REMEDIAL INVESTIGATION REPORT

PACIFIC CITY PARK 600 THIRD AVENUE SOUTHEAST PACIFIC, WASHINGTON



River and Floodplain Management Section Water and Land Resources Division

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PACIFIC CITY PARK 600 THIRD AVENUE SOUTHEAST PACIFIC, WASHINGTON

Prepared for



King County

River and Floodplain Management Section King County Water and Land Resources Division 201 South Jackson Street, Suite 600 Seattle, Washington 98104

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In Conjunction with Aspect Consulting, LLC

January 4, 2019



Prepared for:

King County River and Floodplain Management Section King County Water and Land Resources Division

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Alternate Formats Available.

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CONTENTS

Ab	oreviations		V
Cei	tificate of Lice	ensed Hydrogeologist	vii
Exe	cutive Summa	ary	ix
	Regulatory P	Process	x
1.	Introduction		1
	1.1. General	Site Information	4
	1.2. Report (Organization	4
2.	Site Descript	tion and Background	5
	2.1. Site Hist	tory	5
		2	
	2.3. Site Sett	ting	6
	2.4. Geology	y and Hydrogeology	8
	2.4.1.	Geologic Setting	8
	2.4.2.	Site Geology	11
	2.4.3.	Hydrogeologic Setting	19
	2.4.4.	Site Hydrogeology	19
	2.5. Future S	Site Use	20
	2.6. Ground	water Use	20
3.	Field Investig	gations	25
	3.1. Previous	s Environmental Investigations	25
	3.1.1.	1985 Abandoned Landfill Study	25
	3.1.2.	2010 and 2011 Stormwater Sampling	27
	3.1.3.	2016 Phase II Environmental Site Assessment	27
	3.1.4.	2017 Environmental Investigation	28
	3.2. Site Cha	aracterization	29
	3.2.1.	Contaminants of Potential Concern	29
	3.2.2.	Data Gaps Assessment	33
	3.2.3.	2018 Environmental Investigation	
4.	Conceptual S	Site Model	51
	4.1. Sources	of Chemicals of Potential Concern	51
	4.2. Nature a	and Extent	51

	4.2.1.	Physical Conditions	52
	4.2.2.	Soil	53
	4.2.3.	Groundwater	54
	4.2.4.	Surface Water	57
	4.2.5.	Soil Gas	58
	4.3. Fate and	d Transport	58
	4.4. Prelimin	ary Exposure Assessment	63
5.	Proposed Cle	eanup Standards	65
	5.1. Cleanup	Levels	65
	5.1.1.	Soil	65
	5.1.2.	Groundwater and Surface Water	65
	5.2. Points o	f Compliance	66
	5.2.1.	Soil	66
	5.2.2.	Groundwater	66
	5.3. Terrestri	ial Ecological Evaluation	66
6.	Summary, Co	onclusions, and Recommendations	67
	6.1. Summar	ry	67
	6.2. Conclusi	ions and Recommendations	68
7	References		60

APPENDICES

Appendix A	Soil Boring Logs
Appendix B	Abandoned Landfill Study in King County
Appendix C	Memorandum: Stormwater Quality Sampling Results from the Pacific Right Bank Wetland
Appendix D	Phase II Environmental Site Assessment, Pacific Right Bank Levee Setback Project
Appendix E	Pacific Park/Dumpsite Environmental Investigation Report
Appendix F	Geophysical Investigation Report, Pacific Right Bank Project
Appendix G	Laboratory Analytical Reports
Appendix H	Data Quality Assurance Review Memorandum
Appendix I	Soil Vapor Monitoring Data

TABLES

Table 1.	Summary of Water Level Elevation Data from Monitoring Wells, Pacific City Park Remedial Investigation, Pacific, Washington	73
Table 2.	Summary of Surface Water Sample Results, Pacific City Park Remedial Investigation, Pacific, Washington	75
Table 3.	Summary of Groundwater Sample Results from Push Probes, Pacific City Park Remedial Investigation, Pacific, Washington	77
Table 4.	Proposed Site Screening Levels for Soil, Pacific City Park Remedial Investigation, Pacific, Washington	78
Table 5.	Proposed Site Screening Levels for Groundwater and Surface Water, Pacific City Park Remedial Investigation, Pacific, Washington	81
Table 6.	Summary of Groundwater Sample Results from Monitoring Wells, Pacific City Park Remedial Investigation, Pacific, Washington	85
Table 7.	Summary of Soil Sample Results, Pacific City Park Remedial Investigation, Pacific, Washington	89
Table 8.	Summary of Soil Vapor Monitoring Data, Pacific City Park Remedial	111



FIGURES

Figure 1.	Vicinity Map, Pacific City Park, Pacific, Washington	2
Figure 2.	Site Map, Pacific City Park, Pacific, Washington	3
Figure 3.	Approximate Extent of Historical Dumping and Filling, Pacific City Park, Pacific, Washington.	6
Figure 4.	Wetlands, Stormwater Conveyance Features, and Surface Water Sample Locations, Pacific City Park, Pacific, Washington	9
Figure 5.	Extent of Fill Soil Containing Refuse, and Soil Boring and Monitoring Well Locations, Pacific City Park, Pacific, Washington	13
Figure 6.	Cross Section A-A', Pacific City Park, Pacific, Washington	15
Figure 7.	Cross Section B-B', Pacific City Park, Pacific, Washington	17
Figure 8.	Groundwater Level Contour Map, March 23, 2018, Pacific City Park, Pacific, Washington.	21
Figure 9.	Groundwater Level Contour Map, June 21, 2018, Pacific City Park, Pacific, Washington	22
Figure 10.	Groundwater Level Contour Map, September 26, 2018, Pacific City Park, Pacific, Washington	23
Figure 11.	Soil Vapor Monitoring Locations, Pacific City Park, Pacific, Washington	26
Figure 12.	Extent of Diesel- and Lube-oil Range Petroleum Hydrocarbons in Soil, Pacific City Park, Pacific, Washington	41
Figure 13.	Extent of Lead in Soil, Pacific City Park, Pacific, Washington	43
Figure 14.	Extent of Total Polychlorinated Biphenyls (PCBs) in Soil, Pacific City Park, Pacific, Washington.	45
Figure 15.	Extent of Total Carcinogenic Polycyclic Aromatic Hydrocarbons (cPAHs) in Soil, Pacific City Park, Pacific, Washington	47
Figure 16.	Extent of Soil and Groundwater Contamination Above Site Screening Levels, Pacific City Park, Pacific, Washington	55
Figure 17.	Extent of Total and Dissolved Metals in Groundwater, Pacific City Park, Pacific, Washington	59
Figure 18.	Non-Metals Contaminants of Concern in Groundwater, Pacific City Park,	61

ABBREVIATIONS

Aspect Consulting, LLC

ASTM American Society for Testing and Materials

Bgs below ground surface

BTEX benzene, toluene, ethylbenzene, and xylenes

bTOC below the top of the monitoring well casing

COC contaminant of concern

COPC contaminant of potential concern

cPAHs carcinogenic polycyclic aromatic hydrocarbons

CUL Cleanup level

Ecology Washington Department of Ecology

EPA Environmental Protection Agency

GEM Gas Analyzer & Extraction Monitor

Herrera Environmental Consultants, Inc.

HCID hydrocarbon identification

mg/kg milligrams/kilograms

mg/L milligrams per liter

ml/min milliliters per minute

μg/L micrograms per liter

MCL maximum contaminant level

MTCA Model Toxics Control Act

NWTPH-Dx Northwest total petroleum hydrocarbons, diesel-extended



PAHs polycyclic aromatic hydrocarbons

PCBs polychlorinated biphenyls

RI Remedial Investigation

SSL site screening level

SVOCs semi-volatile organic compounds

S&W Shannon & Wilson, Inc.

TEQ toxic equivalency

TPH total petroleum hydrocarbons

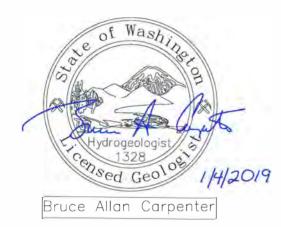
VCP Voluntary Cleanup Program

VOC volatile organic compound

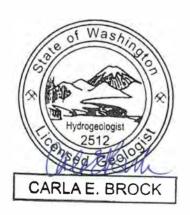
WAC Washington Administrative Code

CERTIFICATE OF LICENSED HYDROGEOLOGIST

This document has been prepared under the supervision of a licensed hydrogeologist.



Bruce Carpenter, LHG	January 4, 2019
Name	Date



Carla Brock, LHG	January 4, 2019	
Name	Date	



EXECUTIVE SUMMARY

This Remedial Investigation (RI) report has been prepared to meet the requirements of the Model Toxics Control Act (MTCA) and regulations implementing it, Washington Administrative Code (WAC) 173-340, and to provide the results of investigations completed to characterize the nature and extent of contamination at the Site. The RI Report has been prepared in general accordance with the Remedial Investigation Checklist Guidance (Ecology 2016).

The Site is located on the existing right bank of the White River in the City of Pacific, Washington, on property that was historically part of the river channel before the construction of a levee and concrete revetment in 1919. The portions of the property located landward of the levee were filled as a King County refuse dump, which was active between approximately 1921 and 1965. The results of investigation activities indicate that portions of the Site were filled with soil and portions of the Site were filled with a combination of soil and refuse. The presence of fill soil alone, in the absence of contaminants of potential concern (COPCs), does not fall under the definition of MTCA as a hazardous substance, and therefore, is not part of the MTCA Site. The Site is defined by any location where refuse is present or where COPCs are present in soil, groundwater, surface water or soil vapor at concentrations exceeding the site screening levels (SSLs). The current Site use includes a seasonal city park and undeveloped natural areas along the White River.

Because of the unknown nature of the fill soil and refuse, initial investigations conducted on the Site included a full analytical suite of COPCs, including total petroleum hydrocarbons (TPH), metals, volatile organic compounds (VOCs), semi-volatile organic compounds (SVOCs), polychlorinated biphenyls (PCBs), herbicides and pesticides. Subsequent phases of investigation focused on characterizing the nature and extent of the primary COPCs, which consist of TPH, metals, PCBs, and total carcinogenic polycyclic aromatic hydrocarbons (cPAHs). The results of investigation work indicate that the fill soil contains ubiquitous concentrations of lead and total cPAHs¹. In areas where refuse was known to have been dumped and/or was observed to be present in subsurface explorations, the results of soil characterization samples identify higher concentrations of total cPAHs and lead, as well as the frequent detection of TPH and PCBs and rare detections of chlorinated VOCs, other SVOCs and pesticides. Laterally, the extent of COPCs in soil are generally defined by the extent of fill. Vertically, the COPCs are present throughout the fill and extend up to 10 feet into the underlying native, alluvial deposits.

The COPCs detected in groundwater include arsenic, lead, total cPAHs and benzene. Arsenic has been detected in groundwater samples collected from across the Site, is naturally occurring in soil and groundwater in Washington state and is likely not present in Site groundwater at concentrations that warrant remedial action. Lead, total cPAHs and benzene have been detected

King County

¹ Total cPAH toxic equivalency (TEQ) concentration calculated in accordance with WAC 173-340-708(8).

above the SSLs in groundwater samples collected to the south-southwest of the refuse, which is hydraulically downgradient based on water level elevations in monitoring wells on the Site. The groundwater data suggest that there may be a low concentration, seasonal and/or diffuse groundwater plume emanating from the refuse or there may be localized groundwater impacts attributable to variations in the fill soil quality. Although groundwater is assumed to discharge to surface water in some areas of the Site, the results of surface water sampling do not indicate the migration of COPCs in groundwater to surface water.

Sufficient information regarding the extent and quality of the fill and refuse, and the associated groundwater impacts, has been collected to allow for the development and evaluation of remedial alternatives, with one exception. Additional investigation is warranted to further evaluate the extent of fill soil and the presence of COPCs in soil and groundwater to the south-southwest of the property boundary. Additional groundwater monitoring wells and soil probes are planned to be installed along the parcel boundary by the apartments immediately west of the Park. A supplemental report to the RI summarizing the results of the additional investigation will be submitted to the Washington State Department of Ecology (Ecology).

REGULATORY PROCESS

The County is seeking an opinion from Ecology through the Voluntary Cleanup Program (VCP) regarding the sufficiency of the RI to meet MTCA requirements. The RI and Ecology's opinion letter will inform the Feasibility Study (FS) which will develop and evaluate cleanup options that will be part of the Pacific Right Bank Flood Protection Project (Project) alternatives to be evaluated in the Draft EIS for the Project.

1. INTRODUCTION

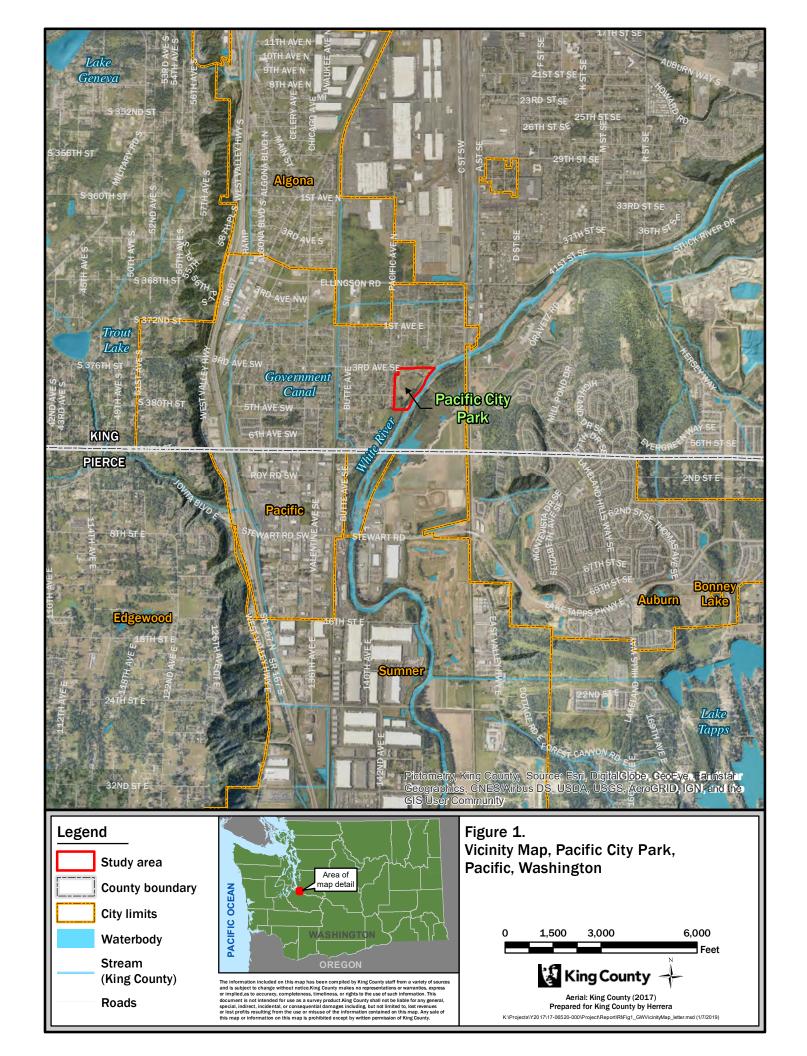
This report presents the results of the RI conducted at the Pacific City Park (Park) located at 600 Third Avenue Southeast in the City of Pacific, Washington (Figure 1). The Park is located on a portion of a 43-acre parcel of land on the right (west) bank of the White River, which is herein referred to as the *Subject Property*, and is currently developed and used as a city park (Figure 2).

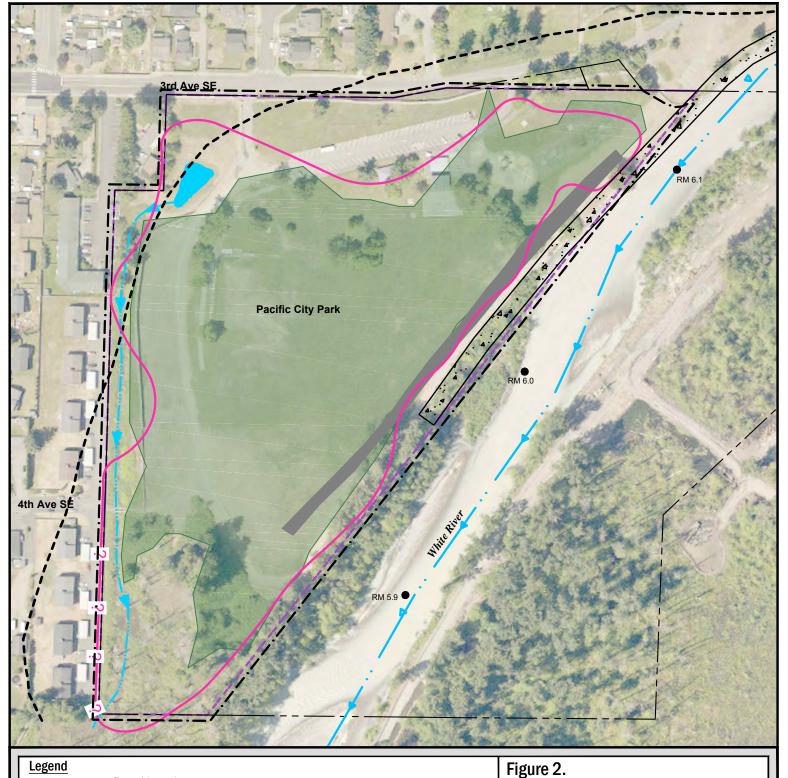
The existing right bank of the White River, which forms the eastern extent of Pacific City Park, consists of a levee and concrete revetment that was constructed in 1919 (S&W 2016). Following construction of the levee, the former river channel located landward of the levee was filled with municipal waste and dredge spoils as an informal dumpsite and city dump until it was closed in 1965. The *study area* for the RI roughly matches the existing Park boundary but extends slightly beyond the Park at the northeast corner. The *Site* depicted on Figure 2 is based on the extent of locations where hazardous substances associated with filling on the landward side of the 1919 levee have been deposited, disposed of, placed, or otherwise come to be located.

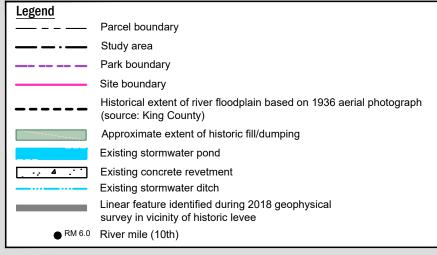
The Site was vacant until 1969 when King County issued a permit to the City of Pacific to use a 21-acre portion of the Subject Property for a park, which opened in 1972 (S&W 2016). In 1990, the City entered into a 30-year lease agreement with King County for continued use of the Subject Property as a park. King County is in the process of evaluating options for increasing flood capacity and reducing flood risk along this section of the river, including consideration of removal of the existing levee and concrete revetment and the construction of a new setback levee.

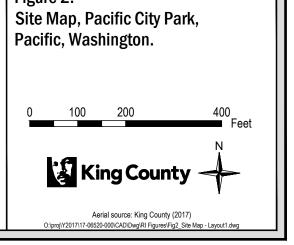
The RI report has been prepared to meet the requirements of MTCA and regulations implementing it: Chapter 173-340 of WAC 173-340. The RI Report has been prepared in general accordance with the Remedial Investigation Checklist Guidance (Ecology 2016).

The purpose of this RI is to collect and evaluate sufficient information to characterize the nature and extent of contamination at the Site to enable the development and selection of a remedial alternative. King County entered the Voluntary Cleanup Program (VCP) in July 2018 and is requesting an opinion from Ecology on the RI report as it moves forward with developing project alternatives that will be analyzed in an Environmental Impact Statement (EIS).









1.1. GENERAL SITE INFORMATION

Site Name	Pacific City Park
Site Address	600 Third Avenue Southeast
Facility Site Identification number (FSID)	2160
King County Assessor's Parcel No.	3621049040
Cleanup Site ID	21
VCP Project Number	TBD
Ecology Site Manager	TBD

The project consultant is Mark Ewbank with Herrera Environmental Consultants, Inc. located at 2200 Sixth Avenue, Suite 1100, Seattle, Washington 98121. Telephone: 206-787-8217, and email: MEwbank@herrerainc.com.

The Subject Property is owned by King County, with a portion leased to the City of Pacific for use as a city park. Chris Brummer, Supervising Engineer with the River and Floodplain Management Section of the King County Water and Land Resources Division, is the Site contact. His office is located at 201 South Jackson Street, Suite 600, Seattle, Washington, 98104. Telephone: 206-477-4655, and email: Chris.Brummer@kingcounty.gov.

1.2. REPORT ORGANIZATION

This RI report has been organized in accordance with Ecology's RI Checklist (Publications No. 16-09-006) dated May 2016 and includes the following:

- Section 2 provides a definition of the Site and property and presents a summary of the background information including the environmental setting, historical use of the property/vicinity, and regulatory involvement.
- Section 3 provides the scope of work and results of the RI, including a summary of the historical environmental studies/actions and screening/cleanup levels used to evaluate the soil, groundwater and soil vapor data collected for the RI to facilitate site characterization.
- Section 4 presents the Conceptual Site Model (CSM) for the Site, including the sources and nature and extent of concentrations of hazardous substances in soil and groundwater at the Site, and a preliminary assessment of potential receptors and exposure pathways.
- Section 5 presents the proposed cleanup standards for future cleanup at the Site, including cleanup levels and points of compliance for soil and groundwater.
- Section 6 presents the summary, conclusions and recommendations.



2. SITE DESCRIPTION AND BACKGROUND

The Site is located approximately 2,000 feet north of the King/Pierce county boundary line (Figure 1). The property is relatively flat, with ground surface elevations ranging from 80 to 87 feet (NAVD88).

Access to the Site is restricted from October through March by a 4- to 8-foot-high temporary flood barrier system (HESCOs) installed in 2009 and extended in 2013. The continuous barrier extends along the northern and western property boundaries, except for sections of the barrier that are removed from April to September to allow public access at two locations on the northern portion of the property, along 3rd Avenue SE.

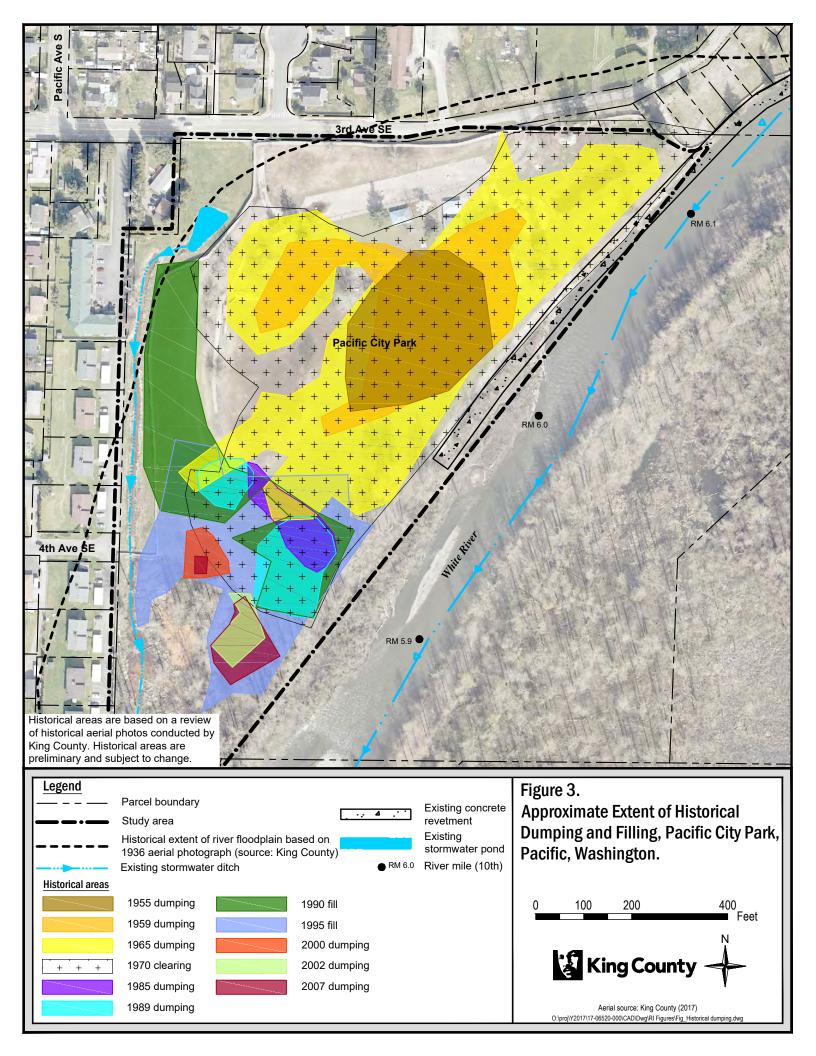
2.1. SITE HISTORY

The Site is located in an area that historically had been occupied by the White River prior to construction of levees in the 1910s and filling of the site from the 1920s through the 1960s. The Site is located in an area that has historically been agricultural and is prone to flooding by the White River. According to historical photos and maps, flood-control measures were taken that included construction of a concrete revetment in 1919 as part of the channelization of the White River along the alignment of the former Stuck River (S&W 2016). The extent of the concrete revetment to the southwest is uncertain; therefore, Figure 2 shows only the portion of the revetment within the project area that is known.

Figures 2 and 3 depict the approximate extent of historical dumping and filling at the Site. Aerial photographs from 1936 and 1944 indicate an orchard occupied the area west of the Site. The Site was closed for use as a City dump in 1965 and was abandoned until 1969 when King County issued a permit to the City of Pacific for a city park that subsequently opened in 1972. Fill soil containing refuse placed in the 1950s and 1960s was covered with additional fill soil as the Park was developed but dumping continued in the southwest portion of the site through the 2000s. In the 1980s, development of the area surrounding the park included placement of fill within the area underlying the four apartments south of 4th Avenue SE and the southeastern half of the area underlying the four apartments north of 4th Avenue SE (S&W 2016). The exact lateral extent of dumping and fill in the southwest corner of the Site is unknown.

In 2009, after significant flooding, the County installed temporary HESCO flood barriers along the right bank of the White River and around the north and west sides of the Park. Additional HESCO barriers were installed in 2013, which extended the barrier upstream of the park (to the BNSF Railway) and added a second tier of HESCOs on the north and west sides of the park.





2.2. SITE USE

The Subject Property is currently used as a public park operated by the City of Pacific under a 30-year lease agreement from King County. Since 2013, the park is accessible April through October, and restricted for use during the winter months when HESCO barriers are fully continuous. The park is relatively flat and includes paved parking, a restroom, picnic shelter, ball field, basketball court, children's play area, and open space along the river. The site is used primarily for passive recreation (including walking, dog walking, visiting the river bank), picnicking, grilling, children's activities in the play area, and an annual city festival (Pacific Days).

2.3. SITE SETTING

The Site is located in the White River valley in south King County and north Pierce County (Figure 1). The White River travels 68 miles from the western foothills of the Cascade Mountains in a general westward flow direction before turning south in the City of Auburn and draining into the Puyallup River near the City of Sumner. A 1914 survey of the Site and adjacent areas shows the former river channel flowing through the approximate central portion of the Site now occupied by the Park (Roberts, W.J. 1920) (Figure 3). The United States Geological Survey (USGS) operates two surface water monitoring stations on the White River in the vicinity of the Site: one located upstream of the Site near A Street (White River near Auburn, WA), and one located approximately 0.2-miles downstream of the Site (White River at Pacific, WA). A review of water level data recorded by the USGS in the White River between October 2012 and the present indicates that surface water elevations are typically highest in the spring and lowest in the late summer and early fall (USGS, 2018). The mean monthly river gage heights in the White River range from 81 to 84 feet NGVD 29 at the A Street gage and from 72 to 74 feet NGVD 29 at the Pacific gage (USGS, 2018).

The White River is classified as a R2UBH wetland (Riverine, Lower Perennial, Unconsolidated Bottom, Permanently Flooded) by the US Fish & Wildlife Service. This classification includes all wetlands and deepwater habitats contained within a river channel, characterized by a low gradient, no tidal influence and water flow all year, at least 25% of the river bottom is covered with particles smaller than stones and a vegetative cover of less than 30%. The White River is also a Class 1 wetland per the 1990 King County Wetlands Inventory (King County, 1991).

Land use surrounding the site includes residential developments adjacent to the north and west, and open space owned by King County to the south and east (KCGIS 2018). Portions of the levee on the left bank of the White River, across the river from the site, were removed as part of the Countyline Levee Setback Project completed by the King County Flood Control District in 2017.

Herrera wetland biologists visited the study area in May, June, July, and October 2018 to conduct wetland delineation work. This delineation work identified six wetlands (Wetlands A through F) located on the Site (see Figure 4), and a seventh wetland (Wetland G) located further



to the southwest in an area not covered by this RI. Wetland A and B are riverine wetlands; Wetlands C, D, and E are depressional wetlands within the active area of Pacific City Park; and Wetland F is a depressional wetland south of the onsite stormwater pond.

A stormwater ditch along the western edge of the site drains south for approximately 2,000 feet before discharging into Wetland A, which drains into the river. The ditch receives stormwater via a series of catch basins located along 3rd Avenue SE, the parking lot in the Park, and the apartments west of the ditch. The water level in the ditch roughly corresponds to groundwater levels in adjacent wells, so the ditch presumably intercepts groundwater that intermingles with surface stormwater conveyed by the ditch.

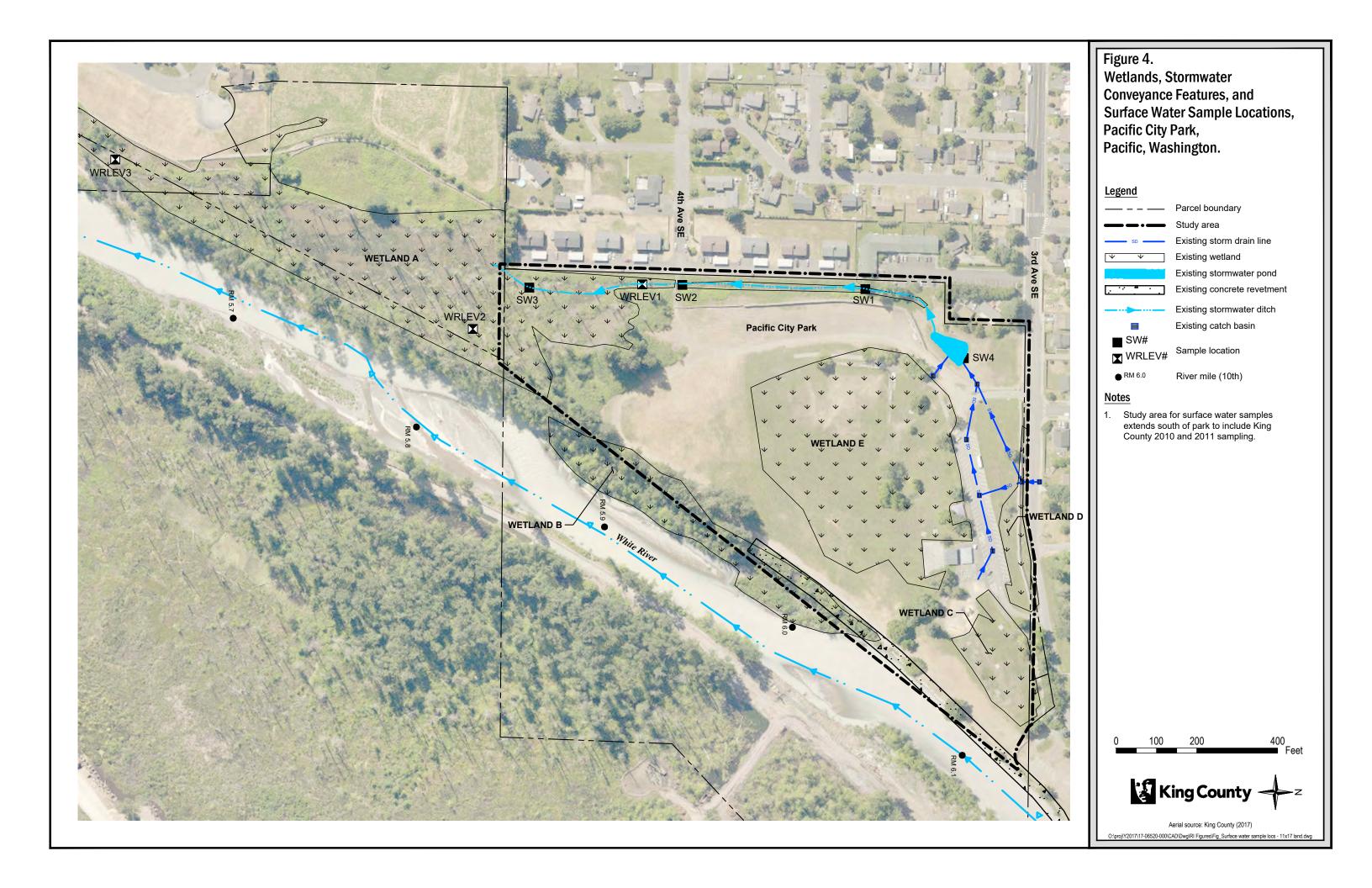
2.4. GEOLOGY AND HYDROGEOLOGY

2.4.1. Geologic Setting

Geological mapping in the project area identifies a thick sequence of post-Fraser-glaciation alluvium deposited over Fraser glacial and pre-Fraser glacial and nonglacial sediments (Mullineaux, 1965). The White River valley wall and valley floor are composed of glacially consolidated sediments from multiple Pleistocene glaciations and interglacial deposition during the Pleistocene (2.6 million to 11,600 years ago). The latest postglacial and Holocene (11,600 years ago to present) geologic processes included erosion and transport of sediment from uplands and drainage headwaters, and transport of sediment from lahars originating on Mount Rainier (Collins and Montgomery, 2010; Zehfuss et. al, 2003). The deep glacial meltwater channel below the project area was subsequently infilled with Holocene (Recent) alluvium. The Holocene alluvium deposited along the White River merged into a continuous alluvial deposit through the regional lowland that now contains (from Seattle southward to Tacoma) the lower reaches of the Duwamish, Green, White, and Puyallup rivers.

The geologic history of the White River valley generally consists of the following processes:

Subglacial Erosion of the White-Green-Duwamish River Valley. Erosion by ice from
the Vashon ice sheet and subglacial meltwater flow created the broad north-to-south
trending marine embayment now occupied by the White, Green, and Duwamish rivers.
Glacially overridden soils and bedrock lie at considerable depth below younger and
weaker, late-glacial and postglacial deposits.



- Postglacial Incision and Sedimentation. Shortly after deglaciation, sea level was several hundred feet lower than at present, and the Puget Lowland was depressed by the weight of the glacial ice. Rapid isostatic rebound (causing uplift of land) combined with low post-glacial sea level allowed the lowland rivers, including the White River, to incise to the lower base level. Several hundred feet of rising sea level continued until about 5,000 years ago, partially resulting from combined effects of melting ice caps and local isostatic rebound following deglaciation. The rise in sea level resulted in marine flooding of the Duwamish, White and Green river valleys. The oldest post-glacial deposits in the valleys include marine clays and muds.
- **Embayment and Valley Filling.** The marine embayment filled with Holocene alluvium derived primarily from multiple lahars (volcanic mudflows) originating on Mount Rainier. The most recent large lahar that impacted the valley, known as the Osceola Mudflow, travelled down the White River valley from Mount Rainier approximately 5,600 years ago. White River deposition in the Puget Lowland has been fan-dominated between Puyallup and Kent. Sedimentation of the White River fan was driven largely by large-scale erosion of postglacial and lahar deposits as the White River carved the confined valley upriver of the marine embayment.

After the initial pulse of diamictic (unsorted) debris from the Osceola Mudflow, sedimentation around the project area occurred as hyperconcentrated flow deposits (sediment rich but water transported and reworked deposits), deltaic deposits, and alluvial deposits. Deposition occurred in several discrete phases, and these deposits are described as tightly packed sand and gravel of volcanic origin with a distinct dark purple-gray color. Other post-Osceola mudflow deposits including hyperconcentrated sands have also been identified within younger Holocene alluvial strata (Zehfuss et al., 2003) in the Lowland.

• Historic Filling and Hydraulic Engineering. Historically, the White River flowed north across its alluvial fan through the present location of Auburn to join the Green River. The Stuck River, which is occupied by the present-day White River, was a distributary of the White River that flowed into the Puyallup River and carried variable amounts of White River flow. Early twentieth century flooding and hydraulic engineering revised flow paths and outlet points of the White and Green rivers, resulting in the system present today where the entire flow of the White River discharges south to the Puyallup River.

2.4.2. Site Geology

Explorations were completed at the Site by Shannon & Wilson in 2015 and 2016 under contract with King County, and additional explorations were completed by Herrera in 2017 under separate contract. Soil borings were advanced using direct push methods to investigate soil and groundwater quality. In addition, geotechnical explorations, including test pits and soil borings, were completed by Shannon & Wilson in 2016 and by Aspect in 2018. The boring logs and test pit logs from these subsurface investigations are included in Appendix A and D, respectively. The geologic conditions at the Site, as observed during these investigation activities, are described below.



Fill Soil

The approximate lateral extent of historical dumping and filling at Pacific City Park was initially estimated based on a review of historic aerial photographs and is depicted on Figure 3 (S&W 2016). Figure 5 depicts the lateral extent of fill soil that contains refuse (i.e., debris) based on explorations conducted in 2016, 2017, and 2018, along with the locations of soil borings and monitoring wells completed at the Site. Cross sections depicting the extent of fill soil are presented in Figures 6 and 7.

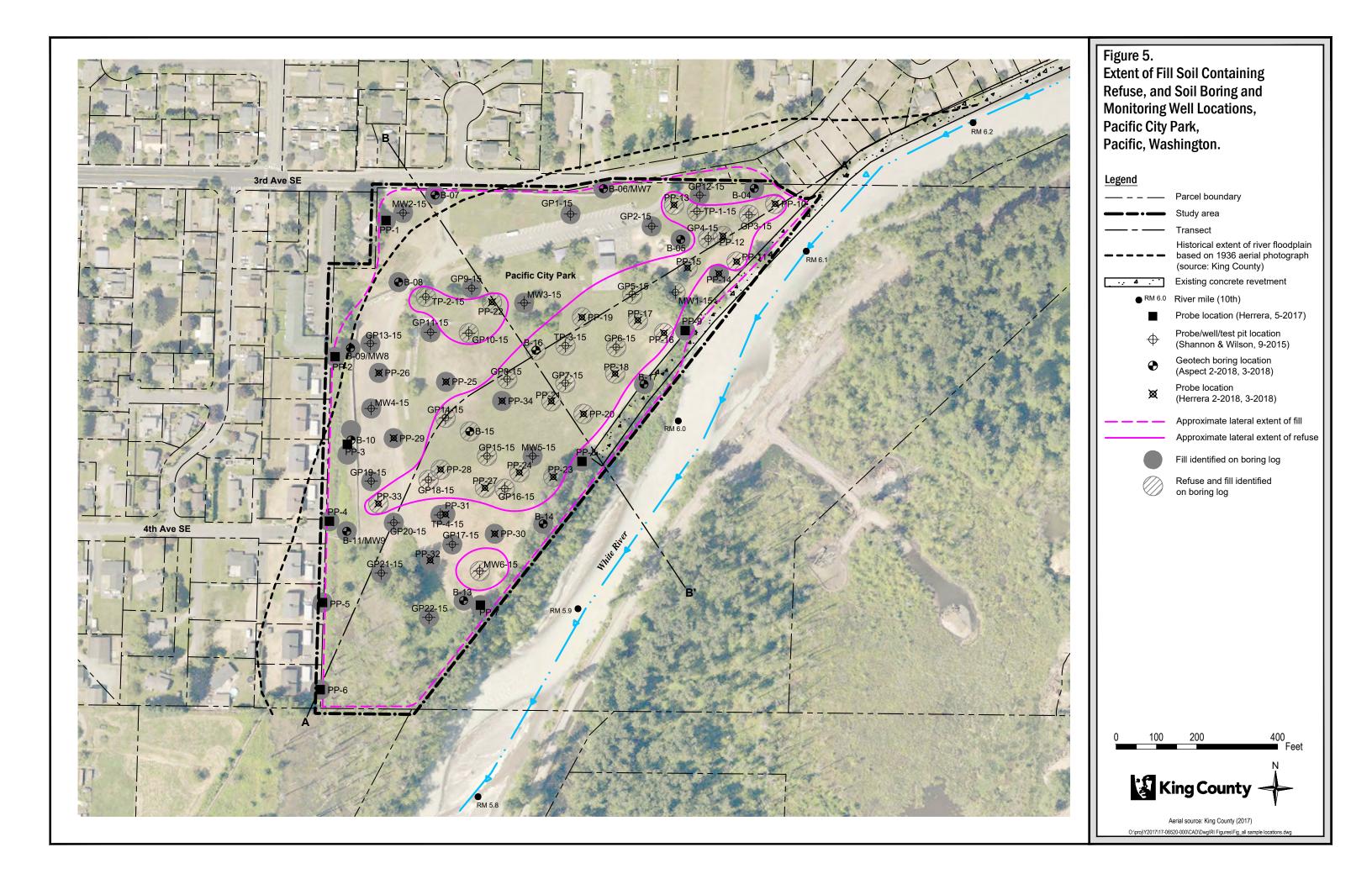
Fill soil containing refuse was encountered at depths ranging from 2 to 20 feet below ground surface (bgs). The refuse includes glass, brick, cement, organics, wood, paper, rubber and ceramics. The refuse is mixed with fill soil consisting of variable amounts of silt, sand, and gravel. The fill soil with refuse is typically thicker in the central/south-central portion of the Site. The fill soil with refuse is covered with approximately 2 to 10 feet of sandy, gravelly fill soil that was reportedly placed following closure of the dumpsite to cover the debris and create a level surface (S&W, 2016). Groundwater levels within the fill soil were observed at depths of approximately 5 feet bgs. Holocene alluvium, described further below, was observed below the fill soil with refuse.

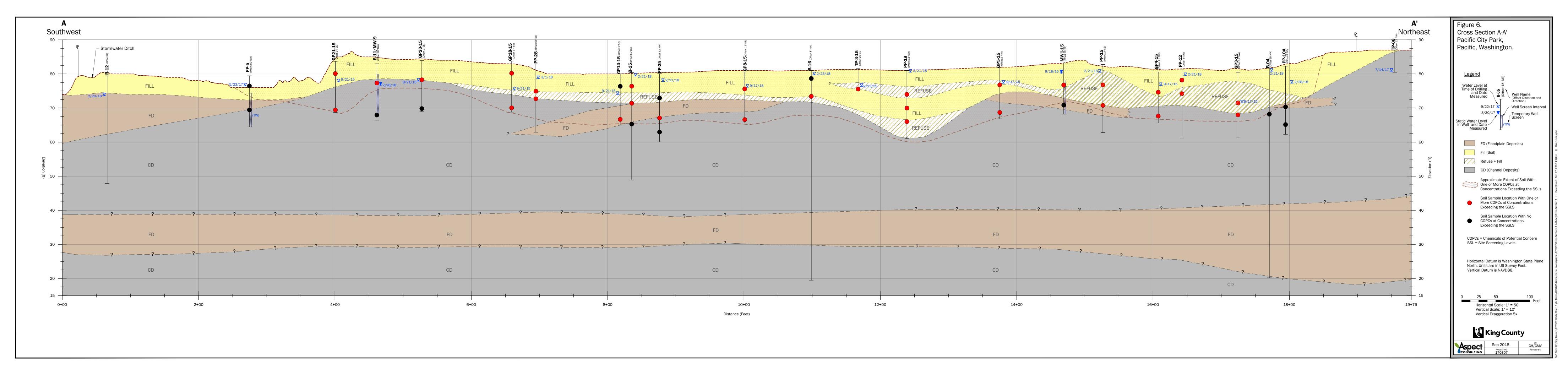
Holocene Alluvium

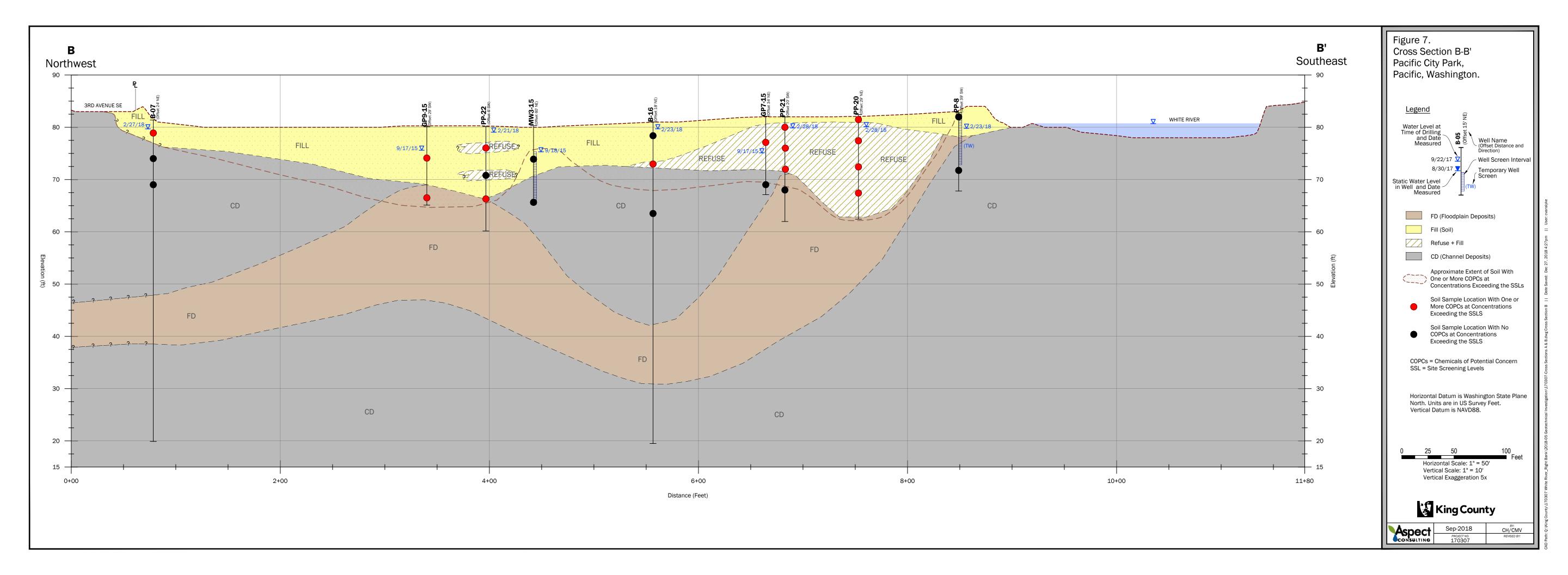
Alluvium consists of floodplain and channel deposits. Floodplain deposits are described as relatively fine-grained soils ranging from silt and clay to silty fine to medium sand. Channel deposits are described as relatively clean sand with gravel, cobbles, and boulders with interbedded fine-grained sediments. The floodplain and channel deposits are not necessarily predictable using a continuous layered stratigraphy model due to past channel meandering through the area and infill processes associated with over-bank flooding events.

Floodplain Deposits. During flood events, suspended-load sediment composed of sand, silt, and clay are deposited on floodplains peripheral to the active river channel. The deposits typically include relatively fine-grained soils ranging from silt and clay to silty fine to medium sand and are typically cross-bedded to laminated, although bedding may be indistinct. These deposits may also contain well-preserved logs, in situ stumps, and other organics and woody debris.

Floodplain deposits are encountered beneath the fill soil or a thin layer of topsoil in all borings, and as deep as 51 feet bgs in boring B-16 (Figure 5). Floodplain deposits are identified based on grain size, fines content, interbedded texture, and density. Floodplain deposit soils consist of moist to wet, brown or gray, sandy, medium plasticity CLAY (CL), low plasticity SILT (ML), fibrous or fine-grained PEAT (PT), and relatively clean sand (containing less than 5 percent silt) to silty sand (SM, SP-SM, SP). The sand size fraction is typically fine to medium and rarely coarse, and soil types are commonly interbedded. Standard Penetration Test (SPT) N-values ranged from 1 to 31 blows per foot (bpf) with an average N-value of 18 bpf, indicating the floodplain deposits were typically stiff/medium dense (Aspect, 2018).







Channel Deposits. Channel deposits are the result of depositional sequences within the
active or historical river channel(s). These deposits typically include relatively clean sand
with gravel, cobbles, and boulders. Interbeds of silt, clay, and peat can be present within
channel deposits, particularly in areas of historical oxbows or other low-energy channel
features. Well-preserved logs and wood debris may be present in the channel deposits.

Channel deposits are encountered both above and beneath the floodplain deposits, and as deep as 61.5 feet bgs in borings B-04, B-07, B-10, B-13, and B-16. Channel deposits are identified based on grain size, fines content, and color, which was commonly a dark purple-gray. Channel deposits generally consist of gray or dark gray, clean to silty SAND (SP/SW, SP-SM/ SW-SM, SM) or clean to sandy GRAVEL (GP). SPT N-values ranged from 2 to greater than 50 bpf, with an average N-value of around 30 bpf, indicating the channel deposits were variable and ranged from loose to very dense and were typically medium dense. Some N-values were overstated due to the presence of coarse gravel. Channel deposit density generally increased with depth.

2.4.3. Hydrogeologic Setting

Groundwater within the White River valley occurs primarily within alluvial deposits and glacial outwash deposits but also may occur in coarse-grained interglacial deposits (Welch et. al., 2015). The upper alluvial aquifer, primarily comprised of alluvial silt, sand and gravel deposits, is present throughout the White River valley and is estimated to range from 46 to 107 feet thick in the Site vicinity (Welch et. al., 2015). Groundwater in the upper alluvial aquifer is generally unconfined. The upper alluvial aquifer is separated from a lower alluvial aquifer by a confining unit comprised of volcanic mudflow-lahar deposits that range from 27 to 52 feet thick in the Site vicinity (Welch et. al., 2015). The horizontal groundwater flow direction within the upper alluvial aquifer is generally in the down-valley direction of the White River. Water level differences between the upper and lower alluvial aquifers indicate the potential for upward groundwater flow in the Site vicinity (Welch et. al, 2015). During late summer and early fall, when contribution from melting snow and glacial ice are lowest, it is likely that baseflow is the main source of water to the White River in the Site vicinity.

2.4.4. Site Hydrogeology

The groundwater levels in nine monitoring wells located at the Site were measured during two sampling events in 2017 and 2018 to range from near the ground surface (0.32 feet below the top of the monitoring well casing [bTOC]) to 6.7 feet bTOC (Table 1). King County has collected continuous groundwater level data in the wells since October 2015. Based on these data and river stage measured at the A Street gage, the general direction of groundwater flow at the Site is inferred to be to the west and southwest and fluctuates with river stage. This indicates that the White River is "losing" water to the upper alluvial aquifer near the Site. Refer to groundwater level contour maps presented in Figures 8, 9, and 10.

Based on slug tests performed during Aspect's 2018 geotechnical investigation, calculated hydraulic conductivities at the Site ranged from 10.4 feet/day to 103.4 feet/per day, resulting in a mean of 42.5 feet per day (Aspect, 2018). Lower hydraulic conductivity (9.8 feet/day to 11.6 feet/day) was observed in monitoring wells that are fully screened within the upper alluvial aquifer than those that are partially screened within the fill soil.

2.5. FUTURE SITE USE

The planned future uses of the Site may include flood protection, restored aquatic and terrestrial habitat, and a public park.

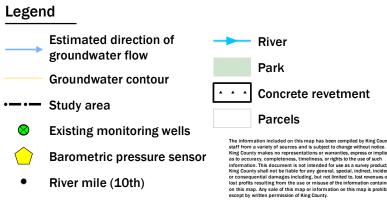
2.6. GROUNDWATER USE

Groundwater at the Site is currently not utilized for potable or non-potable purposes.

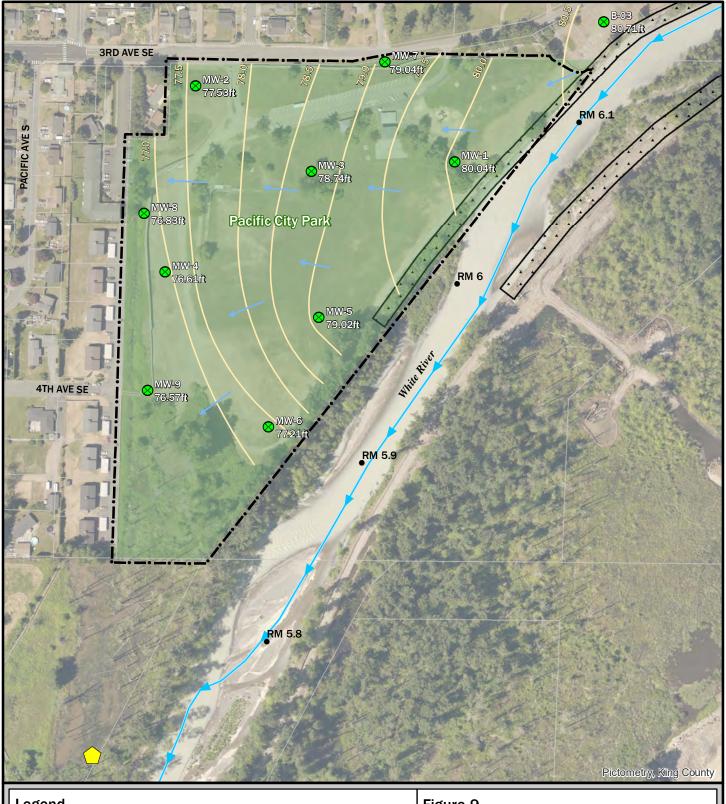
According to the Washington State Department of Health (DOH) Source Water Assessment Program (SWAP) mapping application, the City of Pacific supplies water to customers, including landowners adjacent to the Site and the Site itself, from three municipal water supply wells in a well field approximately 0.6-miles north of the Site (DOH, 2018). The Site is located beyond the 1-year travel time and within the 5-year travel time of the City's water supply wells (DOH, 2018). The Site is not located within the 10-year travel time of any other drinking water wells (DOH, 2018).

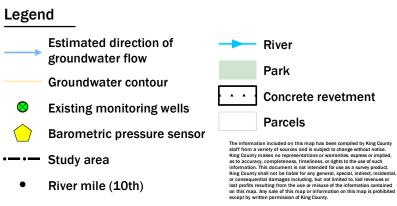
One drinking water well is identified approximately 0.35-miles east-northeast of the Site, on the opposite side of the White River (DOH, 2018). This well is reportedly owned by Danner Corporation, is listed as active, 260 feet deep, with a pumping capacity of approximately 30 gallons per minute (DOH, 2018). The water well is located upgradient of the Site and based on its depth, it is unlikely that it well is hydraulically connected to the upper alluvial aquifer at the Site. No other water supply wells were identified within one mile of the Site (DOH, 2018).













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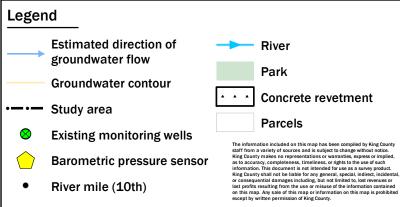
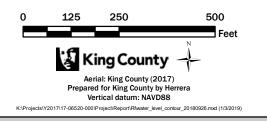


Figure 10. Groundwater Level Contour Map, September 26, 2018, Pacific City Park, Pacific, Washington.



3. FIELD INVESTIGATIONS

3.1. Previous Environmental Investigations

This section describes previous environmental investigations, including methods and results, conducted at the site from 1985 through 2017. The summary of the previous environmental investigations provided in this section compares results to the regulatory criteria that were used to evaluate data at the time of the investigation. Because an evaluation of exposure pathways and applicable screening levels had not yet been completed, the regulatory criteria consisted of the standard MTCA Method A and/or B cleanup levels for soil and groundwater. Site Screening Levels (SSLs) are developed for the RI in Section 3.2 and a comprehensive discussion of the results, for both the previous and current investigations is presented in the Conceptual Site Model, Section 4.

3.1.1. 1985 Abandoned Landfill Study

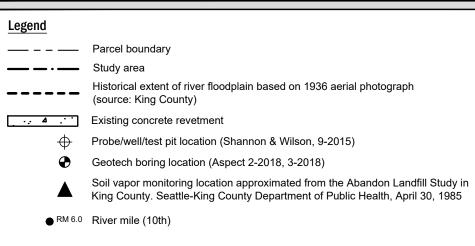
An Abandoned Landfill Study was conducted by the Seattle King County Health Department in 1985 on 20 sites in King County to determine if there were any public health problems associated with these sites (Seattle King County Department of Health, 1985) (Appendix B). The study involved research of geographical and historical data to help guide sampling efforts at the abandoned dumpsites. The sampling efforts consisted of a survey screening of methane and non-specific organics/inorganics and water quality parameters of conductivity, dissolved oxygen, pH, turbidity, and temperature. In the event the initial screening survey revealed any environmental issues, funds would be allocated for additional work.

In October 1984, fifteen boreholes were advanced at the Pacific City Park, and methane and non-specific trace gas concentrations (i.e., organic or inorganic gases with an ionization potential <10 electron volts detectable by photo-ionization detector) were recorded. Low concentrations of methane up to 0.4 percent and low concentrations of non-specific trace gas levels were detected with the exception of 6.2 parts per million detected at one location (Seattle King County Department of Health, 1985). The vapor monitoring data are presented in Table 6 and the monitoring locations are depicted on Figure 11.

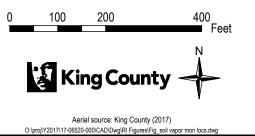
Water quality measurements were recorded at a sample location along the southeast perimeter of the site along the right bank of the White River. A water sample tested for water quality parameters was determined to not be indicative of leachate (Seattle-King County Department of Public Health. 1985). The surface water measurements and laboratory analytical results are summarized in Table 2 and the sampling locations are depicted on Figure 4.











3.1.2. 2010 and **2011** Stormwater Sampling

In 2010 and 2011, King County (2015) collected stormwater samples from the ditch near the end of 4th Avenue SE and from two locations in the wetland located 300 feet and 1,000 feet south of 4th Avenue SE (Appendix C). Water was tested for 6020A metals, pesticides, herbicides, and EPA 8260C VOCs and EPA 8270D SVOCs (2015). Results from all three locations found detectable concentrations of VOCs in surface waters. The laboratory analytical results are summarized in Table 2 and the sampling locations are depicted on Figure 4.

3.1.3. 2016 Phase II Environmental Site Assessment

Beginning in 2015, a Phase II Environmental Site Assessment (ESA) was conducted to evaluate conditions across the Site (S&W 2016) (see Appendix D). Soil and groundwater samples were collected from 28 push-probes advanced to a maximum depth of approximately 15 feet bgs, and six monitoring wells constructed in six of the probe explorations. Soil samples were also analyzed from two of four test pits excavated to depths ranging from 6 to 7 feet bgs (Figure 5).

Soil samples were analyzed for gasoline-, diesel-, and lube oil-range petroleum hydrocarbons; VOCs; RCRA 8 metals; SVOCs, including polycyclic aromatic hydrocarbons (PAHs); herbicides; and pesticides. Groundwater samples were analyzed for gasoline-, diesel-, and lube oil-range petroleum hydrocarbons; PAHs; VOCs; and total and dissolved priority pollutant metals. Tables 6 and 7 summarize soil and groundwater sampling data from the investigation.

The concentrations of contaminants detected in soil and groundwater relative to standard MTCA Method A and/or B cleanup levels, are summarized as follows:

Soil

- Lube oil-range petroleum hydrocarbons were detected in two soil samples at depths of 4.5 and 5.5 feet bgs, above the MTCA Method A cleanup level.
- Three arsenic samples, five cadmium samples, seven lead samples, and two mercury samples were detected above the MTCA Method A cleanup levels, at depths ranging from 4.5 to 12.5 feet bgs.
- No VOCs were detected above the MTCA Method A cleanup levels.
- SVOCs, including total cPAHs² were detected above the MTCA Method A cleanup level at 18 soil sample locations.
- No PCBs were detected above the MTCA cleanup levels.

King County

² Total cPAH toxic equivalency (TEQ) concentration calculated in accordance with WAC 173-340-708(8).

 Low concentrations of herbicides and pesticides were detected in 9 samples, and 3 samples, respectively during the 2016 Phase II ESA, but at several orders of magnitude below MTCA Method B cleanup levels (S&W 2016).

Groundwater

- Total arsenic concentrations exceeded the MTCA Method A cleanup level at eight locations. Dissolved arsenic exceeded the MTCA Method A cleanup level at two locations.
- Total lead concentrations exceeded the MTCA Method A cleanup level at three locations. Dissolved lead was not detected above the MTCA Method A cleanup level.
- No other metals, VOCs or petroleum hydrocarbons were detected above the MTCA cleanup levels.

3.1.4. 2017 Environmental Investigation

In 2017, Herrera advanced nine borings as part of an Environmental Investigation that included soil and groundwater sampling at the park boundary, outside of the fill area (Figure 5) (Herrera 2017) (see Appendix E).

The purpose of the 2017 Environmental Investigation was to determine if soil and/or groundwater contamination was present outside of the fill area and to determine the potential for offsite impacts. Herrera sampled the six existing monitoring wells located on the Site on May 15, 2017 and collected soil and groundwater samples from nine push-probe borings completed at the perimeter of the Site on May 23 and 24, 2017. Six push-probe explorations were completed along the western perimeter of the Site and three along the eastern perimeter.

Shallow and deep soil samples were collected at depth intervals of 0 to 5 feet and 10 to 15 feet in each of the nine borings. The samples were analyzed for gasoline and diesel-range total petroleum hydrocarbons; VOCs; SVOCs, including PAHs; and total and dissolved metals (Table 2). Six soil samples also were analyzed for PCBs based on the detection of lube oil. Groundwater samples were also collected from six existing monitoring wells and each of the nine probe borings. The samples were analyzed for gasoline and diesel-range total petroleum hydrocarbons; VOCs; PAHs; and total and dissolved metals (see Tables 3, 6 and 7).

The concentrations of contaminants detected in soil and groundwater relative to standard MTCA Method A and/or B cleanup levels, are summarized as follows:

- Soil
 - None of the soil concentrations reported for samples collected from nine boring locations exceeded the MTCA cleanup levels for gasoline- and diesel-range total petroleum hydrocarbons, VOCs, SVOCs, total and dissolved metals, or PCBs.



Groundwater

- Benzene and other VOC components of gasoline were detected in one groundwater sample collected from boring PP5, which also had a reported gasoline concentration. The benzene concentration slightly exceeded the MTCA Method A cleanup level. The presence of benzene and gasoline at PP-5 was likely attributable to an isolated spill not associated with the dumpsite, since no gasoline was detected in six monitoring wells or in water samples collected from probes GP-12-15, GP-13-15, and GP-21-15 by S&W in 2015.
- o Concentrations of total PAHs were below the MTCA Method A cleanup level.
- Multiple MTCA exceedances for total metals in groundwater were attributed to extremely turbid samples collected from probe borings. In addition to sample turbidity, elevated arsenic concentrations in groundwater may be due to the release of arsenic derived from emissions associated with historical coal burning that was sequestered in peat (and other organic material), potential uses of arsenic at the nearby orchard, and/or from being located within the plume of the former Asarco lead and copper smelter in Tacoma.
- Arsenic was the only metal detected in groundwater monitoring wells (and only in MW4-15) that exceeded the MTCA Method A cleanup level. Arsenic also was reported above the MTCA cleanup level in two wells (MW2-15 and MW4-15) during the October 2015 groundwater sampling event.

3.2. SITE CHARACTERIZATION

The following sections describe COPCs, SSLs, potential exposure pathways, data gaps, and site characterization activities for each site media including soil, groundwater, surface water, and soil vapor.

3.2.1. Contaminants of Potential Concern

The COPCs for the Site are those that could be potentially associated with, or attributable to, the historical activities conducted at the Site. The COPCs were initially identified based on historical use of the Site as a refuse dump site and the use of adjacent properties as orchards and farmland. In addition, historical photos suggest that treated wood may have been used in bank stabilization (S&W, 2016). Based on these activities, the COPCs include the following contaminant groups:

- TPH in the gasoline-, diesel-, and oil-ranges
- VOCs
- SVOCs, including PAHs

- Metals
- PCBs
- Herbicides and pesticides.

The results of previous investigations and the site characterization have been collectively evaluated to identify the COPCs that are present at the Site. The Site COPCs are chemicals that are reported above the laboratory reporting limits in any media sampled at the Site (soil, groundwater, surface water, and soil gas). SSLs have been developed to evaluate the chemical analytical results for the COPCs as part of this RI. The COPCs and their SSLs are presented on Tables 4 and 5, for soil and groundwater respectively. Because the development of the groundwater SSLs includes consideration of applicable surface water criteria, the SSLs for groundwater and surface water are the same (Table 5).

The development of SSLs is dependent on potential exposure pathways and receptors. The potential exposure pathways and receptors are summarized in Section 3.2.1.1 as the basis for the selection of SSLs in Section 3.2.1.2. Section 3.2.1.3 presents the RI data gaps that were addressed as part of the site characterization and the 2018 environmental exploration activities are summarized in Section 3.2.2.

The results of the site characterization are summarized in Section 3.3.2. The results of the previous investigations and the site characterization together constitute the RI and are the basis for development of the CSM (Section 4).

3.2.1.1. Potential Exposure Pathways

The development of SSLs relies on the identification of current and potential future exposure pathways and receptors. For purposes of developing SSLs, we acknowledge that there is a potential for contaminant migration and groundwater discharge to the river, so we have included consideration of regulatory criteria that are protective of surface water receptors. Because of the site setting, we have also included consideration of regulatory criteria that are protective of ecological receptors. The following exposure pathways and receptors are applicable.

- **Soil/refuse leaching to groundwater:** Contaminants in soil and/or refuse can leach to groundwater by infiltration of precipitation through contaminated soil and refuse or where groundwater is in contact with contaminated soil or refuse.
- **Ingestion of groundwater:** Human receptors have the potential to contact contaminants in groundwater via ingestion.
- Direct contact with soil, refuse, groundwater and surface water: Human, ecological, and aquatic receptors have the potential to contact contaminants in soil, refuse, groundwater, and groundwater that discharges as surface water under current exposure scenarios.

• **Ingestion of surface water and aquatic organisms:** Human receptors have the potential to ingest surface water and aquatic organisms that come into contact with contaminants in groundwater that discharges to surface water and sediment.

SSLs were established for the Site based on these potential exposure scenarios and are described in Section 3.2.3. The potential exposure pathways are evaluated in Section 4 (Conceptual Site Model).

Exposure pathways determined not applicable at the Site include:

• Soil vapor discharge to ambient and indoor air: Landfill gas (i.e., soil vapor) monitoring conducted during the 1985 Abandoned Landfill Study detected low concentrations of methane in boreholes at the site (Figure 11), and data in Appendix I with the highest being 0.4 percent (S&W 2016). Landfill gas monitoring was conducted March 23 and June 21, 2018, at monitoring wells MW-6 and MW-9; and September 23, 2018, at monitoring wells MW-1, MW-6, and MW-9 (Figure 11) and methane was not detected. No VOCs were detected in soil samples above MTCA Method A cleanup levels during the 2016 Phase II ESA (S&W 2016). The sampling and monitoring conducted to date does not indicate a potential exposure pathway from soil vapor to ambient and indoor air.

3.2.1.2. Site Screening Levels

The current and potential future exposure pathways and receptor described above were considered when developing the SSLs for the RI. Data collected during previous investigations and the RI were then evaluated against SSLs to assess the nature and extent of contamination at the Site. The SSLs are intentionally conservative and represent the most stringent of relevant and appropriate criteria for all potential exposure pathways. The SSLs are not cleanup levels. Sitespecific cleanup levels should be developed as part of the Feasibility Study.

3.2.1.2.1. Soil

The SSLs for soil include consideration of the following:

- Standard MTCA Method A and B cleanup levels for the protection of human health for unrestricted land use, through direct contact only
- Calculated MTCA Method B cleanup levels for the protection of groundwater, direct contact and ingestion exposure (saturated soil)
- Ecological Indicator Soil Concentrations for the protection of plants, soil biota and terrestrial wildlife
- Puget Sound Background Metals Concentrations.



Soil concentrations protective of groundwater's highest beneficial use are calculated conservatively using Ecology's variable parameter 3-phase partitioning model (WAC 173-340-747[5]), and using the most stringent groundwater screening level, including potable (drinking) water criteria and criteria protective of surface water quality for the protection of aquatic organisms and human health based on consumption of aquatic organisms. MTCA-default parameter values are used in the 3-phase model. Based on shallow groundwater levels, which often result in surface flooding, the calculated Method B cleanup levels assume that all soil at the Site is saturated.

The soil concentrations generated by this MTCA-default methodology are intentionally conservative and are intended for preliminary data screening in the RI. Soil concentrations above these screening levels may or may not actually be leaching contaminants to groundwater at concentrations of concern. MTCA provides a range of options to further evaluate site-specific soil concentrations protective of groundwater, including use of soil leaching tests and empirical groundwater quality data, as outlined in WAC 173-340-747. The soil-to-groundwater-based soil screening levels may not be considered for a chemical if it can be demonstrated that soil concentrations are protective of groundwater using methods in WAC 173-340-747.

The preliminary soil criteria incorporated into the derivation of soil SSLs to be applied for the RI are included in Table 4 for each contaminant of concern (COC) at the Site.

3.2.1.2.2. Groundwater and Surface Water

SSLs for groundwater include consideration of the following:

- Standard MTCA Method A and B cleanup levels for the protection of human health through direct contact and ingestion for unrestricted land use.
- State and federal maximum contaminant levels for the protection of human health through direct contact and ingestion.
- Applicable state and federal criteria for the protection of surface water quality: protection of human health and aquatic life.

In accordance with MTCA, groundwater screening levels protective of surface water incorporate MTCA surface water cleanup levels including criteria from applicable state and federal laws (WAC 173-340-730). For protection of fresh water quality, screening levels are the most stringent of the following aquatic life criteria and human health criteria for consumption of aquatic organisms under state and federal laws:

- Washington State human health criteria for the consumption of water and organisms, EPA-approved values under Section 303(c) of the Clean Water Act.
- Federal National Recommended Water Quality Criteria pursuant to Section 304(a) of the Clean Water Act.



- Washington State Water Quality Standards (WAC 173-201A-240).
- Standard MTCA Method B surface water cleanup levels based on human consumption of fish (human health only).

The preliminary groundwater criteria incorporated into the derivation of groundwater SSLs to be applied for the RI are included in Table 5 *for* each COC at the Site.

3.2.2. Data Gaps Assessment

This section describes data gaps identified during previous Site investigations (see Section 3.1) and provides a description of the work that was conducted to address them.

3.2.2.1.1. 1985 Abandoned Landfill Study

The 1985 Site investigation by Public Health – Seattle & King County did not characterize the lateral or vertical extent of refuse or potential soil, groundwater, or surface water contamination at the site. These data gaps were addressed in subsequent investigations with soil borings, test pits, monitoring wells, groundwater sampling, and a geophysical survey.

3.2.2.1.2. 2010 and **2011** Stormwater Sampling

No conclusions regarding the source of the VOCs detected in stormwater samples were made by King County (2015). This data gap was addressed in subsequent investigations by collecting surface water and stormwater samples from the ditch and the stormwater pond along the west side of the Site and by collecting groundwater samples from push probes and monitoring wells throughout the Site.

3.2.2.1.3. 2016 Phase II ESA

The 2016 Phase II ESA was conducted to characterize soil and groundwater conditions across the site, and to determine where contaminated soils may be encountered during excavation activities associated with future levee setback work (S&W 2016). The assessment concluded that additional soil and groundwater data was needed to determine if soil and/or groundwater contamination was present outside of the fill area, and to determine the potential for offsite impacts. These data gaps were addressed in subsequent investigations with soil borings, additional monitoring wells, groundwater and surface water sampling, and a geophysical survey.

3.2.2.1.4. 2017 Environmental Investigation

The 2017 Environmental Investigation was conducted to determine if soil and/or groundwater contamination was present outside of the fill areas identified during the 2016 Phase II ESA, and to determine the potential for offsite impacts (Herrera 2017). The report concluded that three additional monitoring wells should be installed for the purpose of collecting samples for metals analyses from permanent wells that produce less turbid water samples than temporary soil borings. The report also concluded that monitoring wells should be resampled during seasonal high and low

water level periods to evaluate effects of seasonal changes in groundwater levels and flow directions on groundwater quality. These data gaps were addressed by installing upgradient monitoring well MW-7, and downgradient monitoring wells MW-8 and MW-9, and collecting quarterly groundwater samples in March, June, and September 2018 from all wells at the Site.

3.2.3. 2018 Environmental Investigation

The 2018 investigation work was conducted to address the data gaps identified and to support development and evaluation of alternatives for permanent flood protection along the right bank of the White River. The specific objectives of the sampling were to:

- Further refine the understanding of the lateral extent and thickness of refuse within the Study Area.
- Evaluate the presence and concentrations of chemical constituents both within the refuse and in soils surrounding the refuse.
- Collect sufficient physical and chemical information about the refuse and surrounding soil conditions to support the development and evaluation of feasible remedial alternatives.
- Evaluate variability in seasonal groundwater fluctuations and flow direction.
- Evaluate groundwater near the refuse for the presence and concentration of chemical constituents.

3.2.3.1. Investigation Methods

The 2018 environmental investigation consisted of the following components.

- January: completed a geophysical survey to characterize the depth and lateral extent of refuse and help target the locations of geotechnical and environmental explorations at the Site.
- February and March: completed 25 push probe environmental borings, and 13 geotechnical borings with environmental sampling, and installed three new monitoring wells (MW-7, MW-8, and MW-9) at three of the geotechnical boring locations.
- March, June, and September: collected groundwater samples from six existing monitoring wells, and the three new monitoring wells.
- June, and September: collected surface water samples at three locations within the stormwater drainage ditch along the west edge of the site, and at a fourth location within the onsite stormwater pond during the September sampling event.
- March, June, and September: monitored landfill gas at Site monitoring wells where the static groundwater level was below the top of the well screen.

3.2.3.1.1. Soil Sampling Methods

The work was conducted in accordance with the Sampling and Analysis Plan (Herrera 2018).

Soil samples from geotechnical borings were collected in 18-inch, split spoon samplers at 2.5-foot intervals from 2.5 feet bgs to the total depth of each boring. The soil was visually inspected for staining and classified in accordance with the ASTM International (ASTM) D2488, Standard Practice for Description and Identification of Soils (Visual-Manual Procedure). Discrete soil samples were collected for potential chemical analysis. The boring logs, including well construction details, are provided in Appendix A.

Soils encountered during drilling were visually inspected and classified in accordance with the Unified Soil Classification System (USCS; ASTM International [ASTM] D2488-09), and a photoionization detector (PID) instrument was used to screen soil for VOCs. Following soil sample collection, each boring was backfilled from the bottom to ground surface with bentonite chips and capped at the surface with soil or cement to match surrounding surface conditions. The soil boring logs are presented in Appendix A.

Sample collection intervals were adjusted in the field depending on the presence or absence of refuse and the homogeneity of the subsurface material. Sample intervals generally included:

- The first sample collected from the 0- to 5-foot depth interval or groundwater interface above the refuse if present
- A second sample from the 5- to 10-foot interval or within refuse
- A third sample from the 10- to 15-foot interval, below refuse (or from the 15- to 20-foot interval at locations within clean waste deep soil grids previously established by Shannon & Wilson [2016]).

A total of 93 soil samples from push probes PP10 through PP24 and geotechnical borings B-04 through B-17 were submitted to OnSite for the following laboratory analyses (see Figure 5):

- TPH identification using Ecology's Northwest total petroleum hydrocarbons, dieselextended (NWTPH-Dx) method
- Total MTCA metals (arsenic, cadmium, chromium, mercury, and lead) using EPA Methods 6010C/7471A.
- Carcinogenic PAHs using EPA Method 8270D/SIM.

If TPH was identified in a sample, the sample was also analyzed for:

 Gasoline-range TPH and benzene, toluene, ethylbenzene and xylenes (BTEX) using Ecology's NWTPH-Gx method/EPA Method 8021B; and/or



- Diesel- and oil-range TPH using Ecology's NWTPH-Dx method; and
- PCBs using EPA Method 8082A (if oil-range TPH is detected).

3.2.3.1.2. Groundwater Sampling Methods

Monitoring wells MW-7, MW-8 and MW-9 drilled and constructed in accordance with the Minimum Standards for Construction and Maintenance of Wells, WAC 173-160. The borings were completed by a Holocene Drilling, Inc. of Puyallup, Washington using hollow stem auger drilling methods. The monitoring wells are each constructed with a 2-inch diameter PVC well casing with an 11-foot long well screen set from 4 feet to 15 feet bgs. Filter pack sand was placed in the annular space between the borehole and well screen. Bentonite chips were placed above the filter pack and each well was completed at the ground surface with flush-mounted steel monument sealed in concrete.

Following installation, the monitoring wells were developed with a submersible pump. Development continued until levels of sand and silt were reduced and water removed from the well was generally of clear quality. Development water from each well was contained in 55-gallon drums, stored temporarily at the site, pending analytical results.

On March 23, June 21, and September 23, 2018, groundwater samples were collected from monitoring wells MW-1 through MW-9, and submitted to Onsite for the following laboratory analyses:

- Gasoline-range TPH and BTEX using Ecology's NWTPH-Gx method/EPA 8021B;
- Diesel- and oil-range TPH using Ecology's NWTPHDx method;
- Total MTCA metals (arsenic, cadmium, chromium, mercury, and lead) using EPA Methods 6010C/200.8/7470A; and
- Carcinogenic PAHs using EPA Method 8270D/SIM.
- Volatile organic compounds using EPA Method 8260C.

If oil-range hydrocarbons were identified in a sample, the sample was also analyzed for PCBs using EPA Method 8280A.

Groundwater samples were collected according to procedures outlined in the Sampling and Analysis Plan (Herrera 2018) as follows:

- The well monument cover was removed, and the condition of the well and surrounding area was inspected.
- Observations were noted in the field notebook and on the well sampling log.



- The well casing plug was removed.
- The depth to groundwater was measured from the top of the well monument to the nearest 0.01 foot using an electronic water level indicator.
- The date, time, and measurements were recorded on the well sampling log.
- Care was taken to ensure that no bubbles or headspace were present for the VOC samples.
- Containers were securely capped, labeled, and placed into a chilled cooler for storage, prior to delivery to the laboratory.
- The date and time of sample collection was recorded on the well sampling log and chain-of-custody form.

The wells were purged of standing water using a low-flow purge method at approximately 0.4 liters per minute with clean, dedicated polyethylene tubing and a submersible pump. Tubing was placed 3 to 5 feet below the top of the water table and opposite the screened zone in the monitoring wells.

During purging, pH, water temperature, dissolved oxygen, specific conductivity, water level, and pumping rate were measured. The amount of water purged, water parameter measurements, and time of collection were recorded on the well sampling log. Recharge occurred quickly for the wells; the water levels remained constant during purging and sampling. Purged water removed during development was placed into 55-gallon drums stored on site.

Samples were collected with the same dedicated polyethylene tubing used to purge the well by pumping directly into sample containers provided by the analytical laboratory.

3.2.3.1.3. Surface Water Sampling Methods

Surface water samples were collected according to King County Standard Operating Procedure (SOP) Sampling Methods for Stream and River Water (SOP #214v3) as follows:

- One-liter unpreserved amber glass bottles were dipped directly into the surface water to fill them, and one of the unpreserved bottles was used to collect and transfer water into other bottles containing preservative.
- The date, time, sample identification, sampler name, and requested analysis were recorded on each sample bottle label and on the chain-of-custody form.
- Field measurements for temperature (degrees centigrade), pH (standard units), dissolved oxygen (milligrams/liter [mg/L]), and specific conductivity (µs/cm), and turbidity (nephelometric turbidity units) were recorded in a log book using a calibrated YSI Pro DSS multi-meter.

- Care was taken to ensure that no bubbles or headspace were present for the VOC samples.
- Containers were securely capped, labeled, and placed into a chilled cooler for storage, prior to delivery to the laboratory.

On June 29, 2018, surface water samples were collected from three locations (SW1 thru SW3) along the stormwater ditch in conjunction with the June 2018 quarterly groundwater sampling event. On October 9, 2018, surface water samples were collected during a rainstorm from the same three locations along the ditch, plus a fourth location (SW4) within the northeast corner of the onsite stormwater pond. The surface water samples were submitted to Onsite for the following laboratory analyses:

- Gasoline-range TPH and BTEX using Ecology's NWTPH-Gx method
- Diesel- and lube oil-range TPH using Ecology's NWTPH-Dx method
- Total MTCA metals (arsenic, cadmium, chromium, mercury, and lead) using EPA Methods 6010C/200.8/7470A
- cPAHs using EPA Method 8270D/SIM
- VOCs using EPA Method 8260C
- Hardness by EPA Method 6010D/SM 2340B.

Samples were held for potential analysis of PCBs pending the results of oil-range TPH analysis, but no oil-range petroleum hydrocarbons were detected so subsequent PCB analysis was not performed.

3.2.3.1.4. Soil Vapor Monitoring Methods

Landfill gas monitoring was performed using a Landtec Gas Analyzer & Extraction Monitor (GEM) 2000 Plus gas analyzer and extraction monitor that was calibrated by the supplier prior to use. During the March 2018 event the wells were purged using the GEM at a rate of approximately 300 milliliters per minute (ml/min), and during the June 2018 event, the wells were purged during monitoring using an Aircheck Sampler pump by SKC, Ltd. at a rate of approximately 3,000 ml/min. Table 8 presents a summary of the gas monitoring conducted in 1985 by King County, as well as the 2018 monitoring conducted by Herrera.

Landfill gas monitoring was conducted March 23 and June 21, 2018, at two locations (MW-6 and MW-9; and September 23, 2018, at three locations (MW-1, MW-6, and MW-9) (Figure 11). These three wells were the only wells where the measured groundwater elevations were below the top of the well screen.

3.2.3.2. Sampling/Analytical Results

The following subsections present the sampling methods and the analytical results by media for the 2018 investigation work described above.

3.2.3.2.1. Geophysical Survey

On January 30 and 31, 2018, a geophysical survey was completed to further delineate the lateral and vertical extent of buried metal debris identified in 2016, and to aid in targeting additional geotechnical and environmental explorations (Duoos, P. 2018; S&W 2016). The survey study area included the open park area and did not extend into the wooded area to the south or beyond the Site boundary. A copy of the geophysical report is included in Appendix F.

Electromagnetic and magnetic geophysical methods were used along a grid established at the Site to determine the general lateral extent of buried refuse, and a magnetometer was also used to search for buried ferrous (iron-bearing) metal.

Areas interpreted to have buried metal debris were categorized as low, moderate, and high anomalous zones. Except for one low metal anomalous zone identified near the northwest part of the Site, most of the buried metal debris anomalies were located near the central and southeastern portions of the site. A long, linear feature was identified along the southeast edge of the Site, parallel to the White River; it is possible that this feature is part of the historic levee.

3.2.3.2.2. Soil Results

A summary of soil analytical results is presented in Table 7. Figures 12 through 15 depict the extent of soil contamination from diesel- and lube-oil range petroleum hydrocarbons, lead, PCBs, and total cPAHs.

Gasoline-range petroleum hydrocarbons were initially detected by method NWTPH-hydrocarbon identification (HCID) in samples at only three locations: PP23, PP24, and PP25. Follow-up analysis detected a concentration of gasoline-range petroleum hydrocarbons in PP25-13 (400 milligrams/kilogram [mg/kg]) that exceeded the SSL of 100 mg/kg; BTEX compounds were not detected above the laboratory reporting limits; however, the RLs for benzene, toluene, and ethyl benzene were higher than the SSLs.

Diesel- and lube-oil range petroleum hydrocarbons were initially detected by method NWTPH-HCID in 25 samples at 18 locations. Follow-up analysis detected diesel-range petroleum hydrocarbons that exceeded the SSL of 200 mg/kg in samples from four locations at concentrations ranging from 400 to 1,800 mg/kg.

Lube oil-range petroleum hydrocarbons exceeded the SSL of 2,000 mg/kg in samples from five locations at concentrations ranging from 3,200 to 12,000 mg/kg. BTEX compounds were not detected above the laboratory reporting limits.



A review of laboratory chromatograms for the analytical results of Ecology Method NWTPH-Dx for quantification of diesel- and oil-range TPH indicates that most of the TPH present in soil at the Site is lube oil. The one exception is diesel-range TPH detected in soil at boring B-05, which appears to have a chromatographic signature similar to weathered diesel fuel. In accordance with Ecology Implementation Memorandum #4, Determining Compliance with Method A Cleanup Levels for Diesel and Heavy Oil, the NWTPH-Dx results for all samples (except B-05-7.5) have been summed to calculate a Total NWTPH-Dx value that is compared to the SSL for lube oil-range TPH of 2,000 mg/kg.

MTCA metals were detected above SSLs in samples from 21 locations. Lead was the metal most frequently detected above the SSL and was therefore used as an indicator of metals contamination across the Site. Arsenic exceeded the SSL of 20 mg/kg in only one sample: PP15-12 (24 mg/kg). Cadmium exceeded the SSL of 1 mg/kg in 18 samples from 13 locations at concentrations ranging from 1.2 to 27 mg/kg. Chromium exceeded the SSL of 48 mg/kg in six samples from five locations at concentrations ranging from 53 to 430 mg/kg. Lead exceeded the SSL of 25 mg/kg in 38 samples from 20 locations at concentrations ranging from 30 to 7,300 mg/kg. And mercury exceeded the SSL of 0.07 mg/kg in four samples from 3 locations at concentrations ranging from 0.41 to 1.2 mg/kg.

Total PCBs were detected above the SSL of 0.33 mg/kg in samples from four locations at concentrations ranging from 0.74 to 1.36 mg/kg.

Concentrations of total cPAHs were detected above the SSL of 0.09 mg/kg in soil samples collected from three locations at concentrations ranging from 0.094 to 1.2 mg/kg.

3.2.3.2.3. Groundwater Results

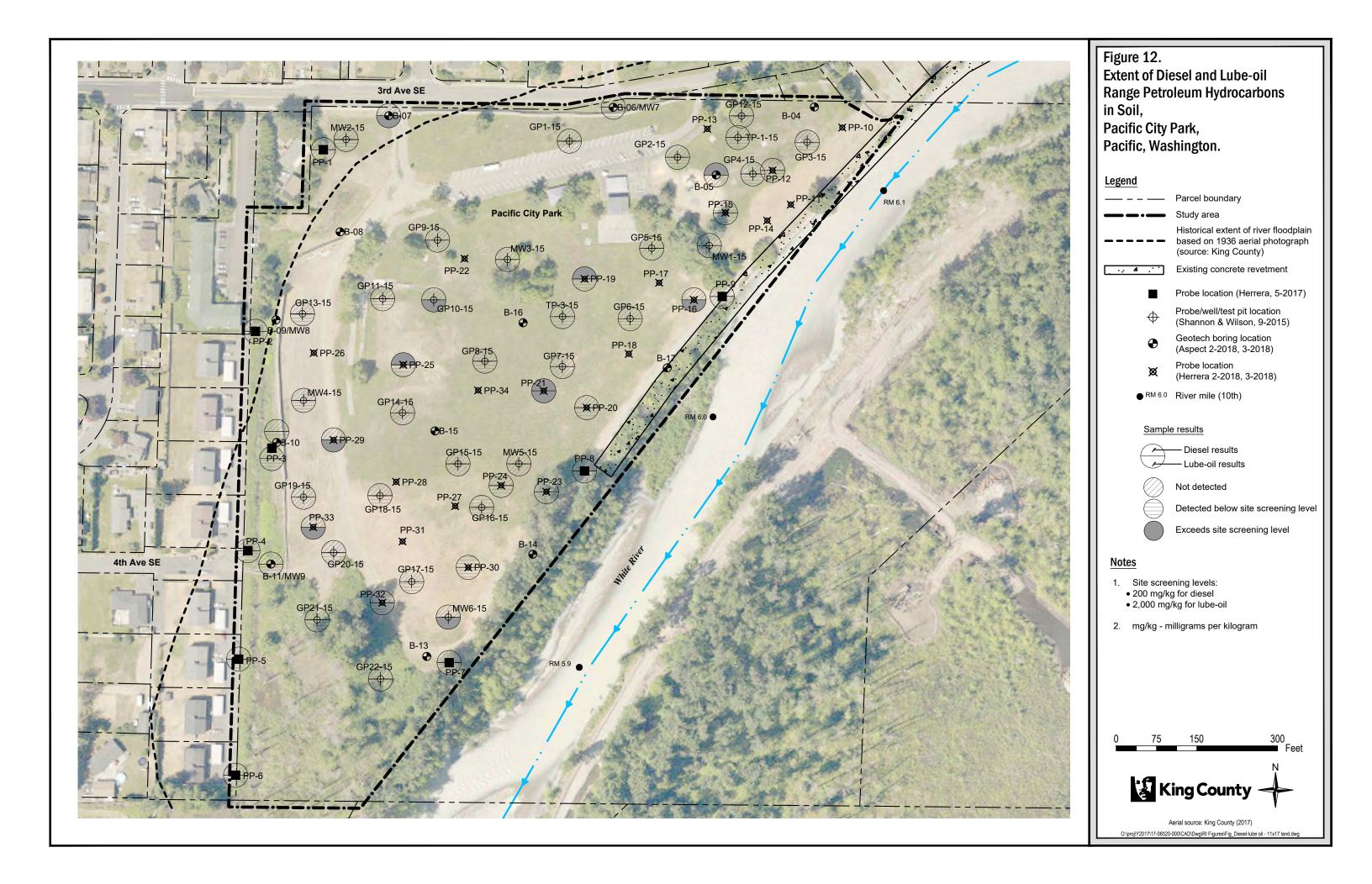
A summary of groundwater analytical results for samples collected from monitoring wells is presented in Table 6. No gasoline-, diesel-, or lube oil-range petroleum hydrocarbons were detected above SSLs in groundwater samples collected from the nine monitoring wells in March, June, and September 2018.

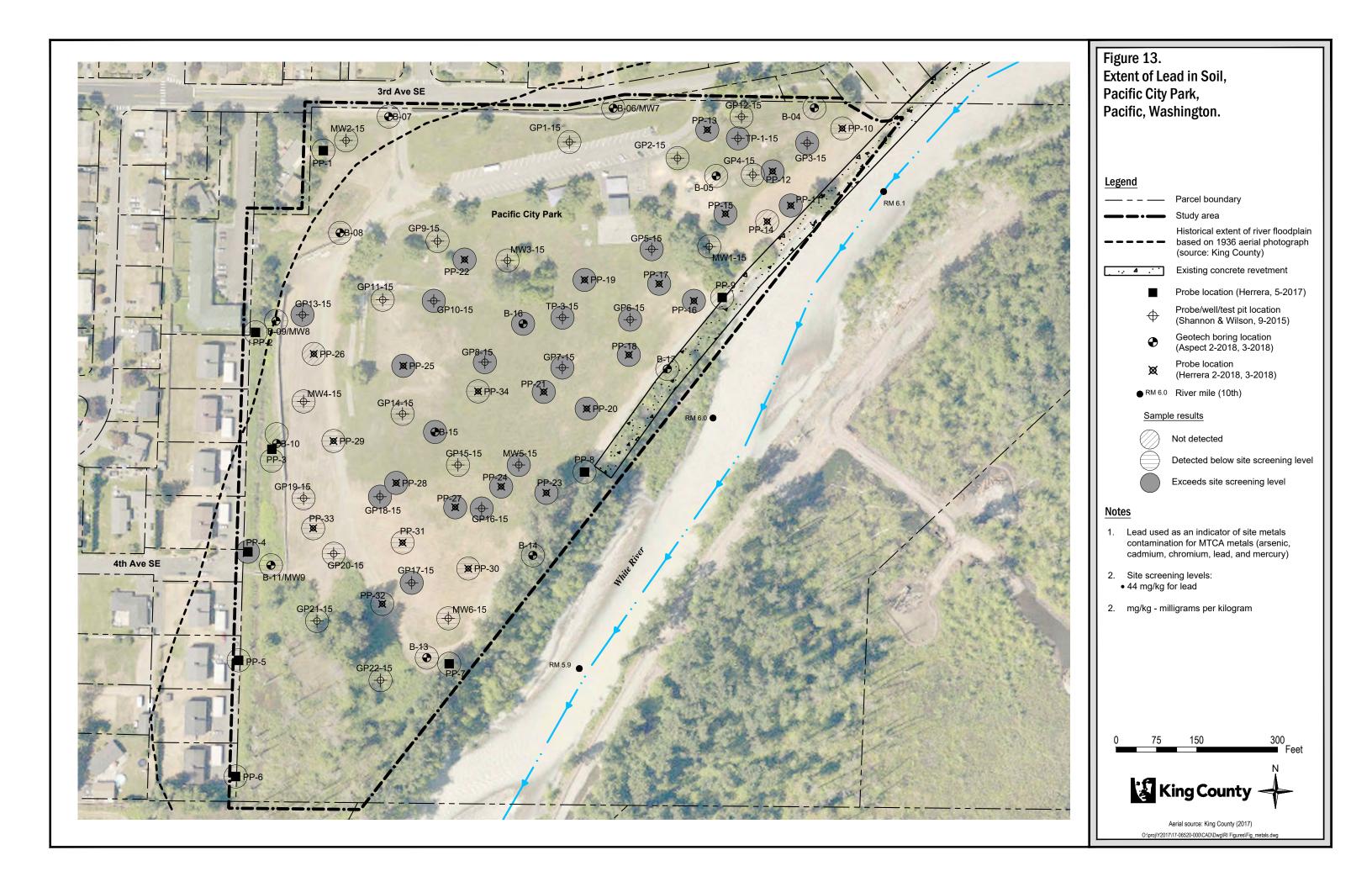
One VOC, (cis) 1,2-dichloroethene, was detected below the SSL in two samples from MW3; 1,4-dichlorobenzene was detected below the SSL in one sample from MW6, and chlorobenzene was detected below the SSL in three samples from MW4. No other VOCs or total PCBs were detected above the SSLs in any of the samples from the wells.

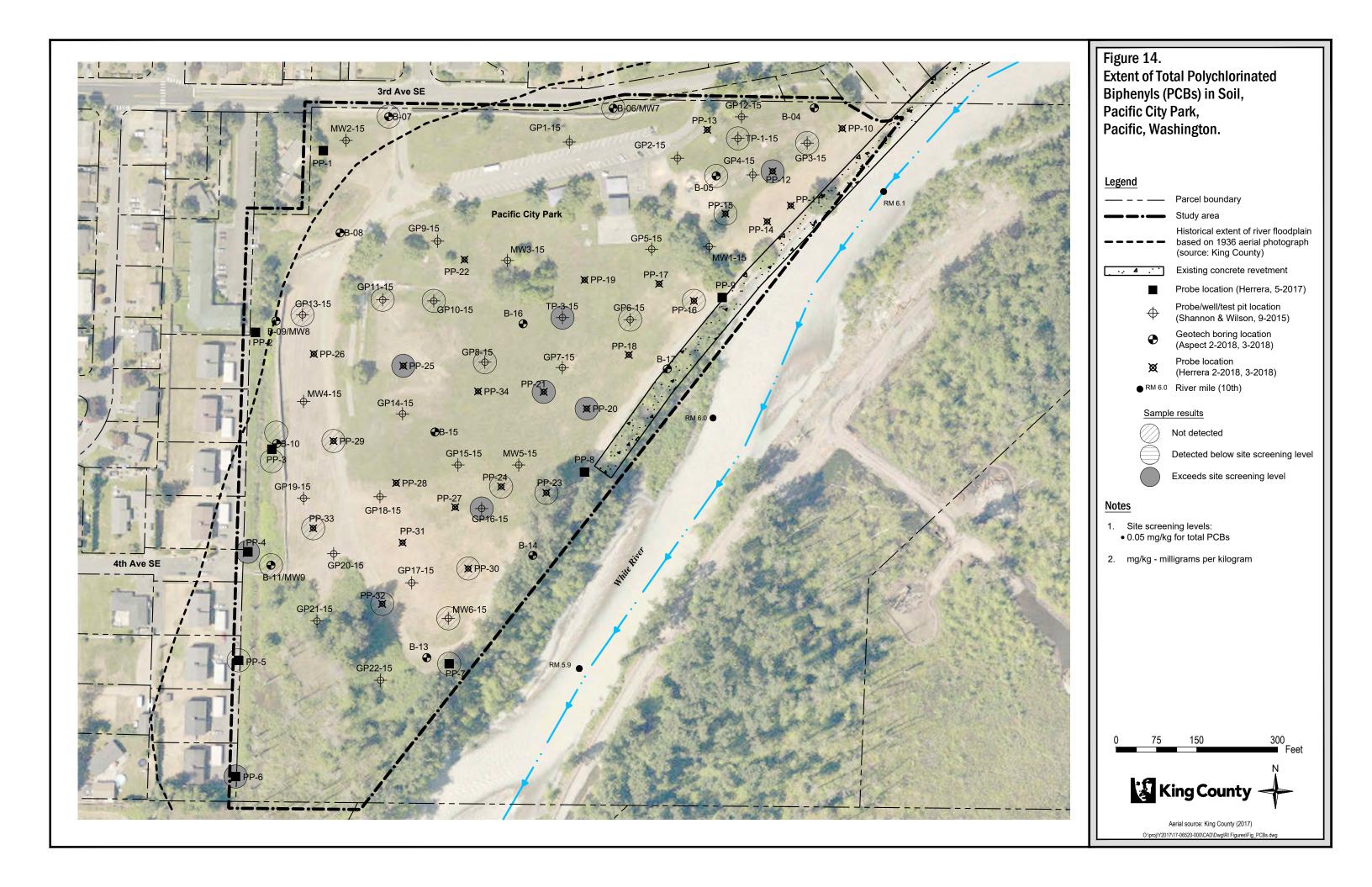
Of the five MTCA metals, total arsenic was detected above the SSL of 3.3 μ g/L in 13 samples from 7 wells at concentrations ranging from 3.4 to 14 μ g/L, and total lead was detected above the SSL of 2.5 μ g/L 3 samples from one well at concentrations ranging from 3.2 to 5.6 μ g/L. Dissolved arsenic was detected above the SSL in four samples from three wells at concentrations ranging from 4.8 to 9.8 μ g/L, and dissolved lead was detected above the SSL in one sample from one well at a concentration of 3.0 μ g/L.

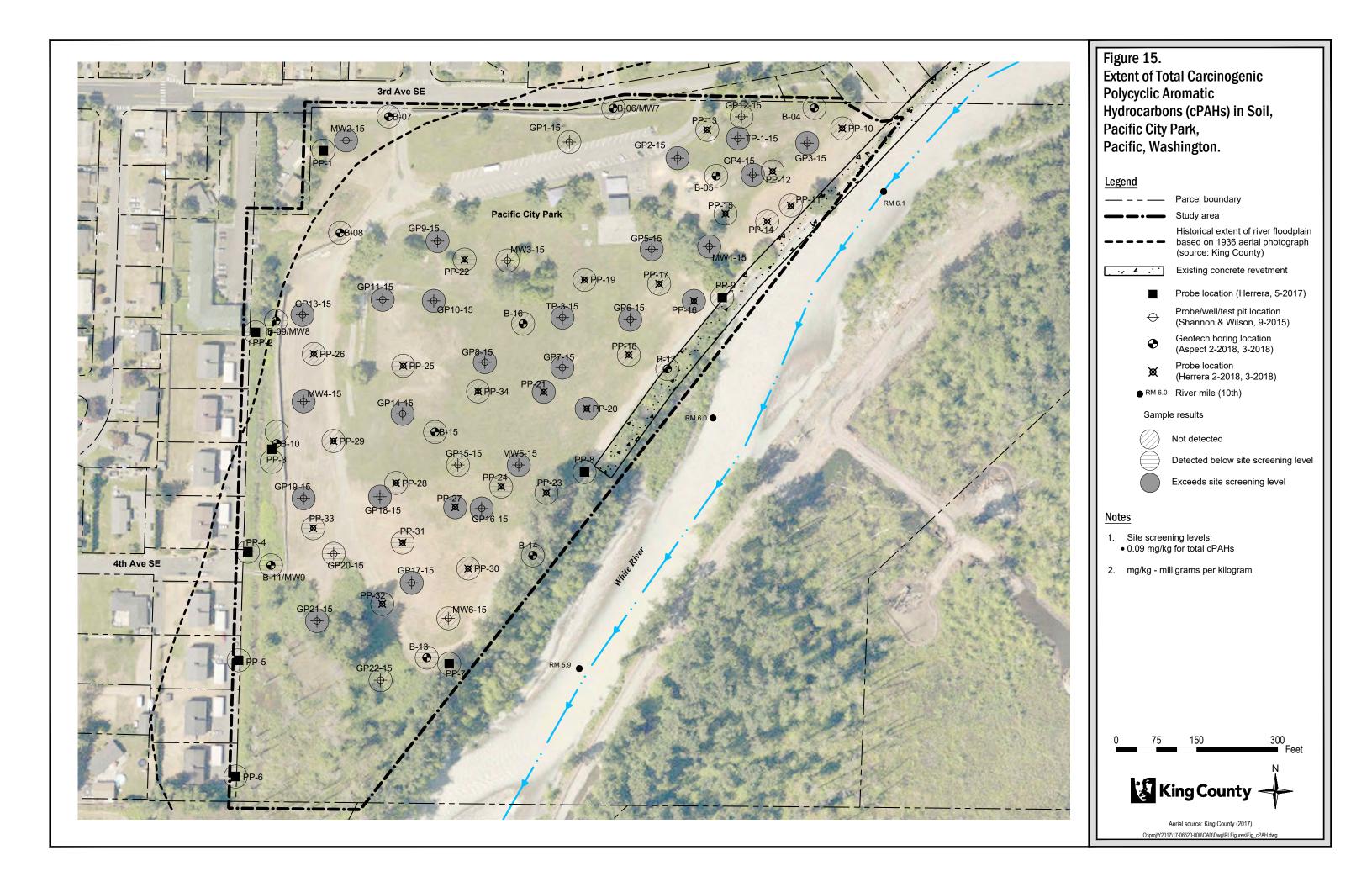
The concentration of total cPAHs exceeded the SSL of 0.015 μ g/L in one sample from MW-6 (0.017 μ g/L).











3.2.3.2.4. Surface Water Results

Surface water analytical results were compared to the groundwater SSLs protective of surface water shown in Table 2. No petroleum hydrocarbons, total MTCA metals, cPAHs, or VOCs were detected above the SSLs in any of the samples.

3.2.3.2.5. Soil Vapor Results

No methane (CH₄) or hydrogen sulfide (H₂S) were detected at MW-6 or MW-9 during the March, June, or September 2018 sampling events. In 1985, only low concentrations of methane (up to 0.4 percent) and low levels of trace gases (up to 6.2 ppm) were detected. Based on these results it does not appear that the Site is producing or releasing landfill gasses.

3.2.3.3. Data Quality Analysis

Laboratory analyses for the 2018 investigation were performed by OnSite Environmental, of Redmond, Washington, an Ecology-accredited laboratory. Laboratory reports and chain of custody forms are provided in Appendix G.

A data quality assurance review was completed by Herrera for all analyses performed (See memorandum included in Appendix H.) Data were validated based on the following:

- Sample custody, preservation, holding times, and completeness
- Laboratory reporting limits
- Method blank analysis
- Laboratory control sample analysis
- Surrogate compound analysis
- Matrix spike analysis
- Laboratory duplicate analysis



4. CONCEPTUAL SITE MODEL

The conceptual site model (CSM) for the Site has been developed based on historical land use information, the results of previous environmental investigations, and results from the RI work completed in 2018 to address data gaps. The CSM is the basis for developing technically feasible cleanup alternatives and selecting a final cleanup action. The CSM is dynamic and may be refined as additional information becomes available for the Site. The following sections describe the components of the CSM including the source of the COPCs, nature and extent of contamination, contaminant fate and transport, and a preliminary exposure assessment.

4.1. Sources of Chemicals of Potential Concern

Historical records indicate that municipal waste was placed on the Site as early as 1920 until approximately 1965. The previous investigations identified fill soil mixed with refuse that consists of pieces of concrete, wood, shards of glass and china, and bits of metal debris. Most of the refuse was reportedly burned. The Site was vacant until it was leased to the City of Pacific in 1969 and subsequently opened as a city park in 1972. Additional fill soil was placed throughout the Site through the 2000s, primarily in the south, southwest, west and north portions of the site.

The sources of COPCs include fill soil containing ubiquitous but low concentrations of metals and total cPAHs, and fill soil mixed with refuse, which analytical testing indicated the presence of TPH, PCBs and VOCs. Suspected sources of COPCs that were historically identified and evaluated included treated wood and possible nearby uses of pesticides and herbicides.

The approximate lateral extent of fill soil and refuse are depicted on Figures 3 and 5. The approximate vertical extent of fill soil and refuse are depicted on cross sections A-A' and B-B' (Figures 6 and 7, respectively).

4.2. NATURE AND EXTENT

The nature and extent of the COPCs at the Site are described based on the sources, physical conditions and analytical data collected during previous investigations and during the 2018 site characterization work. The Site is defined by the extent of hazardous substances in one or more media, at concentrations exceeding the SSLs. The presence of fill soil alone, in the absence of COPCs, does not fall under the definition of MTCA as a hazardous substance, and therefore, is not part of the MTCA Site. However, the presence of municipal waste does qualify as a hazardous substance. The Site is defined by any location where municipal waste and/or COPCs are present in soil, groundwater, surface water or soil vapor at concentrations exceeding the SSLs. The following subsections summarize the known distribution of concentrations of COPCs at the Site.



4.2.1. Physical Conditions

The Site consists of a relatively flat upland area bounded by the White River to the east, a paved road to the north, and a surface water drainage ditch and multi-family residential properties located to the west (Figure 2). During the wet season, typically between October and March, the central and northern portions of the Site are frequently inundated by the surface expression of shallow groundwater. Groundwater depths ranged from 4 to 9 feet bgs at the time of the Phase II investigation, in September 2015. In general, groundwater was encountered between 4 to 6 feet bgs, with deeper groundwater depths encountered in areas of higher relief, such as explorations located on the existing levee and in terraced fill areas in the southern portion of the Site (S&W 2016). The groundwater flow direction suggests that the White River loses surface water to the upper alluvial aquifer, at least seasonally.

Shallow groundwater levels in the unconfined, upper alluvial aquifer are also encountered seasonally in monitoring wells located on the Site and springtime groundwater flow has been observed to the west and southwest (Figures 8, 9, and 10). King County completed an animation of groundwater levels, based on hourly well measurements from October 2015 through October 2016. The direction of groundwater flow is typically to the southwest (parallel to river flow) for most of the year. When water levels rise in winter with the river stage, the gradient increases and has a more westerly component, away from the river (Brummer 2017).

The geology at the Site consists of fill overlying native Holocene alluvium to the maximum depth explored of approximately 62 feet bgs. Fill material was identified as soil with a disturbed appearance and from the presence of unnatural debris (refuse) such as glass shards and bottles, brick, cement, organics, wood, paper, rubber, and ceramic. The fill consists of fill soil and fill soil mixed with refuse. The fill soil consists of variable amounts of silt, sand and gravel that generally directly overlies the Holocene alluvium deposits across the Site, and ranges in thickness from inches to 15 feet. The fill soil mixed with refuse is primarily located beneath the eastern portions of the Park, along the White River, and is up to 20 feet thick in the central and south-central portions of the Site (Figures 6 and 7). The refuse is typically covered by one to four feet of sandy, gravelly fill soil.

The fill is underlain by alluvial overbank deposits consisting mostly of poorly graded sand with gravel to sandy gravel interbedded with backwater lacustrine deposits consisting of silt with organics and interbeds of silty sand.

Figure 5 depicts the estimated lateral extents of fill soil and fill soil with refuse and the transect locations for cross sections that depict subsurface characteristics at the Site. Cross sections A-A' and B-B' are depicted on Figures 6 and 7, respectively.

4.2.2. Soil

The primary COPCs exceeding the SSLs in soil are lead and total cPAHs, which are present in fill soil and refuse throughout the Site (Figures 13 and 15, respectively). Other primary COPCs are present in soil at concentrations exceeding the SSLs sporadically throughout the Site in the fill soil and refuse, including metals (arsenic, barium, cadmium, chromium, mercury and silver), TPH and PCBs (Table 7). For the purposes of the RI, lead is used as an indicator of the presence of all metals in soil at concentrations exceeding the SSLs. Although all metals are not present in soil above SSLs, where lead is reported above the SSLs, the soil samples that contain metals exceeding the SSLs almost always contain lead above the SSLs. The soil data for total cPAHs, lead, TPH and PCBs are depicted on Figures 15, 13, 12, and 14, respectively.

Secondary chemicals that have been detected in soil at the Site at concentrations exceeding the SSLs, include the following:

- 1. Chlorinated VOCs, including tetrachloroethene (PCE) and trichloroethene (TCE);
- 2. Other SVOCs, including non-carcinogenic PAHs, pentachlorophenol (PCP) and phthalate compounds; and
- 3. Pesticides.

Many of these secondary chemicals have been detected in a single soil sample collected at the Site (Table 7.) All of these secondary chemicals have been detected in soil samples collected from areas of the Site where fill soil with refuse has been observed and are co-located with concentrations of total cPAHs and lead in soil. Developing a cleanup action to address exposure risk to human health and the environment from the primary COPCs in the areas of the Site where refuse is located will include those areas where these secondary chemicals have been detected above SSLs, so they are not discussed further in the CSM.

The highest concentrations of total cPAHs and/or lead in soil are generally observed in soil borings where refuse was observed, including the area of borings PP-17, PP-20 and PP-21, where refuse was observed to be up to 20 feet thick (see Figure 5). Relatively high concentrations of total cPAHs and lead are also reported in soil at boring GP10-15, where historical photos suggest refuse disposal in the late 1950s and into the mid-1960s (Table 4). The lateral extent of total cPAHs and lead in soil has not been defined in soil to the west-southwest of borings PP-4, PP-5 and PP-6, where fill soil extends beyond the property boundary and the results of soil samples collected within the fill at the property boundary contain concentrations of one or more of the COPCs above the SSLs. The vertical extent of lead and total cPAHs in soil, depicted on the cross-sections (Figures 6 and 7), is defined by the results of soil samples collected from native Holocene alluvium soils located beneath the fill soil and refuse (Figures 13 and 15). The soil data indicates that COPCs are present above the SSLs in the upper 2- to 8-feet of native soil beneath the fill and refuse.



The locations where TPH was detected in soil above the SSLs are depicted on Figure 12. The distribution of TPH in soil at the Site, at concentrations above the SSLs, is typically limited to shallow depths in the fill soil. Except for borings PP-21 and PP-25, all the total NWTPH-Dx concentrations above the SSLs are reported in fill soil located at depths ranging from 1 foot to 5.5 feet bgs. When deeper, native soil was sampled in these same borings, the concentrations of total NWTPH-Dx are below the SSLs, defining the vertical extent (Table 7). Similarly, the concentrations of total NWTPH-Dx above the SSLs in borings PP-21 and PP-25 are reported at depths of 10 and 13 feet bgs, respectively near the fill soil-native soil interface, and deeper soil samples collected from both borings (at depths of 15 and 17 feet bgs, respectively) do not contain total NWTPH-Dx concentrations above the SSLs (Table 7).

Concentrations of total PCBs in soil are generally below 1 mg/kg, except for soil samples collected from borings PP-21 (1.33 mg/kg) and PP-25 (1.36 mg/kg) (Figure 14). With the exception of boring PP-4, the presence of total PCBs at concentrations above the SSLs in soil generally correlates to the presence of fill soil with refuse.

The extent of fill soil has been inferred based on historical aerial photographs, however the presence and quality of fill soil beyond the property boundaries has not been evaluated. The lateral extent of one or more of the COPCs, at concentrations above the SSLs, in either fill soil or fill soil mixed with refuse is shown on Figure 16. The lateral extent of COPCs in fill soil are sufficiently defined to the north to characterize the quality of the fill soil and to allow for the development and evaluation of remedial alternatives. Further evaluation into the extent and quality of fill soil to the west-southwest is needed to adequately characterize the Site for the purpose of developing and evaluating remedial alternatives.

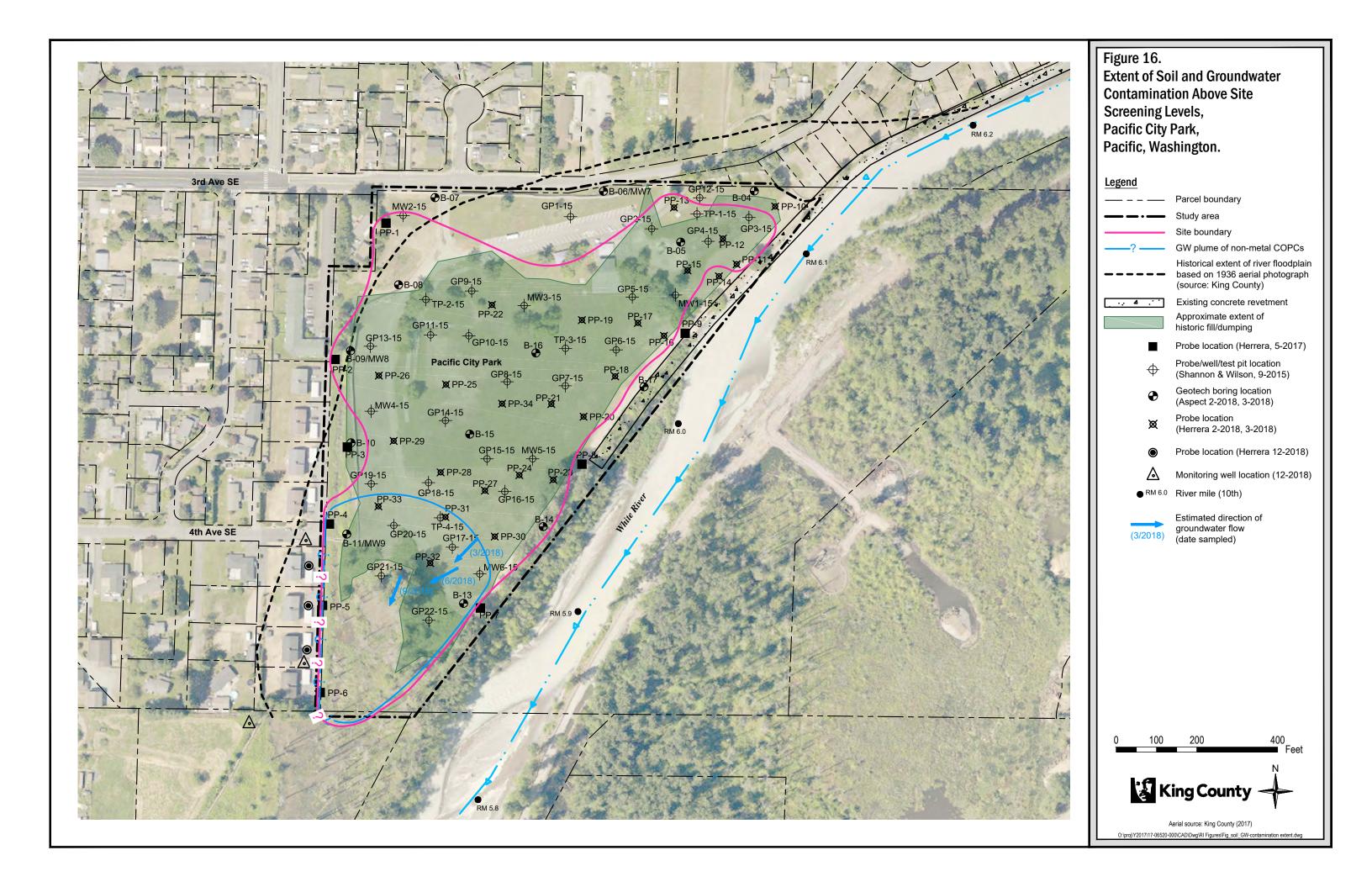
4.2.3. Groundwater

The nature and extent of COPCs in groundwater is based on the results of laboratory analytical results of reconnaissance groundwater samples collected from borings in 2017 (Table 3) and from groundwater samples collected from Site monitoring wells during three separate sampling events conducted between 2015 and 2018 (Table 6)

MTCA (WAC 173-340-720(9)) dictates that groundwater cleanup levels shall be determined by analysis of groundwater samples representative of the groundwater and that analysis shall be conducted on unfiltered samples unless it can be demonstrated that a filtered groundwater sample provides a more representative measure of groundwater quality.

A comparison of total and dissolved metals concentrations detected in groundwater samples collected from the Site suggest that filtered groundwater samples (dissolved metals data) provide a more representative measure of groundwater quality for metals than unfiltered samples (total metals data), for purposes of determining groundwater compliance evaluation under MTCA.





The metals being analyzed for in Site groundwater samples are naturally occurring, inorganic hazardous substances (Ecology, 1994), and even low quantities of suspended solids in a sample can bias high the detected metals concentration in an unfiltered groundwater sample compared to what is dissolved in groundwater. This is evident in the reconnaissance groundwater samples collected from the PP-borings in 2017, where there is a stark difference between the reported concentrations of total metals and dissolved metals in samples collected from temporary borings, where the volume of suspended solids in a sample is typically high. In contrast, there is little difference in the reported concentrations of total metals and dissolved metals in groundwater samples collected from properly developed monitoring wells, where the groundwater turbidity is low.

Assuming that the dissolved metals concentrations are representative of groundwater quality at the Site, the COPCs that have been detected in groundwater at concentrations above the SSLs consist of metals (arsenic and lead), benzene and total cPAHs.

Arsenic has been detected above the SSLs in groundwater samples collected from across the Site (Figure 17). The SSL for arsenic is based on the protection of surface water, adjusted upward in consideration of the laboratory practical quantitation limit, resulting in the SSL of 3.3 micrograms/liter [ug/L] (Table 4). The MTCA Method A cleanup level for arsenic of 5 ug/L is based on natural background concentrations and the state and federal drinking water standard maximum contaminant level (MC) for arsenic is 10 ug/L. Although not yet promulgated, Ecology has proposed a human-health surface water quality standard of 10 ug/L. Dissolved arsenic has been detected in reconnaissance groundwater samples collected from two locations, PP-1 and PP-2 at concentrations above 10 ug/L. None of the groundwater results for samples collected from properly constructed and developed monitoring wells contain dissolved arsenic above 10 ug/L.

Dissolved lead has been detected above the SSLs in groundwater samples collected from well MW-5 and boring PP-4 (Table 6). Benzene and total cPAHs were detected above the SSLs in groundwater samples collected from borings PP-4, PP-5, PP-6 and MW-6 (Figure 18). The nature and extent of dissolved lead, benzene and total cPAHs in groundwater has not been fully defined in the downgradient direction (to the south-southwest).

4.2.4. Surface Water

Results from surface water sampling in the drainage ditch along the west side of the Site indicate there are no SSL exceedances for Site COPCs, except for vinyl chloride detected in two separate surface water samples collected from the same location in December 2010 and January 2011 (Table 2). None of the surface water samples collected from the west drainage ditch in 2018 (locations SW-1, SW-2 and SW-3) had detections of the COPCs above laboratory reporting limits (Table 2). The surface water sample locations are depicted on Figure 4.



4.2.5. Soil Gas

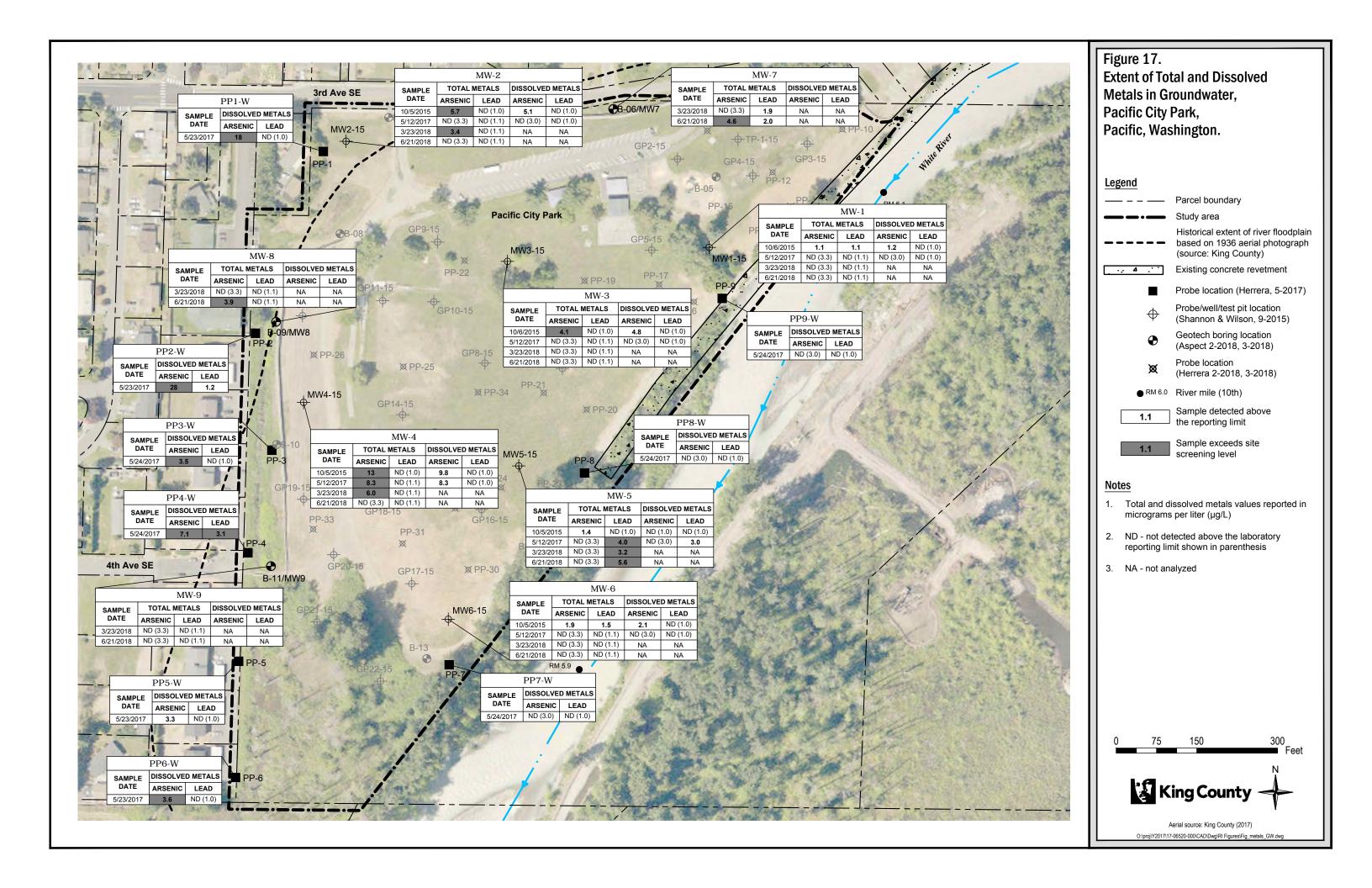
Based on the landfill gas sampling during the 1985 Abandoned Landfill Study and the fill gas monitoring conducted March 23, June 21, and September 23, 2018, during the RI at three monitoring wells (MW-1, MW-6, and MW-9), little to no methane was detected. See Figure 11 for sample locations.

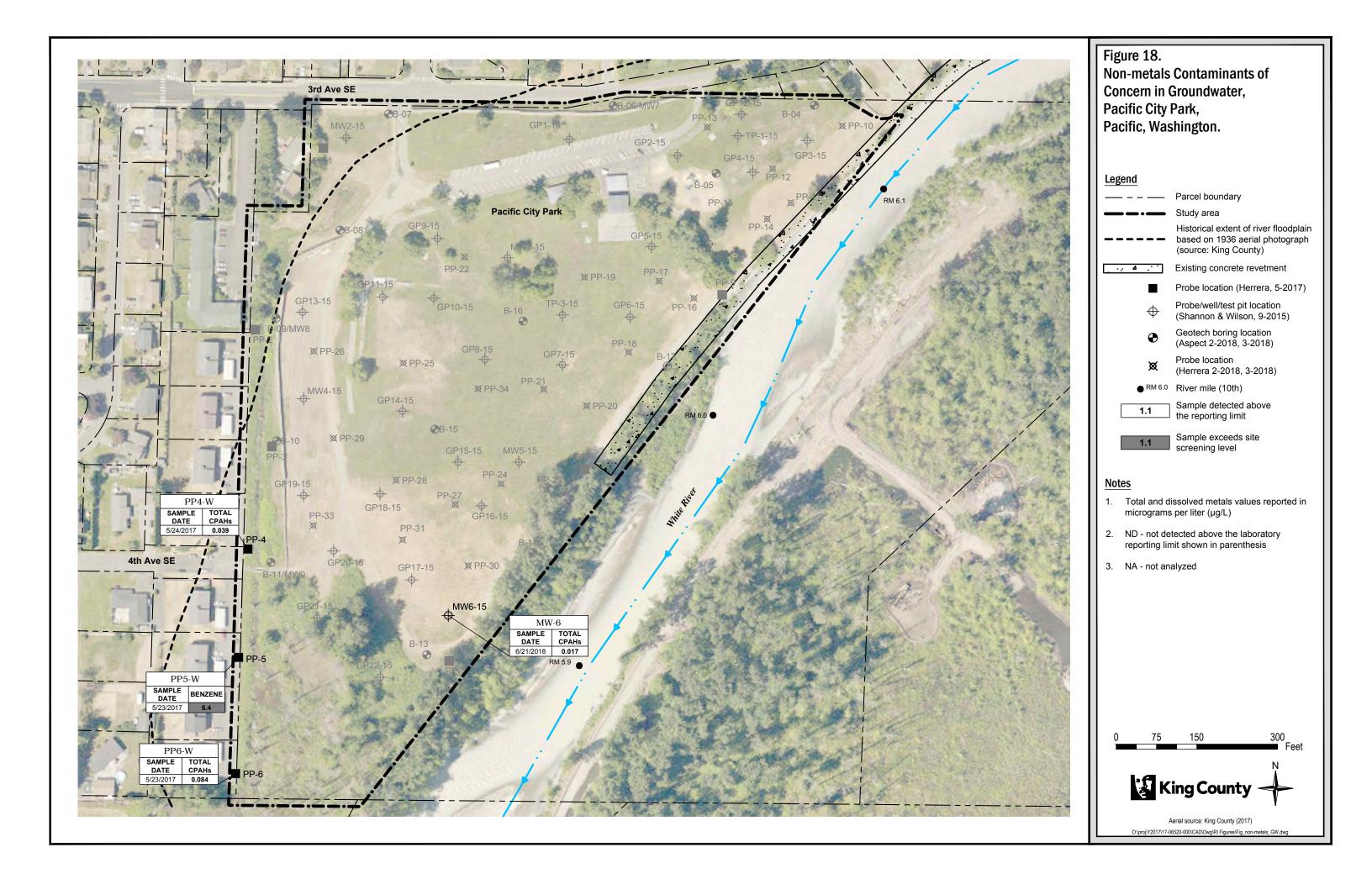
4.3. FATE AND TRANSPORT

Concentrations of several COPCs [including metals (primarily lead and arsenic), total cPAHs, and TPH (primarily lube oil)] exceed SSLs across the Site in native soil, fill soil, and fill soil mixed with refuse. In general, concentrations of COPCs in soil are highest in the east-central portion of the Site where refuse was disposed at thicknesses up to 20 feet bgs. The fill soil mixed with refuse also contains concentrations of other COPCs above the SSLs (chlorinated VOCs, SVOCs and pesticides); however, the presence of these chemicals is limited to a few soil samples and are probably associated with the quality of localized, discrete refuse. In native soil located beneath the refuse, the concentrations of COPCs are typically lower than what is reported in the fill soil with refuse and, the vertical extent is limited to less than 10 feet, suggesting that downward migration of COPCs slowed or ceased because of changes in density between fill and native soils, sorption characteristics, and/or the presence of groundwater.

Although concentrations of total cPAHs and lead are ubiquitous in fill soil and refuse across the Site, there are relatively few detections in groundwater, suggesting that total cPAHs and lead are not readily leaching from the fill soil or the refuse. However, the groundwater data collected at the south-southwest corner of the Site, which is downgradient of the refuse mass but still within the fill soil, suggests that there may be a low concentration, diffuse groundwater plume emanating from the refuse or there may be localized groundwater impacts attributable to variations in the fill soil quality. Groundwater likely discharges as surface water in the drainage ditch, at least seasonally. However, concentrations of total cPAHs, arsenic and lead have not been detected in surface water samples above laboratory reporting limits.

Very low methane concentrations were detected during the landfill gas monitoring for the 1985 Abandoned Landfill Study. No methane was detected during the landfill gas monitoring conducted March 23, June 21, and September 26, 2018 during the RI at monitoring wells (MW-1, MW-6 and/or MW-9). Benzene was detected in one groundwater sample from the southwest corner of the Site above the MTCA Method A cleanup level. However, based on the Site model, the benzene detection appears to be a localized anomaly and due to its proximity to adjacent property dumpsters, and is likely from a recent spill. The sampling and monitoring conducted to date does not indicate a potential exposure pathway from soil vapor to ambient and indoor air.





4.4. PRELIMINARY EXPOSURE ASSESSMENT

An exposure pathway describes the mechanisms by which human or ecological exposure to contaminants can occur assuming no remedial action or protective controls are in place. An exposure pathway is considered complete if a human or ecological receptor can be exposed to a contaminant via that pathway. A preliminary assessment of potential exposure pathways was completed to develop SSLs for evaluation of RI data, as provided in Section 3.2.1.1. All potential receptors and exposure pathways at the Site will be considered when developing the proposed cleanup standards as part of the Feasibility Study (FS). Based on the current and expected future use of the Site, a brief discussion of the receptors and exposure pathways that will be evaluated in the FS is provided below.

Protection of Human Health. The FS will be completed to ensure the protection of human health through the following potential exposure to receptors:

- 1. Direct contact by Site workers during construction and Site maintenance, and park visitors with soil, refuse, groundwater and surface water containing hazardous substances.
- 2. Ingestion of groundwater containing hazardous substances.

The primary COPCs detected above SSLs in soil and groundwater at the Site are total cPAHs and heavy metals, which do not pose a risk to human health via vapor intrusion and inhalation. Because soil at the Site is nearly continuously saturated, there is limited volatilization and preferential pathway.

Protection of Groundwater Discharge to Surface Water. The FS will consider the groundwater to surface water migration pathway to ensure the protection of beneficial use of surface water (human recreational, aquatic, and terrestrial receptors in the White River).

Protection of Terrestrial Ecological Receptors. WAC 173-340-7490 addresses procedures to be followed to ensure protection of terrestrial ecological receptors from exposure to contaminated soil. The FS will consider the protection of terrestrial ecological receptors.

5. PROPOSED CLEANUP STANDARDS

A cleanup standard includes both a cleanup level (chemical- and media-specific concentration of a contaminant that is protective of human health and the environment via all exposure pathways) and a point of compliance (the location where the cleanup level must be attained to achieve protectiveness). The preliminary cleanup levels and points of compliance for the Site are described in the following subsections. Final cleanup standards will be proposed in the draft Cleanup Action Plan for the Site and approved by Ecology.

5.1. CLEANUP LEVELS

Cleanup levels (CULs) are defined by regulatory numeric criteria (contaminant concentrations) that are protective of human health and the environment. Cleanup levels are contaminant-specific and media-specific and are only proposed for hazardous substances that exceed SSLs at the Site. The cleanup levels are used as the basis for developing media-specific remedial action objectives (RAOs) for the cleanup action. As described in Section 3.2.1.2, SSLs were developed for this RI for each hazardous substance that has historically been detected in soil, groundwater, or surface water at the Site.

Final proposed CULs for constituents that exceed SSLs in soil, groundwater and surface water will be defined in the FS and approved by Ecology as part of the development of the Cleanup Action Plan for the Site.

5.1.1. Soil

The preliminary soil cleanup levels are likely to be equivalent to the SSLs developed in 3.2.1.2.2 for the protection of human health and the environment, through complete exposure pathways, including the protection of leaching from soil to groundwater and surface water.

5.1.2. Groundwater and Surface Water

The preliminary groundwater and surface water cleanup levels for non-metal COPCs are likely to be equivalent to the SSLs developed in 3.2.1.2.2 for the protection of surface water. Because metals are naturally occurring in Washington soils and groundwater, the groundwater and surface water cleanup levels for metals may be adjusted based on reasonable and likely exposure scenarios and other relevant requirements. For example, the groundwater cleanup level for arsenic may be adjusted to the MTCA Method A groundwater cleanup level of 5 ug/L, which is based on background concentrations for the state of Washington, or the proposed state of Washington human-health surface water quality standard of 10 ug/L.



5.2. Points of Compliance

5.2.1. Soil

In accordance with MTCA, the point of compliance for direct contact with soil extends to 15 feet bgs, based on a reasonable maximum depth of excavation and assumed placement of excavated soils at the surface where contact occurs. For soil CULs based on leaching to groundwater and the protection of surface water, the soil point of compliance is all depths, above and below the water table.

5.2.2. Groundwater

Under MTCA, the standard point of compliance for groundwater CULs is throughout the site, regardless of whether groundwater is potable (WAC 173-340-720(8)(b)). If it is not practicable to meet groundwater CULs throughout the site within a reasonable restoration time frame, Ecology may approve a conditional point of compliance (WAC 173-340-720(8)(c)). Remedial alternatives will be developed and evaluated in the future FS assuming the standard point of compliance for groundwater.

5.3. TERRESTRIAL ECOLOGICAL EVALUATION

The purpose of a Terrestrial Ecological Evaluation (TEE) is to ensure the protection of land-based plants and animals from potential exposure to contaminated soil and to establish cleanup standards protective of terrestrial ecological receptors, if necessary (WAC 173-340-7490). The Site does not qualify for an exclusion from a TEE, nor does it qualify for a simplified TEE (WAC 173-340-7491 and -7492). To evaluate for the protection of ecological receptors, MTCA Ecological Indicator Soil Concentrations were included in the development of SSLs to evaluate the RI data. Per MTCA, a site-specific TEE will be conducted in consultation with Ecology and will facilitate selection of a cleanup action by developing information necessary to evaluate cleanup action alternatives in the FS.

6. SUMMARY, CONCLUSIONS, AND RECOMMENDATIONS

6.1. SUMMARY

The RI report has been prepared to meet the requirements of the MTCA and regulations implementing it, WAC 173-340, to provide the results of investigations completed to characterize the nature and extent of contamination at the Site. The RI Report has been prepared in general accordance with the Remedial Investigation Checklist Guidance (Ecology 2016).

The Site is located on the existing right bank of the White River in the City of Pacific, Washington, on property that was historically part of the river channel before the construction of a levee and concrete revetment in 1919. The portions of the property located landward of the levee were filled as a King County refuse dump, which was active between approximately 1921 and 1965. The results of investigation activities indicate that portions of the Site were filled with soil and portions of the Site were filled with a combination of soil and refuse. The presence of fill soil alone, in the absence of COPCs, does not fall under the definition of MTCA as a hazardous substance, and therefore, is not part of the MTCA Site. The Site is defined by any location where refuse is present or where COPCs are present in soil, groundwater, surface water or soil vapor at concentrations exceeding the SSLs. The current Site use includes a seasonal city park, undeveloped natural areas along the White River, and temporary flood protection (HESCO) barrier.

Because of the unknown nature of the fill soil and refuse, initial investigations conducted on the Site included a full analytical suite of COPCs, including TPH, metals, VOCs, SVOCs, PCBs, herbicides and pesticides. Subsequent phases of investigation focused on characterizing the nature and extent of the primary COPCs, which consist of TPH, metals, PCBs and total cPAHs. The results of investigative work indicate that the fill soil contains ubiquitous concentrations of lead and total cPAHs. In areas where refuse was known to have been dumped and/or was observed to be present in subsurface explorations, the results of soil characterization samples identify higher concentrations of total cPAHs and lead, as well as the frequent detection of TPH and PCBs and rare detections of chlorinated VOCs, other SVOCs and pesticides. Laterally, the extent of COPCs in soil are generally defined by the extent of fill. Vertically, the COPCs are present throughout the fill and extend up to 10 feet into the underlying native, alluvial deposits.

The COPCs detected in groundwater include arsenic, lead, total cPAHs and benzene. Arsenic has been detected in groundwater samples collected from across the Site, is naturally occurring in soil and groundwater in Washington state and is likely not present in Site groundwater at concentrations that warrant remedial action. Lead, total cPAHs and benzene have been detected

King County

above the SSLs in groundwater samples collected to the south-southwest of the refuse, which is hydraulically downgradient based on water level elevations in monitoring wells on the Site. The groundwater data suggest that there may be a low concentration, seasonal and/or diffuse groundwater plume emanating from the refuse or there may be localized groundwater impacts attributable to variations in the fill soil quality. Although groundwater is assumed to discharge to surface water in some areas of the Site, the results of surface water sampling do not indicate the migration of COPCs in groundwater to surface water.

6.2. Conclusions and Recommendations

The presence of COPCs in soil generally corresponds to the locations of historical placement of fill and refuse (see Figure 16). A remedial action for the Site is likely to include components of removal, treatment and/or containment to mitigate potential risks to human health and the environment associated with the presence of hazardous substances in fill soil and refuse. Sufficient information regarding the extent and quality of the fill and refuse, and the associated groundwater impacts, has been collected to allow for the development and evaluation of remedial alternatives, with one exception. Additional investigation is warranted to further evaluate the extent of fill soil and the presence of COPCs in soil and groundwater to the south-southwest of the property boundary. This would include installing three additional groundwater monitoring wells, one on 4th Avenue SE, just west of probe location PP-4 and MW-9, one to the west of probe location PP-5, and one to the southwest of probe location PP-6. In addition, three explorations are planned to assess soil conditions along the parcel boundary by the apartments immediately west of the Park. The proposed additional explorations are shown in Figure 16.

Quarterly groundwater monitoring of the Site monitoring wells is planned to continue to evaluate seasonal fluctuations in water levels and variability in groundwater quality. The CSM will be revised as additional information is collected.

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TABLES



Table 1. Summary of Water Level Elevation Data from Monitoring Wells, Pacific City Park Remedial Investigation, Pacific, Washington.

Monitoring Well Identification	Measurement Date	Reference Elevation (feet) ^a	Depth to Water (feet)	Water Level Elevation (feet)
	5/12/17		2.33	80.83
	3/23/18	02.46	2.84	80.32
MW-1	6/21/18	83.16	3.12	80.04
	9/26/18		5.8	77.36
	5/12/17		1.37	78.48
N 41 A 2	3/23/18	70.05	1.81	78.04
MW-2	6/21/18	79.85	2.32	77.53
	9/26/18		3.68	76.17
	5/12/17		0.40	79.61
	3/23/18	00.4	0.55	79.46
MW-3	6/21/18	80.1	1.27	78.74
	9/26/18		3.01	77.09
	5/12/17		2.73	78.41
	3/23/18	00.14	3.09	77.05
MW-4	6/21/18	80.14	3.53	76.61
	9/26/18		4.54	75.6
	5/12/17		1.60	79.80
	3/23/18	81.4	2.26	79.14
MW-5	6/21/18		2.38	79.02
	9/26/18		4.8	76.6
	5/12/17		5.71	78.10
NAM 6	3/23/18	02.01	6.65	77.16
MW-6	6/21/18	83.81	6.60	77.21
	9/26/18		8.53	75.28
	3/23/18		0.32	79.50
MW-7 ^b	6/21/18	79.82	0.78	79.04
	9/26/18		2.68	77.14
	3/23/18		2.63	77.32
MW-8 ^b	6/21/18	79.95	3.12	76.83
	9/26/18		4.2	75.75
	3/23/18		5.85	76.74
MW-9 ^b	6/21/18	82.59	6.02	76.57
	9/26/18		6.98	75.61
	3/23/18		5.52	80.60
B-03 ^c	6/21/18	86.12	5.41	80.71
	9/26/18		8.3	77.82

^a Reference elevation is the top of protective casing (North American Vertical Datum 1988 [NAVD 88])

^b MW-7, MW-8, and MW-9 correspond to probe borings B-06, B-09, and B-11, respectively.

^c Standing water level measurement only at boring location B-03. No groundwater sample was collected.

			•	Γable 2. S	Summary	of Surface	Water S	Sample R	esults, Pac	ific City Pa	rk Remed	ial Investi	gation, Pa	cific, Was	shington.					
									Sa	mple Identif	ication									
	,	WRLEV1-Dra	ainage Ditc	h	WR	RLEV2-Upstr	eam Wetla	nd	WR	LEV3-Downs	tream Wetla	and	SV	W1	SV	V2	SI	N3	SW4	
Parameter	12/8/10	1/4/11	2/28/11	3/10/11	12/8/10	1/4/11	2/28/11	3/10/11	12/8/10	1/4/11	2/28/11	3/10/11	6/29/18	10/9/18	6/29/18	10/9/18	6/29/18	10/9/18	10/9/18	SSLa
Field Parameters																				
Temp (°C)	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	14.2	12.2	14.7	12.3	14.3	12.9	12.9	NA
DO (mg/L)	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	1.62	1.2	1.84	0.36	1.12	0.8	0.65	NA
Cond (µS/cm)	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	153.7	204	170.1	208	177.5	151	179	NA
pH (std units)	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	6.73	6.84	6.84	6.74	6.94	6.65	6.69	NA
Turbidity (NTU)	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	13.72	Clear	10.76	Clear	12.33	Clear	Clear	NA
Conventional Paramete	ers (mg/L)																			
TKN	1.75	1.58	0.997	0.772	0.31	3.23	2.54	0.202	1.48	1.63	0.954	1.04	NA	NA	NA	NA	NA	NA	NA	NA
Nitrate+Nitrite Nitrogen	ND (0.01)	ND (0.01)	ND (0.04)	ND (0.04)	ND (0.01)	ND (0.01)	ND (0.01)	ND (0.01)	ND (0.01)	ND (0.01)	ND (0.04)	ND (0.04)	NA	NA	NA	NA	NA	NA	NA	NA
TP	0.136	0.134	0.132	0.134	0.115	0.142	0.070	0.035	0.162	0.205	0.170	0.196	NA	NA	NA	NA	NA	NA	NA	NA
ОР	ND (0.02)	0.049	0.026	0.086	0.007	0.016	0.024	0.025	ND (0.005)	0.014	0.024	0.060	NA	NA	NA	NA	NA	NA	NA	NA
Hardness (mg CaCO3/L)	125	141	86.6	64.6	47.8	39	22	26.9	113	125	90.5	88	66	62	71	51	74	51	45	NA
Petroleum Hydrocarbo	ons (µg/L)																			
GRO	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	ND (100)	ND (100)	ND (100)	ND (100)	ND (100)	ND (100)	ND (100)	1,000
DRO	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	ND (250)	ND (250)	ND (260)	ND (260)	ND (260)	ND (260)	ND (250)	
Lube Oil RO	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	ND (410)	ND (410)	ND (420)	ND (420)	ND (420)	ND (410)	ND (400)	500
Volatile Organic Comp	ounds (µg/	L)		•									•							
Benzene	ND (0.2)	ND (0.2)	NA	NA	ND (0.2)	ND (0.2)	NA	NA	ND (0.2)	ND (0.2)	NA	NA	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	0.44
Toluene	ND (0.2)	ND (0.2)	NA	NA	0.551	0.5	NA	NA	ND (0.2)	ND (0.2)	NA	NA	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	1.1	ND (1.0)	57
Ethylbenzene	ND (0.2)	ND (0.2)	NA	NA	ND (0.2)	ND (0.2)	NA	NA	ND (0.2)	ND (0.2)	NA	NA	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	29
Xylenes	ND (0.2)	ND (0.2)	NA	NA	ND (0.2)	ND (0.2)	NA	NA	ND (0.2)	ND (0.2)	NA	NA	ND (0.40)	ND (0.40)	ND (0.40)	ND (0.40)	ND (0.40)	ND (0.40)	ND (0.40)	1,000
Acetone	ND (2)	ND (2)	NA	NA	ND (2)	ND (2)	NA	NA	ND (4)	ND (2)	NA	NA	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	7,200
(cis)1,2-Dichloroethene	ND (0.2)	ND (0.2)	NA	NA	ND (0.2)	ND (0.2)	NA	NA	ND (0.2)	ND (0.2)	NA	NA	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	16
Chlorobenzene	ND (0.2)	ND (0.2)	NA	NA	ND (0.2)	ND (0.2)	NA	NA	ND (0.2)	ND (0.2)	NA	NA	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	100
Vinyl Chloride	0.0308	0.0371	NA	NA	ND (0.01)	ND (0.01)	NA	NA	ND (0.01)	ND (0.02)	NA	NA	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	0.02
Total Metals (µg/L)		_							_			_					_			_
Arsenic	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	ND (3.3)	ND (3.3)	ND (3.3)	ND (3.3)	ND (3.3)	ND (3.3)	ND (3.3)	3.3
Cadmium	ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	ND (4.4)	ND (4.4)	ND (4.4)	ND (4.4)	ND (4.4)	ND (4.4)	ND (4.4)	4.4
Calcium	37,900	42,600	25,600	18,700	14,300	11,900	6,450	7,790	35,400	39,900	28,500	26,300	NA	NA	NA	NA	NA	NA	NA	NA
Chromium	ND (1)	1.08	ND (1)	ND (1)	0.32	ND (1)	ND (1)	ND (0.2)	ND (1)	ND (1)	ND (1)	ND (1)	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	74
Copper	ND (0.4)	ND (0.4)	ND (2)	ND (2)	ND (0.4)	ND (2)	ND (2)	ND (2)	ND (0.4)	ND (0.4)	ND (2)	ND (2)	NA	NA	NA	NA	NA	NA	NA	1,300
Iron	NA	NA	NA	6,850	NA	NA	NA	2,240	NA	NA	NA	8,580	NA	NA	NA	NA	NA	NA	NA	1,000
Lead	ND (0.1)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.1)	ND (0.5)	ND (0.5)	ND (0.5)	ND (1.1)	ND (1.1)	ND (1.1)	ND (1.1)	ND (1.1)	ND (1.1)	ND (1.1)	2.5
Magnesium	7,370	8,500	5,520	4,330	2,930	2,280	1,440	1,820	6,090	6,200	4,700	5,390	NA	NA	NA	NA	NA	NA	NA	NA
Mercury	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	0.50

		Tab	le 2 (co	ntinued)	. Summar	y of Surfac	e Water	Sample	Results, P	acific City	Park Re	medial I	nvestigat	ion, Pacif	ic, Washin	gton.				
									Sam	ple Identific	ation									
	W	RLEV1-Drain	age Ditch		WRL	.EV2-Upstrea	m Wetlan	d	WRLE	V3-Downstre	eam Wetla	and	SV	V1	SI	W2	S	W3	SW4	1
Parameter	12/8/10	1/4/11	2/28/11	3/10/11	12/8/10	1/4/11	2/28/11	3/10/11	12/8/10	1/4/11	2/28/11	3/10/11	6/29/18	10/9/18	6/29/18	10/9/18	6/29/18	10/9/18	10/9/18	SSLª
Dissolved Metals (µg/L)																				
Chromium	ND (1)	1.12	ND (1)	ND (1)	ND (1)	ND (1)	ND (1)	ND (0.2)	ND (1)	ND (1)	ND (1)	ND (1)	NA	NA	NA	NA	NA	NA	NA	74
Copper	ND (0.4)	ND (2)	ND (2)	ND (2)	ND (0.4)	ND (0.4)	ND (0.4)	ND (0.4)	ND (0.4)	ND (0.4)	ND (0.4)	ND (0.4)	NA	NA	NA	NA	NA	NA	NA	1,300
Lead	ND (0.1)	ND (0.1)	ND (0.5)	ND (0.1)	ND (0.1)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.1)	ND (0.1)	ND (0.1)	ND (0.1)	NA	NA	NA	NA	NA	NA	NA	2.5
Zinc	6.72	4.89	3.42	3.17	ND (2.5)	ND (2.5)	ND (0.5)	2.82	2.5	ND (2.5)	ND (2.5)	ND (2.5)	NA	NA	NA	NA	NA	NA	NA	50
Carcinogenic Polycyclic Aro	matic Hydro	carbons (µg/l	L)																	
Benzo(a) anthracene	ND (0.0094)	ND (0.0096)	NA	NA	ND (0.0094)	ND (0.0095)	NA	NA	ND (0.0094)	ND (0.0095)	NA	NA	ND (0.011)	ND (0.011)	ND (0.010)	ND (0.0099)	ND (0.010)	ND (0.0098)	ND (0.010)	0.01
Benzo(b) fluoranthene	ND (0.0094)	ND (0.0096)	NA	NA	ND (0.0094)	ND (0.0095)	NA	NA	ND (0.0094)	ND (0.0095)	NA	NA	ND (0.011)	ND (0.011)	ND (0.010)	ND (0.0099)	ND (0.010)	ND (0.0098)	ND (0.010)	0.01
Benzo(j,k) fluoranthene	ND (0.0094)	ND (0.0096)	NA	NA	ND (0.0094)	ND (0.0095)	NA	NA	ND (0.0094)	ND (0.0095)	NA	NA	ND (0.011)	ND (0.011)	ND (0.010)	ND (0.0099)	ND (0.010)	ND (0.0098)	ND (0.010)	0.01
Benzo(a) pyrene	ND (0.0094)	ND (0.0096)	NA	NA	ND (0.0094)	ND (0.0095)	NA	NA	ND (0.0094)	ND (0.0095)	NA	NA	ND (0.011)	ND (0.011)	ND (0.010)	ND (0.0099)	ND (0.010)	ND (0.0098)	ND (0.010)	0.01
Chrysene	ND (0.0094)	ND (0.0096)	NA	NA	ND (0.0094)	ND (0.0095)	NA	NA	ND (0.0094)	ND (0.0095)	NA	NA	ND (0.011)	ND (0.011)	ND (0.010)	ND (0.0099)	ND (0.010)	ND (0.0098)	ND (0.010)	0.016
Indeno(1,2,3-cd)pyrene	ND (0.0094)	ND (0.0096)	NA	NA	ND (0.0094)	ND (0.0095)	NA	NA	ND (0.0094)	ND (0.0095)	NA	NA	ND (0.011)	ND (0.011)	ND (0.010)	ND (0.0099)	ND (0.010)	ND (0.0098)	ND (0.010)	0.01
Dibenz(a,h) anthracene	ND (0.0094)	ND (0.0096)	NA	NA	ND (0.0094)	ND (0.0095)	NA	NA	ND (0.0094)	ND (0.0095)	NA	NA	ND (0.011)	ND (0.011)	ND (0.010)	ND (0.0099)	ND (0.010)	ND (0.0098)	ND (0.010)	0.01
Total cPAHs TEQb	ND (0.0071)	ND (0.0072)	NA	NA	ND (0.0071)	ND (0.0072)	NA	NA	ND (0.0071)	ND (0.0072)	NA	NA	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.007)	ND (0.008)	ND (0.007)	ND (0.008)	0.085
Semivolatile Organic Comp	ounds (µg/L)																			
2-Methylphenol	ND (0.024)	ND (0.024)	NA	NA	0.539	0.307	NA	NA	0.113	0.0574	NA	NA	NA		NA		NA			NA
4-Methylphenol	ND (0.047)	ND (0.048)	NA	NA	1.55	0.197	NA	NA	ND (0.047)	ND (0.048)	NA	NA	NA		NA		NA			NA
Acenaphthene	ND (0.0189)	ND (0.0096)	NA	NA	ND (0.0094)	ND (0.0095)	NA	NA	ND (0.0189)	ND (0.0095)	NA	NA	NA		NA		NA			30
Acenaphthylene	ND (0.0094)	ND (0.0096)	NA	NA	ND (0.0094)	ND (0.0095)	NA	NA	ND (0.0094)	ND (0.0095)	NA	NA	NA		NA		NA			NA
Benzoic Acid	1	1.9	NA	NA	1.75	0.909	NA	NA	ND (0.472)	1.55	NA	NA	NA		NA		NA			NA
Benzyl Alcohol	ND (0.094)	ND (0.096)	NA	NA	0.263	0.21	NA	NA	0.225	ND (0.095)	NA	NA	NA		NA		NA			NA
Butyl Benzyl Phthalate	0.102	ND (0.048)	NA	NA	ND (0.047)	0.092	NA	NA	ND (0.047)	ND (0.048)	NA	NA	NA		NA		NA			1.0
Bis(2-Ethylhexyl) phthalate	0.99	NS (0.26)	NA	NA	0.488	0.23	NA	NA	ND (0.472)	1.77	NA	NA	NA		NA		NA			1.0
Diethyl Phthalate	ND (0.024)	ND (0.024)	NA	NA	0.026	ND (0.024)	NA	NA	ND (0.472)	ND (0.024)	NA	NA	NA		NA		NA			NA
Di-n-butyl Phthalate	0.176	0.14	NA	NA	0.134	0.14	NA	NA	0.149	0.126	NA	NA	NA		NA		NA			8
Naphthalene	ND (0.0189)	0.0241	NA	NA	ND (0.0094)	0.013	NA	NA	0.0231	0.0338	NA	NA	NA		NA		NA			4,710

Bold values detected above the reporting limit

Shaded values exceed the site screening level

 μ g/L = micrograms per liter

NA = not analyzed

ND = not detected above laboratory reporting limits shown in parentheses

SSL = site screening levels



^a Refer to Table 7 "Proposed Site Screening Levels for Groundwater and Surface Water" for notes on how each screening level was selected.

b Total carcinogenic polycyclic aromatic hydrocarbon (cPAHs) toxic equivalency (TEQ) concentration was calculated using one-half the reporting limit for compounds that were not detected above the reporting limit.

mg/L = milligrams per liter

	1				stigation, l		n Push Prol Ishington.			
				Sa	ample Location	on				Site
Analytical Parameter	PP1-W	PP2-W	PP3-W	PP4-W	PP5-W	PP6-W	PP7-W	PP8-W	PP9-W	Screening Level ^a
Sample Date				5/23	/2017–5/24/	2017				Levei ^ω (μg/L)
NWTPH-Gx (μg/L)	l .									4 5- 7
Gasoline Range Organics	ND (100)	ND (100)	ND (100)	ND (100)	210	ND (100)	ND (100)	ND (100)	ND (100)	1,000
NWTPH-Dx (mg/L)	140 (100)	140 (100)	145 (100)	140 (100)		140 (100)	140 (100)	140 (100)	140 (100)	1,000
Diesel Range Organics	ND (0.27)	ND (0.37)	ND (0.26)	ND (0.26)	ND (0.27)	ND (0.26)	ND (0.28)	ND (0.26)	ND (0.28)	500
Lube Oil	ND (0.27)	ND (0.60)	ND (0.20)	ND (0.20)	ND (0.43)	ND (0.42)	ND (0.28)	ND (0.20)	ND (0.28)	500
	. ,	ND (0.00)	ND (0.42)	ND (0.41)	ND (0.43)	ND (0.42)	ND (0.44)	ND (0.41)	ND (0.43)	300
Volatile Organic Compou		7-				7.0			7.0	7200
Acetone	6.7	75	5.3	5.6	5.5	7.9	6.1	5.2	7.2	7200
Carbon disulfide	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	0.22	800
Benzene	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	6.4	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	0.44
Toluene	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	5.2	ND (1.0)	ND (1.0)	57
Chlorobenzene	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	0.43	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	100
Xylenes	ND (0.40)	ND (0.40)	ND (0.40)	ND (0.40)	32.57	ND (0.40)	ND (0.40)	ND (0.40)	ND (0.40)	1,000
Isopropylbenzene	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	0.45	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	NA ^c
n-Propylbenzene	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	0.95	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	800
1,3,5-Trimethylbenzene	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	3.1	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	80
1,2,4-Trimethylbenzene	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	12	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	NA ^c
Total Metals (µg/L)		_	_			_	_			_
Antimony	ND (5.6)	ND (5.6)	ND (5.6)	47	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	5.6
Arsenic	36	910	18	110	14	14	110	ND (3.3)	30	3.3
Beryllium	ND (11)	34	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	4
Cadmium	ND (4.4)	24	ND (4.4)	ND (4.4)	ND (4.4)	ND (4.4)	ND (4.4)	ND (4.4)	ND (4.4)	4.4
Chromium	73	2,000	63	240	29	45	210	ND (11)	49	50
Copper	80	4,200	120	460	50	54	570	ND (11)	110	640
Lead	250	2,100	54	2,800	32	55	250	ND (1.1)	61	2.5
Mercury	ND (0.50)	4.3	ND (0.50)	1.2	ND (0.50)	ND (0.50)	0.68	ND (0.50)	ND (0.50)	0.5
Nickel	33	1,800	32	190	ND (22)	46	210	ND (22)	44	100
Selenium	ND (5.6)	53	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	6.0	ND (5.6)	ND (5.6)	50
Silver	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	10
Thallium	ND (5.6)	9.6	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	0.16
Zinc	700	5,500	130	1,000	49	100	520	71	210	120
Dissolved Metals (μg/L)	I		I		<u> </u>	I	I	I	I	
Antimony	ND (5.0)	ND (5.0)	ND (5.0)	7.1	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	5.6
Arsenic	18	28	3.5	7.1	3.3	3.6	ND (3.0)	ND (3.0)	ND (3.0)	3.3
Beryllium	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	273
Cadmium	ND (4.0)	ND (4.0)	ND (4.0)	ND (4.0)	ND (4.0)	ND (4.0)	ND (4.0)	ND (4.0)	ND (4.0)	4.4
Chromium	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	74
Copper	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	1,300
Lead	ND (1.0)	1.2	ND (1.0)	3.1	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	2.5
Mercury	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	0.50
Nickel	ND (20)	ND (20)	ND (20)	ND (20)	ND (20)	ND (20)	ND (20)	ND (20)	ND (20)	52
Selenium	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	5.6
Silver	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	10
Thallium	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	0.216
Zinc	39	ND (25)	ND (25)	ND (25)	ND (25)	ND (25)	ND (25)	62	ND (25)	120
Polycyclic Aromatic Hydro			I		T	I	ı	I	T	
Naphthalene	ND (0.0096)	ND (0.12)	ND (0.0096)	ND (0.10)	0.18	ND (0.097)		ND (0.0095)		160
Benzo(a)anthracene	ND (0.0096)	ND (0.012)	ND (0.0096)	0.013	ND (0.0097)	0.052	` '	`	ND (0.0096)	0.01
Chrysene	ND (0.0096)	0.012	ND (0.0096)	0.021	ND (0.0097)	0.080	` '	ND (0.0095)	` '	0.016
Benzo(b)fluoranthene	ND (0.0096)	0.020	ND (0.0096)	0.036	ND (0.0097)	0.14		ND (0.0095)	0.019	0.01
Benzo(j,k)fluoranthene	ND (0.0096)	ND (0.012)	ND (0.0096)	0.011	ND (0.0097)	0.042		ND (0.0095)		0.01
Benzo(a)pyrene	ND (0.0096)	ND (0.012)	ND (0.0096)	0.024	ND (0.0097)	0.051		ND (0.0095)	0.010	0.01
Indeno(1,2,3-cd)pyrene	ND (0.0096)	0.014	ND (0.0096)	0.040	ND (0.0097)	0.066	` '	ND (0.0095)	0.011	0.01
Dibenz(a,h)anthracene	ND (0.0096)	ND (0.012)	ND (0.0096)	0.011	ND (0.0097)	0.018		ND (0.0095)		0.01
Benzo(g,h,i)perylene	ND (0.0096)	0.020	ND (0.0096)	0.046	ND (0.0097)	0.055	ND (0.0097)	ND (0.0095)	0.012	NAc
Total cPAHs TEQ ^b	ND (0.0072)	0.011	ND (0.0072)	0.039	ND (0.0073)	0.084	ND (0.0073)		0.014	0.015

BOLD values detected above the reporting limit.

Shaded values exceed the site screening level

 μ g/L = micrograms per Liter

ND = not detected above laboratory reporting limits (shown in parentheses)

PP – push probe

a Refer to Table 7 "Proposed Site Screening Levels for Groundwater and Surface Water" for notes on how each screening level was selected.

b Total carcinogenic polycyclic aromatic hydrocarbons (cPAHs) toxic equivalency (TEQ) concentration was calculated using one-half the reporting limit for compounds that were not detected above the reporting limit.

 $^{^{\}rm c}$ $\,$ No Site screening level established for this parameter

	Protection of Human Health – Direct Contact	Listed MTCA	Protection of Groundwater (saturated soil)	MTCA E	tion of Ed Receptor cological Concentra	rs Indicator			
Constituent (by group) ^a	MTCA Method A/B (Unrestricted Land Use) ^c	Method B Protective of Groundwater (Saturated	MTCA Method A/ Calculated Method B ^d	Plants	Soil Biota	Wildlife	Natural Background Metals Concentrations ^e	Practical Quantitati on Limit (PQL)	
Petroleum Hydrocarbons		((- \(- \)	
Gasoline Range Organics									
(no benzene)			100		100	5,000		5	100
Diesel Range Organics			2,000		200	6,000		25	200
Lube Oil Range Organics			2,000					50	2,000
Metals									
Arsenic	20	0.146	0.097	10	60	132	20	10	20
Barium	16,000	82.6	41.3	500		102		2.5	41.3
Cadmium			0.031	4	20	14	1	0.5	1
Chromium		24,000	50	42	42	67	48	0.5	48
Lead	250	150	25	50	500	118	24	5	25
Mercury		0.105	0.026	0.3	0.1	5.5	0.07	0.025	0.07
Selenium	400	0.264	0.030	1	70	0.3	0.78	10	10
Silver	400	0.687	0.086	2			0.61	0.5	0.61
Volatile Organic Compounds (VOCs)	1							
Benzene		0.00174	0.0002					0.001	0.001
Toluene		0.273	0.024	200				0.005	0.024
Ethylbenzene		0.343	0.014					0.001	0.014
Xylenes		0.831	0.52					0.002	0.52
Acetone	72,000	2.07	2.07					0.005	2.07
2-Butanone (methyl ethyl ketone)	48,000	0.266	1.38					0.005	1.38
Carbon Disulfide	8,000	0.266	0.27					0.001	0.27
Cis-1,2-Dichloroethene Chlorobenzene	160	0.00515	0.005		40			0.001	0.005
Methylene Chloride	1,600 0.02	0.0511 0.00148	0.051 0.001		40			0.001	0.051 0.005
p-Isopropyltoluene (cumene)	8,000	0.00148	0.001					0.003	0.229
Styrene	16,000	0.12	0.120	300				0.001	0.120
Tetrachloroethene (PCE)	0.05	0.00276	0.0013	300				0.001	0.0013
Trichloroethene (TCE)	0.03	0.00152	0.0013					0.001	0.001
1,2,4-Trimethylbenzene			0.0002					0.001	
Semi-Volatile Organic Compounds	(SVOCs)								
4-Nitrophenol								0.033	
Benzyl Alcohol								0.017	
Bis(2-ethylhexyl)phthalate	71.4	0.668	0.111					0.033	0.111
Butyl benzylphthalate	526	0.646	0.0140					0.033	0.033
Di-n-butyl Phthalate	8,000	2.97	0.015					0.17	0.17
Di-n-octyl Phthalate	800	13,300	13,312					0.033	800
Naphthalenes ^g	5	0.236	0.236					0.033	0.236
Pentachlorophenol	2.5	0.000878	0.004	3	6	4.5		0.17	0.17
Phenol	24,000	0.757	0.757	70	30			0.033	0.757
p-Cresol	8,000							0.033	8,000
Polycyclic Aromatic Hydrocarbons	(PAHs)								T
Acenaphthene	4,800	4.98	0.156	20				0.0067	0.156
Acenaphthylene								0.0067	
Anthracene	24,000	114	7.134					0.0067	7.134
Benzo(a)anthracene	1.37	0.043	0.004					0.0067	0.0067
Benzo(a)pyrene	0.1	0.116	0.010			12		0.0067	0.010
Benzo(b)fluoranthene	1.37	0.147	0.012					0.0067	0.012
Benzo(j,k)fluoranthene	13.7	1.47	0.012					0.0067	0.012
Chrysene	137	4.77	0.0064					0.0067	0.0067
Dibenz(a,h)anthracene	0.137	0.021	0.018					0.0067	0.018
Fluoranthene	3,200	31.6	0.296		22			0.0067	0.296
Fluorene	3,200	5.12	0.080		30			0.0067	0.080
Indeno(1,2,3-cd)pyrene	1.37	0.416	0.035					0.0067	0.035
Phenanthrene	400	0.000878	0.000					0.0067	0.0067
Pyrene Total cPAHs TEO	2,400	32.8	0.546			12		0.0067	0.546
Total cPAHs TEQ Polychlorinated Biphenyls (PCBs)	0.1	0.0214	0.090			12		0.0101	0.020
PRIVEDIATION RIPHONVIE (DCR)									

	Protection of Human Health – Direct Contact	Listed MTCA	Protection of Groundwater (saturated soil)	MTCA	ction of Eco Receptors Ecological Concentra	s Indicator			
Constituent (by group) ^a	MTCA Method A/B (Unrestricted Land Use) ^c	Method B Protective of Groundwater (Saturated)	MTCA Method A/ Calculated Method B ^d	Plants	Soil Biota Wildlin		Natural Background Metals Concentrations ^e	Practical Quantitati on Limit (PQL)	Proposed Site Soil Screening Level ^t
Herbicides									
2,4-DB	640							0.095	640
2,4,5-T	800							0.095	800
Bentazon	2,400								2,400
Chloramben	1,200								1,200
Dacthal (chlorthal-dimethyl)	800								800
Dalapon	2,400							2.3	2,400
Dinoseb	80							0.095	80
Picloram	5,600								5,600
Silvex (2,4,5-TP)	640							0.095	640
Organochlorine Pesticides									
4,4'-DDD ^h	4.17	0.0168	0.0002			0.75		0.01	0.01
4,4'-DDE ⁱ	2.94	0.0223	0.0004			0.75		0.01	0.01
Cis-chlordane ^j	2.86	0.103	0.0003		1	2.7		0.01	0.01
Endosulfan I ^k	480	0.103	0.0001					0.005	0.005
Endosulfan II ^k	480	0.223	0.0001					0.01	0.01
Endosulfan sulfate		0.223						0.01	
Gamma-chlordane ^j	2.86		0.0003					0.01	0.01
Methoxychlor	400	3.21	0.0016					0.01	0.01

All units in milligrams per kilogram (mg/kg), unless otherwise stated.

Blank cells are intentional and indicate that criteria are not available.

Shaded cells denote the most restrictive criteria that is proposed as the site screening level.

- ^a Cleanup levels have been developed for only those compounds that have been detected at the Site above laboratory reporting limits in any media.
- b Ecological Indicator Soil Concentrations for Protection of Terrestrial Plants and Animals, lowest value of concentrations established protective of plants, soil biota and wildlife (WAC 173-340-7493).
- ^c Model Toxics Control Act (MTCA) Method A or Method B Soil Cleanup Levels for Unrestricted Land Uses (WAC 173-340) for the protection of human health through direct contact.
- d MTCA Method A for petroleum hydrocarbons, calculated Method B for the protection of groundwater discharging to surface water for all other compounds.
- e Natural background values for metals from Natural Background Soil Metals Concentrations in Washington State (Ecology 1994), except arsenic which is from MTCA (WAC 173-340-900, Table 740-1).
- f Proposed site screening level for soil is the most restrictive of MTCA Method A/B and Ecological Indicator Soil Concentration Screening Levels, adjusted for natural background and laboratory POLs.
- ^g Screening level for naphthalenes is a total value for naphthalene, 1-methyl naphthalene, and 2-methyl naphthalene.
- ^h Screening levels for DDD are used for 4,4'-DDD.
- i Screening levels for DDE are used for 4,4'-DDE.
- ^j Screening levels for chlordane are used for cis-chlordane and gamma-chlordane.
- k Screening levels for endosulfan are used for endosulfan I and endosulfan II.

	Table 5. Propo		c.ccimig Ec			541						ga 11011,		g.			
						I			of Surface	Water							
	Protection o as Drinking W			EPA 201	.6 CWA –			Recommend uality Criteria		,	WAC 173-	201A Table	240 ^d	MTCA			
				Criteria Ap Washi	uman Health oplicable to ngton: ^b	Aquat	ction of ic Life –	Protect Human	Health:	Aquat	ction of ic Life –	Human	ction of Health:	Method B Standarde	Practical	Proposed Site	Proposed Site Surface
Constituent (by group) ^a	MTCA Method A/B (unrestricted land use) ^f	Federal MCL ^g	Washington State MCL ^g		Organisms Only	Acute	Chronic	Consum Water and Organisms			Chronic		Organisms Only	Protection of Human Health	Quantitation Limit (PQL)	Groundwater Screening Level ^h	Water Screening Level ⁱ
Total Petroleum Hydrocarbons	(umesuretea iana ase)		June 11102	- Cigamonis	· • • • • • • • • • • • • • • • • • • •	710000	- Cili Cilic	C gamons	<u> </u>	710010		Cigamonio	· · · · · ·		(· ~-)		
Gasoline Range Organics	1,000		T												100	1,000	
Diesel Range Organics	500														250	500	
Lube Oil	500														400	500	
Metals	300				1					ı		1	I		100	1 300	
antimony	6.4	6	6	6	90			5.6	640			12	180	1040	5.0	5.6	5.6
Arsenic	5	10	10	0.018	0.14	340	150	0.018	0.14	360	190	10	100	0.0982	3.3	3.3	3.3
Barium	3,200	2,000	2,000	0.010	0.14	340	130	1,000	0.14	300	150	10	10	0.0302	28	1,000	1,000
beryllium	32	4	4					1,000						273	4	4	273
Cadmium	5	5	5			1.8	0.72							273	4.4	4.4	4.4
Calcium (stormwater only)	j					2.0	0.72								1,000	7	7. 7
Chromium	50	100	100			570	74							243,000	10	50	74
Copper	640	1,300	1,300	1,300				1,300				1,300		2,880	11	640	1,300
Iron (stormwater only)		,	,	,			1,000	,				,		,	10		1,000
Lead	15	15	15			65	2.5								1.1	2.5	2.5
Magnesium (stormwater only)															1,000		
Mercury	2	2	2			1.4	0.77	0.14	0.15	2.1	0.012	0.14	0.15		0.50	0.50	0.50
Nickel	320		100	80	100	470	52	610	4600			150	190	1100	20	52	52
Selenium	80	50	50	60	200		5	170	4,200	20	5	120	480	2,700	5.6	5.6	5.6
Silver	80					3.2								25,900	10	10	10
Thallium	0.16	2	2	1.7	6.3			0.24	0.47			0.24	0.27	0.216	0.20	0.20	0.216
Zinc	4,800			1,000	1,000	120	120	7,400	26,000			2,300	2,900	16,500	50	120	120
Volatile Organic Compounds (VOC	Cs)																
Benzene	5	5	5	0.44	1.6			0.58	16			0.44	1.6	22.7	0.20	0.44	0.44
Toluene	1,000	1,000	1,000	72	130			57	520			180	410	18,900	0.20	57	57
Ethylbenzene	700	700	70	29				68	130			200	270	6,820	0.20	29	29
Xylenes	1,000	10,000	10,000												0.40	1,000	1,000
Acetone	7,200														5.0	7,200	7,200
2-Butanone (methyl ethyl ketone)	4,800														5.0	4,800	
Carbon Disulfide	800														0.20	800	800
Cis-1,2-Dichloroethene	16	70	70												0.20	16	16
Chlorobenzene	160	100	100	100	200			100	800			380	890	5,190	0.2	100	100
isopropylbenzene															0.20		
Methylene Chloride	5	5	5	10	100			20	1,000			16	250	3,600	1.0	5	10
n-propylbenzene	800														0.20	800	

								Protection	n of Surface	Water						
	Protection of	f Groundwat	er				National	Recommend								
	as Drinking W	ater (ingesti	on)	EPA 2016 CW	/A – Effective		Water Q	uality Criteri	a ^c	WAC 173	-201A Table 2	240 ^d	мтса			
				Human Hea Applic Washir Consum	able to	Aquat	ction of ic Life – nwater	Protec Human Consum		Protection of Aquatic Life – Freshwater	Protect Human Consum	Health:	Method B Standarde	Practical Quantitati	Proposed Site Groundwater	Proposed Site Surface Water
Constituent (by group) ^a	MTCA Method A/B (unrestricted land use) ^f	Federal MCL ^g	Washington State MCL ^g		Organisms Only	Acute	Chronic	Water and Organisms	Organisms Only	Acute Chronic	Water and Organisms	Organisms Only		on Limit (PQL)	Screening Level ^h	Screening Level ⁱ
Volatile Organic Compounds (VOC	Cs) (continued)			, ,						1			1		1	
p-Isopropyltoluene (cumene)	800													0.20	800	
Styrene	1,600	100	100											0.20	100	
Tetrachloroethene (PCE)	5	5	5	2.4	2.9			10	29		4.9	7.1	99.6	0.20	2.4	2.4
Trichloroethene (TCE)	0.54	5	5	0.3	0.7			0.6	7		0.38	0.86	12.8	0.20	0.3	0.3
1,2,4-Trimethylbenzene														0.20		
1,3,5-trimethylbenzene	80													0.20	80	
Vinyl Chloride (stormwater only)				0.02	0.18			0.022	1.6		0.02	0.26	3.7	0.02		0.02
Semi-Volatile Organic Compounds	s (SVOCs)															
4-Nitrophenol														5.0		
Benzoic Acid (stormwater only)														5.0		
Benzyl Alcohol														1.0		
Bis(2-ethylhexyl)phthalate	6.25	6	6	0.045	0.046			0.32	0.37		0.23	0.25	3.56	1.0	1.0	1.0
Butyl benzyl phthalate	46.1			0.013	0.013			0.10	0.10		0.56	0.58	8.32	1.0	1.0	1.0
Diethyl Phthalate (stormwater only)				200	200			600	600		4,200	5,000	28,400	1.0		200
Di-n-butyl Phthalate	1,600			8	8			20	30		450	510	2,910	1.0	8	8
Di-n-octyl Phthalate	160													1.0	160	
Naphthalenes ^j	160												4,710	0.01	160	4,710
Pentachlorophenol	0.219	1	1	0.002	0.002	19	15	0.03	0.04		0.046	0.1	1.47	5.0	5.0	5.0
Phenol	2,400			9,000	70,000			4,000	300,000		18,000	200,000	556,000	1.0	2,400	4,000
p-Cresol	800													1.0	800	
Polycyclic Aromatic Hydrocarbons	(PAHs)															
Acenaphthene	960			30	30			70	90		110	110	648	0.10	30	30
Acenaphthylene														0.10		
Anthracene	4,800							300	400		3,100	4,600	25,900	0.10	300	300
Benzo(a)anthracene	0.12			0.00016	0.00016			0.0012	0.0013		0.014	0.021	0.296	0.010	0.010	0.010
Benzo(a)pyrene	0.012	0.2	0.2	0.000016	0.000016			0.00012	0.00013		0.0014	0.0021	0.0296	0.010	0.010	0.010
Benzo(b)fluoranthene	0.12			0.00016	0.00016			0.0012	0.0013		0.014	0.021	0.296	0.010	0.010	0.010
Benzo(j,k)fluoranthene	1.2			0.0016	0.0016			0.012	0.013		0.014	0.21	2.96	0.010	0.010	0.010
Chrysene	12			0.016	0.016			0.12	0.13		1.4	2.1	29.6	0.010	0.016	0.016
Dibenz(a,h)anthracene	0.012			0.000016	0.000016			0.00012	0.00013		0.0014	0.0021	0.0296	0.010	0.010	0.010
Fluoranthene	640			6	6			20	20		16	16	86.4	0.10	6	6
Fluorene	640			10	10			50	70		420	610	3,460	0.10	10	10



								Protectio	n of Surface	Water							
	Protection of as Drinking Wa			EPA 2016 CV	/A – Effective			Recommend uality Criter			WAC 173	-201A Table 2	240 ^d	MTCA			
				Applic Washii	alth Criteria able to ngton: ^b ption of	Aquat	ction of ic Life – nwater		tion of Health: ption of	Aquat	ction of tic Life – hwater	Protect Human Consum	Health:	Method B Standard ^e	Practical Quantitati	Proposed Site Groundwater	Proposed Site Surface Water
Constituent (by group) ^a	MTCA Method A/B (unrestricted land use) ^f	Federal MCL ^g	Washington State MCL ^g	Water and Organisms	Organisms Only	Acute	Chronic	Water and Organisms	Organisms Only	Acute	Chronic	Water and Organisms	Organisms Only	of Human Health	on Limit (PQL)	Screening Level ^h	Screening Level ⁱ
Polycyclic Aromatic Hydrocarbon	s (PAHs) (continued)																
Indeno(1,2,3-cd)pyrene	0.12			0.00016	0.00016			0.0012	0.0013			0.014	0.021	0.296	0.010	0.010	0.010
Phenanthrene															0.10		
Pyrene	480			8	8			20	30			310	460	2,590	0.10	8	8
Total cPAHs TEQ	0.012	0.2	0.2	0.000016	0.000016			0.00012	0.00013			0.0014	0.0021	0.0296	0.085	0.015	0.015
Polychlorinated Biphenyls (PCBs)			<u>'</u>						•			•		'		•	
Total PCBs	0.1	0.5	0.5	0.000007	0.000007		0.014	0.000064	0.000064	2	0.014	0.00017	0.00017	0.000105	0.050	0.050	
Herbicides	<u> </u>							•	•								
2,4-DB	128														0.071	128	
2,4,5-T	160														0.071	160	
Bentazon	480															480	
Herbicides (continued)												_					
Chloramben	240															240	
Dacthal (chlorthal-dimethyl)	160															160	
Dalapon	240	200	200												0.46	200	
Dinoseb	16	7	7												0.047	7	
Picloram	1,120	500	500													500	
Silvex (2,4,5-TP)	128	50	50												0.048	50	
Organochlorine Pesticides															1		
4,4'-DDD ^k	0.365			0.0000079	0.0000079	1.1	0.001	0.00012	0.00012			0.000036	0.000036	0.0005	0.005	0.005	0.005
4,4'-DDE ^l	0.257			0.00000088	0.00000088			0.000018	0.000018			0.000051	0.000051	0.000353	0.005	0.005	0.005
Cis-chlordane ^m	0.25	2	2	0.000022	0.000022	2.4	0.0043	0.00031	0.00032	2.4	0.0043	0.000093	0.000093	0.00132	0.005	0.005	0.005
Gamma-chlordane ^m	0.25	2	2	0.000022	0.000022	2.4	0.0043	0.00031	0.00032	2.4	0.0043			0.00132	0.005	0.005	0.005
Endosulfan I ⁿ	96		1							0.22	0.056			57.6	0.005	0.056	0.056
Endosulfan II ⁿ	96		1							0.22	0.056			57.6	0.005	0.056	0.056
Endosulfan sulfate				9	10			20	40			9.7	10	57.6	0.005	9	9

All units in micrograms per liter (ug/L), unless otherwise stated.

Blank cells are intentional and indicate that criteria are not available.

Shaded cells denote the most restrictive criteria that is proposed as the site screening level.

- ^a Cleanup levels have been developed for only those compounds that have been detected at the Site above laboratory reporting limits. Constituents with SW only indicate that constituent was only detected in surface water samples.
- b Washington State human health criteria for the consumption of water and organisms and organisms only, EPA-approved human health criteria under 40 CFR 131.45.
- c National recommended water quality criteria for the protection of aquatic organisms and protection of human health based on consumption of organisms from Section 304 of the Clean Water Act.
- ^d Water Quality Standards for Surface Waters of the State of Washington, Chapter 173-201A WAC.
- ^e MTCA Method B Standard Surface Water Cleanup Level for the protection of human health; the most stringent value is used.
- f Washington State Model Toxics Control Act (MTCA) Cleanup Regulation Ground Water Method A or Method B Standard Formula Values for protection of human health through ingestion (groundwater as drinking water; WAC 173-340-720).
- ^g Federal and Washington State Maximum Contaminant Level (MCL) for drinking water.

- h Proposed site screening level for groundwater is the most restrictive of MTCA Method A/B and Protection of Surface Water Criteria Screening Levels, adjusted for laboratory PQLs.
- Proposed site screening level for surface water is the most restrictive of Protection of Surface Water Criteria Screening Levels, adjusted for laboratory PQLs.

 Screening level for naphthalenes is a total value for naphthalene, 1-methyl naphthalene, and 2-methyl naphthalene.
- k Screening levels for DDD are used for 4,4'-DDD.
- Screening levels for DDE are used for 4,4'-DDE.
- ^m Screening levels for chlordane are used for cis-chlordane and gamma-chlordane.
- ⁿ Screening levels for endosulfan are used for endosulfan I and endosulfan II.



				rable	6. Sumn	nary of (roundwater	Sample F	Results from M	onitoring \	vells, Pacific C	ity Park	Kemedia	ı ınvestiga	tion, Pa	citic, Was	ningtor	1.			
					T					Analy	rtical Parameter (μg/L)									
		Petrole	um Hydro	arbons			Vol	atile Organ	ic Compounds					Total Metals					Dissolved Met	als	
Sample	Sample								(cis) 1,2-	1,4-Dichloro											
Location	Date	GRO	DRO	Lube Oil	Benzene	Toluene	Ethylbenzene	Xylenes	Dichloroethene	benzene	Chlorobenzene	Arsenic	Cadmium	Chromium	Lead	Mercury	Arsenic	Cadmium	Chromium	Lead	Mercury
Site Screen	-																				
Level (µg/L		1,000	500	500	0.44	57	29	1,000	16		100	3.3	4.4	50	2.5	0.5	3.3	4.4	50	2.5	0.5
	10/6/15	ND (50)	ND (50)	ND (100)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (0.20)	ND (1.0)	1.1	ND (0.20)	ND (0.50)	1.1	ND (0.10)	1.2	ND (0.20)	ND (0.50)	ND (1.0)	ND (0.10)
	5/12/17	ND (100)	ND (260)	ND (420)	ND (0.20)	ND (1.0)	ND (0.20)	ND (0.40)	ND (0.20)	ND (0.20)	ND (0.20)	ND (3.3)	ND (4.4)	ND (11)	ND (1.1)	ND (0.50)		ND (4.0)	ND (10)	ND (1.0)	ND (0.50)
MW-1	3/23/18	ND (100)	ND (260)	ND (410)	ND (0.20)	ND (1.0)	ND (0.20)	ND (0.40)	ND (0.20)	ND (0.20)	ND (0.20)	ND (3.3)	ND (4.4)	ND (11)		ND (0.50)	NA	NA	NA	NA	NA
	6/21/18	ND (100)	ND (260)	ND (410)	ND (0.20)	ND (1.0)	ND (0.20)	ND (0.40)	ND (0.20)	ND (0.20)	ND (0.20)	ND (3.3)	ND (4.4)	ND (11)		ND (0.50)	NA	NA	NA	NA	NA
	9/26/18	ND (100)	ND (270)	ND (430)	ND (0.20)	ND (1.0)	ND (0.20)	ND (0.40)	ND (0.20)	ND (0.20)	ND (0.20)	ND (3.3)	ND (4.4)	ND (11)	ND (11)	ND (0.50)	NA	NA	NA	NA	NA
	10/5/15	ND (50)	ND (50)	ND (100)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (0.20)	ND (1.0)	5.7	ND (0.20)	2.3	ND (1.0)	ND (0.10)	5.1	ND (0.20)	1.6	ND (1.0)	ND (0.10)
	5/12/17	ND (100)	ND (270)	ND (440)	ND (0.20)	ND (1.0)	ND (0.20)	ND (0.40)	ND (0.20)	ND (0.20)	ND (0.20)	ND (3.3)	ND (4.4)	ND (11)	ND (1.1)	ND (0.50)	ND (3.0)	ND (4.0)	ND (10)	ND (1.0)	ND (0.50)
MW-2	3/23/18	ND (110)	ND (260)	ND (420)		ND (1.0)	ND (0.20)	ND (0.40)	ND (0.20)	ND (0.20)	ND (0.20)	3.4	ND (4.4)	ND (11)	ND (1.1)	ND (0.50)	NA	NA	NA	NA	NA
	6/21/18	ND (100)	ND (270)	ND (430)	ND (0.20)	ND (1.0)	ND (0.20)	ND (0.40)	ND (0.20)	ND (0.20)	ND (0.20)	ND (3.3)	ND (4.4)	ND (11)	ND (1.1)	ND (0.50)	NA	NA	NA	NA	NA
	9/26/18	ND (100)	ND (260)	ND (410)	ND (0.20)	ND (1.0)	ND (0.20)	ND (0.40)	ND (0.20)	ND (0.20)	ND (0.20)	4.9	ND (4.4)	ND (11)	ND (11)	ND (0.50)	NA	NA	NA	NA	NA
	10/6/15	ND (50)	ND (50)	ND (100)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (0.20)	ND (1.0)	4.1	ND (0.20)	2.8	ND (1.0)	ND (0.10)	4.8	ND (0.20)	1.5	ND (1.0)	ND (0.10)
	5/12/17	ND (100)	ND (260)	ND (420)	ND (0.20)	ND (1.0)	ND (0.20)	ND (0.40)	ND (0.20)	ND (0.20)	ND (0.20)	ND (3.3)	ND (4.4)	ND (11)	ND (1.1)	ND (0.50)	ND (3.0)	ND (4.0)	ND (10)	ND (1.0)	ND (0.50)
MW-3	3/23/18	ND (100)	ND (260)	ND (410)	ND (0.20)	ND (1.0)	ND (0.20)	ND (0.40)	0.22	ND (0.20)	ND (0.20)	ND (3.3)	ND (4.4)	ND (11)	ND (1.1)	ND (0.50)	NA	NA	NA	NA	NA
	6/21/18	ND (100)	ND (260)	ND (410)	ND (0.20)	ND (1.0)	ND (0.20)	ND (0.40)	ND (0.20)	ND (0.20)	ND (0.20)	ND (3.3)	ND (4.4)	ND (11)	ND (1.1)	ND (0.50)	NA	NA	NA	NA	NA
	9/26/18	ND (100)	ND (260)	ND (410)	ND (0.20)	ND (1.0)	ND (0.20)	ND (0.40)	0.35	ND (0.20)	ND (0.20)	ND (3.3)	ND (4.4)	ND (11)	ND (11)	ND (0.50)	NA	NA	NA	NA	NA
	10/5/15	ND (50)	ND (50)	ND (100)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (0.20)	5.7	13	ND (0.20)	2.7	ND (1.0)	ND (0.10)	9.8	ND (0.20)	1.5	ND (1.0)	ND (0.10)
	5/12/17	ND (100)	ND (260)	ND (420)	ND (0.20)	ND (1.0)	ND (0.20)	ND (0.40)	ND (0.20)	ND (0.20)	1.5	8.3	ND (4.4)	ND (11)	ND (1.1)	ND (0.50)	8.3	ND (4.0)	ND (10)	ND (1.0)	ND (0.50)
MW-4	3/23/18	ND (110)	ND (270)	ND (440)	ND (0.20)	ND (1.0)	ND (0.20)	ND (0.40)	ND (0.20)	ND (0.20)	1.1	6.0	ND (4.4)	ND (11)	ND (1.1)	ND (0.50)	NA	NA	NA	NA	NA
	6/21/18	ND (100)	ND (270)	ND (430)	ND (0.20)	ND (1.0)	ND (0.20)	ND (0.40)	ND (0.20)	ND (0.20)	ND (0.20)	ND (3.3)	ND (4.4)	ND (11)	ND (1.1)	ND (0.50)	NA	NA	NA	NA	NA
	9/26/18	ND (100)	ND (250)	ND (410)	0.22	ND (1.0)	ND (0.20)	ND (0.40)	ND (0.20)	ND (0.20)	4.6	14	ND (4.4)	ND (11)	ND (11)	ND (0.50)	NA	NA	NA	NA	NA
	10/5/15	ND (50)	ND (50)	ND (100)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (0.20)	ND (1.0)	1.4	ND (0.20)	0.52	ND (1.0)	ND (0.10)	ND (1.0)	ND (0.20)	ND (0.5)	ND (1.0)	ND (0.10)
	5/12/17	ND (100)	ND (260)	ND (410)	ND (0.20)	ND (1.0)	ND (0.20)	ND (0.40)	ND (0.20)	ND (0.20)	ND (0.20)	ND (3.3)	ND (4.4)	ND (11)	4.0	ND (0.50)	ND (3.0)	ND (4.0)	ND (10)	3.0	ND (0.50)
MW-5	3/23/18	ND (100)	ND (260)	ND (420)	ND (0.20)	ND (1.0)	ND (0.20)	ND (0.40)	ND (0.20)	ND (0.20)	ND (0.20)	ND (3.3)	ND (4.4)	ND (11)	3.2	ND (0.50)	NA	NA	NA	NA	NA
	6/21/18	ND (100)	ND (260)	ND (410)	ND (0.20)	ND (1.0)	ND (0.20)	ND (0.40)	ND (0.20)	ND (0.20)	ND (0.20)	ND (3.3)	ND (4.4)	ND (11)	5.6	ND (0.50)	NA	NA	NA	NA	NA
	9/26/18	ND (100)	ND (260)	ND (410)	ND (0.20)	ND (1.0)	ND (0.20)	ND (0.40)	ND (0.20)	ND (0.20)	ND (0.20)		ND (4.4)	ND (11)	1.9	ND (0.50)		NA	NA	NA	NA
	10/5/15				ND (1.0)		ND (1.0)	ND (1.0)	ND (1.0)	ND (0.20)	ND (1.0)	1.9	ND (0.20)	0.74	1.5	ND (0.10)		ND (0.20)	ND (0.50)	ND (1.0)	ND (0.10)
					ND (0.20)		ND (0.20)	ND (0.40)	ND (0.20)	ND (0.20)	ND (0.20)	ND (3.3)		ND (11)		ND (0.50)			ND (10)	ND (1.0)	
MW-6					ND (0.20)		ND (0.20)	ND (0.40)	ND (0.20)	ND (0.20)	ND (0.20)	ND (3.3)		ND (11)		ND (0.50)		NA	NA	NA	NA NA
		ND (100)			ND (0.20)		ND (0.20)	ND (0.40)	ND (0.20)	ND (0.20)	ND (0.20)	ND (3.3)		ND (11)		ND (0.50)		NA	NA	NA	NA
					ND (0.20)		ND (0.20)	ND (0.40)	ND (0.20)	0.20	ND (0.20)	4.5	ND (4.4)	ND (11)		ND (0.50)		NA	NA	NA	NA
					ND (0.20)		ND (0.20)	ND (0.40)	ND (0.20)	ND (0.20)	ND (0.20)	ND (3.3)		ND (11)	1.9	ND (0.50)		NA	NA	NA	NA
MW-7					ND (0.20)		ND (0.20)	ND (0.40)	ND (0.20)	ND (0.20)	ND (0.20)	4.6	ND (4.4)	ND (11)	2.0	ND (0.50)		NA NA	NA NA	NA NA	NA NA
1V1 V /					ND (0.20)		ND (0.20)	ND (0.40)		ND (0.20)	ND (0.20)	5.5	ND (4.4)	ND (11)		ND (0.50)		NA NA	NA NA	NA NA	NA NA
	3/20/10	14D (TOO)	140 (200)	1AD (4TO)	140 (0.20)	ND (1.0)	140 (0.20)	110 (0.40)	110 (0.20)	140 (0.20)	140 (0.20)	3.3	110 (4.4)	IND (TT)	IND (TT)	140 (0.30)	INA	INA	INA	11/7	11/7

Table 6 (continued). Summary of Groundwater Sample Results from Monitoring Wells, Pacific City Park Remedial Investigation, Pacific, Washington. Analytical Parameter (µg/L) **Petroleum Hydrocarbons Volatile Organic Compounds Total Metals Dissolved Metals** 1,4-Dichloro Sample Sample (cis) 1,2-Location Date GRO DRO Lube Oil Toluene Ethylbenzene Xylenes Dichloroethene benzene Chlorobenzene Arsenic Cadmium Chromium Lead Mercury Arsenic Cadmium Chromium Lead Mercury Benzene Site Screening 57 1,000 500 500 0.44 29 1,000 100 2.5 0.5 Level (µg/L) 16 3.3 4.4 50 3.3 4.4 50 2.5 0.5 3/23/18 | ND (100) | ND (260) | ND (420) ND (0.20) ND (1.0) ND (0.40) ND (0.20) ND (11) ND (0.50) ND (0.20) ND (0.20) ND (0.20) ND (3.3) ND (4.4) ND (11) NA NA NA NA NA 6/21/18 ND (100) ND (260) ND (410) ND (0.20) ND (1.0) ND (0.20) ND (0.40) ND (0.20) ND (0.20) ND (0.20) 3.9 ND (4.4) ND (11) ND (1.1) ND (0.50) NA MW-8 NA NA NA NA 9/26/18 | ND (100) | ND (260) | ND (410) ND (0.20) ND (1.0) ND (0.20) ND (0.40) ND (0.20) ND (0.20) ND (0.20) ND (3.3) ND (4.4) ND (11) ND (11) ND (0.50) NA NA NA NA NA 3/23/18 ND (100) ND (260) ND (420) ND (0.20) ND (1.0) ND (0.40) ND (11) ND (0.50) ND (0.20) ND (0.20) ND (0.20) ND (0.20) ND (3.3) ND (4.4) ND (11) NA NA NA NA NA MW-9 6/21/18 | ND (100) | ND (260) | ND (410) ND (0.20) ND (1.0) ND (0.20) ND (0.40) ND (0.20) ND (0.20) ND (0.20) ND (3.3) ND (4.4) ND (11) ND (1.1) ND (0.50) NA NA NA NA NA 9/26/18 ND (100) ND (250) ND (410) ND (0.20) ND (1.0) ND (11) ND (0.50) ND (0.20) ND (0.40) ND (0.20) 0.38 3.6 ND (4.4) ND (11) ND (0.20) NA NA NA NA NA

Note: MW7, MW8, and MW9 correspond to borings B06, B09, and B11

BOLD values detected above the reporting limit.

Shaded values exceed site criteria.

cPAHs (TEQ) = Carcinogenic polycyclic aromatic hydrocarbons toxic equivalency

DRO = Diesel range organics

PCBs = Polychlorinated biphenyls

GRO = Gasoline range organics

μg/L = micrograms per liter

NA = not analyzed or not applicable

ND = not detected above laboratory reporting limits shown in parentheses



			Table	e 6 (continued	l). Summary	of Ground	water Sample	e Results, Pa	cific City Par	k Remedial Inv	vestigation, Pa	cific, Washingto	on.		
				Field Parameters	;						Analytical Parame	eter (µg/L)			
										Carcin	ogenic Polycyclic	Aromatic Hydrocai	rbons (cPAHs)		
Sample Location	Sample Date	Temp (°C)	DO (mg/L)	Cond (µS/cm)	pH (std units)	Turbidity (NTU)	Total PCBs	Benzo(a) anthracene	Chrysene	Benzo(b) fluoranthene	Benzo(j,k) fluoranthene	Benzo(a) pyrene	Indeno(1,2,3- cd) pyrene	Dibenz(a,h) anthracene	Total cPAHs (TEQ)
Site Screen	ing Level (μg/L)	NA	NA	NA	NA	NA	0.05	0.01	0.016	0.01	0.01	0.01	0.01	0.01	0.015
	10/6/15	NR	NR	NR	NR	NR	NA	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.076)
	5/12/17	9.0	3.28	98	6.84	Clear	NA	ND (0.0095)	ND (0.0095)	ND (0.0095)	ND (0.0095)	ND (0.0095)	ND (0.0095)	ND (0.0095)	ND (0.0072)
MW-1	3/23/18	6.9	4.67	97	6.94	Clear	NA	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0071)
	6/21/18	11.3	1.69	77	6.79	Clear	NA	ND (0.0095)	ND (0.0095)	ND (0.0095)	ND (0.0095)	ND (0.0095)	ND (0.0095)	ND (0.0095)	ND (0.0072)
	9/26/18	14.2	2.76	113	6.64	Clear	NA	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.0076)
	10/5/15	NR	NR	NR	NR	NR	NA	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.0072)
	5/12/17	11.9	2.47	296	6.58	Clear	NA	ND (0.0099)	ND (0.0099)	ND (0.0099)	ND (0.0099)	ND (0.0099)	ND (0.0099)	ND (0.0099)	ND (0.0075)
MW-2	3/23/18	9.8	0.66	328	6.54	Clear	NA	ND (0.011)	ND (0.011)	ND (0.011)	ND (0.011)	ND (0.011)	ND (0.011)	ND (0.011)	ND (0.0083)
	6/21/18	13.7	3.28	270	6.33	Clear	NA	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.0076)
	9/26/18	15.8	0.23	276	6.30	Clear	NA	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0071)
	10/6/15	NR	NR	NR	NR	NR	NA	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.0072)
	5/12/17	10.9	0.69	332	7.08	Clear	NA	ND (0.0095)	ND (0.0095)	ND (0.0095)	ND (0.0095)	ND (0.0095)	ND (0.0095)	ND (0.0095)	ND (0.0072)
MW-3	3/23/18	8.1	0.50	332	7.01	Clear	NA	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.0076)
	6/21/18	12.8	0.11	281	7.08	Clear	NA	ND (0.0097)	ND (0.0097)	ND (0.0097)	ND (0.0097)	ND (0.0097)	ND (0.0097)	ND (0.0097)	ND (0.0073)
	9/26/18	14.5	0.12	322	6.65	Clear	NA	ND (0.0095)	ND (0.0095)	ND (0.0095)	ND (0.0095)	ND (0.0095)	ND (0.0095)	ND (0.0095)	ND (0.0072)
	10/5/15	NR	NR	NR	NR	NR	NA	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.0072)
	5/12/17	11.5	0.19	348	6.60	Clear	NA	ND (0.0096)	ND (0.0096)	ND (0.0096)	ND (0.0096)	ND (0.0096)	ND (0.0096)	ND (0.0096)	ND (0.0072)
MW-4	3/23/18	9.1	0.55	307	6.15	Clear	NA	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.0076)
	6/21/18	15.4	2.05	309	6.62	Clear	NA	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.0076)
	9/26/18	20.4	1.47	325	6.10	Clear	NA	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.0076)
	10/5/15	NR	NR	NR	NR	NR	NA	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.0072)
	5/12/17	9.5	1.06	156	7.08	Clear	NA	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.0076)
MW-5	3/23/18	6.7	0.47	129	6.69	Clear	NA	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.0076)
	6/21/18	11.6	0.08	126	7.44	Clear	NA	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.0076)
	9/26/18	15.3	0.26	193	6.90	Clear	NA	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0071)
	10/5/15	NR	NR	NR	NR	NR	NA	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.100)	ND (0.0072)
	5/12/17	10.2	0.25	132	6.25	Clear	NA	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.0076)
MW-6	3/23/18	6.9	0.73	161	5.95	Clear	NA	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0071)
	6/21/18	12.4	0.14	154	6.69	Clear	NA	0.014	0.014	0.012	0.012	0.011	0.012	0.011	0.017
	9/26/18	15.1	0.44	341	6.25	Clear	NA	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.0076)
	3/23/18	6.9	0.52	127	6.94	Clear	NA	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0071
MW-7	6/21/18	16.2	0.12	137	6.59	Clear	NA	ND (0.0096)	ND (0.0096)	ND (0.0096)	ND (0.0096)	ND (0.0096)	ND (0.0096)	ND (0.0096)	ND (0.0072
	9/26/18	16.2	0.49	151	6.47	Clear	NA	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0071)

			Table	e 6 (continued). Summary o	of Ground	water Sample	e Results, Pa	cific City Parl	k Remedial Inv	estigation, Pa	cific, Washingto	on.		
				Field Parameters	1						Analytical Param	eter (µg/L)			
										Carcin	ogenic Polycyclic	Aromatic Hydroca	rbons (cPAHs)		
Sample Location	Sample Date	Temp (°C)	DO (mg/L)	Cond (µS/cm)	pH (std units)	Turbidity (NTU)	Total PCBs	Benzo(a) anthracene	Chrysene	Benzo(b) fluoranthene	Benzo(j,k) fluoranthene	Benzo(a) pyrene	Indeno(1,2,3- cd) pyrene	Dibenz(a,h) anthracene	Total cPAHs (TEQ)
Site Screen	ning Level (µg/L)	NA	NA	NA	NA	NA	0.05	0.01	0.016	0.01	0.01	0.01	0.01	0.01	0.015
	3/23/18	10.8	0.45	400	6.62	Clear	NA	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.0076
MW-8	6/21/18	13.4	2.44	384	6.24	Clear	NA	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.010)	ND (0.0076
	9/26/18	16.4	0.61	325	6.56	Clear	NA	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0094)	ND (0.0071)
	3/23/18	10.5	0.42	294	6.22	Clear	NA	ND (0.011)	ND (0.011)	ND (0.011)	ND (0.011)	ND (0.011)	ND (0.011)	ND (0.011)	ND (0.0083
MW-9	6/21/18	11.5	2.65	240	6.58	Clear	NA	ND (0.0098)	ND (0.0098)	ND (0.0098)	ND (0.0098)	ND (0.0098)	ND (0.0098)	ND (0.0098)	ND (0.0074
	9/26/18	14.5	0.60	249	6.41	Clear	NA	ND (0.0098)	ND (0.0098)	ND (0.0098)	ND (0.0098)	ND (0.0098)	ND (0.0098)	ND (0.0098)	ND (0.0074

Note: MW7, MW8, and MW9 correspond to borings B06, B09, and B11

BOLD values detected above the reporting limit.

Shaded values exceed site criteria.

cPAHs (TEQ) = Carcinogenic polycyclic aromatic hydrocarbons toxic equivalency

DRO = Diesel range organics

PCBs = Polychlorinated biphenyls

GRO = Gasoline range organics

μg/L = micrograms per liter

NA = not analyzed or not applicable

ND = not detected above laboratory reporting limits shown in parentheses

NR = not reported



				Table	7. Summ	nary of Soi	I Sample F	Results, Pa	cific City F	Park Reme	dial Inves	stigation,	Pacific, Wa	ashington.	,				
Analytical							•		Sample	Location									
Parameter	GP-:	1-15	GP-	2-15	GP-	3-15	GP-	4-15		GP-5-15		GP-	6-15	GP-	7-15		GP-8-15		Site Screening
Sample Date	9/17	7/15	9/1	7/15	9/17	7/15	9/1	7/15		9/17/15		9/1	7/15	9/1	7/15		9/17/15		Level ^a
Depth (feet)	5.0	13.5	4.5	14.0	9.0	12.5	5.0	13.0	5.0	13.0	13 dup	7.5	14.0	5.0	14.0	5.0	14.5	14.5 dup	(mg/kg)
Petroleum Hydrocarl	bons (mg/	kg)																	
Gasoline Range Organics	ND (2.96)	ND (3.44)	ND (3.19)	ND (3.13)	ND (3.33)	ND (7.99)	ND (2.87)	ND (3.41)	ND (4.31)	ND (3.20)	ND (3.60)	ND (5.40)	ND (3.35)	ND (3.79)	ND (3.74)	ND (3.05)	ND (3.65)	ND (3.25)	100
Diesel Range Organics	ND (23.0)	ND (25.1)	ND (19.8)	ND (21.2	ND (24.4)	ND (38.9)	ND (23.8)	ND (25.0)	ND (23.7)	ND (23.2)	ND (21.9)	ND (33.7)	ND (22.5)	ND (19.1)	ND (22.2)	ND (22.9)	ND (24.9)	ND (24.7)	200
Lube Oil Range Organics	ND (57)	ND (63)	75	ND (53)	ND (61)	275	182	ND (63)	ND (59)	ND (58)	ND (55)	217	ND (56)	ND (48)	ND (56)	119	ND (62)	ND (62)	2,000
Volatile Organic Com	npounds b	y EPA 8260) (mg/kg)																
Benzene	ND (0.012)	ND (0.014)	ND (0.013)	ND (0.013)	ND (0.013)	ND (0.030)	ND (0.012)	ND (0.014)	ND (0.017)	ND (0.013)	ND (0.014)	ND (0.022)	ND (0.013)	ND (0.015)	ND (0.015)	ND (0.012)	ND (0.015)	ND (0.013)	0.001
Toluene	ND (0.012)	ND (0.014)	ND (0.013)	ND (0.013)	ND (0.013)	ND (0.030)	ND (0.012)	ND (0.014)	ND (0.017)	ND (0.013)	ND (0.014)	ND (0.022)	ND (0.013)	ND (0.015)	ND (0.015)	ND (0.012)	ND (0.015)	ND (0.013)	0.024
Ethylbenzene	ND (0.018)	ND (0.021)	ND (0.019)	ND (0.019)	ND (0.013)	ND (0.048)	ND (0.017)	ND (0.020)	ND (0.026)	ND (0.019)	ND (0.022)	ND (0.032)	ND (0.020)	ND (0.023)	ND (0.022)	ND (0.018)	ND (0.022)	ND (0.020)	0.014
Total Xylenes	ND (0.012)	ND (0.014)	ND (0.013)	ND (0.013)	ND (0.013)	ND (0.030)	ND (0.012)	ND (0.014)	ND (0.017)	ND (0.013)	ND (0.014)	ND (0.022)	ND (0.013)	ND (0.015)	ND (0.015)	ND (0.012)	ND (0.015)	ND (0.013)	0.52
Acetone	-	-	-	_	-	-	-	-	-	_	-	-	-	-	_	-	-	-	2.07
2-Butanone	-	_	_	_	-	-	-	-	-	_	_	-	_	-	_	-	_	_	1.38
Carbon Disulfide	-	_	_	_	-	-	-	-	-	_	_	-	_	-	_	-	_	_	0.27
Cis-1,2-Dichloroethene	ND (0.012)	ND (0.014)	ND (0.013)	ND (0.013)	ND (0.013)	ND (0.030)	ND (0.012)	ND (0.014)	ND (0.017)	ND (0.013)	ND (0.014)	ND (0.022)	ND (0.013)	ND (0.015)	ND (0.015)	ND (0.012)	ND (0.015)	ND (0.013)	0.005
Chlorobenzene	ND (0.012)	ND (0.014)	ND (0.013)	ND (0.013)	ND (0.013)	ND (0.030)	ND (0.012)	ND (0.014)	ND (0.017)	ND (0.013)	ND (0.014)	ND (0.022)	ND (0.013)	ND (0.015)	ND (0.015)	ND (0.012)	ND (0.015)	ND (0.013)	0.051
Methylene Chloride	ND (0.012)	ND (0.014)	ND (0.013)	ND (0.013)	ND (0.013)	ND (0.030)	ND (0.012)	ND (0.014)	ND (0.017)	ND (0.013)	ND (0.014)	ND (0.022)	ND (0.013)	ND (0.015)	ND (0.015)	ND (0.012)	ND (0.015)	ND (0.013)	0.005
p-Isopropyltoluene	ND (0.012)	ND (0.014)	ND (0.013)	ND (0.013)	ND (0.013)	ND (0.030)	ND (0.012)	ND (0.014)	ND (0.017)	ND (0.013)	ND (0.014)	ND (0.022)	ND (0.013)	ND (0.015)	ND (0.015)	ND (0.012)	ND (0.015)	ND (0.013)	0.229
Styrene	ND (0.012)	ND (0.014)	ND (0.013)	ND (0.013)	ND (0.013)	ND (0.030)	ND (0.012)	ND (0.014)	ND (0.017)	ND (0.013)	ND (0.014)	ND (0.022)	ND (0.013)	ND (0.015)	ND (0.015)	ND (0.012)	ND (0.015)	ND (0.013)	0.120
Tetrachloroethene	ND (0.012)	ND (0.014)	ND (0.013)	ND (0.013)	ND (0.013)	ND (0.032)	ND (0.012)	ND (0.014)	ND (0.017)	ND (0.013)	ND (0.014)	ND (0.022)	ND (0.013)	0.023	ND (0.015)	ND (0.012)	ND (0.015)	ND (0.013)	0.0013
Trichloroethene	ND (0.012)	ND (0.014)	ND (0.013)	ND (0.013)	ND (0.013)	ND (0.030)	ND (0.012)	ND (0.014)	ND (0.017)	ND (0.013)	ND (0.014)	ND (0.022)	ND (0.013)	ND (0.015)	ND (0.015)	ND (0.012)	ND (0.015)	ND (0.013)	0.001
1,2,4-Trimethylbenzene	ND (0.012)	ND (0.014)	ND (0.013)	ND (0.013)	ND (0.013)	ND (0.030)	ND (0.012)	ND (0.014)	ND (0.017)	ND (0.013)	ND (0.014)	ND (0.022)	ND (0.013)	ND (0.015)	ND (0.015)	ND (0.012)	ND (0.015)	ND (0.013)	NA
Total Metals by EPA	6010D/74	71B (mg/k	g)																
Arsenic	1.7	1.4	3.1	2.9	3.4	102	3.8	1.4	7.6	2.3	4.0	50	2.2	6.1	2.4	12	5.9	3.3	20
Barium	23	15	38	21	24	154	36	17	85	23	19	631	23	148	17	224	50	36	41.3
Cadmium	ND (0.17)	ND (0.18)	0.25	ND (0.20)	ND (0.20)	3.8	0.20	ND (0.21)	1.4	ND (0.19)	ND (0.18)	37	ND (0.19)	0.79	ND (0.18)	0.71	ND (0.20)	ND (0.20)	1
Chromium	9.8	13	18	17	11	143	12	11	20	12	14	115	9.7	23	9.59	17	20	15	48
Lead	1.3	1.2	22	1.4	22	2,780	19	1.3	45	1.9	2.0	2,180	1.5	63	1.5	370	3.2	2.1	<i>2</i> 5
Mercury	ND (0.29)	ND (0.30)	ND (0.23)	ND (0.29)	ND (0.30)	0.55	ND (0.28)	ND (0.32)	ND (0.30)	ND (0.28)	ND (0.27)	9.1	ND (0.30)	2.5	ND (0.28)	ND (0.29)	ND (0.30)	ND (0.29)	0.07
Selenium	1.2	1.3	1.5	1.2	1.1	1.2	1.5	1.1	1.6	1.1	1.4	1.7	1.1	1.3	1.1	1.9	1.5	1.0	10
Silver	ND (0.087)	ND (0.091)	ND (0.084)	ND (0.098)	ND (0.10)	2.5	ND (0.093)	ND (0.11)	0.12	ND (0.095)	ND (0.092)	2.6	ND (0.096)	0.14	ND (0.09)	0.14	ND (0.10)	ND (0.10)	0.61
Polychlorinated Biph	enyls (PCE	Bs) by EPA	8082A (mg	g/kg)															
Total PCBs	-	-	-	_	-	ND (0.20)	-	-	-	_	_	ND (0.17)	-	-	-	ND (0.11)	-	-	0.05
Semi-Volatile Organi	ic Compou	nds by EPA	4 8270D/SI	M (mg/kg)															
Acenapthene	ND (0.089)	ND (0.101)	ND (0.086)	ND (0.093)	ND (0.095)	ND (0.156)	ND (0.091)	ND (0.096)	ND (0.091)	ND (0.094)	ND (0.090)	ND (0.135)	ND (0.097)	ND (0.083)	ND (0.092)	ND (0.095)	ND (0.109)	ND (0.103)	0.156
Acenaphthylene	ND (0.089)	ND (0.101)	ND (0.086)	ND (0.093)	ND (0.095)	ND (0.156)	ND (0.091)	ND (0.096)	ND (0.091)	ND (0.094)	ND (0.090)	ND (0.135)	ND (0.097)	ND (0.083)	ND (0.092)	ND (0.095)	ND (0.109)	ND (0.103)	NA
Anthracene	ND (0.089)	ND (0.101)	0.091	ND (0.093)	ND (0.095)	ND (0.156)	ND (0.091)	ND (0.096)	ND (0.091)	ND (0.094)	ND (0.090)	ND (0.135)	ND (0.097)	ND (0.083)	ND (0.092)	ND (0.095)	ND (0.109)	ND (0.103)	7.134
Benzyl Alcohol	ND (0.111)	ND (0.126)	ND (0.108)	ND (0.116)	ND (0.119)	ND (0.195)	ND (0.113)	ND (0.120)	ND (0.114)	ND (0.117)	ND (0.112)	ND (0.169)	ND (0.122)	ND (0.103)	ND (0.115)	ND (0.119)	ND (0.136)	ND (0.129)	NA



			Tal	ble 7 (con	tinued).	Summar	y of Soil S	ample Res	sults, Pacif	ic City Pa	rk Remed	ial Investi	gation, Pa	cific, Wasł	nington.				
									Sample	Location									
Analytical Parameter	GP-	1-15	GP-	2-15	GP-	3-15	GP-	4-15		GP-5-15		GP-	6-15	GP-	7-15		GP-8-15		Site Screening
Sample Date	9/1	7/15	9/1	7/15	9/17	7/15	9/1	7/15		9/17/15		9/1	7/15	9/1	7/15		9/17/15		Levela
Depth (feet)	5.0	13.5	4.5	14.0	9.0	12.5	5.0	13.0	5.0	13.0	13 dup	7.5	14.0	5.0	14.0	5.0	14.5	14.5 dup	(mg/kg)
Semi-Volatile Organic C	Compounds	by EPA 827	70D/SIM (m	g/kg) (conti	nued)														
Bis(2-Ethylhexyl) Phthalate	ND (0.111)	ND (0.126)	ND (0.108)	ND (0.116)	ND (0.119)	ND (0.195)	ND (0.113)	ND (0.120)	ND (0.114)	ND (0.117)	ND (0.112)	ND (0.169)	ND (0.122)	ND (0.103)	ND (0.115)	ND (0.119)	ND (0.136)	ND (0.129)	0.111
Butyl Benzylphthalate	ND (0.111)	ND (0.126)	ND (0.108)	ND (0.116)	ND (0.119)	ND (0.195)	ND (0.113)	ND (0.120)	ND (0.114)	ND (0.117)	ND (0.112)	ND (0.169)	ND (0.122)	ND (0.103)	ND (0.115)	ND (0.119)	ND (0.136)	ND (0.129)	0.033
Dibutyl Phthalate	ND (0.111)	ND (0.126)	0.13	ND (0.116)	ND (0.119)	ND (0.195)	0.12	ND (0.120)	0.28	ND (0.117)	ND (0.112)	0.17	ND (0.122)	ND (0.103)	ND (0.115)	0.13	ND (0.136)	ND (0.129)	0.17
Di-N-Octyl Phthalate	ND (0.111)	ND (0.126)	ND (0.108)	ND (0.116)	ND (0.119)	ND (0.195)	ND (0.113)	ND (0.120)	ND (0.114)	ND (0.117)	ND (0.112)	ND (0.169)	ND (0.122)	ND (0.103)	ND (0.115)	ND (0.119)	ND (0.136)	ND (0.129)	800
Fluoranthene	ND (0.089)	ND (0.101)	0.13	ND (0.093)	ND (0.095)	ND (0.156)	ND (0.091)	ND (0.096)	ND (0.091)	ND (0.094)	ND (0.090)	0.15	ND (0.097)	ND (0.083)	ND (0.092)	0.11	ND (0.109)	ND (0.103)	0.296
Fluorene	ND (0.089)	ND (0.101)	ND (0.086)	ND (0.093)	ND (0.095)	ND (0.156)	ND (0.091)	ND (0.096)	ND (0.091)	ND (0.094)	ND (0.090)	ND (0.135)	ND (0.097)	ND (0.083)	ND (0.092)	ND (0.095)	ND (0.109)	ND (0.103)	0.080
1-Methylnaphthalene	ND (0.089)	ND (0.101)	ND (0.086)	ND (0.093)	ND (0.095)	ND (0.156)	ND (0.091)	ND (0.096)	ND (0.091)	ND (0.094)	ND (0.090)	ND (0.135)	ND (0.097)	ND (0.083)	ND (0.092)	ND (0.095)	ND (0.109)	ND (0.103)	0.236
2-Methylnaphthalene	ND (0.089)	ND (0.101)	ND (0.086)	ND (0.093)	ND (0.095)	ND (0.156)	ND (0.091)	ND (0.096)	ND (0.091)	ND (0.094)	ND (0.090)	ND (0.135)	ND (0.097)	ND (0.083)	ND (0.092)	ND (0.095)	ND (0.109)	ND (0.103)	0.236
Naphthalene	ND (0.089)	ND (0.101)	ND (0.086)	ND (0.093)	ND (0.095)	ND (0.156)	ND (0.091)	ND (0.096)	ND (0.091)	ND (0.094)	ND (0.090)	ND (0.135)	ND (0.097)	ND (0.083)	ND (0.092)	ND (0.095)	ND (0.109)	ND (0.103)	0.236
4-Nitrophenol	ND (0.553)	ND (0.630)	ND (0.538)	ND (0.578)	ND (0.594)	ND (0.974)	ND (0.566)	ND (0.602)	ND (0.571)	ND (0.586)	ND (0.560)	ND (0.843)	ND (0.608)	ND (0.517)	ND (0.576)	ND (0.594)	ND (0.681)	ND (0.643)	NA
p-Cresol	ND (0.111)	ND (0.126)	ND (0.108)	ND (0.116)	ND (0.119)	ND (0.195)	ND (0.113)	ND (0.120)	ND (0.114)	ND (0.117)	ND (0.112)	ND (0.169)	ND (0.122)	ND (0.103)	ND (0.115)	ND (0.119)	ND (0.136)	ND (0.129)	8,000
Pentachlorophenol	ND (0.111)	ND (0.126)	0.20	ND (0.116)	ND (0.119)	ND (0.195)	ND (0.113)	ND (0.120)	ND (0.114)	ND (0.117)	ND (0.112)	ND (0.169)	ND (0.122)	ND (0.103)	ND (0.115)	ND (0.119)	ND (0.136)	ND (0.129)	0.17
Phenanthrene	ND (0.089)	ND (0.101)	ND (0.086)	ND (0.093)	ND (0.095)	ND (0.156)	ND (0.091)	ND (0.096)	ND (0.091)	ND (0.094)	ND (0.090)	ND (0.135)	ND (0.097)	ND (0.083)	ND (0.092)	ND (0.095)	ND (0.109)	ND (0.103)	0.0067
Phenol	ND (0.221)	ND (0.252)	ND (0.215)	ND (0.231)	ND (0.238)	ND (0.389)	ND (0.226)	ND (0.241)	ND (0.228)	ND (0.234)	ND (0.224)	ND (0.337)	ND (0.243)	ND (0.207)	ND (0.230)	ND (0.237)	ND (0.272)	ND (0.257)	0.757
Pyrene	ND (0.089)	ND (0.101)	0.16	ND (0.093)	ND (0.095)	ND (0.156)	ND (0.091)	ND (0.096)	ND (0.091)	ND (0.094)	ND (0.090)	0.29	ND (0.097)	ND (0.083)	ND (0.092)	0.11	ND (0.109)	ND (0.103)	0.546
Carcinogenic Polycyclic	Aromatic I	Hydrocarbo	ns (cPAHs) k	oy EPA 8270	D/SIM (mg/k	(g)													
Benzo(a)anthracene	ND (0.089)	ND (0.101)	0.14	ND (0.093)	0.11	0.19	0.12	0.11	0.13	0.11	0.10	0.62	ND (0.097)	0.11	ND (0.092)	0.13	ND (0.109)	ND (0.103)	0.0067
Benzo(a)pyrene	ND (0.089)	ND (0.101)	ND (0.086)	ND (0.093)	ND (0.095)	0.28	ND (0.091)	ND (0.096)	ND (0.091)	ND (0.094)	ND (0.090)	0.93	ND (0.097)	ND (0.083)	ND (0.092)	ND (0.095)	0.55	0.23	0.01
Benzo(b)fluoranthene	ND (0.089)	ND (0.101)	ND (0.086)	ND (0.093)	ND (0.095)	ND (0.156)	ND (0.091)	ND (0.096)	ND (0.091)	ND (0.094)	ND (0.090)	0.51	ND (0.097)	ND (0.083)	ND (0.092)	ND (0.095)	ND (0.109)	ND (0.103)	0.012
Benzo(j,k)fluoranthene	ND (0.089)	ND (0.101)	ND (0.086)	ND (0.093)	ND (0.095)	ND (0.156)	ND (0.091)	ND (0.096)	ND (0.091)	ND (0.094)	ND (0.090)	0.19	ND (0.097)	ND (0.083)	ND (0.092)	ND (0.095)	ND (0.109)	ND (0.103)	0.012
Chrysene	ND (0.089)	ND (0.101)	ND (0.086)	ND (0.093)	ND (0.095)	ND (0.156)	ND (0.091)	ND (0.096)	ND (0.091)	ND (0.094)	ND (0.090)	0.25	ND (0.097)	ND (0.083)	ND (0.092)	ND (0.095)	ND (0.109)	ND (0.103)	0.0067
Dibenz(a,h)anthracene	ND (0.089)	ND (0.101)	ND (0.086)	ND (0.093)	ND (0.095)	ND (0.156)	ND (0.091)	ND (0.096)	ND (0.091)	ND (0.094)	ND (0.090)	ND (0.135)	ND (0.097)	ND (0.083)	ND (0.092)	ND (0.095)	ND (0.109)	53	0.018
Indeno(1,2,3-cd)pyrene	ND (0.089)	ND (0.101)	ND (0.086)	ND (0.093)	ND (0.095)	ND (0.156)	ND (0.091)	ND (0.096)	ND (0.091)	ND (0.094)	ND (0.090)	ND (0.135)	ND (0.097)	ND (0.083)	ND (0.092)	ND (0.095)	ND (0.109)	62	0.035
Total cPAHs (TEQ) ^b	ND (0.08)	ND (0.091)	0.087	ND (0.084)	0.092	0.36	0.089	0.093	0.091	0.091	0.087	1.1	ND (0.088)	0.082	ND (0.083)	0.094	0.60	0.27	0.020
Herbicides by EPA 8151	A (mg/kg)																		
2,4-DB	ND (0.029)	ND (0.031)	ND (0.026)	ND (0.030)	ND (0.030)	ND (0.051)	ND (0.029)	ND (0.032)	ND (0.030)	ND (0.030)	ND (0.029)	ND (0.042)	ND (0.030)	ND (0.026)	ND (0.029)	ND (0.030)	ND (0.034)	ND (0.032)	640
2,4,5-T	ND (0.059)		ND (0.052)	ND (0.059)	ND (0.059)	ND (0.102)	ND (0.058)	ND (0.065)	ND (0.061)	ND (0.050	ND (0.058)	ND (0.085)	ND (0.059)	ND (0.051)	ND (0.059)	ND (0.060)	ND (0.068)	ND (0.064)	800
Bentazon		ND (0.063)	ND (0.052)	ND (0.059)	ND (0.059)	ND (0.102)	ND (0.058)	ND (0.065)	ND (0.061)	ND (0.050	ND (0.058)	ND (0.085)	ND (0.059)	ND (0.051)	ND (0.059)	ND (0.060)	ND (0.068)	ND (0.064)	2,400
Chloramben		ND (0.025)	23	ND (0.024)	ND (0.024)	ND (0.041)	25	ND (0.026)	24	ND (0.024)	ND (0.023)	ND (0.034)	ND (0.024)	ND (0.021)	ND (0.023)	26	ND (0.027)	ND (0.025)	1,200
Chlorthal-dimethyl	ND (0.029)	ND (0.031)	ND (0.026)	ND (0.030)	ND (0.030)	ND (0.051)	ND (0.029)	ND (0.032)	ND (0.030)	ND (0.030)	ND (0.029)	ND (0.042)	ND (0.030)	ND (0.026)	ND (0.029)	ND (0.030)	ND (0.034)	ND (0.032)	800
Dalapon	ND (0.024)		ND (0.021)	ND (0.024)	ND (0.024)	ND (0.041)	ND (0.023)	ND (0.026)	ND (0.024)	ND (0.024)	ND (0.023)	ND (0.034)	ND (0.024)	ND (0.021)	ND (0.023)	ND (0.024)	ND (0.027)	ND (0.025)	2,400
Dinoseb	ND (0.059)		ND (0.052)	ND (0.059)	ND (0.059)	ND (0.102)	ND (0.058)	ND (0.065)	ND (0.061)	ND (0.050	ND (0.058)	ND (0.085)	ND (0.059)	ND (0.051)	ND (0.059)	ND (0.060)	ND (0.068)	ND (0.064)	80
Picloram		ND (0.063)	ND (0.052)	ND (0.059)	ND (0.059)	ND (0.102)	ND (0.058)	ND (0.065)	ND (0.061)	ND (0.050	ND (0.058)	ND (0.085)	ND (0.059)	ND (0.051)	ND (0.059)	ND (0.060)	ND (0.068)	ND (0.064)	5,600
Silvex		ND (0.025)	ND (0.021)	ND (0.024)	ND (0.024)	ND (0.041)	ND (0.023)	ND (0.026)	ND (0.024)	ND (0.024)	ND (0.023)	ND (0.034)	ND (0.024)	ND (0.021)	ND (0.023)	ND (0.024)	ND (0.027)	ND (0.025)	640



			Tal	ble 7 (con	tinued).	Summar	y of Soil S	ample Re	sults, Pacif	ic City Pa	k Remedi	ial Investi	gation, Pa	cific, Wash	nington.				
									Sample	Location									
Analytical Parameter	GP-	1-15	GP-2	2-15	GP-	3-15	GP-	4-15		GP-5-15		GP-	6-15	GP-	7-15		GP-8-15		Site Screening
Sample Date	9/1	7/15	9/17	7/15	9/17	7/15	9/1	7/15		9/17/15		9/1	7/15	9/1	7/15		9/17/15		Level ^a
Depth (feet)	5.0	13.5	4.5	14.0	9.0	12.5	5.0	13.0	5.0	13.0	13 dup	7.5	14.0	5.0	14.0	5.0	14.5	14.5 dup	(mg/kg)
Organochlorine Pesticio	les by EPA	8081(mg/kg	g)																
4,4'-DDD	ND (0.023)	ND (0.025)	ND (0.021)	ND (0.023)	ND (0.023)	ND (0.039)	ND (0.023)	ND (0.025)	ND (0.023)	ND (0.024)	ND (0.022)	ND (0.031)	ND (0.023)	0.074	ND (0.023)	ND (0.023)	ND (0.026)	ND (0.025)	0.01
4,4'-DDE	ND (0.023)	ND (0.025)	ND (0.021)	ND (0.023)	ND (0.023)	ND (0.039)	ND (0.023)	ND (0.025)	ND (0.023)	ND (0.024)	ND (0.022)	ND (0.031)	ND (0.023)	ND (0.199)	ND (0.023)	ND (0.023)	ND (0.026)	ND (0.025)	0.01
Cis-Chlordane (alpha)	ND (0.011)	ND (0.012)	ND (0.010)	ND (0.011)	ND (0.011)	ND (0.020)	ND (0.011)	ND (0.013)	ND (0.012)	ND (0.012)	ND (0.011)	ND (0.016)	ND (0.012)	ND (0.010)	ND (0.011)	ND (0.012)	ND (0.013)	ND (0.013)	0.01
Endosulfan I	ND (0.011)	ND (0.012)	ND (0.010)	ND (0.011)	ND (0.011)	ND (0.020)	ND (0.011)	ND (0.013)	ND (0.012)	ND (0.012)	ND (0.011)	ND (0.016)	ND (0.012)	ND (0.010)	ND (0.011)	ND (0.012)	ND (0.013)	ND (0.013)	0.005
Endosulfan II	ND (0.023)	ND (0.025)	ND (0.021)	ND (0.023)	ND (0.023)	ND (0.039)	ND (0.023)	ND (0.025)	ND (0.023)	ND (0.024)	ND (0.022)	ND (0.031)	ND (0.023)	0.063	ND (0.023)	ND (0.023)	ND (0.026)	ND (0.025)	0.01
Endosulfan Sulfate	ND (0.023)	ND (0.025)	ND (0.021)	ND (0.023)	ND (0.023)	ND (0.039)	ND (0.023)	ND (0.025)	ND (0.023)	ND (0.024)	ND (0.022)	ND (0.031)	ND (0.023)	ND (0.199)	ND (0.023)	ND (0.023)	ND (0.026)	ND (0.025)	NA
Gamma-Chlordane	ND (0.011)	ND (0.012)	ND (0.010)	ND (0.011)	ND (0.011)	ND (0.020)	ND (0.011)	ND (0.013)	ND (0.012)	ND (0.012)	ND (0.011)	ND (0.016)	ND (0.012)	ND (0.010)	ND (0.011)	ND (0.012)	ND (0.013)	ND (0.013)	0.01
Methoxychlor	ND (0.057)	ND (0.062)	ND (0.051)	ND (0.058)	ND (0.059)	ND (0.100)	ND (0.058)	ND (0.063)	ND (0.058)	ND (0.060)	ND (0.056)	ND (0.079)	ND (0.058)	ND (0.050)	ND (0.057)	ND (0.057)	ND (0.064)	ND (0.063)	0.01

SEE END OF TABLE 7 FOR COMPLETE LIST OF TABLE NOTES.

BOLD values detected above the reporting limit.

Shaded values exceed the site screening level

Native Soil Fill/Refuse

King County

			Table 7 (c	ontinued).	Summary	of Soil Sam	ple Results,	Pacific City	Park Remed	ial Investiga	tion, Pacific,	Washingto	n.			
								Sample Location	on							
Analytical Parameter	GP-	9-15	GP-1	LO-15	GP-1	11-15	GP-1	L2-15	GP-1	13-15		GP-14-15		GP-1	L5-15	Site
Sample Date	9/1	7/15	9/1	7/15	9/1	7/15	9/1	8/15	9/2	1/15		9/21/15		9/2	1/15	Screening Level ^a
Depth (feet)	5.0	12.5	4.5	13.0	4.5	14.5	4.0	13.5	4.5	13.5	3.5	13.5	13.5 dup	6.0	14.0	(mg/kg)
Petroleum Hydrocarbor	ns (mg/kg)															
Gasoline Range Organics	ND (2.98)	ND (3.30)	ND (4.64)	ND (3.48)	ND (2.85)	ND (6.73)	ND (3.03)	ND (3.34)	ND (3.30)	ND (2.98)	ND (2.98)	ND (2.98)	ND (2.98)	ND (2.98)	ND (2.98)	100
Diesel Range Organics	ND (25)	ND (25)	ND (27)	ND (24)	ND (21)	ND (36)	ND (23)	ND (25)	ND (22)	ND (27)	ND (20)	ND (34)	ND (27)	ND (25)	ND (26)	200
Lube Oil Range Organics	ND (62)	ND (63)	3,840	ND (60)	103	ND (91)	ND (58)	ND (63)	462	ND (68)	ND (50)	ND (86)	ND (66)	ND (64)	ND (65)	2,000
Volatile Organic Compo	unds by EPA 8	3260 (mg/kg)														
Benzene	ND (0.012)	ND (0.013)	ND (0.019)	ND (0.014)	ND (0.011)	ND (0.027)	ND (0.012)	ND (0.013)	ND (0.013)	ND (0.016)	ND (0.014)	ND (0.025)	ND (0.016)	ND (0.018)	ND (0.015)	0.001
Toluene	ND (0.012)	ND (0.013)	ND (0.019)	ND (0.014)	ND (0.011)	ND (0.027)	ND (0.012)	ND (0.013)	ND (0.013)	ND (0.016)	ND (0.014)	ND (0.025)	ND (0.016)	ND (0.018)	ND (0.015)	0.024
Ethylbenzene	ND (0.018)	ND (0.020)	ND (0.028)	ND (0.021)	ND (0.017)	ND (0.040)	ND (0.018)	ND (0.020)	ND (0.020)	ND (0.024)	ND (0.021)	ND (0.037)	ND (0.024)	ND (0.027)	ND (0.022)	0.014
Total Xylenes	ND (0.012)	ND (0.013)	ND (0.019)	ND (0.014)	ND (0.011)	ND (0.027)	ND (0.012)	ND (0.013)	ND (0.013)	ND (0.016)	ND (0.014)	ND (0.025)	ND (0.016)	ND (0.018)	ND (0.015)	0.52
Acetone	_	-	-	_	-	-	_	_	-	_	-	_	_	-	_	2.07
2-Butanone	-	-	-	-	_	_	_	_	-	_	-	_	_	_	_	1.38
Carbon Disulfide	_	_	_	_	_	_	_	_	_	_	_	_	_	_	_	0.27
Cis-1,2-dichloroethene	ND (0.012)	ND (0.013)	ND (0.019)	ND (0.014)	ND (0.011)	ND (0.027)	ND (0.012)	ND (0.013)	ND (0.013)	ND (0.016)	ND (0.014)	ND (0.025)	ND (0.016)	ND (0.018)	ND (0.015)	0.005
Chlorobenzene	ND (0.012)	ND (0.013)	ND (0.019)	ND (0.014)	ND (0.011)	ND (0.027)	ND (0.012)	ND (0.013)	ND (0.013)	ND (0.016)	ND (0.014)	ND (0.025)	ND (0.016)	ND (0.018)	ND (0.015)	0.051
Methylene Chloride	ND (0.012)	ND (0.013)	ND (0.019)	ND (0.014)	ND (0.011)	ND (0.027)	ND (0.012)	ND (0.013)	ND (0.013)	ND (0.016)	ND (0.014)	ND (0.025)	ND (0.016)	ND (0.018)	ND (0.015)	0.005
p-Isopropyltoluene	ND (0.012)	ND (0.013)	ND (0.019)	ND (0.014)	ND (0.011)	ND (0.027)	ND (0.012)	ND (0.013)	ND (0.013)	ND (0.016)	ND (0.014)	ND (0.025)	ND (0.016)	ND (0.018)	ND (0.015)	0.229
Styrene	ND (0.012)	ND (0.014)	ND (0.013)	ND (0.013)	ND (0.013)	ND (0.030)	ND (0.012)	ND (0.014)	ND (0.017)	ND (0.013)	ND (0.014)	ND (0.022)	ND (0.013)	ND (0.015)	ND (0.015)	0.120
Tetrachlorethene	ND (0.012)	ND (0.013)	ND (0.019)	ND (0.014)	ND (0.011)	ND (0.027)	ND (0.012)	ND (0.013)	ND (0.013)	ND (0.016)	ND (0.014)	ND (0.025)	ND (0.016)	ND (0.018)	ND (0.015)	0.0013
Trichloroethene	ND (0.012)	ND (0.013)	0.066	ND (0.014)	ND (0.011)	ND (0.027)	ND (0.012)	ND (0.013)	ND (0.013)	ND (0.016)	ND (0.014)	ND (0.025)	ND (0.016)	ND (0.018)	ND (0.015)	0.001
1,2,4-Trimethylbenzene	ND (0.012)	ND (0.014)	ND (0.013)	ND (0.013)	ND (0.013)	ND (0.030)	ND (0.012)	ND (0.014)	ND (0.017)	ND (0.013)	ND (0.014)	ND (0.022)	ND (0.013)	ND (0.015)	ND (0.015)	NA
Total Metals by EPA 601	LOD/7471B (m	g/kg)					_									
Arsenic	1.9	2.7	39.7	1.2	2.5	3.0	2.8	1.2	3.4	3.3	2.9	7.7	4.6	2.5	1.7	20
Barium	29	59	232	31	33	72	27	18	40	73	36	95	103	42	32	41.3
Cadmium	ND (0.18)	ND (0.22)	8.3	ND (0.19)	ND (0.16)	ND (0.28)	ND (0.20)	ND (0.20)	0.19	ND (0.22)	0.19	ND (0.30)	ND (0.24)	0.39	ND (0.21)	1
Chromium	18	21	314	14	14	22	13	14	19	24	13	23	32	18	12	48
Lead	8.3	3.6	3,320	2.06	40	4.3	3.5	1.3	424	4.3	20	5.2	6.2	15	1.6	25
Mercury	ND (0.28)	ND (0.30)	ND (0.31)	ND (0.30)	ND (0.25)	ND (0.45)	ND (0.29)	ND (0.31)	ND (0.26)	ND (0.33)	ND (0.28)	ND (0.46)	ND (0.31)	ND (0.33)	ND (0.30)	0.07
Selenium	0.74	1.7	0.96	1.0	0.80	1.4	1.4	0.89	0.99	2.1	0.96	3.2	2.5	0.66	0.83	10
Silver	ND (0.091)	ND (0.109)	1.99	ND (0.093)	ND (0.080)	ND (0.14)	ND (0.10)	ND (0.099)	ND (0.091)	ND (0.11)	ND (0.093)	ND (0.15)	ND (0.12)	ND (0.12)	ND (0.11)	0.61
Polychlorinated Biphen	yls (PCBs) by E	PA 8082A (mg	g/kg)			T		1		1		ı	1		T	
Total PCBs	-	-	ND (0.13)	-	ND (0.11)	_	_	_	ND (0.11)	_	-	_	_	-	-	0.05
Semi-Volatile Organic C	ompounds by	EPA 8270D/SI	M (mg/kg)			ı		1		1		T			1	
Acenapthene	ND (0.095)	ND (0.108)	29.2	ND (0.101)	ND (0.085)	ND (0.144)	ND (0.095)	ND (0.100)	ND (0.087)	ND (0.111)	ND (0.092)	ND (0.146)	ND (0.114)	ND (0.117)	ND (0.108)	0.156
Acenaphthylene	ND (0.095)	ND (0.108)	0.13	ND (0.101)	ND (0.085)	ND (0.144)	ND (0.095)	ND (0.100)	ND (0.087)	ND (0.111)	ND (0.092)	ND (0.146)	ND (0.114)	ND (0.117)	ND (0.108)	NA
Anthracene	ND (0.095)	ND (0.108)	106	ND (0.101)	ND (0.085)	ND (0.144)	ND (0.095)	ND (0.100)	ND (0.087)	ND (0.111)	ND (0.092)	ND (0.146)	ND (0.114)	ND (0.117)	ND (0.108)	7.134
Benzyl Alcohol	ND (0.119)	ND (0.135)	ND (0.138)	ND (0.127)	ND (0.106)	ND (0.180)	ND (0.118)	ND (0.124)	ND (0.109)	ND (0.139)	ND (0.115)	ND (0.182)	ND (0.142)	ND (0.146)	ND (0.134)	NA
Bis(2-Ethylhexyl) Phthalate	ND (0.119)	ND (0.135)	1.6	ND (0.127)	ND (0.106)	ND (0.180)	ND (0.118)	ND (0.124)	0.17	ND (0.139)	0.12	0.25	ND (0.142)	0.15	ND (0.134)	0.111



			Table 7 (c	ontinued).	Summary	of Soil Sam	ple Results,	Pacific City	Park Remed	ial Investiga	tion, Pacific	Washingto	n.			
								Sample Location	on							
Analytical Parameter	GP-	9-15	GP-1	.0-15	GP-1	1-15	GP-1	L2-15	GP-1	13-15		GP-14-15		GP-1	L5-15	Site
Sample Date	9/1	7/15	9/17	7/15	9/1 ⁻	7/15	9/1	8/15	9/2:	1/15		9/21/15		9/2	1/15	Screening Level ^a
Depth (feet)	5.0	12.5	4.5	13.0	4.5	14.5	4.0	13.5	4.5	13.5	3.5	13.5	13.5 dup	6.0	14.0	(mg/kg)
Semi-Volatile Organic C	ompounds by	EPA 8270D/S	IM (mg/kg) (co	ntinued)												
Butyl Benzylphthalate	ND (0.119)	ND (0.135)	ND (0.138)	ND (0.127)	ND (0.106)	ND (0.180)	ND (0.118)	ND (0.124)	ND (0.109)	ND (0.139)	ND (0.115)	ND (0.182)	ND (0.142)	ND (0.146)	ND (0.134)	0.033
Di-N-Octyl Phthalate	ND (0.119)	ND (0.135)	ND (0.138)	ND (0.127)	ND (0.106)	ND (0.180)	ND (0.118)	ND (0.124)	ND (0.109)	ND (0.139)	ND (0.115)	ND (0.182)	ND (0.142)	ND (0.146)	ND (0.134)	0.17
Dibutyl Phthalate	ND (0.119)	ND (0.135)	0.18	ND (0.127)	ND (0.106)	ND (0.180)	ND (0.118)	ND (0.124)	ND (0.109)	ND (0.139)	ND (0.115)	ND (0.182)	ND (0.142)	ND (0.146)	ND (0.134)	800
Fluoranthene	ND (0.095)	ND (0.108)	365	0.11	ND (0.085)	0.19	ND (0.095)	ND (0.100)	0.14	ND (0.111)	ND (0.092)	ND (0.146)	ND (0.114)	ND (0.117)	ND (0.108)	0.296
Fluorene	ND (0.095)	ND (0.108)	37.5	ND (0.101)	ND (0.085)	ND (0.144)	ND (0.095)	ND (0.100)	ND (0.087)	ND (0.111)	ND (0.092)	ND (0.146)	ND (0.114)	ND (0.117)	ND (0.108)	0.080
2-Methylnaphthalene	ND (0.095)	ND (0.108)	3.1	ND (0.101)	ND (0.085)	ND (0.144)	ND (0.095)	ND (0.100)	ND (0.087)	ND (0.111)	ND (0.092)	ND (0.146)	ND (0.114)	ND (0.117)	ND (0.108)	0.236
1-Methylnaphthalene	ND (0.095)	ND (0.108)	6.1	ND (0.101)	ND (0.085)	ND (0.144)	ND (0.095)	ND (0.100)	ND (0.087)	ND (0.111)	ND (0.092)	ND (0.146)	ND (0.114)	ND (0.117)	ND (0.108)	0.236
Naphthalene	ND (0.095)	ND (0.108)	4.4	ND (0.101)	ND (0.085)	ND (0.144)	ND (0.095)	ND (0.100)	ND (0.087)	ND (0.00002)	ND (0.00002)	ND (0.00004)	ND (0.110)	ND (0.00003)	ND (0.00002)	0.236
4-Nitrophenol	ND (0.595)	ND (0.68)	1.4	ND (0.633)	ND (0.528)	ND (0.899)	ND (0.592)	ND (0.622)	ND (0.546)	ND (0.693)	ND (0.576)	ND (0.910)	ND (0.711)	ND (0.732)	ND (0.672)	NA
p-Cresol	ND (0.119)	ND (0.135)	ND (0.138)	ND (0.127)	ND (0.106)	ND (0.180)	ND (0.118)	ND (0.124)	ND (0.109)	ND (0.139)	ND (0.115)	ND (0.182)	ND (0.142)	ND (0.146)	ND (0.134)	8,000
Pentachlorophenol	ND (0.119)	ND (0.135)	ND (0.138)	ND (0.127)	ND (0.106)	ND (0.180)	ND (0.118)	ND (0.124)	ND (0.109)	ND (0.139)	ND (0.115)	ND (0.182)	ND (0.142)	ND (0.146)	ND (0.134)	0.17
Phenanthrene	ND (0.095)	ND (0.108)	317	ND (0.101)	ND (0.085)	ND (0.144)	ND (0.095)	ND (0.100)	ND (0.087)	ND (0.111)	ND (0.092)	ND (0.146)	ND (0.114)	ND (0.117)	ND (0.108)	0.0067
Phenol	ND (0.238)	ND (0.270)	ND (0.276)	ND (0.253)	ND (0.211)	ND (0.360)	ND (0.237)	ND (0.249)	ND (0.218)	ND (0.277)	ND (0.231)	ND (0.364)	ND (0.284)	ND (0.293)	ND (0.269)	0.757
Pyrene	ND (0.095)	ND (0.108)	345	ND (0.101)	ND (0.085)	0.19	ND (0.095)	ND (0.100)	0.13	ND (0.111)	ND (0.092)	ND (0.146)	ND (0.114)	ND (0.117)	ND (0.108)	0.546
Carcinogenic Polycyclic	Aromatic Hyd	rocarbons (cP	AHs) by EPA 82	270D/SIM (mg,	/kg)											
Benzo(a)anthracene	0.11	0.12	164	ND (0.10)	0.11	0.20	ND (0.095)	ND (0.092)	0.13	ND (0.111)	ND (0.092)	ND (0.146)	ND (0.114)	ND (0.117)	ND (0.108)	0.0067
Benzo(a)pyrene	ND (0.095)	0.24	91	0.213	ND (0.085)	0.59	ND (0.095)	ND (0.092)	ND (0.087)	0.45	ND (0.092)	0.64	0.54	ND (0.117)	ND (0.108)	0.01
Benzo(b)fluoranthene	ND (0.095)	ND (0.108)	186	ND (0.101)	ND (0.085)	ND (0.144)	ND (0.095)	ND (0.092)	0.23	ND (0.111)	ND (0.092)	ND (0.146)	ND (0.114)	ND (0.117)	ND (0.108)	0.012
Benzo(j,k)fluoranthene	ND (0.095)	ND (0.108)	28	ND (0.101)	ND (0.085)	ND (0.144)	ND (0.095)	ND (0.092)	ND (0.087)	ND (0.111)	ND (0.092)	ND (0.146)	ND (0.114)	ND (0.117)	ND (0.108)	0.012
Chrysene	ND (0.095)	ND (0.108)	193	ND (0.101)	ND (0.085)	ND (0.144)	ND (0.095)	ND (0.092)	ND (0.087)	ND (0.111)	ND (0.092)	ND (0.146)	ND (0.114)	ND (0.117)	ND (0.108)	0.0067
Dibenz(a,h)anthracene	ND (0.095)	ND (0.108)	62	ND (0.101)	ND (0.085)	ND (0.144)	ND (0.095)	ND (0.092)	ND (0.087)	ND (0.111)	ND (0.095)	ND (0.146)	53	ND (0.117)	ND (0.108)	0.018
Indeno(1,2,3-cd)pyrene	ND (0.095)	ND (0.108)	53	ND (0.101)	ND (0.085)	ND (0.144)	ND (0.095)	ND (0.092)	ND (0.087)	ND (0.111)	ND (0.095)	ND (0.146)	62	ND (0.117)	ND (0.108)	0.035
Total cPAHs (TEQ) ^b	0.092	0.25	213	0.25	0.083	0.66	ND (0.86)	ND (0.090)	0.10	0.48	ND (0.083)	0.64	0.58	ND (0.117)	ND (0.108)	0.020
Herbicides by EPA 8151	A (mg/kg)									_			_	_		
2,4-DB	ND (0.030)	ND (0.033)	ND (0.034)	ND (0.032)	ND (0.027)	ND (0.047)	ND (0.031)	ND (0.031)	ND (0.028)	ND (0.034)	ND (0.028)	ND (0.046)	ND (0.036)	ND (0.036)	ND (0.033)	640
2,4,5-T	ND (0.061)	ND (0.067)	ND (0.068)	ND (0.064)	ND (0.054)	ND (0.093)	ND (0.062)	ND (0.062)	ND (0.056)	ND (0.068)	ND (0.057)	ND (0.092)	ND (0.071)	ND (0.073)	ND (0.067)	800
Bentazon	ND (0.061)	ND (0.067)	ND (0.068)	ND (0.064)	ND (0.054)	ND (0.093)	ND (0.062)	ND (0.062)	ND (0.056)	ND (0.068)	ND (0.057)	ND (0.092)	ND (0.071)	ND (0.073)	ND (0.067)	2,400
Chloramben	ND (0.024)	ND (0.027)	ND (0.027)	ND (0.025)	ND (0.022)	ND (0.037)	ND (0.025)	ND (0.025)	ND (0.022)	ND (0.027)	ND (0.023)	ND (0.037)	ND (0.028)	ND (0.029)	ND (0.027)	1,200
Chlorthal-dimethyl	ND (0.030)	ND (0.033)	ND (0.034)	ND (0.032)	ND (0.027)	ND (0.047)	ND (0.031)	ND (0.031)	ND (0.028)	ND (0.034)	ND (0.028)	ND (0.046)	ND (0.036)	ND (0.036)	ND (0.033)	800
Dalapon	ND (0.024)	ND (0.027)	ND (0.027)	ND (0.025)	ND (0.022)	ND (0.037)	ND (0.025)	ND (0.025)	ND (0.022)	ND (0.027)	ND (0.023)	ND (0.037)	ND (0.028)	ND (0.029)	ND (0.027)	2,400
Dinoseb	ND (0.061)	ND (0.067)	ND (0.068)	ND (0.064)	ND (0.054)	ND (0.093)	ND (0.062)	ND (0.062)	ND (0.056)	ND (0.068)	ND (0.057)	ND (0.092)	ND (0.071)	ND (0.073)	ND (0.067)	80
Picloram	ND (0.061)	ND (0.067)	ND (0.068)	ND (0.064)	ND (0.054)	ND (0.093)	ND (0.062)	ND (0.062)	ND (0.056)	ND (0.068)	ND (0.057)	ND (0.092)	ND (0.071)	ND (0.073)	ND (0.067)	5,600
Silvex	ND (0.024)	ND (0.027)	ND (0.027)	ND (0.025)	ND (0.022)	ND (0.037)	ND (0.025)	ND (0.025)	ND (0.022)	ND (0.027)	ND (0.023)	ND (0.037)	ND (0.028)	ND (0.029)	ND (0.027)	640

			Table 7 (c	ontinued).	Summary	of Soil Sam	ple Results,	Pacific City	Park Remedi	ial Investiga	tion, Pacific,	Washingto	n.			
								Sample Location	on							
Analytical Parameter	GP-	9-15	GP-1	LO- 1 5	GP-1	1-15	GP-1	.2-15	GP-1	3-15		GP-14-15		GP-1	L5-15	Site Screening
Sample Date	9/17	7/15	9/1	7/15	9/1	7/15	9/18	3/15	9/21	L/15		9/21/15		9/2:	1/15	Level ^a
Depth (feet)	5.0	12.5	4.5	13.0	4.5	14.5	4.0	13.5	4.5	13.5	3.5	13.5	13.5 dup	6.0	14.0	(mg/kg)
Organochlorine Pesticid	es by EPA 808	31 (mg/kg)														
4,4'-DDD	ND (0.024)	ND (0.027)	ND (0.026)	ND (0.024)	ND (0.022)	ND (0.038)	О	ND (0.024)	ND (0.021)	ND (0.027)	ND (0.021)	ND (0.034)	ND (0.027)	ND (0.027)	ND (0.025)	0.01
4,4'-DDE	ND (0.024)	ND (0.027)	ND (0.026)	ND (0.024)	ND (0.022)	ND (0.038)	ND (0.024)	ND (0.024)	ND (0.021)	ND (0.027)	ND (0.021)	ND (0.034)	ND (0.027)	ND (0.027)	ND (0.025)	0.01
Cis-Chlordane (alpha)	ND (0.012)	ND (0.013)	ND (0.013)	ND (0.012)	ND (0.011)	ND (0.019)	ND (0.012)	ND (0.012)	ND (0.011)	ND (0.014)	ND (0.011)	ND (0.017)	ND (0.013)	ND (0.014)	ND (0.012)	<i>0</i> .01
Endosulfan I	ND (0.012)	ND (0.013)	ND (0.013)	ND (0.012)	ND (0.011)	ND (0.019)	ND (0.012)	ND (0.012)	ND (0.011)	ND (0.014)	ND (0.011)	ND (0.017)	ND (0.013)	ND (0.014)	ND (0.012)	<i>0</i> .005
Endosulfan II	ND (0.024)	ND (0.027)	ND (0.026)	ND (0.024)	0.056	ND (0.038)	ND (0.024)	ND (0.024)	ND (0.021)	ND (0.027)	ND (0.021)	ND (0.034)	ND (0.027)	ND (0.027)	ND (0.025)	0.01
Endosulfan Sulfate	ND (0.024)	ND (0.027)	ND (0.026)	ND (0.024)	0.060	ND (0.038)	ND (0.024)	ND (0.024)	ND (0.021)	ND (0.027)	ND (0.021)	ND (0.034)	ND (0.027)	ND (0.027)	ND (0.025)	NA
Gamma-Chlordane	ND (0.012)	ND (0.013)	ND (0.013)	ND (0.012)	ND (0.011)	ND (0.019)	ND (0.012)	ND (0.012)	ND (0.011)	ND (0.014)	ND (0.011)	ND (0.017)	ND (0.013)	ND (0.014)	ND (0.012)	0.01
Methoxychlor	ND (0.059)	ND (0.066)	ND (0.065)	ND (0.060)	0.127	ND (0.094)	ND (0.060)	ND (0.059)	ND (0.053)	ND (0.067)	ND (0.053)	ND (0.085)	ND (0.066)	ND (0.069)	ND (0.062)	0.01

SEE END OF TABLE 7 FOR COMPLETE LIST OF TABLE NOTES.

BOLD values detected above the reporting limit.

Shaded values exceed the site screening level





			Table	7 (continu	ed). Summ	nary of Soil	Sample Res	ults, Pacific	: City Park I	Remedial In	vestigation	, Pacific, W	ashington.				
								Sample	Location								
Analytical Parameter	GP-1	.6- 1 5		GP-17-15		GP-1	8-15	GP-1	.9-15	GP-2	20-15	GP-2	21-15		GP-22-15		Site Screening
Sample Date	9/2:	1/15		9/21/15		9/2:	L/15	9/2:	1/15	9/2:	1/15	9/2	1/15		9/22/15		Level ^a
Depth (feet)	8.0	13.5	7.0	14.5	7 dup	3.0	14.0	5.0	14.5	5.5	14.0	4.0	14.0	1.5	3.0	3.0 dup	(mg/kg)
Petroleum Hydrocarbo	ns (mg/kg)																
Gasoline Range Organics	ND (4.83)	ND (3.95)	ND (4.2)	ND (3.9)	ND (3.8)	ND (3.5)	ND (3.7)	ND (3.1)	ND (4.7)	ND (3.3)	ND (3.6)	3.7	ND (4.0)	ND (3.5)	ND (3.6)	ND (2.7)	100
Diesel Range Organics	72	ND (25)	ND (27.2)	ND (20.2)	ND (23.9)	ND (20.3)	ND (23.6)	ND (20.7)	ND (24.8)	ND (19.7)	ND (20.8)	ND (23.3)	ND (21.8)	ND (21.4)	ND (20.4)	ND (19.3)	200
Lube Oil Range Organics	427	ND (63)	70	ND (51)	401	302	ND (59)	ND (52)	ND (62)	285	ND (52)	728	ND (55)	ND (53)	ND (51)	ND (48)	2,000
Volatile Organic Compo	ounds by EPA	8260 (mg/k	g)														
Benzene	ND (0.019)	ND (0.016)	ND (0.017)	ND (0.016)	ND (0.015)	ND (0.014)	ND (0.015)	ND (0.012)	ND (0.019)	ND (0.013)	ND (0.014)	ND (0.013)	ND (0.016)	ND (0.014)	ND (0.014)	ND (0.010)	0.001
Toluene	0.022	ND (0.016)	ND (0.017)	ND (0.016)	ND (0.015)	ND (0.014)	ND (0.015)	ND (0.012)	ND (0.019)	ND (0.013)	ND (0.014)	ND (0.013)	ND (0.016)	ND (0.014)	ND (0.014)	ND (0.010)	0.024
Ethylbenzene	ND (0.029)	ND (0.024)	ND (0.025)	ND (0.023)	ND (0.023)	ND (0.021)	ND (0.022)	ND (0.019)	ND (0.028)	ND (0.020)	ND (0.022)	ND (0.020)	ND (0.024)	ND (0.021)	ND (0.022)	ND (0.016)	0.014
Total Xylenes	0.029	ND (0.016)	ND (0.017)	ND (0.016)	ND (0.015)	ND (0.014)	ND (0.015)	ND (0.012)	ND (0.019)	ND (0.013)	ND (0.014)	ND (0.013)	ND (0.016)	ND (0.014)	ND (0.014)	ND (0.010)	0.52
Acetone	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	2.07
2-Butanone	-	-	-	-	-	-	-	-	-	-	_	-	-	-	_	_	1.38
Carbon Disulfide	-	-	_	-	-	_	_	_	-	-	-	-	_	-	_	-	0.27
Cis-1,2-dichloroethene	ND (0.019)	ND (0.016)	ND (0.017)	ND (0.016)	ND (0.015)	ND (0.014)	ND (0.015)	ND (0.012)	ND (0.019)	ND (0.013)	ND (0.014)	ND (0.013)	ND (0.016)	ND (0.014)	ND (0.014)	ND (0.010)	0.005
Chlorobenzene	ND (0.019)	ND (0.016)	ND (0.017)	ND (0.016)	ND (0.015)	ND (0.014)	ND (0.015)	ND (0.012)	ND (0.019)	ND (0.013)	ND (0.014)	ND (0.013)	ND (0.016)	ND (0.014)	ND (0.014)	ND (0.010)	0.051
Methylene Chloride	ND (0.019)	ND (0.016)	ND (0.017)	ND (0.016)	ND (0.015)	ND (0.014)	ND (0.015)	ND (0.012)	0.020	ND (0.013)	0.015	0.014	ND (0.016)	ND (0.014)	ND (0.014)	ND (0.010)	0.005
p-Isopropyltoluene	ND (0.019)	ND (0.016)	ND (0.017)	ND (0.016)	ND (0.015)	ND (0.014)	ND (0.015)	ND (0.012)	ND (0.019)	ND (0.013)	ND (0.014)	0.13	ND (0.016)	ND (0.014)	ND (0.014)	ND (0.010)	0.229
Styrene	ND (0.012)	ND (0.014)	ND (0.013)	ND (0.013)	ND (0.013)	ND (0.030)	ND (0.012)	ND (0.014)	ND (0.017)	ND (0.013)	ND (0.014)	ND (0.022)	ND (0.013)	ND (0.015)	ND (0.015)	ND (0.012)	0.120
Tetrachlorethene	ND (0.019)	ND (0.016)	ND (0.017)	ND (0.016)	ND (0.015)	ND (0.014)	ND (0.015)	ND (0.012)	ND (0.019)	ND (0.013)	ND (0.014)	ND (0.013)	ND (0.016)	ND (0.014)	ND (0.014)	ND (0.010)	0.0013
Trichloroethene	ND (0.019)	ND (0.016)	ND (0.017)	ND (0.016)	ND (0.015)	ND (0.014)	ND (0.015)	ND (0.012)	ND (0.019)	ND (0.013)	ND (0.014)	ND (0.013)	ND (0.016)	ND (0.014)	ND (0.014)	ND (0.010)	0.001
1,2,4-Trimethylbenzene	ND (0.012)	ND (0.014)	ND (0.013)	ND (0.013)	ND (0.013)	ND (0.030)	ND (0.012)	ND (0.014)	ND (0.017)	ND (0.013)	ND (0.014)	ND (0.022)	ND (0.013)	ND (0.015)	ND (0.015)	ND (0.012)	NA
Total Metals by EPA 60	10D/7471B (mg/kg)		ı	ı						ı						
Arsenic	18.5	1.8	8.7	2.0	5.7	5.2	5.6	3.5	2.7	4.0	1.9	2.6	2.8	3.8	3.8	4.0	20
Barium	78	26	52	22	38	47	25	44	38	32	21	40	27	46	44	44	41.3
Cadmium	1.9	ND (0.198)	0.36	ND (0.19)	3.2	0.32	ND (0.18)	ND (0.18)	ND (0.20)	ND (0.17)	ND (0.18)	0.23	ND (0.20)	ND (0.17)	ND (0.19)	ND (0.18)	1
Chromium	57	13	16	12	19	18	56	22	17	19	11	22	14	18	19	19	48
Lead	552	2.2	49	2.6	39	60	5.7	5.4	2.2	19.6	2.3	18	3.4	5.1	4.1	5.5	25
Mercury	ND (0.42)	ND (0.29)	ND (0.31)	ND (0.26)	ND (0.29)	ND (0.24)	ND (0.26)	ND (0.24)	ND (0.28)	ND (0.27)	ND (0.27)	ND (0.25)	ND (0.28)	0.38	ND (0.25)	ND (0.25)	0.07
Selenium	ND (0.63)	0.89	1.5	0.99	1.5	1.1	0.90	1,1	0.86	1.2	0.9	0.85	0.78	1.7	1.6	1.8	10
Silver	0.18	ND (0.10)	ND (0.117)	ND (0.095)	ND (0.102)	ND (0.091)	ND (0.090)	ND (0.088)	ND (0.099)	ND (0.085)	ND (0.092)	ND (0.095)	ND (0.099)	ND (0.084)	ND (0.097)	ND (0.091)	0.61
Polychlorinated Biphen	yls (PCBs) by	, EPA 8082A ((mg/kg)	I	I						I		T		T	<u> </u>	
Total PCBs	0.42	-	-	-	_	-	-	-	-	_	_	-	-	_	-	-	0.05
Semi-Volatile Organic (Compounds b	y EPA 8270D	/SIM (mg/kg))	T-						T-		1		1	I	
Acenapthene	ND (0.133)	ND (0.097)	ND (0.111)	ND (0.091)	ND (0.100)	ND (0.085)	ND (0.095)	ND (0.086)	ND (0.102)	ND (0.085)	ND (0.092)	ND (0.091)	ND (0.092)	ND (0.086)	ND (0.092)	ND (0.090)	0.156
Acenaphthylene	ND (0.133)	ND (0.097)	ND (0.111)	ND (0.091)	ND (0.100)	ND (0.085)	ND (0.095)	ND (0.086)	ND (0.102)	ND (0.085)	ND (0.092)	ND (0.091)	ND (0.092)	ND (0.086)	ND (0.092)	ND (0.090)	NA
Anthracene	ND (0.133)	ND (0.097)	ND (0.111)	ND (0.091)	ND (0.100)	ND (0.085)	ND (0.095)	ND (0.086)	ND (0.102)	ND (0.085)	ND (0.092)	0.13	0.15	ND (0.086)	ND (0.092)	ND (0.090)	7.134
Benzyl Alcohol	ND (0.166)	ND (0.121)	ND (0.139)	ND (0.113)	ND (0.124)	ND (0.107)	ND (0.119)	ND (0.108)	ND (0.128)	ND (0.106)	ND (0.115)	ND (0.114)	ND (0.115)	ND (0.108)	ND (0.115)	ND (0.113)	NA
Bis (2-Ethylhexyl) Phthalate	2.5	ND (0.121)	ND (0.139)	ND (0.113)	1.09	ND (0.107)	0.38	ND (0.108)	ND (0.128)	0.21	ND (0.115)	2.76	ND (0.115)	ND (0.108)	ND (0.115)	ND (0.113)	0.111



			Table	7 (continu	ed). Summ	nary of Soil	Sample Res	sults, Pacific	: City Park I	Remedial In	vestigation	, Pacific, W	ashington.				
								Sample	Location								
Analytical Parameter	GP-1	L6- 1 5		GP-17-15		GP-1	L8-15	GP-1	.9-15	GP-2	20-15	GP-2	21-15		GP-22-15		Site
Sample Date	9/2:	1/15		9/21/15		9/2:	1/15	9/2:	1/15	9/2:	1/15	9/2	1/15		9/22/15		Screening Level ^a
Depth (feet)	8.0	13.5	7.0	14.5	7 dup	3.0	14.0	5.0	14.5	5.5	14.0	4.0	14.0	1.5	3.0	3.0 dup	(mg/kg)
Semi-Volatile Organic C	Compounds b	y EPA 8270D	/SIM (mg/kg)	(continued)													
Butyl Benzylphthalate	0.17	ND (0.121)	ND (0.139)	ND (0.113)	ND (0.124)	ND (0.107)	ND (0.119)	0.27	ND (0.128)	ND (0.106)	ND (0.115)	0.18	ND (0.115)	ND (0.108)	ND (0.115)	ND (0.113)	0.033
Dibutyl Phthalate	ND (0.166)	ND (0.121)	ND (0.139)	ND (0.113)	ND (0.124)	ND (0.107)	ND (0.119)	ND (0.108)	ND (0.128)	ND (0.106)	ND (0.115)	ND (0.114)	ND (0.115)	ND (0.108)	ND (0.115)	ND (0.113)	0.17
Di-N-Octyl Phthalate	ND (0.166)	ND (0.121)	ND (0.139)	ND (0.113)	ND (0.124)	0.17	ND (0.119)	ND (0.108)	ND (0.128)	ND (0.106)	ND (0.115)	ND (0.114)	ND (0.115)	ND (0.108)	ND (0.115)	ND (0.113)	800
Fluoranthene	0.15	ND (0.097)	ND (0.111)	ND (0.091)	0.13	0.11	ND (0.095)	ND (0.086)	ND (0.102)	0.096	ND (0.092)	0.25	0.13	ND (0.086)	ND (0.092)	ND (0.090)	0.296
Fluorene	ND (0.133)	ND (0.097)	ND (0.111)	ND (0.091)	ND (0.100)	ND (0.085)	ND (0.095)	ND (0.086)	ND (0.102)	ND (0.085)	ND (0.092)	ND (0.091)	ND (0.092)	ND (0.086)	ND (0.092)	ND (0.090)	0.080
2-Methylnaphthalene	ND (0.133)	ND (0.097)	ND (0.111)	ND (0.091)	ND (0.100)	ND (0.085)	ND (0.095)	ND (0.086)	ND (0.102)	ND (0.085)	ND (0.092)	ND (0.091)	ND (0.092)	ND (0.086)	ND (0.092)	ND (0.090)	0.236
1-Methylnaphthalene	ND (0.133)	ND (0.097)	ND (0.111)	ND (0.091)	ND (0.100)	ND (0.085)	ND (0.095)	ND (0.086)	ND (0.102)	ND (0.085)	ND (0.092)	ND (0.091)	ND (0.092)	ND (0.086)	ND (0.092)	ND (0.090)	0.236
Naphthalene	ND (0.133)	ND (0.097)	ND (0.111)	ND (0.091)	ND (0.100)	ND (0.085)	ND (0.095)	ND (0.086)	ND (0.102)	ND (0.085)	ND (0.092)	ND (0.091)	ND (0.092)	ND (0.086)	ND (0.092)	ND (0.090)	0.236
4-Nitrophenol	ND (0.831)	ND (0.605)	ND (0.694)	ND (0.566)	ND (0.622)	ND (0.533)	ND (0.596)	ND (0.539)	ND (0.638)	ND (0.530)	ND (0.575)	ND (0.571)	ND (0.575)	ND (0.540)	ND (0.575)	ND (0.565)	NA
p-Cresol	ND (0.166)	ND (0.121)	ND (0.139)	ND (0.113)	ND (0.124)	ND (0.107)	ND (0.119)	ND (0.108)	ND (0.128)	ND (0.106)	ND (0.115)	ND (0.114)	ND (0.115)	ND (0.108)	ND (0.115)	ND (0.113)	8,000
Pentachlorophenol	ND (0.166)	ND (0.121)	ND (0.139)	ND (0.113)	ND (0.124)	ND (0.107)	ND (0.119)	ND (0.108)	ND (0.128)	ND (0.106)	ND (0.115)	ND (0.114)	ND (0.115)	ND (0.108)	ND (0.115)	ND (0.113)	0.17
Phenanthrene	ND (0.133)	ND (0.097)	ND (0.111)	ND (0.091)	ND (0.100)	ND (0.085)	ND (0.095)	ND (0.086)	ND (0.102)	ND (0.085)	ND (0.092)	0.11	0.13	ND (0.086)	ND (0.092)	ND (0.090)	0.0067
Phenol	ND (0.333)	ND (0.242)	ND (0.278)	ND (0.226)	ND (0.249)	ND (0.213)	ND (0.238)	ND (.216)	ND (0.255)	ND (0.212)	ND (0.230)	0.57	ND (0.230)	ND (0.216)	ND (0.230)	ND (0.226)	0.757
Pyrene	0.19	ND (0.097)	0.13	ND (0.091)	0.14	0.12	ND (0.095)	ND (0.086)	ND (0.102)	0.090	ND (0.092)	0.23	0.11	ND (0.086)	ND (0.092)	ND (0.090)	0.546
Carcinogenic Polycyclic	Aromatic Hy	ydrocarbons ((cPAHs) by EP	A 8270D/SIM	(mg/kg)												
Benzo(a)anthracene	0.21	ND (0.097)	0.19	ND (0.091)	ND (0.100)	0.16	ND (0.095)	ND (0.086)	ND (0.102)	0.13	ND (0.092)	0.16	0.12	ND (0.086)	ND (0.092)	ND (0.090)	0.0067
Benzo(a)pyrene	ND (0.133)	ND (0.097)	ND (0.111)	ND (0.091)	ND (0.100)	ND (0.085)	0.17	ND (0.086)	0.28	ND (0.085)	ND (0.092)	ND (0.091)	ND (0.092)	ND (0.086)	ND (0.092)	ND (0.090)	0.01
Benzo(b)fluoranthene	ND (0.133)	ND (0.097)	ND (0.111)	ND (0.091)	ND (0.100)	ND (0.085)	0.19	0.17	ND (0.102)	ND (0.085)	ND (0.092)	ND (0.091)	ND (0.092)	ND (0.086)	ND (0.092)	ND (0.090)	0.012
Benzo(j,k)fluoranthene	ND (0.133)	ND (0.097)	ND (0.111)	ND (0.091)	ND (0.100)	ND (0.085)	ND (0.095)	ND (0.086)	ND (0.102)	ND (0.085)	ND (0.092)	ND (0.091)	ND (0.092)	ND (0.086)	ND (0.092)	ND (0.090)	0.012
Chrysene	ND (0.133)	ND (0.097)	ND (0.111)	ND (0.091)	ND (0.100)	0.095	ND (0.095)	ND (0.086)	ND (0.102)	ND (0.085)	ND (0.092)	0.23	ND (0.092)	ND (0.086)	ND (0.092)	ND (0.090)	0.0067
Dibenz(a,h)anthracene	ND (0.133)	ND (0.097)	ND (0.111)	ND (0.091)	ND (0.100)	ND (0.085)	ND (0.009)	ND (0.086)	ND (0.102)	ND (0.085)	ND (0.092)	ND (0.091)	ND (0.092)	ND (0.086)	ND (0.092)	ND (0.090)	0.018
Indeno(1,2,3-cd)pyrene	ND (0.133)	ND (0.097)	ND (0.111)	ND (0.091)	ND (0.100)	ND (0.085)	ND (0.009)	ND (0.086)	ND (0.102)	ND (0.085)	ND (0.092)	ND (0.091)	ND (0.092)	ND (0.086)	ND (0.092)	ND (0.090)	0.035
Total cPAHs (TEQ) ^b	0.13	ND (0.088)	0.10	ND (0.082)	ND (0.090)	0.089	0.54	0.091	0.32	0.086	ND (0.083)	0.096	0.091	ND (0.078)	ND (0.083)	ND (0.081)	0.020
Herbicides by EPA 8151	.A (mg/kg)																
2,4-DB	ND (0.041)	ND (0.031)	ND (0.035)	ND (0.029)	ND (0.031)	ND (0.027)	ND (0.030)	ND (0.027)	ND (0.032)	ND (0.027)	ND (0.029)	ND (0.029)	ND (0.029)	ND (0.027)	ND (0.029)	ND (0.028)	640
2,4,5-T	ND (0.082)	ND (0.063)	ND (0.071)	ND (0.057)	ND (0.063)	ND (0.055)	ND (0.059)	ND (0.055)	ND (0.064)	ND (0.054)	ND (0.058)	ND (0.059)	ND (0.059)	ND (0.055)	ND (0.058)	ND (0.056)	800
Bentazon	ND (0.082)	ND (0.063)	ND (0.071)	ND (0.057)	ND (0.063)	ND (0.055)	ND (0.059)	ND (0.055)	ND (0.064)	ND (0.054)	ND (0.058)	ND (0.059)	ND (0.059)	ND (0.055)	ND (0.058)	ND (0.056)	2,400
Chloramben	ND (0.033)	ND (0.025)	ND (0.028)	ND (0.023)	ND (0.025)	ND (0.022)	ND (0.024)	ND (0.022)	ND (0.026)	ND (0.022)	ND (0.023)	ND (0.023)	ND (0.024)	ND (0.022)	ND (0.023)	ND (0.022)	1,200
Chlorthal-dimethyl	ND (0.041)	ND (0.031)	ND (0.035)	ND (0.029)	ND (0.031)	ND (0.027)	ND (0.030)	ND (0.027)	ND (0.032)	ND (0.027)	ND (0.029)	ND (0.029)	ND (0.029)	ND (0.027)	ND (0.029)	ND (0.028)	800
Dalapon	ND (0.033)	ND (0.025)	ND (0.028)	ND (0.023)	ND (0.025)	ND (0.022)	ND (0.024)	ND (0.022)	ND (0.026)	ND (0.022)	ND (0.023)	ND (0.023)	ND (0.024)	ND (0.022)	ND (0.023)	ND (0.022)	2,400
Dinoseb	ND (0.082)	ND (0.063)	ND (0.071)	ND (0.057)	ND (0.063)	ND (0.055)	ND (0.059)	ND (0.055)	ND (0.064)	ND (0.054)	ND (0.058)	ND (0.059)	ND (0.059)	ND (0.055)	ND (0.058)	ND (0.056)	80
Picloram	ND (0.082)	ND (0.063)	ND (0.071)	ND (0.057)	ND (0.063)	ND (0.055)	ND (0.059)	ND (0.055)	ND (0.064)	ND (0.054)	ND (0.058)	ND (0.059)	ND (0.059)	ND (0.055)	ND (0.058)	ND (0.056)	5,600
Silvex	ND (0.033)	ND (0.025)	ND (0.028)	ND (0.023)	ND (0.025)	ND (0.022)	ND (0.024)	ND (0.022)	ND (0.026)	ND (0.022)	ND (0.023)	ND (0.023)	ND (0.024)	ND (0.022)	ND (0.023)	ND (0.022)	640



			Table	7 (continu	ed). Summ	ary of Soil	Sample Res	sults, Pacific	City Park F	Remedial In	vestigation	, Pacific, W	ashington.				
								Sample	Location								
Analytical Parameter	GP-1	L6-15		GP-17-15		GP-1	18-15	GP-1	L9-15	GP-2	20-15	GP-2	21-15		GP-22-15		Site Screening
Sample Date	9/2	9/21/15 9/21/15 13.5 7.0 14.5 7 dup				9/2	1/15	9/2	1/15	9/2	1/15	9/2	1/15		9/22/15		Level ^a
Depth (feet)	8.0	13.5	7.0	14.5	7 dup	3.0	14.0	5.0	14.5	5.5	14.0	4.0	14.0	1.5	3.0	3.0 dup	(mg/kg)
Organochlorine Pestici	des by EPA 8	081 (mg/kg)															
4,4'-DDD	ND (0.029)	ND (0.025)	ND (0.027)	ND (0.021)	ND (0.025)	ND (0.021)	ND (0.022)	ND (0.020)	ND (0.023)	ND (0.019)	ND (0.023)	ND (0.022)	ND (0.021)	ND (0.021)	ND (0.020)	ND (0.022)	0.01
4,4'-DDE	ND (0.029)	ND (0.025)	ND (0.027)	ND (0.021)	ND (0.025)	ND (0.021)	ND (0.022)	ND (0.020)	ND (0.023)	ND (0.019)	ND (0.023)	ND (0.022)	ND (0.021)	ND (0.021)	ND (0.020)	ND (0.022)	0.01
Endosulfan I	ND (0.015)	ND (0.012)	ND (0.013)	ND (0.011)	ND (0.013)	ND (0.010)	ND (0.011)	ND (0.010)	ND (0.011)	ND (0.010)	ND (0.011)	ND (0.011)	ND (0.010)	ND (0.011)	ND (0.010)	ND (0.011)	0.01
Endosulfan II	ND (0.029)	ND (0.025)	ND (0.027)	ND (0.021)	ND (0.025)	ND (0.021)	ND (0.022)	ND (0.020)	ND (0.023)	ND (0.019)	ND (0.023)	ND (0.022)	ND (0.021)	ND (0.021)	ND (0.020)	ND (0.022)	0.005
Cis-Chlordane (alpha)	ND (0.015)	ND (0.012)	ND (0.013)	ND (0.011)	ND (0.013)	ND (0.010)	ND (0.011)	ND (0.010)	ND (0.011)	ND (0.010)	ND (0.011)	ND (0.011)	ND (0.010)	ND (0.011)	ND (0.010)	ND (0.011)	0.01
Endosulfan Sulfate	ND (0.029)	ND (0.025)	ND (0.027)	ND (0.021)	ND (0.025)	ND (0.021)	ND (0.022)	ND (0.020)	ND (0.023)	ND (0.019)	ND (0.023)	ND (0.022)	ND (0.021)	ND (0.021)	ND (0.020)	ND (0.022)	NA
Gamma-Chlordane	ND (0.015)	ND (0.012)	ND (0.013)	ND (0.011)	ND (0.013)	ND (0.010)	ND (0.011)	ND (0.010)	ND (0.011)	ND (0.010)	ND (0.011)	ND (0.011)	ND (0.010)	ND (0.011)	ND (0.010)	ND (0.011)	0.01
Methoxychlor	ND (0.073)	ND (0.062)	ND (0.067)	ND (0.052)	ND (0.064)	ND (0.052)	ND (0.054)	ND (0.051)	ND (0.057)	ND (0.047)	ND (0.057)	ND (0.056)	ND (0.052)	ND (0.053)	ND (0.0500)	ND (0.054)	0.01

BOLD values detected above the reporting limit.

Shaded values exceed the site screening level

Native Soil
Fill/Refuse

King County

			Table 7	(continued).	Summary	of Soil Sam	ple Results,	Pacific City	Park Remed	ial Investiga	tion, Pacific	, Washingto	n.			
								Sample Locati	on							
Analytical Parameter	MW-	1-15	MW-	-2-15	MW	-3-15	MW-	-4-15	MW	-5-15		MW-6-15		TP-1-15	TP-3-15	Site Screening
Sample Date	9/18	3/15	9/18	3/15	9/1	8/15	9/22	2/15	9/2	2/15		9/22/15		9/23/15	9/23/15	Level ^a
Depth (feet)	6.0	12.0	3.0	13	6.5	15.0	6.0	14.5	5.0	12.5	5.5	5.5 dup	14.0	5.5	6.0	(mg/kg)
Petroleum Hydrocarbor	ns (mg/kg)															
Gasoline Range Organics	ND (2.7)	ND (3.3)	ND (3.3)	ND (5.5)	ND (3.88)	ND (3.69)	ND (2.66)	ND (4.89)	ND (3.79)	ND (3.57)	ND (3.59)	ND (4.29)	ND (3.88)	ND (5.83)	ND (3.66)	100
Diesel Range Organics	ND (24.0)	ND (23.9)	ND (21.9)	ND (37.8)	ND (23)	ND (26)	ND (22)	ND (29)	ND (24)	ND (24)	ND (20)	ND (20)	ND (20)	ND (27)	ND (24)	200
Lube Oil Range Organics	ND (60)	ND (60)	137	ND (95)	ND (59)	ND (64)	1,250	ND (74)	ND (61)	ND (61)	ND (50)	4,080	ND (51)	895	267	2,000
Volatile Organic Compo	ounds by EPA	8260 (mg/kg)														
Benzene	ND (0.011)	ND (0.013)	ND (0.013)	ND (0.022)	ND (0.016)	ND (0.015)	ND (0.011)	ND (0.020)	ND (0.015)	ND (0.014)	ND (0.014)	ND (0.016)	ND (0.017)	ND (0.023)	ND (0.015)	0.001
Toluene	ND (0.011)	ND (0.013)	ND (0.013)	ND (0.022)	ND (0.016)	ND (0.015)	ND (0.011)	ND (0.020)	ND (0.015)	ND (0.014)	ND (0.014)	ND (0.016)	ND (0.017)	ND (0.023)	ND (0.015)	0.024
Ethylbenzene	ND (0.016)	ND (0.020)	ND (0.020)	ND (0.033)	ND (0.023)	ND (0.022)	ND (0.016)	ND (0.029)	ND (0.023)	ND (0.021)	ND (0.022)	ND (0.023)	ND (0.026)	ND (0.035)	ND (0.022)	0.014
Total Xylenes	ND (0.011)	ND (0.013)	ND (0.013)	ND (0.022)	ND (0.016)	ND (0.015)	ND (0.011)	ND (0.020)	ND (0.015)	ND (0.014)	ND (0.014)	ND (0.016)	ND (0.017)	ND (0.023)	ND (0.015)	0.52
Acetone	-	-	_	-	-	_	-	_	-	-	_	_	_	-	_	2.07
2-Butanone	-	-	_	-	-	_	_	_	_	_	_	_	_	-	_	1.38
Carbon Disulfide	-	-	_	_	-	_	_	_	_	_	_	_	_	_	_	0.27
Cis-1,2-dichloroethene	ND (0.011)	ND (0.013)	ND (0.013)	ND (0.022)	ND (0.016)	ND (0.015)	ND (0.011)	ND (0.020)	ND (0.015)	ND (0.014)	ND (0.014)	ND (0.016)	ND (0.017)	ND (0.023)	ND (0.015)	0.005
Chlorobenzene	ND (0.011)	ND (0.013)	ND (0.013)	ND (0.022)	ND (0.016)	ND (0.015)	ND (0.011)	ND (0.020)	ND (0.015)	ND (0.014)	ND (0.014)	ND (0.016)	ND (0.017)	ND (0.023)	ND (0.015)	0.051
Methylene Chloride	ND (0.011)	ND (0.013)	ND (0.013)	ND (0.022)	ND (0.016)	ND (0.015)	ND (0.011)	ND (0.020)	ND (0.015)	ND (0.014)	ND (0.014)	ND (0.016)	ND (0.017)	ND (0.023)	ND (0.015)	0.005
p-Isopropyltoluene	ND (0.011)	ND (0.013)	ND (0.013)	ND (0.022)	ND (0.016)	ND (0.015)	ND (0.011)	ND (0.020)	ND (0.015)	ND (0.014)	ND (0.014)	ND (0.016)	ND (0.017)	ND (0.023)	ND (0.015)	0.229
Styrene	ND (0.012)	ND (0.014)	ND (0.013)	ND (0.013)	ND (0.013)	ND (0.030)	ND (0.012)	ND (0.014)	ND (0.017)	ND (0.013)	ND (0.014)	ND (0.022)	ND (0.013)	ND (0.015)	ND (0.015)	0.120
Tetrachlorethene	ND (0.011)	ND (0.013)	ND (0.013)	ND (0.022)	ND (0.016)	ND (0.015)	ND (0.011)	ND (0.020)	ND (0.015)	ND (0.014)	ND (0.014)	ND (0.016)	ND (0.017)	ND (0.023)	ND (0.015)	0.0013
Trichloroethene	ND (0.011)	ND (0.013)	ND (0.013)	ND (0.022)	ND (0.016)	ND (0.015)	ND (0.011)	ND (0.020)	ND (0.015)	ND (0.014)	ND (0.014)	ND (0.016)	ND (0.017)	ND (0.023)	ND (0.015)	0.001
1,2,4-Trimethylbenzene	ND (0.012)	ND (0.014)	ND (0.013)	ND (0.013)	ND (0.013)	ND (0.030)	ND (0.012)	ND (0.014)	ND (0.017)	ND (0.013)	ND (0.014)	ND (0.022)	ND (0.013)	ND (0.015)	ND (0.015)	NA
Total Metals by EPA 603	10D/7471B (m	ıg/kg)														
Arsenic	3.7	2.7	3.1	8.1	3.8	1.6	4	4.5	4.6	2.2	3.5	3.3	3.9	5.2	17	20
Barium	31	20	28	74	17	27	45	80	98	30	39	29	28	78	315	41.3
Cadmium	0.44	ND (0.20)	ND (0.17)	ND (0.28)	ND (0.21)	ND (0.23)	ND (0.19)	ND (0.25)	5.4	ND (0.19)	ND (0.17)	ND (0.19)	ND (0.17)	0.61	1.4	1
Chromium	17	14	15	22	12	12	26	21	24	15	18	12	13	22	30	48
Lead	41.1	3.4	23	4.9	1.3	1.5	19	3.8	56	7.7	12	14	2.0	88	364	25
Mercury	ND (0.26)	ND (0.28)	ND (0.24)	ND (0.44)	ND (0.30)	ND (0.32)	ND (0.28)	ND (0.34)	ND (0.27)	ND (0.28)	ND (0.27)	ND (0.29)	ND (0.28)	ND (0.35)	ND (0.30)	0.07
Selenium	1.1	1.0	0.89	2.35	0.86	0.98	1.2	2.5	1.6	1.3	1.5	1.1	1.4	1.0	1.2	10
Silver	ND (0.094)	ND (0.10)	ND (0.084)	ND (0.14)	ND (0.10)	ND (0.11)	ND (0.094)	ND (0.13)	ND (0.092)	ND (0.094)	ND (0.087)	ND (0.095)	ND (0.087)	0.20	0.47	0.61
Polychlorinated Biphen	yls (PCBs) by	EPA 8082A (m	ıg/kg)													
Total PCBs	-	-	_	-	-	_	-	_	-	_	_	ND (0.11)	_	ND (0.13)	0.23	0.05
Semi-Volatile Organic C	Compounds by	EPA 8270D/S	SIM (mg/kg)													
Acenapthene	ND (0.094)	ND (0.098)	ND (0.091)	ND (0.152)	ND (0.099)	ND (0.108)	0.093	ND (0.123)	ND (0.096)	ND (0.097)	ND (0.089)	ND (0.095)	ND (0.094)	ND (0.110)	ND (0.098)	0.156
Acenaphthylene	ND (0.094)	ND (0.098)	ND (0.091)	ND (0.152)	ND (0.099)	ND (0.108)	ND (0.092)	ND (0.123)	ND (0.096)	ND (0.097)	ND (0.089)	ND (0.095)	ND (0.094)	ND (0.110)	ND (0.098)	NA
Anthracene	ND (0.094)	ND (0.098)	ND (0.091)	ND (0.152)	ND (0.099)	ND (0.108)	0.84	ND (0.123)	ND (0.096)	ND (0.097)	ND (0.089)	ND (0.095)	ND (0.094)	ND (0.110)	ND (0.098)	7.134
Benzyl Alcohol	ND (0.118)	ND (0.122)	ND (0.114)	ND (0.190)	ND (0.124)	ND (0.135)	ND (0.114)	ND (0.154)	ND (0.120)	ND (0.121)	ND (0.111)	ND (0.118)	ND (0.119)	ND (0.138)	ND (0.122)	NA
Bis (2-Ethylhexyl) Phthalate	ND (0.118)	ND (0.122)	0.12	0.33	ND (0.124)	ND (0.135)	ND (0.114)	ND (0.154)	ND (0.120)	ND (0.121)	ND (0.111)	ND (0.118)	ND (0.119)	ND (0.138)	ND (0.122)	0.111



			Table 7	(continued)	. Summary	of Soil Sam	ple Results,	Pacific City	Park Remed	ial Investiga	tion, Pacific,	, Washingto	n.			
								Sample Location	on							
Analytical Parameter	MW-	-1-15	MW-	-2-15	MW-	-3-15	MW-	-4-15	MW	-5-15		MW-6-15		TP-1-15	TP-3-15	Site
Sample Date	9/18	8/15	9/18	8/15	9/18	8/15	9/2:	2/15	9/2	2/15		9/22/15		9/23/15	9/23/15	Screening Level ^a
Depth (feet)	6.0	12.0	3.0	13	6.5	15.0	6.0	14.5	5.0	12.5	5.5	5.5 dup	14.0	5.5	6.0	(mg/kg)
Semi-Volatile Organic C	ompounds by	y EPA 8270D/S	SIM (mg/kg) (c	ontinued)												
Butyl Benzylphthalate	ND (0.118)	ND (0.122)	ND (0.114)	ND (0.190)	ND (0.124)	ND (0.135)	ND (0.114)	ND (0.154)	ND (0.120)	ND (0.121)	ND (0.111)	ND (0.118)	ND (0.119)	ND (0.138)	ND (0.122)	0.033
Dibutyl Phthalate	ND (0.118)	ND (0.122)	ND (0.114)	ND (0.190)	ND (0.124)	ND (0.135)	0.14	ND (0.154)	0.14	0.15	0.13	0.16	0.14	0.31	0.14	0.17
Di-N-Octyl Phthalate	ND (0.118)	ND (0.122)	ND (0.114)	ND (0.190)	ND (0.124)	ND (0.135)	ND (0.114)	ND (0.154)	ND (0.120)	ND (0.121)	ND (0.111)	ND (0.118)	ND (0.119)	ND (0.138)	ND (0.122)	800
Fluoranthene	ND (0.094)	ND (0.098)	0.16	ND (0.152)	ND (0.099)	ND (0.108)	0.47	ND (0.123)	ND (0.096)	ND (0.097)	ND (0.089)	0.11	ND (0.094)	ND (0.110)	0.15	0.296
Fluorene	ND (0.094)	ND (0.098)	ND (0.091)	ND (0.152)	ND (0.099)	ND (0.108)	0.18	ND (0.123)	ND (0.096)	ND (0.097)	ND (0.089)	ND (0.095)	ND (0.094)	ND (0.110)	ND (0.098)	0.080
1-Methylnaphthalene	ND (0.094)	ND (0.098)	ND (0.091)	ND (0.152)	ND (0.099)	ND (0.108)	0.18	ND (0.123)	ND (0.096)	ND (0.097)	ND (0.089)	ND (0.095)	ND (0.094)	ND (0.110)	ND (0.098)	0.236
2-Methylnaphthalene	ND (0.094)	ND (0.098)	ND (0.091)	ND (0.152)	ND (0.099)	ND (0.108)	0.16	ND (0.123)	ND (0.096)	ND (0.097)	ND (0.089)	ND (0.095)	ND (0.094)	ND (0.110)	ND (0.098)	0.236
Naphthalene	ND (0.094)	ND (0.098)	ND (0.091)	ND (0.152)	ND (0.099)	ND (0.108)	ND (0.092)	ND (0.123)	ND (0.096)	ND (0.097)	ND (0.089)	ND (0.095)	ND (0.094)	ND (0.110)	ND (0.098)	0.236
4-Nitrophenol	ND (0.590)	ND (0.612)	ND (0.570)	ND (0.952)	ND (0.618)	ND (0.618)	0.80	ND (0.0445)	ND (0.036)	ND (0.038)	ND (0.553)	ND (0.594)	ND (0.588)	0.041	0.049	NA
p-Cresol	ND (0.118)	ND (0.122)	ND (0.114)	ND (0.190)	ND (0.124)	ND (0.135)	ND (0.114)	ND (0.154)	ND (0.120)	ND (0.121)	ND (0.111)	ND (0.118)	ND (0.119)	ND (0.138)	ND (0.122)	8,000
Pentachlorophenol	ND (0.118)	ND (0.122)	ND (0.114)	ND (0.190)	ND (0.124)	ND (0.135)	ND (0.114)	ND (0.154)	ND (0.120)	ND (0.121)	ND (0.111)	ND (0.118)	ND (0.119)	ND (0.138)	ND (0.122)	0.17
Phenanthrene	ND (0.094)	ND (0.098)	ND (0.091)	ND (0.152)	ND (0.099)	ND (0.108)	0.82	ND (0.123)	ND (0.096)	ND (0.097)	ND (0.089)	ND (0.095)	ND (0.094)	ND (0.110)	ND (0.098)	0.0067
Phenol	ND (0.236)	ND (0.245)	ND (0.228)	ND (0.381)	ND (0.247)	ND (0.270)	ND (0.229)	ND (0.308)	ND (0.241)	ND (0.241)	ND (0.221)	ND (0.235)	ND (0.238)	ND (0.276)	ND (0.244)	0.757
Pyrene	ND (0.094)	ND (0.098)	0.15	ND (0.152)	ND (0.099)	ND (0.108)	0.72	ND (0.123)	ND (0.096)	ND (0.097)	ND (0.089)	ND (0.095)	ND (0.094)	ND (0.110)	0.14	0.546
Carcinogenic Polycyclic	Aromatic Hy	drocarbons (cl	PAHs) by EPA 8	3270D/SIM (mg	g/kg)											
Benzo(a)anthracene	0.13	ND (0.098)	0.14	ND (0.152)	ND (0.099)	ND (0.108)	0.33	ND (0.123)	0.13	ND (0.097)	ND (0.089)	ND (0.095)	ND (0.094)	0.16	ND (0.098)	0.0067
Benzo(a)pyrene	ND (0.094)	ND (0.098)	ND (0.091)	0.52	ND (0.099)	ND (0.108)	0.20	0.27	0.17	ND (0.097)	ND (0.089)	ND (0.095)	ND (0.094)	ND (0.110)	ND (0.098)	0.01
Benzo(b)fluoranthene	ND (0.094)	ND (0.098)	0.26	ND (0.152)	ND (0.099)	ND (0.108)	0.35	ND (0.123)	ND (0.096)	ND (0.097)	ND (0.089)	ND (0.095)	ND (0.094)	ND (0.110)	0.22	0.012
Benzo(j,k)fluoranthene	ND (0.094)	ND (0.098)	0.11	ND (0.152)	ND (0.099)	ND (0.108)	ND (0.092)	ND (0.123)	ND (0.096)	ND (0.097)	ND (0.089)	ND (0.095)	ND (0.094)	ND (0.110)	ND (0.098)	0.012
Chrysene	ND (0.094)	ND (0.098)	ND (0.091)	ND (0.152)	ND (0.099)	ND (0.108)	0.30	ND (0.123)	ND (0.096)	ND (0.097)	ND (0.089)	ND (0.095)	ND (0.094)	ND (0.110)	ND (0.098)	0.0067
Dibenz(a,h)anthracene	ND (0.094)	ND (0.098)	ND (0.091)	ND (0.152)	ND (0.099)	ND (0.108)	ND (0.008)	ND (0.123)	ND (0.096)	ND (0.097)	ND (0.089)	ND (0.095)	ND (0.094)	ND (0.110)	ND (0.098)	0.018
Indeno(1,2,3-cd)pyrene	ND (0.094)	ND (0.098)	ND (0.091)	ND (0.152)	ND (0.099)	ND (0.108)	ND (0.008)	ND (0.123)	ND (0.096)	ND (0.097)	ND (0.089)	ND (0.095)	ND (0.094)	ND (0.110)	ND (0.098)	0.035
Total cPAHs (TEQ) ^b	0.093	ND (0.089)	0.10	0.58	ND (0.089)	ND (0.098)	0.28	0.27	0.18	ND (0.087)	ND (0.080)	ND (0.085)	ND (0.086)	0.10	0.099	0.020
Herbicides by EPA 8151	A (mg/kg)															
2,4-DB	ND (0.028)	ND (0.030)	ND (0.028)	ND (0.047)	ND (0.031)	ND (0.034)	ND (0.028)	ND (0.037)	ND (0.030)	ND (0.032)	ND (0.027)	0.022	0.019	0.024	ND (0.015)	640
2,4,5-T	ND (0.056	ND (0.059	ND (0.056	ND (0.094	ND (0.063)	ND (0.068)	ND (0.056)	ND (0.074)	ND (0.061)	ND (0.063)	ND (0.054)	0.071	ND (0.028)	ND (0.034)	ND (0.031)	800
Bentazon	ND (0.056	ND (0.059	ND (0.056	ND (0.094	ND (0.056	ND (0.068)	ND (0.056)	ND (0.074)	ND (0.061)	ND (0.063)	ND (0.054)	0.071	ND (0.028)	ND (0.034)	ND (0.031)	2,400
Chloramben	ND (0.023)	ND (0.024)	ND (0.022)	ND (0.038)	ND (0.025)	ND (0.027)	ND (0.022)	ND (0.030)	ND (0.024)	ND (0.025)	ND (0.022)	0.029	0.021	0.077	0.11	1,200
Chlorthal-dimethyl	ND (0.028)	ND (0.030)	ND (0.028)	ND (0.047)	ND (0.031)	ND (0.034)	ND (0.028)	ND (0.037)	ND (0.030)	ND (0.032)	ND (0.027)	0.022	0.019	0.024	ND (0.015)	800
Dalapon	ND (0.023)	ND (0.024)	ND (0.022)	ND (0.038)	ND (0.025)	ND (0.027)	ND (0.022)	ND (0.030)	ND (0.024)	ND (0.025)	ND (0.022)	0.029	0.021	0.077	0.11	2,400
Dinoseb	ND (0.056	ND (0.059	ND (0.056	ND (0.094	ND (0.056	ND (0.068)	ND (0.056)	ND (0.074)	ND (0.061)	ND (0.063)	ND (0.054)	0.040	0.028	0.045	0.032	80
Picloram	ND (0.056)	ND (0.059)	ND (0.056)	ND (0.094)	ND (0.063)	ND (0.068)	ND (0.056)	ND (0.074)	ND (0.061)	ND (0.063)	ND (0.054)	0.040	0.028	0.045	0.032	5,600
Silvex	ND (0.023	ND (0.024	ND (0.022	ND (0.038	ND (0.031)	ND (0.034)	ND (0.028)	ND (0.037)	ND (0.030)	ND (0.032)	ND (0.027)	0.022	0.019	0.024	ND (0.015)	640



			Table 7	(continued)	. Summary	of Soil Sam	ple Results,	Pacific City	Park Remed	ial Investiga	tion, Pacific,	Washingto	n.			
								Sample Location	n							
Analytical Parameter	MW-	1-15	MW-	2-15	MW-	3-15	MW-	4-15	MW-	-5-15		MW-6-15		TP-1-15	TP-3-15	Site Screening
Sample Date	9/18	3/15	9/18	3/15	9/18	3/15	9/2	2/15	9/22	2/15		9/22/15		9/23/15	9/23/15	Level ^a
Depth (feet)	6.0	12.0	3.0	13	6.5	15.0	6.0	14.5	5.0	12.5	5.5	5.5 dup	14.0	5.5	6.0	(mg/kg)
Organochlorine Pesticio	des by EPA 80	81 (mg/kg)														
4,4'-DDD	ND (0.024)	ND (0.023)	ND (0.022)	ND (0.035)	ND (0.024)	ND (0.025)	ND (0.020)	ND (0.030)	ND (0.023)	ND (0.024)	ND (0.022)	ND (0.023)	ND (0.021)	ND (0.024)	ND (0.023)	0.01
4,4'-DDE	ND (0.024)	ND (0.023)	ND (0.022)	ND (0.035)	ND (0.024)	ND (0.025)	ND (0.020)	ND (0.030)	ND (0.023)	ND (0.024)	ND (0.022)	ND (0.023)	ND (0.021)	ND (0.024)	ND (0.023)	0.01
Cis-Chlordane (alpha)	ND (0.012)	ND (0.012)	ND (0.011)	ND (0.018)	ND (0.012)	ND (0.013)	ND (0.010)	ND (0.015)	ND (0.012)	ND (0.012)	ND (0.011)	ND (0.012)	ND (0.011)	ND (0.012)	ND (0.012)	0.01
Endosulfan I	ND (0.012)	ND (0.012)	ND (0.011)	ND (0.018)	ND (0.012)	ND (0.013)	ND (0.010)	ND (0.015)	ND (0.012)	ND (0.012)	ND (0.011)	ND (0.012)	ND (0.011)	ND (0.012)	ND (0.012)	0.005
Endosulfan II	ND (0.024)	ND (0.023)	ND (0.022)	ND (0.035)	ND (0.024)	ND (0.025)	ND (0.020)	ND (0.030)	ND (0.023)	ND (0.024)	ND (0.022)	ND (0.023)	ND (0.021)	ND (0.024)	ND (0.023)	0.01
Endosulfan Sulfate	ND (0.024)	ND (0.023)	ND (0.022)	ND (0.035)	ND (0.024)	ND (0.025)	ND (0.020)	ND (0.030)	ND (0.023)	ND (0.024)	ND (0.022)	ND (0.023)	ND (0.021)	ND (0.024)	ND (0.023)	NA
Gamma-Chlordane	ND (0.012)	ND (0.012)	ND (0.011)	ND (0.018)	ND (0.012)	ND (0.013)	ND (0.010)	ND (0.015)	ND (0.012)	ND (0.012)	ND (0.011)	ND (0.012)	ND (0.011)	ND (0.012)	ND (0.012)	0.01
Methoxychlor	ND (0.059)	ND (0.058)	ND (0.054)	ND (0.088)	ND (0.059)	ND (0.063)	ND (0.049)	ND (0.075)	ND (0.058)	ND (0.060)	ND (0.056)	ND (0.058)	ND (0.053)	ND (0.059)	ND (0.058)	0.01

BOLD values detected above the reporting limit.

Shaded values exceed the site screening level





			Tal	ole 7 (cont	tinued). S	Summary	of Soil Sar	nple Resu	lts, Pacific	City Park	Remedial	Investigat	tion, Pacif	ic, Washin	gton.				
									Sample	Location									
Analytical Parameter	PI	P1	PI	2	PI	P3	P	P4	P	P5	PI	P6	PI	P7	PI	P8	PI	> 9	Site Screening
Sample Date	5/23,	/2017	5/23/	2017	5/24/	/2017	5/24,	/2017	5/23,	/2017	5/23,	/2017	5/24/	/2017	5/24,	/2017	5/24/	2017	Level ^a
Depth (feet)	2.0	10	2.5	10	2.5	10	3.0	10	2.5	10	2.0	10	0	10	0.5	10	0	10	(mg/kg)
Petroleum Hydrocarbon	s by Metho	ds NWTPH-	Gx and NWT	PH-Dx (mg/	/kg)														
Gasoline	ND (6.9)	ND (8.0)	ND (7.8)	ND (9.0)	ND (6.6)	ND (7.0)	ND (7.4)	ND (7.1)	ND (7.0)	ND (8.1)	ND (5.7)	ND (9.7)	ND (6.6)	ND (7.8)	ND (6.1)	ND (9.6)	ND (5.6)	ND (7.0)	100
Diesel Range Organics	ND (31)	ND (33)	ND (33)	ND (36)	ND (31)	ND (31)	ND (54)	ND (55)	ND (68)	ND (34)	ND (29)	ND (37)	ND (31)	ND (31)	ND (28)	ND (36)	ND (26)	ND (31)	200
Lube Oil Range Organics	ND (62)	ND (67)	ND (66)	ND (73)	140	ND (62)	500	570	620	ND (67)	62	ND (75)	63	ND (63)	ND (57)	ND (71)	ND (52)	ND (62)	2,000
Volatile Organic Compo	unds by EP	A 8260C (mg	J/kg)																
Benzene	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.002)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	0.001
Toluene	0.019	0.020	0.025	0.019	ND (0.006)	0.016	0.016	0.011	ND (0.006)	0.016	0.010	0.008	ND (0.007)	0.022	0.014	0.021	0.014	0.015	0.024
Ethylbenzene	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.002)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	0.014
Total Xylenes	ND (0.002)	ND (0.003)	ND (0.003)	ND (0.003)	ND (0.002)	ND (0.003)	ND (0.002)	ND (0.003)	ND (0.002)	ND (0.003)	ND (0.002)	ND (0.003)	ND (0.003)	ND (0.003)	ND (0.003)	ND (0.003)	ND (0.003)	ND (0.003)	0.52
Acetone	ND (0.011)	0.051	0.071	0.088	0.058	0.028	ND (0.011)	0.051	ND (0.012)	0.046	0.011	0.10	ND (0.014)	0.019	0.027	0.016	ND (0.013)	0.015	2.07
2-Butanone	ND (0.006)	0.012	0.019	0.023	0.013	ND (0.006)	ND (0.006)	0.007	ND (0.006)	0.013	ND (0.005)	0.025	ND (0.007)	ND (0.007)	ND (0.007)	ND (0.007)	ND (0.006)	ND (0.006)	1.38
Carbon Disulfide	ND (0.002)	ND (0.002)	ND (0.001)	ND (0.002)	ND (0.002)	ND (0.002)	ND (0.002)	ND (0.001)	ND (0.006)	0.002	ND (0.001)	ND (0.002)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	0.27
Cis-1,2-dichloroethene	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.002)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	0.005
Chlorobenzene	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.002)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	0.051
Methylene Chloride	ND (0.011)	ND (0.013)	ND (0.012)	ND (0.014)	ND (0.012)	ND (0.012)	ND (0.011)	ND (0.013)	ND (0.012)	ND (0.013)	ND (0.001)	ND (0.015)	ND (0.014)	ND (0.013)	ND (0.013)	ND (0.014)	ND (0.013)	ND (0.012)	0.005
p-Isopropyltoluene	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.069)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.002)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.057)	ND (0.001)	0.229
Styrene	0.019	0.020	0.025	0.019	ND (0.006)	0.016	0.016	0.011	ND (0.006)	0.016	0.010	0.008	ND (0.007)	0.022	0.014	0.021	0.014	0.015	0.120
Tetrachloroethene	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.002)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	0.0013
Trichloroethene	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.002)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	0.001
1,2,4-Trimethylbenzene	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.001)	ND (0.002)	0.012	ND (0.001)	NA				
Total Metals by EPA 601	LOD/7471B ((mg/kg)																	
Arsenic	ND (12)	ND (13)	ND (13)	ND (15)	ND (13)	ND (12)	ND (13)	ND (12)	ND (13)	ND (13)	ND (12)	ND (15)	ND (12)	ND (13)	ND (11)	ND (14)	ND (10)	ND (12)	20
Cadmium	ND (0.62)	ND (0.67)	ND (0.66)	ND (0.73)	ND (0.63)	ND (0.62)	ND (0.63)	ND (0.62)	ND (0.65)	ND (0.67)	ND (0.58)	ND (0.74)	ND (0.61)	ND (0.63)	ND (0.57)	ND (0.71)	ND (0.52)	ND (0.62)	1
Chromium	20	13	11	11	27	9.5	29	13	23	13	29	14	12	11	13	13	25	14	48
Lead	8.9	ND (6.7)	ND (6.6)	ND (7.3)	25	ND (6.2)	84	ND (6.2)	27	ND (6.7)	9.7	ND (7.4)	ND (6.1)	ND (6.3)	ND (5.7)	ND (7.1)	ND (5.2)	ND (6.2)	25
Mercury	ND (0.31)	ND (0.33)	ND (0.33)	ND (0.36)	ND (0.31)	ND (0.31)	ND (0.32)	ND (0.31)	ND (0.33)	ND (0.34)	ND (0.29)	ND (0.37)	ND (0.31)	ND (0.31)	ND (0.28)	ND (0.36)	ND (0.26)	ND (0.31)	0.07
Polychlorinated Bipheny	yls (PCBs) b	y EPA 8082A	(mg/kg)																
Total PCBs	-	_	_	-	ND (0.063)	-	0.18	ND (0.062)	ND (0.065)	_	0.12	_	ND (0.061)	_	_	-	-	-	0.05
Semi-Volatile Organic C	ompounds	by EPA 8270	D/SIM (mg/	kg)															
Acenapthene	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.010)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.008)	ND (0.010)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.010)	ND (0.007)	ND (0.008)	0.156
Acenaphthylene	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.010)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.008)	ND (0.010)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.010)	ND (0.007)	ND (0.008)	NA
Anthracene	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.010)	ND (0.008)	ND (0.008)	0.009	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.008)	ND (0.010)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.010)	ND (0.007)	ND (0.008)	7.134
Benzyl Alcohol	ND (0.210)	ND (0.220)	ND (0.220)	ND (0.240)	ND (0.210)	ND (0.210)	ND (0.210)	ND (0.210)	ND (0.220)	ND (0.220)	ND (0.190)	ND (0.250)	ND (0.200)	ND (0.210)	0.70	ND (0.240)	ND (0.170)	ND (0.210)	NA
Bis (2-Ethylhexyl) Phthalate	0.052	ND (0.045)	ND (0.044)	ND (0.048)	ND (0.042)	ND (0.041)	ND (0.042)	ND (0.041)	ND (0.044)	ND (0.045)	ND (0.039)	ND (0.050)	ND (0.041)	ND (0.042)	ND (0.038)	ND (0.047)	ND (0.035)	ND (0.041)	0.111
Butyl Benzylphthalate	ND (0.041)	ND (0.045)	ND (0.044)	ND (0.048)	ND (0.042)	ND (0.041)	ND (0.042)	ND (0.041)	ND (0.044)	ND (0.045)	ND (0.039)	ND (0.050)	ND (0.041)	ND (0.042)	ND (0.038)	ND (0.047)	ND (0.035)	ND (0.041)	0.033
Dibutyl Phthalate	ND (0.21)	ND (0.22)	ND (0.22)	ND (0.24)	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.22)	ND (0.22)	ND (0.19)	ND (0.25)	ND (0.20)	ND (0.21)	ND (0.19)	ND (0.24)	ND (0.17)	ND (0.21)	0.17

Remedial Investigation Report—Pacific City Park

			Tal	ble 7 (con	tinued). S	Summary o	of Soil Sar	nple Resu	lts, Pacific	City Park	Remedial	Investiga	tion, Pacif	ic, Washin	gton.				
									Sample	Location									
Analytical Parameter	P	P1	PI	P2	PI	23	Р	P4	PI	P5	P	P6	PI	P7	PI	P8	P	P9	Site Screening
Sample Date	5/23,	/2017	5/23/	/2017	5/24/	2017	5/24,	/2017	5/23,	2017	5/23,	/2017	5/24,	/2017	5/24/	/2017	5/24,	/2017	Level ^a
Depth (feet)	2.0	10	2.5	10	2.5	10	3.0	10	2.5	10	2.0	10	0	10	0.5	10	0	10	(mg/kg)
Semi-Volatile Organic C	Compounds	by EPA 8270	D/SIM (mg/	kg) (continu	ıed)														
Di-N-Octyl Phthalate	ND (0.041)	ND (0.045)	ND (0.044)	ND (0.048)	ND (0.042)	ND (0.041)	ND (0.042)	ND (0.041)	ND (0.044)	ND (0.045)	ND (0.039)	ND (0.050)	ND (0.041)	ND (0.042)	ND (0.038)	ND (0.047)	ND (0.035)	ND (0.041)	800
Fluoranthene	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.010)	ND (0.008)	ND (0.008)	0.042	ND (0.008)	ND (0.009)	ND (0.009)	0.013	ND (0.010)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.010)	ND (0.007)	ND (0.008)	0.296
Fluorene	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.010)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.008)	ND (0.010)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.010)	ND (0.007)	ND (0.008)	0.080
1-Methylnaphthalene	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.010)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.008)	ND (0.010)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.010)	ND (0.007)	ND (0.008)	0.236
2-Methylnaphthalene	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.010)	ND (0.008)	ND (0.008)	0.017	0.010	ND (0.009)	ND (0.009)	ND (0.010)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.010)	ND (0.008)	0.236
Naphthalene	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.010)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.008)	ND (0.010)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.010)	ND (0.007)	ND (0.008)	0.236
4-Nitrophenol	ND (0.041)	ND (0.045)	ND (0.044)	ND (0.048)	ND (0.042)	ND (0.041)	ND (0.042)	ND (0.041)	ND (0.044)	ND (0.045)	ND (0.039)	ND (0.050)	ND (0.041)	ND (0.042)	ND (0.038)	ND (0.047)	ND (0.035)	ND (0.041)	NA
p-Cresol	ND (0.041)	ND (0.045)	ND (0.044)	ND (0.048)	ND (0.042)	ND (0.041)	ND (0.042)	ND (0.041)	ND (0.044)	ND (0.045)	ND (0.039)	ND (0.050)	ND (0.041)	ND (0.042)	ND (0.038)	ND (0.047)	ND (0.035)	ND (0.041)	8,000
Pentachlorophenol	ND (0.21)	ND (0.22)	ND (0.22)	ND (0.24)	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.22)	ND (0.22)	ND (0.19)	ND (0.25)	ND (0.20)	ND (0.21)	ND (0.19)	ND (0.24)	ND (0.17)	ND (0.21)	0.17
Phenanthrene	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.010)	ND (0.008)	ND (0.008)	0.028	0.011	ND (0.009)	ND (0.009)	ND (0.008)	ND (0.010)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.010)	ND (0.007)	ND (0.008)	0.0067
Phenol	ND (0.041)	ND (0.045)	ND (0.044)	ND (0.048)	ND (0.042)	ND (0.041)	ND (0.042)	ND (0.041)	ND (0.044)	ND (0.045)	ND (0.039)	ND (0.050)	ND (0.041)	ND (0.042)	ND (0.038)	ND (0.047)	ND (0.035)	ND (0.041)	0.757
Pyrene	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.010)	ND (0.008)	ND (0.008)	0.046	ND (0.008)	ND (0.009)	ND (0.009)	0.015	ND (0.010)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.010)	ND (0.007)	ND (0.008)	0.546
Carcinogenic Polycyclic	Aromatic H	ydrocarbons	(cPAHs) by	EPA 8270D	/SIM (mg/kg)													
Benzo(a)anthracene	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.010)	ND (0.008)	ND (0.008)	0.029	ND (0.008)	ND (0.009)	ND (0.009)	0.008	ND (0.010)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.010)	ND (0.007)	ND (0.008)	0.0067
Benzo(a)pyrene	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.010)	ND (0.008)	ND (0.008)	0.034	ND (0.008)	ND (0.009)	ND (0.009)	0.012	ND (0.010)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.010)	ND (0.007)	ND (0.008)	0.01
Benzo(b)fluoranthene	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.010)	ND (0.008)	ND (0.008)	0.042	ND (0.008)	ND (0.009)	ND (0.009)	0.016	ND (0.010)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.010)	ND (0.007)	ND (0.008)	0.012
Benzo(j,k)fluoranthene	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.010)	ND (0.008)	ND (0.008)	0.014	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.008)	ND (0.010)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.010)	ND (0.007)	ND (0.008)	0.012
Chrysene	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.010)	ND (0.008)	ND (0.008)	0.037	0.014	ND (0.009)	ND (0.009)	0.012	ND (0.010)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.010)	ND (0.007)	ND (0.008)	0.0067
Dibenz(a,h)anthracene	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.010)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.008)	ND (0.010)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.010)	ND (0.007)	ND (0.008)	0.018
Indeno(1,2,3-cd)pyrene	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.010)	ND (0.008)	ND (0.008)	0.027	ND (0.008)	ND (0.009)	ND (0.009)	0.009	ND (0.010)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.010)	ND (0.007)	ND (0.008)	0.035
Total cPAHs (TEQ)b	ND (0.006)	ND (0.007)	ND (0.007)	ND (0.007)	ND (0.006)	ND (0.006)	0.046	0.006	ND (0.007)	ND (0.007)	0.016	ND (0.008)	ND (0.006)	ND (0.006)	ND (0.006)	ND (0.007)	ND (0.005)	ND (0.006)	0.020

BOLD values detected above the reporting limit.

Shaded values exceed the site screening level





			Table 7 (c	ontinued). S	ummary of S	oil Sample R	esults, Pacifi	c City Park Re	emedial Inve	stigation, Pac	ific, Washing	jton.			
Analytical							Sample	Location							
Parameter	PP	10	PP	11		PP12		PP13	PF	P1 4	PP	15	PP	16	Site Screening
Sample Date	2/28	8/18	2/28	8/18		2/21/18		2/21/18	2/28	8/18	2/2:	1/18	2/28	3/18	Level ^a
Depth (feet)	11	17	5	10	2	7	10	10	5	12	6	12	1	11	(mg/kg)
Petroleum Hydrocarb	ons by Method	NWTPH-Dx (mg	g/kg)												
Diesel Range Organics	-	-	-	_	ND (31)	ND (130)	ND (32)	-	-	_	81	-	ND (320)	-	200
Lube Oil Range Organics	-	_	-	_	69	800	82	-	-	_	550	_	3,200	-	2,000
Volatile Petroleum Pr	oducts Includin	g Gasoline, Benz	zene, Toluene, E	thylbenzene, an	nd Total Xylenes	(BTEX) by NW	TPH-GX (mg/kg)							
Gasoline	-	_	-	_	-	_	_	-	-	_	-	_	ND (5.3)	-	100
Benzene	-	_	-	_	-	_	_	-	-	_	-	_	ND (0.020)	-	0.001
Toluene	-	_	-	_	-	_	_	-	-	_	-	_	ND (0.053)	_	0.024
Ethylbenzene	-	-	-	-	-	-	_	-	-	_	-	-	ND (0.053)	_	0.014
Total Xylenes	-	-	-	_	-	_	_	-	-	_	-	_	ND (0.11)	_	0.52
Total Metals by EPA 6	6010D/7471B (m	ng/kg)													
Arsenic	ND (16)	ND (12)	19	ND (13)	ND (12)	16	ND (13)	ND (12)	ND (12)	ND (13)	ND (15)	24	ND (11)	ND (12)	20
Cadmium	ND (0.82)	ND (0.62)	1.7	0.72	ND (0.62)	4.1	ND (0.64)	3.5	ND (0.62)	ND (0.65)	0.79	1.0	ND (0.53)	1.2	1
Chromium	22	15	39	36	15	31	25	29	12	11	22	34	20	24	48
Lead	15	ND (6.2)	840	480	8.8	500	30	330	ND (6.2)	ND (6.5)	75	270	17	200	25
Mercury	ND (0.41)	ND (0.31)	0.41	1.2	ND (0.31)	0.76	ND (0.32)	ND (0.31)	ND (0.31)	ND (0.32)	ND (0.38)	ND (0.30)	ND (0.27)	ND (0.31)	0.07
Polychlorinated Biphe	enyls (PCBs) by	EPA 8082A (mg	/kg)					_							
Total PCBs	-	-	-	_	ND (0.062)	0.29	ND (0.064)	-	-	_	ND (0.076)	_	ND (0.053)	_	0.05
Carcinogenic Polycycl	lic Aromatic Hy	drocarbons (cPA	NHs) by EPA 827	OD/SIM (mg/kg)										
Benzo(a)anthracene	ND (0.011)	ND (0.008)	ND (0.009)	ND (0.008)	0.008	0.010	ND (0.009)	ND (0.008)	ND (0.008)	ND (0.009)	ND (0.010)	ND (0.008)	0.20	0.12	0.0067
Benzo(a)pyrene	ND (0.011)	ND (0.008)	ND (0.009)	ND (0.008)	0.015	0.011	ND (0.009)	ND (0.008)	ND (0.008)	ND (0.009)	ND (0.010)	ND (0.008)	0.058	0.10	0.01
Benzo(b)fluoranthene	ND (0.011)	ND (0.008)	ND (0.009)	ND (0.008)	0.016	0.014	ND (0.009)	ND (0.008)	ND (0.008)	ND (0.009)	ND (0.010)	ND (0.008)	0.086	0.15	0.012
Benzo(j,k)fluoranthene	ND (0.011)	ND (0.008)	ND (0.009)	ND (0.008)	ND (0.008)	ND (0.010)	ND (0.009)	ND (0.008)	ND (0.008)	ND (0.009)	ND (0.010)	ND (0.008)	ND (0.036)	0.049	0.012
Chrysene	ND (0.011)	ND (0.008)	ND (0.009)	ND (0.008)	0.014	0.013	ND (0.009)	ND (0.008)	ND (0.008)	ND (0.009)	ND (0.010)	ND (0.008)	0.060	0.11	0.0067
Dibenz(a,h)anthracene	ND (0.011)	ND (0.008)	ND (0.009)	ND (0.008)	ND (0.008)	ND (0.010)	ND (0.009)	ND (0.008)	ND (0.008)	ND (0.009)	ND (0.010)	ND (0.008)	ND (0.036)	0.015	0.018
Indeno(1,2,3-cd)pyrene	ND (0.011)	ND (0.008)	ND (0.009)	ND (0.008)	0.012	ND (0.010)	ND (0.009)	ND (0.008)	ND (0.008)	ND (0.009)	ND (0.010)	ND (0.008)	0.036	0.069	0.035
Total cPAHs (TEQ) ^b	ND (0.008)	ND (0.006)	ND (0.007)	ND (0.006)	0.020	0.015	ND (0.007)	ND (0.006)	ND (0.006)	ND (0.007)	ND (0.008)	ND (0.060)	0.094	0.14	0.020

BOLD values detected above the reporting limit.

Shaded values exceed the site screening level

Native Soil Fill/Refuse

King County

January 2019 Remedial Investigation Report—Pacific City Park

			Table 7 (con	tinued). Sumi	mary of Soil Sa	ample Results	, Pacific City P	ark Remedial	Investigation	Pacific, Wash	ington.			
Analytical							Sample Location	l						
Parameter		PP17			PP18			PP19			PP	20		Site Screening
Sample Date		2/28/18			2/28/18			2/21/18			2/28	8/18		Level ^a
Depth (feet)	1	5	10	3	5	10	7	10	15	2	5	10	15	(mg/kg)
Petroleum Hydrocarb	ons by Method N	NWTPH-Dx (mg/	kg)											
Diesel Range Organics	-	_	_	_	-	-	400	_	_	ND (29)	ND (200)	ND (32)	ND (58)	200
Lube Oil Range Organics	-	_	_	_	-	_	370	_	_	380	1,300	95	320	2,000
Volatile Petroleum Pr	oducts Including	Gasoline, Benze	ne, Toluene, Ethy	lbenzene, and To	tal Xylenes (BTEX	() by NWTPH-GX	(mg/kg)							
Gasoline	-	_	_	_	-	-	-	_	_	-	_	-	_	100
Benzene	-	_	_	_	_	_	-	_	-	_	_	_	_	0.001
Toluene	_	_	_	-	_	_	-	_	_	_	_	_	_	0.024
Ethylbenzene	-	-	-	-	_	-	-	-	-	-	-	-	-	0.014
Total Xylenes	_	_	_	-	-	-	-	_	_	_	_	_	_	0.52
Total Metals by EPA 6	010D/7471B (mg	g/kg)												
Arsenic	19	ND (12)	ND (12)	ND (12)	ND (20)	ND (13)	ND (13)	ND (12)	ND (12)	ND (12)	ND (15)	ND (13)	ND (14)	20
Cadmium	4.0	4.5	1.6	2.4	3.1	1.2	ND (0.63)	0.95	3.2	ND (0.58)	3.7	ND (0.64)	ND (0.70)	1
Chromium	57	59	24	63	39	40	17	30	17	15	53	22	15	48
Lead	7,300	380	29	130	230	97	82	340	96	28	630	270	330	<i>2</i> 5
Mercury	ND (0.32)	ND (0.30)	ND (0.29)	ND (0.31)	ND (0.49)	ND (0.33)	ND (0.32)	ND (0.29)	ND (0.30)	ND (0.29)	ND (0.37)	ND (0.32)	ND (0.35)	0.07
Polychlorinated Biphe	enyls (PCBs) by E	PA 8082A (mg/k	g)											
Total PCBs	-	_	_	-	_	-	-	_	_	ND (0.058)	0.74	ND (0.064)	ND (0.070)	0.05
Carcinogenic Polycycl	ic Aromatic Hyd	rocarbons (cPAH	s) by EPA 8270D,	/SIM (mg/kg)										
Benzo(a)anthracene	ND (0.009)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.013)	ND (0.009)	ND (0.008)	ND (0.008)	ND (0.008)	0.0080	0.23	ND (0.043)	ND (0.046)	0.0067
Benzo(a)pyrene	ND (0.009)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.013)	ND (0.009)	ND (0.008)	ND (0.008)	ND (0.008)	0.013	0.22	ND (0.043)	ND (0.046)	0.01
Benzo(b)fluoranthene	ND (0.009)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.013)	ND (0.009)	ND (0.008)	ND (0.008)	ND (0.008)	0.023	0.20	ND (0.043)	ND (0.046)	0.012
Benzo(j,k)fluoranthene	ND (0.009)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.013)	ND (0.009)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.0077)	ND (0.098)	ND (0.043)	ND (0.046)	0.012
Chrysene	ND (0.009)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.013)	ND (0.009)	ND (0.008)	ND (0.008)	ND (0.008)	0.019	0.37	ND (0.043)	ND (0.046)	0.0067
Dibenz(a,h)anthracene	ND (0.009)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.013)	ND (0.009)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.0077)	ND (0.098)	ND (0.043)	ND (0.046)	0.018
Indeno(1,2,3-cd)pyrene	ND (0.009)	ND (0.008)	ND (0.008)	ND (0.008)	ND (0.013)	ND (0.009)	ND (0.008)	ND (0.008)	ND (0.008)	0.013	0.12	ND (0.043)	ND (0.046)	0.035
Total cPAHs (TEQ) ^b	ND (0.007)	ND (0.006)	ND (0.006)	ND (0.006)	ND (0.010)	ND (0.007)	ND (0.006)	ND (0.006)	ND (0.006)	0.018	0.29	ND (0.032)	ND (0.035)	0.020

BOLD values detected above the reporting limit.

Shaded values exceed the site screening level

Native Soil
Fill/Refuse



			Table 7 (con	tinued). Sumi	mary of Soil S	ample Results	, Pacific City P	ark Remedial	Investigation	Pacific, Wash	nington.			
Analytical							Sample Location	ı						
Parameter		PI	P21			PP22			PI	23		PP	24	Site Screening
Sample Date						2/21/18			2/2	8/18		2/28	8/18	Level ^a
Depth (feet)	2	6	10	15	4	9	13	2	5	10	15	7	10	(mg/kg)
Petroleum Hydrocarb	ons by Method	NWTPH-Dx (mg/	'kg)											
Diesel Range Organics	57	150	1,800	110	_	_	_	ND (29)	_	_	_	ND (35)	_	200
Lube Oil Range Organics	540	960	10,000	1,200	-	_	_	200	_	_	_	130	-	2,000
Volatile Petroleum Pr	oducts Including	g Gasoline, Benze	ne, Toluene, Ethy	lbenzene, and To	tal Xylenes (BTE	X) by NWTPH-GX	(mg/kg)							
Gasoline	_	_	39	_	_	_	_	ND (13)	_	_	_	ND (7.5)	_	100
Benzene	_	_	ND (0.024)	_	-	_	_	ND (0.025)	_	_	_	ND (0.020)	_	0.001
Toluene	_	_	ND (0.12)	_	-	_	_	ND (0.13)	_	_	_	ND (0.075)	-	0.024
Ethylbenzene	_	_	ND (0.12)	_	-	_	_	ND (0.13)	_	_	_	ND (0.075)	-	0.014
Total Xylenes	-	_	ND (0.24)	_	-	_	_	ND (0.26)	_	_	_	ND (0.15)	_	0.52
Total Metals by EPA 6	6010D/7471B (m	g/kg)												
Arsenic	ND (13)	17	ND (16)	ND (14)	ND (13)	ND (12)	ND (12)	ND (12)	ND (11)	ND (13)	ND (15)	ND (14)	ND (12)	20
Cadmium	0.90	27	6.2	ND (1.4)	ND (0.64)	ND (0.61)	2.4	ND (0.59)	ND (0.55)	ND (0.67)	ND (0.76)	2.3	ND (0.58)	1
Chromium	30	74	23	20	9.9	9.9	36	19	15	13	18	18	10	48
Lead	740	2,800	180	ND (14)	ND (6.4)	ND (6.1)	270	460	97	ND (6.7)	220	480	ND (5.8)	25
Mercury	0.87	ND (0.51)	ND (0.410	ND (0.68)	ND (0.32)	ND (0.31)	ND (0.29)	ND (0.29)	ND (0.27)	ND (0.34)	ND (0.38)	ND (0.35)	ND (0.29)	0.07
Polychlorinated Biphe	enyls (PCBs) by E	PA 8082A (mg/k	g)					_			_			
Total PCBs	0.52	0.27	1.33	ND (0.14)	-	_	_	ND (0.059)	_	_	_	ND (0.069)	_	0.05
Carcinogenic Polycycl	lic Aromatic Hyd	rocarbons (cPAH	ls) by EPA 8270D	/SIM (mg/kg)										
Benzo(a)anthracene	0.076	0.14	0.90	ND (0.018)	0.056	ND (0.004)	ND (0.008)	0.012	0.18	ND (0.009)	ND (0.010)	ND (0.018)	ND (0.008)	0.0067
Benzo(a)pyrene	0.062	0.13	0.85	ND (0.018)	0.064	ND (0.004)	ND (0.008)	0.011	0.044	ND (0.009)	ND (0.010)	ND (0.018)	ND (0.008)	0.01
Benzo(b)fluoranthene	0.11	0.22	1.3	ND (0.018)	0.088	ND (0.004)	ND (0.008)	0.017	0.10	ND (0.009)	ND (0.010)	0.031	ND (0.008)	0.012
Benzo(j,k)fluoranthene	0.033	0.077	0.30	ND (0.018)	0.029	ND (0.004)	ND (0.008)	ND (0.0078)	ND (0.015)	ND (0.009)	ND (0.010)	ND (0.018)	ND (0.008)	0.012
Chrysene	0.097	0.18	1.0	ND (0.018)	0.086	ND (0.004)	ND (0.008)	0.017	0.044	ND (0.009)	ND (0.010)	0.035	ND (0.008)	0.0067
Dibenz(a,h)anthracene	0.010	0.024	0.19	ND (0.018)	0.017	ND (0.004)	ND (0.008)	ND (0.0078)	0.019	ND (0.009)	ND (0.010)	ND (0.018)	ND (0.008)	0.018
Indeno(1,2,3-cd)pyrene	0.040	0.11	0.71	ND (0.018)	0.051	ND (0.004)	ND (0.008)	0.0082	0.026	ND (0.009)	ND (0.010)	ND (0.018)	ND (0.008)	0.035
Total cPAHs (TEQ) ^b	0.090	0.19	1.2	ND (0.014)	0.089	ND (0.006)	ND (0.006)	0.0059	0.078	ND (0.007)	ND (0.008)	0.016	ND (0.006)	0.020

BOLD values detected above the reporting limit.

Shaded values exceed the site screening level

Native Soil Fill/Refuse

January 2019

King County

Remedial Investigation Report—Pacific City Park

			Table 7 (c	ontinued). S	Summary of S	Soil Sample R	esults, Pacifi	c City Park R	emedial Inve	stigation, Pac	ific, Washing	gton.			
							Sample	Location							
Analytical Parameter		PP25		PP	26	PP	27	PF	28		PP29		PP	30	Site Screening
Sample Date		2/21/18		2/2:	1/18	2/28	3/18	3/1	/18		3/1/18		2/28	8/18	Level ^a
Depth (feet)	7	13	17	11	17	7	10	8	10	3	5	10	5	10	(mg/kg)
Petroleum Hydrocarbo	ons by Method	NWTPH-Dx (m	g/kg)												
Diesel Range Organics	_	1,400	49	-	_	-	_	_	_	ND (280)	_	_	71	_	200
Lube Oil Range Organics	-	1,200	130	-	-	-	-	_	-	5,900	-	_	630	-	2,000
Volatile Petroleum Pro	ducts Includin	g Gasoline, Ben	zene, Toluene, I	Ethylbenzene, aı	nd Total Xylenes	s (BTEX) by NW	TPH-GX (mg/kg)							
Gasoline	-	400	ND (15)	-	_	-	_	_	_	-	_	_		_	100
Benzene	-	ND (0.026)	ND (0.15)	-	-	-	-	_	-	-	-	_		-	0.001
Toluene	_	ND (0.13)	ND (0.15)	-	_	-	_	-	_	-	_	_		_	0.024
Ethylbenzene	_	ND (0.13)	ND (0.15)	_	_	-	_	_	_	_	_	_		_	0.014
Total Xylenes	_	ND (0.13)	ND (0.15)	_	_	_	_	_	_	-	_	_		_	0.52
Total Metals by EPA 6	010D/7471B (n	ng/kg)													
Arsenic	ND (14)	ND (17)	ND (19)	ND (12)	ND (15)	ND (15)	ND (12)	ND (13)	ND (12)	ND (11)	ND (13)	ND (13)	ND (13)	ND (11)	20
Cadmium	ND (0.68)	1.9	ND (0.93)	ND (0.58)	ND (0.77)	0.88	ND (0.61)	ND (0.63)	ND (0.62)	ND (0.56)	ND (0.66)	ND (0.66)	0.83	ND (0.57)	1
Chromium	13	38	28	21	15	28	12	25	23	25	15	16	16	14	48
Lead	8.8	140	ND (9.3)	ND (5.8)	ND (7.7)	180	ND (6.1)	49	6.8	8.8	ND (6.5)	ND (6.6)	31	ND (5.7)	25
Mercury	ND (0.34)	ND (0.42)	ND (0.47)	ND (0.29)	ND (0.39)	ND (0.36)	ND (0.30)	ND (0.31)	ND (0.31)	ND (0.28)	ND (0.32)	ND (0.33)	ND (0.32)	ND (0.28)	0.07
Polychlorinated Biphe	nyls (PCBs) by	EPA 8082A (mg	/kg)												
Total PCBs	-	1.36	ND (0.093)	-	_	-	_	_	_	ND (0.056)	_	_	ND (0.064)	_	0.05
Carcinogenic Polycycli	c Aromatic Hy	drocarbons (cPA	AHs) by EPA 827	70D/SIM (mg/kg	J)										
Benzo(a)anthracene	ND (0.009)	ND (0.011)	ND (0.012)	ND (0.008)	ND (0.010)	0.036	ND (0.008)	ND (0.008)	0.011	ND (0.038)	ND (0.009)	ND (0.009)	ND (0.085)	ND (0.008)	0.0067
Benzo(a)pyrene	ND (0.009)	ND (0.011)	ND (0.012)	ND (0.008)	ND (0.010)	0.038	ND (0.008)	ND (0.008)	0.010	ND (0.038)	ND (0.009)	ND (0.009)	ND (0.085)	ND (0.008)	0.01
Benzo(b)fluoranthene	ND (0.009)	ND (0.011)	ND (0.012)	ND (0.008)	ND (0.010)	0.049	ND (0.008)	ND (0.008)	0.0084	ND (0.038)	ND (0.009)	ND (0.009)	ND (0.085)	ND (0.008)	0.012
Benzo(j,k)fluoranthene	ND (0.009)	ND (0.011)	ND (0.012)	ND (0.008)	ND (0.010)	0.014	ND (0.008)	ND (0.008)	ND (0.0083)	ND (0.038)	ND (0.009)	ND (0.009)	ND (0.085)	ND (0.008)	0.012
Chrysene	ND (0.009)	0.012	ND (0.012)	ND (0.008)	ND (0.010)	0.051	ND (0.008)	ND (0.008)	0.011	ND (0.038)	ND (0.009)	ND (0.009)	ND (0.085)	ND (0.008)	0.0067
Dibenz(a,h)anthracene	ND (0.009)	ND (0.011)	ND (0.012)	ND (0.008)	ND (0.010)	ND (0.0097)	ND (0.008)	ND (0.008)	ND (0.0083)	ND (0.038)	ND (0.009)	ND (0.009)	ND (0.085)	ND (0.008)	0.018
Indeno(1,2,3-cd)pyrene	ND (0.009)	ND (0.011)	ND (0.012)	ND (0.008)	ND (0.010)	0.025	ND (0.008)	ND (0.008)	ND (0.0083)	ND (0.038)	ND (0.009)	ND (0.009)	ND (0.085)	ND (0.008)	0.035
Total cPAHs (TEQ) ^b	ND (0.007)	0.008	ND (0.009)	ND (0.006)	ND (0.008)	0.051	ND (0.006)	ND (0.006)	0.013	ND (0.029)	ND (0.007)	ND (0.007)	ND (0.064)	ND (0.006)	0.020

BOLD values detected above the reporting limit.

Shaded values exceed the site screening level

Native Soil
Fill/Refuse



			Table 7 (cont	inued). Sum	mary of Soil S	ample Results	, Pacific City P	ark Remedial	Investigation	, Pacific, Wash	ington.			
Analytical							Sample Location							
Parameter	PF	931		PP32			PP33		PP34		B-04	B-05		Site Screening
Sample Date	3/1	_/18		3/1/18			3/1/18		2/2	1/18	2/21/18	2/20	0/18	Level ^a
Depth (feet)	3	11	4	7	10	3	5	10	8	15	12.5	7.5	12.5	(mg/kg)
Petroleum Hydrocarbons by Method NWTPH-Dx (mg/kg)														
Diesel Range Organics	-	_	-	ND (60)	_	ND (750)	_	-	-	_	-	440	-	200
Lube Oil Range Organics	-	_	_	650	_	12,000	_	_	-	-	-	ND (220)	_	2,000
Volatile Petroleum Products Including Gasoline, Benzene, Toluene, Ethylbenzene, and Total Xylenes (BTEX) by NWTPH-GX (mg/kg)														
Gasoline	-	_	-	_	ND (13)	ND (12)	_	-	-	_	-	-	_	100
Benzene	-	_	_	_	ND (0.026)	ND (0.024)	_	_	_	_	_	-	_	0.001
Toluene	-	_	_	_	ND (0.13)	ND (0.12)	_	_	-	-	_	-	_	0.024
Ethylbenzene	-	_	-	_	ND (0.13)	ND (0.12)	_	_	-		_	-	_	0.014
Total Xylenes	-	_	-	_	ND (0.26)	ND (0.24)	_	_	-		_	-	_	0.52
Total Metals by EPA 6010D/7471B (mg/kg)														
Arsenic	ND (11)	ND (12)	ND (12)	ND (12)	ND (12)	ND (11)	ND (12)	ND (13)	ND (12)	_	ND (13)	ND (12)	ND (13)	20
Cadmium	ND (0.56)	ND (0.62)	ND (0.62)	ND (0.61)	ND (0.61)	ND (0.56)	ND (0.59)	ND (0.67)	ND (0.61)	-	ND (0.64)	ND (0.60)	ND (0.63)	1
Chromium	21	12	16	15	20	17	19	13	12		16	430	11	48
Lead	12	ND (6.2)	12	15	ND (6.1)	8.1	7.3	ND (6.7)	6.1	-	ND (6.4)	ND (6.0)	ND (6.3)	25
Mercury	ND (0.28)	ND (0.31)	ND (0.31)	ND (0.31)	ND (0.31)	ND (0.28)	ND (0.30)	ND (0.34)	ND (0.30)	_	ND (0.32)	ND (0.30)	ND (0.32)	0.07
Polychlorinated Biphe	enyls (PCBs) by E	PA 8082A (mg/k	g)											
Total PCBs	-	_	_	ND (0.061)	ND (0.061)	ND (0.056)	_	_	-	-	_	ND (0.060)	_	0.05
Carcinogenic Polycyc	lic Aromatic Hyd	rocarbons (cPAH	s) by EPA 8270D/	SIM (mg/kg)										
Benzo(a)anthracene	ND (0.008)	ND (0.008)	0.010	0.038	ND (0.008)	ND (0.075)	ND (0.008)	ND (0.009)	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.008)	ND (0.008)	0.0067
Benzo(a)pyrene	ND (0.008)	ND (0.008)	0.011	0.036	ND (0.008)	ND (0.075)	ND (0.008)	ND (0.009)	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.008)	ND (0.008)	0.01
Benzo(b)fluoranthene	0.0093	ND (0.008)	0.024	0.051	ND (0.008)	0.14	ND (0.008)	ND (0.009)	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.008)	ND (0.008)	0.012
Benzo(j,k)fluoranthene	ND (0.008)	ND (0.008)	ND (0.008)	0.013	ND (0.008)	ND (0.075)	ND (0.008)	ND (0.009)	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.008)	ND (0.008)	0.012
Chrysene	ND (0.008)	ND (0.008)	0.015	0.055	ND (0.008)	0.19	ND (0.008)	ND (0.009)	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.008)	ND (0.008)	0.0067
Dibenz(a,h)anthracene	ND (0.008)	ND (0.008)	ND (0.008)	0.009	ND (0.008)	ND (0.075)	ND (0.008)	ND (0.009)	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.008)	ND (0.008)	0.018
Indeno(1,2,3-cd)pyrene	ND (0.008)	ND (0.008)	0.0095	0.025	ND (0.008)	ND (0.075)	ND (0.008)	ND (0.009)	ND (0.008)	ND (0.009)	ND (0.009)	ND (0.008)	ND (0.008)	0.035
Total cPAHs (TEQ) ^b	0.006	ND (0.006)	0.016	0.050	ND (0.006)	0.068	ND (0.006)	ND (0.007)	ND (0.006)	ND (0.007)	ND (0.006)	ND (0.006)	ND (0.006)	0.020

BOLD values detected above the reporting limit.

Shaded values exceed the site screening level

Native Soil Fill/Refuse

January 2019

King County

			Table 7 (cont	inued). Sum	mary of Soil	Sample Resul	ts, Pacific Pa	rk City Park I	Remedial I	nvestigation,	Pacific, Wa	shington.			
							Sample L	ocation							
Analytical Parameter		B-06			B-07			B-08 B-		В-(B-09 B-10		В	-11	Site Screening
Sample Date		2/22/18			2/27/18			2/26/18		2/22	/18	2/26/18	2/26/18		Level ^a
Depth (feet)	2.5	10	12.5	2.5	7.5	12.5	5	7.5	12.5	5	15	7.5	2.5	12.5	(mg/kg)
Petroleum Hydrocarb	ons by Method	NWTPH-Dx (mg	ı/kg)												
Diesel Range Organics	ND (31)	ND (31)	ND (32)	ND (150)	ND (310)	ND (33)	_	_	-	-	_	ND (30)	ND (33)	_	200
Lube Oil Range Organics	150	ND (61)	130	4,400	1,800	180	_	_	-	-	-	88	380	-	2,000
Volatile Petroleum Pro	um Products Including Gasoline, Benzene, Toluene, Ethylbenzene, and Total Xylenes (BTEX) by NWTPH-GX (mg/kg)														
Gasoline	-	_	_	-	-	_	_	_	-	_	-	-	_	_	100
Benzene	_	_	_	_	_	_	_	_	-	_	_	_	-	_	0.001
Toluene	_	_	_	-	_	_	_	_	_	_	_	_	_	_	0.024
Ethylbenzene	-	_	_	-	-	_	-	_	-	_	_	_	-	_	0.014
Total Xylenes	-	_	_	-	-	_	-	_	-	_	_	_	-	_	0.52
Total Metals by EPA 6	010D/7471B (m	ıg/kg)													
Arsenic	ND (12)	ND (13)	ND (13)	ND (12)	ND (12)	ND (13)	ND (14)	_	ND (13)	ND (15)	ND (16)	ND (12)	ND (13)	ND (12)	20
Cadmium	0.75	ND (0.61)	ND (0.64)	ND (0.58)	ND (0.62)	ND (0.66)	ND (0.69)	_	ND (0.67)	ND (0.77)	ND (0.78)	ND (0.59)	ND (0.65)	ND (0.60)	1
Chromium	13	13	11	15	14	10	16	_	18	16	24	8.0	16	13	48
Lead	20	ND (6.1)	ND (6.4)	6.0	ND (6.2)	ND (6.6)	6.9	-	ND (6.7)	20	ND (7.8)	ND (5.9)	15	ND (6.0)	25
Mercury	ND (0.31)	ND (0.31)	ND (0.32)	ND (0.29)	ND (0.31)	ND (0.33)	ND (0.35)	_	ND (0.34)	ND (0.39)	ND (0.39)	ND (0.29)	ND (0.32)	ND (0.30)	0.07
Polychlorinated Biphe	nyls (PCBs) by l	EPA 8082A (mg/	kg)												
Total PCBs	ND (0.062)	_	ND (0.064)	ND (0.058)	ND (0.062)	ND (0.066)	_	_	_	_	_	ND (0.059)	ND (0.26)	_	0.05
Carcinogenic Polycycl	ic Aromatic Hyd	drocarbons (cPA	Hs) by EPA 8270	D/SIM (mg/kg)											
Benzo(a)anthracene	ND (0.008)	ND (0.008)	ND (0.009)	ND (0.039)	ND (0.041)	ND (0.009)	ND (0.009)	_	ND (0.009)	ND (0.010)	ND (0.010)	ND (0.008)	0.016	ND (0.008)	0.0067
Benzo(a)pyrene	ND (0.008)	ND (0.008)	ND (0.009)	ND (0.039)	ND (0.041)	ND (0.009)	ND (0.009)	_	ND (0.009)	ND (0.010)	ND (0.010)	ND (0.008)	0.019	ND (0.008)	0.01
Benzo(b)fluoranthene	ND (0.008)	ND (0.008)	ND (0.009)	ND (0.039)	ND (0.041)	ND (0.009)	ND (0.009)	-	ND (0.009)	0.011	ND (0.010)	ND (0.008)	0.053	ND (0.008)	0.012
Benzo(j,k)fluoranthene	ND (0.008)	ND (0.008)	ND (0.009)	ND (0.039)	ND (0.041)	ND (0.009)	ND (0.009)	-	ND (0.009)	ND (0.010)	ND (0.010)	ND (0.008)	0.014	ND (0.008)	0.012
Chrysene	ND (0.008)	ND (0.008)	ND (0.009)	ND (0.039)	ND (0.041)	ND (0.009)	ND (0.009)	-	ND (0.009)	0.012	ND (0.010)	ND (0.008)	0.035	ND (0.008)	0.0067
Dibenz(a,h)anthracene	ND (0.008)	ND (0.008)	ND (0.009)	ND (0.039)	ND (0.041)	ND (0.009)	ND (0.009)	-	ND (0.009)	ND (0.010)	ND (0.010)	ND (0.008)	ND (0.009)	ND (0.008)	0.018
Indeno(1,2,3-cd)pyrene	ND (0.008)	ND (0.008)	ND (0.009)	ND (0.039)	ND (0.041)	ND (0.009)	ND (0.009)	-	ND (0.009)	ND (0.010)	ND (0.010)	ND (0.008)	0.017	ND (0.008)	0.035
Total cPAHs (TEQ) ^b	ND (0.006)	ND (0.006)	ND (0.007)	ND (0.029)	ND (0.031)	ND (0.007)	ND (0.007)	_	ND (0.007)	0.008	ND (0.008)	ND (0.006)	0.030	ND (0.006)	0.020

BOLD values detected above the reporting limit.

Shaded values exceed the site screening level

Native Soil
Fill/Refuse



	Table 7 (continued). Summary of Soil Sample Results, Pacific Park City Park Remedial Investigation, Pacific, Washington.												
Analytical						Sample	Location						
Parameter	B-	13	B-	14		B-15			B-16		B-17		Site Screening
Sample Date	2/2	3/18	2/21	L/18	2/21/18				2/23/18		2/20/18		Level ^a
Depth (feet)	7.5	15	5	10	5	7.5	15	2.5	7.5	17.5	10	15	(mg/kg)
Petroleum Hydrocarb	ons by Method N	WTPH-Dx (mg/kg))										
Diesel Range Organics	-	_	_	-	_	_	_	-	_	_	-	_	200
Lube Oil Range Organics	-	_	-	_	-	-	-	_	_	_	-	_	2,000
Volatile Petroleum Pr	oleum Products Including Gasoline, Benzene, Toluene, Ethylbenzene, and Total Xylenes (BTEX) by NWTPH-GX (mg/kg)												
Gasoline	-	_	-	_	_	_	_	-	_	_	-	_	100
Benzene	-	_	-	_	-	-	-	-	_	_	-	_	0.001
Toluene	_	_	-	_	-	-	_	_	_	_	-	_	0.024
Ethylbenzene	-	-	-	_	-	-	-	-	-	-	-	-	0.014
Total Xylenes	-	_	-	-	-	_	_	-	-	_	-	_	0.52
Total Metals by EPA 6010D/7471B (mg/kg)													
Arsenic	ND (17)	ND (13)	ND (13)	ND (12)	ND (17)	ND (15)	ND (13)	ND (14)	ND (12)	ND (13)	ND (12)	ND (11)	20
Cadmium	ND (0.84)	ND (0.66)	ND (0.67)	ND (0.62)	1.3	ND (0.76)	ND (0.67)	ND (0.71)	1.6	ND (0.66)	ND (0.61)	ND (0.54)	1
Chromium	24	8.9	23	17	35	31	17	18	28	16	23	9.6	48
Lead	12	ND (6.6)	31	33	75	49	ND (6.7)	9.2	68	ND (6.6)	ND (6.1)	ND (5.4)	25
Mercury	ND (0.42)	ND (0.33)	ND (0.34)	ND (0.31)	ND (0.43)	ND (0.38)	ND (0.34)	ND (0.36)	ND (0.29)	ND (0.33)	ND (0.31)	ND (0.27)	0.07
Polychlorinated Bipho	enyls (PCBs) by EP	A 8082A (mg/kg)					_						_
Total PCBs	-	_	-	-	-	-	_	-	_	_	-	_	0.05
Carcinogenic Polycyc	lic Aromatic Hydro	ocarbons (cPAHs) l	by EPA 8270D/SIM	(mg/kg)									
Benzo(a)anthracene	ND (0.011)	ND (0.009)	ND (0.009)	ND (0.008)	ND (0.012)	ND (0.010)	ND (0.009)	ND (0.010)	ND (0.008)	ND (0.009)	ND (0.008)	ND (0.007)	0.0067
Benzo(a)pyrene	ND (0.011)	ND (0.009)	ND (0.009)	ND (0.008)	ND (0.012)	ND (0.010)	ND (0.009)	ND (0.010)	ND (0.008)	ND (0.009)	ND (0.008)	ND (0.007)	0.01
Benzo(b)fluoranthene	0.014	ND (0.009)	ND (0.009)	ND (0.008)	0.014	ND (0.010)	ND (0.009)	ND (0.010)	ND (0.008)	ND (0.009)	ND (0.008)	ND (0.007)	0.012
Benzo(j,k)fluoranthene	ND (0.011)	ND (0.009)	ND (0.009)	ND (0.008)	ND (0.012)	ND (0.010)	ND (0.009)	ND (0.010)	ND (0.008)	ND (0.009)	ND (0.008)	ND (0.007)	0.012
Chrysene	0.016	ND (0.009)	ND (0.009)	ND (0.008)	ND (0.012)	ND (0.010)	ND (0.009)	ND (0.010)	ND (0.008)	ND (0.009)	ND (0.008)	ND (0.007)	0.0067
Dibenz(a,h)anthracene	ND (0.011)	ND (0.009)	ND (0.009)	ND (0.008)	ND (0.012)	ND (0.010)	ND (0.009)	ND (0.010)	ND (0.008)	ND (0.009)	ND (0.008)	ND (0.007)	0.018
Indeno(1,2,3-cd)pyrene	ND (0.011)	ND (0.009)	ND (0.009)	ND (0.008)	ND (0.012)	ND (0.010)	ND (0.009)	ND (0.010)	ND (0.008)	ND (0.009)	ND (0.008)	ND (0.007)	0.035
Total cPAHs (TEQ) ^b	0.009	ND (0.007)	ND (0.007)	ND (0.006)	0.010	ND (0.008)	ND (0.007)	ND (0.007)	ND (0.006)	ND (0.067)	ND (0.006)	ND (0.006)	0.020

BOLD values detected above the reporting limit.

Shaded values exceed the site screening level



January 2019

ND = not detected above laboratory reporting limits (shown in parentheses)

– = not analyzed or not applicable

^a Refer to Table 6 "Proposed Site Screening Levels for Soil" for notes on how each screening level was selected.

b Total carcinogenic polycyclic aromatic hydrocarbons (cPAHs) toxicity equivalency (TEQ) concentration was calculated using one-half the reporting limit for compounds that were not detected above the reporting limit. mg/kg = milligrams per kilogram

	Table 8. Summary of Soil Vapor Monitoring Data, Pacific City Park Remedial Investigation, Pacific, Washington.																					
		Sample Identification ^a																				
	1	2	3	4	5	6	7	8	9	10	а	b	c	d	е	MW6	MW9	MW6	MW9	MW1	MW6	MW9
Sample Date								10/23/	1984							3/2	3/18	6/21	/18		9/26/18	
Parameter	Parameter																					
Methane (% Vol	Trace	0	0	0	0.3	0	0	0	NA	0	0.4	Trace	0.2	Trace	Trace	0.0	0.0	0.0	0.0	0.0	0.0	0.0
Trace Gas ^b (ppm)	0.1	0.3	-0.1	0	6.2	0	0	NA	0	-0.1	0	0	0	0.1	0	NA	NA	NA	NA	NA	NA	NA
H2S (ppm)	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	0.0	0.0	0.0	0.0	0.0	0.0	0.0

^a Samples 1 through 10, and a through e by King County 1984. Samples MW6 and MW9 by Herrera 2018.

NA = not analyzed

ppm = parts per million

H₂S = hydrogen sulfide

b Trace gases include any organic or inorganic gases with an ionization potential <10.2 electron volts (eV) detected by photo-ionization detector (PID).

APPENDIX A

Soil Boring Logs





 Boring ID
 PP1

 Total depth
 15 feet

 Sheet
 1
 of
 1

Project name	Pacific Park	Drilling Cont	ractor ESN	Drilling method P	ush-probe rig
Project number	15-05986-040	Location	NW corner of park	Sampling method	5 ft core with plastic liner
Client King	County	~ 50 feet w	vest of MW-2	Air monitoring (Y/N)	Yes
HEC rep.	Bruce Carpenter	Date M	ay 23, 2017	Instrument(s)	Photoionization detector

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group	Soil description
(11)		,		,	ML	Grass/Brown sandy SILT, trace of gravel, FILL, damp
				1		D. OHT.
	5 C .		_	2		Brown SILT, trace of sand, FILL, damp
0	5-foot	20/50	<u>▼</u> 1.8	2		Static water level measured at 1.8 feet.
0	core with	20/50	1.8	3		Soil sample PP1-2 collected at 9:20
	liner			3	SM	Brown silty SAND, trace of gravel, FILL,wet
	inici			4	5111	Biown sinty states, trace of graver, files, wer
				5		
					ML	Dark Brown-gray sandy SILT, wet
				6		
	5-foot			7		
0	core	100		- 0		
	with liner			8	SM	D. d. D 11-1-14-14-14-14-14-14-14-14-14-14-14-14
	imer			9	SIVI	Dark Brown-black silty SAND, wet
				9		
				10	PT	Dark Brown-gray PEAT, damp
					SM	Soil sample PP1-10 collected at 9:45
				11		Dark Brown-black silty SAND, wet
						·
	5-foot			12		
0	core	100			ML	8" Dark Brown SILT, trace of sand, wet
	with			13	SM	Dark Brown-black, silty SAND, wet
	liner			1.4) (T	D OHT 1
				14	ML MI /DT	Brown-gray SILT, damp
				15	ML/PT ML	Brown SILT, and PEAT, damp Brown-gray SILT, damp
				13	1711	Set temporary screen from 5 feet to 15 feet.
						Purged approximately 1-gallon prior to sample collection.
						Collected water sample PP1-W at 10:00.
						Backfilled borehole with bentonite chips.



 Boring ID
 PP2

 Total depth
 15 feet

 Sheet
 1
 of
 1

Project name Pacific Park	Drilling Contractor ESN	Drilling method _ Push-probe rig
Project number <u>15-05986-040</u>	Location SE corner of Megan's Court	Sampling method 5 ft core with plastic liner
Client King County		Air monitoring (Y/N) Yes
HEC rep. Bruce Carpenter	Date May 23, 2017	Instrument(s) Photoionization detector

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group	Soil description
	5-foot			1 2	ML	Grass/Brown gravelly sandy SILT, FILL, damp Brown sandy SILT, FILL, damp
0.1	core with liner	60	<u>▽</u> 2.5	3		Soil sample PP2-2.5 collected at 10:35 Groundwater encountered during drilling Dark Brown sandy SILT, wet
	5-foot			5 6 7		
0	core with liner	50		8	SM	Dark Brown silty SAND, wet
				10	ML	Dark Brown sandy SILT, trace of gravel, wet Soil sample PP2-10 collected at 10:50
	5-foot			11	SM	Dark Brown-black silty SAND, wet
0	core with liner	100		13	MH MH/PT	Greenish-gray silty CLAY, with red-brown mottling, damp Occasional 1-inch peat lenses
				14		
						Initially set temporary screen from 5 feet to 15 feet, raised screen to 2 feet to 12 feet. Very slow yielding, took 2.5 hours to collect water Sample. Purged approximately 1-quart prior to sample collection. Began water sample PP2-W collection at 11:00. Backfilled borehole with bentonite chips.



 Boring ID
 PP3

 Total depth
 15 feet

 Sheet
 1
 of
 1

Project name Pacific Park	Drilling Contractor ESN	Drilling method Push-probe rig
Project number <u>15-05986-040</u>	Location East side of ditch	Sampling method 5 ft core with plastic liner
Client King County	~ 190 feet north of PP4	Air monitoring (Y/N) Yes
HEC rep. Bruce Carpenter	Date May 24, 2017	Instrument(s) Photoionization detector

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group	Soil description
\(\frac{1}{1}\)		j			ML	Grass/Brown gravelly sandy SILT, FILL, damp
0	5-foot core with	40/35	<u>▼</u>	2 3	ML	Brown gravelly SILT, FILL, damp Brown gravelly SILT, FILL, wet. Soil sample PP3-2.5 collected at
	liner		3.1		1412	10:05 Static water level measured at 3.1 feet.
				5		
				6	ML	Brown sandy SILT, FILL, wet
	5-foot			7		
0	core with liner	50/55		8	SP	Dark Brown-black fine to medium SAND, trace of gravel, wet
	inici			9		
				10		Soil sample PP3-10 collected at 10:15
	5-foot	0/100		12		
0	core 0/100 with liner		100			
				14		
				15		Greenish-gray sandy SILT, wet
						Set temporary screen from 5 feet to 15 feet. Purged approximately 1-gallon prior to sample collection. Collected water sample PP3-W at 10:00 Backfilled borehole with bentonite chips.



 Boring ID
 PP4

 Total depth
 15 feet

 Sheet
 1
 of
 1

Project name Pacific Park	Drilling C	Contractor ESN	Drilling method P	ush-probe rig
Project number <u>15-05986-040</u>	Location	East end of 4 th Avenue SE in	Sampling method	5 ft core with plastic liner
Client King County	grass		Air monitoring (Y/N)	Yes
HEC rep. Bruce Carpenter	Date	May 24, 2017	Instrument(s)	Photoionization detector

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group	Soil description
M I Z	5-foot	,		1 2	SM SW	Grass/Dark Brown gravelly silty SAND, some concrete, FILL, damp 4" zone of medium silty SAND, FILL, damp
0	core with liner	50	<u>▼</u> 3.2	3	ML	Brown gravelly sandy SILT, red-brown mottling, FILL, damp. Soil sample PP4-3 collected at 8:30
				5	SW	Static water level measured at 3.2 feet. Dark Brown-black gravelly SAND, FILL, wet
				6	SM	Gray-black gravelly silty SAND, piece of asphalt, FILL, wet
0	5-foot core with liner	55		8	ML	Dark Brown-black sandy SILT, wet
	iner			9	SP	Dark Brown-black fine to medium SAND, wet
				11		Soil sample PP4-10 collected at 8:40
0	5-foot core with	100		12	ML	Dark Brown SILT, wet
	liner			1.4	SM	Dark Brown silty SAND, wet
				15	ML	Dark Brown SILT, organic material, damp
						Set temporary screen from 5 feet to 15 feet. Purged approximately 1-gallon prior to sample collection. Collected water sample PP4-W at 9:00 Backfilled borehole with bentonite chips.



Boring ID _	PP5
Total depth	15 feet
Sheet 1	of 1

Project name	Pacific Park	Drilling Cont	ractor ESN	Drilling method P	ush-probe rig
Project number	15-05986-040	Location	~ 215 feet north of PP6,	Sampling method	5 ft core with plastic liner
Client King	County	east side o	f HESCOs	Air monitoring (Y/N)	Yes
HEC rep.	Bruce Carpenter	Date Ma	ay 23, 2017	Instrument(s)	Photoionization detector

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group	Soil description
0	5-foot	50		1 2	SM	Grass/Topsoil, Dark Brown gravelly silty SAND, piece of wire, FILL, damp
0	core with liner	30	<u>▼</u> 3.0	3	SW	Soil sample PP5-2.5 collected at 15:40 Dark Brown-gray gravelly SAND, FILL, damp Static water level measured at 3.0 feet.
				5	PT	Brown PEAT, wet
				6	ML	Dark Brown-black sandy SILT, wet
0.3	5-foot core with liner	60		7 8 9	SW	Dark Brown-black fine to medium SAND, trace of silt, organic material, wet
				10		Dark Brown-black gravelly fine to coarse SAND, trace of silt, wet
0	5-foot core with liner	40		11 12 13	SW	Soil sample PP5-10 collected at 15:50 Dark Brown-black fine to medium SAND, trace of gravel, wet
	inici			14		Dark Brown-black, fine to coarse gravelly SAND, wet
						Set temporary screen from 5 feet to 15 feet. Purged approximately 1-gallon prior to sample collection. Collected water sample PP5-W at 16:00 Backfilled borehole with bentonite chips.



Boring ID	PP6
Total depth	15 feet
Sheet 1	of 1

Project name	Pacific Park	Drilling Cont	ractor ESN	Drilling method Pu	ısh-probe rig
Project numbe	r <u>15-05986-040</u>	Location	SW corner of park, east side	Sampling method	5 ft core with plastic liner
Client King	County	of HESCO	es .	Air monitoring (Y/N)	Yes
HEC rep.	Bruce Carpenter	Date Ma	ay 23, 2017	Instrument(s)	Photoionization detector

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group	Soil description
(ррііі)		recovery	(leet)	1	GW	Grass/Brown sandy GRAVEL, FILL, damp
0.1	5-foot core with liner	40	<u>▼</u> 3.3	3 4	SM	Soil sample PP6-2 collected at 14:35 Brown gravelly silty SAND, piece of a ball, FILL, damp Static water level measured at 3.3 feet .
0	5-foot core with liner	60		6 7 8 9	SM	Dark Brown-black silty SAND, organic material, wet
0	5-foot core with liner	100		11 12 13	ML SP	Soil sample PP6-10 collected at 14:45 Dark Brown sandy SILT, organic material, wet Dark Brown-black, medium SAND, trace of silt, wet
				15	ML	Dark Brown sandy SILT, wet Set temporary screen from 2 feet to 12 feet. Purged approximately 1-gallon prior to sample collection. Collected water sample PP6-W at 15:00 Backfilled borehole with bentonite chips.



 Boring ID
 PP7

 Total depth
 15 feet

 Sheet
 1
 of
 1

Project name Pacific Park	Drilling Contractor ESN	Drilling method Push-probe rig
Project number <u>15-05986-040</u>	Location SE corner of park, ~ 125 feet	Sampling method 5 ft core with plastic liner
Client King County	south of MW-6	Air monitoring (Y/N) Yes
HEC rep. Bruce Carpenter	Date May 24, 2017	Instrument(s) Photoionization detector

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group	Soil description
0	5-foot core with liner	40/30/30	<u>▼</u> 1.8	1 2 3 4 5	SM	Dark Brown silty SAND, organic material, FILL, damp Soil sample PP7-0 collected at 14:30 Soil sample PP10-0 (duplicate sample, false time-14:35) Static water level measured at 1.8 feet Dark Brown sandy GRAVEL, trace of silt, cobbles, wet .
0	5-foot core with liner	75		6 7 8 9 10	SM	Dark Brown silty gravelly SAND, wet
0	5-foot core with liner	60		11 12 13 14		Soil sample PP7-10 collected at 14:40
						Set temporary screen from 0 feet to 10 feet. Purged approximately 1-gallon prior to sample collection. Collected water sample PP7-W at 14:50 Backfilled borehole with bentonite chips.



 Boring ID
 PP8

 Total depth
 15 feet

 Sheet
 1
 of
 1

Project name	Pacific Park	Drilling Co	ntractor <u>ESN</u>	Drilling method P	ush-probe rig
Project numbe	r <u>15-05986-040</u>	Location	~125 feet east-southeast of	Sampling method	5 ft core with plastic liner
Client King	County	MW-5		Air monitoring (Y/N)	Yes
HEC rep.	Bruce Carpenter	Date I	May 24, 2017	Instrument(s)	Photoionization detector

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group	Soil description								
W. F		,	,	1	ML	Dark Brown sandy gravelly SILT, some wood, FILL, damp Soil sample PP8-0.5 collected at 13:20								
0	5-foot core with	20/30		3	SM ▼	Brown silty gravelly SAND, FILL, damp wet								
	liner			4	3.1	Static water level measured at 3.1 feet .								
				5	GW	Dark Gray-black, fine to coarse sandy GRAVEL, wet								
0	5-foot core	30		7										
	with liner	with										9		
				10		Soil sample PP8-10 collected at 13:30								
	5-foot	5-foot			11		cobbles							
0	core with liner	40	40	13										
				14										
						Set temporary screen from 0 feet to 10 feet. Purged approximately 1-gallon prior to sample collection. Collected water sample PP8-W at 13:10 Collected duplicate water sample PP10-W (false time 13:45) Backfilled borehole with bentonite chips.								



 Boring ID
 PP9

 Total depth
 15 feet

 Sheet
 1
 of
 1

Project name Pacific Park	Drilling Contr	ractor ESN	Drilling method Po	ush-probe rig
Project number <u>15-05986-040</u>	Location	~100 feet south-southeast of	Sampling method	5 ft core with plastic liner
Client King County	MW-1		Air monitoring (Y/N)	Yes
HEC rep. Bruce Carpenter	Date Ma	ay 24, 2017	Instrument(s)	Photoionization detector

PID (ppm)	Sample type, interval	%	Water level	Depth (feet, BGS)	Soil	Soil description
(ррпп)	ilitervai	recovery	(feet)	BG3)	group	Gravel at surface
				1		Soil sample PP9-0 collected at 11:50
	·				GW	Brown-dark brown sandy GRAVEL, FILL, damp
0	5-foot core	20		2		
0	with	20		3		
	liner					wet
				4	<u>▼</u>	Static water level measured at 4.0 feet
				-	4.0	
				5	GW	Dark Brown fine to coarse sandy GRAVEL, small cobbles, wet
				6	0 "	Dark Brown line to coarse sailty GRAVEL, sman coooles, wet
	5-foot	• •		7		
0	core with	20		8		
	liner			0		
	inici			9		
				10		C-111- DD0 10114-1 -4 12:00
				11		Soil sample PP9-10 collected at 12:00
				- 11	SW	Dark Red-brown fine to medium SAND, wet
	5-foot			12	GW	Dark Red-brown fine to coarse sandy GRAVEL, wet
0	core	80		10		
	with liner			13		
	IIIICI			14		
					SW	Dark Brown-black fine to medium SAND, wet
				15	GW	Dark Brown-black sandy GRAVEL, wet
						Set temporary screen from 3 feet to 13 feet. Purged approximately 1-gallon prior to sample collection.
						Collected water sample PP9-W at 12:10
						Backfilled borehole with bentonite chips.



 Boring ID
 PP10

 Total depth
 8 feet

 Sheet
 1
 of
 1

Project name _Pacific Park	Drilling Contractor ESN	Drilling method Push-probe rig
Project number <u>17-06520-000</u>	Location NE corner of park	Sampling method 5 ft core with plastic liner
Client King County		Air monitoring (Y/N) Yes
HEC rep. Bruce Carpenter	Date <u>2/21/2018</u>	Instrument(s) Photoionization detector

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group	Soil description
V 6/		,	, <i>/</i>	1	GW	Grass/topsoil Brown sandy GRAVEL, trace of silt, small cobbles, FILL, damp
0	5-foot core	35		2		
	with liner			3		
			<u>▼</u> 5.0	5		
0	3-foot core	25	5.0	6	ML	Brown gravelly SILT, trace of sand, cobbles, FILL, wet Soil sample PP10-6 collected at 9:40
	with liner			7		
				8		Backfilled borehole with bentonite chips. Unable to penetrate beyond 6 feet in first borehole. Unable to penetrate beyond 8 feet in second borehole.



 Boring ID
 PP10A

 Total depth
 20 feet

 Sheet
 1
 of
 1

Project name Pacific Park	Drilling Contractor ESN	Drilling method Push probe
Project number <u>17-06520-000</u>	Location NE corner of park.	Sampling method 5 foot-core with plastic liner
Client King County	Immediately adjacent to PP10.	Air monitoring (Y/N) Yes
HEC rep. Bruce Carpenter	Date <u>2/28/2018</u>	Instrument(s) PID

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group	Soil description
0	5-foot core with liner	25		1 2 3 4	ML	Grass/Brown sandy SILT, trace of gravel, cobbles, brick fragment, FILL, damp
0	5-foot core with liner	55	5.0	6 7 8 9	CL GW	Dark Brown gravelly fine to coarse SAND, FILL, wet Brown gravelly silty CLAY, FILL, wet Gray sandy GRAVEL, pieces of concrete, wood, FILL, damp
0	5-foot core with liner	80		11 12 13 14	ML SW	Wood Gray-brown, sandy SILT, organic material, trace of gravel, wet Soil sample PP10-11 collected at 8:40 Gray-Dark brown, fine to coarse SAND, trace of gravel, wet Gray-Dark brown, sandy GRAVEL, cobbles, wet
0	5-foot core with liner	100		16 17 18 19 20	SW	Dark Brown-gray, fine to coarse SAND, trace of gravel, wet Soil sample PP10-17 collected at 9:08 Wood/thin silt zone (1")
						Backfilled borehole with bentonite chips.



 Boring ID
 PP11

 Total depth
 20 feet

 Sheet
 1
 of
 1

Project name Pacific Park	Drilling Contractor _ ESN	Drilling method Push probe
Project number <u>17-06520-000</u>	Location 35 feet northeast of levee	Sampling method 5 foot-core with plastic liner
Client King County	50 feet northeast of PP14	Air monitoring (Y/N) Yes
HEC rep. Bruce Carpenter	Date <u>2/28/2018</u>	Instrument(s) PID

PID	Sample type,	%	Water level	Depth (feet,	Soil	
(ppm)	interval	recovery	(feet)	BGS)	group	Soil description
					SW	Grass, Brown silty gravelly fine to coarse SAND, cobbles, FILL, damp
				1		
	5-foot			2		
0	core	30				Soil sample PP11-2 collected at 9:15
	with			3		
	liner					
				4	-	
			_	5		
			<u>▼</u> 5.0	3	GW	Soil sample PP11-5 collected at 9:20
				6		Dark Brown sandy GRAVEL, glass fragments, plastic, brick fragments,
						FILL wet
	5-foot	20		7		
0	core with	20		8		
	liner			0		
				9		
				10	-	G 1
				11	GW	Soil sample PP11-10 collected at 9:25 Dark Brown sandy GRAVEL, glass fragments, plastic, brick fragments,
				11	- 0 **	FILL wet
	5-foot			12		
0	core	40				
	with			13		
	liner					
				14	-	
				15	-	
				13		Soil sample PP11-15 collected at 9:25
				16		
	_				=	
	5-foot			17		
=	core with	=		18		
	liner			10	1	
				19]	
				20	-	Destablish to a shall saids beautiful to
						Backfilled borehole with bentonite chips.



 Boring ID
 PP12

 Total depth
 20 feet

 Sheet
 1
 of
 1

Project name Pacific Park	Drilling Contractor ESN	Drilling method _ Push probe
Project number <u>17-06520-000</u>	Location East of standing water, north	Sampling method 5 foot-core with plastic liner
Client King County	end of park	Air monitoring (Y/N) Yes
HEC rep. Bruce Carpenter	Date 2/21/2018	Instrument(s) PID

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group	Soil description
			<u>▼</u>	1	SW	Grass, Brown gravelly SAND, trace of silt, cobbles, FILL, wet
0	5-foot core with	30	<u>▼</u> 1.7	3		Soil sample PP12-2 collected at 9:50
	liner			4		Small piece of thin metal wire, plastic
			-	5	SM	Brown silty SAND, trace of cobbles and gravel, glass, plastic, brick
	5 C4			6		fragments, organic material, thin root fibers, wood, FILL, wet
0	5-foot core with	40		8		Soil sample PP12-7 collected at 10:00
	liner			9		
				10	SW	Soil sample PP12-10 collected at 10:10 Gray-brown fine to coarse SAND, wet
0	5-foot core	50		12		
	with liner			13		
			-	15		
				16	CW	Soil sample PP12-16 collected at 10:20
0	5-foot core with	100		17	SW	Dark Brown-gray fine to medium SAND, with a trace of silt, wet
	liner			19		
			_	20		Backfilled borehole with bentonite chips.



 Boring ID
 PP13

 Total depth
 20 feet

 Sheet
 1
 of
 1

Project name Pacific Park	Drilling Contractor ESN	Drilling method Push probe
Project number <u>17-06520-000</u>	Location West end of standing water	Sampling method 5 foot-core with plastic liner
Client King County		Air monitoring (Y/N) Yes
HEC rep. Bruce Carpenter	Date <u>2/21/2018</u>	Instrument(s) PID

PID	Sample type,	%	Water level	Depth (feet,	Soil	
(ppm)	interval	recovery	(feet)	BGS)	group	Soil description
			<u>▼</u> 0.3	1	ML	Grass, Brown fine sandy SILT, FILL, wet
			0.3	1		
	5-foot		-	2		
0	core	35				Soil sample PP13-2 collected at 10:30
	with			3		
	liner			4		
				5		
				(SW	Dark Brown-gray fine to medium SAND, FILL, wet
				6		Soil sample PP13-6 collected at 10:40
	5-foot			7		Son sample 11 13 0 conceied at 10.10
0	core	15				
	with			8		
	liner			9		
				,		
				10		
					SW	Soil sample PP13-10 collected at 10:50
				11		Dark Gray medium to coarse SAND, trace of gravel, cobbles, piece
	5-foot			12		of glass, FILL, wet
0	core	45		12		
	with			13		
	liner					
				14		
				15		
				13		Soil sample PP13-15 collected at 11:00
				16		Dark Gray fine to coarse SAND, with a trace of gravel, wood
						fragments, wet
	5-foot	100		17	SW	
0	core with	100	-	18		
	liner			10		
				19		
				20		
			-	20		Backfilled borehole with bentonite chips.
						Backfined outchole with bemonite chips.

PID = Photoionization detector



 Boring ID
 PP14

 Total depth
 20 feet

 Sheet
 1
 of
 1

Project name Pacific Park	Drilling Contractor ESN	Drilling method Push probe
Project number <u>17-06520-000</u>	Location 120 feet southwest of PP11,	Sampling method 5 foot-core with plastic liner
Client King County	30 feet northwest of levee.	Air monitoring (Y/N) Yes
HEC rep. Bruce Carpenter	Date 2/28/2018	Instrument(s) PID

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group	Soil description
V .		-	, ,	,	SW	Grass, Brown silty gravelly fine to coarse SAND, cobbles, FILL,
				1		damp
						Soil sample PP14-1 collected at 9:50
	5-foot			2		
0	core	20				
	with			3		
	liner					
				4		
				5		
				3		C-111- DD14 1114-1-4 0.55
				6	SW	Soil sample PP14-1 collected at 9:55 Dark Brown-gray, fine to coarse SAND, trace of gravel, FILL, damp
				U	SW	Dark Brown-gray, fine to coarse SAND, trace of graver, Fill, damp
	5-foot			7		
0	core	20		/		
	with	20		8		
	liner			0		
	inici			9		
						Water encountered in core sample at 9 feet.
				10		
				11		
					GW	Dark Brown-gray, fine to coarse sandy GRAVEL, cobbles, wet
	5-foot			12	SW	Dark Brown-gray, fine to coarse SAND, trace of gravel, wet
0	core	80				Soil sample PP14-12 collected at 10:00
	with			13		
	liner					
				14		
					ML	Brown-gray, fine sandy SILT, wet
				15	GW	Brown-gray, fine to coarse sandy GRAVEL, cobbles, wet
				16		
						Soil sample PP14-16 collected at 10:05
	5-foot	100		17	ML	Brown fine to medium sandy SILT, wet
0	core	100		10		
	with			18	CW	Constitution for the constitution of the CDAVIDIA 111
	liner			19	GW	Gray-brown, fine to coarse sandy GRAVEL, cobbles, wet
				19		
				20		
			}	20		Backfilled borehole with bentonite chips.
						Savanned obtenior with contonic emps.



 Boring ID
 PP15

 Total depth
 20 feet

 Sheet
 1
 of
 1

Project name Pacific Park	Drilling Contractor	ESN	Drilling method _ F	ush probe
Project number <u>17-06520-000</u>	Location ~55 fee	t SE of Geotech Boring	Sampling method	5 foot-core with plastic liner
Client King County	B-05		Air monitoring (Y/N)	Yes
HEC rep. Bruce Carpenter	Date 2/21/2018		Instrument(s)	PID

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group	Soil description
5-foot 0 core with liner		40	<u>▼</u> 2.3	1	SM	Grass, Brown silty SAND, with a trace of gravel, FILL, damp
				2	SM	Brown gravelly silty SAND, brick fragments, FILL, wet
				3	SM	
				5		Soil sample PP15-4 collected at 11:40 Proven grow gifty SAND, with some college, piece of gloss, thin piece.
		20		6		Brown-gray silty SAND, with some cobbles, piece of glass, thin piece of wire, brick fragments, rubber, FILL, wet Soil sample PP15-6 collected at 11:55
0	5-foot core with			8		
	liner			9		
		50		10	SM	Brown-gray silty SAND, with glass, wire, brick fragments, FILL, wet
	5-foot			11		
0 0	core with liner			13	SM	Soil sample PP15-12 collected at 12:05 Brown-gray silty SAND, wet
	inter			14		
		90		15	SW	Gray gravelly SAND, with a trace of silt, wet
0	5-foot core with liner			17		Soil sample PP15-16 collected at 12:15
				18		
				20		
						Backfilled borehole with bentonite chips.



 Boring ID
 PP16

 Total depth
 20 feet

 Sheet
 1
 of
 1

Project name Pacific Park	Drilling Contractor ESN	Drilling method Push probe
Project number <u>17-06520-000</u>	Location 50 feet west of PP16	Sampling method 5 foot-core with plastic liner
Client King County		Air monitoring (Y/N) Yes
HEC rep. Bruce Carpenter	Date <u>2/28/2018</u>	Instrument(s) PID

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group	Soil description
				1	SM	Grass, Brown silty fine to medium SAND, FILL, damp Soil sample PP16-1 collected at 10:30
0	5-foot core	30		2	SW	Brown-gravelly fine to coarse SAND, FILL, damp
	with liner			3		
				4		
			_	5		N
				6		No recovery – three attempts
	5-foot core	0		7		
	with liner			8		
				9		
				10	CM	
				11	GM	Dark Brown – gray, fine to coarse sandy GRAVEL, trace of silt, wet
	5-foot			12	SM	Soil sample PP16-11 collected at 10:45
0	core	55		12	GM	Dark Gray-brown, silty SAND, wet Dark Gray – brown fine to coarse sandy GRAVEL, wet
Ü	with liner	33		13	GIVI	Dark Gray Grown fine to course sandy Grove EE, wet
	IIIIei			14		
				15		
						Soil sample PP16-15 collected at 10:50
				16	SW	Dark Gray-brown gravelly fine to coarse SAND, piece of glass, wet
0	5-foot core	80		17	GW	Gray-brown fine to coarse sandy GRAVEL, wet
	with			18		
	liner			19	SW	Gray-brown gravelly fine to coarse SAND, wet
				20	GW	Gray-brown fine to coarse sandy GRAVEL, wet
						Backfilled borehole with bentonite chips.



 Boring ID
 PP17

 Total depth
 20 feet

 Sheet
 1
 of
 1

Project name	Pacific Park	Drilling Co	ontractor ESN	Drilling method Pu	ush probe
Project number	17-06520-000	Location	50 feet west of PP16	Sampling method	5 foot-core with plastic liner
Client King (County			Air monitoring (Y/N)	Yes
HEC repE	Bruce Carpenter	Date	2/28/2018	Instrument(s)	PID

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group	Soil description
(PPIII)	morvan	10001019	(1001)	200)	SM	Grass, Brown silty SAND, trace of gravel, FILL, damp
				1		
	7 C .			2	CIV	Soil sample PP17-1 collected at 11:00
0	5-foot	30		2	GW	Brown-dark gray, fine to coarse sandy GRAVEL, glass, plastic, brick, rubber, FILL, damp-wet
0	core with	30	•	3	<u> </u>	rubber, Filel, damp-wet
	liner		<u>▼</u> 2.7			
				4		
				5		Soil sample PP17-5 collected at 11:10
				6	GW	Brown-dark gray, fine to coarse sandy GRAVEL, glass, tile, brick,
					0,,	wood, FILL, wet
	5-foot			7		
0	core	20				
	with			8		
	liner			9		
				,		
				10	SM	Gray-brown silty SAND, wet
						Soil sample PP17-5 collected at 11:20
				11	GW	Dark Gray-brown fine to coarse sandy GRAVEL, cobbles, wet
	5-foot			12		
0	core	50		12		
	with			13		
	liner					
				14		
				15		
				13		
				16		
0	5-foot	100		17		
0	core with	100		18		
	liner			10		Soil sample PP17-18 collected at 11:28
				19	SM	Dark Gray-brown, silty fine SAND, wet
						wood
				20		D 1611 11 11 11 11 11 11 11 11 11 11 11 1
						Backfilled borehole with bentonite chips.



 Boring ID
 PP18

 Total depth
 20 feet

 Sheet
 1
 of
 1

Project name Pacific Par	Drilling Contract	or ESN	Drilling method P	ush probe
Project number 17-0652	0-000 Location Ap	proximately 40 feet west of	Sampling method	5 foot-core with plastic liner
Client King County	levee and soutl	n of PP16	Air monitoring (Y/N)	Yes
HEC rep. Bruce Carpe	nter Date <u>2/28/2</u>	018	Instrument(s)	PID

PID	Sample type,	%	Water	Depth (feet,	Soil	Cail decariation
(ppm)	interval	recovery	(feet)	BGS)	group SM	Soil description Grass, Brown silty SAND, FILL, wet
				1	SIVI	Glass, Blown sitty SAND, FILE, wet
				1	SW	Dark Gray-brown, gravelly fine to medium SAND, FILL, damp
	5-foot		_	2	5 **	Dark Gray-Grown, graverry fine to inequalif SAND, 1 feb., damp
0	core	35	<u>▼</u> 2.2			
	with			3		
	liner					Soil sample PP18-3 collected at 11:50
				4	SM	Light Brown silty SAND, with a trace of gravel, wet, FILL
						·
				5	SW	Light Brown-gray gravelly SAND, FILL, wet
						Soil sample PP18-5 collected at 12:00
				6	SM	Light Brown fine to medium sandy SILT, concrete, FILL, wet
	5-foot			7		
0	core	30				
	with			8		
	liner					
				9		
				1.0		
			_	10	CW	G '1
				1.1	GW	Soil sample PP18-10 collected at 12:05
				11		Dark Gray -brown fine to coarse sandy GRAVEL, glass, brick
	5-foot			12		fragments, FILL, wet
0	core	30		12		
U	with	30		13		
	liner			13		
	111101			14		
				15		
					ML	Soil sample PP18-15 collected at 12:10
				16		Dark Gray-brown fine sandy SILT, wet
	5-foot			17		
0	core	50				
	with			18		
	liner					
				19		
				20		W/ 1
			-	20		Wood Dealt Strad house also with house with a bine.
						Backfilled borehole with bentonite chips.



 Boring ID
 PP19

 Total depth
 20 feet

 Sheet
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 1

Project name _	Pacific Park	Drilling Con	tractor ESN	Drilling method Pu	ush probe
Project number	17-06520-000	Location	~30 feet south of the	Sampling method	5 foot-core with plastic liner
Client King C	County	Basketball	court	Air monitoring (Y/N)	Yes
HEC rep. B	ruce Carpenter	Date 2/	/21/2018	Instrument(s)	PID

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group	Soil description
W.I.		,	<u>▼</u> 0.5	1	ML	Grass, Brown sandy SILT, with a trace of gravel, FILL, wet
0	5-foot core with liner	35		3		Soil sample PP19-3 collected at 12:20
	mer			5		Soil sample 11 19-3 conceied at 12.20
	5.0			6	SM	Brown silty SAND, with a trace of gravel, pieces of concrete, FILL, wet
0	5-foot core with liner	40		8		Soil sample PP19-7 collected at 12:30
				9	GW	Gray sandy GRAVEL, with cobbles, FILL, wet
0	5-foot core with liner	50		11 12 13	GW	Soil sample PP19-10 collected at 12:40 Gray sandy GRAVEL, FILL, wet
				15	CIV	G 11
0	5-foot	60		16	SW	Soil sample PP19-15 collected at 12:50 Gray-brown fine to medium SAND, brick fragments, FILL, wet
V	with liner			18	GW	Gray sandy GRAVEL, brick fragments, glass, FILL, wet
				20		Backfilled borehole with bentonite chips.



 Boring ID
 PP20

 Total depth
 20 feet

 Sheet
 1
 of
 1

Project name _F	Pacific Park	Drilling Conti	ractor ESN	Drilling method Pu	ush probe
Project number	17-06520-000	Location	50 feet northwest of levee,	Sampling method	5 foot-core with plastic liner
Client King C	ounty	110 feet no	rth of PP8.	Air monitoring (Y/N)	Yes
HEC rep. B	ruce Carpenter	Date 2/2	28/2018	Instrument(s)	PID

PID (ppm)	Sample type, interval 5-foot core with liner	% recovery 30	Water level (feet) V 2.3	Depth (feet, BGS) 1 2 3 4	Soil group SM SM	Soil description Grass, Brown fine to medium silty SAND, FILL, damp Dark Brown fine to medium silty SAND, glass, plastic, trace of gravel, FILL, wet Soil sample PP20-2 collected at 12:50
0	5-foot core with liner	35		6 7 8 9	SM ML	Soil sample PP20-5 collected at 12:55 Dark Brown silty fine to medium SAND, plastic, glass, trace of gravel FILL, wet Dark Gray-brown silty fine sandy SILT, trace of gravel, FILL, wet
0	5-foot core with liner	30		11 12 13 14	SM	Soil sample PP20-10 collected at 13:05 Dark Gray-brown silty fine to coarse SAND, trace of gravel, glass, FILL, wet
0	5-foot core with liner	30		16 17 18 19	SM	Soil sample PP20-15 collected at 13:10 Dark Gray-brown silty fine to coarse SAND, trace of gravel, glass, FILL, wet
						Backfilled borehole with bentonite chips.



 Boring ID
 PP21

 Total depth
 20 feet

 Sheet
 1
 of
 1

Project name Pacific Parl	Drilling Contract	or ESN	Drilling method _ F	Push probe
Project number 17-0652	-000 Location Ap	proximately 80 feet west of	Sampling method	5 foot-core with plastic liner
Client King County	PP20 and 160	eet west of levee.	Air monitoring (Y/N)	Yes
HEC rep. Bruce Carpe	nter Date 2/28/2	018	Instrument(s)	PID

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group SM	Soil description Topsoil/Brown silty fine SAND, FILL, damp
0	5-foot core with liner	30	<u>▼</u> 2.1	3 4	SW	Dark Brown, gravelly fine to medium SAND, trace of silt, glass, FILL, wet Soil sample PP21-2 collected at 13:30 Wood -2"
0	5-foot core with liner	30		6 7 8 9	GW	Wood -4" Dark Brown-black fine to coarse sandy GRAVEL, trace of silt, wood, Soil sample PP21-6 collected at 13:35 plastic, glass, FILL, wet
0	5-foot core with liner	20		11 12 13 14	SM	Soil sample PP21-10 collected at 13:45 Dark Brown silty SAND, trace of gravel, wood, wet
0	5-foot core with liner	20		16 17 18 19 20	SM	Soil sample PP21-15 collected at 13:50 Dark Brown silty SAND, peat, wet Wood
						Backfilled borehole with bentonite chips.



 Boring ID
 PP22

 Total depth
 20 feet

 Sheet
 1
 of
 1

Project name Pa	acific Park	Drilling Cont	tractor ESN	Drilling method P	ush probe
Project number	17-06520-000	Location	North end of Baseball field	Sampling method	5 foot-core with plastic liner
Client King Co	ounty			Air monitoring (Y/N)	Yes
HEC rep. Bru	uce Carpenter	Date 2/	21/2018	Instrument(s)	PID

	Sample		Water	Depth		
PID (nnm)	type,	%	level (feet)	(feet, BGS)	Soil	Soil description
(ppm)	interval	recovery	(leet)	663)	group SM	Grass, Brown-gray silty SAND, with some gravel, FILL, damp
			▼	1	21.1	States, 210 mil gray emily entres, militarious graves, 1122, aminp
			<u>▼</u> 1.1			
	5-foot			2		
0	core	50				
	with			3	G) (
	liner			4	SM	Dark Brown silty SAND, with some gravel, plastic, wood, FILL, wet
				4		Soil sample PP22-4 collected at 13:40
				5		Son sample 11 22-4 conceiled at 13.40
					SW	Dark Gray fine to coarse SAND, with a trace of silt, FILL, wet
				6		, , , , , , , , , , , , , , , , , , , ,
	5-foot			7		
0	core	100				
	with			8		
	liner			9	SW	Soil sample PP22-9 collected at 13:50
					5 **	Dark Gray gravelly SAND, with a trace of silt, glass fragments, FILL,
				10		wet
					SW	Dark Gray gravelly SAND, with a trace of silt, FILL, wet
				11		
	- 0			- 10		
	5-foot	100		12		
0	core with	100		13		
	liner			13	GW	Soil sample PP22-13 collected at 13:55
				14		Dark Gray sandy GRAVEL, piece of plastic, FILL, wet
					ML/PT	Light Brown-gray SILT, with a trace of clay, peat, wet
				15		
				1.5	ML	Light Brown-gray SILT, with a trace of clay, fine sand, wet
				16		
	5-foot			17		
0	core	100		1 /		Soil sample PP22-17 collected at 14:05
	with	100		18		Some sample II BB I / Controlled at I 1100
	liner					
				19]	
			4	20		D 169 11 11 24 1 4 2 12
						Backfilled borehole with bentonite chips.
		l	1	l .	l .	I .



 Boring ID
 PP23

 Total depth
 20 feet

 Sheet
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Project name Pacific Park	Drilling Contractor ESN	Drilling method _ Push probe
Project number <u>17-06520-000</u>	Location 80 feet east of PP24, adjacent	Sampling method 5 foot-core with plastic liner
Client King County	to levee.	Air monitoring (Y/N) Yes
HEC rep. Bruce Carpenter	Date <u>2/28/2018</u>	Instrument(s) PID

S-foot S	PID	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group	Soil description
S-foot core with liner S-foot core so learned at 14:15 Dark Gray-brown, fine to medium SAND, trace of gravel Dark Gray-brown, gravelly fine to coarse, SAND, wet so learned at 14:20 Soil sample PP23-10 collected at 14:15 Dark Gray-brown, gravelly fine to coarse, SAND, wet soil sample PP23-15 collected at 14:20 Soil sample PP23-15 collected at 14:20 Dark Gray-brown, gravelly fine to coarse sandy GRAVEL, two pice soil sample PP23-15 collected at 14:20 Wet Dark Brown-gray fine to coarse SAND, trace of gravel, wet with liner S-foot lample PP23-15 collected at 14:20 Dark Brown-gray fine to coarse SAND, trace of gravel, wet wet lample PP23-15 collected at 14:20 Dark Brown-gray fine to coarse SAND, trace of gravel, wet lample PP23-15 collected at 14:20					1	SP	Grass, Brown silty fine SAND, FILL, damp
S-foot S-foot Soil sample PP23-2 collected at 14:00					1	SW	Light Brown-gray gravelly SAND, trace of silt, rubber, wood, plastic.
with liner 2.7		5-foot			2	5	
Iiner Soil sample PP23-5 collected at 14:05	0	core	30	<u>▼</u>			
Soil sample PP23-5 collected at 14:05 Swith Swith Swith Swith Swith Single Swith Swith Swith Swith Single Swith Swi				2.7	3		
Soil sample PP23-5 collected at 14:05 Light Brown-gray gravelly SAND, trace of silt, plastic, FILL, damp-wet Soil sample PP23-10 collected at 14:15 Light Brown-gray gravelly SAND, trace of silt, plastic, FILL, damp-wet Soil sample PP23-10 collected at 14:15 Dark Gray-brown, fine to medium SAND, trace of gravel Dark Gray-brown, gravelly fine to coarse, SAND, wet Soil sample PP23-10 collected at 14:15 Dark Gray-brown, fine to medium SAND, trace of gravel Dark Gray-brown, gravelly fine to coarse, SAND, wet Soil sample PP23-10 collected at 14:15 Dark Gray-brown, fine to coarse sandy GRAVEL, two pieces of silt, plastic, FILL, damp-wet Soil sample PP23-10 collected at 14:15 Dark Gray-brown, fine to coarse sandy GRAVEL, two pieces of silt, plastic, FILL, damp-wet Soil sample PP23-10 collected at 14:15 Dark Gray-brown, fine to coarse sandy GRAVEL, two pieces of silt, plastic, FILL, damp-wet Soil sample PP23-10 collected at 14:15 Dark Gray-brown, fine to coarse sandy GRAVEL, two pieces of silt, plastic, FILL, damp-wet					4		
Soil sample PP23-5 collected at 14:05 Light Brown-gray gravelly SAND, trace of silt, plastic, FILL, damp-wet Soil sample PP23-5 collected at 14:05 Light Brown-gray gravelly SAND, trace of silt, plastic, FILL, damp-wet Soil sample PP23-10 collected at 14:15 Dark Gray-brown, fine to medium SAND, trace of gravel Dark Gray-brown, gravelly fine to coarse, SAND, wet Soil sample PP23-10 collected at 14:15 Dark Gray-brown, fine to coarse, SAND, wet Dark Gray-brown, gravelly fine to coarse sandy GRAVEL, two pieces Soil sample PP23-15 collected at 14:20 wet Soil sample PP23-15 collected at 14:15 Dark Gray-brown, fine to coarse sandy GRAVEL, two pieces Soil sample PP23-15 collected at 14:20 wet Dark Brown-gray fine to coarse SAND, trace of gravel, wet Soil sample PP23-15 collected at 14:05 Dark Brown-gray fine to coarse SAND, trace of gravel, wet					5		
5-foot core with liner Sw Light Brown-gray gravelly SAND, trace of silt, plastic, FILL, damp-wet				1			Soil sample PP23-5 collected at 14:05
5-foot core with liner 13 SW Dark Gray-brown, fine to medium SAND, trace of gravel Dark Gray-brown, gravelly fine to coarse, SAND, wet 15 GW Dark Gray-brown, fine to coarse sandy GRAVEL, two piece Soil sample PP23-15 collected at 14:20 wet SW Dark Brown-gray fine to coarse SAND, trace of gravel, wet 17 Dark Brown-gray fine to coarse SAND, trace of gravel, wet					6	SW	
0 core with liner 9 Soil sample PP23-10 collected at 14:15 11 SW Dark Gray-brown, fine to medium SAND, trace of gravel 13 SW Dark Gray-brown, gravelly fine to coarse, SAND, wet 14 SW Dark Gray-brown, fine to coarse sandy GRAVEL, two piece Soil sample PP23-15 collected at 14:20 wet 5-foot core with SHOW Dark Brown-gray fine to coarse SAND, trace of gravel, wet 0 Soil sample PP23-10 collected at 14:15 Dark Gray-brown, fine to coarse sandy GRAVEL, two piece Soil sample PP23-15 collected at 14:20 Wet Dark Brown-gray fine to coarse SAND, trace of gravel, wet							
with liner Solity Solity	:	5-foot			7		
liner Soil sample PP23-10 collected at 14:15			30				
Soil sample PP23-10 collected at 14:15 11 SW Dark Gray-brown, fine to medium SAND, trace of gravel 12 SW Dark Gray-brown, gravelly fine to coarse, SAND, wet 13 SW Dark Gray-brown, gravelly fine to coarse, SAND, wet 15 GW Dark Gray-brown, fine to coarse sandy GRAVEL, two piece Soil sample PP23-15 collected at 14:20 wet 16 SW Dark Brown-gray fine to coarse SAND, trace of gravel, wet 17 SW Dark Brown-gray fine to coarse SAND, trace of gravel, wet					- 8		
Soil sample PP23-10 collected at 14:15 Dark Gray-brown, fine to medium SAND, trace of gravel 12		liner			9		
5-foot ore with liner 90 11 SW Dark Gray-brown, fine to medium SAND, trace of gravel 12 Dark Gray-brown, gravelly fine to coarse, SAND, wet 13 SW Dark Gray-brown, gravelly fine to coarse, SAND, wet 15 GW Dark Gray-brown, fine to coarse sandy GRAVEL, two piece Soil sample PP23-15 collected at 14:20 wet SW Dark Brown-gray fine to coarse SAND, trace of gravel, wet 17 Dark Brown-gray fine to coarse SAND, trace of gravel, wet					10		
5-foot with liner 13 SW Dark Gray-brown, gravelly fine to coarse, SAND, wet 15 GW Dark Gray-brown, fine to coarse sandy GRAVEL, two piece Soil sample PP23-15 collected at 14:20 wet 5-foot core of with 18 SW Dark Brown-gray fine to coarse SAND, trace of gravel, wet							Soil sample PP23-10 collected at 14:15
0 core with liner 90 13 SW Dark Gray-brown, gravelly fine to coarse, SAND, wet 14 Dark Gray-brown, fine to coarse sandy GRAVEL, two piece Soil sample PP23-15 collected at 14:20 wet SHOW Dark Brown-gray fine to coarse SAND, trace of gravel, wet 5-foot ore with 18					11	SW	Dark Gray-brown, fine to medium SAND, trace of gravel
0 core with liner 90 13 SW Dark Gray-brown, gravelly fine to coarse, SAND, wet 14 Dark Gray-brown, fine to coarse sandy GRAVEL, two piece Soil sample PP23-15 collected at 14:20 wet SW Dark Brown-gray fine to coarse SAND, trace of gravel, wet 5-foot ore with 18		5 0			10		
with liner 13			00		12		
liner 14 15 GW Dark Gray-brown, fine to coarse sandy GRAVEL, two piece Soil sample PP23-15 collected at 14:20 wet 5-foot 0 core 60 with 18			90		13	SW	Dark Gray brown grayally fine to coarse SAND wat
15 GW Dark Gray-brown, fine to coarse sandy GRAVEL, two piece Soil sample PP23-15 collected at 14:20 wet Sw Dark Brown-gray fine to coarse SAND, trace of gravel, wet 5-foot ore with 18					13		Dark Gray-orown, graverry line to coarse, SAND, wet
15 GW Dark Gray-brown, fine to coarse sandy GRAVEL, two piece Soil sample PP23-15 collected at 14:20 wet Soll sample PP23-15 collected at 14:20 wet SW Dark Brown-gray fine to coarse SAND, trace of gravel, wet 17 0 core 60 with 18		inici			14		
Soil sample PP23-15 collected at 14:20 wet Sw Dark Brown-gray fine to coarse SAND, trace of gravel, wet ocore 60 with 18							
5-foot 0 core with 16 wet 5-foot 17 Dark Brown-gray fine to coarse SAND, trace of gravel, wet 18					15	GW	Dark Gray-brown, fine to coarse sandy GRAVEL, two pieces of glass
5-foot 0 core with SW Dark Brown-gray fine to coarse SAND, trace of gravel, wet 17 18							
0 core with 17 18					16		
0 core 60 18	0 core			1.7	SW	Dark Brown-gray fine to coarse SAND, trace of gravel, wet	
with 18		60		1/			
		60	60	18	-		
							10
					19	GW	Dark Brown-gray, fine to coarse sandy GRAVEL, cobbles, wet
20					20		
Backfilled borehole with bentonite chips.							Backfilled borehole with bentonite chips.



 Boring ID
 PP24

 Total depth
 20 feet

 Sheet
 1
 of
 1

Project name Pacific Park	Drilling Contractor _ ESN	Drilling method _ Push probe
Project number 17-06520-0	0 Location Approximately 70 feet NW o	Sampling method 5 foot-core with plastic liner
Client King County	PP23 and 80 feet NW of levee.	Air monitoring (Y/N) Yes
HEC rep. Bruce Carpente	Date <u>2/28/2018</u>	Instrument(s) PID

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group SM	Soil description Grass/Brown silty SAND, FILL, damp
				1	SIVI	Grass/Brown Sirty SAND, FILL, damp
0.1	5-foot core	50	<u>▼</u> 2.3	2		Soil sample PP24-2 collected at 14:30
0.1	with liner	30	2.3	3	GW	Dark Brown-black sandy GRAVEL, glass, FILL, wet
	imer			4		
				5		
					SW	Brown-gray, gravelly SAND, glass, rubber, ceramic shards, FILL, wet
				6		
	5-foot	40		7		G '' 1 PP015 11 11 11 15
0.2	core with	40		8	GW	Soil sample PP24-7 collected at 14:35 Dark Gray-black fine to coarse sandy GRAVEL, rubber, glass, plastic,
	liner			9		FILL, wet
				10		
				11	GW	Soil sample PP24-10 collected at 14:45 Dark Gray fine to coarse sandy GRAVEL, with cobbles, wet
	5-foot	50		12		
0	with	50		13		
	liner			14		
				15		
				16	GW	Dark Gray, fine to coarse sandy GRAVEL, with cobbles, wet Sample was too coarse to collect soil sample
	5-foot			17		
0	core with	25		18		
	liner			19		
				20		
				20		Backfilled borehole with bentonite chips.



 Boring ID
 PP25

 Total depth
 20 feet

 Sheet
 1
 of
 1

Project name Pa	cific Park	Drilling Conti	actor ESN	Drill	illing method Po	ush probe
Project number	17-06520-000	Location	Behind pitcher's	mound San	mpling method	5 foot-core with plastic liner
Client King Cou	ınty			Air	monitoring (Y/N)	Yes
HEC rep. Brue	ce Carpenter	Date	21/2018	Inst	strument(s)	PID

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group	Soil description
				1	SM	Grass, Brown silty SAND, FILL, damp
	5-foot		<u>▼</u> 2.3	2	SP	Light Gray medium SAND, FILL, damp
0	core	85	2.3		SM	Light Brown-gray gravelly silty SAND, FILL, wet
	with liner			3		Soil sample PP25-3 collected at 14:20
	IIIICI			4		Soft sample 11 23-3 conceicd at 14.20
					SM	Light Brown-gray silty SAND, with a trace of gravel, FILL, wet
				5		
					SM	Light Brown-gray silty SAND, FILL, wet
				6		
	5-foot			7		
0	core	100				Soil sample PP25-7 collected at 14:25
	with			8		
	liner			0		
				9		
				10		
			1		SM	Light Brown-gray silty SAND, FILL, wet
				11		
	5 C 4			10		
0	5-foot core	50		12		
	with	30		13		
	liner				SM	Soil sample PP25-13 collected at 14:30
				14		Dark Brown-black silty SAND, Styrofoam, plastic, hydrocarbon odor,
				1.5		FILL, wet
			-	15		
				16		
	5-foot			17	.	
0	core	90		10	ML	Soil sample PP25-17 collected at 14:45
	with liner			18		Gray-brown sandy SILT, wet
				19		
			1	20		D 160 11 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1
						Backfilled borehole with bentonite chips.



 Boring ID
 PP26

 Total depth
 20 feet

 Sheet
 1
 of
 1

Project name	Pacific Park	Drilling Cor	ntractor	ESN	Drilling method P	rush probe
Project number	17-06520-000	Location	West of	Baseball field	Sampling method	5 foot-core with plastic liner
Client King	County				Air monitoring (Y/N)	Yes
HEC rep.	Bruce Carpenter	Date 2	/21/2018		Instrument(s)	PID

PID (nnm)	Sample type,	%	Water level	Depth (feet,	Soil	Soil description
(ppm)	interval	recovery	(feet)	BGS)	group SM	Soil description Grass, Light brown silty SAND, with a trace of gravel, FILL, damp
				1		
	5-foot		_	2	SM	Gray-brown gravelly silty SAND, FILL, wet
0	core with	65	<u>▼</u> 2.9	3		
	liner			4		Soil sample PP26-3 collected at 14:55
				5	SM	Gray-brown silty SAND, FILL, wet
				6		
	5-foot			7		
0	core with	25		8	SM	Soil sample PP26-7 collected at 15:05 Gray-brown silty gravelly SAND, FILL, wet
	liner				SIVI	Gray-brown sinty graverry SAND, FILL, wet
				9		
			-	10		
				11		
	5-foot			12	GW	Soil sample PP26-11 collected at 15:10 Dark gray sandy GRAVEL, with cobbles, FILL, wet
0	core	60				Dark gray sandy GRAVEL, with coopies, FILL, wet
	with liner			13		
	inici			14		
				15		
				16	GW	Dark gray sandy GRAVEL, with cobbles, FILL, wet
0	5-foot core	100		17	ML	Soil sample PP26-17 collected at 15:15
	with	100		18	1.112	Light Brown fine sandy SILT, with a trace of clay, some wood, wet
	liner			19		
				20		
			-	20		Backfilled borehole with bentonite chips.



 Boring ID
 PP27

 Total depth
 20 feet

 Sheet
 1
 of
 1

Project name Pacific Park	Drilling Contractor ESN	Drilling methodPush probe
Project number <u>17-06520-000</u>	Location 80' NW of levee and 70'SE	Sampling method 5 foot-core with plastic liner
Client King County	Of PP23	Air monitoring (Y/N) Yes
HEC rep. Bruce Carpenter	Date <u>2/28/2018</u>	Instrument(s) PID

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group	Soil description								
VI I		,		1	SM	Grass, Brown fine to medium silty SAND, FILL, wet								
0	5-foot core with	40		3										
	liner		<u>▼</u> 4.7	4	SW	Soil sample PP27-4 collected at 15:45 Brown fine to coarse gravelly SAND, FILL, wet								
			4.7	6	GW	Dark Brown to black, wood, plastic, glass, fine to coarse sandy GRAVEL, FILL, wet								
0	5-foot core with	50		7 8	ML	Dark Brown sandy SILT, FILL, wet Soil sample PP27-7 collected at 15:50								
	liner											9	GW	Dark Gray-black, fine to coarse sandy GRAVEL, FILL, wet
				10	CW	Soil sample PP27-10 collected at 16:00								
0	5-foot core with	50		11 12 13	GW	Dark Gray, fine to coarse sandy GRAVEL, glass, FILL, wet								
	liner			14										
	5-foot			16	GW	Soil sample PP27-15 collected at 16:10 Dark Gray, fine to coarse sandy GRAVEL, glass, wet								
0	core with liner	75		18										
				20		Destrilled housheds with houtswite abin-								
						Backfilled borehole with bentonite chips.								



 Boring ID
 PP28

 Total depth
 20 feet

 Sheet
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Project name Pacific Park	Drilling Contractor ESN	Drilling method
Project number <u>17-06520-000</u>	Location 110' northwest of PP27	Sampling method 5 foot-core with plastic liner
Client King County		Air monitoring (Y/N) Yes
HEC rep. Bruce Carpenter	Date 3/1/2018	Instrument(s) PID

PID	Sample type,	%	Water	Depth (feet,	Soil	
(ppm)	interval	recovery	(feet)	BGS)	group	Soil description Grass, Brown SILT, FILL, damp
				1	ML SW	Brown-gray, gravelly fine to medium SAND, FILL, damp-wet
				1	5 W	Brown-gray, graverry fine to medium SAND, FIEL, damp-wet
	5-foot			2		
0	core	45				
	with			3		
	liner					Soil sample PP28-3 collected at 10:30
			<u>▼</u>	4	GW	Brown-gray, fine to medium sandy GRAVEL, FILL, wet
			<u>▼</u> 4.3			
				5	ML	Brown-gray, fine to medium sandy SILT, wood, FILL, wet
					SW	Dark gray, gravelly fine to coarse SAND, FILL, wet
				6		
	<i>5.6.</i>			7		
0.2	5-foot	25		7		
0.2	core with	23		8		
	liner			0		Soil sample PP28-8 collected at 10:35
	inici			9		Plastic, glass, wood
						Timeter, genes, week
				10		
			1			Soil sample PP28-10 collected at 10:40
				11	SW	Gray gravelly fine to coarse SAND, wet
	5-foot			12		
0	core	40				
	with			13		
	liner			14	CW	Constructor and CDAVEL and
				14	GW	Gray fine to coarse sandy GRAVEL, wet
				15		
			1	13		Soil sample PP28-15 collected at 10:45
				16	SW	Gray-brown, fine to coarse SAND, wet
	5-foot			17	1	
0	core	60				
	with			18	GW	Gray fine to coarse sandy GRAVEL, wet
	liner					
				19		
				20		
			4	20		Destruited beautiful and the second of the s
						Backfilled borehole with bentonite chips.
	1		1	1	1	



 Boring ID
 PP29

 Total depth
 20 feet

 Sheet
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Project name Pacific Park	Drilling Contractor _ ESN	Drilling method Push probe
Project number <u>17-06520-000</u>	Location East of gravel road, west of	Sampling method 5 foot-core with plastic liner
Client King County	right field	Air monitoring (Y/N) Yes
HEC rep. Bruce Carpenter	Date <u>3/1/2018</u>	Instrument(s) PID

PID	Sample type,	%	Water	Depth (feet,	Soil				
(ppm)	interval	recovery	(feet)	BGS)	group SM	Soil description Grass, Light Brown silty fine SAND, FILL, damp			
				1	SIVI	Grass, Light Brown sitty fine SAND, FILL, damp			
				1					
	5-foot			2	GW	Light Gray sandy GRAVEL, FILL, damp			
0	core	50			J	Eight Gray bandy Gratt Ez, 1122, damp			
	with			3	GW	Gray-brown, fine to coarse sandy GRAVEL, FILL, damp-wet			
	liner		▼			Soil sample PP29-3 collected at 11:40			
			<u>▼</u> 3.1	4					
					GM	Light Brown silty sandy GRAVEL, FILL, wet			
				5	SW	Dark Brown gravelly fine to coarse SAND, FILL, wet			
						Soil sample PP29-5 collected at 11:45			
				6	GW	Gray fine to coarse sandy GRAVEL, FILL, wet			
	5-foot			7	SW	Gray fine to coarse SAND, FILL, wet			
0	core	80				, ,			
	with			8					
	liner								
				9	GW	Gray fine to coarse sandy GRAVEL, FILL, wet			
				10	OL	Gray brown, SILT, organic material, damp			
			1	10	OL	Soil sample PP29-10 collected at 11:50			
							11	SW	Gray silty fine to coarse SAND, wet
					~	Stary stary and to country starts, were			
	5-foot			12					
0	core	80							
	with			13	ML	Gray brown clayey SILT, trace of sand, wet			
	liner								
				14					
			1	15					
				1.0		Soil sample PP29-15 collected at 11:55			
				16	ML	Gray clayey SILT, wet			
	F.C. ,			17					
	5-foot	90		17	SM	Gray silty fine SAND, wet			
0	core with	90		18	SIVI	Gray Sitty lifte SAIND, wet			
	liner			10					
	iner			19	ML	Gray sandy SILT, wet			
					SM	Gray silty fine SAND, wet			
				20					
						Backfilled borehole with bentonite chips.			



 Boring ID
 PP30

 Total depth
 20 feet

 Sheet
 1
 of
 1

Project name Pacific Park	Drilling Contractor ESN	Drilling method Push probe
Project number <u>17-06520-000</u>	Location Approximately 100 feet west	Sampling method 5 foot-core with plastic liner
Client King County	O levee, south of PP23	Air monitoring (Y/N) Yes
HEC rep. Bruce Carpenter	Date <u>2/28/2018</u>	Instrument(s) PID

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group SM	Soil description Grass, Brown silty fine SAND, FILL, damp									
	5-foot			1 2	SIVI	Grass, Brown siny fine SAND, FILL, damp									
0	core with liner	40		3		Soil sample PP30-3 collected at 15:05									
				4	SW	Dark Gray-Brown gravelly, fine to coarse SAND, cobbles, 4" zone of iron oxidation, FILL, damp									
			<u>▼</u> 5.3	5	SM	Dark Brown silty fine SAND, FILL, damp									
			5.3	6	SW	Soil sample PP30-5 collected at 15:10 Dark Gray-brown gravelly fine to coarse SAND, glass, plastic, FILL, wet									
0	5-foot core	20		7		wet									
	with liner			8											
			9												
				10		Soil sample PP30-10 collected at 15:20									
		40		11	GW	Dark Gray-brown fine to coarse sandy GRAVEL, Cobbles, wet									
0	5-foot core		40	40	40		12								
	with liner														
				14											
				15		C 1 1 PP20 15 11 4 1 4 15 25									
						16	SW	Soil sample PP30-15 collected at 15:25 Dark Gray gravelly fine to coarse SAND, piece of glass, cobbles, wet							
0	5-foot core	40		17											
	with liner	. •	.0	- 🗸		- 3		.0	.0		18				
				19											
			-	20		Backfilled borehole with bentonite chips.									
						Backfined objetiole with bentomic emps.									



 Boring ID
 PP31

 1Total depth
 20 feet

 Sheet
 1
 of
 1

Project name	Pacific Park	Drilling Cor	ntractor ESN	Drilling method P	ush probe
Project number	17-06520-000	Location	120 feet northwest of PP30	Sampling method	5 foot-core with plastic liner
Client King	County	levee		Air monitoring (Y/N)	Yes
HEC rep.	Bruce Carpenter	Date 3	3/1/2018	Instrument(s)	PID

	Sample		Water	Depth		
PID	type,	%	level	(feet,	Soil	
(ppm)	interval	recovery	(feet)	BGS)	group	Soil description
					ML	Grass, Brown SILT, FILL, damp
				1	SW	Light Brown fine to medium gravelly SAND, FILL, damp
	5-foot	60		2	SW	Light Brown-gray gravelly, fine to coarse SAND, FILL, damp
0	core	60				
	with liner			3		Soil sample PP31-3 collected at 10:00
				4	SW	Light Brown-gray fine to coarse sandy GRAVEL, FILL, damp
			<u>▼</u> 5.5	5	SM	Dark Brown-gray silty gravelly fine SAND, FILL, damp
						Soil sample PP31-5 collected at 10:05
				6	SM	Gray-brown silty fine to medium SAND, trace of gravel, FILL, damp
	5-foot			7		
2.8	core	40			SM	Black silty SAND, wood, FILL, damp
	with liner			8		Predominantly wood
	IIIIei			9		Fredominantly wood
				10		
			-	10		Wood
				11	SW	Gray fine to medium SAND, wet
						Soil sample PP31-11 collected at 10:10
	5-foot			12		
0	core	70				
	with			13		
	liner					
				14		
				15		
			1	13		Call samula DD21 15 callected at 10:15
				16	SW	Soil sample PP31-15 collected at 10:15 Gray fine to coarse SAND, trace of gravel, wet
				10		Gray fine to coarse SAND, trace of graver, wet
	5-foot			17		
0	core	40		- /		
	with			18	GW	Gray fine to coarse sandy GRAVEL, wet
	liner					·
				19		
				20		
			-	20		D 1011 11 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1
						Backfilled borehole with bentonite chips.



 Boring ID
 PP32

 Total depth
 20 feet

 Sheet
 1
 of
 1

Project name Pacific Park	Drilling Contractor ESN	Drilling method Push probe
Project number <u>17-06520-000</u>	Location 160' northwest of PP7 and	Sampling method 5 foot-core with plastic liner
Client King County	200' northwest of levee	Air monitoring (Y/N) Yes
HEC rep. Bruce Carpenter	Date 3/1/2018	Instrument(s) PID

PID	Sample type,	%	Water level	Depth (feet,	Soil	
(ppm)	interval	recovery	(feet)	BGS)	group	Soil description
				1	ML SM	Grass, Brown SILT, FILL, damp Light Brown silty gravelly SAND, FILL, damp
				1	SIVI	Light brown sitty graverry SAND, Filel, damp
	5-foot			2		
0	core	40				
	with			3		
	liner			4	ML	Dark Brown-gray fine sandy SILT, FILL, damp
				4		Soil sample PP32-4 collected at 9:20
				5		Soil sample FF32-4 collected at 9.20
			-		SM	Light Brown-gray silty gravelly SAND, some wood, FILL, damp
				6		
	5-foot		<u>▼</u> 7.4	7		
0	core	35	7.4	0	CM	Soil sample PP32-7 collected at 9:30
	with liner			8	SM	Light Brown-gray silty gravelly SAND, some wood, FILL, wet
	IIIICI			9		
				10		
						Soil sample PP32-10 collected at 9:35
				11	GW	Brown-gray sandy GRAVEL, wet
	5-foot			12		
0	core	25		12		
	with	23		13		
	liner					
				14		
			-	15		C-il
				16	SW	Soil sample PP32-15 collected at 9:40 Dark gray fine to coarse SAND, trace of gravel, wet
				10	5 **	Dark gray fine to coarse SAIVD, trace of graver, wet
	5-foot			17		
0	core	60			GW	Dark gray fine to coarse sandy GRAVEL, wet
	with			18		
	liner			10		
				19		
				20		
			1			Backfilled borehole with bentonite chips.
						1



 Boring ID
 PP33

 Total depth
 20 feet

 Sheet
 1
 of
 1

Project name Pacific Park	Drilling Contra	actor <u>ESN</u>	Drilling methodP	Push probe
Project number 17-06520-0	0 Location	Southwest portion of the park.	Sampling method	5 foot-core with plastic liner
Client King County			Air monitoring (Y/N)	Yes
HEC rep. Bruce Carpente	Date 3/1/	/2018	Instrument(s)	PID

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group	Soil description
(ррііі)	Intorval	recovery	(ICCI)	ВОО)	SM	Grass, Gray-brown silty fine to medium SAND, FILL, damp
				1		
	5-foot			2		
0	core	50			SW	Gray-brown, fine to medium gravelly SAND, burned waste, plastic,
	with		<u>▼</u> 3.6	3		FILL, damp-wet
	liner		3.6	4		Soil sample PP33-3 collected at 11:05
				- +	GM	Gray-brown, silty sandy GRAVEL, FILL, wet
				5		
				(Soil sample PP33-5 collected at 11:10
				6		
	5-foot			7		
0	core	30			SM	Gray-brown silty fine SAND, FILL, wet
	with liner			8		
	IIIICI			9	GW	Gray-brown fine to coarse sandy GRAVEL, FILL, wet
]	
				10		C 1 1 PP22 10 11 4 1 411 17
				11	SW	Soil sample PP33-10 collected at 11:15 Gray fine to coarse SAND, trace of gravel, FILL, wet
				11	5,,,	Gray fine to course of fixe, face of graves, fines, wet
	5-foot			12		
0	core with	80		13		
	liner			13		
				14		
				1.5	ML	Gray fine sandy SILT, wet
				15	SM	Soil sample PP33-15 collected at 11:20; Gray fine silty SAND, wet
				16	ML	Gray fine sandy SILT, wet
				1-	g; ;	
0	5-foot core	100		17	SM	Gray fine silty SAND, wet
	with	100		18		
	liner				ML	Gray-brown, clayey SILT, wet
		ı	19	Gray-brown sandy SILT, wet		
				20	-	Gray-brown sandy SIL1, wet Gray clayey sandy SILT, wet
				-		Backfilled borehole with bentonite chips.



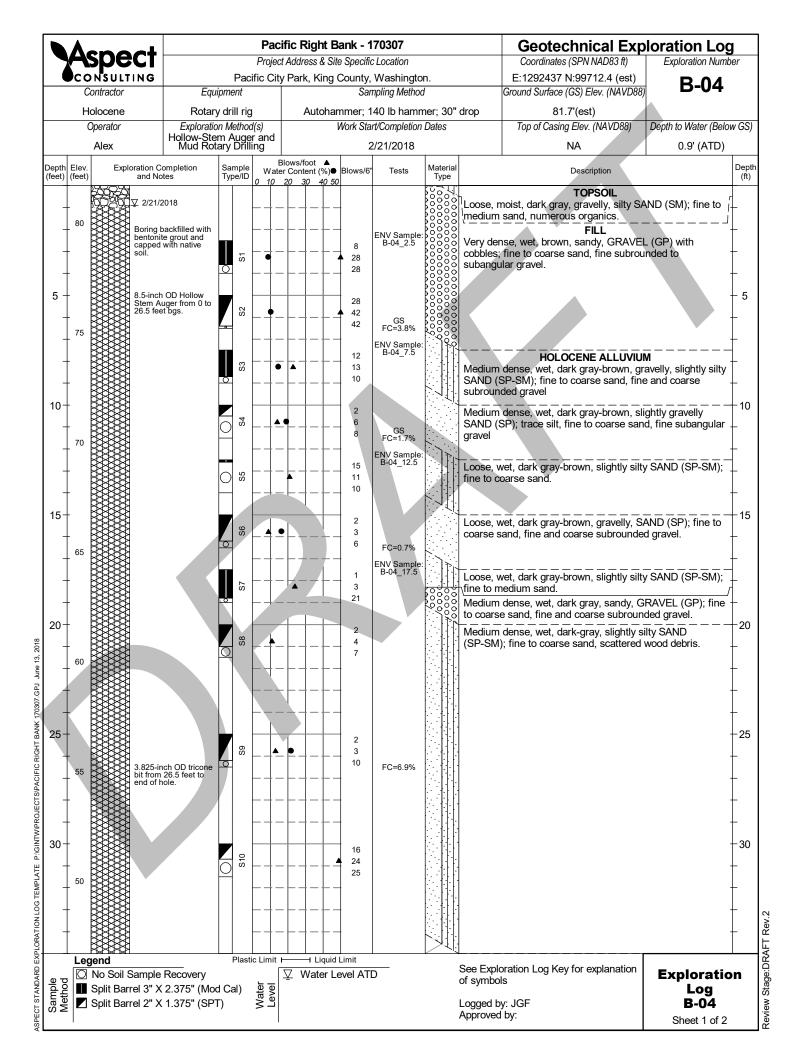
 Boring ID
 PP34

 Total depth
 20 feet

 Sheet
 1
 of
 1

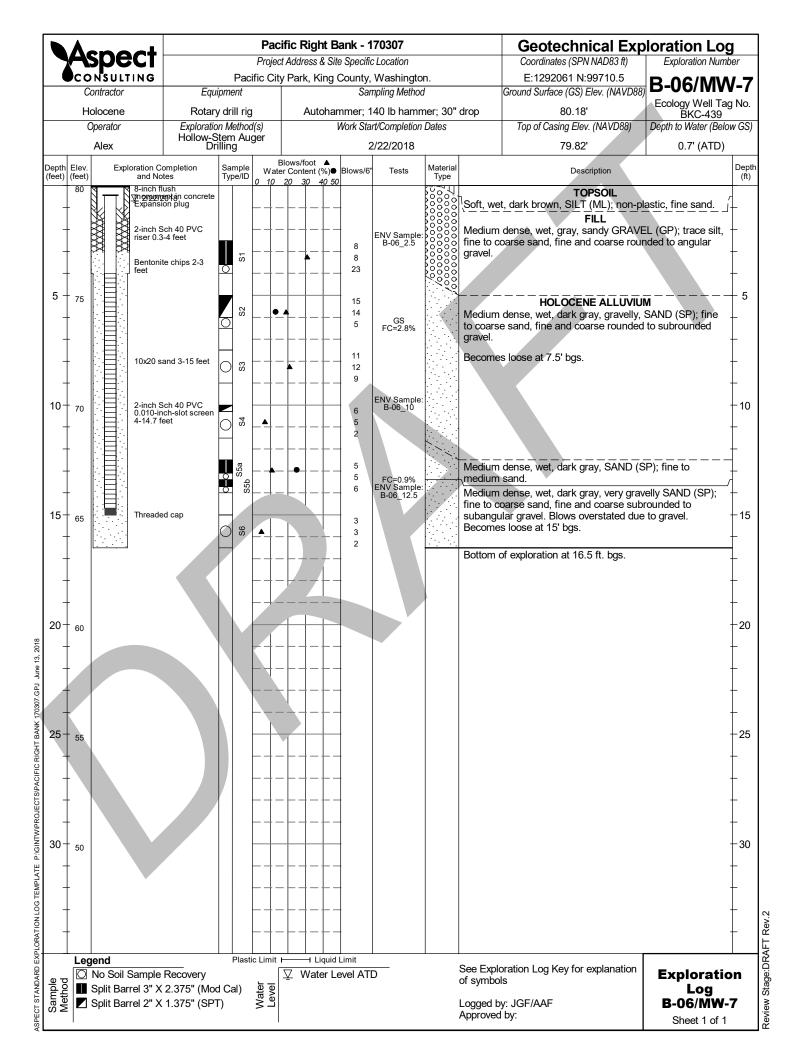
Project name Pacific Park	Drilling Contractor ESN	Drilling method Push probe	
Project number <u>17-06520-000</u>	Location Between geotech	borings B15 Sampling method 5 foot-core with plastic liner	
Client King County	and B16	Air monitoring (Y/N) Yes	
HEC rep. Bruce Carpenter	Date <u>2/21/2018</u>	Instrument(s) PID	

PID (npm)	Sample type, interval	% rocovery	Water level (feet)	Depth (feet, BGS)	Soil	Soil description
(ppm)	interval	recovery		BG3)	group SM	Grass, Light brown silty gravelly SAND, FILL, wet
			<u>▼</u> 0.9	1	5111	States, Eight ete wit entry graverity et in 12, 1122, wet
	5-foot			2		
0	core	40				
	with			3		
	liner			4		
				4	SW	Soil sample PP34-4 collected at 15:40
				5	5 **	Light Brown-gray gravelly SAND, FILL, wet
					SM	Light Brown-gray silty SAND, FILL, wet
				6		
	5-foot			7		
0	core	35				
	with			8		G '1 1 DD24 0 11 4 1 4 15 50
	liner			9		Soil sample PP34-8 collected at 15:50
				7		
				10		
					SW	Brown-gray fine to medium SAND, FILL, wet
				11		
	5-foot	4.0		12		
0	core	40		13		
	with liner			13	GW	Soil sample PP34-13 collected at 16:00
	IIIICI			14	G W	Brown sandy GRAVEL, FILL, wet
				11		Brown sundy Grant BE, TiBE, wet
				15	PT	Brown PEAT, 4-inch zone, wet
					ML	Soil sample PP34-15 collected at 16:05
				16		Light Brown-gray SILT, wet
	- 0					
	5-foot	100		17	CI	Class CH T and
0	core with	100		18	CL	Clayey SILT, wet
	liner			10	-	
				19		
				-	1	
				20		
						Backfilled borehole with bentonite chips.
			l			



	٨	cpost						-		170307		Geotechnical Exploratio	
	7	spect		_	_					fic Location		, , , , , , , , , , , , , , , , , , , ,	tion Number
•		ON SULTING Contractor	Equ			tic Cit	y Par	k, King		y, Washingto ampling Metho		E:1292437 N:99712.4 (est) Ground Surface (GS) Elev. (NAVD88)	-04
		Holocene						Autob		140 lb hamı		Ground Surface (GS) Liev. (NAVD00)	
		Operator	Rotar Exploration	on M	lethod	(s)		<i>r</i> autOH		tart/Completion	-		ater (Below GS
		Alex	Hollow-Ste Mud Rot	m /	Auger Drilli	and				2/21/2018			' (ATD)
nth	Elev						l Blows/	foot 🔺			Material		Dep
et)	(feet	and No		Ty	mple pe/ID	Wate 0 10	er Con 20	tent (%) 30 40	● Blows/6	6" Tests	Type	Description	(ft
-	45				S11	-	-		8 14			No recovery, gravel slough. No recovery	
0+	- - 40 -			7	S12	 			9 13 18	FC=35.4%		Dense, wet, gray, very silty SAND (SM); non-plastic sand, laminations of varying silt content, rare wood fragments and peat laminations.	5, fine 40
5 +	- 35 -			7	813		•		10 12 18			Dense, wet, gray, slightly silty SAND (SP-SM); low-plasticity, fine to medium sand, trace coarse rogravel, scattered clay and peat laminations.	45 unded
	30				\$14	1			5 5 6			Stiff, wet, gray, SILT (ML); non-plastic, trace fine so with scattered peat and silty sand laminations. Stiff, wet, green, CLAY (CL); low to medium plastic	
- - -	25			Z	S15	4	•		18 16 12			Medium dense, wet, gray, silty SAND (SM); fine to medium sand, scattered slightly silty sand lamination	
) -	20				S16		•		46 30 24			Very dense, wet, dark gray, slightly silty SAND (SP- fine to coarse sand. Bottom of exploration at 61.5 ft. bgs.	-SM); -60
5-	- - - 15					_							- - 65 -
Method		gend No Soil Sample Split Barrel 3" X Split Barrel 2" X	2.375" (Mo	d C		Water		⊣ Liqui Water	d Limit	TD TD		L	oration .og

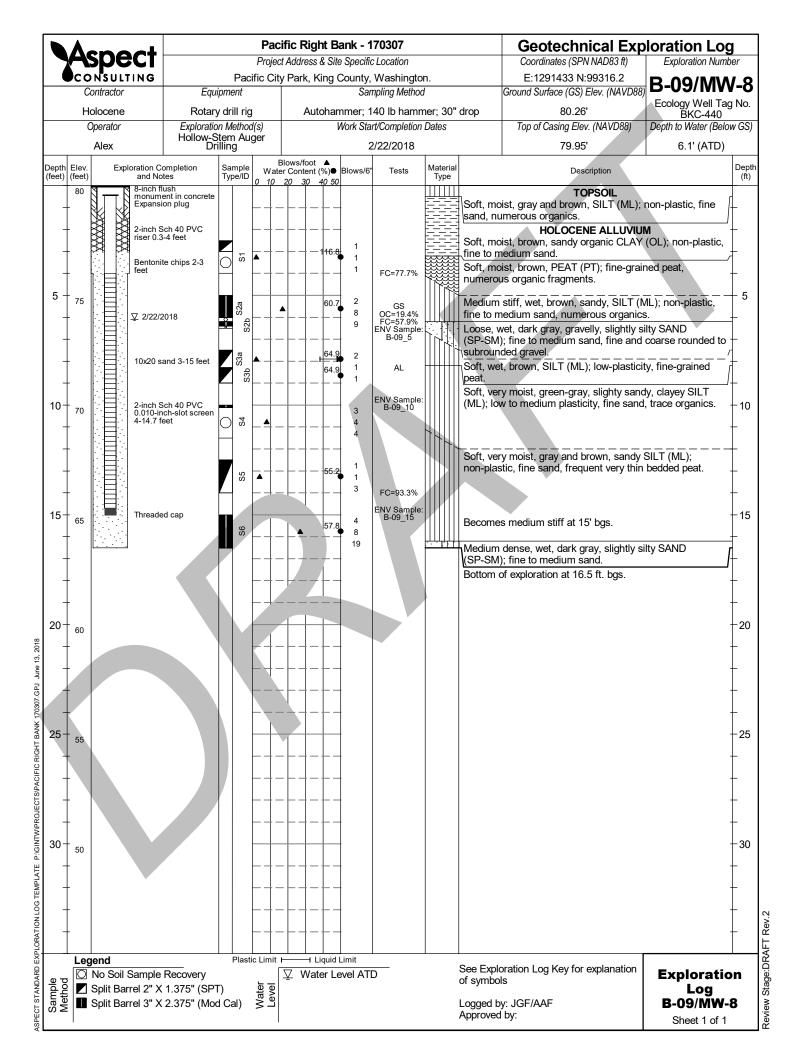
		<u>.</u>					-	Bank -				Geotechnical Exp		
	spec				-				ic Location			Coordinates (SPN NAD83 ft)	Exploration Num	ber
	ONSULTIN		quipm		ific Ci	ty Par	rk, King		, Washingto Impling Methor			E:1292254 N:99584.8 (est) Ground Surface (GS) Elev. (NAVD88)	B-05	
							Λι. 4 L		, 0		dran	` ' ' '		
	Holocene Operator	Explora	ary di ation M		<u> </u>		Autona		140 lb hamr art/Completion		arop	81.8'(est) Top of Casing Elev. (NAVD88)	Depth to Water (Belo	w G.S
	Alex	Hollow	-Ster	n Au	ger				2/20/2018	Datoo		NA	1.4' (ATD)	00
						Blows	/foot ▲		7/20/2010			IVA	1.4 (A1D)	1_
epth Elevet) (feet		Completion Notes	Sa Ty	mple pe/ID	Wa 0 10	ter Con	ntent (%)	● Blows/6	6" Tests	Material Type		Description		Dep (ft)
	a 8.5- Steet finds	g backfilled with nite grout and ad with native with OD Hollow Auger from 0 gs to end of		S10 S9 S8 S7 S6a S5 S4 S3 S2			Liqui		ENV Sample B-05_2.5 GS FC=3.9% ENV Sample B-05_7.5 ENV Sample B-05_17.5	00000	Very stift medium Medium fine to cogravel, research Medium coarse signavel. Embedium coarse signavel. Embedium medium medium numeror Medium (SP-SM) Medium (SP-SM) Medium (SP-SM) Medium medium numeror medium numero	dense, wet, brown, sandy, silty parse sand, fine and coarse ang metal sheet at 3' bgs. Inse, wet, gray, slightly silty SANI sand. HOLOCENE ALLUVIU wet, gray, sandy GRAVEL (GP); he and coarse subrounded to any dense, wet, dark gray, gravelly sand, fine and coarse subroundes slows overstated due to gravel. Idense, wet, dark gray, SAND (Soarse sand, fine and coarse sub oarse sand, fine and coarse sub oarse sand, fine and coarse sub oarse, wet, dark gray, gravelly, if the to coarse sand, fine and coarse, wet, dark gray, slightly so ngular gravel. Idense, wet, dark gray, slightly so if the to coarse sand, fine and coarse sub organics. Idense, wet, dark gray, slightly so if the to coarse sand, fine and coarse sub organics. Idense, wet, dark gray, slightly so if the to coarse sand, fine and coarse sub organics.	mon-plastic, fine to GRAVEL (GM); ular to rounded D (SP-SM); fine to M fine to coarse gular gravel. SAND (SP); fine to d to angular SP) with cobbles; rounded to angular M); fine to tions and lenses of slightly silty SAND oarse subrounded iity SAND	
pod T	egend No Soil Sam Split Barrel 3 Split Barrel 2	X 2.375" (M			Water Level	∇		d Limit Level AT	ΓD		See Explored by Approved	y: JGF	Exploration Log B-05 Sheet 1 of 1	on



	A.	- A	4									170307			Geotechnical Exp		
7		spe	CT									c Location			Coordinates (SPN NAD83 ft)	Exploration Num	
•		NSULT	ING		u iln		ific C	ity Pa	rk, K	ing (_	, Washingto			E:1291646 N:99695.2 (est)	B-07	
		ontractor	Rotary drill rig Autohammer; Exploration Method(s) Work St Hollow-Stem Auger and									mpling Method		alua ··	Ground Surface (GS) Elev. (NAVD88)		
		Olocene Operator			•				Aut			140 lb namn art/Completion		drop	81.4'(est) Top of Casing Elev. (NAVD88)	Depth to Water (Belo	ow GS
		Alex		Hollow-St	em /	Auae	er ánc	i		,		2/27/2018	Daics		NA	1.7' (ATD)	, vv GG
				l	T			Blows	s/foot	_		2/2//2016				1.7 (ATD)	T
	Elev. (feet)	Expl	loration C and No	ompletion ites		mple pe/ID		ater Co		(%)●	Blows/6	" Tests	Material Type		Description		Dept (ft)
			and No.	2018 2018 2018 2018 2018 2018 2018 2018								ENV Sample: B-07_2.5 ENV Sample: B-07_7.5 GS FC=2.7% ENV Sample: B-07_12.5	Type	Loose, v sand, tra Loose, v coarse s gravel, co Medium SAND (3 subroun Become Loose, v trace or coarse s trace or coars	TOPSOIL wet, dark brown, silty SAND (SM ace fine gravel. FILL wet, brown, sandy GRAVEL (GP sand, fine and coarse subrounde concrete fragments, trace organic HOLOCENE ALLUVIUI dense, wet, dark gray, very grav SP-SM); fine to coarse sand, fine ded to subangular gravel, trace of se fine to medium sand with no gravet, dark gray, GRAVEL (GP); fine to gravet. Wet, dark gray, SAND (SP); fine to wet, da	ity SAND carse rounded to	(ft)
1	-]_						stiff, wet, gray and brown, SILT t from cuttings.	(ML); silt, clay,	Ţ
	Loc					Plact	ic Lim	 		iquid I	limit						
od bot				Recovery 2.375" (M	od C			∇			evel AT	D		See Explo of symbo	oration Log Key for explanation ls	Exploration Log	on
Method				1.375" (SF		,	Water	ž						Logged b		B-07	
2 ر	1			,	′		_	1						Approved			

Managt			ific Right Bank - 17		Geotechnical Exp	loration Log
Aspect		-	t Address & Site Specific I		Coordinates (SPN NAD83 ft)	Exploration Number
Contractor	Equipmer		y Park, King County, V Sam	oling Method	E:1291646 N:99695.2 (est) Ground Surface (GS) Elev. (NAVD88)	B-07
Holocene	Rotary drill			0 lb hammer; 30" drop		
Operator	Exploration Mea	ethod(s)		/Completion Dates	Top of Casing Elev. (NAVD88)	Depth to Water (Below GS)
Alex	Hollow-Stem Au Mud Rotary D	uger and Drilling	2/	27/2018	NA	1.7' (ATD)
pth Elev. Exploration et) (feet) and N	Completion Sam Type	npie Wate	Blows/foot ▲ Blows/6" 20 30 40 50	Tests Material Type	Description	Dep (ft)
45	0	S	3 3 3 3		edium stiff, wet, gray and brown, SILT d peat from cuttings. (continued)	(ML); silt, clay,
	0	S12	12 15 15		edium dense, wet, gray, silty SAND (Sedium sand, scattered laminations of v	
5-	O	813	36 50/5"	ိုဝိဒ္ဓဝိဒ္ဓါ coa	ry dense, wet, dark gray, sandy GRAN arse sand, fine and coarse rounded to ows elevated due to gravel.	/EL (GP); fine to angular gravel. ——45
- - - - 30	0	41.8	10 18 50/5"	00000000000000000000000000000000000000	comes very dense at 50' bgs.	- - 50 - - -
5 - 25	0	315	44 A 31 17	. (SF	ense, wet, dark gray, slightly gravelly, s P-SM); predominantly fine to medium arse rounded to subangular gravel.	slightly silty SAND sand, fine and
)-		9	18	Vei	ry dense, wet, dark gray, silty SAND (SM); non-plastic,
20		816	30 30		e to medium sand, scattered thin beds	s or silt.
5-						- - 65 -
Legend No Soil Sample Split Barrel 3" Split Barrel 2"	e Recovery K 2.375" (Mod Cal		⊢ i Liquid Limit ☑ Water Level ATD	of s	Exploration Log Key for explanation symbols ged by: JGF proved by:	Exploration Log B-07 Sheet 2 of 2

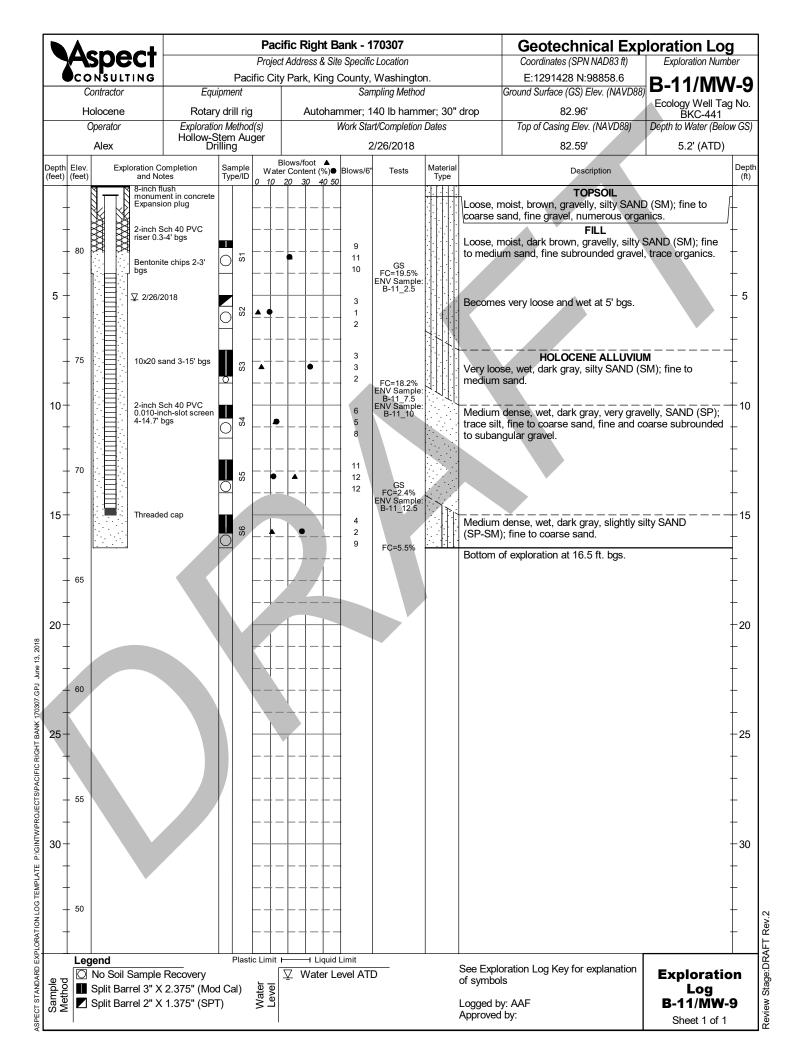
	\cn							_			170307			Geotechnical Exp		
Y	/ spo					-				•	c Location		_	Coordinates (SPN NAD83 ft)	Exploration Num	ber
	CONSU Contracto		E^	uipm		ific C	City Pa	ark, K	ing (Washingto			E:1291556 N:99480.2 (est) Ground Surface (GS) Elev. (NAVD88)	B-08	
	Holocen		Rota	•				Λ. ı+.	ohon		140 lb hamn		dron	80.5'(est)		
	Operator		Explorat	•		•	+	Auto			rt/Completion		urop	Top of Casing Elev. (NAVD88)	Depth to Water (Bel	ow GS
	Alex		Hollow-	Ster	n Au	ger			•		2/26/2018	24100		NA	4.4' (ATD)	
							Blow	s/foot	A						4.4 (////2)	
epth Ele eet) (fe		xploration C and No		Ty	mple pe/ID	0 1		ontent (30 4		Blows/6	Tests	Material Type		Description		Dep (ft)
5 - 7 - 7 - 7 - 7 - 7 - 7 - 7 - 7 - 7 -	75 55 55 55 56 56 56 56 56 56 56 56 56 56	bentoni capped soil. 2/26/2 8.5-inch Stem A	packfilled with le grout and with native 2018 a OD Hollow uger from 0 to end of		S8 S7a S6 S5 S4 S3b S2 S1				58:8	111 8 8 2 4 5 6 4 2	OC=2.7% ENV Sample B-08_5 ENV Sample B-08_7.5 GS FC=1.9% AL OC=3.5% ENV Sample B-08_12.5 OC=4.2% FC=84.8% ENV Sample B-08_17.5	7,000000000000 (000000000000000000000000	Sand, tra Medium trace sil subangu Loose, v (SP-SM subroun Become Very loo silt, fine subroun Become Very loo (SP-SM Medium non-plas peat lam Stiff, we sand. Medium (SP-SM Loose, v medium Become	TOPSOIL wet, dark brown, silty SAND (SM ace fine gravel. FILL I dense, wet, brown-gray, sandy it, fine to coarse sand, fine and coallar gravel. HOLOCENE ALLUVIUI wet, dark brown, gravelly, slightly i); fine to medium sand, fine rour ided gravel, trace organics. ses dark gray at 7.5' bgs. ses, wet, dark gray, very gravelly to coarse sand, fine and coarse ided gravel. ses loose at 10' bgs. Ty moist, gray, SILT (ML); low play initiations. Ty moist, dark gray, slightly i); fine to coarse sand. The stiff, very moist, brown-gray, SII stiff, very moist, brown-gray, SII stiff, very moist, brown-gray, SII stiff, very moist, dark gray, slightly initiations. The stiff of the sand, trace organications. The stiff of the sand organications or the sand, trace organications. The stiff of the sand organications or the sand organications. The stiff of the sand organications or the sand organications. The stiff of the sand organications or the sand organications or the sand organications. The stiff of the sand organications or the sand organications	GRAVEL (GP); parse rounded to M silty SAND ded to SAND (SP); trace rounded to asticity, frequent v silty SAND T (ML); cs, frequent thin on-plastic, fine silty SAND SM); fine to	
pod I	Split B	arrel 3" X	Recovery 2.375" (Mo			Water oil	∇		iquid I	Limit Evel AT	<u> </u> D		of symbo	by: JGF/AAF	Exploration Log B-08 Sheet 1 of 1	



	cna	▗ ▐▄						170307			Geotechnical Exp	ioration Log	<u> </u>
Y	spec			-				c Location			Coordinates (SPN NAD83 ft)	Exploration Numb	per
	ONSULTIN Contractor		Pa Equipment		y Park, I	King (, Washingto mpling Method			E:1291438 N:99086.4 (est) Ground Surface (GS) Elev. (NAVD88)	B-10	
	Holocene		-quipineni tary drill i		۸	itohan		140 lb hamn		dron	80.2'(est)		
	Operator		ation Meth	•	Au			nt/Completion		шор	Top of Casing Elev. (NAVD88)	Depth to Water (Belo	w GS
	Alex	Hollow-S	Stem Aud Rotary Di	ger and		-		2/26/2018			NA NA	2.7' (ATD)	,
				. 6	 Blows/foot	i 🔺						2.7 (7112)	D
epth Elev eet) (feet		ion Completion nd Notes	Samp Type/I	חו ייימויי	er Content 20 30			" Tests	Material Type		Description		Dept (ft)
5 - 75 - 75 - 75 - 75 - 75 - 75 - 75 - 7	Bo bela soi	nd Notes 2/26/2018 ring backfilled wintonite grout and pped with native	th to S s s s s s s s s s s s s s s s s s s		20 30		17 23 10 2 2 1 1 8 9 7 5 10 13 12 11 13	GS OC=0.6% FC=2.4% ENV Sample: B-10_7.5 GS FC=1.4% ENV Sample: B-10_12.5 GS FC=9.6% ENV Sample: B-10_17.5	Туре	Loose, vsand, nu Medium gravelly, coarse sorganics Medium coarse sorganics Medium (SP-SM coarse sorganics)	TOPSOIL wet, dark brown, silty SAND (SM) umerous organics. FILL dense, moist, dark brown and be silty SAND (SM); fine to coarse subrounded to angular gravel, asp HOLOCENE ALLUVIUI use, wet, brown-gray. slightly grave erbedded with soft, SILT (ML); lo um sand, fine subrounded gravel, wet, dark gray, very gravelly, SAN sand, fine and coarse subrounded	rown, very sand, fine and ohalt fragments. Velly, silty SAND by plasticity, fine wood fragments. ID (SP); fine to d gravel, rare SAND (SP); fine to d gravel. Ity SAND -plasticity, fine to	- 10 - 15 - 20 - 25
		nple Recovery 3" X 2.375" (N	/	Water Level		Liquid later Le	Limit evel AT	D		See Exploof symbo	oration Log Key for explanation	Exploration Log	- - - on

•	Λ	cpost							_		170307			Geotechnical Exp	loration Lo	g
7	7	spect			_						c Location			Coordinates (SPN NAD83 ft)	Exploration Num	ber
_		ON SULTING Contractor	Equ	iinm		ific C	City F	Park,	King		, Washingto mpling Method		-	E:1291438 N:99086.4 (est) Ground Surface (GS) Elev. (NAVD88)	B-10	
								۸.	ıt ah a							
		Holocene Operator	Rotar Exploration					Al			140 lb hamr art/Completion		агор	80.2'(est) Top of Casing Elev. (NAVD88)	Depth to Water (Belo	ow GS
		Alex	Hollow-Ste Mud Rot	m /	Auge	r and	t				2/26/2018			NA NA	2.7' (ATD)	,
nth	Elevi			1			Blov	ws/foc	ot 🔺			Material			2.7 (1115)	Dep
eet)	Elev (feet	Exploration C and No	ompletion	Ту	mple pe/ID	0 10	ater C 0 20	Conten	nt (%) € 40 5	Blows/6	" Tests	Material Type		Description ret, dark gray, slightly gravelly, s	lightly cilty SAND	(ft)
	-				S11		•	_ 4	<u> </u>	15 18	GS FC=9.1%		(SW-SM) gravel.	et, dank gray, signity gravelly, s ; fine to coarse sand, fine roun	ded to subrounded	
- 04 - -	40			0	S12					7 8 9			Becomes	medium dense at 40' bgs.		-40 -
- 5- - -	35			0	S13	_				5 3 2			Becomes (SP-SM);	loose, wet, dark gray, slightly s fine to coarse sand, fining dow	iilty SAND nhole.	+ +45 +
- 50 - - -	30			0	\$14					17 11 10	K		Medium c	dense, wet, dark gray, gravelly so coarse sand, fine and coarse ar gravel.	SAND (SP); trace subrounded to	-50 -
- 55 - - -	25				S15					14 12 10			Medium of sand.	dense, wet, brown-gray, silty SA	ND (SM); fine	-55
60-	20				S16					12 15 14				fine to medium sand at 60' bgs	s.	-60 -
- - 65 -	15													. 3		- - -65
-	- -						· — · — · —	-	-							 - - -
Sample Method		gend No Soil Sample Split Barrel 3" X Split Barrel 2" X	2.375" (Mo	d C		Water min	∇		Liquid ater L	Limit Level AT	<u> </u>		See Explor of symbols Logged by Approved I	: JGF/AAF	Exploration Log B-10 Sheet 2 of 2	

	Λ.	o 10							_		170307			Geotechnical Exp		
7		spe					-				ic Location			Coordinates (SPN NAD83 ft)	Exploration Num	
(N SULT	ING	_			ific C	ity Pa	rk, Kin		, Washingto			E:1290986 N:98067.3 (est)	B-100)
		Contractor			quipm						mpling Metho			Ground Surface (GS) Elev. (NAVD88)		
		lolocene Operator		Explora	,	rill rig	<u> </u>		Autoh		140 lb hami art/Completion		drop	77.5'(est) Top of Casing Elev. (NAVD88)	Depth to Water (Belo	ow GS
	,	Alex		Hollow-S Mud R	tem /	Auge	r and				2/27/2018	Dales				W GC
1		1			otar y	ווווט	ing		/foot ▲		2/2//2016			NA NA	8.8' (ATD)	1
	Elev. (feet)		loration Co and No	ompletion tes	Sa Ty	mple pe/ID	W a	ter Cor	ntent (%)	● Blows/6	5" Tests	Material Type		Description		Dep (ft
5	75 70 65 55 50 45 60		bentonit capped soil. 8.5-inch Stem At 10 feet I	oackfilled with e grout and with native OD Hollow uger from 0 to ogs.		S10 S9 S8 S7 S6 S5 S4 S3 S2				1 0 1 1 2.2 1 1 2 2 1 1 3 3 3 3 3 4 4 6 7 5 6 8 8 4 4 4 7 7 5 6 6 8 8 14 15 7 7 9 9	FC=29.7% GS FC=39% FC=40%		Numero Very loc SAND ((ML); no organics Loose, vorganics Soft, we Loose, vorganics Soft, we Loose, vorganics Soft, we Loose, vorganics Soft, we low plas Stiff, we dense, sono-plas organics Medium interbed	wet, dark brown, silty SAND (SMs. et, brown, PEAT (PT); coarse-grawet, dark gray, very silty SAND (Seganics. es with frequent thinly bedded silt ace organics. et, brown, slightly sandy, SILT (Mc.) interbedded slightly silty SAND (SP-SM) and stic to low plasticity, fine to medical dense, wet, dark gray, silty SAND (SP-sand, frequent thinly bedded sand, frequent thinly bedded sand, frequent thinly bedded sand.	M y, slightly silty oft, brown, SILT numerous); fine sand, trace sined peat. SM); fine sand, t, organic silt, and it, organic silt, and with medium silty SAND (SM); um sand, trace ID (SM) SM); fine to	
Method				Recovery 1.375" (SI	PT)	Plast	ic Limi	∇		id Limit Level A	D D	+ +1-19	See Explored of symbol	oration Log Key for explanation	Exploration Log	on
Jet 3	-	,	• •	- (2	,		Water Level						Logged b		B-100	
								1					Approved	1 by		



Pacific						_					Geotechnical Exploration Log		
1	hect			•			•	c Location		Coordinates (SPN NAD83 ft) Exploration Num			
	NSULTING ontractor	Faui		citic Cit	y Park	, King (Washingto			E:1291356 N:98461.5 (est) Ground Surface (GS) Elev. (NAVD88) B-12		
	olocene		Equipment Sampling Method Rotary drill rig Autohammer; 140 lb hammer; 30"								79.4'(est)		
	perator	Exploration Method(s)			Work Start/Completion Dates						Top of Casing Elev. (NAVD88) Depth to Water (Below		
	Alex	Hollow-S Dri	tem Au Illing	ıger			2	2/20/2018			NA NA	5.8' (ATD)	
	Exploration Co and No	ompletion tes	Sample Type/ID	Wate		ent (%)●		Tests	Material Type		Description		
apth Elev. (feet) - 75 - 75 - 70 - 70 - 65 60 60 60 60	and No Boring bentonit capped soil soil 8.5-inch Stem At	ackfilled with e grout and with native OD Hollow Iger from 0 to end of		Wate o 10	er Conte	oot Aent (%) • 0 40 50		OC=2.7% FC=5.1% FC=5.3%		Very loo fine to conumeror very loo interbed coarses very loo numeror very loo (SP-SM gravel. Loose, very loose,	TOPSOIL bist, gray-brown, SILT (ML); low	MD (SM); subangular gravel, but any subangular gravel, but any subangular gravel, but agments. (SM); fine sand, but any sand. RAVEL (GP); fine to subangular but any subangular but an	Dept (ft)
45										Bottom	of exploration at 31.5 ft. bgs.		<u>+</u> +
Leg			Plas	tic Limit		- - Liquid		<u> </u>	1	See Evol	oration Log Key for evolunation		1
	No Soil Sample Split Barrel 2" X		-)	Water Level	ΔM	Vater Le	evel AT	D		of symbol Logged b Approved	y: JGF	Exploration Log B-12	on

	chast			cific Right Bar			Geotechnical Exploration Log		
	spect		•	ect Address & Site S	•		Coordinates (SPN NAD83 ft)	Exploration Numb	er
	ONSULTING	Finit		ity Park, King Co	ounty, Washingto		E:1291717 N:98690.9 (est)	B-13	
	Contractor	Equipn		A 1	Sampling Method		Ground Surface (GS) Elev. (NAVD88)		
	lolocene Operator	Rotary d			mer; 140 lb hamr ork Start/Completion	· · · · · · · · · · · · · · · · · · ·	84.5'(est) Top of Casing Elev. (NAVD88)	Depth to Water (Below GS	
`	Alex	Hollow-Stem Mud Rotary	Auger and	770	2/23/2018	Duico	NA		w oo
			y Drilling	Blows/foot ▲	2/23/2016		NA NA	6.7' (ATD)	Т
epth Elev. eet) (feet)				iter Content (%)● Bl	lows/6" Tests	Material Type	Description		Dep (ft)
55 - 80 55 - 75 110 - 75 115 - 70 15 - 65 10 - 55	Boring to bentonit capped soil. 8.5-inch 21.5 fee	ackfilled with e grout and with native OD Hollow uper from 0 to t bgs. 018	S10 S9a S8a S7 S6 S5a S4 S3 S2 S1 S6 S6b S6b S6b S6b S6b S6b S6b S6b S6b		5 4 4 4 8 10 114 4 5 3 GS OC=8.5% FC=36.8% ENV Sample B-13_15.5 6 4 5 ENV Sample B-13_15 2 1 1 8 13 12 FC=9.7% ENV Sample B-13_17.5 16 32 31 6 6 11 8 8 21 1 7 11 8 FC=4.6%	Loose, (GP-Gi gravel, occord) Becom gravel. Mediur SAND gravel, occord fine an occord occor	FILL moist, brown, sandy slightly silty M); fine to coarse sand, fine subatrace wood fragments, strong iror les medium dense at 5' bgs. Blows les wet at 6.7' bgs. HOLOCENE ALLUVIUI m dense, wet, brown, slightly graw (SM); fine to medium sand, trace numerous organics. lense, wet, dark gray, GRAVEL (G d coarse gravel, trace organics (w losse, wet, dark gray, gravelly, slightly); fine to coarse sand, fine round trace organics. In dense, wet, dark gray, silty SAN sand, trace fine subrounded gravel lium sand, fine subrounded gravel m dense, wet, dark gray, slightly gravell ium sand, fine subrounded gravel m dense, wet, dark gray, slightly gravell ium sand, fine subrounded gravel m dense, wet, dark gray, slightly gravell ium sand, fine subrounded gravel m dense, wet, dark gray, very sand gitt, medium to coarse sand, fine an gular to angular gravel.	Melly, very silty fine and coarse P) with cobbles; ood fragments). Ty silty SAND ed to angular Ty SAND (SP); fine to el.	- 5 - 10 - 15 20 25
	gend		Plastic Limi			See Evr	ploration Log Key for explanation		1
	No Soil Sample Split Barrel 2" X Split Barrel 3" X	1.375" (SPT)	Cal) Mater	☑ Water Lev	el ATD	of symb	ools	Exploration Log B-13	'n

	snact	Pacific Right Bank - 170307											Geotechnical Exploration Log Coordinates (SPN NAD83 ft) Exploration Number			
Aspect		Project Address & Site Specific Location Pacific City Park, King County, Washington.											Coordinates (SPN NAD83 ft, E:1291717 N:98690.9 (e:			
	Contractor	Equ	uipme				airt, 1 t	9		npling Method			Ground Surface (GS) Elev. (NAV		3	
	lolocene	Rotar	-	_			Aut		-	40 lb hamr)" drop	84.5'(est)			
(Operator	Exploration Hollow-Ste	em A	luae	r and			١		rt/Completion	Dates		Top of Casing Elev. (NAVD88			
	Alex	Mud Ro	tary T	Dřilli	ing		s/foot		1	2/23/2018		1	NA NA	6.7' (AT	D)	
pth Elev. et) (feet)	Exploration C and No	ompletion otes	Sar Typ	mple pe/ID	W a	ater Co	s/100t ontent (30	(%)●	Blows/6'	Tests	Mater Type	ial :	Description		Dept (ft)	
				S11		•	_		17 17			Dense,	wet, dark gray, SAND (SP); to sand, trace fine subrounded	trace silt, fine to		
1			0	0)			7		18			·	r saira, trace fine subroarrace	a graver.	Ī	
Ť					\vdash	- -	7-	<u> </u>	Ī						Ť	
†					\vdash	- -	1-	-	1						†	
45					\vdash		-	-	-						+	
+ "				2a					6			Become	es loose at 41' bgs.		-40	
+				S12a S12b			4-	-	4 2				n stiff, wet, gray, sandy SILT	(ML): low to medium	+	
+				S	-+		4-	-	_			plasticity	y, fine sand.	(IVIL), IOW to medium	' 	
1						_	- ↓	L.							1	
1					$ldsymbol{ldsymbol{ldsymbol{ldsymbol{ld}}}$	$_{L}$	\rfloor	L_{-}								
40																
Ť				S13a		A		1	3 5						+45	
†				S13a S13b	\vdash		1-	 	9			Medium	dense, wet, dark gray, slight	tly silty SAND	7	
+					\vdash		-	-				Subrour	I); trace silt, fine to medium s nded gravel.	and, trace fine	+	
+					\vdash		4-	-	-						+	
-							4-	ļ.,	,						+	
35															-50	
1				S14					17 18			Become	es dense at 50' bgs.			
Ī									19			-[-]				
†					\vdash			7	1						Ť	
†					\vdash	-F	7	-	1		\	.[]			†	
30			M		\vdash										+	
+ 30						1			15			[]			-55	
+		· ·	Z	S15	-4	_ •	, _	_	19 24						+	
↓ .						_	4_	<u> </u>	24	FC=7.2%		.[-]			1	
						LL	1_	\perp							1	
												:[-]				
25			М					Γ								
				S16		Y			20 21			:[]			+60	
†				S	\vdash	- -	7-	<u> </u>	24			1				
+ 1					\vdash		-	-	1			Bottom	of exploration at 61.5 ft. bgs.		+	
† `					$\vdash \dashv$		4-	<u> </u>	-						+	
+					$\vdash \downarrow$	$ \mid$	4-	<u> </u>	-						+	
20					\Box	\perp		_							-65	
1						$_{L}$	1_	L_]	
T					\Box		7-								†	
†					-	- -	7-	† -	†						†	
15					\vdash		1-	-	†						+	
	gend			Plasti	c Limi	t 		iquid	 Limit		1					
	No Soil Sample				_ _	Ţ			evel ATI)		See Expl of symbo	loration Log Key for explanati ols	Exploia		
	Split Barrel 2" X Split Barrel 3" X			۱اد	Water Level	3						Logged b		Log B-13	ł	
≥ ■	Spir Dairei J A	. Z. 01 0 (IVIO	,u Oc	<i>ا</i> ا	> -	1						Approved	d by:	Sheet 2 d		

ΔV	cuvet					_	Bank - 1				Geotechnical Exp		
_	spect			-			Site Specifi				Coordinates (SPN NAD83 ft)	Exploration Num	
	ONSULTING Contractor	Equip		ıfic Ci	ty Par	k, Kin		, Washingto			E:1291958 N:98958.7 (est) Ground Surface (GS) Elev. (NAVD88)	B-14	
	Holocene	Rotary				Autoh		140 lb hamn		dron	83.9'(est)		
	Operator	Exploration	Method	d(s)		7 tatoi		art/Completion		чор	Top of Casing Elev. (NAVD88)	Depth to Water (Beld	ow GS
	Alex	Hollow-Ste Drill	em Aug ing	ger			:	2/21/2018			NA	3.8' (ATD)	
epth Elev	Exploration C	ompletion S	Sample		ter Con		● Blows/6	" Tests	Material Type		Description		Dep
- 80 5 - 75 10 - 70 15 - 65 20 - 60 25 - 55	Boring benton capped soil.	packfilled with the grout and with native 2018 DOD Hollow uger from 0 to end of	S10 S9 S8 S7 S6a S5 S4 S3 S2 S1	0 10	20	Lent (%)		ENV Sample B-14_5 ENV Sample B-14_10 FC=70.1% ENV Sample B-14_20 FC=70.1%	Type 0000 (Dense, I fine to coroad base Loose, I with coborounded fragmen Loose, I to coars gravel. Loose, I to coars gravel. Loose, I to coars gravel, so the coars gravel, to coars gravel.	FILL moist, brown, sandy, GRAVEL (Coarse sand, fine and coarse subsel levee crest. wet, dark gray-brown, gravelly, silbles; fine to medium sand, fine all to angular gravel, scattered organics. wet, brown, gravelly, slightly silty sibles; fine to medium sand, fine all to angular gravel, scattered organits. Wet, brown, sandy GRAVEL (GP) e sand, fine and coarse subrounded the sand, fine and coarse subrounded the sand, fine and coarse sand, fine subarticatered organics. Mense, wet, dark gray, slightly silty sand, fine and coarse sand. In gray, slightly sandy SILT (ML); and coarse sand sand. In gray, slightly sandy SILT (ML); and coarse sand sand. In gray, slightly sandy SILT (ML); and coarse sand sand. In gray, slightly sandy SILT (ML); and coarse sand sand sand sand sand sand sand sand	ty SAND (SM) and coarse anics. SAND (SP-SM) and coarse anics, PVC with cobbles; fine ded to angular wightly silty SAND and angular lty SAND RAVEL (GP); fine to subangular O (SP-SM); fine to non-plastic, fine d thin peat	5
thod [gend No Soil Sample Split Barrel 2" >			Water Level	\(\sum_{\subset}\)		id Limit Level AT	D		See Explored by Logged by		Exploration Log B-14	on

	۸	'n	~ +						_			70307			Geotechnical Exp	loration Log	g
X	<u>_/;</u>	he	CT			_	-				•	c Location			Coordinates (SPN NAD83 ft)	Exploration Numb	ber
		NSULT ontractor	ING	Fo	uipm		itic (ity Pa	ark, K	ung (Washingto			E:1291732 N:99109.4 (est) Ground Surface (GS) Elev. (NAVD88)	B-15	
		olocene		Rota			1		Aut	ohan		40 lb hamn		drop	80.4'(est)		
		perator		Explorat	ion N	1etho	d(s)				,	rt/Completion	,		Top of Casing Elev. (NAVD88)	Depth to Water (Belo	w GS
		Alex		Hollow-	Ster Orillin	n Au າg	ger				2	2/21/2018			NA	1' (ATD)	
epth E		Expl	oration C	ompletion otes		ample pe/ID		ater Co	s/foot ontent	(%)●	Blows/6	Tests	Materia Type		Description	1	Dept (ft)
5 -	75		bentoni capped soil. 8.5-inch Stem A	2018 Dackfilled with the grout and with native OD Hollow uger from 0 to end of	C	S2 S1					2 2 1	ENV Sample B-15_5		numero Become Very loc	see, moist, brown, silty SAND (St us organics, from cuttings. es wet at 1' bgs. use to loose, wet, silty SAND (SM including: ABS pipe fragments, p erial.	l); predominantly	- - - - 5
0+	70					S5 S4 S3					7 9 13 12 8 7 4 4 6	ENV Sample B-15_7.5		(SP-SM gravel. Medium fine to o gravel. Medium (ML); no	HOLOCENE ALLUVIUI wet, dark gray, slightly gravelly, s); fine to coarse sand, fine round dense, wet, dark gray, very grav coarse sand, fine and coarse roun stiff, moist, brown and gray, slig on-plastic, fine to medium sand, s	lightly silty SAND led to subangular relly, SAND (SP); nded to subangular htly sandy SILT	- - - - - 10
5-	65				7	S7 S6	4		-	•	3 2 3 5 10 12	FC=92.3% ENV Sample B-15_15 ENV Sample B-15_18.5		 Medium	slightly silty sand. I dense, very moist, gray, very silt nedium sand, scattered thin silt a ons.	ty SAND (SM); nd peat	- -15 - - - - - - -
25-	55					S9a S8					9 12 6 12 10	GS FC=3.7%		withGR	dense, wet, dark gray, SAND (S AVEL (GP); fine to coarse sand, I to subangular gravel.	SP) interbedded fine and coarse	- - - - -25
30 -	50				0	S10a S10b			A		6 12 14			Bottom	of exploration at 31.5 ft. bgs.		- - -30 - - -
Method		No Soil S Split Bar	rel 2" X	Recovery 1.375" (SR 2.375" (Me			Water In Inc.	∇		iquid I	Limit evel AT	<u> </u>		See Explored by See Explored b	by: JGF	Exploration Log B-15 Sheet 1 of 1	on .

	haat		Pau	ific Right Bank - 1	10301		Geotechnical Exp		
1/2	haci		-	t Address & Site Specific			Coordinates (SPN NAD83 ft)	Exploration Numb	per
	NSULTING ntractor	Equip		y Park, King County, San	Washington. Inpling Method		E:1291896 N:99309.9 (est) Ground Surface (GS) Elev. (NAVD88)	B-16	
	locene	Rotary			40 lb hammer; 30	l" dron	81'(est)		
	perator	Exploration		· ·	rt/Completion Dates	чиор	Top of Casing Elev. (NAVD88)	Depth to Water (Belo	w GS)
	Alex	Hollow-Stem Mud Rota	n Auger ánd Irv Drilling		2/23/2018		NA NA	1.3' (ATD)	,
epth Elev.	Exploration C and No	ompletion	Sample Wate	Blows/foot ▲ er Content (%)● Blows/6"	Tests Materia		Description	- ()	
Elev. (feet) (fe	and No	ch OD tricone 21.5 feet to ole.	Sample Wate	67 2 4 6 6 3 0 0 1 1 7 9 9 3 3 2 2 2 3 3 2 5 5 10 10 6 6	OC=6.6% ENV Sample: B-16_2.5 GS ENV Sample: B-16_12.5 ENV Sample: B-16_17.5 ENV Sample: B-16_17.5	Soft, mowithorgath broken is Become wood frage in the second wood in the second wood in the second wood wood wood wood wood wood wood w	FILL Dist, dark brown, sandy SILT (ML anic SILT (OL); low to medium pla	ND (SM); fine to ace organics and ND (SP); fine to do angular Shtly gravelly SILT ine subrounded EL (GP); medium SAND (SM); fine Sand, fine and	Dept (ft)
ਜੂ ਨੂੰ Ⅲ s	end No Soil Sample Split Barrel 3" X Split Barrel 2" X	2.375" (Mod	Cal) $\frac{1}{2}$ $\frac{1}{2}$	⊢	<u> </u>	See Expl of symbo		Exploration Log B-16	on

		٨.							_		- 170307					Geotechnica	I Exp	loration Log	g
1	X	1	spec ¹				-				cific Location					Coordinates (SPN NAD	83 ft)	Exploration Num	ber
_			SULTING	I	guipn		ific Ci	ty Par	k, Kin		nty, Washing					E:1291896 N:99309.9		B-16	
			ontractor						A		Sampling Met		~			Ground Surface (GS) Elev. (ΝΑΥΔΟΘ)		
			Olocene Operator	Rot Explora		rill rig			Autoh		r; 140 lb hai Start/Complet				arop	81'(est) Top of Casing Elev. (NA	VD88)	Depth to Water (Belo	ري الار (25
			•	Hollow-S Mud F	Stem	Auge	r and			VVOIK	•		ales	•			v D00)		w GS
	1		Alex		Rotary	/ Drilli			foot ▲		2/23/2018	3		1		NA NA		1.3' (ATD)	
epth eet)	(fe	lev. eet)	Exploration and	Completion Notes	S: Ty	ample /pe/ID	Wat	ter Con	tent (%)	● Blow <i>50</i>			Mate Typ	rial e F1.	Modium	Description		oliabthy oilty SAND	Dept (ft)
- - 40 - - -	- - -	40				S12			6	5.2 2 2 11	FC=6.4 ⁹	%			(SP-SM) coarse s	dense, wet, dark gray, g); predominantly fine to r subrounded to subangula ist, brown, sandy, organ ous PEAT (PT); non-plas	nedium ir gravel	sand, fine and (continued)	+ 40
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APPENDIX B

Abandoned Landfill Study in King County



ABANDONED LANDFILL STUDY IN KING COUNTY

SEATTLE-KING COUNTY DEPARTMENT OF PUBLIC HEALTH
APRIL 30, 1985



TABLE OF CONTENTS

	PAGE
ACKNOWLED GEMENTS	2
EXECUTIVE SUMMARY	3
CHAPTER I. INTRODUCTION	
 BACKGROUND INFORMATION 	4
• HISTORICAL CONCERNS	6
• FIGURE I	10
• TABLE I	. 11
CHAPTER II. METHODS AND MATERIALS	12
CHAPTER III. RESULTS	16
 HISTORICAL AND FIELD DATA 	
• MAPS AND TABLES	
• OTHER SITES	
CHAPTER IV. CONCLUSIONS AND RECOMMENDATIONS	135
GLOSSARY	145
REFERENCES	146

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Greg Bishop Project Co-Leader

Wayne L. Turnberg Project Co-Leader

Karen VanDusen Historical Research

EXECUTIVE SUMMARY

The King County Abandoned Landfill Survey was conducted from October through December 1984 by the Health Department's Environmental Health Division at the request of the King County Council. The twenty sites were selected during a September 17, 1984 meeting between the Seattle-King County Department of Public Health, the Solid Waste Division. (Subsequent to this meeting three additional sites were added to the list.) The primary objective of the survey was to determine if any public health problems existed at the predetermined sites.

The Health Department researched the geographical and historical data on each site with input from various city and county agencies. This information was used by the field staff for guidance in determining where the sampling efforts should be focused.

Due to the limited time and resources available for the study, the staff decided to conduct a primary survey utilizing general sampling parameters. In the event the primary survey revealed any environmental health problems, then funds would be requested for more detailed followup work. The parameters chosen as criteria for the primary survey were:

- 1. Gas Methane and non-specific organics/inorganics exclusive of methane;
- 2. Water Conductivity, dissolved oxygen, pH, turbidity and temperatures.

The report presents the results of the preliminary survey including the field and historical data for each site. Of the twenty-three sites studied, eleven were found to have no significant environmental health problems. Better leachate or methane control are recommended at four sites. It is further recommended that additional water and/or soil samples be taken and analyzed at eight sites for primary organics and inorganics that may be toxic to humans.

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CHAPTER I INTRODUCTION

- BACKGROUND INFORMATION
- · HISTORICAL CONCERNS
- FIGURE I
- · TABLE I

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BACKGROUND INFORMATION

Land use is a major concern for any governmental entity, but particularly for those located in urban areas. In King County the increasing demand for residential, commercial and industrial sites causes both the public and private sectors to look for appropriate, effective and efficient means to use and/or develop remaining vacant land. Reclaiming land through the use of sanitary landfills which are later developed for many different uses has been an acceptable pattern of land use development for years. At least one-hundred-fifty acres of land were so reclaimed in the City of Seattle. However, since the complicated community impacts of New York's "Love Canal," a long forgotten hazardous waste landfill, became known in the 1970's, both public and private citizens have become increasingly sensitive to the legacy of long accepted waste disposal practices. It has become increasingly clear that some sites previously used for landfills are not appropriate for certain types of development due to the potential risk they present for chronic exposure to hazardous materials for users or developers of the site. Other sites have negatively impacted ground and/or surface waters. As a result, several federal, state and local regulations have been passed in an effort to change long accepted patterns of waste disposal and studies have been undertaken or are underway to document existing or abandoned sites which pose either environmental and/or human health hazards.

Against this national background of activity and concern, in September 1984 the King County Council requested the Environmental Health Division of the Seattle-King County Health Department to conduct a preliminary study of the abandoned landfills in King County. In undertaking this study, the Department focused on these questions:

1. Where are the abandoned sanitary landfills in King County?

- 2. Have abandoned landfills at the currently known sites undergone enough stabilization to allow development to safely occur on site?
- 3. Is there any evidence to suggest that the abandoned site may contain materials known to be toxic to human health which could impact human health during site development or due to chronic exposure from later use of the site?
- 4. Is there any evidence to suggest that the site may be negatively impacting the surrounding community via ground/surface water contamination or methane migration?

Both field studies and a search of site specific historical records were conducted in order to answer these questions. In addition, generalized information about waste disposal practices, provided a background perspective regarding the significance of the problem in this County.

HISTORICAL CONCERNS

A preliminary review of abandoned landfills in King County suggests a picture of "probably anything" located "almost anyplace." Garbage was defined by the 1931 Superintendent of the Garbage Division of the Health and Sanitation Department, C.L. Murray, as, "Everything that is wasted from the home, the business house, manufacturing plant...We collect this material without any separation..."

The material collected went to sanitary fills at convenient spots throughout the area. Low lands such as tide lands, marsh or dry ravines were favorite locations. It was not unusual for property owners to lease land to the County for a garbage dump in order to level the property, making it more desirable for development. In addition, the County was plagued by promiscuous dumping on vacant property and dead-end streets.

As noted in a 1970 Public Works document, <u>King County Solid Waste Disposal</u> - <u>For 2020 Vision, Vol. I</u>, "Historically the major objective of solid waste disposal was to get waste out of residential neighborhoods into rural dumping areas at a minimum of cost to the taxpayer. These objectives, shared by both officials and citizens, were conducive neither to operating disposal sites in a sanitary manner nor to preventing environmental pollution. However, 'the out of sight, out of mind' attitude toward solid waste disposal was shared across the nation, and any system which moved solid waste away from citizens' homes was generally deemed an adequate system!

One of the major concerns left from the legacy of solid waste disposal practices in the past are the number of abandoned sites at forgotten locations, long since developed for other uses. Another concern is the traditional practice of mixing commercial/industrial and residential wastes. According to a 1980 Battelle report on the "Identification of Hazardous Waste Disposal Sites and

Management Practices in Region 10," it was not unusual for industries to use regional landfills for waste disposal. This is underscored by a report released by CH₂M Hill in August 1969 entitled, <u>Seattle Area Oil Waste Disposal Facility Study</u>. This report notes that the <u>closing of landfills</u> to oil waste disposal due to water pollution and the cessation of open burning of oil due to air pollution made it difficult to properly and legally dispose of oily wastes from shipyards, industries and service stations. These wastes were noted to contain crude, diesel, lubricating, cutting, cooling, cooking or any other form of mineral or vegetable oil.

Thus the composition of the waste disposal of the several landfills which have been abandoned is of concern from a historical perspective.

Another and well known problem associated with solid waste landfills is the contamination of ground and/or surface waters due to leachate. Closed landfills without leachate interception equipment can contribute to ground water contamination. A draft EIS on the 1982 Comprehensive Solid Waste Management Plan for King County notes that leachate from landfills is the major environmental problem currently caused by solid waste disposal in King County. Unfortunately few, if any of the abandoned sites studied in this report had leachate interception and thus not only current but past landfills as well may be contributing to ground water pollution.

Many of the abandoned landfill sites in King County come as a result of a new strategy in waste disposal. The expanding build-up of residential areas in the 1950's began to crowd out potential future County landfill disposal sites. Old sites were reaching their planned capacities. At the same time the construction of Interstate 5 forced the closure of three large County landfills. Two of these handled 75% of the solid waste in King County. In 1960 the shift from using

several landfill sites to the use of transfer stations in conjunction with large remote (at that time) disposal sites was adopted. Eventually seven transfer stations eliminated fifteen of the previously used open dumps in the County. What happened to those closed sites and their existing conditions is the concern of this report.



ABANDONED LANDFILL SITES

On September 17, 1984 a meeting was held between representatives of the Seattle-King County Department of Public Health and King County's Solid Waste division to determine the abandoned landfills to be targeted for this study. A list of twenty sites was organized based on the information available at that time. Subsequent to this meeting three additional sites were added to the list. All sites are presented in Table I and Figure 1.



CHAPTER II METHODS AND MATERIALS



FIGURE 1

KING COUNTY ABANDONED LANDFILL LOCATIONS

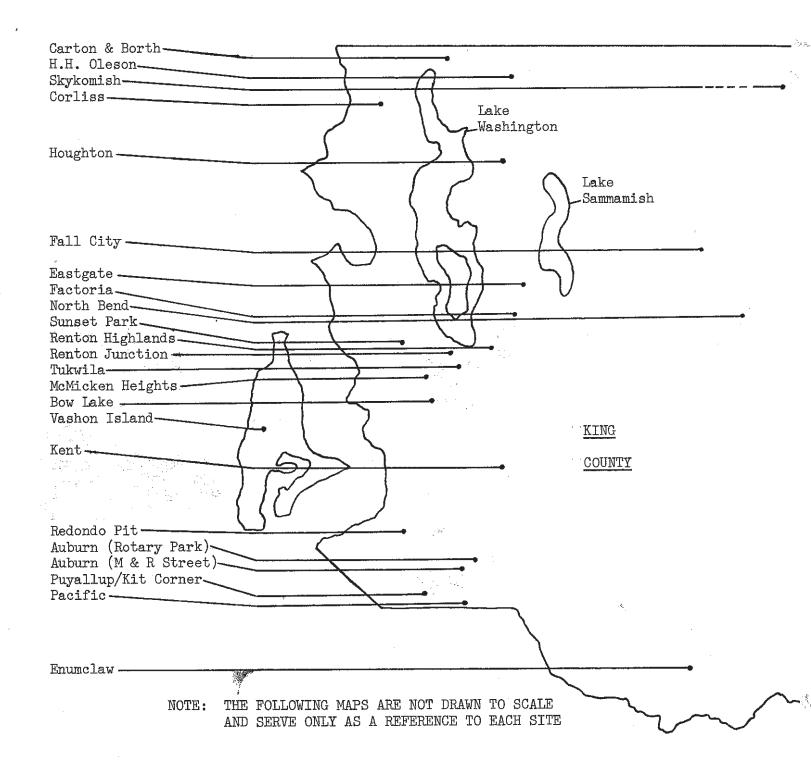


TABLE I

KING COUNTY ABANDONED LANDFILL LOCATIONS

Auburn (M & R Street) - Northwest of R Street and 25th Street S.E.

Auburn (Rotary Park) - Northwest of 27th Street S.E. and Alpine St. S.E.

Bow Lake - Northeast of S. 188th St. and Interstate-5

Carton & Borth - Northwest of 71st Ave. N.E. and N.E. 186th Street

Corliss - Northeast of Corliss Ave. N. and N. 163rd Street

Eastgate - Northeast of 156th Ave. S.E. and I-90, north end of Bellevue Airstrip

Enumclaw - Southeast of 284th Ave. S.E. and S.E. 448th Street

Factoria Pit - Northwest of 135th Ave S.E. and S.E. 40th Street

Fall City - Old Dump Rd. at first road bend

H.H. Oleson - Northeast of N.E. 172nd St. and 152 Pl. N.E.

Houghton - Northwest of N.E. 60th St. and 120th Ave. N.E.

Kent - Northeast of Maple St. and Tilden Avenue

McMicken Heights - Between the S. 175th St. dead-end and Interstate-5

North Bend - West of the Middle Fork Road, about 1 mile north of the Y turn from Edgewick Road

Pacific - South of 3rd Ave. S.E. at the White River

Puyallup/Kit Corner - Northeast of S. 360th St. and Interstate-5

Redondo Pit - Southwest of S. Dash Point Road and Pacific Highway South

Renton Highlands - Southwest of N.E. 3rd St. and Jefferson Avenue N.E.

Renton Junction - Between Monster Road and the Green River

Skykomish - North of the Stephen's Pass Highway just east of Skykomish

Sunset Park - Southwest of 18th Ave. S. and S. 136th Street

Tukwila - South of 62nd Ave. S. and S. 151st Street

Vashon Island - West of 130th Ave. S.W. across from the present landfill

METHODS AND MATERIALS

OBJECTIVE

It is the objective of this study to identify obvious and potential problems at the King County abandoned landfills that could adversely impact the public's health and safety.

General test parameters were selected for this preliminary assessment to:

- Identify sites with problematic methane off-gasing;
- Identify site areas that may have been used for hazardous waste dumping;
 - 3. Locate potential or obvious leachate seepage problems.

TEST PARAMETERS

The test parameters selected to identify problematic off-gasing include measurements of methane gas and non-specific trace gases. These are described as follows:

- 1. Methane Gas a final by-product of anaerobic organic decomposition, methane is explosive when concentrated in the range of 4% to 18% per volume of atmoshere.
- 2. Trace Gas for the purposes of this study, trace gas includes any organic or inorganic gas with an ionization potential of ≤ 10.2 eV. Methane, with an ionization potential of 12.98 eV would not be detected. Sites exhibiting trace gas levels greater than ambient air levels are suspect as potential areas of hazardous waste dumping.

The parameters selected to identify surface leachate seepage include pH, temperature, dissolved oxygen, electrolytic conductivity and turbidity. These are described as follows:

- 1. pH typical leachate values have been reported to range from 3.7 to 8.5 with a median value of 5.8. Typical surface water values encountered in King County range between 6.5 to 8.5.
- 2. Temperature fluctuates seasonally (electrolytic conductivity increases with temperature at a rate of approxiately 2% per degree C).
- 3. Dissolved Oxygen dependent upon the physical, chemical, and biological activities in the water. Water quality tends to decrease with low dissolved oxygen values.
- 4. Electrolytic Conductivity the ability for a water sample to carry an electric current. It is dependent upon the total concentration of the ionized substances dissolved in the water and temperature. Typical sanitary landfill leachate values have been reported to range from 0.1 to 1.2 milliohm/cm. Surface waters in King County typically exhibit conductivity values between 0.05 to 0.15 milliohm/cm.
- 5. Turbidity measurement of suspended solids in a water sample reported in ppm. To approximate NTU's the following is given: $ppm \times 2.1 = NTU's$.

SAMPLING INSTRUMENTATION

Gas measurements were made from one-half inch diameter bore holes sunk three feet into the ground at locations throughout each landfill area. Each hole was capped for a minimum of twenty minutes prior to testing to allow landfill gases to reach a state of equilibrium within the test holes. Methane gas and trace gas levels were monitored using the following instrumentation:

1. Methane Gas - Gas-Pointer Combustible Gas indicator, a product of Bacharach Instrument Company of Pittsburgh, Pennsylvania and Mountain View, California, a division of AMBAC Industries, Inc.

Sample gas was collected from the test hole through a suction tube, processed directly by the meter, and reported as percentage methane per volume of atmosphere.

2. Trace Gas - Model PI 101 Photoionization Analyzer, a product of HNU Systems, Inc. of Newton, Massachusetts.

Sample gas was collected from the test hole through a suction tube designed to draw gas at a flow rate of three-hundred to seven-hundred centimeters per minute past a photoionization sensor. The sample is directly processed by the instrument, recording trace organic and inorganic gases with ionization potentials of less than 10.2eV as parts per million.

Molecules with higher ionization potentials such as those forming the major components of air, would go undetected. Methane, with an ionization potential of 12.98 eV, would not register as part of the trace gases.

Water samples were tested in the field using a Model U-7 Water Quality Checker, a product of Horiba Instruments, Inc. of Irvine, California. The instrument was field calibrated using a standarized solution prior to testing, at a frequency of one per hour of operation. Field water samples were then placed into the unit's water test chamber for direct analysis by the sensor probes and read-out on digital display. The sensor probes were thoroughly rinsed with distilled water before and after each water testing.

HISTORICAL REVIEW METHODS

Since field tests give information about a site at only one point in time, it was appropriate to gain historical information on the twenty-three abandoned disposal sites. This was to provide a broader perspective of what might potentially be the site's characteristics. Specifically data gathering focused on past and present uses of the site in question; any engineering information about the site, including topography and soil studies; information pertaining to waste disposal practices at the site; and any information on known or suspected problems. This effort relied on available secondary data sources. Official records were few. However, files on some of the sites available from the King County Solid Waste Division contained many valuable references as did those from the King County Parks Department for sites subsequently developed for County parks. Anecdotal recollections, historical documents and maps, newspaper clippings, environmental impact statements and specialized studies done for development purposes, or citizen's advisory committee reports formed the bulk of the written report.

CHAPTER III RESULTS

HOLLING HE	Page
AUBURN (M & R STREET SITE)	116
AUBURN (ROTARY PARK SITE)	20
Bow Lake	23
CARTON & BORTH	29
CORLISS LANDFILL	34
EASTGATE ABANDONED LANDFILL	49
Enumclaw	47
FACTORIA PIT (SUNSET RAVINE PARK)	52
FALL CITY ABANDONED LANDFILL	56
H.H. OLESON	61
Houghton	66
KENT ABANDONED LANDFILL (MILL CREEK CANYON PARK)	72
McMicken Heights	78
North Bend	83
PACIFIC CITY	86
Puyallup/Kit Corner	91
REDONDO PIT	98
RENTON HIGHLANDS	103
RENTON JUNCTION (MONSTER ROAD)	108
Skykomish	114
Sunset Park	120
Tukwila	126
Vashon Island	130
OTHER SITES	133

AUBURN (M & R STREET SITE)

During the time period from the early 1930's through the 1940's King County operated a landfill in the area bordered approximately by 21st and 25th Northeast and M & R Streets in Auburn. The exact closure date is unknown, but since the A.L. Wolfe Addition was platted and built upon by the mid-1950's, it can be assumed that the site closed no later than the early 1950's.

PAST AND PRESENT USE

The area in question was originally operated by King County and later annexed by the City of Auburn. The site was at one time a gravel pit which was subsequently reclaimed by the garbage fill. After the closure, the area was platted as the A.L. Wolfe Addition and is now a large single-family residential area.

SITE/ENGINEERING INFORMATION

Since closure of the landfill and the subsequent housing development there have been numerous settling problems. Particularly along 21st, 22nd and 23rd Streets inadequate compaction and settling prior to on-site development has played havoc with housing stability, water lines, streets and sidewalks. At one point, F.H.A. became involved and required that basements be dug under existing houses to increase stability of the structures.

The area has been described as including some of the old river bottom of the White River and as a relatively shallow fill area.

WASTE DISPOSAL PRACTICES

Generally the landfill is remembered as a place where one could bury and get rid of anything. However, Auburn has been primarily a bedroom community in a rural

setting so that at this time there is no evidence to support anything but residential garbage and rubbish as being deposited on site. The one exception is the apparent existence of a toothbrush factory in the area and the occasional finding of toothbrushes when digging post holes in residential back yards!

The fill was estimated to have been six to eight feet in depth.

SUSPECTED PROBLEMS

With the exception of the long standing land compaction/settling problem in this area, there is no historical information to suggest need for any further investigation.

FIELD RESULTS

On November 26, 1984 four bore holes were tested within the M and R Street site for the presence of methane and non-specific trace gases. Methane was not detected above a trace level from any of the test holes. Trace gas concentrations were low with a range of 0 to +0.3 ppm relative to ambient air levels (Table II).

No surface water was observed on or about the site. Street settling was observed within the vicinity of the former fill.

FIGURE 2
AUBURN ABANDONED LANDFILL: M AND R STREET

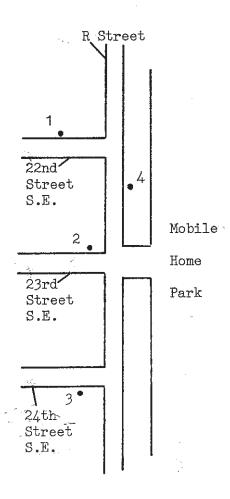


TABLE II

METHANE AND TRACE GAS CONCENTRATIONS

AUBURN ABANDONED LANDFILL: M and R STREETS

Site	Methane (%)	Trace Gas (ppm)*
1	Trace	0
2	Trace	0
3	0	.3
4	Trace	0

^{*}Reading represents change from ambient air level

AUBURN (ROTARY PARK SITE)

During the investigation of the main Auburn landfill, the existence of another landfill in the area was discovered. This was on the old Mead property, which is now part of the Auburn Rotary Park.

According to personal accounts this operated for several years until about 1965. In contrast to the main landfill this did not operate as a garbage dump but was primarily used for old cars, rocks, tires and general rubbish. Putrescible waste was not dumped here.

It is estimated that this site may have been thirty feet deep. Upon closure it was covered with one to two feet of top soil.

Since this site took everything except garbage it may be a site of long term concern, if "everything" included crankcase oil, pesticide cans and/or other potentially hazardous waste. However, no files, documents, or personal accounts suggest any problems at the site.

FIELD RESULTS

Twelve test holes were examined for methane and trace gas concentrations at the Rotary Park site on November 5, 1984 (Table III).

Methane gas levels were not observed exceeding trace concentrations throughout the site. Non-specific trace gas levels were observed ranging between 0 to +5.2 ppm relative to ambient air concentrations. Of note, the highest trace gas levels of +2.8, +1.9, and +5.2 ppm were located within the southeast corner of the park.

No waste was observed on or around the former landfill site, though a relatively high perched water table was noted within the southeast park quadrant.

FIGURE 3
AUBURN ABANDONED LANDFILL: ROTARY PARK

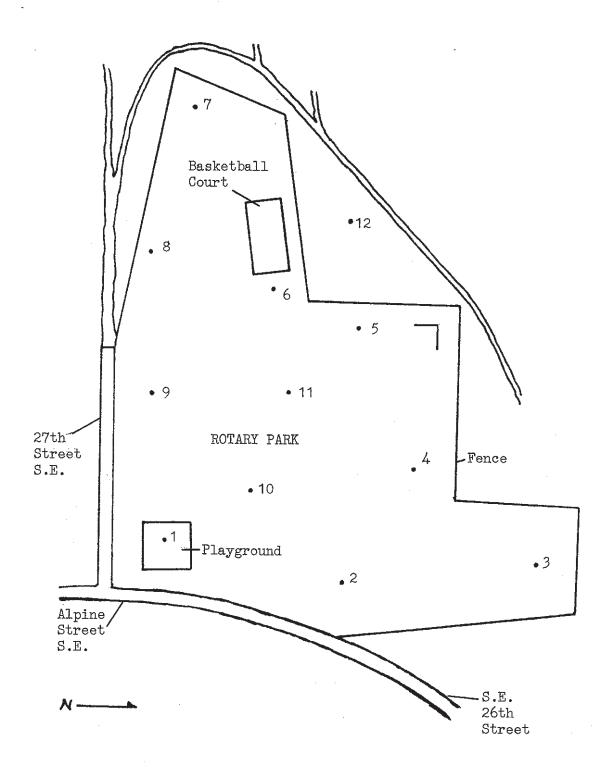


TABLE III

METHANE AND TRACE GAS CONCENTRATIONS

AUBURN ABANDONED LANDFILL: ROTARY PARK

Site	Methane (%)	Trace Gas (ppm)*
1	Trace	2.8
2	0	.5
3	Trace	1.9
4	0	.5
5	Trace	.7
6	Trace	.2
7	Trace	.2
8	Trace	0
9	Trace	.4
10	0	5.2
11	Trace	0
12	0	0

^{*}Reading represents change from ambient air level

BOW LAKE

Located in Section 35T-23N-R4E, the old Bow Lake landfill originally consisted of approximately 14 acres bordered on the west by Military Road and extending from slightly north of South 186th to the Orillia Exit road. It is typically identified as being at 188th and Military, the site of the old access road.

PAST AND PRESENT USE

The landfill was located in a predominantly rural area at the time of its operation. Old maps show a "dump" at the northern portion of the landfill site as early as in 1943. In the late 1950's the construction of Interstate 5 necessitated closure of the site as the freeway ran diagonally through this tract. In 1961 a transfer station was constructed on the site. This station became outdated and was redtagged by the Department of Labor and Industries in 1970. The "new" and existing transfer station was opened for a cost of \$1.7 million in the fall of 1978. At that time it was serving an estimated population of 125,000 and transferring at least 60,000 tons of refuse annually to the Cedar Hills landfill site. It remains today the major transfer point for South King County.

SITE/ENGINEERING INFORMATION

There is a distinct slope of the old landfill site to the southeast corner.

Early contour maps indicate a two hundred foot difference in elevation between the northwestern portion and the southeastern portion of the site.

Today the transfer site is bordered by freeway and freeway access roads on the west and south and private property on the north and east.

WASTE DISPOSAL PRACTICES

Records indicate that the Bow Lake landfill was the largest landfill in the County during the 1950's when it received $160,000 \text{ yd}^3$ of refuse annually. It served a large geographic area of South King County, running from the Seattle City boundary to Puget Sound, bordering Renton and Kent.

The toe of the fill appears to have moved from northwest to southeast as filling occurred through the years.

Interestingly a small incinerator was installed at the Bow Lake site on an experimental basis in 1955. Unfortunately this proved to be too smokey and incapable of handling the large items of refuse or garbage. In addition, it burned at too low a temperature to ensure complete combustion, so the incinerator was closed.

SUSPECTED PROBLEMS

Comments from private owners, the City of Tukwila and personal observation make it clear that the major problem at this abandoned landfill is uncontrolled leachate. It has apparently been a chronic problem for years.

A secondary concern is what might be <u>in</u> the site. Since the dump operated through the years when there was less effort and no regulation to separate waste types and since it served an area where there were commercial and industrial developments such as Sea-Tac and Boeing, there is the possibility that some potentially hazardous material could be on site. No records have been found to document this supposition. The site is listed on the EPA Region 10 ERRIS list.

FIELD RESULTS

The former Bow Lake Landfill was tested for methane and non-specific trace gas concentrations on October 22, 1984 (Table IV).

Of the seventeen bore holes tested for methane, four were observed within or above the explosive range of methane gas (4% to 18%). Peak methane levels of 30% and 35% were found at the northend of the site.

Trace gas levels ranged between 0 to +0.8 ppm relative to ambient air concentrations. Of note, four of the five positive trace gas readings corresponded to test holes which exhibited the highest levels of methane gas.

Leachate with an oily sheen and characteristic discoloration was observed seeping from the site from the southeast quadrant. Water parameters did not indicate it to be highly concentrated. These data are found in Table V.

Test parameters of a water sample retrieved from the transfer station storm drain did not indicate leachate contamination.

FIGURE 4
BOW LAKE ABANDONED LANDFILL

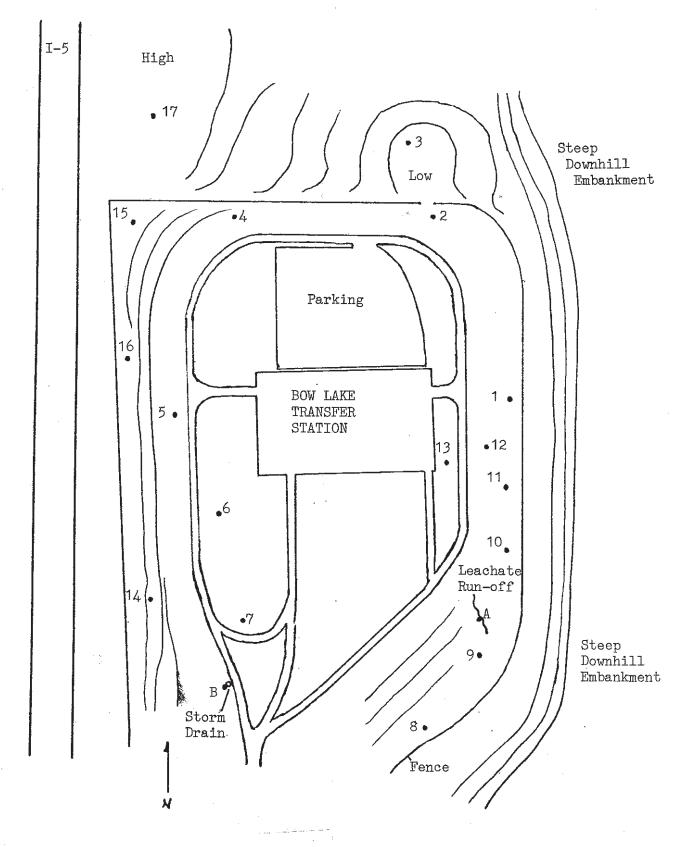


TABLE IV

METHANE AND TRACE GAS CONCENTRATIONS

BOW LAKE ABANDONED LANDFILL

Site	Methane (%)	Trace Gas (ppm)*
1	0	0
2	35	0.3
3	30	0.7
4	0	0.1
5	0	0
6	5	0.8
7	5	0.7
8	0 (Meter Box)	
9	0	0
10	0	0
11	0	0
12	0	0
13	2.4	0
14	0	0
15	0	0
16	0	0
17	0	

^{*}Reading represents change from ambient air level

TABLE V
SURFACE WATER PARAMETERS
BOW L'AKE ABANDONED LANDFILL

	Site A (1)	Site B (2)
рН	6.6	6.6
Temperature	13.1	15.5
Dissolved	1.6	4.2
Electrolytic	0.5	0.5
Turbidity	110	2

⁽¹⁾ Flowing leachate stream

⁽²⁾ Standing water in storm drain

CARTON & BORTH

The Carton and Borth former landfill is situated between 68th Avenue Northeast and 71st Avenue Northeast. The southwesterly quarter of the site is currently occupied by the WesMar Industrial Park. The northwesterly quarter is presently under new building construction.

This site was formally known as the Loveless and Dillon Site. A Seattle <u>Times</u> article dated August 3, 1971 reported that the County Council approved an unclassified use permit for a sanitary landfill by Loveless and Dillon, Inc. This was described as a 12.81 acre landfill on the east side of 68th Avenue Northeast, Kenmore and about three hundred thirty feet north of the Tolt River pipeline.

Since it was a private operation, the County regulated and inspected the site. Health Department officials remember getting complaints about the site regarding stench, which may have been an indication of inadequate cover operations. In addition a stream, which is currently diverted around the perimeter of the property, once ran through the property. The Health Department had responded to complaints regarding the off-colored appearance and odor associated with this stream.

The site was primarily used for rubbish such as wood and stumps, demolition materials, and oil from roads. A minimal amount of garbage was apparently received. It operated for only a short time for the purpose of reclaiming the site for the commercial development that currently exists.

FIELD RESULTS

On October 30, 1984 nine test bore holes were placed in the easterly half of the Carton and Borth site for methane and trace gas monitoring. These data appear in Table VI. Access for testing of the westerly half of the fill was denied by its current property owners.

Methane gas was observed in low concentrations throughout the site with a peak reading of 5%. All trace gas observations were neutral relative to ambient air concentrations.

One water sample was retrieved from the stream passing around the north to east perimeter of the fill. These data are presented in Table VII. Leachate contamination was not indicated by the test parameters.

FIGURE 5
CARTON & BORTH ABANDONED LANDFILL

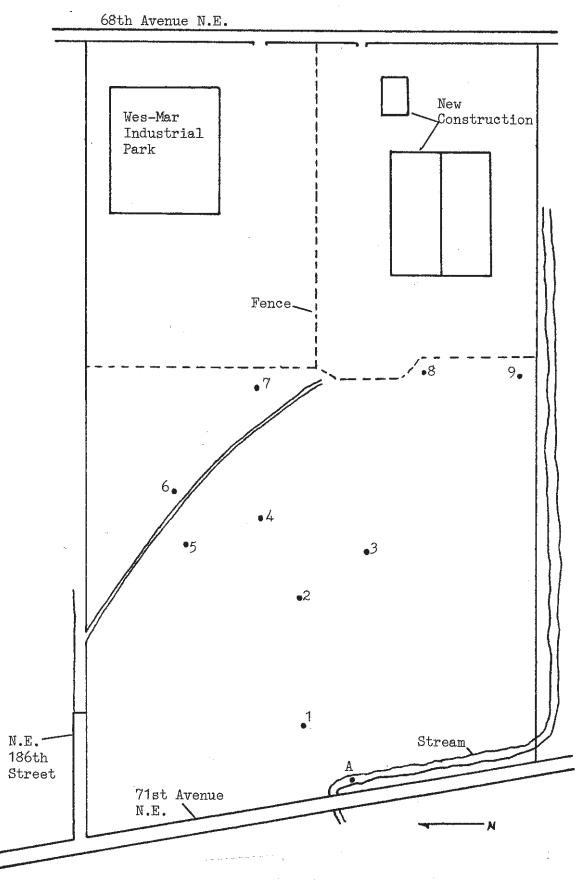


TABLE VI

METHANE AND TRACE GAS CONCENTRATIONS

CARTON & BORTH ABANDONED LANDFILL

Site	Methane (%)	Trace Gas (ppm)*
1	Trace	0.1
2	Trace	0
3	Trace	0
4	Trace	0
5	. 0	0
6	Trace	0
7	5	0
8	1.2	0
9	0	0
	•	

^{*}Reading represents change from ambient air level

TABLE VII SURFACE WATER PARAMETERS CARTON & BORTH ABANDONED LANDFILL

	Site A (1)
рН	6.1
Temperature	9.7
Dissolved	7.8
Electrolytic	0.2
Turbidity	1

⁽¹⁾ Flowing water from creek

CORLISS LANDFILL

The north end of King County was serviced for several years by a landfill located in the vicinity of 1st Avenue Northeast and Corliss Avenue between approximately Northeast 163rd and Northeast 165th. It operated from the 1940's until closed by the construction of Interstate 5 in June 1959. The landfill site today is divided midway by a fence extending east to west. The northerly half is the site of the Northeast Transfer Station. The southerly half is an undeveloped tract.

PAST AND PRESENT USE

Prior to 1940 the site was essentially rural, undeveloped property. Between about 1946 and 1959 King County used the area for a sanitary landfill.

Construction of Interstate 5 necessitated excavation of the earliest part of the waste fill on the southeasterly portion of the site. The freeway now runs over that section. The north transfer station was constructed on the more recently used portion of the landfill. The remaining area consists of the undeveloped McCormick Park at Northeast 165th and Corliss, which the King County Parks Department acquired in 1967. This tract is also currently being considered by Metro as a potential site for their North Operating Base.

SITE/ENGINEERING INFORMATION

Data derived from Washington State Department of Transportation and King County explorations reported in a 1984 EIS prepared by Metro describe the site as an area providing surface drainage for Ronald Bog and perimeter areas north of this site. A large peat bog is found adjacent to the landfill site and was mined about the same time the solid waste operation existed. In addition, peat is found under portions of the landfill site.

Soil studies undertaken in 1977 in this area showed that garbage and rubbish extended from a depth of two feet to fourteen feet. Other soil explorations on site note that the maximum fill recorded is thirty seven feet. Motor oil smell was encountered at three feet in one soil boring reported for the site.

The site soils are characterized as fill, peat, loose sand and soft silt deposit, glacial advance sand and glacial till. The ground water is encountered anywhere from four to seven feet in the recorded borings. Groundwater flows generally to the southwest across the site. An early map of the landfill area characterizes the southeastern portion as a "swamp" and shows the toe of the fill adjacent to a creek - possibly Thornton Creek.

WASTE DISPOSAL PRACTICES

The sanitary fill at Northeast 165th and 1st Avenue Northeast handled about 60,000 cubic yards of refuse annually and was the primary north end disposal site. Over 900 cars and trucks were reported to come to the dump on weekends. The nearest County fill after the closure in June 1959 was the site at Houghton.

The earliest portion of the fill at 1st Northeast was lost to the freeway. Included in this portion, in 1947, at the northeastern edge, was an authorized septic tank dump. Plans were underway by 1960 to use part of the remaining landfill site for a transfer station as it exists today.

SUSPECTED PROBLEMS

Since closure of the landfill site, the land has remained undeveloped except for construction of the transfer station.

Concerns about differential settlement, ground water and the natural drainage in the area have made planners cautious regarding development.

In addition, since the site was one of two that handled the majority of the waste produced in King County for several years, it is possible hazardous materials could have been placed on site. The site is mentioned on the EPA ERRIS listing. However, the 1984 Metro Draft Supplemental Environmental Impact Statement for the North Operating Base notes that, "Municipal waste material samples were collected and analyzed from four borings drilled in the filled area. The chemical analyses indicate that the waste material would not likely be considered hazardous or dangerous."

FIELD RESULTS

The Corliss Abandoned Landfill is divided midway by a fence east to west. Testing of the south sector was conducted on October 11, 1984 for methane and non-specific trace gas (Table VIII). Gas sampling of the north sector and all surface water tests were completed on October 12, 1984.

Methane gas was observed at levels within or above the explosive range for methane gas (4% to 18%) from five of twenty test holes located without pattern throughout the site.

Non-specific trace gas was observed near that of ambient air levels with the exception of test hole #11 (0.8 ppm), #14 (3.8 ppm), #18 (9.2 ppm) and #19 (0.8 ppm).

A stream which runs along the site's westerly perimeter was sampled from locations upstream and downstream from the former fill. Test parameters did not indicate any obvious leachate impact upon the stream. These data appear in Table IX.

FIGURE 6
CORLISS ABANDONED LANDFILL

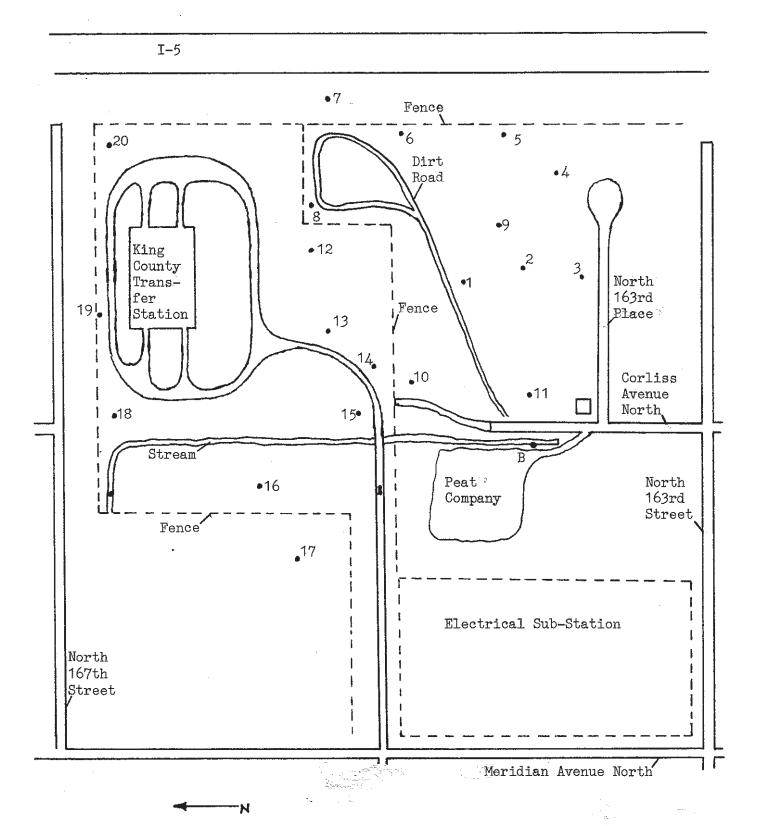


TABLE VIII
METHANE AND TRACE GAS CONCENTRATIONS
CORLISS ABANDONED LANDFILL

Site	Methane (%)	Trace Gas (ppm)*
1	3.2	0.1
2	Trace	0.2
3	0	0.2
4	0	0.1
5	Trace	0.1
6	30	0.1
7	10	0.2
8	1.2	0.2
9	Trace	0.2
10	0.8	0.2
11	27	0.8
12	Trace	0.1
13	Trace	0
14	42	3.8
15	PO 500 VIII	0.1
16	5	0
17	0	0.1
18	3	9.2
19	Trace	0.8
20	Trace	0

^{*} Reading represents change from ambient air level

TABLE IX SURFACE WATER PARAMETERS CORLISS ABANDONED LANDFILL

	(1) Site A	Site B (2)
рН	6.5	6.6
Temperature	14.9	14.5
Dissolved	6.5	6.0
Electrolytic	0.1	0.1
Turbidity	4	3

⁽¹⁾ Stream water upstream from landfill(2) Stream water downstream from landfill

EASTGATE ABANDONED LANDFILL

The north end of the Bellevue Airfield was the site of a former landfill which was closed and covered in 1964. Solid Waste Division files show the landfill operation in this area in 1953 when it was referred to as the Factoria Garbage Dump. It apparently began in about 1951.

Little historical documentation has been found regarding the specific operations of this site. Personal anecdotes reveal that at one time a gravel pit was adjacent to the site.

Very few operational problems were remembered. It was estimated to have been filled to a depth of six to twenty-eight feet. The site was a burn dump and consequently became a problem for the airfield due to smoke reduced visibility.

The general area was purchased by Cabot, Cabat & Forbes and subsequently by Boeing (the current owners) and the Bellevue School District. The site is part of the I-90/Bellevue Business Park.

SUSPECTED PROBLEMS

The site is currently surrounded by controversy. Residents in the Phantom Lake Area north of the old landfill have become concerned about the development of the site and its impact on the lake.

Drainage going through the garbage has caused leachate to be a problem. This has necessitated costly interventions by private owners, including an ongoing methane monitoring program conducted by Boeing. Entranco Engineers recently conducted a study of the landfill leachate to evaluate "whether landfill leachate constitutes significant and dangerous or damaging contamination to Phantom Lake."

This was done by assessing priority pollutant data from two sampling stations

located on the I-90 Bellevue Business Park Property. Given the limitations of small sample size, the results included: 1) "Landfill leachate contains certain heavy metal and organic priority pollutants at detectable levels." 2) "Of the organic pollutants measured in the landfill leachate, the insecticide chlordane was the only compound found at concentrations higher than expected." The report cautions that the leachate could have adverse impacts on ground water and that concentrations of the leachate pollutants could vary with higher concentrations possible.

FIELD RESULTS

Nineteen bore holes were tested for methane and non-specific trace gases at the Eastgate Abandoned Landfill on November 20-21, 1984. These data appear in Table X. Methane was observed ranging from 0% to 10% on the site. Only two test holes registered methane in the explosive range (4% to 18%).

Non-specific trace gas levels were low relative to ambient air levels with only one test hole exhibiting a level of +0.6 ppm.

Of note, the Boeing Company, current owners of most of the site property, has installed several off-site methane test wells dug to approximately 30 feet to measure off-site gas migration. It has been their experience to observe dramatic fluctuations in methane levels with changing barometric pressure. High pressure systems have been associated with low methane levels, and low pressure systems with high levels.

Our observations of this phenomenon appear in Table XI. Two test wells sampled on different days under different barometric conditions yielded dramatically different results. Methane levels were observed fluctuating from trace to 40% from test well A, and 4% to 70% from test well B.

Leachate has historically been a problem with this site. At one time a leachate collection system was installed at the north end of the fill which channeled into a stream and eventually Phantom Lake.

Samples obtained from the settling pond, the leachate recovery stream by the fill and by the lake did not indicate concentrated leachate contamination. Surface water from a run-off stream located directly on the landfill exhibited parameters of good water quality. These data are presented in Table XII.

FIGURE 7

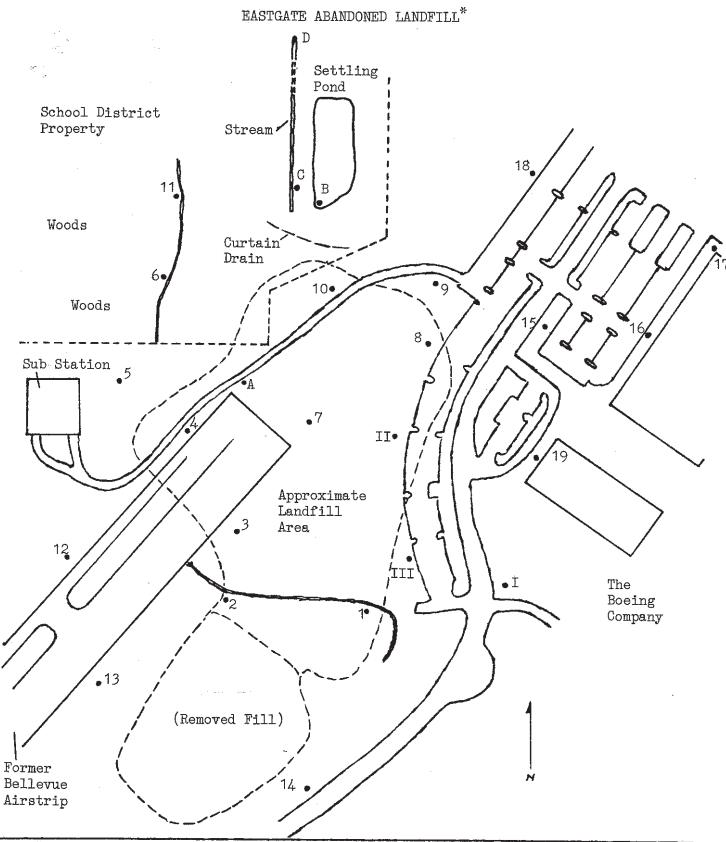


TABLE X
METHANE AND TRACE GAS CONCENTRATIONS
EASTGATE ABANDONED LANDFILL

Site	Methane (%)	Trace Gas (ppm)*	
1	0.2	0	
2	0	0	
3	0.2	0	
4	10	0.1	
5	0.2	0	
6	0.2	0	
7	Trace	0	
8	Trace	0	
9	5	0	
10	Trace	0.6	
11	Trace	0	
12	Trace	0	
13	Trace	0	
14	0	0	
15	0	0	
16	Water Table	Water Table	
17	0.2	0	
18	0.2	0	
19	0.4	0	

^{*} Reading represents change from ambient air level

TABLE XI

METHANE CONCENTRATIONS IN TEST WELLS

EASTGATE ABANDONED LANDFILL

<u>Well</u>	Sampling	Methane (%)*
I .	11-21-84	Trace
d e	11-29-84	40
II	11-20-84	4
: -	11-29-84	70
III	11-29-84	24

^{*} Variations due to barometric pressure differences

TABLE XII SURFACE WATER PARAMETERS EASTGATE ABANDONED LANDFILL

	Site A(1)	Site B(2)	Site C(3)	Site D(4)
рН	5.7	5.9	6.2	6.9
Temperature	8.4	9.7	9.4	9.4
Dissolved Oxygen ppm	7.5	7.8	2.4	9.6
Electrolytic Conductivity m~v/cm	.05	.15	.2	.05
Turbidity ppm	2	2	160	1

Drainage ditch on landfill surface Settling pond Drainage stream from leachate collection system Drainage stream from leachate collection system

ENUMCLAW

The current sanitary landfill on Southeast 440th in Enumclaw, operating since 1939, is generally referenced as the "old" landfill. The Sonneson family deeded this area to the City of Enumclaw on January 5, 1939 for the purposes of a garbage dump. However, there was an earlier landfill which was abandoned in 1939 when the Sonneson property became available.

The landfill was located on Roosevelt Avenue to the east of the Farman's Pickle Factory adjacent to the land currently occupied by Pete's Pool at the King County Park. The County leases land to the Enumclaw Golf Course on the site of the old landfill and early maps of the abandoned site show that the fairway parallels the approximate site of the old dump.

Since this dump has been closed nearly fifty years, preliminary data searching revealed no existing records on the site's specific operation.

However, it might be assumed that the old site served the area serviced by the current landfill. This would include the city of Enumclaw as well as some private residents of King County and commercial refuse haulers. The property in the area is primarily residential and Enumclaw itself is a predominantly rural town. It is unlikely that any hazardous material found its way to the landfill unless there could have been wastes from any wood manufacturing existing at that time. The current site is on the EPA ERRIS list, indicating a potential possibility of hazardous waste materials in that site. The earlier site of current concern was closed before the major chemical developments and subsequent waste generation which occurred during and after World War II. Thus, while the existing site may well contain hazardous materials, there is little to suggest the same for the old site.

FIELD RESULTS

On December 8, 1984 nine bore holes were tested at the Enumclaw site for methane gas concentrations. These data appear in Table XIII.

Methane gas levels were observed in low concentrations (trace to 0.2%) from test holes located along the east and west perimeter of the fill, indicating its boundaries.

Non-specific trace gas levels were not monitored at this site.

Surface water was not observed on the site. However, a nearby stream was sampled and exhibited parameters of good water quality (Table XIV).

FIGURE 8
ENUMCLAW ABANDONED LANDFILL

S.E. 448th Street 3 •1 Enumclaw Golf Course Boise Creek Office

TABLE XIII

METHANE GAS CONCENTRATIONS
ENUMCLAW ABANDONED LANDFILL

<u>Site</u>	Methane (%)
1	0
2	Trace
3	0.2
4	0.2
5	0.2
6	0.2
7	Trace
8	0
9	. 0
	:

TABLE XIV

SURFACE WATER PARAMETERS ENUMCLAW ABANDONED LANDFILL

рН	٠	٠	•	•	•	•	•	•		•		5.8
Temperature O _C	•	•	•	•	•		•	•	•	•	•	6.5
Dissolved . Oxygen ppm	•	•	•	•	•	•	•	•	•	•	•	10.2
Electrolytic Conductivity m \(\forall / \text{cm}\)	•	•	٠		•	•	•	•	•	•	•	0.2
Turbidity .	•											42

FACTORIA PIT (SUNSET RAVINE PARK)

South of I-90 on 136th Avenue Southeast is the site of another abandoned landfill in the City of Bellevue. It was closed in about 1951 when operations were moved to the northend of the Bellevue Airfield.

Official information on this site is sketchy. It was apparently the original landfill site in the general area and operated as a burn dump. The site remains undeveloped. It is listed on the EPA ERRIS report.

FIELD RESULTS

On December 1, 1984 eleven bore holes were tested for methane and non-specific trace gases at the Factoria abandoned landfill site. These data appear in Table XV.

Methane was observed ranging from 0% to 0.2% indicating the stability of the site. Trace gas levels were all observed undifferentiated from ambient air levels.

Upstream and downstream water samples obtained from a stream, passing adjacent to the site indicated no change in water quality (Table XVI).

FIGURE 9
FACTORIA ABANDONED LANDFILL

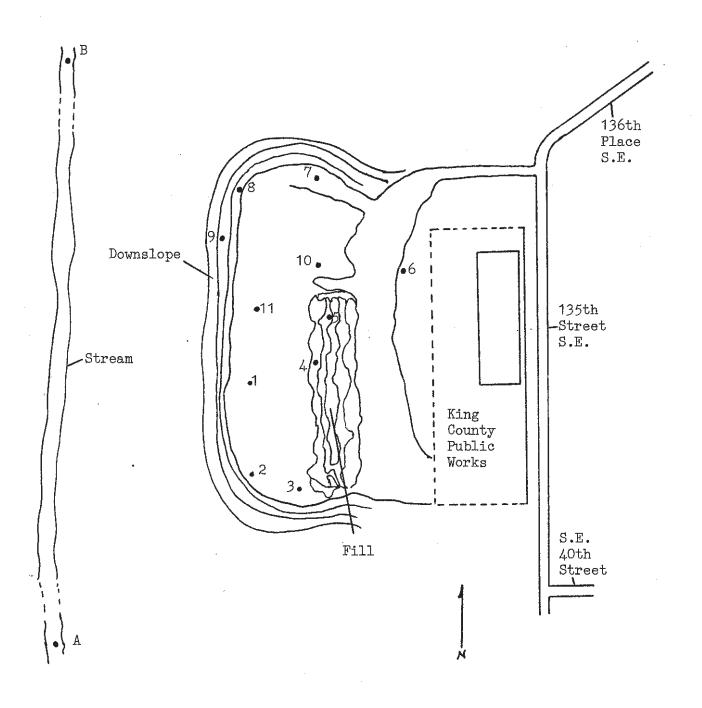


TABLE XV

METHANE AND TRACE GAS CONCENTRATIONS
FACTORIA PIT ABANDONED LANDFILL

Site	Methane (%)	Trace Gas (ppm)*
1	Trace	0
2	Trace	0
3	Trace	0
4	0	0
5	Trace	0
6	0.2	0
7	Trace	0
8	Trace	0
9.	0.2	0
10	0	0
11	0	0

^{*} Reading represents change from ambient air level

TABLE XVI SURFACE WATER PARAMETERS FACTORIA PIT ABANDONED LANDFILL

	Site A (1)	Site B (2)
рН	5.9	6.0
Temperature	9.4	9.0
Dissolved	9.7	10.1
Electrolytic	0.15	0.15
Turbidity ppm	3	9

⁽¹⁾ Upstream from landfill(2) Downstream from landfill

FALL CITY ABANDONED LANDFILL

Located off the Fall City-Duvall Road is a turn to Southeast 39th Place, the "old Dump Road." Following this road about 0.3 mile beyond two gate posts on the road brings one to the site of a small former landfill. This operated several years and was closed sometime in the early 1960's.

The site was located on Weyerhauser property and privately operated. It was approximately three hundred feet deep and one hundred to one hundred fifty feet in width, located above Rutherford Slough on the top of a steep embankment. Vehicles pulled in and dumped material over the hillside. It was a burn dump.

It was predominantly a community dump and being in a very rural part of King County it is doubtful hazardous materials are on site. The one reasonable exception is the potential presence of pesticides in a rural area. In addition, it is suspected that septic tank pumpers may have dumped on site. However, the practice of burning on site has eliminated any pollutants of this nature as a chronic problem.

When it was closed, Weyerhauser was to clean the area up and close it off.

Today there is no visual evidence that a landfill once operated here. It is now overgrown with brush and tall grasses and remains vacant.

FIELD RESULTS

On-site gas testing at the Fall City site was run on October 20, 1984 for methane and non-specific trace gases. These data are presented in Table XVII. Methane was observed in only trace concentrations in each of the five test bore holes.

Trace gas levels were neutral to the ambient air with the exception of a low ± 0.2 ppm reading of hole $\pm 3.$

No off-site leaching problems were directly observed. Leachate contamination was not indicated by the test parameters of a slough sample, though an oily sheen was observed on the slough surface (Table XVIII).

Recent illegal dumping of household refuse was observed on the site in small quantities.

FIGURE 10
FALL CITY ABANDONED LANDFILL

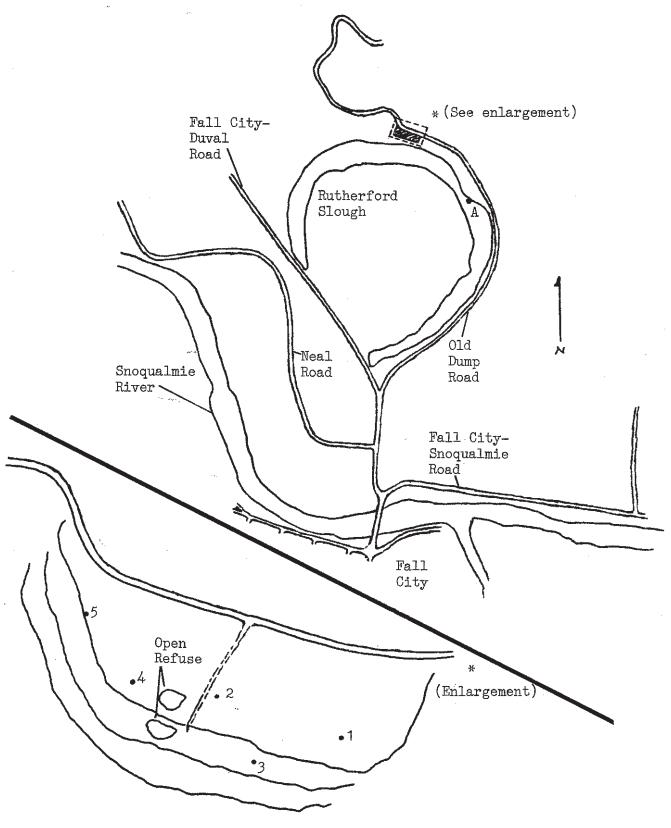


TABLE XVII

METHANE AND TRACE GAS CONCENTRATIONS
FALL CITY ABANDONED LANDFILL

<u>Site</u>	Methane (%)	Trace Gas (ppm)*
1	Trace	0
2	Trace	0
3	Trace	0.2
4	Trace	0
5	Trace	0

^{*} Reading represents change from ambient air level

TABLE XVIII

SURFACE WATER PARAMETERS FALL CITY ABANDONED LANDFILL

рН		•			•		•	. •		•				•	6.3
Temperature O _C		•	•	•	•	•	•	•	•	•	•	•	•		7.3
Dissolved . Oxygen ppm	1		•	•	•	•		•	•	•	•	•		•	8.4
Electrolytic Conductivity m v /cm		•		•		•		•			•	•	•	•	0.1
Turbidity .		•			•	•	٠		•	•		•	•	•	8

The H.H. Oleson landfill operated from about 1972 until about 1979 as a means of reclaiming an existing abandoned excavation site. It consisted of about 10.8 acres according to the legal description.

PAST AND PRESENT USE

In the early 1970's H.H. Sand and Gravel Company requested the permit to operate a demolition landfill on the site of their former excavation pit. It was primarily a reclamation effort. A time extension for the operation was requested in 1977. At that time the site was described as slightly wooded on the last portion, with the excavation on the west portion currently being filled. Thus the property apparently was filled from east to west.

SITE/ENGINEERING INFORMATION

Maps of the site show that drainage was an anticipated problem. Settling basins were planned along 152nd Avenue Northeast so that drainage swales could have water channeled off property to drains and/or ditch on 152nd Avenue Northeast.

The completed fill was anticipated to change the topography of the site considerably. The final fill slopes gradually to 152nd Northeast (Van Brocklin Road).

WASTE DISPOSAL PRACTICES

According to a S.E.P.A. Application at the time the landfill began, the operation was to follow the standard sanitary fill procedures, alternating layers of sanitary fill and soil. It was estimated that a thirty mile radius around the fill site could be anticipated to contribute waste.

It appears that the fill depth varied from ten to fifteen feet to seventy to seventy five feet in some of the deepest sections.

SUSPECTED PROBLEMS

The type of waste suspected at this site consists of demolition waste. The site is on the EPA ERRIS list.

In addition, it appears that ground and/or surface waters may be subject to leachate.

It was anticipated in the SEPA checklist prepared for this operation that the fill would produce compaction variability, causing some portions of the site to be unsuitable for construction.

FIELD RESULTS

On January 11, 1985 nine bore holes were tested at the former Oleson Landfill for methane and trace gas emissions. These data are presented in Table XIX.

Methane gas levels were not observed above trace levels throughout the site indicating minimal organic decomposition. Trace gas levels were at an equilibrium with ambient air levels.

A water sample was retrieved from a small gravel pit pond located immediately south of the fill. Evidence of leachate contamination was not indicated by the test data (Table XX).

FIGURE 11
OLESON ABANDONED LANDFILL

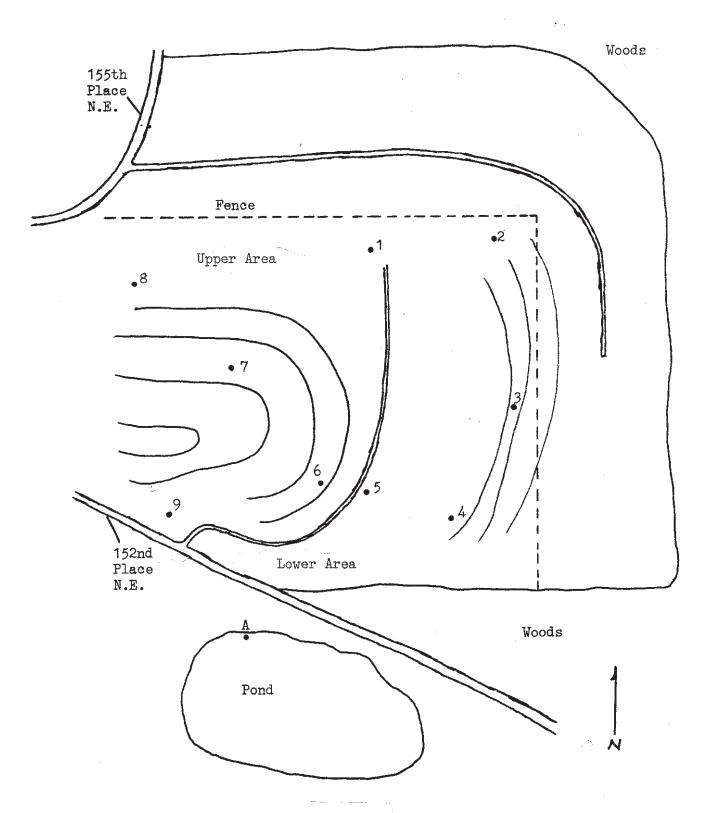


TABLE XIX

METHANE AND TRACE GAS CONCENTRATIONS

OLESON ABANDONED LANDFILL

		ንካ
Site	Methane (%)	Trace Gas (ppm) (1,2)
	0	0
2	Trace	0.1
3	Trace	0.2
4	Trace	. 0
5	.0	0
6	Trace	0
7	Trace	0
8	0	0
9	0	0

⁽¹⁾ Trace gas measurements made using an 11.2 eV HNU probe(2) Reading represents change from ambient air level

TABLE XX SURFACE WATER PARMETERS OLESON ABANDONED LNADFILL

•			<u>Site (A)</u> (1)
pH		 	5.0
Temperature °C	13.	 	1.8
Dissolved Oxygen (ppm)		 	13.9
Electrolytic Conductivity mu/cm		 • • •	0.35
Turbidity (ppm)		 • • •	34

⁽¹⁾ Gravel pit pond

HOUGHTON

The Houghton Transfer Station located near Bridle Trails State Park is the site of an old abandoned landfill. The landfill operated between approximately Northeast 60th Steet and Northeast 67th bounded by 117th Avenue Northeast and 120th Avenue Northeast. It was used from at least 1945 until closed in 1965.

PAST AND PRESENT USE

As a part of a predominantly rural area, the site was apparently undeveloped until used as a landfill operation beginning in the 1940's. It operated as a landfill until closed in about April 1965. By this time the site was noted to be leveled, harrowed, fertilized and seeded, with a portion of the site adjacent to the current transfer station still being used to dump hot ashes. This transfer station was in operation at the site by 1970 and that operation, plus open space, characterizes the site today.

SITE/ENGINEERING INFORMATION

The site is not perfectly level but is characterized by a gentle slope with the elevations in the general vicinity of the transfer station being greater than those along Northeast 67th. Fill depth varies on site from approximately two feet to over twenty feet. The area around 119th between Northeast 66th and Northeast 67th was characterized as swamp. High water tables and consequent saturated garbage could be anticipated throughout the site.

The final cover consisted of 8,000 cubic yards of top soil brought on site by the Thos. Scalzo Company in 1970. In 1972, due to heavy rains, an additional 2,000 cubic yards of impervious cover material were placed on site and compacted over the area to seal the garbage off from further moisture infiltration. It was replanted with new cover material characterized as having a high transpiration rate.

WASTE DISPOSAL PRACTICES

According to 1963 correspondence, operations at Houghton included garbage dumping, an "old dump area," a salvage operation and an excavation area being used by the State for fill material for highway work.

Burning was practiced in the area of the fill now occupied by the Transfer Station.

After the Corliss/1st Avenue Northeast landfill site was closed in 1959, the refuse was hauled to Houghton for disposal.

The oldest portion of the site, referred to in correspondence as "the old dump area" was east of 119th Northeast and south of Northeast 63rd Street. The site apparently filled from that location northerly toward the area of the existing transfer station.

SUSPECTED PROBLEMS

While Houghton was the receiving site for much King County waste for several years, the characteristics of the waste are not documented. The site is on the EPA ERRIS list.

Problems with impounding water in a low spot in the older section of the landfill have occurred as have leachate problems. Since part of the area has high ground water, it can be anticipated that leachate may be an ongoing problem, as it has been historically.

FIELD RESULTS

On October 8-9, 1984 twenty one bore holes were tested at the Houghton site for methane and non-specific gas. These data are presented in Table XXI.

Methane gas was observed within the explosive concentration range of 4% to 18% in eight of twenty one bore holes tested. Of these, seven were concentrated in the southwest section of the fill.

Trace gas levels ranged between -0.1 to +1.6 ppm relative to ambient air concentrations. Positive trace gas readings generally corresponded with areas of higher methane levels.

Site examination for leachate production, particularly along the west perimeter, were made on October 9, November 8, and December 1, 1984. No leachate or surface water was observed leaving the site.

Water was sampled from a storm run-off stream within the vicinity of the fill's southwest corner on November 8, 1984. Leachate contamination was not indicated. On January 15, 1985 a second water sample was taken from a dug well located along the site's west perimeter. The sample had a strong stagnation off-odor but did not exhibit test parameters indicating heavy leachate contamination. It was noted by the well's owner that the well had a history of containing iron-orange colored water during summer months which would kill grass if used for lawn watering. A third sample was retrieved that day from a swamp located at the west perimeter of the former fill. The influence of leachate was not observed.

FIGURE 12

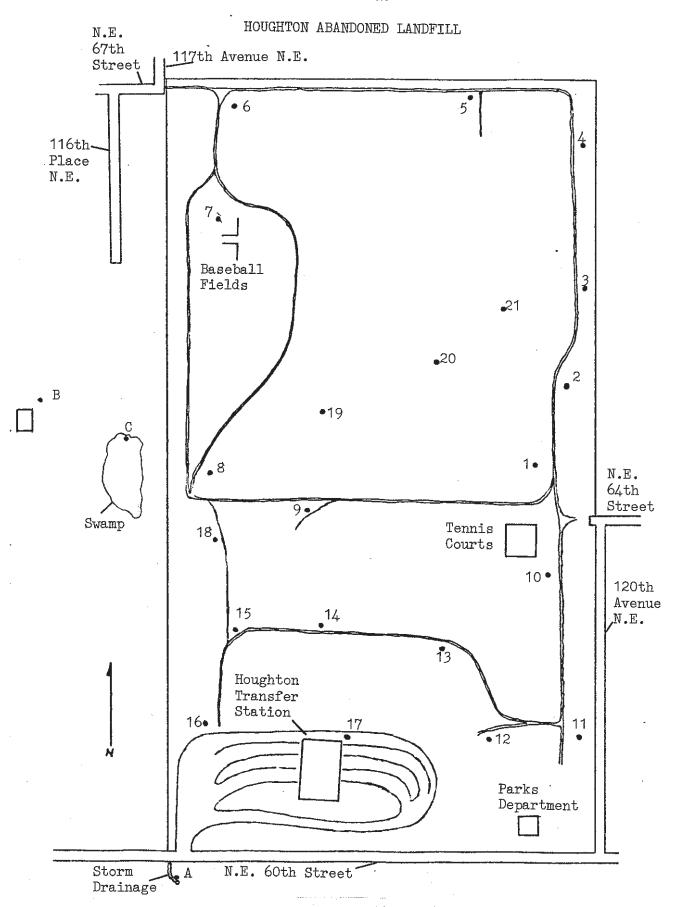


TABLE XXI
METHANE AND TRACE GAS CONCENTRATIONS
HOUGHTON ABANDONED LANDFILL

Site	Methane (%)	Trace Gas (ppm)*
1	0	0
2	Trace	0
3	0	0
4	Trace	0
5	Trace	0.2
6	11	0.8
7	0.4	-0.1
8	18	1.0
9	11	0.4
10	0	0
11	0	0
12	Trace	0
13	1.2	0.2
14	Trace	0
15	12	1.6
16	<u>≥</u> 4	0.4
17	12	0.8
18	11	0.2
19	≥ 4	0
20	0.4	. 0
21	0.7	0

^{*} Reading represents change from ambient air level

TABLE XXII SURFACE WATER PARAMETERS HOUGHTON ABANDONED LANDFILL

	Site A (1)	Site B (2)	Site C (3)
рН	6.4	6.4	6.3
Temperature	10.8	10.7	4.6
Dissolved	8.5	1.6	5.6
Electrolytic	0.4	0.4	0.0
Turbidity	55	22	29

Water from an off-site culvert
 Dug well 45 feet to static water
 Swamp surface water

KENT ABANDONED LANDFILL (MILL CREEK CANYON PARK)

The City of Kent operated a landfill at Woodland Way and Maple Street, this was closed in about 1961 with the opening of the Kent-Highlands operation by the City of Seattle. The former landfill stands today as a vacant clearing abutting the steep slopes of Canyon Park.

PAST AND PRESENT USE

There is no indication that the site in question had any major use until it was part of the landfill operation. A portion may have been part of a general pit operation earlier. Since its closure it has been left largely in a natural state and serves the community as a passive park.

SITE/ENGINEERING INFORMATION

Discussions with individuals who were in some way affiliated with the landfill operation provided some infomation regarding conditions on site. The fill on site is apparently deep, characterized as being 200 or more feet high. Cover fill was deep, and described as Class A or Class B bank run. Top soil was probably put on top of that, with the last topping being pit run from a cut about 100 feet deep from the area and hence probably similar to the original site conditions. The site is quite rocky at present.

During closure a berm was put around the site to control drainage.

WASTE DISPOSAL PRACTICES

This site served primarily the City of Kent. Excerpts from a 1956 document entitled <u>Community Report of Kent</u>, <u>Washington</u>, characterize the disposal operation as follows:

"The Kent City dump is an uncontrolled dump with 24 to 48 yards of municipal refuse being dumped daily as well as frequent dumping by individuals and a contractor.

"20 of the 32 persons interviewed having individual disposal use the Kent dump, 2 of the total 32 use other dumps, 1 uses a ravine and the remainder use burial, incineration and/or animal or fowl feeding.

"By using the term 'uncontrolled dump' reference is made to the fact that there is no dumping charge, no controlled burning, irregular burial, very loose supervision and little attempt to control scavenging. The height and width of the dump is only limited by the topography of the area. This dump is a public nuisance to nearby residences by reason of odors and smoke. The Kent dump is a definite health hazard both from the potentials of insect and rodent carriers of disease and direct disease contraction by scavengers. If in the future an air pollution problem should arise, such an open burning dump would contribute to air pollution. It is recommended that:

"'A sanitary method of refuse disposal be used and the Kent dump be closed and covered. Incineration and a sanitary landfill are the two accepted methods of sanitary refuse disposal by a community.'"

The type of waste deposited on site can be surmised from the characteristics of the community served. The above report notes that:

"Kent is a growing community in area and population. Its 1955 census of 3,675 in the incorporated limits has been expanded to well over 4,000 by annexations since that time. The population of the immediate area is considerably greater, however, as evidenced by the unusually large business district for a city of this size. The entire Kent School District has a population of over 12,000 persons in residence. A properous community, Kent has an assessed valuation of over four million dollars.

"The City is surrounded by numerous truck farms that contribute in large part to the economy of the community. Several sizable food processing industries provide seasonal employment for the community, as do the farms themselves. Although other industries are moving into the Kent area, the farms and [dependent] processing plants will continue to make an important contribution to the economy and will provide sanitation problems inherent to food industries."

The "other industries" mentioned are of concern as there began to be a shift in the industry type to manufacturing and non-seasonal industries about this time.

New establishments working in plastics, steel, and wood are reported. One plant is noted to discharge acid and neutralizer into a local creek and another is reported as handling a product causing dermatitis among employees. Knowing the general

policy of solid waste disposal typical during that time period, it is likely some potentially hazardous materials found their way to the Kent dump. However, most operators of the time remember the waste as being primarily residential waste.

SUSPECTED PROBLEMS

There are no specific records site characteristics documented which indicate any problems since the site was abandoned. With the exception of the possibility that some small amount of industrial waste may have been deposited there, it is unlikely problems should be anticipated.

FIELD RESULTS

On November 2, 1984 nine bore holes wree tested at the former Kent Landfill for methane and non-specific trace gases. Methane gas was observed only at trace levels. Non-specific trace gas tests were neutral relative to ambient air (Table XXII).

Surface water samples were retrieved from a stream at the bottom of Canyon Park at locations upstream and downstream from the fill site. No changes in water parameters were noted. No characteristics of leachate contamination were observed. These data are presented in Table XXIV.

FIGURE 13
KENT ABANDONED LANDFILL

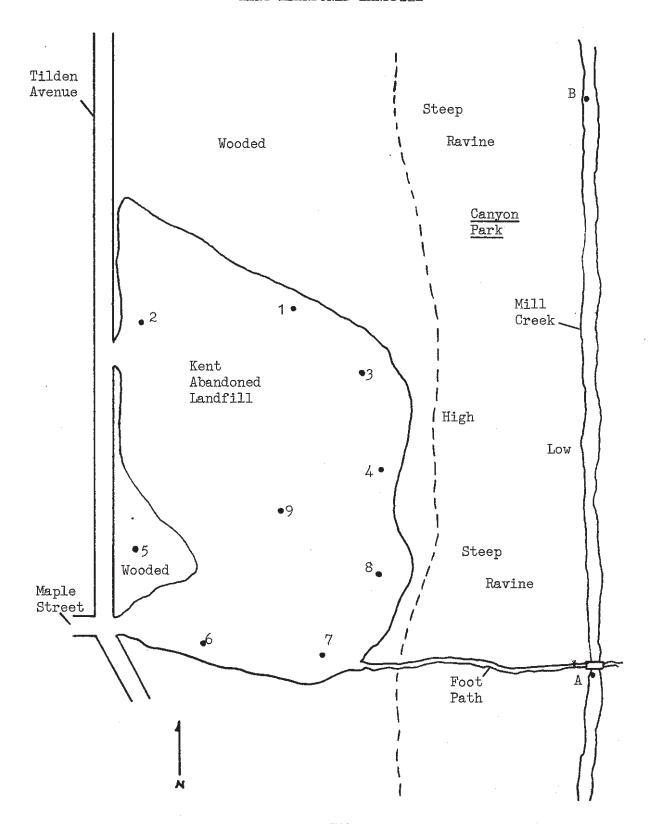


TABLE XXIII

METHANE AND TRACE GAS CONCENTRATIONS

KENT ABANDONED LANDFILL

Site	Methane (%)	Trace Gas (ppm)*
1	Trace	0
2	0	0
3	Trace	0
4	0	0
5	0	0
6		0
7	Trace	. 0
8	Trace	0
9	Trace	0

^{*} Reading represents change from ambient air level

TABLE XXIV SURFACE WATER PARAMETERS KENT ABANDONED LANDFILL

	Site A (1)	Site B (2)
рн	6.2	6.3
Temperature	9.5	9.7
Dissolved	9.3	9.4
Electrolytic	0.2	0.2
Turbidity	65	66

⁽¹⁾ Stream water upstream from landfill(2) Stream water downstream from landfill

McMICKEN HEIGHTS

In 1942 a garbage disposal site was located in the South King County area on the hillside southwest of the City of Tukwila. Bordered on the west by 51st Avenue South, and the east by the area currently occupied by Interstate 5 at the location of 53rd Avenue South, the south by South 176th and on the north by South 173rd, the site is a steep ravine where the bordering roads deadend.

No documents regarding this site were located with the exception of the old map showing its location. Long time residents of the area do not recall the dump, referred to as the McMicken Heights garbage disposal site.

The site was probably a small residential site which closed sometime in the 1940's and was entered on 51st Avenue South.

Current land use in the area is residential on all sides, exclusive of the freeway. The disposal site itself remains undeveloped except for a trail system through the area. Some residents recall mention of a "gun club" at one time, so it is possible the site may have been used for target practice sometime after the fill closed and before residential development began. Plat maps indicate the site is approximately 13.95 acres and list it as the Castillo Land Co. property.

Unauthorized neighborhood dumping of rubbish is still done at the northern edge of the site.

FIELD RESULTS

Nine bore holes were tested at this site for methane and non-specific trace gas concentrations on November 16, 1984. Results appear in Table XXV. Methane levels were low as measured from the nine test sites, ranging between 0% to 0.2% gas. Trace gas levels never exceeded ambient air concentration.

No surface water was observed directly on-site. Surface water run-off streams were sampled at east and west points of the north canyon and displayed no evidence of leachate pollution (Table XXVI).

FIGURE 14
MeMICKEN HEIGHTS ABANDONED LANDFILL

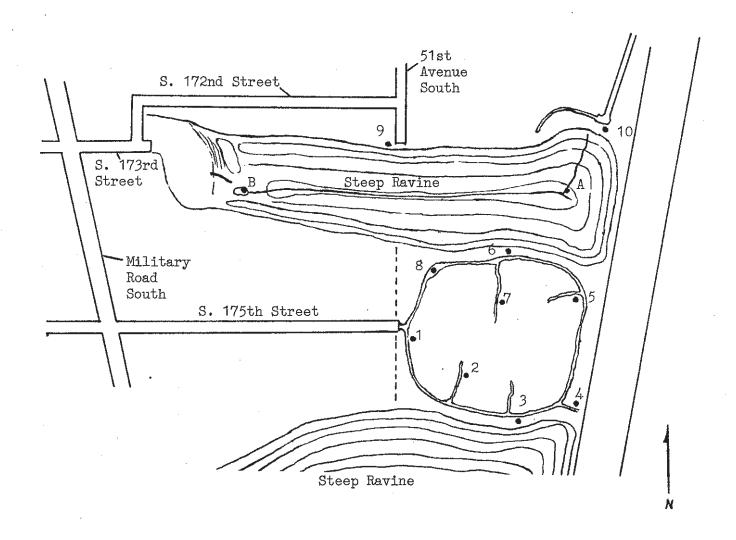


TABLE XXV

METHANE AND TRACE GAS CONCENTRATIONS

McMICKEN HEIGHTS ABANDONED LANDFILL

Site	Methane (%)	Trace Gas (ppm)*
1	0	0
2	Trace	0
3	0	0
4	Trace	0
5	Trace	0
6	Trace	0
7	Trace	0
8	0.2	0
9	Trace	0
10	Trace	0

^{*} Reading represents change from ambient air level

TABLE XXVI

SURFACE WATER PARAMETERS McMICKEN HEIGHTS ABANDONED LANDFILL

	Site A (1)	Site B (2)
рН	5.6	6.1
Temperature	10.4	10.3
Dissolved	10.1	9.9
Electrolytic	0.2	0.4
Turbidity	5.	2
·		

Flowing water from storm drain run-off Pooled water from storm drain run-off

NORTH BEND

The North Bend Abandoned Landfill stands today as a small clearing along the west side of the Middle Fork Road, about one mile north of the "Y" turn from Edgewick Road located on the east side of North Bend. There are several well-defined paths extending into the new-growth woods. A gravel pit is across the road and the Snoqualmie River some distance to the north.

This approximately two acre landfill was city operated during the 1950's. The Seattle-King County Health Department baited the area but did not have any other major responsibilities at the site. Being in a small rural area, disposal of any hazardous material there seems unlikely.

FIELD RESULTS

On October 20, 1984 ten bore holes were tested at this site for methane and non-specific trace gas concentrations. These data appear in Table XXVII.

Methane levels were observed ranging between 0% to 0.4% from the ten test bore holes. All non-specific trace gas readings were neutral relative to ambient air concentrations.

No surface water was observed on or about the site.

FIGURE 15 NORTH BEND ABANDONED LANDFILL

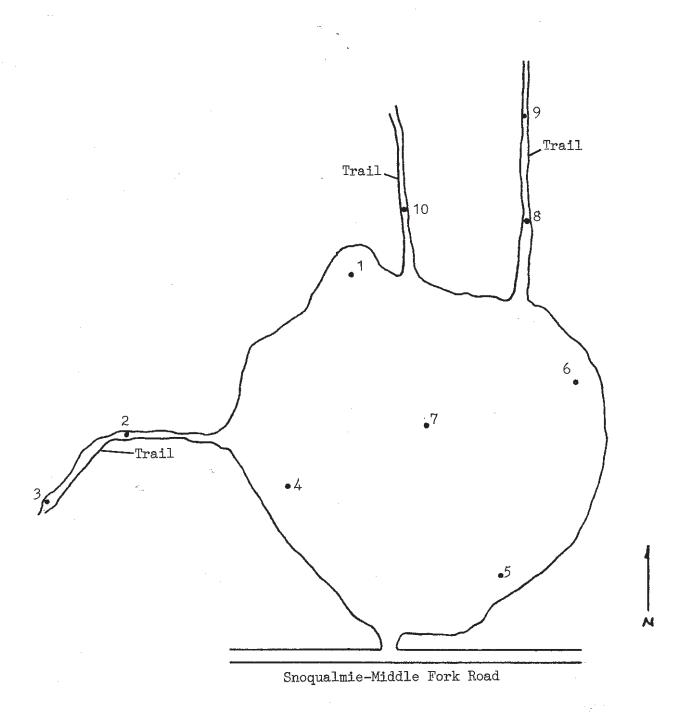


TABLE XXVII

METHANE AND TRACE GAS CONCENTRATIONS

NORTH BEND ABANDONED LANDFILL

Site	Methane (%)	Trace Gas (ppm)*
1	Trace	0.1
2	Trace	0
3	Trace	0
4	Trace	0
5	Trace	0
6	Trace	0
7	Trace	0
8	Trace	0
9	Trace	0
10	.4	0

^{*} Reading represents change from ambient air level

PACIFIC CITY

The King County refuse dump in the City of Pacific was located on an approximately seventy two acre site at the location of the present Stuck River Park. Encompassing both sides of the Stuck (White) River the site was large and old. Early maps show "garbage dump" operations as early as 1921 and possibly even in 1914. It was closed about 1961 with refuse being routed to the Puyallup Cutoff landfill.

The legal description of the site is:

E 495 FEET OF W 660 FEET OF NW 1/4 OF SE 1/4 LESS NORTH 200 FEET OF W 100 FEET AND LESS THE PORTION DEEDED TO KING COUNTY 12/14/14, TAX LOT 40 SECTION 36, TOWNSHIP 21 NORTH, RANGE 4 EAST

PAST AND PRESENT USE

There is no information available on the land use prior to the landfill. Since it is located in the traditional farm belt of South King County it can be assumed it was, as it still is predominantly, rural and agricultural. Apparently flooding from the Stuck River was a chronic problem in the early 1900's. Several documents exist showing deeds transferring lands needed for Stuck River improvements. These included concrete bulkheads and rechanneling the river for the benefit of both King and Pierce Counties.

On March 7, 1966 the City of Pacific was granted permission to use about twenty one acres of the abandoned site for park purposes for a period of twenty five years. The City proceeded with site development and discovered not all of the land contemplated for park development had been properly described in the 1966 King County Resolution No. 31548. Thus on August 1, 1969, King County granted the City the permission to use the land originally intended for park purposes. The "new" city park was developed and dedicated in September 1972.

The remaining area adjacent to the landfill is zoned either multi-family or single family residential or industrial property.

SITE/ENGINEERING INFORMATION AND WASTE DISPOSAL PRACTICES

Nothing regarding the use or type of waste disposed was found in the existing records.

Individuals who remember the operation of the landfill recall that it was ten to twelve feet deep. Water was typically encountered, at least at twelve feet and the toe of the fill appeared to border the river. Most of the refuse was burned.

SUSPECTED PROBLEMS

No major problems are anticipated at this site <u>unless</u>, given the rural character of the region, pesticides or their containers may have been dumped here.

No records exist to substantiate that. This landfill is mentioned in the EPA ERRIS list.

FIELD RESULTS

On October 23, 1984 fifteen bore holes were tested at the Pacific site for methane and non-specific trace gas concentrations. These data appear in Table XXVIII.

Methane gas levels were observed only at low levels throughout the site with a peak reading of 0.4%.

Non-specific trace gas levels were also non-significant (Range = -0.1 to +0.3 ppm relative to ambient air) with the exception of the 5.6 ppm reading of hole #5. This test was made within the center of a dead patch of grass. Five test holes (a - e) placed within forty feet of hole #5 yielded non-significant levels of methane or trace gases.

A water sample retrieved from the adjacent White River was not indicative of leachate Table XXIX.

FIGURE 16
PACIFIC ABANDONED LANDFILL

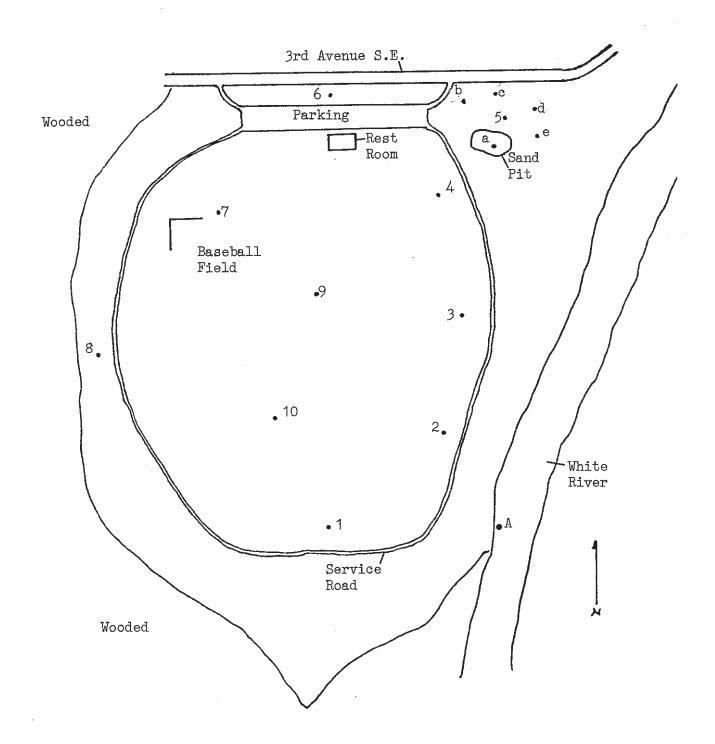


TABLE XXIX

SURFACE WATER PARAMETERS PACIFIC ABANDONED LANDFILL

											Site A (1)
рН	•	•	•		•		•		•		5.1
Temperature o _C	•		÷	•	•	٠	•	•	•	•	8.5
Dissolved . Oxygen ppm	•		•	•	•	•	•	•	•	•	11.6
Electrolytic Conductivity m~v/cm	•	•	•	•	•	•	•	•	•	•	0.2
Turbidity .	•	•	•	•	•	•	•	•	•	•	5

TABLE XXVIII METHANE AND TRACE GAS CONCENTRATIONS PACIFIC ABANDONED LANDFILL

Site	Methane (%)	Trace Gas (ppm) (1)
1	Trace	0.1
2	0	0.3
3	0	-0.1
4	0	0
5(2)	0.3	6.2
6	0	0
7	0	0
8	0	
9		0
10	0	-0.1
a	0.4	0
Б	Trace	0
С	0.2	0
d	Trace	0.1
е	Trace	0

⁽¹⁾ Reading represents change from ambient air level(2) Test hole centered an area of dead grass

PUYALLUP/KIT CORNER

King County operated a 30.45 acre sanitary landfill in Section 20, T21N, R4E, immediately east of I-5 and one-quarter mile north of South 360th Street.

PAST AND PRESENT USE

This was apparently tax title property deeded to King County for operation of a "garbage dump." Old maps of the section show a King County dump site here in 1947. The original parcel of land consisted of forty acres, but 9.55 acres were deeded to the State of Washington Department of Highways in 1959, reducing the landfill portion to 30.45 acres. Aerial maps of the area taken in May, 1970 show the landfill as closed.

It is currently vacant land being surplused by the County. Brush and high grass cover the site which is returning to its natural state. Bike trails can be seen on site but the original frontage access road to the site has been closed.

SITE/ENGINEERING INFORMATION AND WASTE DISPOSAL PRACTICES

This was one of the larger King County sites. As the Bow Lake, Renton Junction, and Pacific sites closed, refuse was hauled to this site. It operated until shortly after the opening of the Cedar Hills landfill. It is estimated that thirty to forty feet of refuse fill the site.

Operational maps indicate that the borrow site for cover was located at the south edge of the property. This was also one of the last areas filled. The northern edge of the property was also the site of some of the older fill area, but was expanded and raised with new fill in later stages of operation. The same, ie. raising the fill, appears to have occurred in the central portion of the older fill area.

A paved service road fronted the west side of the site and a dirt service road appears to have run around the northern, eastern and southern edges. A drainage ditch was on the southeastern and southern edges of the fill which appear to be the lower elevations.

SUSPECTED PROBLEMS

Since this site received refuse from a large geographic area and operated for about twenty five years, it seems likely that almost anything could be found here. The site is mentioned on the EPA ERRIS list. The landfill's proximity to the major Tacoma industrial area raises some suspicions regarding the possible unauthorized disposal of hazardous materials at some time during its operation. However, the presence of two nearby landfills in Pierce County at that time probably make this a remote possibility.

The cover fill, apparently put on the site when closed with no provisions made for methane release, make likely a potential for methane build up on site and additional land settlement.

FIELD RESULTS

For the purposes of this study, methane gas was tested from twenty-two test areas located throughout the Puyallup/Kit Corner site on October 27, 1984 which was a day of heavy rainfall. These data appear in Table XXX. Non-specific trace gases were not measured at that time because of that test equipment's sensitivity to excessive moisture.

Methane was observed at levels within or above the explosive range for methane gas (4% to 18%) in ten test holes, located throughout the site with peak levels reaching 68% to 72%. Fissures and areas of dead vegetation were frequently

encountered. Methane gas was measured at 50% directly from the fissure at test site #22. Readings at five test holes were confounded by the high water table.

Trace gases were measured from six bore holes on December 1, 1984. Levels were non-significant, ranging from -0.2 to +0.2 ppm relative to the ambient air levels (Table XXXI).

During examination of the site for leachate problems a metal drainage conduit was found along the west side of the site's access road which drained into a surface water collection stream. A discoloration characteristic of leachate was noted on the drainage conduit. Test parameters of the stream water under the conduit on October 27 did not indicate a leachate problem at that time, though a slightly acidic pH (5.3) was noted.

Water was also sampled from an east perimeter stream. Leachate contamination was not indicated at that location. Surface water data are presented in Table XXXII.

FIGURE 17

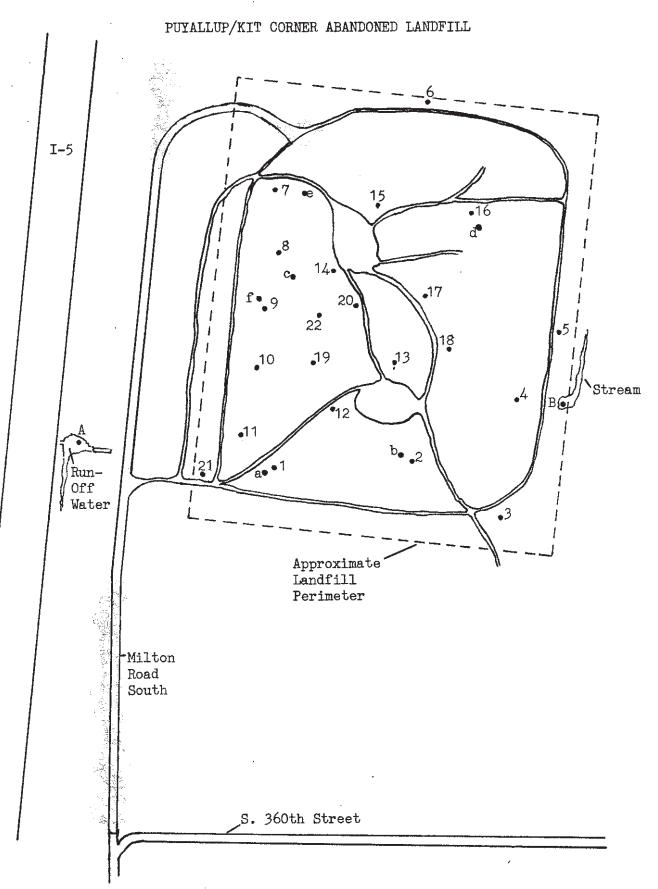


TABLE XXX METHANE GAS CONCENTRATIONS PUYALLUP/KIT CORNER ABANDONED LANDFILL

	Site	Methane (%)
	1	0.8
	2	60
	3	Trace
	4	72
	5	25
	6	Trace
	7	Water Table
	8	Water Table
	9	70
1	10	18
•	11	Water Table
•	12	Water Table
	13	5
٠	14	12
√r r (Å=	15	68
	16	Water Table
	17	Trace
	18	5
\	19	3
	20	47
	21	Trace
	22	50

TABLE XXXI

TRACE GAS CONCENTRATIONS
PUYALLUP/KIT CORNER ABANDONED LANDFILL

Site	Trace Gas (ppm)*
a	-0.2
b	0
С	0
d	0.2
e	-0.2
f	0

^{*}Reading represents change from ambient air level

TABLE XXXII SURFACE WATER PARAMETERS

PUYALLUP/KIT CORNER ABANDONED LANDFILL

	Site A (1)	$\underline{\text{Site B}}$ (2)
рН	5.3	6.0
Temperature	8.9	8.2
Dissolved	10.2	9.5
Electrolytic Conductivity m· v /cm	0.1	0.1
Turbidity ppm	12	175

⁽¹⁾ Run-off stream surface water(2) Stream surface water, east perimeter

REDONDO PIT

The Redondo Pit referred to a large gravel pit located at the intersection of the Pacific Highway South and South Dash Point Road a few blocks north of Federal Way High School. It was a nineteen acre site.

The Redondo Pit was operated by the King County Department of Public Works as a gravel pit for several years. During the 1940's the Port of Seattle, the United States Navy, and the County used the site as an oil dump. Bilge oil, crankcase oil, and road oil were dumped there. It was common to burn the oil off until residents and/or regulations curtailed the practice. In 1969, when the gravel pit was exhausted, the land was transferred to the King County Parks for use as a community park site. Final plans for the park were made in 1973 and the current Sacajawea Park was completed.

King County Park Department officials report running into bunker oil during development. They excavated what they could find, using earth fill from the Federal Way School District's Sacajawea Junior High School site. Today the area is a three level athletic/recreational facility with track and football/soccer activities on the lower level, baseball and tennis on the middle, and general playground/picnic area on the upper section.

FIELD RESULTS

This site was tested on November 7, 1984. Methane gas levels were not observed above 0.6% from the twenty bore holes tested. Trace gases were noted with peak readings above ambient air levels of +1.0 ppm in the lower section, +8/8 ppm in the center, and +1.2 ppm in the upper section. A ground water table was observed in test holes located on the lower section (Table XXXIII).

On-site structures (restrooms) exhibited no indication of problematic ground settling due to the fill.

No surface water was observed on or about the fill area. However, water samples were obtained from run-off sewer drains located on the lower tier (site A) and the upper tire (site B). These data are presented in Table XXXIV. Parameters of leachate contamination were not observed in the water samples.

FIGURE 18
REDONDO PIT ABANDONED LANDFILL

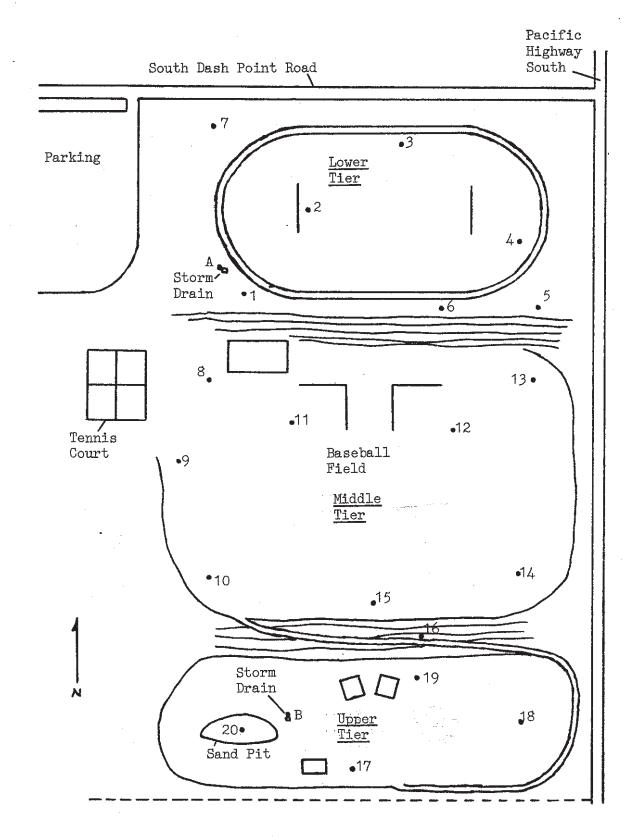


TABLE XXXIII

METHANE AND TRACE GAS CONCENTRATIONS
REDONDO PIT ABANDONED LANDFILL

Site	Methane (%)	Trace Gas (ppm)*
1	0	1.0
2	0	0.8
3	0	0
4	0	0
5	0	1.0
6	Water Table	Water Table
7	0	0
8	0	0
9	0.6	0.8
10	Trace	
11	0	0
12	. 0	0
13	Trace	0
14	0	0.4
15	Trace	8.8
16	Trace	
17	Trace	1.2
18	Trace	1.0
19	Trace	0.8
20	Trace	0.2

^{*} Reading represents change from ambient air level

TABLE XXXIV

SURFACE WATER PARAMETERS REDONDO PIT ABANDONED LANDFILL

$\underline{\text{Site A}}^{(1)}$	(2)
pH 5.9 6.0	
Temperature 8.8 9.7	
Dissolved 8.4 3.5 Oxygen ppm	
Electrolytic 0.1 0.05 Conductivity m·v/cm	
Turbidity 13 8	

⁽¹⁾ Storm-drain water(2) Storm-drain water

RENTON HIGHLANDS

The Renton Highlands abandoned landfill was on about 11.2 acres. On the south side of Northeast 3rd Street slightly east of Mt. Olivet Cemetery, and west of the Southeast District Office of the Seattle-King County Health Department.

PAST AND PRESENT USE

Operating since the 1940's, correspondence dated June 7, 1951, described the site as follows:

"The area of approximately 12 acres, privately owned, lies directly south of the Renton Highland Housing Project and just east of the Renton City limits, a former gravel pit with plenty of cover material. It is 1 1/2 miles from the Renton City Hall, a minimum trucking distance.

"This site appears to be a good one for dumping, is in need of fill to round out a broken area and has a probable expectancy period of 10 to 20 years. It has about the best possible soil condition, good proximity to collection area, but also good visual and wind isolation from thickly settled parts of the city."

Unfortunately, perhaps, in the thirty years since this correspondence the site did not stay "isolated." The Renton Highlands has seen substantial growth in the ensuing years. The dump site was closed by the late 1960's. Today the land is still undeveloped private property.

In addition, it appears that another site operated slightly north of that location. Correspondence dated January 31, 1949 from D.L. Evans, County Road Engineer to the County Commissioners, notes the following:

"The attached application from the City of Renton to purchase Lots 1, 2, 3, 8, 9 and 10, Section 7, Rainier Acres, to be used for garbage disposal site.... This location is a portion of an old gravel pit which has been worked out and is of no further value to the County for this purpose...."

This area, just west of Renton Vocational Institute is currently used by the Housing Authority for residential dwellings.

SITE/ENGINEERING INFORMATION AND WASTE DISPOSAL PRACTICES

According to the 1951 correspondence noted earlier, "The soil type in this area is Everett gravelly sandy loam, a very deep deposit of gravel, very well drained with little if any possibility of horizontal seepage."

The site was a large, major fill in the area which is remembered as receiving "everything." Fires were common with smoke drifting into the neighborhood adjacent to the site.

It is thought that a portion of Northeast 4th and/or the housing project in that general area may actually be on some of the older fill.

In 1951 it was noted that major material being dumped was ash from the housing project which "has no fumes, smoke, or odor."

SUSPECTED PROBLEMS

Although no records regarding waste disposal are available the existence of several industries, including Boeing and Paccar, in Renton during the life span of this fill raise the question of possible hazardous materials disposal. The site is mentioned in the EPA ERRIS list.

FIELD RESULTS

On November 14, 1984 eight bore holes were tested for methane and non-specific gas at this site along the north side of Northeast 3rd Street. Three bore holes were tested on January 25, 1985 for only methane gas along the south side of Northeast 3rd Street. Methane gas was not observed above trace levels from any of the test holes. Trace gas levels were all observed at equilibrium with ambient air conditions.

A water sample was retrieved from the Mt. Olivet Creek on January 25, 1985. Signs of leachate contamination were not indicated.

FIGURE 19
RENTON HIGHLANDS ABANDONED LANDFILL

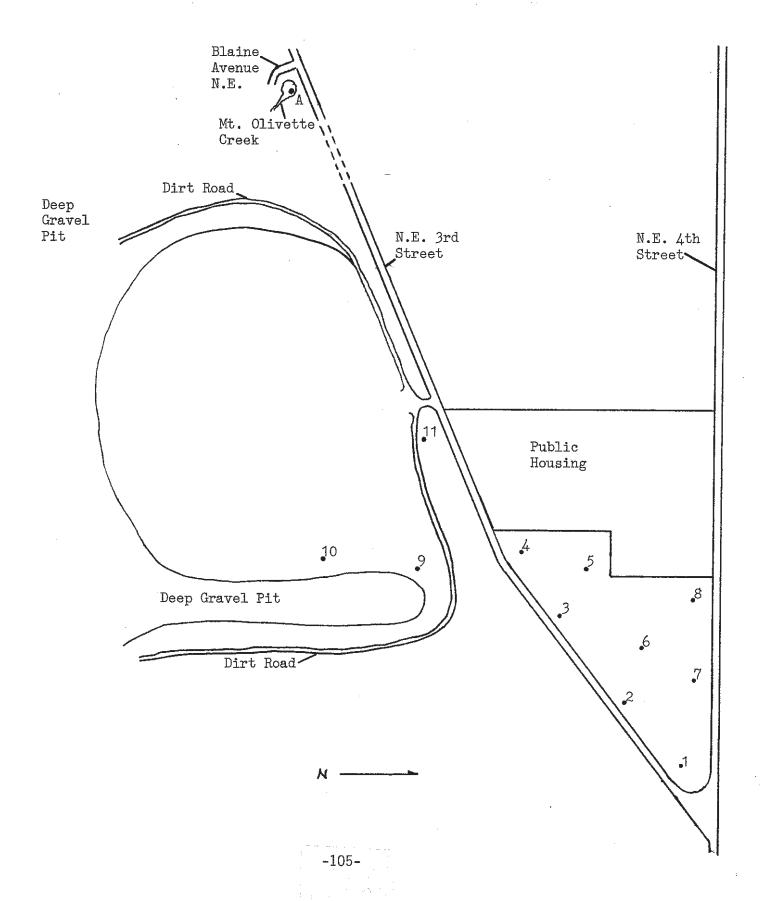


TABLE XXXV

METHANE AND TRACE GAS CONCENTRATIONS
RENTON HIGHLANDS ABANDONED LANDFILL

,		/1)
<u>Site</u>	Methane (%)	Trace Gas (ppm) (1)
1.	Trace	0
2	0	0
3	Trace	0
4	0	0
5	0	0
6	Trace	0
7	0	0
8	0	0
9	0	
10	0	
11	0	

⁽¹⁾ Reading represents change from ambient air level

TABLE XXXVI

SURFACE WATER PARAMETERS

RENTON HIGHLANDS ABANDONED LANDFILL

													Site (A)
рН	٠	٠	•	•	•	•		•	•		•	•	6.7
Temperature °C	•	•	•	•	•	•	•	•	•		•	•	8.3
Dissolved . Oxygen (ppm)	•	•	•	•	•	•	•	•	•	•	•	•	9.8
Electrolytic Conductivity mr/cm	•	•	•	•	•	•	•	•	•	•	•	•	0.2
Turbidity . (ppm)	•	•	•	•	•	•	٠	•	•	٠	٠	•	4

RENTON JUNCTION (MONSTER ROAD)

King County operated a refuse disposal site at the Renton Junction near Longacres from approximately 1946 to about 1961. It operated on land leased from the Northern Pacific Railway Company described as:

"Those portions of Lots 31 and 32 of Interurban Addition to Seattle, according to the recorded plat thereof, lying northeasterly of a line parallel with and distant 36 feet northeasterly, measured at right angles, from the center line of the most northeasterly main track as now constructed across said lots; together with the southwesterly one-half of the original channel of the White River which attached thereto when the channel of said river was relocated and constructed along the southwesterly side of the Railway Company's tracks."

Additional land was obtained from Mr. Fred Nelson of Renton for the operation described as: "that portion of government Lot 6, Section 24, Township 23N, Range 4EWM, lying between the westerly right of way line of the Steel Hill County road No. 24-23-4-1 and the centerline of the old channel of the Green River."

PAST AND PRESENT USE

As noted by the legal description, part of this property was once the old river channel which became part of the Northern Pacific Railway right of way and private rural property. After its operation as a sanitary landfill the land was surplused and in 1979 purchased for commercial use. The site is currently used by a decorative rock company and is used for storage of crushed rock and gravel associated with that operation.

Across the Green River lies Fort Dent Park, southwest of the site is the Riverview Nursery and to the east is the Metro Secondary Sewage Treatment Plant.

SITE/ENGINEERING INFORMATION AND WASTE DISPOSAL PRACTICES

No specific geological or hydrological information was found regarding the site other than it was part of the old river channel. It may be surmised that the

base soil is clay, sand and gravel overlain by fill forty to fifty feet deep and topped off with a relatively impermeable layer and topsoil.

Once the landfill operation started there are records showing numerous complaints regarding the site. Nuisance conditions prevailed in the warm weather. Fire and smoke were reported night and day. Inadequate fill dirt for the cover of the operation was noted, as were problems of dumping sewage and oils on site. The County had a contract for the maintenance of this dump with a private concern in effect until December 31, 1957. This contract called for bulldozing and compaction twice weekly with the top side being covered with eighteen inches of dirt. It also required that the garbage be deposited in lifts or layers not to exceed twelve feet in depth after initial compaction. There was no earth available on the site and all cover material had to be brought in by trucks. The only earth available within hauling distance had been of a hardpan type with a heavy clay concentration.

The site was used not only by the County, but according to copies of agreements, also by the City of Renton to dispose of a portion of their garbage and refuse.

The seriousness of the fires at the south end of the dump operations in 1950 prompted correspondence between the King County Fire Marshal and the Health and Sanitation Department, who ran the landfill, to confine burning to the north area of the landfill.

The landfill was closed effective December 27, 1957 with directions for refuse to be taken instead to either the Bow Lake Fill at South 188th and Military or the landfill at South 352nd Street about one-half mile east of the Puyallup cut off highway, both still in operation at that time.

However, the Health Department was requested to continue filling operations as plans to discontinue the fill were described as leaving the fill in an unusable

condition. The fill was reopened, using County equipment to conduct operations, which continued for about two years.

SUSPECTED PROBLEMS

Due to the proximity of the old landfill site to the Green River, it is conceivable that leachate may reach the river.

The site is recorded on the EPA ERRIS list. The site's proximity to several industries in South King County make it possible that some potentially hazardous materials, including oil, were dumped at the site during its years of operation.

FIELD RESULTS

The Renton Junction Abandoned Landfill was tested for methane and trace gas emissions on January 10, 1985. These data are presented in Table XXXVI. Methane gas levels were observed ranging between 17% and 33% from test holes located at the northerly half of the former fill. Lower levels of methane gas (3% to 5%) were observed within the southerly section.

Trace gas levels were observed ranging between -4.8 ppm to 0 ppm relative to the ambient air.

A surface water sample was retrieved along the shoreline of the Green River immediately adjacent to the former landfill. Leachate contamination was not indicated (Table XXXVII).

FIGURE 20
RENTON JUNCTION ABANDONED LANDFILL

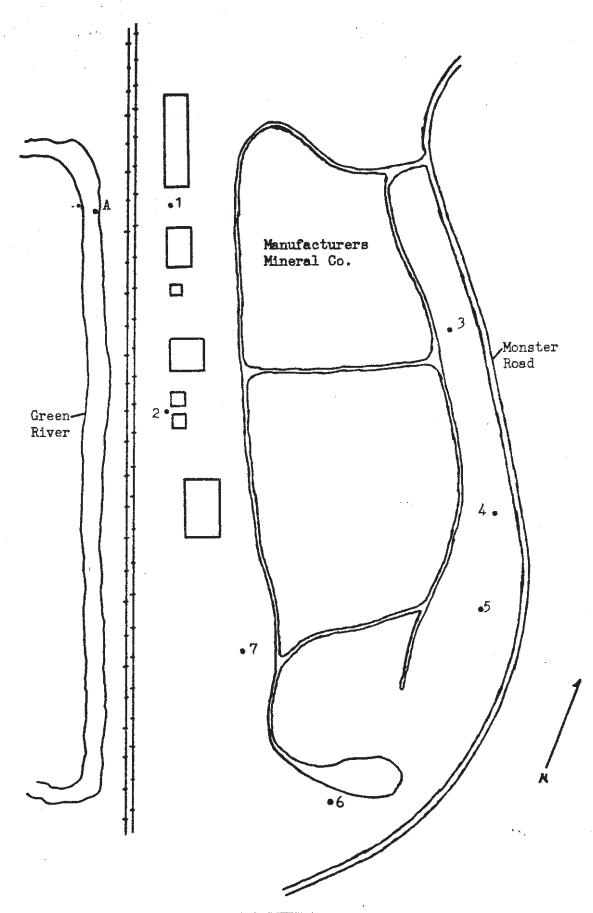


TABLE XXXVII METHANE AND TRACE GAS CONCENTRATIONS RENTON JUNCTION ABANDONED LANDFILL

<u>Site</u>	Methane (%)	Trace Gas (ppm)
1	33	-2.7
2	30	-4.6
3	17	-4.8
4	4	0
5	3	0
6	3	0
7	5	0

⁽¹⁾ Trace gas measurements made with 11.2 eV HNU probe(2) Reading represents change from ambient air level

TABLE XXXVIII

SURFACE WATER PARAMETERS

RENTON JUNCTION ABANDONED LANDFILL

													Site (A) (1)
рН			•		•	•	•	•	•		•	•	5.6
Temperature °C	•	•	•	•	•	•	•	•	•	•	•	•	5.2
Dissolved . Oxygen (ppm)	•	•	•	•	•	•	•	•	•	•	•	•	9.8
Electrolytic Conductivity mu/cm		•	•	•	•	•	•	•	•	•	•	•	0.2
Turbidity . (ppm)	•	•	•	•	•	•	•	•	•	•	•	•	1

⁽¹⁾ Green River surface water

SKYKOMISH

The town of Skykomish operated a landfill outside the corporated limits of the town between the old Cascade Highway, south of the Railroad right of way east of town from about 1946 until 1979. The parcel of land is described in Sec 25, T 26 N, R11EWM, as Gov't Lot 6 lying E of the W line of # 1/2 of SW 1/4 of Sw 1/4 produced N across Gov't Lot 6 and South of Great Northern Rwy Co. right of way and northerly of the County road, westerly of line in boundary line agreement, King Co. recording #7812070786 EXCEPT portions deeded to King County under record #4627707 and #7503170254.

In 1979 this was transferred to the County for a county operated drop box station.

The waste disposal practices, site conditions and problems of the landfill were discussed in depth in a report entitled "Solid Waste Disposal Alternatives for the Skykomish Area - a Feasibility Study" prepared by the King County Solid Waste Division in May 1978. This assisted in the decision to develop the current transfer station found on site. Excerpts from that report follow.

SITE/ENGINEERING INFORMATION

"Geology data for the Skykomish area is scarce. Soil studies done by the U.S.D.A. - Soil Conservation Service in King County do not cover Skykomish.

"During the winter of 1976-77, the King County Department of Public Works took some soil boring samples in connection with a project to renovate the Old Cascade Highway around Skykomish. The current landfill is located a few hundred feet north of the Old Cascade Highway and one mile east of the Town's business district.

"The samples taken in the general area of the landfill went no deeper than six-plus feet. The soils were composed of sand and gravels with cobbles up to 8 inches in diameter.

"The gravel pit excavations made in former years which now comprise the landfill site, show similar material at all exposed depths (up to 15 feet deep)."

WASTE DISPOSAL PRACTICES

"The town of Skykomish [now] operates an open burning (landfill). This has been the solid waste disposal practice in the Skykomish area for as far back as long-time residents can remember. The property which the Town now uses as a landfill was originally purchased from the Northwestern Improvement Company in 1946. At that time, the Town's landfilling practice consisted of digging a trench with dimensions of about 30' x 100' x 12'. The garbage would be dumped in the trench and burned once or twice a week. Once or twice a year the Town would obtain a backhoe and/or bulldozer and compact the waste in the trench and cover it with about three feet of soil. When one trench was full, they would dig another and repeat the process.

"The Town's current practice is quite similar. They now use an area method of fill rather than digging trenches. The burning of the solid waste is done within the proscriptions of PSAPCA's Resolution No. 353 in that rubber tires, waste oil, asphalt roofing or flooring materials, dead animals or demolition material are not burned. The burned residue is compacted in an area left from a gravel excavation operation. The compacted waste is still covered only at infrequent intervals."

SUSPECTED PROBLEMS

- "...Assuming that the problems of inadequate cover material could be solved, the present landfill site would still have a leachate problem...At the Skykomish landfill, leachate is produced when rain or melting snow percolates through the buried waste.
- "...The Skykomish landfill is located within 100 yards of the South Fork of the Skykomish River. Thus the possibility of water pollution from leachate exists.
- "Visual inspection of the Skykomish River near the landfill reveals no obvious leachate outfall.

"The lack of obvious leachate problems does not however make the Skykomish landfill a conforming site. Leachate is being produced, although it's simply not traceable at present...."

The closure of the site and the development of the drop box station eliminated most of these problems.

FIELD RESULTS

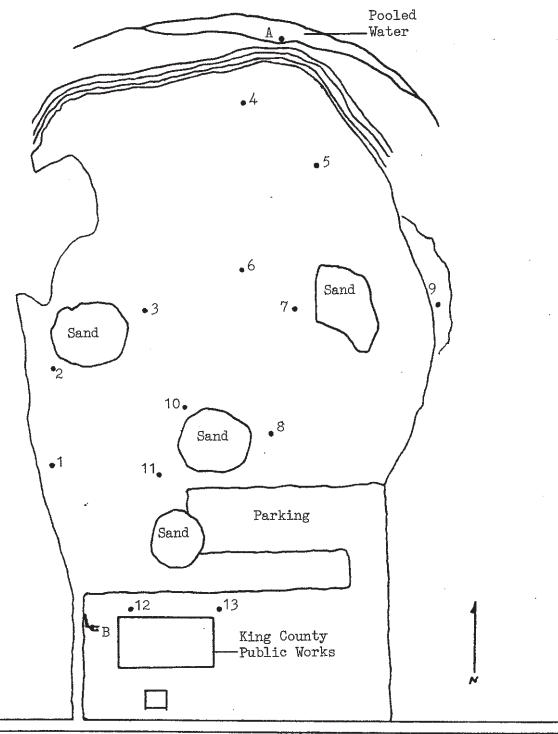
On November 3, 1984 thirteen bore holes were tested for methane and non-specific trace gases at the Skykomish site. These data are presented in Table XXXVIII.

Throughout the site, only low levels of methane gas were observed with a peak reading of 0.8%. Trace gas levels were all neutral relative to the ambient air concentrations with a range of -0.1 to +0.1 ppm.

Heavy intermittent fainfall and a four inch snow cover created surface water pooling on the site which confounded three test hole results.

Surface waters were tested from a pool at the northern base of the fill and from a storm water run-off stream at the southern entrance (Table XXXIX). Leachate contamination was not indicated by the test parameters, although weather conditions may have masked such a problem.

FIGURE 21
SKYKOMISH ABANDONED LANDFILL



Stephen's Pass Highway

TABLE XXXIX

METHANE AND TRACE GAS CONCENTRATIONS

SKYKOMISH ABANDONED LANDFILL

Site	Methane (%)	Trace Gas (ppm)*
1	0	0
2	Water Table	Water Table
3	0.8	-0.1
4	0.4	0
5	Trace	0
6	0.4	0
7	Water Table	Water Table
8	Trace	0
9	0	0.1
10	Trace	0.1
11	Trace	0.1
12	0	0
13	Water Table	Water Table

^{*} Reading represents change from ambient air level

TABLE XL SURFACE WATER PARAMETERS SKYKOMISH ABANDONED LANDFILL

	Site A (1)	Site B (2)
рН	6.6	7.1
Temperature	5.1	4.6
Dissolved	8.1	10.9
Electrolytic Conductivity m ***/cm	0.4	0.4
Turbidity	13	17

⁽¹⁾ Standing pooled water at north base of landfill(2) Flowing run-off stream

SUNSET PARK

The southeast corner of Sunset Park at South 140th and 18th Avenue South is the general location of an abandoned oil dump site.

PAST AND PRESENT USE

King County Park and the Sunset Shops of the Public Works Department have facilities on the east edge of the Sunset Park. Directly south of these facilities is the site of the old oil dump. It apparently operated from about 1936 to 1941 or 1942. The Navy and possibly the Port of Seattle are remembered as using the dump facility. The site was two to three acres in size and the dump from ten to twelve feet in depth. Only liquid waste was deposited here and at least 98% of that was oil waste.

A number of precautions have been taken at the site since closure of the site. These have included fencing, storm drainage, and the installation of a skimmer and baffle on the drainage inlet into Tub Lake. Oil problems remain however as the oil bubbles up in Tub Lake. There is a high water table in the area and an oily sheen can at times be seen on high water adjacent to Tub Lake.

SITE/ENGINEERING INFORMATION AND POTENTIAL PROBLEMS

There is some information indicating that this site was originally a gravel pit. A 1980 report detailing the plans for the "North Sea Tac Park" prepared by Jongigan, Gerrard and McNeal, Inc. describes the general geological and hydrological conditions surrounding this area. In that report they note that South Park is part of the Miller Creek Drainage Basin, with Miller Creek itself <u>not</u> being a distinctive stream at this time, but rather characterized as narrow channelizations or shallow murky areas.

That report further notes that there are particular problems at this site including,

"pollution and debris caused by garbage dumped into the stream channel and fertilizers from playfields...a serious on-site source of pollution is found adjacent to Tub Lake...at present the oil leaches through the soil into the surrounding area. The light oil slick intermittently visible in Tub Lake has been attributed to the abandoned dump."

Soils in the area are poorly drained. As noted in the 1980 study of the area, the soils are derived from glacial deposits with the most common parts being Alderwoods, Everetts, and Indianola.

"The remainder of the soils at North Sea-Tac Park were formed in glacial depressions or clayey alluvium and are very poorly drained. Derived from vegetation in varying degrees of decomposition, the soils are acidic with a high organic content. Orcas and Seattle Muck (Or and Sk) are characterized by thick layers of peat. Organic soils are inappropriate for construction because of high compressibility. Because the water table is at or near the surfaces of these soils, they are unsuitable for development of any kind."

SUSPECTED PROBLEMS

In essence the site remains of concern due to the oil deposits. Even after forty years, the effects of the oil dump are still visible.

FIELD RESULTS

Sunset Park is geographically divided into two distinct elevations with the northerly two-thirds at a higher elevation than the southerly one-third. These areas are used for recreational/athletic activities. At the south perimeter of hte park encroaches the marsh/swamp lands of Tub Lake. Access to Tub Lake is restricted by fences and "warning/danger" signs.

The Phase I field evaluation of the former landfill at Sunset Park was conducted on October 18, 1984. Sixteen bore holes were examined for methane and non-specific trace gas levels (Table XL). Methane was observed at low concentrations (trace to 0.4%) from all test holes located within the northerly

two-thirds of the park. Higher levels within the explosive range for methane gas were observed from four to six bore holes located within the southerly third of the site.

Non-specific trace gas levels were all non-significant with the exception of a 2.8 ppm level relative to ambient air concentrations at bore hole #2.

Surface water was sampled from a roadside stream at the southeastern corner of the fill and from the marsh water of Tub Lake. These data are presented in Table XLI. The test parameters did not indicate leachate contamination.

FIGURE 22
SUNSET PARK ABANDONED LANDFILL

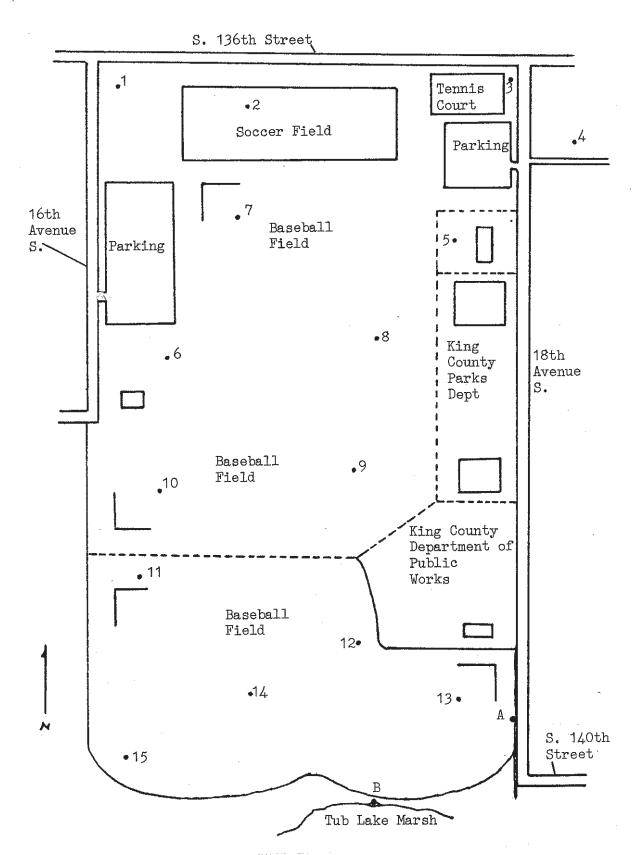


TABLE XLI
METHANE AND TRACE GAS CONCENTRATIONS
SUNSET PARK ABANDONED LANDFILL

Site	Methane (%)	Trace Gas (ppm)*
1	Trace	0
2	Trace	2.8
3	0.2	0.1
4	Trace	0
5	0.4	0
6	Trace	0
7	Trace	0
8	Trace	0
9	Trace	0
. 10	Trace	0
11	Trace	0
12	12	0
13	14	0
14	15	-0.1
15	10	0
16	1.2	0

^{*} Reading represents change from ambient air level

TABLE XLII

SURFACE WATER PARAMETERS SUNSET PARK ABANDONED LANDFILL

	Site A (1)	Site B (2)
рН	6.1	6.2
Temperature $^{\circ}$ C	9.8	14
Dissolved	3.9	10.1
Electrolytic Conductivity m v /cm	0.3	0.5
Turbidity	7	4

⁽¹⁾ Flowing water from road ditch(2) Stagnant water from swamp

TUKWILA

An old abandoned landfill exists at the end of 62nd Avenue South at the turn of South 153rd in Tukwila.

The site is located in the middle of a residential area but at the time of its operation, there was very little housing on the hill. Tukwila was an agricultural region and the landfill is remembered as a disposal site for household trash and garbage. Broken bottles and waste from the dairy which operated in the valley prior to the 1960's were also deposited there. It is thought that the site was discontinued in the mid-1940's. No documents on the site have been found.

FIELD RESULTS

This former landfill was tested for methane and trace-gas levels on October 29, 1984. Because of its small size, only three bore holes were placed for testing. Methane gas was observed only at trace levels indicating the stability of the site. Trace gas levels wer not observed above ambient air levels (Table XLII).

One water sample was retrieved from the swamp adjacent to the former fill. Water stagnation was indicated by a low dissolved oxygen level (1.2 ppm) and a relatively high electrolytic conductivity reading (0.5 milliohm/em). These data appear in Table XLIII. Leachate contamination was not indicated.

FIGURE 23
TUKWILA ABANDONED LANDFILL

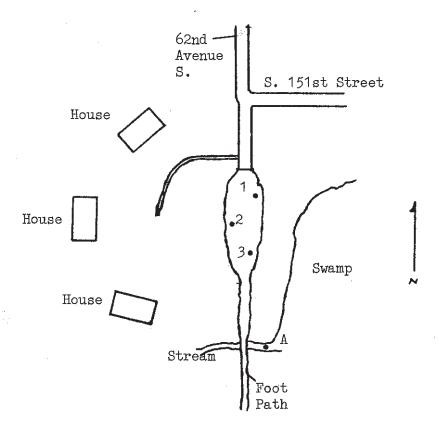


TABLE XLIII METHANE AND TRACE GAS CONCENTRATIONS TUKWILA ABANDONED LANDFILL

<u>Site</u>	Methane (%)	<u>Trace Gas (ppm)</u> *
1	Trace	0
2	Trace	0
3	Trace	0

^{*} Reading represents change from ambient air level

TABLE XLIV SURFACE WATER PARAMETERS TUKWILA ABANDONED LANDFILL

	Site A (1)
рН	6.0
Temperature	7.6
Dissolved	1.2
Electrolytic	0.5
Turbidity	17

⁽¹⁾ Marsh/swamp water

VASHON ISLAND

Across the road from the existing Vashon Island disposal site is the location of the original Island garbage dump.

No documents regarding the original site were located. It is believed that this site closed over forty years ago. It is remembered primarily as a site where residents dumped waste "off the bank." Age and location make it an unlikely problem area.

FIELD RESULTS

This site was tested on November 17, 1984 for methane and trace gases. These data are presented in Table XLIV. Methane levels ranged from betwen 0% to trace in the nine bore holes tested indicating the age and stability of the site. Trace gas levels did not differ from the ambient air concentrations. No surface water was evident on or around the site. Of note, the wooded downhill slopes were littered with old bottles and decomposed household refuse.

FIGURE 24

VASHON ISLAND ABANDONED LANDFILL

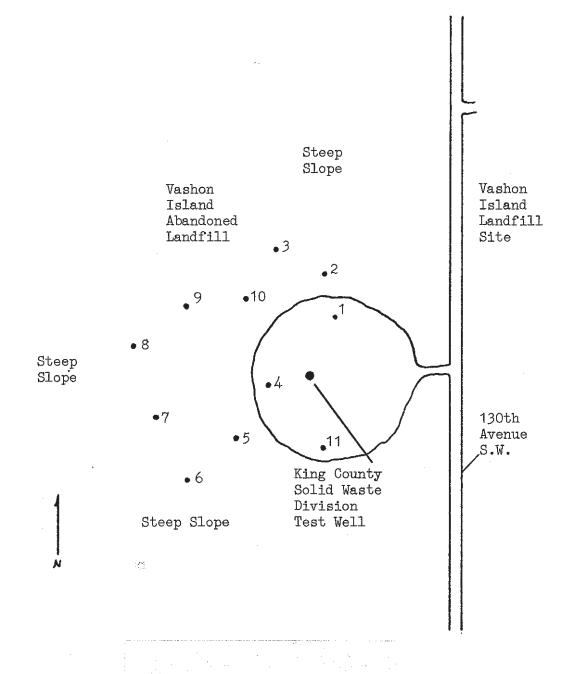


TABLE XLV

METHANE AND TRACE GAS CONCENTRATIONS

VASHON ISLAND ABANDONED LANDFILL

Site	Methane (%)	Trace Gas (ppm)*
1	Trace	0
2	0	0
3	Trace	0
4	Trace	0
5	0	0
6	0	0
7	0	0
8	0	0
9	0	0
10	Trace	0
11	0	0

^{*} Reading represents change from ambient air level

OTHER SITES

While the preliminary study presented in the previous sections focused on known abandoned sites identified for the Environmental Health Division, during the course of the study the existence of other abandoned sites came to light. One of these, Rotary Park in Auburn, discussed in the previous sections, had operated until the mid-1960's. However, there were no records readily available for the Auburn site. This was also true for the other newly identified sites. Their existence became known either by the passing mention of a private citizen or by seeing a reference made to a site buried in a report which focused on another site or issue. They are mentioned here to demonstrate that this study does focus on the generally known and accepted disposal sites in King County. But there indeed may be others that are lost in time and in the memories of those who lived and worked throughout the County in earlier years.

Some of the sites identified during the course of researching historical references were additional and unsuspected dump sites in Renton. During the 1930's and 1940's a garbage site was in operation in the area of the old Renton Coal Mining Company. It was in the canyon area on the east side of Interstate 405, currently seen as a blackberry-covered hill below Cedar Avenue South, slightly north of South 8th Street. All materials were dumped here and burning was common.

Another site from the 1940's and 1950's is remembered along the Cedar River industrial waterway near where it makes its entrance to Lake Washington, situated between the Boeing Renton Plant and the Renton Airport. The Cedar River tract and river park area are at that location today. Fires were common at this site which reportedly received "all kinds of stuff."

In addition, bottles were found about 1969 when trenches were dug at the site of the Renton City Hall parking lot. The one time existence of a dump at that location was a suggested explanation. However, early historical pictures at the Renton Historical Society's Museum at the Old Renton Fire Station show the presence of a glass and bottle manufacturing operation in the general vicinity. It is possible that the parking lot area may have been the site of that industry's waste.

Another site referenced in some reports is "Bellefield," a third Bellevue site, distinguished from the Factoria and Bellevue Air Field sites. The Bellefield site was located on the area of what is now the parking lot of the Bellefield Office Park located between the Mercer Slough and 112th Southeast in Bellevue. According to the 1974 document, Environmental Management for the Metropolitan Area, Part IV, Solid Waste, this fill accepted rubbish, street sweepings, tires, demolition waste and industrial waste. The King County Solid Waste Management Plan of 1976 reported the Bellefield site had problems with on-site and off-site surface runoff central, leachate and gas venting. It also did not meet the minimum functional standards required for either daily cover or for a clean and sanitary site. The site apparently closed sometime after 1976.

Sites as these have been developed and used to meet community's land use needs. No documented problems have arisen to suggest that they present any existing hazards, but the fact that they, at one time, were landfill sites should be common, not forgotten, consumer knowledge.



CHAPTER IV CONCLUSIONS AND RECOMMENDATIONS

	Page
Preface	135
AUBURN (M & R STREET SITE)	136
Auburn (Rotary Park Site)	136
Bow Lake	136
CARTON & BORTH	137
CORLISS LANDFILL	137
Eastgate Abandoned Landfill	138
Enumclaw	138
FACTORIA PIT (SUNSET RAVINE PARK)	139
FALL CITY ABANDONED LANDFILL	139
H.H. OLESON	139
Ноиднтом	139
Kent Abandoned Landfill (Mill Creek Canyon Park)	143
McMicken Heights	140
North Bend	140
Pacific City	141
Puyallup/Kit Corner	141
Redondo Pit	141
Renton Highlands	142
RENTON JUNCTION (MONSTER ROAD)	142
Skykomish	142
Sunset Park	143
Pukwila	143
VASHON	144

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PREFACE

The Conclusions and Recommendations that follow represent the opinion of the Seattle-King County Department of Public Health. Statements are based on the field and historic data presented earlier in this study.

The Seattle-King County Department of Public Health will submit a budget proposal to the King County Budget Office for a follow-up study of the eight abandoned landfill sites recommended for soil and/or water sampling. This project, if funded, will be conducted by the Health Department in 1986.

AUBURN ABANDONED LANDFILL (M & R STREET SITE)

This site revealed only slight readings of methane and non-specific organic/inorganic off-gasing. No surface water was available for sampling. Settling of streets, sidewalks and building foundations was readily apparent at the site. No significant environmental health problems were observed at the site and no further study is recommended.

AUBURN ABANDONED LANDFILL (ROTARY PARK)

This site exhibited elevated off-gasing of non-specific organic/inorganic gases concentrated in the eastern half of the site. Methane off-gasing was only noted at trace levels. No surface water was readily available for sampling.

It is recommended that soil and water samples be taken and analyzed for primary organics and inorganics that may be toxic to humans.

BOW LAKE ABANDONED LANDFILL

This site exhibited pockets of off-gasing of both methane and non-specific organics/inorganics. The elevated gas levels are consistent with the general type and volume of waste taken at the former landfill. Leachate was observed draining down and off the southeast portion of the property.

It is recommended:

- 1. The above noted leachate be intercepted and managed in an environmentally sound manner within the confines of the County transfer station property so as to prevent air, soil, and/or water contamination.
- 2. No further building construction take place on the property until the site has been stabilized.

3. That soil and water samples be taken and analyzed for primary organics and inorganics that may be toxic to humans.

CARTON & BORTH ABANDONED LANDFILL

This site demonstrated virtually no evidence of non-specific organic/inorganic off-gasing and slight evidence of methane generation within the center of
the site.

It is recommended:

- That methane monitoring be periodically conducted in proximity to the commercial structures adjacent to and west of the former landfill to insure that methane migration is not occurring in the building substructures.

CORLISS LANDFILL (McCORMICK PARK)

This site exhibited off-gasing of both methane and non-specific organics/inorganics. Explosive or greater levels of methane were found clustered through the central portions of the former fill. Water samples taken both up and downstream of the site in Thornton Creek revealed no evidence of leachate.

It is recommended:

- 1. No building construction take place on the property until the site has been stabilized.
- 2. Gas flaring or similar technology be utilized to reduce high concentrations of trapped methane within the fill.
- 3. Soil and water samples be taken and analyzed for priority organics and inorganics that may be toxic to humans.

EASTGATE ABANDONED LANDFILL

This site includes the old Bellevue Airfield and a small portion of Bellevue School District property. Little evidence of non-specific organic/inorganic off-gasing was observed. Methane gas was detected both on the former landfill and migrating to the east/northeast of the landfill. Leachate has been historically observed leaving the site via a surface ditch and draining into Phantom Lake.

It is recommended:

- 1. Gas flaring or similar technology be utilized to reduce high concentrations of trapped methane within the fill.
- 2. No further building construction take place until the landfill site has been stabilized.
- 3. Boeing continue methane migration monitoring, including specific checks of buildings in the path of potential migration.
- 4. The above noted leachate be intercepted and managed in an environmentally sound manner within the confines of the property so as to prevent air, soil, and/or water contamination.

ENUMCLAW ABANDONED LANDFILL

This site exhibited trace levels of methane off-gasing. The down gradient surface stream did not reveal evidence of leachate. Given the field data and the age of the former landfill, no environmental health problems are evidenced and no further study is warranted.

FACTORIA PIT (SUNSET RAVINE PARK)

This site demonstrated little evidence to suggest an environmental health risk. Since surface water revealed no evidence of leachate, and off-gasing was insignificant, no further study is warranted.

FALL CITY ABANDONED LANDFILL

This site revealed no evidence of methane off-gasing and one slightly elevated non-specific organic/inorganic reading. The water sample did not reveal high conductivity but the water had an oily appearance.

No significant environmental health problems were observed at the site and no further study is warranted.

H.H. OLESON

Methane and trace gas levels were observed at low levels throughout the site.

The off-site water sample tested did not indicate a leachate contamination problem.

Given the depth of the fill and its relatively recent deposit, further ground settling may be expected. It is therefore recommended that no building construction take place over the fill until the property has settled.

HOUGHTON ABANDONED LANDFILL

This site exhibited elevated levels of methane off-gasing particularly in the southwest section and only low levels of non-specific organic/inorganic gases were observed. Obvious signs of leachate were not observed during site inspections. However, leachate problems have historically been encountered.

It is recommended:

- 1. No additional building construction take place on the landfill until it has stabilized.
- 2. Soil and water samples be taken and analyzed for priority organics and inorganics that may be toxic to humans.
- 3. Gas flaring or similar technology be utilized to reduce high concentrations of trapped methane within the fill.

KENT ABANDONED LANDFILL (MILL CREEK CANYON PARK)

This site revealed no evidence of non-specific organic/inorganic off-gasing and very little evidence of methane off-gasing. The lack of significant environmental health problems at the site make further study unnecessary.

McMICKEN HEIGHTS ABANDONED LANDFILL

This site revealed no evidence of non-specific organic/inorganic off-gasing and very little evidence of methane off-gasing. No significant environmental health problems were observed at the site and it is recommended that no further study be done.

NORTH BEND ABANDONED LANDFILL

This site demonstrated very little evidence to suggest an environmental health risk. Only one methane and one non-specific organic/inorganic reading were even slightly positive. No further study seems warranted.

PACIFIC ABANDONED LANDFILL (PARK SITE)

This site revealed low level readings of methane and non-specific organic/inorganic off-gasing. The water sample from the adjacent White River did not indicate the presence of leachate. No significant environmental health problems were observed at the site and no further study is warranted.

PUYALLUP/KIT CORNER ABANDONED LANDFILL

This site demonstrated consistently elevated levels of methane off-gasing. Although both surface water samples failed to detect leachate parameters. A stained half round corrugated pipe was observed draining from the landfill area into the freeway storm drain.

- It is recommended that:
- 1. No building construction take place until the site has been stabilized.
- 2. Gas flaring or similar technology be utilized to reduce high concentrations of trapped methane within the fill.
- 3. Soil and surface/ground water samples be taken and analyzed for priority organics and inorganics that may be toxic to humans.

REDONDO PIT (SACAJAWEA COUNTY PARK)

This site exhibited elevated off-gasing of non-specific organic/inorganic gases, with the peak level observed in the middle terraced portion of the park. No corresponding significant levels of methane gas were noted at the site. This finding is consistent with the understanding that the site was utilized as a waste oil disposal site by the Navy and the Port of Seattle during and before World War II. No leachate was detected in the storm waste water system utilized by the park.

It is recommended that:

- Soil samples be taken and analyzed for priority organics and inorganics that may be toxic to humans.

RENTON HIGHLANDS ABANDONED LANDFILL

This site revealed no evidence of non-specific organic/inorganic off-gasing and only occasional traces of methane off-gasing. Mt. Olivet Creek was sampled and revealed no evidence of leachate. No significant environmental health problems were observed at the site and recommend that no further study is warranted.

RENTON JUNCTION ABANDONED LANDFILL

This site demonstrated relatively high levels of methane off-gasing with five of the seven test holes being within or above the explosive range. No evidence of non-specific organic/inorganic gases were observed exceeding background levels.

It is recommended:

- 1. That all existing building construction be properly vented and periodically monitored to insure that methane is not accumulating in any substructures.
- 2. That no further building construction take place over the former landfill site until it has been stabilized.

SKYKOMISH ABANDONED LANDFILL

This site revealed little evidence of methane or non-specific organic/inorganic off-gasing. No significant environmental health problems were observed at the site and no further study is warranted.

SUNSET PARK ABANDONED LANDFILL

This site exhibited active off-gasing of methane in the far southern portion of the park immediately adjacent to the Tub Lake property owned by the Port of Seattle. The non-specific organic/inorganic gases were at background levels in all but three samples and hence were not significant. The surface water sample at the border of the park and the lake property exhibited as slightly elevated specific conductivity.

It is recommended that:

- 1. No building construction take place at the southern portion of the park property until the site has been stabilized.
- 2. Water and soil sampling be conducted on Port of Seattle property (Tub Lake) and analyzed for priority organics and inorganics that may be toxic to humans.

TUKWILA ABANDONED LANDFILL

This site revealed no evidence of non-specific organic/inorganic off-gasing and very little evidence of methane off-gasing. Water sample parameters raise questions regarding the quality of on-site water and it is recommended:

- That surface water samples be taken and analyzed for priority organics and inorganics that may be toxic to humans.

VASHON ISLAND ABANDONED LANDFILL

This site exhibited no significant evidence of methane or non-specific organic/inorganic off-gasing. The monitoring well of King County Solid Waste division located on the former landfill was most recently sampled on September 12, 1984. The results indicate elevated conductivity but no corresponding high levels of heavy metals. No significant environmental health problems exist at the site and it is recommended that no further study is warranted.



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GLOSSARY

- 1. Building Construction: Refers to a dwelling or habitable structure.
- 2. <u>Stabilized</u>: Refers to the cessation or appropriate control of methane and trace gas generation.

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Personal Discussions With:

Lauren Anderson, City of Bellevue Maxine Anderson, City of Tukwila Leonard Chapman, City of Auburn Hayes Evans, Seattle-King County Department of Public Health and King County Public Works (Retired) Robert Howell, Tacoma-Pierce County Health Department Robert Hughes, City of Kent Bob Jacobs, King County Parks Department Jack Lange, City of Auburn (Retired) Don Mosley, City of Enumclaw Municipality of Metropolitan Seattle (METRO), Water Quality Section Ron Nelson, City of Renton Pat Nevins, City of Auburn Pat O'Flagherty, JRB Associates Leon Owen, King County Public Works Doug Pierce, Tacoma-Pierce County Health Department Norm Regan, City of Enumclaw Byron Sneva, City of Tukwila John Throap, Auburn private citizen Ernie Tonda, Renton Historical Society A.K. VanDusen, King County Public Works (Retired) Don Wickstrom, City of Kent Paul Wolfe, Auburn private citizen

APPENDIX C

Memorandum: Stormwater Quality Sampling Results from the Pacific Right Bank Wetland





Water and Land Resources Division

Memorandum

Department of Natural Resources and Parks King Street Center 201 South Jackson Street, Suite 600 Seattle, WA 98104-3855

206-296-6519 Fax 206-296-0192 TTY Relay: 711

February 18, 2015

TO: Chris Brummer, Senior Engineer, White River Basin, River and Floodplain Management Section (RFMS)

CC: Jeanne Stypula, Supervising Engineer, White River and Technical Services, RFMS

FM: Sarah McCarthy, Senior Ecologist, White River Basin, RFMS Sevin Bilir, Environmental Scientist III, STSS

RE: Stormwater quality sampling results from the Pacific Right Bank Wetland

Background and Purpose

In 2010 and 2011, we collected stormwater samples from three locations in the vicinity of Pacific City Park, landward of the Pacific City Park Levee (Figure 1). This work was done to support a feasibility study for the future levee setback project on the right bank of the Lower White River. The purpose of this stormwater sampling was twofold:

- 1. Determine whether the quality of stormwater entering the wetland poses a concern for aquatic life, and
- 2. Determine whether analytes detected in the wetland indicate impacts from an abandoned dumpsite located in the vicinity of Pacific City Park that was active between ca. 1920 through the late-1960s

Methods

Sample locations were selected by RFMS personnel to identify potential impacts from offsite sources (WRLEV1) and to evaluate water quality conditions in two wetland habitats (WRLEV2 and WRLEV3) within the study area. WRLEV1 is located in a stormwater drainage ditch receiving water from residential properties to the west and north of the wetland. The northern wetland sampling location (WRLEV2) is located south of and adjacent to the Pacific City Park, approximately 300 feet downstream of where the drainage ditch discharges into the wetland. The southern wetland sampling location (WRLEV3) is located in the wetland approximately 1,000 feet south of WRLEV2.

Field parameters, including dissolved oxygen (DO), pH, water temperature, turbidity and conductivity, were measured prior to sample collection at all three sample locations during storm events on December 8, 2010, February 28, 2011, and March 10, 2011. A baseline (non-storm) collection was completed on January 04, 2011 at all three sampling locations.

Field parameters were measured, and surface water samples were collected using standard practice field sampling protocols and equipment, as described in *Standard Operating Procedure for Sampling Methods for Stream and River Water* (SOP #214v3). Samples were transported to the King County Environmental Laboratory (KCEL) in Seattle for analysis of a suite of parameters (Table 1).

Results

Field water quality parameter results and laboratory detected analytes are presented in Tables 2 and 3, respectfully. Laboratory results from the KCEL may be qualified as less than the method detection limit (MDL) or the reporting detection limit (RDL). Values less than the MDL indicate a parameter is not present in the sample above this level. Values greater than the MDL, but less than the RDL indicate confidence that the parameter is present, but at low enough levels to decrease reliability of the numerical result. Values greater than the RDL are considered reliable after undergoing laboratory Quality Assurance/Quality Control (QA/QC). For most of results, the data passed all KCEL internal QA/QC checks for accuracy and completeness with a few exceptions related to exceedances of preservation time limits for TKN, and manganese, semi-volatile organics and zinc detected in blanks.

Discussion

Results from this initial testing should be shared with the consultant preparing the Phase I Environmental Site Assessment (ESA) for the property. This will assist them in providing recommendations for a Phase II ESA for the study area.

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Attachments

Figure 1: Water Quality Sampling Locations

Table 1: Laboratory Methods and Analytes Tested, 2010 and 2011.

Table 2. Field Water Quality Results, 2010 and 2011.

Table 3. Reported Detected Analytes, 2010 and 2011

Appendices: KCEL Analytical Reports attached in FINAL document



Table 1. Laboratory Methods and Analytes Tested, 2010 and 2011.

	Baseline	Storm Events				
	1/4/2011	12/8/2010	2/28/2011	3/10/2011		
CV EPA 351.2						
Total Kjeldahl Nitrogen	\checkmark	\checkmark	\checkmark	\checkmark		
CV SM4500-NO3-F						
Nitrite + Nitrate Nitrogen	\checkmark	\checkmark	$\sqrt{}$	\checkmark		
CV SM4500-P-B,F						
Total Phosphorus	\checkmark	\checkmark	$\sqrt{}$	\checkmark		
CV SM4500-P-F						
Orthophosphate Phosphorus	$\sqrt{}$	\checkmark	$\sqrt{}$	$\sqrt{}$		
MT EPA 200.8*SW846 6020A						
Cadmium, Dissolved, ICP-MS	$\sqrt{}$	\checkmark	$\sqrt{}$	$\sqrt{}$		
Cadmium, Total, ICP-MS	$\sqrt{}$	\checkmark	$\sqrt{}$	$\sqrt{}$		
Calcium, Total, ICP-MS	$\sqrt{}$	\checkmark	$\sqrt{}$	$\sqrt{}$		
Chromium, Dissolved, ICP-MS	\checkmark	\checkmark	$\sqrt{}$	\checkmark		
Chromium, Total, ICP-MS	$\sqrt{}$	$\sqrt{}$	$\sqrt{}$			
Copper, Dissolved, ICP-MS	\checkmark	\checkmark	$\sqrt{}$	$\sqrt{}$		
Copper, Total, ICP-MS	$\sqrt{}$	$\sqrt{}$	$\sqrt{}$			
Iron, Total, ICP-MS				$\sqrt{}$		
Lead, Dissolved, ICP-MS	$\sqrt{}$	$\sqrt{}$	$\sqrt{}$			
Lead, Total, ICP-MS	$\sqrt{}$	$\sqrt{}$	$\sqrt{}$			
Magnesium, Total, ICP-MS	$\sqrt{}$	\checkmark	$\sqrt{}$	$\sqrt{}$		
Zinc, Dissolved, ICP-MS	$\sqrt{}$	$\sqrt{}$	$\sqrt{}$			
Zinc, Total, ICP-MS	$\sqrt{}$	$\sqrt{}$	$\sqrt{}$			
MT EPA 200.8/SW846 6020A*SM2340B						
Hardness, Calc	\checkmark	\checkmark	$\sqrt{}$	\checkmark		
OR SW846 3520C*SW846 8270D SIM						
Chlorpyrifos	$\sqrt{}$	\checkmark				
Diazinon	\checkmark	\checkmark				
Disulfoton	$\sqrt{}$	\checkmark				
Malathion	$\sqrt{}$	\checkmark				
Parathion-Ethyl	$\sqrt{}$	\checkmark				
Parathion-Methyl	\checkmark	\checkmark				
Phorate	$\sqrt{}$	\checkmark				
OR SW846 3535A/8151A*8270D SIM						
2,4,5-T	$\sqrt{}$	\checkmark				
2,4,5-TP (Silvex)	$\sqrt{}$	\checkmark				
2,4-D	\checkmark	$\sqrt{}$				
2,4-DB	$\sqrt{}$	\checkmark				
Dichloroprop	$\sqrt{}$	\checkmark				
Dinoseb	\checkmark	\checkmark				
MCPA	\checkmark	\checkmark				
MCPP	\checkmark	\checkmark				

Table 1. Laboratory Methods and Analytes Test, 2010 and 2011. (continued)

	Baseline	Storm Events				
	1/4/2011	12/8/2010	2/28/2011	3/10/2011		
OR SW846 5030B*SW846 8260C						
1,1,1,2-Tetrachloroethane		\checkmark				
1,1,1-Trichloroethane	$\sqrt{}$	$\sqrt{}$				
1,1,2,2-Tetrachloroethane	$\sqrt{}$	$\sqrt{}$				
1,1,2-Trichloroethane	$\sqrt{}$	$\sqrt{}$				
1,1,2-Trichloroethylene		$\sqrt{}$				
1,1-Dichloroethane		$\sqrt{}$				
1,1-Dichloroethylene		$\sqrt{}$				
1,2,3-Trichloropropane		$\sqrt{}$				
1,2-Dibromo-3-chloropropane		$\sqrt{}$				
1,2-Dibromoethane		$\sqrt{}$				
1,2-Dichlorobenzene		$\sqrt{}$				
1,2-Dichloroethane		$\sqrt{}$				
1,2-Dichloropropane	V	, V				
1,4-Dichlorobenzene	, v	, J				
2-Butanone (MEK)	, J	J				
2-Hexanone	J	J				
4-Methyl-2-Pentanone (MIBK)	V	J				
Acetone	, J	J				
Acrylonitrile	J	J				
Benzene	J	J				
Bromochloromethane	J.	J				
Bromodichloromethane	1	1				
Bromoform	1	2/				
Bromomethane	1	2/				
Carbon Disulfide	\ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \	2/				
Carbon Tetrachloride	2/	2/				
Chlorobenzene	2/	2/				
Chlorodibromomethane	N al	N al				
	N al	N al				
Chloroethane	N I	.,				
Chloroform	N I	\ .1				
Chloromethane	N I	N I				
Cis-1,2-Dichloroethylene	N I	N I				
Cis-1,3-Dichloropropene	N /	N I				
Dibromomethane	N I	N I				
Dichlorodifluoromethane	N ,	V				
Ethylbenzene	V	V				
Methylene Chloride	V	V				
O-Xylene	V	V				
Styrene	V	√				
Tetrachloroethylene	V	√				
Toluene	√,	√,				
Trans-1,2-Dichloroethylene	√,	√,				
Trans-1,3-Dichloropropene	√	√ ,				
Trichlorofluoromethane	√,	√,				
Vinyl Acetate	√,	√,				
Vinyl Chloride	√	√				

Table 1. Laboratory Methods and Analytes Tested, 2010 and 2011. (continued)

	Baseline	Storm Events				
	1/4/2011	12/8/2010	2/28/2011	3/10/2011		
OR SW846 3520C*SW846 8270D						
1,2,4-Trichlorobenzene		\checkmark				
1,2-Dichlorobenzene	$\sqrt{}$	\checkmark				
1,2-Diphenylhydrazine	$\sqrt{}$	\checkmark				
1,3-Dichlorobenzene	$\sqrt{}$	\checkmark				
1,4-Dichlorobenzene	$\sqrt{}$	$\sqrt{}$				
2,4,5-Trichlorophenol	$\sqrt{}$	$\sqrt{}$				
2,4,6-Trichlorophenol	$\sqrt{}$	\checkmark				
2,4-Dichlorophenol	$\sqrt{}$	\checkmark				
2,4-Dimethylphenol	$\sqrt{}$	\checkmark				
2,4-Dinitrophenol	$\sqrt{}$	\checkmark				
2,4-Dinitrotoluene	$\sqrt{}$	\checkmark				
2,6-Dinitrotoluene	\checkmark	\checkmark				
2-Chloronaphthalene	$\sqrt{}$	\checkmark				
2-Chlorophenol	\checkmark	\checkmark				
2-Methylnaphthalene		$\sqrt{}$				
2-Methylphenol		V				
2-Nitroaniline		$\sqrt{}$				
2-Nitrophenol	V	$\sqrt{}$				
3,3'-Dichlorobenzidine		V				
3-Methylphenol		$\sqrt{}$				
3-Nitroaniline	V	√ √				
4,6-Dinitro-O-Cresol	, ,	, V				
4-Bromophenyl Phenyl Ether	V	√ √				
4-Chloro-3-Methylphenol	V	V				
4-Chloroaniline	V	√ √				
4-Chlorophenyl Phenyl Ether	V	√ √				
4-Methylphenol	, v	V				
4-Nitroaniline	V	√ √				
4-Nitrophenol	V	$\sqrt{}$				
Acenaphthene	V	√ √				
Acenaphthylene	, ,	, V				
Aniline	V	√ √				
Anthracene	V	√ √				
Benzo(a)anthracene	√ √	√ √				
Benzo(a)pyrene		$\sqrt{}$				
Benzo(b)fluoranthene	V	√ √				
Benzo(g,h,i)perylene	V	√ √				
Benzo(k)fluoranthene	V	V				
Benzoic Acid	, V	, √				
Benzyl Alcohol	, V	, V				
Benzyl Butyl Phthalate	, √	, V				
Bis(2-Chloroethoxy)Methane	, √	, V				
Bis(2-Chloroethyl)Ether	, V	į				
Bis(2-Chloroisopropyl)Ether	, √	, V				
Bis(2-Ethylhexyl)Phthalate	, V	, V				
Dista Entry in Chyrji Titriaiate	V	٧				

Table 1. Laboratory Methods and Analytes Tested, 2010 and 2011. (continued)

	Baseline	Storm Events				
	1/4/2011	12/8/2010	2/28/2011	3/10/2011		
OR SW846 3520C*SW846 8270D						
Caffeine	\checkmark	$\sqrt{}$				
Carbazole	\checkmark	$\sqrt{}$				
Chrysene	\checkmark	$\sqrt{}$				
Coprostanol	$\sqrt{}$	$\sqrt{}$				
Dibenzo(a,h)anthracene	\checkmark	$\sqrt{}$				
Dibenzofuran	\checkmark	$\sqrt{}$				
Diethyl Phthalate	\checkmark	$\sqrt{}$				
Dimethyl Phthalate	\checkmark	$\sqrt{}$				
Di-N-Butyl Phthalate	\checkmark	$\sqrt{}$				
Di-N-Octyl Phthalate	\checkmark	$\sqrt{}$				
Fluoranthene	\checkmark	$\sqrt{}$				
Fluorene	\checkmark	$\sqrt{}$				
Hexachlorobenzene	\checkmark	$\sqrt{}$				
Hexachlorobutadiene	\checkmark	$\sqrt{}$				
Hexachlorocyclopentadiene	\checkmark	$\sqrt{}$				
Hexachloroethane	\checkmark	$\sqrt{}$				
Indeno(1,2,3-Cd)Pyrene	\checkmark	$\sqrt{}$				
Isophorone	\checkmark	$\sqrt{}$				
Naphthalene	$\sqrt{}$	$\sqrt{}$				
Nitrobenzene	\checkmark	$\sqrt{}$				
N-Nitrosodimethylamine	\checkmark	$\sqrt{}$				
N-Nitrosodi-N-Propylamine	\checkmark	$\sqrt{}$				
N-Nitrosodiphenylamine	\checkmark	$\sqrt{}$				
Pentachlorophenol	\checkmark	$\sqrt{}$				
Phenanthrene	\checkmark	$\sqrt{}$				
Phenol	\checkmark	$\sqrt{}$				
Pyrene	\checkmark	$\sqrt{}$				
Pyridine	√	√				

Note: Baseline indicates non-storm collection

Table 2. Field Water Quality Results, 2010 and 2011.

	Baseline		Storm Event	s							
	1/4/2011*	12/8/2010**	2/28/2011	3/10/2011							
WRLEV1 - Drainage Ditch											
Dissolved oxygen	5.86	3.45	3.14	0.7							
pH	6.5	6.51	6.48	6.52							
Temperature	4.54	8.63	5.11	9.18							
Turbidity	2.22	10.2	12.5	7.43							
Conductivity	408.9	173.4	238.9	356.3							
WRLEV2 - Upstream Wetland											
Dissolved oxygen	1.66	0.95	1.81	2.7							
pH	6.35	6.51	6.38	6.41							
Temperature	4.89	7.43	5.97	10.1							
Turbidity	12.3	2.2	2.41	5.1							
Conductivity	102.3	NC	NC	146							
	WRLEV3 - Do	ownstream W	etland								
Dissolved oxygen	6.01	4.91	5.11	3.63							
pH	6.54	6.75	6.66	6.62							
Temperature	3.38	8.71	4.45	8.29							
Turbidity	30.6	44.8	41.1	4.95							
Conductivity	288.3	NC	NC	289.4							

Note:

NC not collected

*WRLEV1 Iron oxide odor, orange color and sheen present 1/4/2011

Note: Baseline indicates non-storm collection

^{*}WRLEV2 Sulfide odor 1/4/2011

^{**}WRLEV1 Algae and sheen present, low flow

^{**}WRLEV2 Sulfur odor, algae present

^{**}WRLEV3 Orange color, suspended algae, some flow

Table 3. Reported Detected Analytes, 2010 and 2011

			Total Kjeldahl Nitrogen	Nitrite + Nitrate Nitrogen	Total Phosphorus	Orthophosphate Phosphorus	Calcium, Total, ICP-MS	Chromium, Dissolved, ICP- MS	Chromium, Total, ICP-MS	Copper, Dissolved, ICP- MS	Copper, Total, ICP-MS	Iron, Total, ICP- MS	Lead, Dissolved, ICP-MS	Lead, Total, ICP- MS	Magnesium, Total, ICP-MS	Disso CP-MS	Zinc, Total, ICP- MS	Hardness, Calc	2-Methylphenol	4-Methylphenol	Acenaphthene	Acenaphthylene	Benzoic Acid	Benzyl Alcohol	Benzyl Butyl Phthalate	Bis(2- Ethylhexyl)Phtha late	Diethyl Phthalate	Di-N-Butyl Phthalate	Naphthalene	Acetone	Toluene	Vinyl Chloride
		Units	mg/L	mg/L	mg/L	mg/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	mg CaCO3/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Event	Date																															
WRLEV1 -	Drainage D																															
Baseline	1/4/2011	Result	1.58		0.134	0.049	42600	1.12	1.08	0.41	<mdl< td=""><td></td><td></td><td></td><td>8500</td><td>4.89</td><td>5.2</td><td>141</td><td></td><td></td><td></td><td></td><td>1.9</td><td></td><td></td><td>0.26</td><td></td><td></td><td>0.0241</td><td></td><td></td><td>0.0371</td></mdl<>				8500	4.89	5.2	141					1.9			0.26			0.0241			0.0371
		Qual	SH	<mdl< td=""><td></td><td></td><td></td><td></td><td></td><td><rdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td><mdl< td=""><td></td><td></td><td></td><td></td><td><mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td></td><td><mdl< td=""><td><mdl< td=""><td><rdl,b< td=""><td><mdl< td=""><td>В</td><td></td><td><mdl< td=""><td><mdl< td=""><td></td></mdl<></td></mdl<></td></mdl<></td></rdl,b<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<></td></mdl<>						<rdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td><mdl< td=""><td></td><td></td><td></td><td></td><td><mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td></td><td><mdl< td=""><td><mdl< td=""><td><rdl,b< td=""><td><mdl< td=""><td>В</td><td></td><td><mdl< td=""><td><mdl< td=""><td></td></mdl<></td></mdl<></td></mdl<></td></rdl,b<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<>	<mdl< td=""><td></td><td><mdl< td=""><td><mdl< td=""><td></td><td></td><td></td><td></td><td><mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td></td><td><mdl< td=""><td><mdl< td=""><td><rdl,b< td=""><td><mdl< td=""><td>В</td><td></td><td><mdl< td=""><td><mdl< td=""><td></td></mdl<></td></mdl<></td></mdl<></td></rdl,b<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<>		<mdl< td=""><td><mdl< td=""><td></td><td></td><td></td><td></td><td><mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td></td><td><mdl< td=""><td><mdl< td=""><td><rdl,b< td=""><td><mdl< td=""><td>В</td><td></td><td><mdl< td=""><td><mdl< td=""><td></td></mdl<></td></mdl<></td></mdl<></td></rdl,b<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<>	<mdl< td=""><td></td><td></td><td></td><td></td><td><mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td></td><td><mdl< td=""><td><mdl< td=""><td><rdl,b< td=""><td><mdl< td=""><td>В</td><td></td><td><mdl< td=""><td><mdl< td=""><td></td></mdl<></td></mdl<></td></mdl<></td></rdl,b<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<>					<mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td></td><td><mdl< td=""><td><mdl< td=""><td><rdl,b< td=""><td><mdl< td=""><td>В</td><td></td><td><mdl< td=""><td><mdl< td=""><td></td></mdl<></td></mdl<></td></mdl<></td></rdl,b<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<>	<mdl< td=""><td></td><td><mdl< td=""><td></td><td><mdl< td=""><td><mdl< td=""><td><rdl,b< td=""><td><mdl< td=""><td>В</td><td></td><td><mdl< td=""><td><mdl< td=""><td></td></mdl<></td></mdl<></td></mdl<></td></rdl,b<></td></mdl<></td></mdl<></td></mdl<></td></mdl<>		<mdl< td=""><td></td><td><mdl< td=""><td><mdl< td=""><td><rdl,b< td=""><td><mdl< td=""><td>В</td><td></td><td><mdl< td=""><td><mdl< td=""><td></td></mdl<></td></mdl<></td></mdl<></td></rdl,b<></td></mdl<></td></mdl<></td></mdl<>		<mdl< td=""><td><mdl< td=""><td><rdl,b< td=""><td><mdl< td=""><td>В</td><td></td><td><mdl< td=""><td><mdl< td=""><td></td></mdl<></td></mdl<></td></mdl<></td></rdl,b<></td></mdl<></td></mdl<>	<mdl< td=""><td><rdl,b< td=""><td><mdl< td=""><td>В</td><td></td><td><mdl< td=""><td><mdl< td=""><td></td></mdl<></td></mdl<></td></mdl<></td></rdl,b<></td></mdl<>	<rdl,b< td=""><td><mdl< td=""><td>В</td><td></td><td><mdl< td=""><td><mdl< td=""><td></td></mdl<></td></mdl<></td></mdl<></td></rdl,b<>	<mdl< td=""><td>В</td><td></td><td><mdl< td=""><td><mdl< td=""><td></td></mdl<></td></mdl<></td></mdl<>	В		<mdl< td=""><td><mdl< td=""><td></td></mdl<></td></mdl<>	<mdl< td=""><td></td></mdl<>	
Storm	12/8/2010	Result	1.75		0.136	0.019	37900	0.74	8.0						7370	6.72	5.54	125			0.015		1		0.102	0.99		0.176	0.015			0.0308
		Qual		<mdl< td=""><td></td><td><rdl< td=""><td></td><td><rdl< td=""><td><rdl< td=""><td><mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td><mdl< td=""><td></td><td></td><td></td><td></td><td><mdl< td=""><td><mdl< td=""><td><rdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td></td><td></td><td><mdl< td=""><td>В</td><td><rdl< td=""><td></td><td><mdl< td=""><td></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<></td></rdl<></td></rdl<></td></mdl<>		<rdl< td=""><td></td><td><rdl< td=""><td><rdl< td=""><td><mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td><mdl< td=""><td></td><td></td><td></td><td></td><td><mdl< td=""><td><mdl< td=""><td><rdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td></td><td></td><td><mdl< td=""><td>В</td><td><rdl< td=""><td></td><td><mdl< td=""><td></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<></td></rdl<></td></rdl<>		<rdl< td=""><td><rdl< td=""><td><mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td><mdl< td=""><td></td><td></td><td></td><td></td><td><mdl< td=""><td><mdl< td=""><td><rdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td></td><td></td><td><mdl< td=""><td>В</td><td><rdl< td=""><td></td><td><mdl< td=""><td></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<></td></rdl<>	<rdl< td=""><td><mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td><mdl< td=""><td></td><td></td><td></td><td></td><td><mdl< td=""><td><mdl< td=""><td><rdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td></td><td></td><td><mdl< td=""><td>В</td><td><rdl< td=""><td></td><td><mdl< td=""><td></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<>	<mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td><mdl< td=""><td></td><td></td><td></td><td></td><td><mdl< td=""><td><mdl< td=""><td><rdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td></td><td></td><td><mdl< td=""><td>В</td><td><rdl< td=""><td></td><td><mdl< td=""><td></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<>	<mdl< td=""><td></td><td><mdl< td=""><td><mdl< td=""><td></td><td></td><td></td><td></td><td><mdl< td=""><td><mdl< td=""><td><rdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td></td><td></td><td><mdl< td=""><td>В</td><td><rdl< td=""><td></td><td><mdl< td=""><td></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<>		<mdl< td=""><td><mdl< td=""><td></td><td></td><td></td><td></td><td><mdl< td=""><td><mdl< td=""><td><rdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td></td><td></td><td><mdl< td=""><td>В</td><td><rdl< td=""><td></td><td><mdl< td=""><td></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<>	<mdl< td=""><td></td><td></td><td></td><td></td><td><mdl< td=""><td><mdl< td=""><td><rdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td></td><td></td><td><mdl< td=""><td>В</td><td><rdl< td=""><td></td><td><mdl< td=""><td></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<>					<mdl< td=""><td><mdl< td=""><td><rdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td></td><td></td><td><mdl< td=""><td>В</td><td><rdl< td=""><td></td><td><mdl< td=""><td></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<></td></mdl<></td></mdl<>	<mdl< td=""><td><rdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td></td><td></td><td><mdl< td=""><td>В</td><td><rdl< td=""><td></td><td><mdl< td=""><td></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<></td></mdl<>	<rdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td></td><td></td><td><mdl< td=""><td>В</td><td><rdl< td=""><td></td><td><mdl< td=""><td></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<>	<mdl< td=""><td></td><td><mdl< td=""><td></td><td></td><td><mdl< td=""><td>В</td><td><rdl< td=""><td></td><td><mdl< td=""><td></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<>		<mdl< td=""><td></td><td></td><td><mdl< td=""><td>В</td><td><rdl< td=""><td></td><td><mdl< td=""><td></td></mdl<></td></rdl<></td></mdl<></td></mdl<>			<mdl< td=""><td>В</td><td><rdl< td=""><td></td><td><mdl< td=""><td></td></mdl<></td></rdl<></td></mdl<>	В	<rdl< td=""><td></td><td><mdl< td=""><td></td></mdl<></td></rdl<>		<mdl< td=""><td></td></mdl<>	
Storm	2/28/2011	Result		0.012	0.132	0.0263	25600	0.44	0.5		0.55			0.12	5520	3.07	3.42	86.6	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
		Qual	SH	<rdl< td=""><td></td><td></td><td></td><td><rdl< td=""><td><rdl< td=""><td><mdl< td=""><td><rdl< td=""><td></td><td><mdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td></td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<></td></mdl<></td></rdl<></td></mdl<></td></rdl<></td></rdl<></td></rdl<>				<rdl< td=""><td><rdl< td=""><td><mdl< td=""><td><rdl< td=""><td></td><td><mdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td></td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<></td></mdl<></td></rdl<></td></mdl<></td></rdl<></td></rdl<>	<rdl< td=""><td><mdl< td=""><td><rdl< td=""><td></td><td><mdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td></td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<></td></mdl<></td></rdl<></td></mdl<></td></rdl<>	<mdl< td=""><td><rdl< td=""><td></td><td><mdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td></td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<></td></mdl<></td></rdl<></td></mdl<>	<rdl< td=""><td></td><td><mdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td></td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<></td></mdl<></td></rdl<>		<mdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td></td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<></td></mdl<>	<rdl< td=""><td></td><td></td><td></td><td></td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<>					NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Storm	3/10/2011	Result	0.772	0.025	0.134	0.0857	18700	0.35	0.43	0.45		6850		0.17	4330	3.17	3.35	64.6	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
		Qual	SH	<rdl< td=""><td></td><td></td><td></td><td><rdl< td=""><td><rdl< td=""><td><rdl< td=""><td><rdl< td=""><td></td><td><mdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td></td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<></td></mdl<></td></rdl<></td></rdl<></td></rdl<></td></rdl<></td></rdl<>				<rdl< td=""><td><rdl< td=""><td><rdl< td=""><td><rdl< td=""><td></td><td><mdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td></td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<></td></mdl<></td></rdl<></td></rdl<></td></rdl<></td></rdl<>	<rdl< td=""><td><rdl< td=""><td><rdl< td=""><td></td><td><mdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td></td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<></td></mdl<></td></rdl<></td></rdl<></td></rdl<>	<rdl< td=""><td><rdl< td=""><td></td><td><mdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td></td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<></td></mdl<></td></rdl<></td></rdl<>	<rdl< td=""><td></td><td><mdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td></td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<></td></mdl<></td></rdl<>		<mdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td></td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<></td></mdl<>	<rdl< td=""><td></td><td></td><td></td><td></td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<>					NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
WRLEV2 –	Upstream \																															
Baseline	1/4/2011	Result	0.323		0.142	0.0156	11900	0.33	0.37		0.67		0.13	0.27	2280	1.5	1.4	39	0.307	0.197			0.909	0.21	0.092	0.23		0.14	0.013		0.5	——
		Qual	SH	<mdl< td=""><td></td><td></td><td></td><td><rdl< td=""><td><rdl< td=""><td><mdl< td=""><td><rdl< td=""><td></td><td><rdl< td=""><td><rdl< td=""><td></td><td><rdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td><mdl< td=""><td><mdl< td=""><td></td><td></td><td><rdl< td=""><td><rdl,b< td=""><td><mdl< td=""><td>В</td><td><rdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""></mdl<></td></mdl<></td></rdl<></td></mdl<></td></rdl,b<></td></rdl<></td></mdl<></td></mdl<></td></rdl<></td></rdl<></td></rdl<></td></rdl<></td></rdl<></td></mdl<></td></rdl<></td></rdl<></td></mdl<>				<rdl< td=""><td><rdl< td=""><td><mdl< td=""><td><rdl< td=""><td></td><td><rdl< td=""><td><rdl< td=""><td></td><td><rdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td><mdl< td=""><td><mdl< td=""><td></td><td></td><td><rdl< td=""><td><rdl,b< td=""><td><mdl< td=""><td>В</td><td><rdl< td=""><td><mdl< td=""><td></td><td><mdl< 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Storm	12/8/2010	Result	0.31		0.115	0.00689	14300	0.26	0.32						2930	0.65	0.6	47.8	0.539	1.55			1.75	0.263		0.488	0.026	0.134			0.551	——
		Qual		<mdl< td=""><td></td><td></td><td>2.172</td><td><rdl< td=""><td><rdl< td=""><td><mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td><rdl< td=""><td></td><td><rdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td><mdl< td=""><td><mdl< td=""><td></td><td></td><td><mdl< td=""><td>B2</td><td><rdl< td=""><td>В</td><td><mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""></mdl<></td></mdl<></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<></td></rdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<></td></rdl<></td></mdl<>			2.172	<rdl< td=""><td><rdl< td=""><td><mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td><rdl< td=""><td></td><td><rdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td><mdl< td=""><td><mdl< td=""><td></td><td></td><td><mdl< td=""><td>B2</td><td><rdl< td=""><td>В</td><td><mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""></mdl<></td></mdl<></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<></td></rdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<></td></rdl<>	<rdl< td=""><td><mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td><rdl< td=""><td></td><td><rdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td><mdl< td=""><td><mdl< td=""><td></td><td></td><td><mdl< td=""><td>B2</td><td><rdl< td=""><td>В</td><td><mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""></mdl<></td></mdl<></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<></td></rdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<>	<mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""><td><rdl< td=""><td></td><td><rdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td><mdl< td=""><td><mdl< td=""><td></td><td></td><td><mdl< td=""><td>B2</td><td><rdl< td=""><td>В</td><td><mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""></mdl<></td></mdl<></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<></td></rdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<>	<mdl< td=""><td></td><td><mdl< td=""><td><rdl< td=""><td></td><td><rdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td><mdl< td=""><td><mdl< td=""><td></td><td></td><td><mdl< td=""><td>B2</td><td><rdl< td=""><td>В</td><td><mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""></mdl<></td></mdl<></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<></td></rdl<></td></rdl<></td></mdl<></td></mdl<>		<mdl< td=""><td><rdl< td=""><td></td><td><rdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td><mdl< td=""><td><mdl< td=""><td></td><td></td><td><mdl< td=""><td>B2</td><td><rdl< td=""><td>В</td><td><mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""></mdl<></td></mdl<></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<></td></rdl<></td></rdl<></td></mdl<>	<rdl< td=""><td></td><td><rdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td><mdl< td=""><td><mdl< td=""><td></td><td></td><td><mdl< td=""><td>B2</td><td><rdl< td=""><td>В</td><td><mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""></mdl<></td></mdl<></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<></td></rdl<></td></rdl<>		<rdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td><mdl< td=""><td><mdl< td=""><td></td><td></td><td><mdl< td=""><td>B2</td><td><rdl< td=""><td>В</td><td><mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""></mdl<></td></mdl<></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<></td></rdl<>	<rdl< td=""><td></td><td></td><td></td><td><mdl< td=""><td><mdl< td=""><td></td><td></td><td><mdl< td=""><td>B2</td><td><rdl< td=""><td>В</td><td><mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""></mdl<></td></mdl<></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<>				<mdl< td=""><td><mdl< td=""><td></td><td></td><td><mdl< td=""><td>B2</td><td><rdl< td=""><td>В</td><td><mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""></mdl<></td></mdl<></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<>	<mdl< td=""><td></td><td></td><td><mdl< td=""><td>B2</td><td><rdl< td=""><td>В</td><td><mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""></mdl<></td></mdl<></td></mdl<></td></rdl<></td></mdl<></td></mdl<>			<mdl< td=""><td>B2</td><td><rdl< td=""><td>В</td><td><mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""></mdl<></td></mdl<></td></mdl<></td></rdl<></td></mdl<>	B2	<rdl< td=""><td>В</td><td><mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""></mdl<></td></mdl<></td></mdl<></td></rdl<>	В	<mdl< td=""><td><mdl< td=""><td></td><td><mdl< td=""></mdl<></td></mdl<></td></mdl<>	<mdl< td=""><td></td><td><mdl< td=""></mdl<></td></mdl<>		<mdl< td=""></mdl<>
Storm	2/28/2011	Result			0.0699	0.0243	6450	0.2	0.25	.1.451	0.47		0.21	0.29	1440	. MDI	0.82	22	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
		Qual	SH	<mdl< td=""><td>0.0045</td><td>0.0054</td><td>7700</td><td><rdl< td=""><td><rdl< td=""><td><mdl< td=""><td><rdl< td=""><td>00.40</td><td><rdl< td=""><td><rdl< td=""><td>1000</td><td><mdl< td=""><td><rdl< td=""><td>00.0</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<></td></mdl<></td></rdl<></td></rdl<></td></rdl<></td></mdl<></td></rdl<></td></rdl<></td></mdl<>	0.0045	0.0054	7700	<rdl< td=""><td><rdl< td=""><td><mdl< td=""><td><rdl< td=""><td>00.40</td><td><rdl< td=""><td><rdl< td=""><td>1000</td><td><mdl< td=""><td><rdl< td=""><td>00.0</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<></td></mdl<></td></rdl<></td></rdl<></td></rdl<></td></mdl<></td></rdl<></td></rdl<>	<rdl< td=""><td><mdl< td=""><td><rdl< td=""><td>00.40</td><td><rdl< td=""><td><rdl< td=""><td>1000</td><td><mdl< td=""><td><rdl< td=""><td>00.0</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<></td></mdl<></td></rdl<></td></rdl<></td></rdl<></td></mdl<></td></rdl<>	<mdl< td=""><td><rdl< td=""><td>00.40</td><td><rdl< td=""><td><rdl< td=""><td>1000</td><td><mdl< td=""><td><rdl< td=""><td>00.0</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<></td></mdl<></td></rdl<></td></rdl<></td></rdl<></td></mdl<>	<rdl< td=""><td>00.40</td><td><rdl< td=""><td><rdl< td=""><td>1000</td><td><mdl< td=""><td><rdl< td=""><td>00.0</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<></td></mdl<></td></rdl<></td></rdl<></td></rdl<>	00.40	<rdl< td=""><td><rdl< td=""><td>1000</td><td><mdl< td=""><td><rdl< td=""><td>00.0</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<></td></mdl<></td></rdl<></td></rdl<>	<rdl< td=""><td>1000</td><td><mdl< td=""><td><rdl< td=""><td>00.0</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<></td></mdl<></td></rdl<>	1000	<mdl< td=""><td><rdl< td=""><td>00.0</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<></td></mdl<>	<rdl< td=""><td>00.0</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<>	00.0	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Storm	3/10/2011	Result	0.202	4MDI	0.0345	0.0254	7790	-MDI	4MD1	4MDI		2240	0.11	0.33	1820	2.82	6.48	26.9	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
WDI EVO	Dannatus	Qual	SH	<mdl< td=""><td></td><td></td><td></td><td><mdl< td=""><td><mdl< td=""><td><mdl< td=""><td><rdl< td=""><td></td><td><rdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td></td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<></td></rdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<>				<mdl< td=""><td><mdl< td=""><td><mdl< td=""><td><rdl< td=""><td></td><td><rdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td></td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<></td></rdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<>	<mdl< td=""><td><mdl< td=""><td><rdl< td=""><td></td><td><rdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td></td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<></td></rdl<></td></rdl<></td></mdl<></td></mdl<>	<mdl< td=""><td><rdl< td=""><td></td><td><rdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td></td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<></td></rdl<></td></rdl<></td></mdl<>	<rdl< td=""><td></td><td><rdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td></td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<></td></rdl<></td></rdl<>		<rdl< td=""><td><rdl< td=""><td></td><td></td><td></td><td></td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<></td></rdl<>	<rdl< td=""><td></td><td></td><td></td><td></td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td></rdl<>					NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
WKLEV3 -	Downstrea				0.205	0.0120	20000	0.27	0.52						6200	1 E	2.2	105	0.0574				1.55			1.77		0.426	0.0220			0.015
Baseline	1/4/2011	Result Qual	1.63 SH	<mdl< td=""><td>0.205</td><td>0.0139</td><td>39900</td><td>0.37 <rdl< td=""><td>0.53</td><td><mdl< td=""><td><n di<="" td=""><td></td><td><mdl< td=""><td><n di<="" td=""><td>6200</td><td>1.5</td><td>2.3</td><td>125</td><td>0.0574</td><td><mdl< td=""><td>-MDI</td><td><n di<="" td=""><td>1.55</td><td><mdl< td=""><td><mdl< td=""><td></td><td><n di<="" td=""><td>0.126</td><td>0.0338</td><td><nd1< td=""><td><md1< td=""><td>0.015</td></md1<></td></nd1<></td></n></td></mdl<></td></mdl<></td></n></td></mdl<></td></n></td></mdl<></td></n></td></mdl<></td></rdl<></td></mdl<>	0.205	0.0139	39900	0.37 <rdl< td=""><td>0.53</td><td><mdl< td=""><td><n di<="" td=""><td></td><td><mdl< td=""><td><n di<="" td=""><td>6200</td><td>1.5</td><td>2.3</td><td>125</td><td>0.0574</td><td><mdl< td=""><td>-MDI</td><td><n di<="" td=""><td>1.55</td><td><mdl< td=""><td><mdl< td=""><td></td><td><n di<="" td=""><td>0.126</td><td>0.0338</td><td><nd1< td=""><td><md1< td=""><td>0.015</td></md1<></td></nd1<></td></n></td></mdl<></td></mdl<></td></n></td></mdl<></td></n></td></mdl<></td></n></td></mdl<></td></rdl<>	0.53	<mdl< td=""><td><n di<="" td=""><td></td><td><mdl< td=""><td><n di<="" td=""><td>6200</td><td>1.5</td><td>2.3</td><td>125</td><td>0.0574</td><td><mdl< td=""><td>-MDI</td><td><n di<="" td=""><td>1.55</td><td><mdl< td=""><td><mdl< td=""><td></td><td><n di<="" td=""><td>0.126</td><td>0.0338</td><td><nd1< td=""><td><md1< td=""><td>0.015</td></md1<></td></nd1<></td></n></td></mdl<></td></mdl<></td></n></td></mdl<></td></n></td></mdl<></td></n></td></mdl<>	<n di<="" td=""><td></td><td><mdl< td=""><td><n di<="" td=""><td>6200</td><td>1.5</td><td>2.3</td><td>125</td><td>0.0574</td><td><mdl< td=""><td>-MDI</td><td><n di<="" td=""><td>1.55</td><td><mdl< td=""><td><mdl< td=""><td></td><td><n di<="" td=""><td>0.126</td><td>0.0338</td><td><nd1< td=""><td><md1< td=""><td>0.015</td></md1<></td></nd1<></td></n></td></mdl<></td></mdl<></td></n></td></mdl<></td></n></td></mdl<></td></n>		<mdl< td=""><td><n di<="" td=""><td>6200</td><td>1.5</td><td>2.3</td><td>125</td><td>0.0574</td><td><mdl< td=""><td>-MDI</td><td><n di<="" td=""><td>1.55</td><td><mdl< td=""><td><mdl< td=""><td></td><td><n di<="" td=""><td>0.126</td><td>0.0338</td><td><nd1< td=""><td><md1< td=""><td>0.015</td></md1<></td></nd1<></td></n></td></mdl<></td></mdl<></td></n></td></mdl<></td></n></td></mdl<>	<n di<="" td=""><td>6200</td><td>1.5</td><td>2.3</td><td>125</td><td>0.0574</td><td><mdl< td=""><td>-MDI</td><td><n di<="" td=""><td>1.55</td><td><mdl< td=""><td><mdl< td=""><td></td><td><n di<="" td=""><td>0.126</td><td>0.0338</td><td><nd1< td=""><td><md1< td=""><td>0.015</td></md1<></td></nd1<></td></n></td></mdl<></td></mdl<></td></n></td></mdl<></td></n>	6200	1.5	2.3	125	0.0574	<mdl< td=""><td>-MDI</td><td><n di<="" td=""><td>1.55</td><td><mdl< td=""><td><mdl< td=""><td></td><td><n di<="" td=""><td>0.126</td><td>0.0338</td><td><nd1< td=""><td><md1< td=""><td>0.015</td></md1<></td></nd1<></td></n></td></mdl<></td></mdl<></td></n></td></mdl<>	-MDI	<n di<="" td=""><td>1.55</td><td><mdl< td=""><td><mdl< td=""><td></td><td><n di<="" td=""><td>0.126</td><td>0.0338</td><td><nd1< td=""><td><md1< td=""><td>0.015</td></md1<></td></nd1<></td></n></td></mdl<></td></mdl<></td></n>	1.55	<mdl< td=""><td><mdl< td=""><td></td><td><n di<="" td=""><td>0.126</td><td>0.0338</td><td><nd1< td=""><td><md1< td=""><td>0.015</td></md1<></td></nd1<></td></n></td></mdl<></td></mdl<>	<mdl< td=""><td></td><td><n di<="" td=""><td>0.126</td><td>0.0338</td><td><nd1< td=""><td><md1< td=""><td>0.015</td></md1<></td></nd1<></td></n></td></mdl<>		<n di<="" td=""><td>0.126</td><td>0.0338</td><td><nd1< td=""><td><md1< td=""><td>0.015</td></md1<></td></nd1<></td></n>	0.126	0.0338	<nd1< td=""><td><md1< td=""><td>0.015</td></md1<></td></nd1<>	<md1< td=""><td>0.015</td></md1<>	0.015
				<ndl< td=""><td>0.460</td><td>0.0026</td><td>25400</td><td></td><td><rdl< td=""><td>< IVIDL</td><td><mdl< td=""><td></td><td><ndl< td=""><td><mdl< td=""><td>6000</td><td><rdl< td=""><td><rdl< td=""><td>112</td><td>0.112</td><td><ividl< td=""><td><mdl< td=""><td><ivid l<="" td=""><td>0.25</td><td></td><td><ndl< td=""><td>B2 0.25</td><td><mdl< td=""><td>B</td><td>0.0001</td><td></td><td><mdl< td=""><td><rdl< td=""></rdl<></td></mdl<></td></mdl<></td></ndl<></td></ivid></td></mdl<></td></ividl<></td></rdl<></td></rdl<></td></mdl<></td></ndl<></td></mdl<></td></rdl<></td></ndl<>	0.460	0.0026	25400		<rdl< td=""><td>< IVIDL</td><td><mdl< td=""><td></td><td><ndl< td=""><td><mdl< td=""><td>6000</td><td><rdl< td=""><td><rdl< td=""><td>112</td><td>0.112</td><td><ividl< td=""><td><mdl< td=""><td><ivid l<="" td=""><td>0.25</td><td></td><td><ndl< td=""><td>B2 0.25</td><td><mdl< td=""><td>B</td><td>0.0001</td><td></td><td><mdl< td=""><td><rdl< td=""></rdl<></td></mdl<></td></mdl<></td></ndl<></td></ivid></td></mdl<></td></ividl<></td></rdl<></td></rdl<></td></mdl<></td></ndl<></td></mdl<></td></rdl<>	< IVIDL	<mdl< td=""><td></td><td><ndl< td=""><td><mdl< td=""><td>6000</td><td><rdl< td=""><td><rdl< td=""><td>112</td><td>0.112</td><td><ividl< td=""><td><mdl< td=""><td><ivid l<="" td=""><td>0.25</td><td></td><td><ndl< td=""><td>B2 0.25</td><td><mdl< td=""><td>B</td><td>0.0001</td><td></td><td><mdl< td=""><td><rdl< td=""></rdl<></td></mdl<></td></mdl<></td></ndl<></td></ivid></td></mdl<></td></ividl<></td></rdl<></td></rdl<></td></mdl<></td></ndl<></td></mdl<>		<ndl< td=""><td><mdl< td=""><td>6000</td><td><rdl< td=""><td><rdl< td=""><td>112</td><td>0.112</td><td><ividl< td=""><td><mdl< td=""><td><ivid l<="" td=""><td>0.25</td><td></td><td><ndl< td=""><td>B2 0.25</td><td><mdl< td=""><td>B</td><td>0.0001</td><td></td><td><mdl< td=""><td><rdl< td=""></rdl<></td></mdl<></td></mdl<></td></ndl<></td></ivid></td></mdl<></td></ividl<></td></rdl<></td></rdl<></td></mdl<></td></ndl<>	<mdl< td=""><td>6000</td><td><rdl< td=""><td><rdl< td=""><td>112</td><td>0.112</td><td><ividl< td=""><td><mdl< td=""><td><ivid l<="" td=""><td>0.25</td><td></td><td><ndl< td=""><td>B2 0.25</td><td><mdl< td=""><td>B</td><td>0.0001</td><td></td><td><mdl< td=""><td><rdl< td=""></rdl<></td></mdl<></td></mdl<></td></ndl<></td></ivid></td></mdl<></td></ividl<></td></rdl<></td></rdl<></td></mdl<>	6000	<rdl< td=""><td><rdl< td=""><td>112</td><td>0.112</td><td><ividl< td=""><td><mdl< td=""><td><ivid l<="" td=""><td>0.25</td><td></td><td><ndl< td=""><td>B2 0.25</td><td><mdl< td=""><td>B</td><td>0.0001</td><td></td><td><mdl< td=""><td><rdl< td=""></rdl<></td></mdl<></td></mdl<></td></ndl<></td></ivid></td></mdl<></td></ividl<></td></rdl<></td></rdl<>	<rdl< td=""><td>112</td><td>0.112</td><td><ividl< td=""><td><mdl< td=""><td><ivid l<="" td=""><td>0.25</td><td></td><td><ndl< td=""><td>B2 0.25</td><td><mdl< td=""><td>B</td><td>0.0001</td><td></td><td><mdl< td=""><td><rdl< td=""></rdl<></td></mdl<></td></mdl<></td></ndl<></td></ivid></td></mdl<></td></ividl<></td></rdl<>	112	0.112	<ividl< td=""><td><mdl< td=""><td><ivid l<="" td=""><td>0.25</td><td></td><td><ndl< td=""><td>B2 0.25</td><td><mdl< td=""><td>B</td><td>0.0001</td><td></td><td><mdl< td=""><td><rdl< td=""></rdl<></td></mdl<></td></mdl<></td></ndl<></td></ivid></td></mdl<></td></ividl<>	<mdl< td=""><td><ivid l<="" td=""><td>0.25</td><td></td><td><ndl< td=""><td>B2 0.25</td><td><mdl< td=""><td>B</td><td>0.0001</td><td></td><td><mdl< td=""><td><rdl< td=""></rdl<></td></mdl<></td></mdl<></td></ndl<></td></ivid></td></mdl<>	<ivid l<="" td=""><td>0.25</td><td></td><td><ndl< td=""><td>B2 0.25</td><td><mdl< td=""><td>B</td><td>0.0001</td><td></td><td><mdl< td=""><td><rdl< td=""></rdl<></td></mdl<></td></mdl<></td></ndl<></td></ivid>	0.25		<ndl< td=""><td>B2 0.25</td><td><mdl< td=""><td>B</td><td>0.0001</td><td></td><td><mdl< td=""><td><rdl< td=""></rdl<></td></mdl<></td></mdl<></td></ndl<>	B2 0.25	<mdl< td=""><td>B</td><td>0.0001</td><td></td><td><mdl< td=""><td><rdl< td=""></rdl<></td></mdl<></td></mdl<>	B	0.0001		<mdl< td=""><td><rdl< td=""></rdl<></td></mdl<>	<rdl< td=""></rdl<>
Storm	12/8/2010	Result	1.48	<nd1< td=""><td>0.162</td><td>0.0036 <rdl< td=""><td>35400</td><td>0.26</td><td>0.46</td><td><mdi< td=""><td><n di<="" td=""><td></td><td><n di<="" td=""><td><n di<="" td=""><td>6090</td><td>2.5</td><td>2.5 <rdl< td=""><td>113</td><td>0.113</td><td><mdi< td=""><td>0.017</td><td>-MDI</td><td></td><td>0.225</td><td><ndi< td=""><td></td><td>0.027 <rdl< td=""><td></td><td>0.0231</td><td>2.6 <rdl< td=""><td><mdi< td=""><td>-MDI</td></mdi<></td></rdl<></td></rdl<></td></ndi<></td></mdi<></td></rdl<></td></n></td></n></td></n></td></mdi<></td></rdl<></td></nd1<>	0.162	0.0036 <rdl< td=""><td>35400</td><td>0.26</td><td>0.46</td><td><mdi< td=""><td><n di<="" td=""><td></td><td><n di<="" td=""><td><n di<="" td=""><td>6090</td><td>2.5</td><td>2.5 <rdl< td=""><td>113</td><td>0.113</td><td><mdi< td=""><td>0.017</td><td>-MDI</td><td></td><td>0.225</td><td><ndi< td=""><td></td><td>0.027 <rdl< td=""><td></td><td>0.0231</td><td>2.6 <rdl< td=""><td><mdi< td=""><td>-MDI</td></mdi<></td></rdl<></td></rdl<></td></ndi<></td></mdi<></td></rdl<></td></n></td></n></td></n></td></mdi<></td></rdl<>	35400	0.26	0.46	<mdi< td=""><td><n di<="" td=""><td></td><td><n di<="" td=""><td><n di<="" td=""><td>6090</td><td>2.5</td><td>2.5 <rdl< td=""><td>113</td><td>0.113</td><td><mdi< td=""><td>0.017</td><td>-MDI</td><td></td><td>0.225</td><td><ndi< td=""><td></td><td>0.027 <rdl< td=""><td></td><td>0.0231</td><td>2.6 <rdl< td=""><td><mdi< td=""><td>-MDI</td></mdi<></td></rdl<></td></rdl<></td></ndi<></td></mdi<></td></rdl<></td></n></td></n></td></n></td></mdi<>	<n di<="" td=""><td></td><td><n di<="" td=""><td><n di<="" td=""><td>6090</td><td>2.5</td><td>2.5 <rdl< td=""><td>113</td><td>0.113</td><td><mdi< td=""><td>0.017</td><td>-MDI</td><td></td><td>0.225</td><td><ndi< td=""><td></td><td>0.027 <rdl< td=""><td></td><td>0.0231</td><td>2.6 <rdl< td=""><td><mdi< td=""><td>-MDI</td></mdi<></td></rdl<></td></rdl<></td></ndi<></td></mdi<></td></rdl<></td></n></td></n></td></n>		<n di<="" td=""><td><n di<="" td=""><td>6090</td><td>2.5</td><td>2.5 <rdl< td=""><td>113</td><td>0.113</td><td><mdi< td=""><td>0.017</td><td>-MDI</td><td></td><td>0.225</td><td><ndi< td=""><td></td><td>0.027 <rdl< td=""><td></td><td>0.0231</td><td>2.6 <rdl< td=""><td><mdi< td=""><td>-MDI</td></mdi<></td></rdl<></td></rdl<></td></ndi<></td></mdi<></td></rdl<></td></n></td></n>	<n di<="" td=""><td>6090</td><td>2.5</td><td>2.5 <rdl< td=""><td>113</td><td>0.113</td><td><mdi< td=""><td>0.017</td><td>-MDI</td><td></td><td>0.225</td><td><ndi< td=""><td></td><td>0.027 <rdl< td=""><td></td><td>0.0231</td><td>2.6 <rdl< td=""><td><mdi< td=""><td>-MDI</td></mdi<></td></rdl<></td></rdl<></td></ndi<></td></mdi<></td></rdl<></td></n>	6090	2.5	2.5 <rdl< td=""><td>113</td><td>0.113</td><td><mdi< td=""><td>0.017</td><td>-MDI</td><td></td><td>0.225</td><td><ndi< td=""><td></td><td>0.027 <rdl< td=""><td></td><td>0.0231</td><td>2.6 <rdl< td=""><td><mdi< td=""><td>-MDI</td></mdi<></td></rdl<></td></rdl<></td></ndi<></td></mdi<></td></rdl<>	113	0.113	<mdi< td=""><td>0.017</td><td>-MDI</td><td></td><td>0.225</td><td><ndi< td=""><td></td><td>0.027 <rdl< td=""><td></td><td>0.0231</td><td>2.6 <rdl< td=""><td><mdi< td=""><td>-MDI</td></mdi<></td></rdl<></td></rdl<></td></ndi<></td></mdi<>	0.017	-MDI		0.225	<ndi< td=""><td></td><td>0.027 <rdl< td=""><td></td><td>0.0231</td><td>2.6 <rdl< td=""><td><mdi< td=""><td>-MDI</td></mdi<></td></rdl<></td></rdl<></td></ndi<>		0.027 <rdl< td=""><td></td><td>0.0231</td><td>2.6 <rdl< td=""><td><mdi< td=""><td>-MDI</td></mdi<></td></rdl<></td></rdl<>		0.0231	2.6 <rdl< td=""><td><mdi< td=""><td>-MDI</td></mdi<></td></rdl<>	<mdi< td=""><td>-MDI</td></mdi<>	-MDI
		Qual	0.054	<mdl< td=""><td>0.17</td><td></td><td>20500</td><td><rdl< td=""><td><rdl< td=""><td><mdl< td=""><td><mdl 0.45</mdl </td><td></td><td><mdl< td=""><td><mdl< td=""><td>4700</td><td>1.2</td><td></td><td>90.5</td><td>NIA</td><td><mdl< td=""><td><rdl< td=""><td><mdl< td=""><td><rdl< td=""><td>NIA</td><td><mdl< td=""><td><rdl,b< td=""><td></td><td>В</td><td>NΙΛ</td><td></td><td><mdl< td=""><td><mdl NA</mdl </td></mdl<></td></rdl,b<></td></mdl<></td></rdl<></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<></td></rdl<></td></mdl<>	0.17		20500	<rdl< td=""><td><rdl< td=""><td><mdl< td=""><td><mdl 0.45</mdl </td><td></td><td><mdl< td=""><td><mdl< td=""><td>4700</td><td>1.2</td><td></td><td>90.5</td><td>NIA</td><td><mdl< td=""><td><rdl< td=""><td><mdl< td=""><td><rdl< td=""><td>NIA</td><td><mdl< td=""><td><rdl,b< td=""><td></td><td>В</td><td>NΙΛ</td><td></td><td><mdl< td=""><td><mdl NA</mdl </td></mdl<></td></rdl,b<></td></mdl<></td></rdl<></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<></td></rdl<>	<rdl< td=""><td><mdl< td=""><td><mdl 0.45</mdl </td><td></td><td><mdl< td=""><td><mdl< td=""><td>4700</td><td>1.2</td><td></td><td>90.5</td><td>NIA</td><td><mdl< td=""><td><rdl< td=""><td><mdl< td=""><td><rdl< td=""><td>NIA</td><td><mdl< td=""><td><rdl,b< td=""><td></td><td>В</td><td>NΙΛ</td><td></td><td><mdl< td=""><td><mdl NA</mdl </td></mdl<></td></rdl,b<></td></mdl<></td></rdl<></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<></td></rdl<>	<mdl< td=""><td><mdl 0.45</mdl </td><td></td><td><mdl< td=""><td><mdl< td=""><td>4700</td><td>1.2</td><td></td><td>90.5</td><td>NIA</td><td><mdl< td=""><td><rdl< td=""><td><mdl< td=""><td><rdl< td=""><td>NIA</td><td><mdl< td=""><td><rdl,b< td=""><td></td><td>В</td><td>NΙΛ</td><td></td><td><mdl< td=""><td><mdl NA</mdl </td></mdl<></td></rdl,b<></td></mdl<></td></rdl<></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<></td></mdl<>	<mdl 0.45</mdl 		<mdl< td=""><td><mdl< td=""><td>4700</td><td>1.2</td><td></td><td>90.5</td><td>NIA</td><td><mdl< td=""><td><rdl< td=""><td><mdl< td=""><td><rdl< td=""><td>NIA</td><td><mdl< td=""><td><rdl,b< td=""><td></td><td>В</td><td>NΙΛ</td><td></td><td><mdl< td=""><td><mdl NA</mdl </td></mdl<></td></rdl,b<></td></mdl<></td></rdl<></td></mdl<></td></rdl<></td></mdl<></td></mdl<></td></mdl<>	<mdl< td=""><td>4700</td><td>1.2</td><td></td><td>90.5</td><td>NIA</td><td><mdl< td=""><td><rdl< td=""><td><mdl< td=""><td><rdl< td=""><td>NIA</td><td><mdl< td=""><td><rdl,b< td=""><td></td><td>В</td><td>NΙΛ</td><td></td><td><mdl< td=""><td><mdl NA</mdl </td></mdl<></td></rdl,b<></td></mdl<></td></rdl<></td></mdl<></td></rdl<></td></mdl<></td></mdl<>	4700	1.2		90.5	NIA	<mdl< td=""><td><rdl< td=""><td><mdl< td=""><td><rdl< td=""><td>NIA</td><td><mdl< td=""><td><rdl,b< td=""><td></td><td>В</td><td>NΙΛ</td><td></td><td><mdl< td=""><td><mdl NA</mdl </td></mdl<></td></rdl,b<></td></mdl<></td></rdl<></td></mdl<></td></rdl<></td></mdl<>	<rdl< td=""><td><mdl< td=""><td><rdl< td=""><td>NIA</td><td><mdl< td=""><td><rdl,b< td=""><td></td><td>В</td><td>NΙΛ</td><td></td><td><mdl< td=""><td><mdl NA</mdl </td></mdl<></td></rdl,b<></td></mdl<></td></rdl<></td></mdl<></td></rdl<>	<mdl< td=""><td><rdl< td=""><td>NIA</td><td><mdl< td=""><td><rdl,b< td=""><td></td><td>В</td><td>NΙΛ</td><td></td><td><mdl< td=""><td><mdl NA</mdl </td></mdl<></td></rdl,b<></td></mdl<></td></rdl<></td></mdl<>	<rdl< td=""><td>NIA</td><td><mdl< td=""><td><rdl,b< td=""><td></td><td>В</td><td>NΙΛ</td><td></td><td><mdl< td=""><td><mdl NA</mdl </td></mdl<></td></rdl,b<></td></mdl<></td></rdl<>	NIA	<mdl< td=""><td><rdl,b< td=""><td></td><td>В</td><td>NΙΛ</td><td></td><td><mdl< td=""><td><mdl NA</mdl </td></mdl<></td></rdl,b<></td></mdl<>	<rdl,b< td=""><td></td><td>В</td><td>NΙΛ</td><td></td><td><mdl< td=""><td><mdl NA</mdl </td></mdl<></td></rdl,b<>		В	NΙΛ		<mdl< td=""><td><mdl NA</mdl </td></mdl<>	<mdl NA</mdl
Storm	2/28/2011	Result	0.954 SH	0.011 <rdl< td=""><td>0.17</td><td>0.024</td><td>28500</td><td>0.27 <rdl< td=""><td>0.41</td><td><mdl< td=""><td>0.45 <rdl< td=""><td></td><td><mdl< td=""><td>0.1 <rdl< td=""><td>4700</td><td>1.3</td><td>2</td><td>90.5</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA NA</td></rdl<></td></mdl<></td></rdl<></td></mdl<></td></rdl<></td></rdl<>	0.17	0.024	28500	0.27 <rdl< td=""><td>0.41</td><td><mdl< td=""><td>0.45 <rdl< td=""><td></td><td><mdl< td=""><td>0.1 <rdl< td=""><td>4700</td><td>1.3</td><td>2</td><td>90.5</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA NA</td></rdl<></td></mdl<></td></rdl<></td></mdl<></td></rdl<>	0.41	<mdl< td=""><td>0.45 <rdl< td=""><td></td><td><mdl< td=""><td>0.1 <rdl< td=""><td>4700</td><td>1.3</td><td>2</td><td>90.5</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA NA</td></rdl<></td></mdl<></td></rdl<></td></mdl<>	0.45 <rdl< td=""><td></td><td><mdl< td=""><td>0.1 <rdl< td=""><td>4700</td><td>1.3</td><td>2</td><td>90.5</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA NA</td></rdl<></td></mdl<></td></rdl<>		<mdl< td=""><td>0.1 <rdl< td=""><td>4700</td><td>1.3</td><td>2</td><td>90.5</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA NA</td></rdl<></td></mdl<>	0.1 <rdl< td=""><td>4700</td><td>1.3</td><td>2</td><td>90.5</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA</td><td>NA NA</td></rdl<>	4700	1.3	2	90.5	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA NA
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Note: Baseline indicates non-storm collection

NA = not analyzed

MDL - method detection limit

RDL - reporting detection limit

Qual- qualifiers

B - qualifier indicates result is equal to or less than 5 times the concentration found in the method blank.

B2 - qualifier indicates result is greater than 5 times but is equal to or less than 5 times the concentration found in the method blank.

SH - qualifier indicates samples not preserved within 15 minutes of collection of the last sample into the sample composite sampler.

APPENDIX D

Phase II Environmental Site Assessment, Pacific Right Bank Levee Setback Project

This Site Assessment will be provided separately on a flash drive.



APPENDIX E

Pacific Park/Dumpsite Environmental Investigation Report



PACIFIC PARK/DUMPSITE ENVIRONMENTAL INVESTIGATION

600 3RD AVENUE SOUTHEAST, PACIFIC, WASHINGTON

Prepared for River and Floodplain Management Section King County Water and Land Resources Division

Prepared by Herrera Environmental Consultants, Inc.



Note:

Some pages in this document have been purposely skipped or blank pages inserted so that this document will copy correctly when duplexed.

PACIFIC PARK/DUMPSITE ENVIRONMENTAL INVESTIGATION

600 3RD AVENUE SOUTHEAST, KING COUNTY, WASHINGTON

Prepared for
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King County Water and Land Resources Division
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CONTENTS

Certificate of Licensed Hydrogeologist					
Introduction	1				
Site Background Conditions	1				
Site Conditions	3				
Physical Setting	3				
Geology	3				
Field Investigation	7				
Utility Locate	7				
Sampling Activities	7				
Investigation Results	9				
Subsurface Conditions	9				
Analytical Results	10				
Soil Sampling	10				
Groundwater Sampling	11				
Conclusions	15				
Subsurface Conditions	15				
Soil Sampling Results	16				
Groundwater Sampling Results	16				
Recommendations	17				
Limitations	19				
References	21				
Tables	23				
Figures	33				



APPENDICES

Appendix	A Field Notes				
Appendix	a B Probe Logs				
Appendix	C Photographic Log				
Appendix	D Laboratory Analytical Reports				
Appendix	E Data Quality Assurance Review				
TABLES	S				
Table 1.	Summary of Soil Sample Results, Pacific Park/Dumpsite, Site Investigation, Pacific, Washington	25			
Table 2.	ble 2. Summary of Groundwater Sample Results, Pacific Park/Dumpsite, Site Investigation, Pacific, Washington				
Figuri	ES				
Figure 1.	Vicinity Map, Pacific Park, Pacific, Washington	35			
Figure 2.	Site Map, Pacific Park, Pacific, Washington	36			
Figure 3.	Sample Location Map, Pacific Park, Pacific, Washington	37			
Figure 4.	Water Level Contour Map, October 6, 2016, Pacific Park, Pacific, Washington	38			
Figure 5.	Water Level Contour Map, March 19, 2016, Pacific Park, Pacific, Washington	39			
Figure 6.	Water Level Contour Map, May 12, 2017, Pacific Park, Pacific, Washington	40			
Figure 7.	MTCA Metals Exceedances in Groundwater Sampling Results, Pacific Park, Pacific, Washington	41			



CERTIFICATE OF LICENSED HYDROGEOLOGIST

This document has been prepared under the supervision of a licensed hydrogeologist.



Bruce A. Carpenter	September 18, 2017		
Signature	Date		



INTRODUCTION

This report summarizes field activities, including soil and groundwater sampling results for the Pacific Park/Dumpsite (Site) located at 600 Third Avenue Southeast in Pacific, Washington (Figure 1). The Site consists of 21 acres located along the right bank of the White River. The Site is located on a 43-acre parcel owned by King County since 1921 and used as an informal dumpsite and as a city dump until it was closed in 1965 to refuse dumping. The Site remained abandoned until King County leased the 21-acre portion of the property to the City of Pacific in 1969 for use as a park that opened in 1972 (S&W, 2016). It remains closed for flood control purposes from approximately October through March. A site map is provided in Figure 2.

On February 14, 2017, Herrera Environmental Consultants, Inc., (Herrera) entered into an agreement with the King County River and Floodplain Management Section as part of King County Contract Number E00375E15, Work Order Number E00375-40. Herrera developed a Sampling and Analysis Plan (SAP) describing the field activities to be conducted for this investigation (Herrera 2017). This report focusses on soil and groundwater sampling at the park boundary, outside of the fill area, to evaluate potential offsite impacts.

SITE BACKGROUND CONDITIONS

In 2016, a Phase II Environmental Site Assessment (ESA) was conducted to evaluate conditions across the Site (S&W, 2016). Soil and groundwater samples were collected from 28 push-probes and six monitoring wells constructed from six of the probes. Additionally, four test pits were excavated. Sampling locations were selected based on historical records, including aerial photographs. The report presented the following conclusions for soil and groundwater:

- Contaminants revealed in soil in excess of the Model Toxics control Act (MTCA) regulatory cleanup criteria included the following:
 - Lube oil-range petroleum hydrocarbons were detected in two soil samples at depths of 4.5 and 5.5 feet below ground surface (bgs), above the MTCA Method A cleanup criterion.
 - Three samples, five samples, seven samples, and two samples were detected above the MTCA Method A cleanup criteria for arsenic, cadmium, lead, and mercury, respectively, at depths ranging from 4.5 to 12.5 feet bgs.
 - No volatile organic compounds (VOCs) were detected above the MTCA Method A cleanup criteria.



- Semi-volatile organic compounds (SVOCs), including carcinogenic polycyclic aromatic hydrocarbons (cPAHs), were detected above the MTCA Method A cleanup criterion at 18 soil sample locations. The MTCA Method A cleanup level for cPAHs is based on the total of cPAH compounds. The Toxics Equivalency Factor (TEF) for cPAHs is used to adjust cPAH concentrations to benzo(a)pyrene. If a particular cPAH compound is not detected, one half of the detection limit is used in the TEF calculation. At some locations, no cPAHs were reported detected, but due to a high detection limit, the TEF calculation triggered a MTCA exceedance. If detection limits are higher than normal, the cPAH total could exceed MTCA cleanup levels with no reported detections.
- No polychlorinated biphenyls (PCBs) were detected above the MTCA cleanup criterion.
- Contaminants revealed in groundwater in excess of the MTCA regulatory cleanup criteria included the following:
 - Total arsenic concentrations exceeded the MTCA Method A cleanup criterion at eight locations. Dissolved arsenic exceeded the MTCA Method A cleanup criterion at two locations.
 - Total lead concentrations exceeded the MTCA Method A cleanup criterion at three locations. Dissolved lead was not detected above the MTCA Method A cleanup criterion.
 - No other metals, VOCs or petroleum hydrocarbons were detected above the MTCA Method A cleanup criteria.
- Contaminants appear to be confined to the dumpsite and do not appear to pose a threat
 to the public, based on current land use. Additional soil and groundwater sampling was
 recommended around the perimeter of the Site to determine if contamination had
 migrated off site.
- An operations and maintenance plan should be implemented that would limit ground disturbance to less than 1 foot bgs.

In 2010 and 2011, King County (2015) collected stormwater samples from the ditch near the end of 4th Avenue SE and from two locations in the wetland located 300 feet and 1,000 feet south of 4th Avenue SE. Water was tested for 6020A metals, pesticides, herbicides, and EPA 8260C VOCs and EPA 8270D SVOCs. Results from all three locations found detectable concentrations of VOCs in surface waters. No conclusions regarding the source of the contaminants were made by King County (2015).



SITE CONDITIONS

PHYSICAL SETTING

The Site is located approximately 2,000 feet north of the King/Pierce County line. The property is relatively flat. Ground surface elevation for the Site ranges from 80 to 87 feet (NAVD88), based on 2016 King County Water and Land Resources Division lidar and orthoimagery acquisition completed by Tetra Tech.

The Site is located in an area that historically had been occupied by the White River prior to construction of levees in the 1910s and filling of the site from the 1920s through the 1960s. It is subject to flooding by the White River and seasonal inundation by ponded stormwater and shallow groundwater. According to historical photos and maps, flood-control measures were taken that included construction of a concrete revetment in 1919 as part of the channelization of the White River along the alignment of the former Stuck River (S&W 2016).

The 1936 and 1944 aerial photographs indicate an orchard occupied the area west of the Site. Fill was placed in the 1980s within the area underlain by the four apartments south of 4th Avenue SE, the southeastern half of the area underlain by the four apartments north of 4th Avenue SE, and the extreme southeastern corner of the Megan's Court Apartments.

A stormwater ditch drains south along the western edge of the Site. The ditch receives stormwater from residential properties located along 3rd Avenue SE, Spencer Court, and the apartments west of the ditch and also from the parking lot in the park. The water level in the ditch roughly corresponds to groundwater levels in adjacent wells, so the ditch presumably intercepts groundwater, which intermingles with surface stormwater conveyed by the ditch.

Access to the site is restricted from October through March by a continuous HESCO barrier system installed for flood control purposes along the northern and western property boundaries. A section of the barrier is removed from April to September to allow public access at two locations on the northern portion of the property, along 3rd Avenue SE.

Geology

Regional Geology

The Puget Lowland area has been subjected to six or more major glaciations during the past 2 million years. The last ice covering the Site to a thickness estimated to be 3,000 feet, receded about 13,500 years ago, leaving a series of north-trending ridges and valleys. These deep valleys



were partially or completely filled with recessional glacial deposits and recent Holocene deposits (S&W 2016).

Subsequent to the most recent glaciation, a series of lahars and mudflows associated with volcanism from Mt. Rainier occurred, which contributed to the deposition of sediments in the White River valley. The most significant of these volcanic events was the Osceola mudflow that occurred 5,600 years ago and diverted the White River and initiated the deposition of the White River alluvial fan. During the Sumerland age volcanism approximately 2,500 years ago, the White and Green rivers extended 20 miles north to Elliott Bay. During a large flood that occurred in 1906, the White River was re-diverted to the south and avulsed to the Stuck River.

Site Geology

The Geologic Map of Auburn 1:24,000 Quadrangle (Mullineaux, 1965) maps the Site as artificial fill (af) with the surrounding area as alluvium (Qaw). The map indicates that the alluvium consists of mostly gravel and sand deposits of the White River and boulder-cobble and pebble-cobble gravel and sand in the White River valley.

Shannon and Wilson 2016 Investigation

Fill material was encountered in each of the geoprobe explorations across the Site during the Phase II ESA (S&W 2016). Fill material was identified based on soil with a disturbed appearance and from the presence of unnatural debris such as glass shards and bottles, brick, cement, organics, wood, paper, rubber, and ceramic. Fill consisted of gravelly silt, silty sand, and sandy gravel to gravelly sand. Based on observations from the geoprobes and test pit explorations, fill is generally 2 to 12 feet thick across the Site, with fill being thicker in the central/south central portion of the Site.

The fill is overlain by sandy gravelly fill placed during the period after the dumpsite was closed in order to cover the waste and level out the site for development of a park. Based on the boring and test pit logs, the thickness of this fill ranges from 1.8 to 10 feet.

The fill is underlain by alluvial overbank deposits consisting mostly of poorly graded sand with gravel to sandy gravel interbedded with backwater lacustrine deposits consisting of silt with organics and interbeds of silty sand. Alluvium and lacustrine deposits were encountered from 2 to 12 feet bgs. Large cobbles to small boulders were not observed in the geoprobe explorations due to sampler size limitations, but were encountered in the test pit explorations. Fill material and alluvial deposits were loose, causing caving and collapse that limited excavation depths in some of the test pits. Loose soils and refuse debris limited the recovery of the geoprobe core samples at some locations (S&W 2016).



Site Hydrogeology

Shannon and Wilson 2016 Investigation

Groundwater was encountered in each of the geoprobe and test pit explorations, with the exception of the southernmost test pit and geoprobe that was advanced to characterize the soil mounds in the southern portion of the site. Groundwater depths ranged from 4 to 9 feet bgs at the time of the Phase II investigation, in September 2015. In general, groundwater was encountered between 4 to 6 feet bgs, with deeper groundwater depths encountered in areas of higher relief, such as explorations located on the existing levee and in terraced fill areas in the southern portion of the Site (S&W 2016).

Groundwater observed in test pits was rapid flowing, causing caving and collapse that limited the excavation with the exception of the southernmost test pit. Soils with generally high-permeability were observed in the test pits, which facilitated a rapid movement of groundwater into the excavation (S&W 2016).

A sketch of a water level contour map for the Site was completed by King County, based on October 2015 water level data. Groundwater flow was to the southwest and follows the gradient of the river (Brummer 2017a).

King County completed an animation of groundwater levels, based on measurements from October 2015 through October 2016. The direction of groundwater flow is typically to the southwest for most of the year. When water levels rise in winter with the river stage, the gradient increases, and flow is toward the west (Brummer 2017b).



FIELD INVESTIGATION

Herrera staff collected groundwater samples from the six monitoring wells located across the Site and soil and groundwater samples from nine push-probe borings completed around the perimeter of the Site as part of this investigation. Sampling was performed according to procedures described in the SAP. The locations of the borings were selected to determine whether contamination identified on site during the Phase II ESA has migrated off site.

UTILITY LOCATE

The toll-free underground utility location service ("one call") was contacted prior to conducting probe drilling, and paint markings were made in the public right-of-way to designate water, sewer, gas, electric, and other communication lines. A private locating service (APS of North Bend, Washington) identified underground piping around the perimeter of the property on May 23, 2017.

SAMPLING ACTIVITIES

On May 15, 2017, two Herrera representatives collected water samples from the six existing monitoring wells, MW-1 through MW-6. The water samples were hand delivered to OnSite Environmental, Inc. (OnSite) of Redmond, Washington for the following analyses:

- Gasoline-range total petroleum hydrocarbons (TPH) by method Northwest TPH-Gx
- Diesel-range TPH by Method NWTPH-Dx) with silica gel cleanup
- VOCs by EPA Method 8260C
- PAHs by EPA Method 8270D/SIM
- Total and Dissolved Priority Pollutant Metals by EPA Methods 200.8/7470A

The samples were collected by the low-flow purge method described in the SAP. The dissolved metals samples were filtered in the field through a 0.45-micron filter.

On May 23 and 24, ESN Northwest provided a track mounted Bobcat push-probe rig to complete nine borings (PP1 to PP9) to 15 feet bgs at locations shown on Figure 3. Two soil samples were collected at each location, including one from the 0- to 5-foot depth interval or groundwater interface, and the second from the 10- to 15-foot interval. Groundwater samples also were collected from the borings. Samples were hand delivered to Onsite for analysis.



Soil and groundwater samples were analyzed for the same list of analyses provided above. Soil samples also were analyzed for the entire suite of SVOCs, which includes PAHs and total metals analyses (which includes the Resource Conversation, and Recovery Act (RCRA) metals arsenic, barium, cadmium, chromium, lead, mercury, selenium, and silver).



INVESTIGATION RESULTS

Field notes are provided in Appendix A, and probe boring logs are in Appendix B. GPS coordinates for each boring location were recorded and used to produce Figure 3. Photographic documentation of the field investigation is provided in Appendix C. Laboratory reports and chain-of-custody records are provided in Appendix D, and a Data Quality Assurance Review is provided in Appendix E.

SUBSURFACE CONDITIONS

Fill material was encountered at depths ranging from 2 feet bgs in boring PP7, consisting of sandy gravel to 7 feet bgs in borings PP3 and PP4. The fill material was identified by the presence of man-made material, characteristics of the soil core sample material such as layering, and findings from previous borings completed at the site. Concrete was found near ground surface and asphalt was found at 5 feet bgs in boring PP4, a piece of wire was found near ground surface in boring PP5, and a piece of a plastic ball was found at 2 feet bgs in boring PP6. Despite the presence of some man-made material, refuse was not found in any of the nine borings. Additionally, no sheens or petroleum odors were observed during sample collection.

Silt and sand with occasional gravel were encountered predominantly beneath the fill material in the six borings, PP1 through PP6, completed along the western perimeter of the property. A 6-inch zone of peat was observed at 9.5 feet bgs in boring PP1 and at 4.5 feet bgs in boring PP5. Thin layers of peat were found at depths of 9.5 and 14.5 feet bgs in boring PP1 and at 13 feet bgs in boring PP2. Occasional thin zones of organic material also were encountered in borings PP4, PP5, and PP6. Greenish gray silty clay was observed at 12 feet bgs in boring PP2, with some thin peat layers at 12 feet bgs.

Coarse alluvial material, including sand, gravel, and cobbles were observed beneath the levee fill in borings PP7, PP8, and PP9, completed on the levee along the eastern perimeter of the Site. No silt, clay, peat, other organic material, or any man made material was observed beneath the fill.

Groundwater was measured in the nine borings at depths ranging from 1.8 feet bgs in borings PP1, PP2, and PP7 to 4 feet bgs in boring PP9 and from 0.4 feet bgs in well MW-3 to 5.71 feet bgs in well MW-6. Water level contour maps were completed for the highest and lowest water levels observed during the data collection period, which ranged from the date of monitoring well installation in October 2015 to the date of the groundwater sampling event on May 12, 2017. Water level data loggers deployed in monitoring wells MW-2 through MW-6 in October 2015, provided data used to complete Figure 4, representing the lowest water levels measured during the data collection period, which occurred on October 6, 2016. Figure 5 depicts a contour map, based on the highest water level data over the monitoring period, which occurred on



March 19, 2017. Figure 6 depicts a contour map, based on water levels measured during recent groundwater sampling on May 12, 2017.

Water level data fluctuated by approximately 4.5 feet bgs during the data collection period. The groundwater flow direction is toward the southwest during periods of lower flow during the end of the summer and toward the west to southwest during periods of higher flow during the winter.

ANALYTICAL RESULTS

Soil Sampling

A summary of soil analytical results is presented in Table 1 and summarized as follows:

- Gasoline-range total petroleum hydrocarbons and diesel-range petroleum hydrocarbons were not detected in any of the 18 samples above reporting limits.
- Lube oil-range petroleum hydrocarbons were detected in shallow soil samples collected from borings PP3, PP4, PP5, PP6, and PP7 at concentrations ranging from 62 to 620 mg/kg, well below MTCA Method A cleanup level 2,000 mg/kg. Lube oil was detected in only one deep sample collected from boring PP4 at a concentration of 570 mg/kg. Fill material was observed in Boring PP4 10 to a depth of 7 feet bgs and a piece of asphalt was observed at 5 feet bgs.
- Six VOC compounds were detected, including acetone, carbon disulfide, 2-butanone, toluene, styrene, and 1,2,4-trimethylbenzene. All of the detections were below MTCA Method A or B cleanup levels. Acetone, carbon disulfide, 2-Butanone, and toluene are common laboratory contaminants and may have been introduced into the samples at the laboratory. Styrene and 1,2,4-trimethylbenzene were detected at one location. Trimethlybenzene was detected in the sample from boring PP4 that contained fill material, including a piece of asphalt.
- Two SVOCs, including benzyl alcohol and bis(2-ethylhexyl)phthalate were detected in fill
 material from borings PP1 and PP8, respectively, at low concentrations below MTCA
 Method A or B cleanup levels.
- PAHs were detected in three samples, including the shallow and deep samples from boring PP4 and the shallow sample from boring PP6 at concentrations below the MTCA Method A cleanup level, 0.1 mg/kg total cPAHs. Lube oil was reported at each of these three sample locations.
- Barium, chromium, and lead were the only total metals detected. Barium and chromium
 were detected in all of the soil samples at concentrations well below MTCA cleanup
 levels. Chromium concentrations were typical of background conditions reported for



Puget Sound ranging from a minimum of 12 to a maximum of 235 mg/kg and median of 22 mg/kg (Ecology 1994). Lead was detected in five samples all collected from shallow sample locations. With the exception of the shallow sample collected from boring PP4, all of the sample results are typical of background concentrations in Puget Sound ranging from 4.7 to 30 mg/kg and well below the MTCA Method A cleanup level of 250 mg/kg.

 Six samples with reported lube oil detections also were analyzed for PCBs to comply with MTCA sampling requirements. PCBs were often mixed with lube oil prior to the 1976 EPA Toxic Substance Control Act regulation. Only two of the soil samples had reported detections of PCBs, including the shallow samples collected from borings PP4 and PP6.
 None of the PCBs exceeded the MTCA Method A cleanup level.

Groundwater Sampling

A summary of groundwater sampling analytical results is presented in Table 2, and detected concentrations above MTCA are provided on Figure 7 for metals analyses. The groundwater sampling results are presented according to samples collected from the monitoring wells and from the probe borings.

Monitoring Wells

- Gasoline-range total petroleum hydrocarbons, diesel-range petroleum hydrocarbons, and lube oil-range petroleum hydrocarbons were not detected in any of the six samples above the reporting limits.
- Chlorobenzene was the only VOC detected at a concentration of 1.5 μg/L in well MW-4.
 It also was the only VOC detected during the previous sampling effort in October 2015.
- No PAHs were detected in any of the six samples.
- Three of the 13 priority pollutant metals sampled in the six monitoring wells were detected, including arsenic, lead, and zinc. In most cases there was correlation between the total and dissolved detected concentrations. Arsenic was the only metal that exceeded the MTCA Method A cleanup level of 5 μg/L with an identical reported detection of 8.3 μg/L for both the total and dissolved sample collected from well MW-4. Arsenic was reported at respective total and dissolved concentrations of 13 and 9.8 μg/L during the October 2015 sampling effort. Total and dissolved arsenic also was detected at respective concentrations of 5.7 and 5.1 μg/L during the October 2015 monitoring effort in well MW-2, but it was not detected in the May 2017.



Probe Borings

- Gasoline-range total petroleum hydrocarbons were detected in samples collected from one of the nine boring locations, PP5, at a concentration of 210 μg/L, less than the MTCA Method A cleanup level of 800 μg/L.
- Diesel-range petroleum hydrocarbons, and lube oil-range petroleum hydrocarbons were not detected in any of the nine probe boring samples.
- Acetone, a common laboratory contaminant was detected in all nine probe boring samples at concentrations well below the MTCA cleanup level. Benzene and other VOCs found in gasoline were detected in the sample collected from PP5, which had a reported gasoline concentration. The reported benzene concentration of 6.4 μg/L slightly exceeds the MTCA Method A cleanup level of 5 μg/L. No gasoline was detected in the six monitoring wells located on the Site or in the three probe water samples collected in 2015. Two other VOCs, toluene and carbon disulfide were detected in the water samples collected from boring PP7, and PP9, respectively, at concentrations well below MTCA cleanup levels.
- Two PAHs, including benzo(b)fluoranthene and dibenz(a,h)anthracene exceeded the MTCA Method B, cancer criteria. No Total PAHs exceeded the MTCA Method A cleanup criterion of 0.100 µg/L.
- Multiple MTCA exceedances occurred for total metals as shown in Table 2 and on Figure 7. Ten of the 13 priority pollutant metals analyzed in the sample collected from boring PP2 exceeded MTCA cleanup levels. The sample collected from PP2 was extremely turbid, the well yield was very slow (it took over 2.5 hours to collect the sample), and much of the sample was suspended sediment that was acidified from the sample preservatives when placed into the sample container most likely resulting in elevated metals concentrations. Only two metals, arsenic and lead were detected in the dissolved sample from PP2, and only arsenic exceeded the MTCA cleanup level. Arsenic was the only MTCA exceedance for the remaining dissolved metals analyses from water samples collected from PP1 and PP4. According to MTCA guidance, 173-340-720 (9) (b), "metals analyses shall be conducted on unfiltered ground water samples, unless it can be demonstrated that a filtered water sample provides a more representative measure of ground water quality." Typically probe samples are used for screening and to guide well installation locations for long-term monitoring. Based on this MTCA guidance, the dissolved sample results are considered more representative of the groundwater quality.
- Arsenic may become sequestered in peat and other organic material from emissions
 associated with coal burning and other anthropogenic activities (Mikutta and Rothwell
 2016). As the peat decays it slowly releases arsenic into groundwater. This phenomenon
 may be occurring at Pacific Park, based on elevated levels of dissolved arsenic reported
 in the six probes along the western boundary of the Site, which contain peat and organic
 material. Dissolved arsenic was not reported in groundwater in the three probes



- completed along the eastern boundary of the Site and screened in coarser material that did not contain peat.
- Historically, arsenic use has been associated with orchards. Borings PP2, PP4, PP5, and PP6 were located in an area that was used as an orchard in the 1930s and 1940s. The elevated arsenic concentrations detected in groundwater samples collected from borings PP2, PP4, PP5, and PP6 may be attributed to arsenic used in the orchards.
- A recent publication by Ecology (2016) suggests the MTCA Method A cleanup level for arsenic of 5 μg/L should be raised to 10 or 15 μg/L, based on natural background arsenic concentrations in Washington State. In Puget Sound lowlands, reported background arsenic concentrations ranged from 0.8 to 76 μg/L with an upper tolerance limit (UTL) of 8 μg/L. If the arsenic level is raised from 5 μg/L to 10 μg/L, which is the Federal (EPA) groundwater standard for arsenic, there would be only one dissolved arsenic exceedance.



CONCLUSIONS

The purpose of this Environmental Investigation is to determine if soil and/or groundwater contamination is present outside of the fill area to determine the potential for offsite impacts. Herrera sampled the six existing monitoring wells located on the Site on May 15, 2017, and collected soil and groundwater samples from nine push-probe borings completed at the perimeter of the Site on May 23 and 24. Six probes were completed along the western perimeter of the Site and three along the eastern perimeter.

Herrera's May 2017 Investigation complements the 2015 field effort reported in the 2016 Phase II ESA completed by Shannon and Wilson, which included soil samples collected from 28 push-probe borings and groundwater samples from six probe borings and six monitoring wells across the Site.

SUBSURFACE CONDITIONS

During the 2016 Phase II investigation, fill material was encountered in each of the 28 probe explorations generally 2 to 12 feet thick across the Site. It consisted of silt, sand, and gravel with debris such as glass shards and bottles, brick, cement, organics, wood, paper, rubber, and ceramics. This fill was overlain by sandy gravel ranging from 1.8 to 10 feet thick, placed after the dumpsite was closed to level out the site; this fill was thicker in the central/south central portion of the site.

During the perimeter investigation, fill material was encountered at depths ranging from 2 feet bgs to 7 feet bgs. Despite the presence of some man-made material such as pieces of concrete, asphalt, plastic wire, plastic ball, refuse was not found in any of the nine borings. Typically, the fill consisted of silt, sand, and gravel. Additionally, no sheens or petroleum odors were observed during sample collection.

Groundwater levels fluctuated by approximately 4.5 feet bgs during the data collection period between October 2015 and May 2016 across the Site. Groundwater was measured in the nine probe borings at depths ranging from 1.8 feet bgs to 5.7 feet bgs in May 2017.

The groundwater flow direction is toward the southwest during periods of low flow in the river during the end of the summer and shifts toward the west to southwest during periods of high flow in the river during the winter.



SOIL SAMPLING RESULTS

During the 2016 Phase II investigation, 28 soil samples were collected from the 0- to 10-foot interval and from the 10- to 15-foot interval. Lube oil-range petroleum hydrocarbons were detected above the MTCA Method A cleanup criterion in two soil samples at depths of 4.5 and 5.5 feet bgs,. Three samples, five samples, seven samples, and two samples were detected above the MTCA Method A cleanup criteria for arsenic, cadmium, lead, and mercury, respectively, at depths ranging from 4.5 to 12.5 feet bgs. SVOCs, including cPAHs were detected above MTCA Method A cleanup criterion at 18 sample locations.

During this May 2017 investigation, soil samples were collected at depth intervals of 0 to 5 feet and 10 to 15 feet in each of the nine borings. The samples were analyzed for gasoline and diesel-range total petroleum hydrocarbons, VOCs, SVOCs, and total and dissolved metals. Six soil samples also were analyzed for PCBs based on detection of lube oil. None of the soil sample analyses exceeded MTCA cleanup levels.

GROUNDWATER SAMPLING RESULTS

Groundwater samples were collected from six existing monitoring wells and each of the nine probe borings. The samples were analyzed for gasoline and diesel-range total petroleum hydrocarbons, VOCs, PAHs, and total and dissolved metals.

Benzene and other VOCs, which are components of gasoline, were detected in the sample collected from PP5, which also had a reported gasoline concentration. The reported benzene concentration of 6.4 μ g/L slightly exceeds the MTCA Method A cleanup level of 5 μ g/L. No gasoline was detected in the six monitoring wells located on the Site or in the three probe water samples collected in 2015. The presence of benzene and gasoline at this location is likely attributable to an isolated spill not associated with the dumpsite.

Two PAHs, including benzo(b)fluoranthene and dibenz(a,h)anthracene exceeded the MTCA Method B, cancer criteria in the sample collected from PP6. No Total PAHs exceeded the MTCA Method A cleanup criterion of 0.100 μ g/L. The reported PAH detections may be attributable to the turbid groundwater sample.

Arsenic was the only metal detected in the groundwater monitoring wells (and only in MW-4) that exceeded the MTCA Method A cleanup level of 5 μ g/L. Arsenic also was reported above the MTCA cleanup level during the October 2015 groundwater sampling event in two wells, MW-2 and MW-4.

Multiple MTCA exceedances for total metals occurred in groundwater samples collected from the probe borings due to extremely turbid samples. In addition to sample turbidity, elevated arsenic levels in groundwater may be due to the release of arsenic sequestered in peat (and other organic material) that was derived from emissions associated with historical coal burning and the former Asarco lead and copper smelter in Tacoma. The elevated levels of arsenic may also be attributed to the orchard that historically occupied this area.



RECOMMENDATIONS

Based on a review of previous work and the findings of this investigation, Herrera offers the following recommendations:

- Install two monitoring wells on the east side of the drainage ditch adjacent to PP-2 and PP-4, and one upgradient monitoring well on the north side of the property for purposes of collecting samples for metals analyses from permanent wells that produce less turbid water samples. The upgradient well would provide background water quality conditions.
- Resample monitoring wells during seasonal high and low water level periods, typically in March and October, respectively, to evaluate effects of seasonal changes in groundwater levels and flow directions on ground water quality.



LIMITATIONS

This report has been prepared for exclusive use by King County. The analyses and conclusions included in this report are based on conditions encountered at the time of the field investigation, as well as professional experience and judgment. Herrera cannot be responsible for interpretation by others of the data contained in this report.

Herrera's services were performed with due diligence in a manner consistent with that level of care and skill ordinarily exercised by members of the profession currently practicing under similar conditions in the area. No other warranty, express or implied, is made.



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TABLES



			Т	able 1. Si	ummary o	of Soil Sai	mple Resu	ults, Pacif	ic Park/D	umpsite, S	Site Invest	igation, Pa	acific, Was	hington.					
									Sam	ple Location	1								MTCA
Analytical Parameter	PP1 PP2		2	PP3 PP4			P4	PP5		Р	PP6 PP7		PP7	PP8			PP9	Method A	
Sample Date	5/23	/2017	5/23/	2017	5/24,	/2017	5/24	/2017	5/23	/2017	5/23	/2017	5/24	1/2017	5/24	4/2017	5/24	1/2017	Cleanup Level
Depth (feet)	2.0	10	2.5	10	2.5	10	3.0	10	2.5	10	2.0	10	0	10	0.5	10	0	10	(mg/kg)
Gasoline Range Organics by	NWTPH-Gx ((mg/kg)																	
Gasoline	ND (6.9)	ND (8.0)	ND (7.8)	ND (9.0)	ND (6.6)	ND (7.0)	ND (7.4)	ND (7.1)	ND (7.0)	ND (8.1)	ND (5.7)	ND (9.7)	ND (6.6)	ND (7.8)	ND (6.1)	ND (9.6)	ND (5.6)	ND (7.0)	100
	<u>'</u>					•	Dies	el Range Org	ganics by N\	NTPH-Dx (m	g/kg)	•							
Diesel Range Organics	ND (31)	ND (33)	ND (33)	ND (36)	ND (31)	ND (31)	ND (54)	ND (55)	ND (68)	ND (34)	ND (29)	ND (37)	ND (31)	ND (31)	ND (28)	ND (36)	ND (26)	ND (31)	2,000
Lube Oil Range Organics	ND (62)	ND (67)	ND (66)	ND (73)	140	ND (62)	500	570	620	ND (67)	62	ND (75)	63	ND (63)	ND (57)	ND (71)	ND (52)	ND (62)	2,000
Volatile Organic Compounds	by EPA 8260	OC (mg/kg)																	
Acetone	ND (0.011)	0.051	0.071	0.088	0.058	0.028	ND (0.011)	0.051	ND (0.012)	0.046	0.011	0.10	ND (0.014)	0.019	0.027	0.016	ND (0.013)	0.015	72,000 ^a
Carbon Disulfide	ND (0.0016)	ND (0.0018)	ND (0.0012)	ND (0.0016)	ND (0.0017)	ND (0.0017)	ND (0.0016)	ND (0.0013)	ND (0.0058)	0.0024	ND (0.00090)	ND (0.0015)	ND (0.0014)	ND (0.0013)	ND (0.0013)	ND (0.0014)	ND (0.0013)	ND (0.0012)	8,000°
2-Butanone	ND (0.0057)	0.012	0.019	0.023	0.013	ND (0.0061)	ND (0.0057)	0.0068	ND (0.0058)	0.013	ND (0.0045)	0.025	ND (0.0071)	ND (0.0065)	ND (0.0065)	ND (0.0070)	ND (0.0063)	ND (0.0062)	48,000ª
Toluene	0.019	0.020	0.025	0.019	ND (0.0060)	0.016	0.016	0.011	ND (0.0058)	0.016	0.0096	0.0083	ND (0.0071)	0.022	0.014	0.021	0.014	0.015	7.00
Styrene	ND (0.0011)	ND (0.0013)	ND (0.0012)	ND (0.0014)	ND (0.0012)	ND (0.0012)	ND (0.0011)	ND (0.0013)	ND (0.0012)	ND (0.0013)	ND (0.00090)	ND (0.0015)	0.012	ND (0.0013)	ND (0.0013)	ND (0.0014)	ND (0.0013)	ND (0.0012)	16,000ª
1,2,4-Trimethylbenzene	ND (0.0011)	ND (0.0013)	ND (0.0012)	ND (0.0014)	ND (0.069)	ND (0.0012)	ND (0.0011)	0.0024	ND (0.0012)	ND (0.0013)	ND (0.00090)	ND (0.0015)	ND (0.0014)	ND (0.0013)	ND (0.0013)	ND (0.0014)	ND (0.057)	ND (0.0012)	_
Semivolatile Organic Compo				,	,	,	,		,	,	,	, ,	,	1	,	,	, ,	,	
Benzyl Alcohol	ND (0.21)	ND (0.22)	ND (0.22)	ND (0.24)	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.22)	ND (0.22)	ND (0.19)	ND (0.25)	ND (0.20)	ND (0.21)	0.70	ND (0.24)	ND (0.17)	ND (0.21)	800ª
Bis(2-ethylhexyl)phthalate	0.052	ND (0.045)	ND (0.044)	ND (0.048)	ND (0.042)	ND (0.041)	ND (0.042)	ND (0.041)	ND (0.044)	ND (0.045)	ND (0.039)	ND (0.050)	ND (0.041)	ND (0.042)	ND (0.038)	ND (0.047)	ND (0.035)	ND (0.041)	1,600ª
Polycyclic Aromatic Hydroca	rhons by FD/	 ∆8270D/SIM	l (ma/ka)	l	(0.0 12)	(0.011)	(0.012)	(0.011)	(0.011)	(0.015)	(0.033)		<u> </u>		(0.050)		(0.033)		
2-Methylnaphthalene	ND (0.0082)	ND (0.0089)	ND (0.0088)	ND (0.0097)	ND (0.0084)	ND (0.0082)	0.017	0.010	ND (0.0087)	ND (0.0090)	ND (0.0077)	ND (0.0099)	ND (0.0082)	ND (0.0084)	ND (0.0076)	ND (0.0095)	ND (0.0070)	ND (0.0083)	320 ^a
Phenanthrene	ND (0.0082)	ND (0.0089)	ND (0.0088)	ND (0.0097)	ND (0.0084)	ND (0.0082)	0.028	0.011	ND (0.0087)	ND (0.0090)	ND (0.0077)	ND (0.0099)	ND (0.0082)	ND (0.0084)	ND (0.0076)	ND (0.0095)	ND (0.0070)	ND (0.0083)	_
Anthracene	ND (0.0082)	ND (0.0089)	ND (0.0088)	ND (0.0097)	ND (0.0084)	ND (0.0082)	0.0089	ND (0.0082)	ND (0.0087)	ND (0.0090)	ND (0.0077)	ND (0.0099)	ND (0.0082)	ND (0.0084)	ND (0.0076)	ND (0.0095)	ND (0.0070)	ND (0.0083)	24,000°
Fluoranthene	ND (0.0082)	ND (0.0089)	ND (0.0088)	ND (0.0097)	ND (0.0084)	ND (0.0082)	0.042	ND (0.0082)	ND (0.0087)	ND (0.0090)	0.013	ND (0.0099)	ND (0.0082)	ND (0.0084)	ND (0.0076)	ND (0.0095)	ND (0.0070)	ND (0.0083)	3,200ª
Pyrene	ND (0.0082)	ND (0.0089)	ND (0.0088)	ND (0.0097)	ND (0.0084)	ND (0.0082)	0.046	ND (0.0082)	ND (0.0087)	ND (0.0090)	0.015	ND (0.0099)	ND (0.0082)	ND (0.0084)	ND (0.0076)	ND (0.0095)	ND (0.0070)	ND (0.0083)	2,400ª
Benzo(a)anthracene	(0.0082) ND (0.0082)	ND (0.0089)	ND (0.0088)	ND (0.0097)	ND (0.0084)	ND (0.0082)	0.029	ND (0.0082)	ND (0.0087)	(0.0090) ND (0.0090)	0.0084	ND (0.0099)	ND (0.0082)	ND (0.0084)	ND (0.0076)	ND (0.0095)	ND (0.0070)	ND (0.0083)	1.37 ^b
Chrysene	(0.0082) ND (0.0082)	(0.0089) ND (0.0089)	(0.0088) ND (0.0088)	(0.0097) ND (0.0097)	ND (0.0084)	ND (0.0082)	0.037	0.014	(0.0087) ND (0.0087)	(0.0090) ND (0.0090)	0.012	(0.0099) ND (0.0099)	ND (0.0082)	ND (0.0084)	(0.0076) ND (0.0076)	(0.0093) ND (0.0095)	(0.0070) ND (0.0070)	(0.0083) ND (0.0083)	137 ^b



			Tab	le 1 (cont	inued). S	ummary o	of Soil Sa	mple Resu	ılts, Pacif	ic Park/D	umpsite, S	ite Investi	gation, Pa	cific, Wash	ington.				
									Samp	le Location	1								МТСА
Analytical Parameter	P	P1	P	P2	Р	P3	Р	P4	PI	P5	P	P6	ı	PP7	Р	P8	P	P9	Method A
Sample Date	5/23,	/2017	5/23,	/2017	5/24	/2017	5/24,	/2017	5/23,	/2017	5/23,	/2017	5/24/2017		5/24/2017		5/24/2017		Cleanup Level
Depth (feet)	2.0	10	2.5	10	2.5	10	3.0	10	2.5	10	2.0	10	0	10	0.5	10	0	10	Level (mg/kg)
Polycyclic Aromatic Hydroca	arbons by EPA	\8270D/SIN	l (mg/kg) (c	ontinued)					•			•		-				•	
Benzo(b)fluoranthene	ND (0.0082)	ND (0.0089)	ND (0.0088)	ND (0.0097)	ND (0.0084)	ND (0.0082)	0.042	ND (0.0082)	ND (0.0087)	ND (0.0090)	0.016	ND (0.0099)	ND (0.0082)	ND (0.0084)	ND (0.0076)	ND (0.0095)	ND (0.0070)	ND (0.0083)	1.37 ^b
Benzo(j,k)fluoranthene	ND (0.0082)	ND (0.0089)	ND (0.0088)	ND (0.0097)	ND (0.0084)	ND (0.0082)	0.014	ND (0.0082)	ND (0.0087)	ND (0.0090)	ND (0.0077)	ND (0.0099)	ND (0.0082)	ND (0.0084)	ND (0.0076)	ND (0.0095)	ND (0.0070)	ND (0.0083)	13.7 ^b
Benzo(a)pyrene	ND (0.0082)	ND (0.0089)	ND (0.0088)	ND (0.0097)	ND (0.0084)	ND (0.0082)	0.034	ND (0.0082)	ND (0.0087)	ND (0.0090)	0.012	ND (0.0099)	ND (0.0082)	ND (0.0084)	ND (0.0076)	ND (0.0095)	ND (0.0070)	ND (0.0083)	0.100
Indeno(1,2,3-cd)pyrene	ND (0.0082)	ND (0.0089)	ND (0.0088)	ND (0.0097)	ND (0.0084)	ND (0.0082)	0.027	ND (0.0082)	ND (0.0087)	ND (0.0090)	0.0092	ND (0.0099)	ND (0.0082)	ND (0.0084)	ND (0.0076)	ND (0.0095)	ND (0.0070)	ND (0.0083)	1.37 ^b
Benzo(g,h,i)perylene	ND (0.0082)	ND (0.0089)	ND (0.0088)	ND (0.0097)	ND (0.0084)	ND (0.0082)	0.032	ND (0.0082)	ND (0.0087)	ND (0.0090)	0.011	ND (0.0099)	ND (0.0082)	ND (0.0084)	ND (0.0076)	ND (0.0095)	ND (0.0070)	ND (0.0083)	-
Total cPAHs (TEF) ^c	0.0062	0.0067	0.0066	0.0073	0.0063	0.0062	0.046	0.0063	0.0066	0.0068	0.016	0.0075	0.0062	0.0063	0.0057	0.0072	0.0053	0.0063	0.100
Total Metals by EPA 6010C/	7471B (mg/kg	g)																	
Arsenic	ND (12)	ND (13)	ND (13)	ND (15)	ND (13)	ND (12)	ND (13)	ND (12)	ND (13)	ND (13)	ND (12)	ND (15)	ND (12)	ND (13)	ND (11)	ND (14)	ND (10)	ND (12)	20
Barium	62	30	31	21	80	26	78	38	47	28	53	34	31	22	22	37	62	28	16,000ª
Cadmium	ND (0.62)	ND (0.67)	ND (0.66)	ND (0.73)	ND (0.63)	ND (0.62)	ND (0.63)	ND (0.62)	ND (0.65)	ND (0.67)	ND (0.58)	ND (0.74)	ND (0.61)	ND (0.63)	ND (0.57)	ND (0.71)	ND (0.52)	ND (0.62)	2.00
Chromium	20	13	11	11	27	9.5	29	13	23	13	29	14	12	11	13	13	25	14	2,000
Lead	8.9	ND (6.7)	ND (6.6)	ND (7.3)	25	ND (6.2)	84	ND (6.2)	27	ND (6.7)	9.7	ND (7.4)	ND (6.1)	ND (6.3)	ND (5.7)	ND (7.1)	ND (5.2)	ND (6.2)	250
Mercury	ND (0.31)	ND (0.33)	ND (0.33)	ND (0.36)	ND (0.31)	ND (0.31)	ND (0.32)	ND (0.31)	ND (0.33)	ND (0.34)	ND (0.29)	ND (0.37)	ND (0.31)	ND (0.31)	ND (0.28)	ND (0.36)	ND (0.26)	ND (0.31)	2.00
Selenium	ND (12)	ND (13)	ND (13)	ND (15)	ND (13)	ND (12)	ND (13)	ND (12)	ND (13)	ND (13)	ND (12)	ND (15)	ND (12)	ND (13)	ND (11)	ND (14)	ND (10)	ND (12)	400 ^a
Silver	ND (1.2)	ND (1.3)	ND (1.3)	ND (1.5)	ND (1.3)	ND (1.2)	ND (1.3)	ND (1.2)	ND (1.3)	ND (1.3)	ND (1.2)	ND (1.5)	ND (1.2)	ND (1.3)	ND (1.1)	ND (1.4)	ND (1.0)	ND (1.2)	400 ^a
Polychlorinated Biphenyls b	y EPA 8082A	(mg/kg)																	
Aroclor 1254	-	_	_	_	ND (0.063)	_	ND (0.063)	ND (0.062)	ND (0.065)	_	0.12	_	ND (0.061)	-	_	_	_	_	_
Aroclor 1260	-	_	-	-	ND (0.063)	_	0.18	ND (0.062)	ND (0.065)	-	ND (0.058)	_	ND (0.061)	-	_	-	_	-	-
Total PCBs	-	-	_	_	ND (0.063)	_	0.18	ND (0.062)	ND (0.065)	_	0.12	_	ND (0.061)	-	-	-	_	-	1.00

BOLD values detected above the reporting limit.

cPAHs = carcinogenic polycyclic aromatic hydrocarbons

mg/kg = milligrams per kilogram

September 2017

MTCA = Model Toxic s Control Act (Chapter 173-340 Washington Administrative Code [WAC]).

ND = not detected above laboratory reporting limits shown in parentheses

Pacific Park/Dumpsite Environmental Investigation

^a MTCA Method B, non cancer

b MTCA Method B, cancer

^c Total cPAHs (TEF) was calculated using ½ the reporting limit for compounds that were not detected above the reporting limit.

^{– =} not analyzed or not applicable

							Sa	mple Locatio	n							
			Monitoring W	/ell Samples				mpic zocatio	···	Pı	ısh Probe San	nnles				
Analytical Parameter	MW1	MW2	MW3	MW4	MW5	MW6	PP1-W	PP2-W	PP3-W	PP4-W	PP5-W	PP6-W	PP7-W	PP8-W	PP9-W	MTCA
Sample Date	101002	141442	5/12/2		WW	141440	112 **	112 **	113 **	117 00	5/23-24/201		117 **	110 11	113 W	Method A Cleanup Level
NWTPH-Gx (µg/L)	I		3/12/2	-017							3/23 2 1/201	,				
Gasoline Range Organics	ND (100)	ND (100)	ND (100)	ND (100)	ND (100)	ND (100)	ND (100)	ND (100)	ND (100)	ND (100)	210	ND (100)	ND (100)	ND (100)	ND (100)	800
NWTPH-Dx (mg/L)	140 (100)	140 (100)	140 (100)	110 (100)	110 (100)	110 (100)	140 (100)	110 (100)	110 (100)	140 (100)	210	110 (100)	140 (100)	110 (100)	140 (100)	
Diesel Range Organics	ND (0.26)	ND (0.27)	ND (0.26)	ND (0.26)	ND (0.26)	ND (0.26)	ND (0.27)	ND (0.37)	ND (0.26)	ND (0.26)	ND (0.27)	ND (0.26)	ND (0.28)	ND (0.26)	ND (0.28)	500
Lube Oil	ND (0.20)	ND (0.27)	ND (0.42)	ND (0.42)	ND (0.20)	ND (0.41)	ND (0.43)	ND (0.60)	ND (0.42)	ND (0.20)	ND (0.27)	ND (0.42)	ND (0.28)	ND (0.41)	ND (0.28)	500
Volatile Organic Compound		110 (0.77)	110 (0.72)	ND (0.72)	ND (0.71)	ND (0.41)	ND (0. 1 3)	140 (0.00)	14D (0.42)	ND (0. 1 1)	110 (0.43)	110 (0.42)	IND (U.TT)	ND (0.71)	ND (0.43)	
Acetone	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	6.7	75	5.3	5.6	5.5	7.9	6.1	5.2	7.2	7,200ª
Carbon disulfide	ND (5.0)	ND (3.0)	ND (3.0)	ND (0.20)	ND (0.20)	ND (3.0)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	0.22	800°
Benzene	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	6.4	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	5.0
Toluene	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	5.2	ND (1.0)	ND (1.0)	1,000
Chlorobenzene	ND (0.20)	ND (0.20)	ND (0.20)	1.5	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	0.43	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	160°
Xylenes ^c	ND (0.40)	ND (0.40)	ND (0.40)	ND (0.40)	ND (0.40)	ND (0.40)	ND (0.40)	ND (0.40)	ND (0.40)	ND (0.40)	32.57	ND (0.40)	ND (0.40)	ND (0.40)	ND (0.40)	1,000
Isopropylbenzene	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	0.45	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	_
n-Propylbenzene	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	0.95	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	800°
1,3,5-Trimethylbenzene	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	3.1	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	80°a
1,2,4-Trimethylbenzene	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	12	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	_
Total Metals (µg/L)	•		•					•					•			
Antimony	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	47	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	6.4 ^a
Arsenic	ND (3.3)	ND (3.3)	ND (3.3)	8.3	ND (3.3)	ND (3.3)	36	910	18	110	14	14	110	ND (3.3)	30	5.0
Beryllium	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	34	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	<i>32</i> ^a
Cadmium	ND (4.4)	ND (4.4)	ND (4.4)	ND (4.4)	ND (4.4)	ND (4.4)	ND (4.4)	24	ND (4.4)	ND (4.4)	ND (4.4)	ND (4.4)	ND (4.4)	ND (4.4)	ND (4.4)	5.0
Chromium	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	73	2,000	63	240	29	45	210	ND (11)	49	50
Copper	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	80	4,200	120	460	50	54	570	ND (11)	110	640 ^a
Lead	ND (1.1)	ND (1.1)	ND (1.1)	ND (1.1)	4.0	ND (1.1)	250	2,100	54	2,800	32	55	250	ND (1.1)	61	15
Mercury	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	4.3	ND (0.50)	1.2	ND (0.50)	ND (0.50)	0.68	ND (0.50)	ND (0.50)	2.0
Nickel	ND (22)	ND (22)	ND (22)	ND (22)	ND (22)	ND (22)	33	1,800	32	190	ND (22)	46	210	ND (22)	44	<i>320</i> ^a
Selenium	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	53	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	6.0	ND (5.6)	ND (5.6)	80ª
Silver	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	ND (11)	80 ^a
Thallium	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	9.6	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	ND (5.6)	0.16 ^a
Zinc	160	42	170	80	37	ND (28)	700	5,500	130	1,000	49	100	520	71	210	<i>4,800</i> ^a



September 2017

Pacific Park/Dumpsite Environmental Investigation

Table 2 (continued). Summary of Groundwater Sample Results, Pacific Park/Dumpsite, Site Investigation, Pacific, Washington.																
							Sa	mple Locatio	on							
			Monitoring W	/ell Samples						Р	ush Probe Sam	nples				MTCA
Analytical Parameter	MW1	MW2	MW3	MW4	MW5	MW6	PP1-W	PP2-W	PP3-W	PP4-W	PP5-W	PP6-W	PP7-W	PP8-W	PP9-W	MTCA Method A
Sample Date		•	5/12/2	2017				5/23-24/2017							Cleanup Level	
Dissolved Metals (μg/L) (co	ntinued)															
Antimony	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	7.1	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	6.4ª
Arsenic	ND (3.0)	ND (3.0)	ND (3.0)	8.3	ND (3.0)	ND (3.0)	18	28	3.5	7.1	3.3	3.6	ND (3.0)	ND (3.0)	ND (3.0)	5.0
Beryllium	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	32ª
Cadmium	ND (4.0)	ND (4.0)	ND (4.0)	ND (4.0)	ND (4.0)	ND (4.0)	ND (4.0)	ND (4.0)	ND (4.0)	ND (4.0)	ND (4.0)	ND (4.0)	ND (4.0)	ND (4.0)	ND (4.0)	5.0
Chromium	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	50
Copper	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	640 ^a
Lead	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	3.0	ND (1.0)	ND (1.0)	1.2	ND (1.0)	3.1	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	15
Mercury	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	2.0
Nickel	ND (20)	ND (20)	ND (20)	ND (20)	ND (20)	ND (20)	ND (20)	ND (20)	ND (20)	ND (20)	ND (20)	ND (20)	ND (20)	ND (20)	ND (20)	<i>320</i> ^a
Selenium	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	80 ^a
Silver	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	ND (10)	80 ^a
Thallium	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	0.16 ^a
Zinc	170	ND (25)	140	37	35	ND (25)	39	ND (25)	ND (25)	ND (25)	ND (25)	ND (25)	ND (25)	62	ND (25)	<i>4,800</i> ^a
Polycyclic Aromatic Hydroc	arbons (µg/L)															
Naphthalene	ND (0.095)	ND (0.099)	ND (0.095)	ND (0.096)	ND (0.10)	ND (0.10)	ND (0.096)	ND (0.12)	ND (0.096)	ND (0.10)	0.18	ND (0.097)	ND (0.097)	ND (0.095)	ND (0.096)	160
Benzo(a)anthracene	ND (0.0095)	ND (0.0099)	ND (0.0095)	ND (0.0096)	ND (0.010)	ND (0.010)	ND (0.0096)	ND (0.012)	ND (0.0096)	0.013	ND (0.0097)	0.052	ND (0.0097)	ND (0.0095)	ND (0.0096)	0.12 ^b
Chrysene	ND (0.0095)	ND (0.0099)	ND (0.0095)	ND (0.0096)	ND (0.010)	ND (0.010)	ND (0.0096)	0.012	ND (0.0096)	0.021	ND (0.0097)	0.080	ND (0.0097)	ND (0.0095)	ND (0.0096)	<i>12</i> ^b
Benzo(b)fluoranthene	ND (0.0095)	ND (0.0099)	ND (0.0095)	ND (0.0096)	ND (0.010)	ND (0.010)	ND (0.0096)	0.020	ND (0.0096)	0.036	ND (0.0097)	0.14	ND (0.0097)	ND (0.0095)	0.019	0.12 ^b
Benzo(j,k)fluoranthene	ND (0.0095)	ND (0.0099)	ND (0.0095)	ND (0.0096)	ND (0.010)	ND (0.010)	ND (0.0096)	ND (0.012)	ND (0.0096)	0.011	ND (0.0097)	0.042	ND (0.0097)	ND (0.0095)	ND (0.0096)	1.2 ^b
Benzo(a)pyrene	ND (0.0095)	ND (0.0099)	ND (0.0095)	ND (0.0096)	ND (0.010)	ND (0.010)	ND (0.0096)	ND (0.012)	ND (0.0096)	0.024	ND (0.0097)	0.051	ND (0.0097)	ND (0.0095)	0.0098	0.100
Indeno(1,2,3-cd)pyrene	ND (0.0095)	ND (0.0099)	ND (0.0095)	ND (0.0096)	ND (0.010)	ND (0.010)	ND (0.0096)	0.014	ND (0.0096)	0.040	ND (0.0097)	0.066	ND (0.0097)	ND (0.0095)	0.011	0.12 ^b
Dibenz(a,h)anthracene	ND (0.0095)	ND (0.0099)	ND (0.0095)	ND (0.0096)	ND (0.010)	ND (0.010)	ND (0.0096)	ND (0.012)	ND (0.0096)	0.011	ND (0.0097)	0.018	ND (0.0097)	ND (0.0095)	ND (0.0096)	<i>0.012</i> ^b
Benzo(g,h,i)perylene	ND (0.0095)	ND (0.0099)	ND (0.0095)	ND (0.0096)	ND (0.010)	ND (0.010)	ND (0.0096)	0.020	ND (0.0096)	0.046	ND (0.0097)	0.055	ND (0.0097)	ND (0.0095)	0.012	_
Total cPAHs (TEF) ^d	0.0072	0.0075	0.0072	0.0072	0.0076	0.0076	0.0072	0.011	0.0072	0.039	0.0073	0.084	0.0073	0.0072	0.014	0.100

BOLD values detected above the reporting limit.

Shaded values exceed MTCA criteria.

- a MTCA Method B, non cancer
- b MTCA Method B, cancerc Xylenes is the sum of m-,p-, and o-xylene.
- d Total cPAHs (TEF) was calculated using ½ the reporting limit for compounds that were not detected above the reporting limit.

– = not analyzed or not applicable

cPAHs = Carcinogenic polycyclic aromatic hydrocarbons

 μ g/L = micrograms per liter

mg/L = milligrams per kilogram

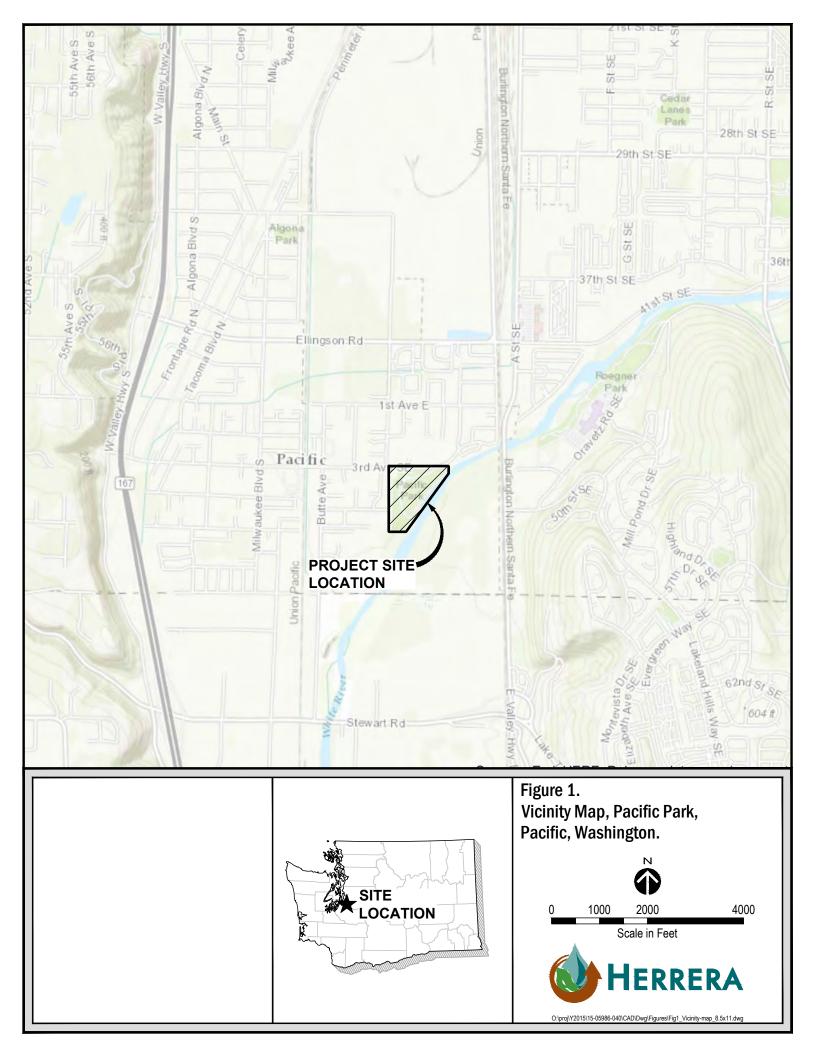
MTCA = Model Toxic s Control Act (Chapter 173-340 Washington Administrative Code [WAC]).

ND = not detected above laboratory reporting limits shown in parentheses

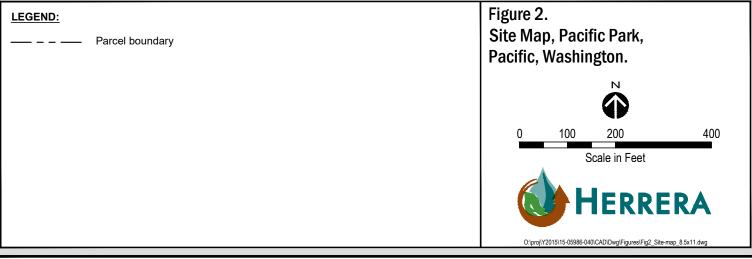
HERRERA

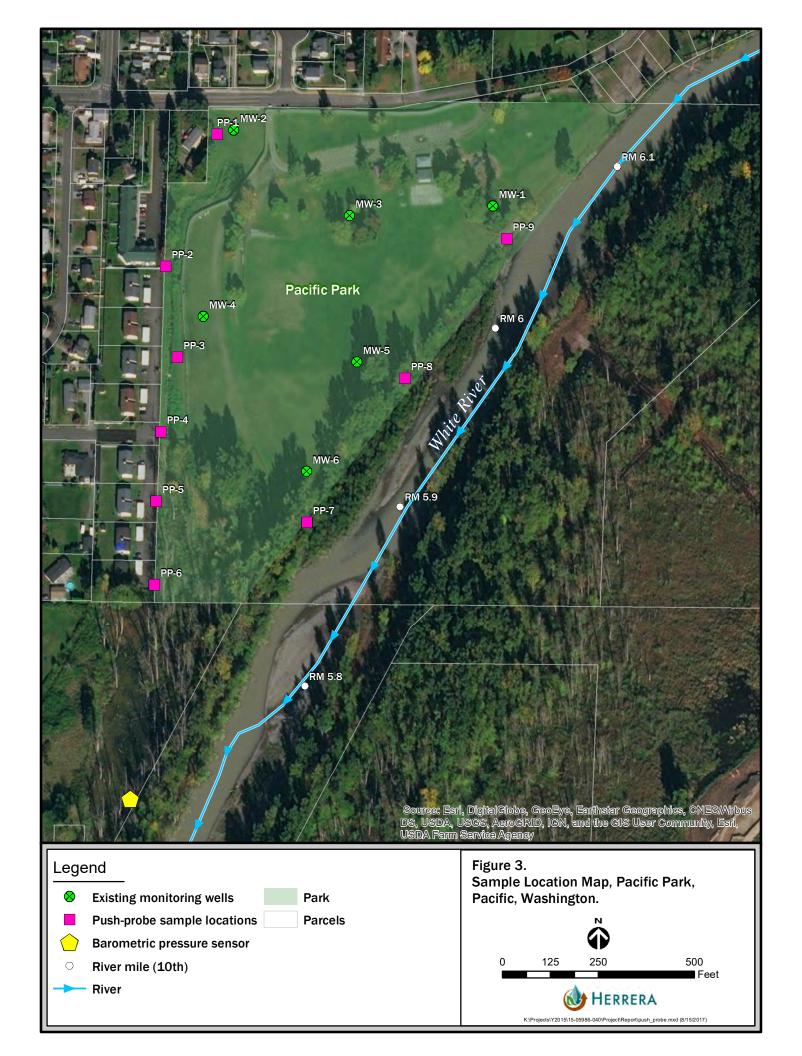
FIGURES

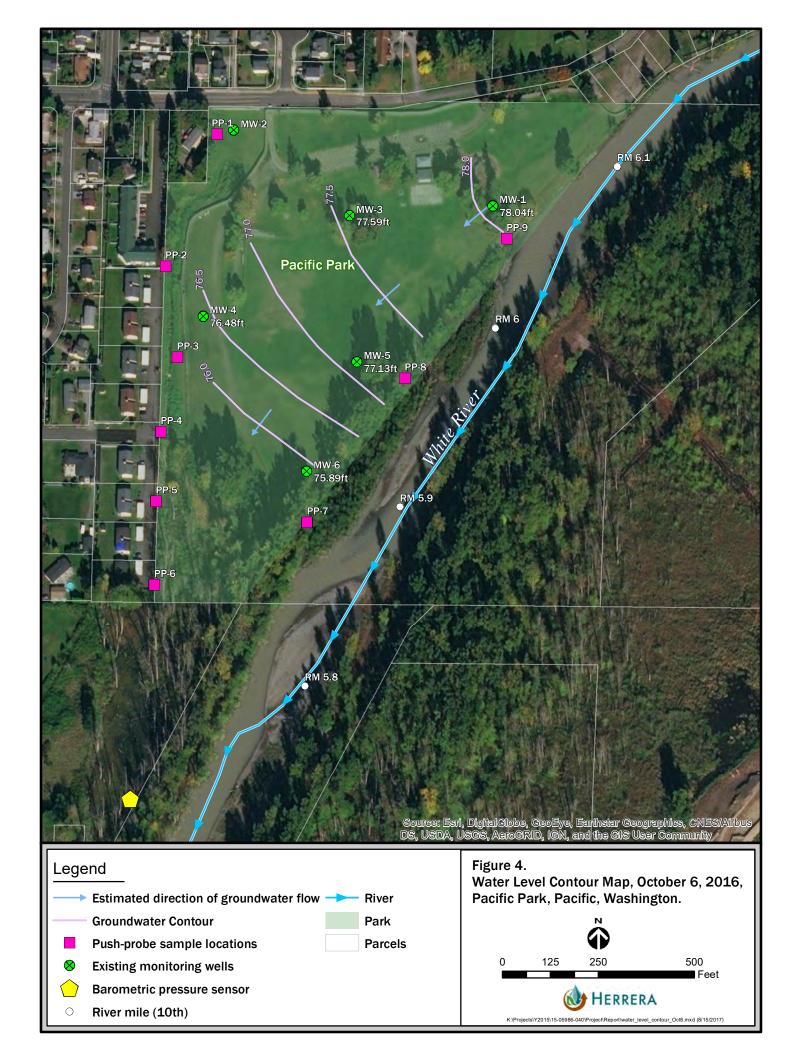


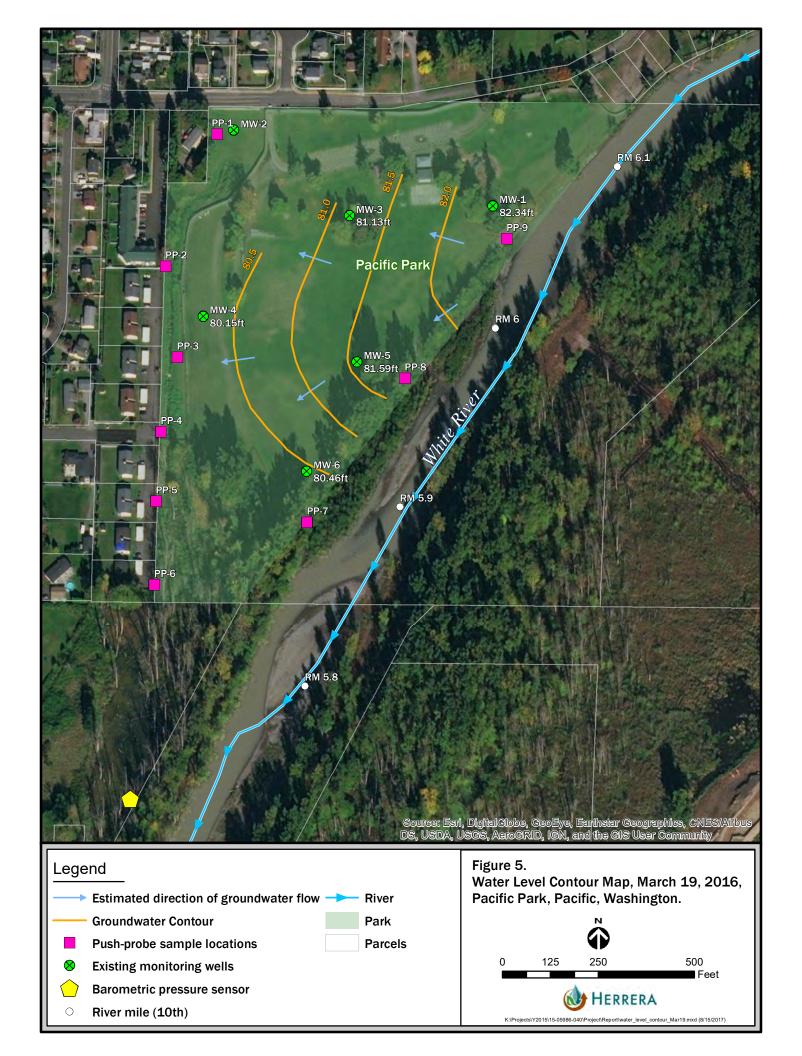


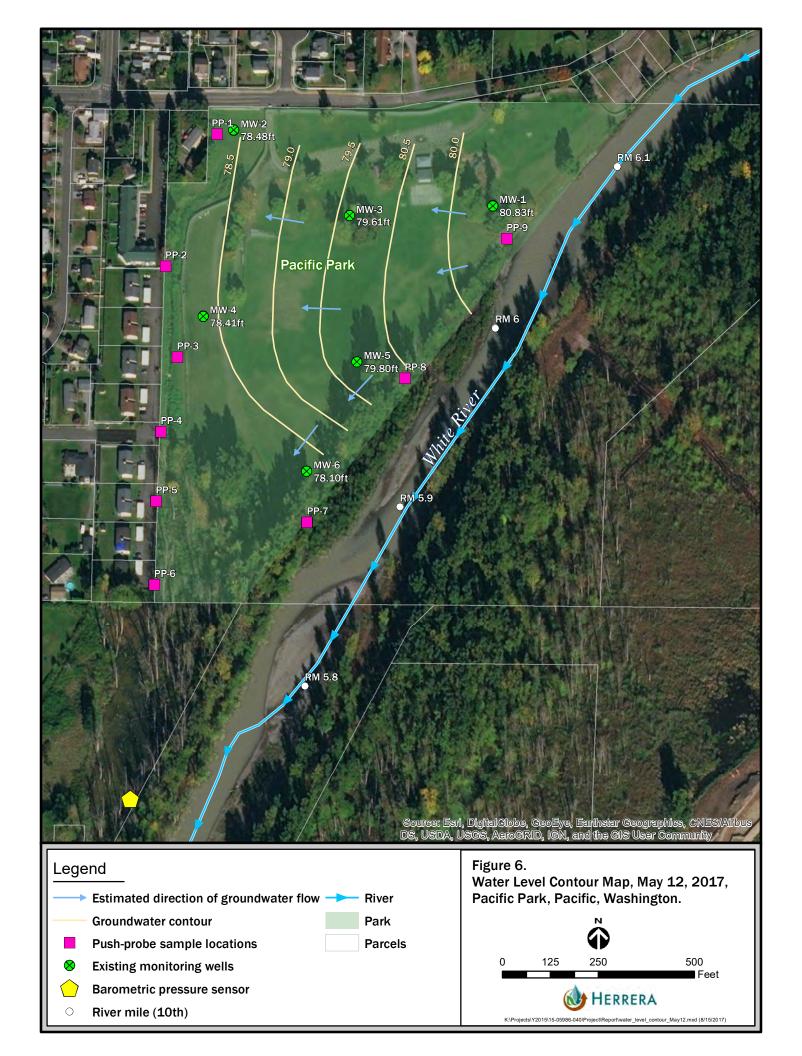


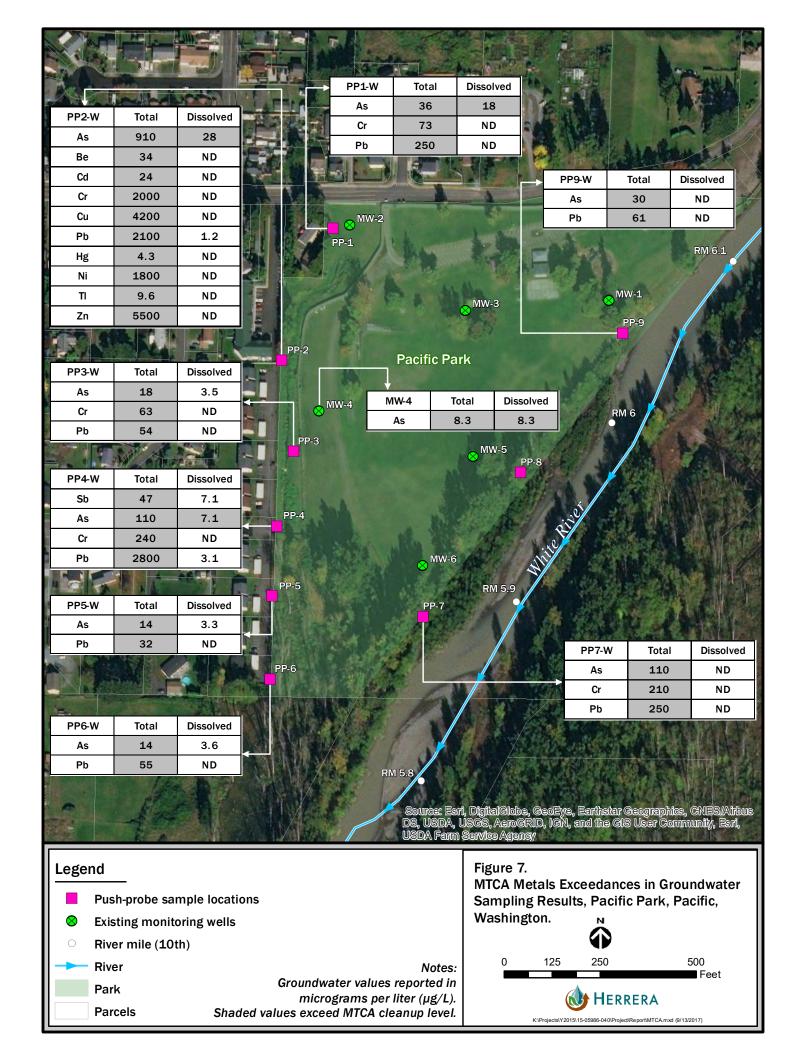












APPENDIX A

Field Notes





Project No.: 15-08/80 Site: Pacific Pour	Well No.: MWI Date: 5/12/17
Well Depth: 15 Screen Length: 10	Well Diameter: 2" Casing Type: PUC
Sampling Device Peristaltic Tubing Type:	polyfluline Water Level: 2.33
Measuring Point: top of well Other Info:	Flow 0.4 L/m year
Sampling Personnel: 8B G	

Time	рН	Temp.	Cond.	Dis.O ₂	Turb.	[]Conc.			Notes
11:55			(00,8					sturt	purging
12:00	7.81	9.4	39.3	5.99	89.3				1
12:05	7.04	9.2	18.5	4.52	42.5		5		
12:10	6.94	9.1	98.0	3.80	27,2		cleen	w	Myss od
12:15	636	9.0	1777	3.4	14.0	•			
12:20	4.84	9.0	97.5	3,28	10.0	٠			
12:22								Star-	Samplin
12:35								WLS	2,33
								,	
								1	,
						*			
						\ \ \			
		1				1			
	wecter	Jevels	alvos	al slight	Alar he	t sta	ble a	10	ell was
	DINNER	ed	11	0	4	00 0 100		7 001	7-1-1

Type of Samples Collected for Laboratory Analysis:

5 VOAS for GK VOCS, 2x 600 ml Amber Dx, 2x IL Amber PAts, 2x500 ml porty
Wall Casing Volumes:

T3D PP matals Well Casing Volumes:

 $\text{Gal/Ft} \quad 1\frac{1}{4}\text{"} = 0.077 \quad 1\frac{1}{2}\text{"} = 0.10 \quad 2\text{"} = 0.16 \quad 2\frac{1}{2}\text{"} = 0.24 \quad 3\text{"} = 0.37 \quad 3\frac{1}{2}\text{"} = 0.50 \quad 4\text{"} = 0.65 \quad 6\text{"} = 1.46$



Project No.: 15-05986 Site: NW 2 Well No.: MWZ Date: 5/12/17
Well Depth: Screen Length: 10.0 Well Diameter: 2.0 Casing Type: PVC
Sampling Device Peristaltic Tubing Type: Polyethylene Water Level: 1.37
Measuring Point: Nom top of Other Info: Flan 0,4 L/n 4cel
Sampling Personnel: Geoval 1. Briama

Time	pН	Temp.	Cond.	Dis.O ₂	Turb.	[]Conc.			Notes
10:45								Star	t purge
10:47	8.02	11.7	391.4 45	4,67 h	\$ 10,6				1
0150	7,18	11.8	342	3,82	19.0			WL-	= 1.4(
0155	10.76	12.0	319	3.04	52.0		clear	1	iless, odori
1:00	4.46	12.0	306	2.63	980				1.43
11:05	458	11.9	294	247	1352				
11:10	¢						Gum	pla	ollection
					,				
			,						
-					4				
				-	-			-	,
la.	ater 1	levels	cha spen	slight	h hi	vt stad	A nA	124	l was
r r	unje	1	//	0	0	· V - 100	- MA	0.00	C V WY
Wolfea	to	CIANA	TN	= mw	DUP	010:00	2 140		1

Type of Samples Collected for Laboratory Analysis:

for Laboratory Analysis:

\[
\langle VOCE, 2x 500 m; Amber Dx, 2x | Lamber PAHs, 2x 500 m; poly
\]

Well Casing Volumes:

\[
\text{Volumes:}
\]

Gal/Ft $1\frac{1}{4}$ " = 0.077 $1\frac{1}{2}$ " = 0.10 2" = 0.16 $2\frac{1}{2}$ " = 0.24 3" = 0.37 $3\frac{1}{2}$ " = 0.50 4" = 0.65 6" = 1.46



Project No.: 13-05986 -040 Site: Pacific Park	Well No.: _MW3 Date: _S-12(17
	Well Diameter: 2" Casing Type: PVC
Sampling Device PONSTATIC Tubing Type:	olyethy 1 se Water Level: 0.40
Measuring Point: Top of well other Info:	Flow 0,4 c/m brel
Sampling Personnel: BB G1	

Time	рH	Temp.	Cond.	Dis.O ₂	Turb.	[]Conc.		T	Natar
13:15	P	Temp.	Cona.	D13.02	Turb.	Sta	rb.	NA/4	Notes 0,29
	0.12	1	101	1	00.	Pu	rging	WL=	0,2
13:20	7.43	12.4	291.1	4.25	28.1		,		
13:25	7.15	12.0	310.4	1.94	20.9	clea	V Gold	114 S	adonesa
13:30	7.10	N.Le	321.3	1.10	16.3		tar	ins	odones
13:35	7.08	11.2	328.1	0.79	33.0		1	Mouz	wout
	4.03	10.9	331.9	0.69	87.1				
13:43						S	art	sam	ding
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						7)			
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-									
							-		·
	Water	- level	demon	rl Sisan	Hy bu	et stabi	pas	NELL	1200
	Pwan	e.	1/		3				- V- V

Type of Samples Collected for Laboratory Analysis:

S VOAS for Gx VOCs, 2x SOO my Amber Dx, 2x IL Amber PAHS, 2x 500 poly
Well Casing Volumes: T3D PP metals

Gal/Ft $1\frac{1}{4}$ " = 0.077 $1\frac{1}{2}$ " = 0.10 2" = 0.16 $2\frac{1}{2}$ " = 0.24 3" = 0.37 $3\frac{1}{2}$ " = 0.50 4" = 0.65 6" = 1.46



Project No.: 15-05986 Project No.: 15-05986 Paulic Paule Well No.: MW-4 Date: 5/12/17
Well Depth: 15 Screen Length: 10 Well Diameter: 2" Casing Type: FUC
Sampling Device penstaltic Tubing Type: powerlulere Water Level: 2,73
Measuring Point: top of Well Other Info. The purge vate
Sampling Personnel: BB G 0.4 L/M

Time	pН	Temp.	Cond.	Dis.O ₂	Turb.	[]Conc.			Notes
14:27							STENCE	puro	e.
14:30	11.47	11.3	31097	219	12.4			= 2.7	
14:35	V.53	11.5	367.8		30.2				
14:40	4.58	11.5	355.10		42.7		WL	=2:	18
14:45	10:00	N.5	348:2	0.19	449				
14148						4	start	Sam	pling
									*
						-			
ζ.,									
					*				
								-	**
	Wafer	1400	el dry	peil shi	Mathe. 1	out s	toble	as we	Of Lus
	PIM	R.P.D.	V	1					4

Type of Samples Collected for Laboratory Analysis:

5 VOAS for GX VOCS, 2x 500 mL Amber DX, 2x 12 Amber PAHS, 2 x 500 mL poly
Well Casing Volumes:
TBD PProtection

 $\text{Gal/Ft} \qquad 1\frac{1}{4}\text{"} = 0.077 \qquad 1\frac{1}{2}\text{"} = 0.10 \qquad 2\text{"} = 0.16 \qquad 2\frac{1}{2}\text{"} = 0.24 \qquad 3\text{"} = 0.37 \qquad 3\frac{1}{2}\text{"} = 0.50 \qquad 4\text{"} = 0.65 \qquad 6\text{"} = 1.46$



Project No.:	15-05	166 O Site: 1	Paulie	Park 1	Well No.:	NWS	Da	te: 5	12/17		
Well Depth:	_15	Screen	Length:	10 V	Vell Diamet	er: 2	Ca		e: PVC		
Sampling De	evice Pe	vistalti	Tubir	ng Type: pd	netrul	ene W	ater Leve	d:	40		
Measuring P	oint: to	of well	Other	Info: 0	wain	at	0,41	In	rate		
Measuring Point: top of well other Info: pwying at 0.4 L/m vate Sampling Personnel: BB GI											
Marie Tourist Control of the Control											
Time.	pН	Temp.	Cond.	Dis.O ₂	Turb.	[]Conc.			Notes		
12:15	SW	De.	ws/ch	ma/L	NTU			STOM	+ powar		
12117	7,53	9150	165	m 6.87	113			Part .	· 1 11 1		
12:20	7,20	9,5	157	6.78	10	-	WL=	1,65			
12125	7.12	9.6	156	0.98	1.2						
12:30	7.09	9.6	155	1,03	1,2		,1 =	1.1.5	7		
12:35	7.08	9,5	156	1.06		reach	DAA '	tibi	10		
12:40	^	ollect	1	Samo)	0-/1	y epopoli	1	100	1 11		
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	Wute	y levi	e dro	iped Sh	ighter	1, bin	Sta	ble a	& well		
	war	1 puni		V	/	/ '					
		* 1					,				
vpe of Sample	es Collecte	ed for Labor	ratory Anal	veie.					94 X		

5 VOAS POV GX VOCS, 2 x 500 ML Amber Dx, 2 x 1L Amber PAH3, 2 x 500 ML poly
Well Casing Volumes: T3D PP metals Well Casing Volumes:

 $\text{Gal/Ft} \quad 1\frac{1}{4}\text{"} = 0.077 \quad 1\frac{1}{2}\text{"} = 0.10 \quad 2\text{"} = 0.16 \quad 2\frac{1}{2}\text{"} = 0.24 \quad 3\text{"} = 0.37 \quad 3\frac{1}{2}\text{"} = 0.50 \quad 4\text{"} = 0.65 \quad 6\text{"} = 1.46$



	(5-059	186	_						
Project No.:	-04	O Site: 3	Parilic ¹	Pourle v	Veli No.: -	MW-LO		te: <u>5</u>	1217
Well Depth:	15'	Screen	Length:	(0' N	/ell Diamet	ter: 2"	Ca	sing Type	= PVC
Sampling Device Pristaltic Tubing Type: Polyelly lene Water Level: 5,71									
Measuring Point: Top of MMMM Other Info: water lands have visen Wire									
Compling Decomposition									
sampling Personnel: BB (4) and the river, Pwal at 0.44 / Mrin.									
Time	pН	Temp.	Cond.	Dis.O ₂	Turb.	[]Conc.			Notes
13:25	Sin	°C-	us/cm	mg/L	NTIL		St	mit	Pural
13:30	6.44	10.6	134	1,04	3.8			0.4	L/120
13:35	6.29	10.6	134	0.48	2,6	(5,85	-WL	clean	Ciplor lass
13:40	6,26	10.2	33	0.28	0.7			oday	V 10
13:45	6,25	10.2	133	0,27	0,6€	- Hert	2		
13:50	6.25	10.2	132	0.25		collea	tol.	Sa	imple.
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	Wa	ten 1	م الاسا	table a	1 1111	00 100	1 011	5000	
			VIC 37	YWORL II	4 1/10	CE MAR	PVU	JECL	
								-	
pe of Sample	e Collage	ed for Labor	aton/ Anali	usia:					
4 PER PROPERTY AND A			aithe villa	U N I N					

5 VOAS for Gx/VOCs, 2x500 ml Amber Dx, 2x IL Amber PAH, 2x500 ml poly

Well Casing Volumes:

Gal/Ft $1\frac{1}{4}$ " = 0.077 $1\frac{1}{2}$ " = 0.10 2" = 0.16 $2\frac{1}{2}$ " = 0.24 3" = 0.37 $3\frac{1}{2}$ " = 0.50 4" = 0.65 6" = 1.46

HERRERA

25-12-17 PACIFIC PARK
Beorge Ithan + Brianna Blands 8:00 Am moite
Weather 55 Sunny rain earlier this morning roccasion al showers
Health + Safety Briefing
Calibrated multi meters
Pro P\$5 # 1. 758 4 mmHg
DI - U.Zuslam DH 7: 7: U 1000 : 1003 US LM PH 10: 10.11 DO: 104.81. PH 4: 4,42
DI: 4.3 NS cm >+7: 6.99
DO: 101.4 1. PH 4: 3.89
Harry trouble califorating ptt 10:45 Started purying MWZ 11:10: Sample time MWZ 10:00: Sample time MWZ 10:00: Sample time MW Sup 10:00 AM
1000: Sample time and sup 10:00 Am

5-12-17	PACIFIC PARK.
11:55	MWI start purge Collected Sample
12:15	MWS Start puge Collected sample,
13:15	MW3 start proge collected samples
13:25 18:50	MW6 Start Purge Collected Sample
14:27	Start purge MW-4 Collected sample,
Purge Wa drum	ter put in St gallan de sop by dimpsters, Labeled top withe contents and
-1	Usite
	2

Rete in the Rain.

5/23/17 Sonny 60's Bruce Corpenter and Ryan - APS stility to cater. Walked site and cleared barehole 8:00 Chris B. and Brianna asrive at site. Meet of Sean UC re. Vac/HESCO. 9:00 APS Locator - Ryan left site 10:05 Sel 01.8' Stem set 5:15' Backfilled brehole 4/ besture! "pre w/ 105 lot 10:20 March to PP Cocation in Megamis 10:35 Collected Sil Sample PP2-2.5@2.5 10:50 Collected six south PP2-10010 screen get 5.15'insticle, moved screen up from 3 to 13. Apr. mgp stopped by doing saw 13:30 Very sonto recharge. Confair

Rite in the Rain

Rete in the Rain

5/24/17 Paarie Park 7:25 Arrive at sote Cloudy, Sois with breeze 7:30 Call brote PIT 7.45 Dave to PP.4 B: 00 Meet Brame + Pichod. Esw Daller. 8:15 Begin drilling PP4. Divided to 5'- con code a hunter March are 5 and redilled come. 8:30 Soil Sauce PP-4-3 9:40 Soil sample PALF-co 9:00 Began collection water Surple 9:20 Buchfilled trackete 9:25 Moved do Latin PP.3. 9:35 Begin direlling PP3. Por recovery 0'0-5' Por recovery @ 10-15' Set screen from 5'- 15! Began Collecty water surfle 10:00. Movedares 2' and redilled to 15' to get additional sometimes soil. 10:05 Soil Sample PP3- 20 10:15 Sail Super PP3-10 10:30 Moved divil in off site

second dill tods, Suplere boaded Na onto trailer and moved to PAG location 11:15 Began drilly at PP-9. For recovery due to coarse gravel Collected two O-S' core samples and combined sample into contained 11:50 Soil Souple PP9-0 (composite from (2) 0-5 core samples is 220% recovery 12:00 SNI Sample PP9-10 17:30 Began collecting water Sample. 12:30 Bal filled brochole moved to Cocation PP-B. 12:45 Begin dilling @ PPB. 1 0-51, 20% regnery drillar seems location and had 30% recorny. Combined sample. 13:10 water fample begin collection 13:105vil Sample 888-05 13:30 Sail Sample PP8-10. 13:45 Dupticate water same PP10-W collected tran PP8. 14.10 Moved to west Location PP7. 14:15 Begin dilling - PP7 14:30 Co locked soil Sample PP7-0

from 0-5' @ PP7. Depleate Souple from PP7-0 laboled as PP10-0 @ 15:0 141.40 Sil sample PP7-10. Screen Set from D-10. 15.50 Water Sample. 15:10 Barlifolled brocke, left site to 15:30 Burn left site. 15:40 Dilles and I Cat site

Rite in the Rain

APPENDIX B

Probe Logs





 Boring ID
 PP1

 Total depth
 15 feet

 Sheet
 1
 of
 1

Project name Pacific Park	Drilling Contractor ESN	Drilling method Push-probe rig
Project number <u>15-05986-040</u>	Location NW corner of park	Sampling method 5 ft core with plastic liner
Client King County	~ 50 feet west of MW-2	Air monitoring (Y/N) Yes
HEC rep. Bruce Carpenter	Date May 23, 2017	Instrument(s) Photoionization detector

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group ML	Soil description Grass/Brown sandy SILT, trace of gravel, FILL, damp
0	5-foot core with liner	20/50	<u>▼</u> 1.8	1 2 3 4	SM	Brown SILT, trace of sand, FILL, damp Static water level measured at 1.8 feet. Soil sample PP1-2 collected at 9:20 Brown silty SAND, trace of gravel, FILL,wet
0	5-foot core with liner	100		6 7 8 9	ML SM	Dark Brown-gray sandy SILT, wet Dark Brown-black silty SAND, wet
0	5-foot core with	100	-	10 11 12 13	PT SM ML SM	Dark Brown-gray PEAT, damp Soil sample PP1-10 collected at 9:45 Dark Brown-black silty SAND, wet 8" Dark Brown SILT, trace of sand, wet Dark Brown-black, silty SAND, wet
	liner			14	ML ML/PT ML	Brown-gray SILT, damp Brown-gray SILT, damp Brown-gray SILT, damp Brown-gray SILT, damp Set temporary screen from 5 feet to 15 feet. Purged approximately 1-gallon prior to sample collection.
						Collected water sample PP1-W at 10:00. Backfilled borehole with bentonite chips.



 Boring ID
 PP2

 Total depth
 15 feet

 Sheet
 1
 of
 1

Project name _	Pacific Park	Drilling Co	ntractor ESN	Drilling method Pu	sh-probe rig
Project number	15-05986-040	Location	SE corner of Megan's Court	Sampling method	5 ft core with plastic liner
Client King C	County			Air monitoring (Y/N)	Yes
HEC rep. E	Bruce Carpenter	Date I	May 23, 2017	Instrument(s)	Photoionization detector

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group	Soil description
W.1. /		,			ML	Grass/Brown gravelly sandy SILT, FILL, damp
				1		Brown sandy SILT, FILL, damp
	5-foot			2		Brown Sandy St. 1, 11LL, damp
0.1	core	60	$\frac{\nabla}{2.5}$			
	with		2.5	3		Soil sample PP2-2.5 collected at 10:35
	liner			4		Groundwater encountered during drilling Dark Brown sandy SILT, wet
				5		
				6		
	5-foot			7		
0	core with	50		8	SM	Dark Brown silty SAND, wet
	liner			0		
				9		
				10	ML	Dark Brown sandy SILT, trace of gravel, wet
				10		Soil sample PP2-10 collected at 10:50
				11	SM	Dark Brown-black silty SAND, wet
	5 C .			10		
0	5-foot core	100		12	MH	Greenish-gray silty CLAY, with red-brown mottling, damp
	with	100		13	14111	Greenish gray shey C2711, what red Grown mouning, damp
	liner				MH/PT	Occasional 1-inch peat lenses
				14		
				15		
						Initially set temporary screen from 5 feet to 15 feet, raised screen to 2 feet to 12 feet. Very slow yielding, took 2.5 hours to collect water Sample. Purged approximately 1-quart prior to sample collection. Began water sample PP2-W collection at 11:00. Backfilled borehole with bentonite chips.



 Boring ID
 PP3

 Total depth
 15 feet

 Sheet
 1
 of
 1

Project name Pacific Park	Drilling Contractor ESN	Drilling method Push-probe rig
Project number <u>15-05986-040</u>	Location East side of ditch	Sampling method 5 ft core with plastic liner
Client King County	~ 190 feet north of PP4	Air monitoring (Y/N) Yes
HEC rep. Bruce Carpenter	Date May 24, 2017	Instrument(s) Photoionization detector

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group	Soil description
W 1 /				1	ML	Grass/Brown gravelly sandy SILT, FILL, damp
0	5-foot core	40/35	_	2		Brown gravelly SILT, FILL, damp
	with liner		3.1	3	ML	Brown gravelly SILT, FILL, wet. Soil sample PP3-2.5 collected at 10:05 Static water level measured at 3.1 feet.
	iner		3.1	4		Static water level measured at 3.1 feet.
				6		
	5-foot			7	ML	Brown sandy SILT, FILL, wet
0	core with liner	50/55		8	SP	Dark Brown-black fine to medium SAND, trace of gravel, wet
	inici			9		
				10		Soil sample PP3-10 collected at 10:15
0	5-foot core	0/100		12		
	with liner	0/100		13		
				14		
				15		Greenish-gray sandy SILT, wet Set temporary screen from 5 feet to 15 feet.
						Purged approximately 1-gallon prior to sample collection. Collected water sample PP3-W at 10:00 Backfilled borehole with bentonite chips.



Boring ID _	PP4		
Total depth	15 feet		
Sheet 1	of 1		

Project name Pacific Park	Drilling C	Contractor ESN	Drilling method F	Push-probe rig
Project number <u>15-05986-040</u>	Location	East end of 4th Avenue SE in	Sampling method	5 ft core with plastic liner
Client King County	grass		Air monitoring (Y/N)	Yes
HEC rep. Bruce Carpenter	Date	May 24, 2017	Instrument(s)	Photoionization detector

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group	Soil description
M I Z	5-foot	,		1 2	SM SW	Grass/Dark Brown gravelly silty SAND, some concrete, FILL, damp 4" zone of medium silty SAND, FILL, damp
0	core with liner	50	<u>▼</u> 3.2	3	ML	Brown gravelly sandy SILT, red-brown mottling, FILL, damp. Soil sample PP4-3 collected at 8:30
				5	SW	Static water level measured at 3.2 feet. Dark Brown-black gravelly SAND, FILL, wet
				6	SM	Gray-black gravelly silty SAND, piece of asphalt, FILL, wet
0	5-foot core with liner	55		8	ML	Dark Brown-black sandy SILT, wet
	iner			9	SP	Dark Brown-black fine to medium SAND, wet
				11		Soil sample PP4-10 collected at 8:40
0	5-foot core with	100		12	ML	Dark Brown SILT, wet
	liner			1.4	SM	Dark Brown silty SAND, wet
				15	ML	Dark Brown SILT, organic material, damp
						Set temporary screen from 5 feet to 15 feet. Purged approximately 1-gallon prior to sample collection. Collected water sample PP4-W at 9:00 Backfilled borehole with bentonite chips.



Boring ID _	PP5
Total depth	15 feet
Sheet 1	of 1

Project name	Pacific Park	Drilling Cont	ractor ESN	Drilling method P	ush-probe rig
Project number	15-05986-040	Location	~ 215 feet north of PP6,	Sampling method	5 ft core with plastic liner
Client King	County	east side o	f HESCOs	Air monitoring (Y/N)	Yes
HEC rep.	Bruce Carpenter	Date Ma	ay 23, 2017	Instrument(s)	Photoionization detector

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group	Soil description
0	5-foot core	50		2	SM	Grass/Topsoil, Dark Brown gravelly silty SAND, piece of wire, FILL, damp
	with liner	30	<u>▼</u> 3.0	3	SW	Soil sample PP5-2.5 collected at 15:40 Dark Brown-gray gravelly SAND, FILL, damp Static water level measured at 3.0 feet.
				5	PT	Brown PEAT, wet
	5-foot			6	ML	Dark Brown-black sandy SILT, wet
0.3	core with liner	60		8	SW	Dark Brown-black fine to medium SAND, trace of silt, organic material, wet
				10		Dark Brown-black gravelly fine to coarse SAND, trace of silt, wet
0	5-foot core with liner	40		11 12 13	SW	Soil sample PP5-10 collected at 15:50 Dark Brown-black fine to medium SAND, trace of gravel, wet
				14		Dark Brown-black, fine to coarse gravelly SAND, wet
						Set temporary screen from 5 feet to 15 feet. Purged approximately 1-gallon prior to sample collection. Collected water sample PP5-W at 16:00 Backfilled borehole with bentonite chips.



Boring ID	PP6
Total depth	15 feet
Sheet 1	of 1

Project name	Pacific Park	Drilling Cont	ractor ESN	Drilling method Pu	ısh-probe rig
Project numbe	r <u>15-05986-040</u>	Location	SW corner of park, east side	Sampling method	5 ft core with plastic liner
Client King	County	of HESCO	es .	Air monitoring (Y/N)	Yes
HEC rep.	Bruce Carpenter	Date Ma	ay 23, 2017	Instrument(s)	Photoionization detector

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group	Soil description
(ррііі)		recovery	(leet)	1	GW	Grass/Brown sandy GRAVEL, FILL, damp
0.1	5-foot core with liner	40	<u>▼</u> 3.3	3 4	SM	Soil sample PP6-2 collected at 14:35 Brown gravelly silty SAND, piece of a ball, FILL, damp Static water level measured at 3.3 feet .
0	5-foot core with liner	60		6 7 8 9	SM	Dark Brown-black silty SAND, organic material, wet
0	5-foot core with liner	100		11 12 13	ML SP	Soil sample PP6-10 collected at 14:45 Dark Brown sandy SILT, organic material, wet Dark Brown-black, medium SAND, trace of silt, wet
				15	ML	Dark Brown sandy SILT, wet Set temporary screen from 2 feet to 12 feet. Purged approximately 1-gallon prior to sample collection. Collected water sample PP6-W at 15:00 Backfilled borehole with bentonite chips.



 Boring ID
 PP7

 Total depth
 15 feet

 Sheet
 1
 of
 1

Project name Pacific Park	Drilling Contractor ESN	Drilling method Push-probe rig
Project number <u>15-05986-040</u>	Location SE corner of park, ~ 125 feet	Sampling method 5 ft core with plastic liner
Client King County	south of MW-6	Air monitoring (Y/N) Yes
HEC rep. Bruce Carpenter	Date May 24, 2017	Instrument(s) Photoionization detector

PID (ppm)	Sample type, interval	% recovery	Water level (feet)	Depth (feet, BGS)	Soil group	Soil description
0	5-foot core with liner	40/30/30	<u>▼</u> 1.8	1 2 3 4	SM	Dark Brown silty SAND, organic material, FILL, damp Soil sample PP7-0 collected at 14:30 Soil sample PP10-0 (duplicate sample, false time-14:35) Static water level measured at 1.8 feet Dark Brown sandy GRAVEL, trace of silt, cobbles, wet
0	5-foot core with liner	75		6 7 8 9	SM	Dark Brown silty gravelly SAND, wet
0	5-foot core with liner	60		11 12 13 14		Soil sample PP7-10 collected at 14:40
						Set temporary screen from 0 feet to 10 feet. Purged approximately 1-gallon prior to sample collection. Collected water sample PP7-W at 14:50 Backfilled borehole with bentonite chips.



 Boring ID
 PP8

 Total depth
 15 feet

 Sheet
 1
 of
 1

Project name Pacific F	ark Drilling Co	ontractor ESN	Drilling method Pu	ush-probe rig
Project number 15-05	986-040 Location	~125 feet east-southeast of	Sampling method	5 ft core with plastic liner
Client King County	_ MW-5		Air monitoring (Y/N)	Yes
HEC rep. Bruce Ca	penter Date	May 24, 2017	Instrument(s)	Photoionization detector

	Sample		Water	Depth		
PID	type,	%	level	(feet,	Soil	Soil description
(ppm)	interval	recovery	(feet)	BGS)	group ML	Dark Brown sandy gravelly SILT, some wood, FILL, damp
				1		Soil sample PP8-0.5 collected at 13:20
	5.0.4			2	CM	D 'l' - II CAND FILL I
0	5-foot core	20/30		2	SM	Brown silty gravelly SAND, FILL, damp
	with	20/30		3	<u>▼</u> 3.1	wet
	liner			4	3.1	Static water level measured at 3.1 feet
				4		·
				5		
					GW	Dark Gray-black, fine to coarse sandy GRAVEL, wet
				6		
	5-foot			7		
0	core	30		0		
	with liner			8		
	111101			9		
				10		
				10		Soil sample PP8-10 collected at 13:30
				11		cobbles
	5 C4			12		
0	5-foot core	40		12		
	with			13		
	liner			1.4		
				14		
				15		
						Set temporary screen from 0 feet to 10 feet. Purged approximately 1-gallon prior to sample collection.
						Collected water sample PP8-W at 13:10
						Collected duplicate water sample PP10-W (false time 13:45)
						Backfilled borehole with bentonite chips.



 Boring ID
 PP9

 Total depth
 15 feet

 Sheet
 1
 of
 1

Project name Pacific Pa	k Drilling Cor	ntractor ESN	Drilling method P	Push-probe rig
Project number 15-059	6-040 Location	~100 feet south-southeast of	Sampling method	5 ft core with plastic liner
Client King County	MW-1		Air monitoring (Y/N)	Yes
HEC rep. Bruce Carp	enter Date N	May 24, 2017	Instrument(s)	Photoionization detector

PID (ppm)	Sample type, interval	%	Water level	Depth (feet, BGS)	Soil	Soil description
(ррпп)	ilitervai	recovery	(feet)	BG3)	group	Gravel at surface
				1		Soil sample PP9-0 collected at 11:50
					GW	Brown-dark brown sandy GRAVEL, FILL, damp
0	5-foot core	20		2		
0	with	20		3		
	liner					wet
				4	<u>▼</u>	Static water level measured at 4.0 feet
				-	4.0	
				5	GW	Dark Brown fine to coarse sandy GRAVEL, small cobbles, wet
				6	0 "	Dark Brown line to coarse sailty GRAVEL, sman coooles, wet
	5-foot	• •		7		
0	core with	20		8		
	liner			0		
				9		
				10		C-111- DD0 10114-1 -4 12:00
				11		Soil sample PP9-10 collected at 12:00
				- 11	SW	Dark Red-brown fine to medium SAND, wet
	5-foot			12	GW	Dark Red-brown fine to coarse sandy GRAVEL, wet
0	core	80		10		
	with liner			13		
	IIIICI			14		
					SW	Dark Brown-black fine to medium SAND, wet
				15	GW	Dark Brown-black sandy GRAVEL, wet
						Set temporary screen from 3 feet to 13 feet. Purged approximately 1-gallon prior to sample collection.
						Collected water sample PP9-W at 12:10
						Backfilled borehole with bentonite chips.

APPENDIX C

Photographic Log



PACIFIC PARK/DUMPSITE ENVIRONMENTAL INVESTIGATION— PHOTOGRAPHIC LOG

Photo Number	Photo Description
1	PP1, looking northwest
2	PP2, looking south
3	PP3, looking north
4	PP4, looking south
5	PP5, looking east
6	PP5, looking east
7	PP6, looking east
8	PP6, looking east
9	PP7, looking south
10	PP8, looking south
11	PP9, looking south



1 PP1, looking northwest



3 PP3, looking north



2 PP2, looking south



4 PP4, looking south





5 PP5, looking east



7 PP6, looking east



6 PP5, looking east



8 PP6, looking east





9 PP7, looking south



11 PP9, looking south



APPENDIX D

Laboratory Analytical Reports





14648 NE 95th Street, Redmond, WA 98052 • (425) 883-3881

May 23, 2017

Bruce Carpenter Herrera Environmental Consultants, Inc. 2200 6th Avenue, Suite 1100 Seattle, WA 98121

Re: Analytical Data for Project 15-05986-040

Laboratory Reference No. 1705-184

Dear Bruce:

Enclosed are the analytical results and associated quality control data for samples submitted on May 12, 2017.

The standard policy of OnSite Environmental, Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

David Baumeister Project Manager

Enclosures

Project: 15-05986-040

Case Narrative

Samples were collected on May 12, 2017 and received by the laboratory on May 12, 2017. They were maintained at the laboratory at a temperature of 2°C to 6°C.

Please note that any and all soil sample results are reported on a dry-weight basis, unless otherwise noted below.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

Project: 15-05986-040

NWTPH-Gx

Matrix: Water
Units: ug/L (ppb)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW1					
Laboratory ID:	05-184-01					
Gasoline	ND	100	NWTPH-Gx	5-19-17	5-19-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	92	61-118				
Client ID:	MW2					
Laboratory ID:	05-184-02					
Gasoline	ND	100	NWTPH-Gx	5-19-17	5-19-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	84	61-118				
Client ID:	MW3					
Laboratory ID:	05-184-03					
Gasoline	ND	100	NWTPH-Gx	5-19-17	5-19-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	85	61-118				
Client ID:	MW4					
Laboratory ID:	05-184-04					
Gasoline	ND	100	NWTPH-Gx	5-19-17	5-19-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	86	61-118				
Client ID:	MW5					
Laboratory ID:	05-184-05					
Gasoline	ND	100	NWTPH-Gx	5-19-17	5-19-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	91	61-118				
Client ID:	MW6					
Laboratory ID:	05-184-06					
Gasoline	ND	100	NWTPH-Gx	5-19-17	5-19-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	82	61-118				

Project: 15-05986-040

NWTPH-Gx

Matrix: Water
Units: ug/L (ppb)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW DUP					
Laboratory ID:	05-184-07					
Gasoline	ND	100	NWTPH-Gx	5-19-17	5-19-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	84	61-118				
Client ID:	Trip Blank					
Laboratory ID:	05-184-09					
Gasoline	ND	100	NWTPH-Gx	5-19-17	5-19-17	
Surrogate:	Percent Recovery	Control Limits	_			
Fluorobenzene	84	61-118				

Project: 15-05986-040

NWTPH-Gx QUALITY CONTROL

Matrix: Water Units: ug/L (ppb)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0519W1					
Gasoline	ND	100	NWTPH-Gx	5-19-17	5-19-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	85	61-118				

Analyte	Res	sult	Spike	Level	Source Result	Pero Reco		Recovery Limits	RPD	RPD Limit	Flags
DUPLICATE											
Laboratory ID:	05-18	34-01									
	ORIG	DUP									
Gasoline	ND	ND	NA	NA		N/	Ą	NA	NA	30	
Surrogate:											
Fluorobenzene						92	85	61-118			

Project: 15-05986-040

NWTPH-Dx

Matrix: Water Units: mg/L (ppm)

Offics. Hig/L (ppiff)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW1					
Laboratory ID:	05-184-01					
Diesel Range Organics	ND	0.26	NWTPH-Dx	5-15-17	5-15-17	X1
Lube Oil Range Organics	ND	0.42	NWTPH-Dx	5-15-17	5-15-17	X1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	109	50-150				
Client ID:	MW2					
Laboratory ID:	05-184-02					
Diesel Range Organics	ND	0.27	NWTPH-Dx	5-15-17	5-15-17	X1
Lube Oil Range Organics	ND	0.44	NWTPH-Dx	5-15-17	5-15-17	X1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	107	50-150				
Client ID:	MW3					
Laboratory ID:	05-184-03					
Diesel Range Organics	ND	0.26	NWTPH-Dx	5-15-17	5-18-17	X1
Lube Oil Range Organics	ND	0.42	NWTPH-Dx	5-15-17 5-15-17	5-18-17	X1 X1
Surrogate:	Percent Recovery	Control Limits	NWII II-DX	0-10-17	0-10-17	XI
o-Terphenyl	91	50-150				
Client ID:	MW4					
Laboratory ID:	05-184-04					
Diesel Range Organics	ND	0.26	NWTPH-Dx	5-15-17	5-15-17	X1
Lube Oil Range Organics	ND -	0.42	NWTPH-Dx	5-15-17	5-15-17	X1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	86	50-150				
Client ID:	MW5					
Laboratory ID:	05-184-05					
Diesel Range Organics	ND	0.26	NWTPH-Dx	5-15-17	5-15-17	X1
Lube Oil Range Organics	ND	0.41	NWTPH-Dx	5-15-17	5-15-17	X1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	80	50-150				
Client ID:	MW6					
Laboratory ID:	05-184-06					
Diesel Range Organics	ND	0.26	NWTPH-Dx	5-15-17	5-15-17	X1
Lube Oil Range Organics	ND ND	0.20	NWTPH-Dx	5-15-17 5-15-17	5-15-17 5-15-17	X1 X1
Surrogate:	Percent Recovery	Control Limits	INVVIIII-DA	J-1J-11	U-1U-11	Λ1
o-Terphenyl	91	50-150				
o respiration	<i>3</i> i	00 100				

Date of Report: May 23, 2017 Samples Submitted: May 12, 2017 Laboratory Reference: 1705-184 Project: 15-05986-040

NWTPH-Dx

Matrix: Water Units: mg/L (ppm)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW DUP					
Laboratory ID:	05-184-07					
Diesel Range Organics	ND	0.27	NWTPH-Dx	5-15-17	5-15-17	X1
Lube Oil Range Organics	ND	0.43	NWTPH-Dx	5-15-17	5-15-17	X1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	93	50-150				

Project: 15-05986-040

NWTPH-Dx QUALITY CONTROL

Matrix: Water
Units: mg/L (ppm)

Result	POI	Method	Date Prepared	Date Analyzed	Flags
rtoun		ou.iou		7.11.0.19200	· iago
MB0515W2					
ND	0.25	NWTPH-Dx	5-15-17	5-15-17	X1
ND	0.40	NWTPH-Dx	5-15-17	5-15-17	X1
Percent Recovery 123	Control Limits 50-150				
	ND ND Percent Recovery	MB0515W2 ND 0.25 ND 0.40 Percent Recovery Control Limits	MB0515W2 ND 0.25 NWTPH-Dx ND 0.40 NWTPH-Dx Percent Recovery Control Limits	Result PQL Method Prepared MB0515W2 ND 0.25 NWTPH-Dx 5-15-17 ND 0.40 NWTPH-Dx 5-15-17 Percent Recovery Control Limits 5-15-17	Result PQL Method Prepared Analyzed MB0515W2 ND 0.25 NWTPH-Dx 5-15-17 5-15-17 ND 0.40 NWTPH-Dx 5-15-17 5-15-17 Percent Recovery Control Limits 5-15-17 5-15-17

Analyta	Por	oul4	Snika	Loval	Source Result	Percent	Recovery	RPD	RPD Limit	Elogo
Analyte	Kes	sult	Бріке	Level	Result	Recovery	Limits	KPU	Limit	Flags
DUPLICATE										
Laboratory ID:	05-19	97-01								
	ORIG	DUP								
Diesel Fuel #2	18.9	7.28	NA	NA		NA	NA	89	NA	X1
Lube Oil	1.31	0.688	NA	NA		NA	NA	62	NA	X1
Surrogate:										
o-Terphenyl						149 94	50-150			

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

Matrix: Water Units: ug/L

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW1					
Laboratory ID:	05-184-01					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Chloromethane	ND	1.0	EPA 8260C	5-16-17	5-16-17	
Vinyl Chloride	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromomethane	ND	0.40	EPA 8260C	5-16-17	5-16-17	
Chloroethane	ND	1.0	EPA 8260C	5-16-17	5-16-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Acetone	ND	5.0	EPA 8260C	5-16-17	5-16-17	
lodomethane	ND	2.1	EPA 8260C	5-16-17	5-16-17	
Carbon Disulfide	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Methylene Chloride	ND	1.0	EPA 8260C	5-16-17	5-16-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Vinyl Acetate	ND	1.0	EPA 8260C	5-16-17	5-16-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Butanone	ND	5.0	EPA 8260C	5-16-17	5-16-17	
Bromochloromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Chloroform	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Benzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Trichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Dibromomethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromodichloromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Chloroethyl Vinyl Ether	ND	11	EPA 8260C	5-16-17	5-16-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	5-16-17	5-16-17	
Toluene	ND	1.0	EPA 8260C	5-16-17	5-16-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-16-17	5-16-17	

Project: 15-05986-040

VOLATILES EPA 8260C page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW1					
Laboratory ID:	05-184-01					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Tetrachloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Hexanone	ND	2.0	EPA 8260C	5-16-17	5-16-17	
Dibromochloromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Chlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Ethylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
m,p-Xylene	ND	0.40	EPA 8260C	5-16-17	5-16-17	
o-Xylene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Styrene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromoform	ND	1.0	EPA 8260C	5-16-17	5-16-17	
Isopropylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1,2,2-Tetrachloroethane	ND	0.27	EPA 8260C	5-16-17	5-16-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
n-Propylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
tert-Butylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
sec-Butylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
n-Butylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	5-16-17	5-16-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Naphthalene	ND	1.3	EPA 8260C	5-16-17	5-16-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	115	77-129				
Toluene-d8	105	80-127				

4-Bromofluorobenzene

80-125

93

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

Matrix: Water Units: ug/L

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW2					
Laboratory ID:	05-184-02					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Chloromethane	ND	1.0	EPA 8260C	5-16-17	5-16-17	
Vinyl Chloride	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromomethane	ND	0.40	EPA 8260C	5-16-17	5-16-17	
Chloroethane	ND	1.0	EPA 8260C	5-16-17	5-16-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Acetone	ND	5.0	EPA 8260C	5-16-17	5-16-17	
Iodomethane	ND	2.1	EPA 8260C	5-16-17	5-16-17	
Carbon Disulfide	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Methylene Chloride	ND	1.0	EPA 8260C	5-16-17	5-16-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Vinyl Acetate	ND	1.0	EPA 8260C	5-16-17	5-16-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Butanone	ND	5.0	EPA 8260C	5-16-17	5-16-17	
Bromochloromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Chloroform	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Benzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Trichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Dibromomethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromodichloromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Chloroethyl Vinyl Ether	ND	11	EPA 8260C	5-16-17	5-16-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	5-16-17	5-16-17	
Toluene	ND	1.0	EPA 8260C	5-16-17	5-16-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-16-17	5-16-17	

Project: 15-05986-040

VOLATILES EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW2					
Laboratory ID:	05-184-02					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Tetrachloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Hexanone	ND	2.0	EPA 8260C	5-16-17	5-16-17	
Dibromochloromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Chlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Ethylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
m,p-Xylene	ND	0.40	EPA 8260C	5-16-17	5-16-17	
o-Xylene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Styrene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromoform	ND	1.0	EPA 8260C	5-16-17	5-16-17	
Isopropylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1,2,2-Tetrachloroethane	ND	0.27	EPA 8260C	5-16-17	5-16-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
n-Propylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
tert-Butylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
sec-Butylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
n-Butylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	5-16-17	5-16-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Naphthalene	ND	1.3	EPA 8260C	5-16-17	5-16-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	110	77-129				

4-Bromofluorobenzene

Toluene-d8

80-127

80-125

107

91

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

Matrix: Water Units: ug/L

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW3					
Laboratory ID:	05-184-03					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Chloromethane	ND	1.0	EPA 8260C	5-16-17	5-16-17	
Vinyl Chloride	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromomethane	ND	0.40	EPA 8260C	5-16-17	5-16-17	
Chloroethane	ND	1.0	EPA 8260C	5-16-17	5-16-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Acetone	ND	5.0	EPA 8260C	5-16-17	5-16-17	
Iodomethane	ND	2.1	EPA 8260C	5-16-17	5-16-17	
Carbon Disulfide	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Methylene Chloride	ND	1.0	EPA 8260C	5-16-17	5-16-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Vinyl Acetate	ND	1.0	EPA 8260C	5-16-17	5-16-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Butanone	ND	5.0	EPA 8260C	5-16-17	5-16-17	
Bromochloromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Chloroform	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Benzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Trichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Dibromomethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromodichloromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Chloroethyl Vinyl Ether	ND	11	EPA 8260C	5-16-17	5-16-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	5-16-17	5-16-17	
Toluene	ND	1.0	EPA 8260C	5-16-17	5-16-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-16-17	5-16-17	

Project: 15-05986-040

VOLATILES EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW3					
Laboratory ID:	05-184-03					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Tetrachloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Hexanone	ND	2.0	EPA 8260C	5-16-17	5-16-17	
Dibromochloromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Chlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Ethylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
m,p-Xylene	ND	0.40	EPA 8260C	5-16-17	5-16-17	
o-Xylene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Styrene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromoform	ND	1.0	EPA 8260C	5-16-17	5-16-17	
Isopropylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1,2,2-Tetrachloroethane	ND	0.27	EPA 8260C	5-16-17	5-16-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
n-Propylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
tert-Butylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
sec-Butylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
n-Butylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	5-16-17	5-16-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Naphthalene	ND	1.3	EPA 8260C	5-16-17	5-16-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	108	77-129				

Dibromofluoromethane 108 77-129
Toluene-d8 105 80-127
4-Bromofluorobenzene 91 80-125



Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW4					
Laboratory ID:	05-184-04					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Chloromethane	ND	1.0	EPA 8260C	5-16-17	5-16-17	
Vinyl Chloride	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromomethane	ND	0.40	EPA 8260C	5-16-17	5-16-17	
Chloroethane	ND	1.0	EPA 8260C	5-16-17	5-16-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Acetone	ND	5.0	EPA 8260C	5-16-17	5-16-17	
Iodomethane	ND	2.1	EPA 8260C	5-16-17	5-16-17	
Carbon Disulfide	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Methylene Chloride	ND	1.0	EPA 8260C	5-16-17	5-16-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Vinyl Acetate	ND	1.0	EPA 8260C	5-16-17	5-16-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Butanone	ND	5.0	EPA 8260C	5-16-17	5-16-17	
Bromochloromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Chloroform	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Benzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Trichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Dibromomethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromodichloromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Chloroethyl Vinyl Ether	ND	11	EPA 8260C	5-16-17	5-16-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	5-16-17	5-16-17	
Toluene	ND	1.0	EPA 8260C	5-16-17	5-16-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-16-17	5-16-17	

Project: 15-05986-040

VOLATILES EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW4					
Laboratory ID:	05-184-04					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Tetrachloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Hexanone	ND	2.0	EPA 8260C	5-16-17	5-16-17	
Dibromochloromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Chlorobenzene	1.5	0.20	EPA 8260C	5-16-17	5-16-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Ethylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
m,p-Xylene	ND	0.40	EPA 8260C	5-16-17	5-16-17	
o-Xylene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Styrene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromoform	ND	1.0	EPA 8260C	5-16-17	5-16-17	
Isopropylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1,2,2-Tetrachloroethane	ND	0.27	EPA 8260C	5-16-17	5-16-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
n-Propylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
tert-Butylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
sec-Butylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
n-Butylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	5-16-17	5-16-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Naphthalene	ND	1.3	EPA 8260C	5-16-17	5-16-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	110	77-129				
-						

4-Bromofluorobenzene

Toluene-d8

80-127

80-125

105

91

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW5					
Laboratory ID:	05-184-05					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Chloromethane	ND	1.0	EPA 8260C	5-16-17	5-16-17	
Vinyl Chloride	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromomethane	ND	0.40	EPA 8260C	5-16-17	5-16-17	
Chloroethane	ND	1.0	EPA 8260C	5-16-17	5-16-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Acetone	ND	5.0	EPA 8260C	5-16-17	5-16-17	
Iodomethane	ND	2.1	EPA 8260C	5-16-17	5-16-17	
Carbon Disulfide	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Methylene Chloride	ND	1.0	EPA 8260C	5-16-17	5-16-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Vinyl Acetate	ND	1.0	EPA 8260C	5-16-17	5-16-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Butanone	ND	5.0	EPA 8260C	5-16-17	5-16-17	
Bromochloromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Chloroform	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Benzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Trichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Dibromomethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromodichloromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Chloroethyl Vinyl Ether	ND	11	EPA 8260C	5-16-17	5-16-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	5-16-17	5-16-17	
Toluene	ND	1.0	EPA 8260C	5-16-17	5-16-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-16-17	5-16-17	

Project: 15-05986-040

VOLATILES EPA 8260C page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW5					
Laboratory ID:	05-184-05					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Tetrachloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Hexanone	ND	2.0	EPA 8260C	5-16-17	5-16-17	
Dibromochloromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Chlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Ethylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
m,p-Xylene	ND	0.40	EPA 8260C	5-16-17	5-16-17	
o-Xylene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Styrene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromoform	ND	1.0	EPA 8260C	5-16-17	5-16-17	
Isopropylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1,2,2-Tetrachloroethane	ND	0.27	EPA 8260C	5-16-17	5-16-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
n-Propylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
tert-Butylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
sec-Butylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
n-Butylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	5-16-17	5-16-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Naphthalene	ND	1.3	EPA 8260C	5-16-17	5-16-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	115	77-129				
Toluene-d8	102	80-127				

4-Bromofluorobenzene

80-125

97

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW6					
Laboratory ID:	05-184-06					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Chloromethane	ND	1.0	EPA 8260C	5-16-17	5-16-17	
Vinyl Chloride	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromomethane	ND	0.40	EPA 8260C	5-16-17	5-16-17	
Chloroethane	ND	1.0	EPA 8260C	5-16-17	5-16-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Acetone	ND	5.0	EPA 8260C	5-16-17	5-16-17	
Iodomethane	ND	2.1	EPA 8260C	5-16-17	5-16-17	
Carbon Disulfide	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Methylene Chloride	ND	1.0	EPA 8260C	5-16-17	5-16-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Vinyl Acetate	ND	1.0	EPA 8260C	5-16-17	5-16-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Butanone	ND	5.0	EPA 8260C	5-16-17	5-16-17	
Bromochloromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Chloroform	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Benzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Trichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Dibromomethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromodichloromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Chloroethyl Vinyl Ether	ND	11	EPA 8260C	5-16-17	5-16-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	5-16-17	5-16-17	
Toluene	ND	1.0	EPA 8260C	5-16-17	5-16-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-16-17	5-16-17	

Project: 15-05986-040

VOLATILES EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW6					
Laboratory ID:	05-184-06					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Tetrachloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Hexanone	ND	2.0	EPA 8260C	5-16-17	5-16-17	
Dibromochloromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Chlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Ethylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
m,p-Xylene	ND	0.40	EPA 8260C	5-16-17	5-16-17	
o-Xylene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Styrene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromoform	ND	1.0	EPA 8260C	5-16-17	5-16-17	
Isopropylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1,2,2-Tetrachloroethane	ND	0.27	EPA 8260C	5-16-17	5-16-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
n-Propylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
tert-Butylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
sec-Butylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
n-Butylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	5-16-17	5-16-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Naphthalene	ND	1.3	EPA 8260C	5-16-17	5-16-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	110	77-129				

 Dibromofluoromethane
 110
 77-129

 Toluene-d8
 104
 80-127

 4-Bromofluorobenzene
 95
 80-125



Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW DUP					
Laboratory ID:	05-184-07					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Chloromethane	ND	1.0	EPA 8260C	5-16-17	5-16-17	
Vinyl Chloride	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromomethane	ND	0.40	EPA 8260C	5-16-17	5-16-17	
Chloroethane	ND	1.0	EPA 8260C	5-16-17	5-16-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Acetone	ND	5.0	EPA 8260C	5-16-17	5-16-17	
Iodomethane	ND	2.1	EPA 8260C	5-16-17	5-16-17	
Carbon Disulfide	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Methylene Chloride	ND	1.0	EPA 8260C	5-16-17	5-16-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Vinyl Acetate	ND	1.0	EPA 8260C	5-16-17	5-16-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Butanone	ND	5.0	EPA 8260C	5-16-17	5-16-17	
Bromochloromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Chloroform	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Benzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Trichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Dibromomethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromodichloromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Chloroethyl Vinyl Ether	ND	11	EPA 8260C	5-16-17	5-16-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	5-16-17	5-16-17	
Toluene	ND	1.0	EPA 8260C	5-16-17	5-16-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-16-17	5-16-17	

Project: 15-05986-040

VOLATILES EPA 8260C page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW DUP					
Laboratory ID:	05-184-07					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Tetrachloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Hexanone	ND	2.0	EPA 8260C	5-16-17	5-16-17	
Dibromochloromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Chlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Ethylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
m,p-Xylene	ND	0.40	EPA 8260C	5-16-17	5-16-17	
o-Xylene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Styrene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromoform	ND	1.0	EPA 8260C	5-16-17	5-16-17	
Isopropylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1,2,2-Tetrachloroethane	ND	0.27	EPA 8260C	5-16-17	5-16-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
n-Propylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
tert-Butylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
sec-Butylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
n-Butylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	5-16-17	5-16-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Naphthalene	ND	1.3	EPA 8260C	5-16-17	5-16-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	112	77-129				
Talvana dO	400	00.407				

 Dibromofluoromethane
 112
 77-129

 Toluene-d8
 102
 80-127

 4-Bromofluorobenzene
 93
 80-125



Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	Trip Blank					
Laboratory ID:	05-184-09					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Chloromethane	ND	1.0	EPA 8260C	5-16-17	5-16-17	
Vinyl Chloride	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromomethane	ND	0.40	EPA 8260C	5-16-17	5-16-17	
Chloroethane	ND	1.0	EPA 8260C	5-16-17	5-16-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Acetone	ND	5.0	EPA 8260C	5-16-17	5-16-17	
Iodomethane	ND	2.1	EPA 8260C	5-16-17	5-16-17	
Carbon Disulfide	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Methylene Chloride	ND	1.0	EPA 8260C	5-16-17	5-16-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Vinyl Acetate	ND	1.0	EPA 8260C	5-16-17	5-16-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Butanone	ND	5.0	EPA 8260C	5-16-17	5-16-17	
Bromochloromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Chloroform	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Benzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Trichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Dibromomethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromodichloromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Chloroethyl Vinyl Ether	ND	11	EPA 8260C	5-16-17	5-16-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	5-16-17	5-16-17	
Toluene	ND	1.0	EPA 8260C	5-16-17	5-16-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-16-17	5-16-17	

Project: 15-05986-040

VOLATILES EPA 8260C page 2 of 2

Analyte Result PQL Method Prepared Analyzed Flags					Date	Date	
Laboratory ID:	Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
1,1,2-Trichloroethane ND 0.20 EPA 8260C 5-16-17 5-16-17 Tetrachloroethene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,3-Dichloropropane ND 0.20 EPA 8260C 5-16-17 5-16-17 2-Hexanone ND 0.20 EPA 8260C 5-16-17 5-16-17 1,1-2-Dibromoethane ND 0.20 EPA 8260C 5-16-17 5-16-17 Chlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Ethylsenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Styrene ND 0.20 EPA 8260C 5-16-17 5-16-17 Styrene ND 0.20 EPA 8260C 5-16-17 5-16-17	Client ID:	Trip Blank					
Tetrachloroethene	Laboratory ID:	05-184-09					
1,3-Dichloropropane ND 0.20 EPA 8260C 5-16-17 5-16-17 2-Hexanone ND 2.0 EPA 8260C 5-16-17 5-16-17 Dibromochloromethane ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Dibromoethane ND 0.20 EPA 8260C 5-16-17 5-16-17 Chlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,1,1,2-Tetrachloroethane ND 0.20 EPA 8260C 5-16-17 5-16-17 Ethylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Thylene ND 0.40 EPA 8260C 5-16-17 5-16-17 Styrene ND 0.20 EPA 8260C 5-16-17 5-16-17 Styrene ND 0.20 EPA 8260C 5-16-17 5-16-17 Isopropylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Isopropylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17	1,1,2-Trichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Hexanone ND 2.0 EPA 8260C 5-16-17 5-16-17 Dibromochloromethane ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Dibromoethane ND 0.20 EPA 8260C 5-16-17 5-16-17 1,1,1,2-Tetrachloroethane ND 0.20 EPA 8260C 5-16-17 5-16-17 1,1,1,2-Tetrachloroethane ND 0.20 EPA 8260C 5-16-17 5-16-17 Ethylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Ethylbenzene ND 0.40 EPA 8260C 5-16-17 5-16-17 Syrlene ND 0.20 EPA 8260C 5-16-17 5-16-17 Styrene ND 0.20 EPA 8260C 5-16-17 5-16-17 Bromoform ND 1.0 EPA 8260C 5-16-17 5-16-17 Bromobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Bromobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 <t< td=""><td>Tetrachloroethene</td><td>ND</td><td>0.20</td><td>EPA 8260C</td><td>5-16-17</td><td>5-16-17</td><td></td></t<>	Tetrachloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Dibromochloromethane ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Dibromoethane ND 0.20 EPA 8260C 5-16-17 5-16-17 Chlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,1,1,2-Tetrachloroethane ND 0.20 EPA 8260C 5-16-17 5-16-17 Ethylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Ethylbenzene ND 0.40 EPA 8260C 5-16-17 5-16-17 O-Xylene ND 0.20 EPA 8260C 5-16-17 5-16-17 Styrene ND 0.20 EPA 8260C 5-16-17 5-16-17 Bromoform ND 1.0 EPA 8260C 5-16-17 5-16-17 Isopropylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Bromobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Horopylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17	1,3-Dichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dibromoethane ND 0.20 EPA 8260C 5-16-17 5-16-17 Chlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,1,1,2-Tetrachloroethane ND 0.20 EPA 8260C 5-16-17 5-16-17 Ethylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 mp-Xylene ND 0.40 EPA 8260C 5-16-17 5-16-17 o-Xylene ND 0.20 EPA 8260C 5-16-17 5-16-17 Styrene ND 0.20 EPA 8260C 5-16-17 5-16-17 Bromoform ND 0.20 EPA 8260C 5-16-17 5-16-17 Bromobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Bromobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Bromobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Horrich Stylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17	2-Hexanone	ND	2.0	EPA 8260C	5-16-17	5-16-17	
Chlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,1,1,2-Tetrachloroethane ND 0.20 EPA 8260C 5-16-17 5-16-17 Ethylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 mp-Xylene ND 0.40 EPA 8260C 5-16-17 5-16-17 o-Xylene ND 0.20 EPA 8260C 5-16-17 5-16-17 Styrene ND 0.20 EPA 8260C 5-16-17 5-16-17 Bromoform ND 1.0 EPA 8260C 5-16-17 5-16-17 Bromobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Bromobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,1,2,2-Tetrachloroethane ND 0.20 EPA 8260C 5-16-17 5-16-17 1,1,2,3-Triholroforopane ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2,3-Triholroforolulene ND 0.20 EPA 8260C 5-16-17 5-16-17	Dibromochloromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1,1,2-Tetrachloroethane ND 0.20 EPA 8260C 5-16-17 5-16-17 Ethylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 m,p-Xylene ND 0.40 EPA 8260C 5-16-17 5-16-17 o-Xylene ND 0.20 EPA 8260C 5-16-17 5-16-17 Styrene ND 0.20 EPA 8260C 5-16-17 5-16-17 Bromoform ND 1.0 EPA 8260C 5-16-17 5-16-17 Isopropylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Bromobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Bromobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Interpopylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1-2,3-Trichloropropane ND 0.20 EPA 8260C 5-16-17 5-16-17 1-2,3-Trichloropropane ND 0.20 EPA 8260C 5-16-17 5-16-17	1,2-Dibromoethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Ethylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 m,p-Xylene ND 0.40 EPA 8260C 5-16-17 5-16-17 o-Xylene ND 0.20 EPA 8260C 5-16-17 5-16-17 Styrene ND 0.20 EPA 8260C 5-16-17 5-16-17 Bromoform ND 1.0 EPA 8260C 5-16-17 5-16-17 Isopropylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Bromobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Bromobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,1,2,2-Tetrachloroethane ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2,3-Trichloropropane ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2,3-Trichloropropane ND 0.20 EPA 8260C 5-16-17 5-16-17 2-Chlorotoluene ND 0.20 EPA 8260C 5-16-17 5-16-17	Chlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
m.p-Xylene ND 0.40 EPA 8260C 5-16-17 5-16-17 o-Xylene ND 0.20 EPA 8260C 5-16-17 5-16-17 Styrene ND 0.20 EPA 8260C 5-16-17 5-16-17 Bromoform ND 1.0 EPA 8260C 5-16-17 5-16-17 Isopropylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Bromobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,1,2,2-Tetrachloroethane ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2,3-Trichloropropane ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2,3-Trichloropropane ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2,3-Trichloropropane ND 0.20 EPA 8260C 5-16-17 5-16-17 2-Chlorotoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 4-Chlorotoluene ND 0.20 EPA 8260C 5-16-17 5-16-17	1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
o-Xylene ND 0.20 EPA 8260C 5-16-17 5-16-17 Styrene ND 0.20 EPA 8260C 5-16-17 5-16-17 Bromoform ND 1.0 EPA 8260C 5-16-17 5-16-17 Isopropylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Bromobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Bromobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Bromobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,1,2,2-Tetrachloroethane ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2,3-Trichloropropane ND 0.20 EPA 8260C 5-16-17 5-16-17 n-Propylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 -Chlorotoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,3,5-Trimethylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17	Ethylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Styrene ND 0.20 EPA 8260C 5-16-17 5-16-17 Bromoform ND 1.0 EPA 8260C 5-16-17 5-16-17 Isopropylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Bromobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,1,2,2-Tetrachloroethane ND 0.27 EPA 8260C 5-16-17 5-16-17 1,2,3-Trichloropropane ND 0.20 EPA 8260C 5-16-17 5-16-17 n-Propylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 2-Chlorotoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 4-Chlorotoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 4-Chlorotoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 4-Chlorotoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,3-5-Trimethylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17<	m,p-Xylene	ND	0.40	EPA 8260C	5-16-17	5-16-17	
Bromoform ND 1.0 EPA 8260C 5-16-17 5-16-17 Isopropylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Bromobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,1,2,2-Tetrachloroethane ND 0.27 EPA 8260C 5-16-17 5-16-17 1,2,3-Trichloropropane ND 0.20 EPA 8260C 5-16-17 5-16-17 n-Propylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 2-Chlorotoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 4-Chlorotoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 4-Chlorotoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,3,5-Trimethylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2,4-Trimethylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,3-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17	o-Xylene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Isopropylbenzene	Styrene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,1,2,2-Tetrachloroethane ND 0.27 EPA 8260C 5-16-17 5-16-17 1,2,3-Trichloropropane ND 0.20 EPA 8260C 5-16-17 5-16-17 n-Propylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 2-Chlorotoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 4-Chlorotoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 4-Chlorotoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,3,5-Trimethylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2,4-Trimethylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Q-Trimethylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,4-Dichlorobenzene ND 0.20 EPA 8260C <	Bromoform	ND	1.0	EPA 8260C	5-16-17	5-16-17	
1,1,2,2-Tetrachloroethane ND 0.27 EPA 8260C 5-16-17 5-16-17 1,2,3-Trichloropropane ND 0.20 EPA 8260C 5-16-17 5-16-17 n-Propylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 2-Chlorotoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 4-Chlorotoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 4-Chlorotoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,3,5-Trimethylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2,4-Trimethylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Lichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,3-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,4-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Dichlorobenzene ND 0.20 EPA 8260C	Isopropylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2,3-Trichloropropane ND 0.20 EPA 8260C 5-16-17 5-16-17 n-Propylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 2-Chlorotoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 4-Chlorotoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,3,5-Trimethylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 tert-Butylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2,4-Trimethylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 sec-Butylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,3-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,4-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Dibromo-3-chloropropane ND 0.20 EPA 8260C	Bromobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
n-Propylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 2-Chlorotoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 4-Chlorotoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,3,5-Trimethylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 tert-Butylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2,4-Trimethylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 sec-Butylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,3-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,3-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,4-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Dibromo-3-chloropropane ND 1.0 EPA 8260C	1,1,2,2-Tetrachloroethane	ND	0.27	EPA 8260C	5-16-17	5-16-17	
2-Chlorotoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 4-Chlorotoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,3,5-Trimethylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 tert-Butylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2,4-Trimethylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 sec-Butylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,3-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,3-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,4-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Dibromo-3-chloropropane ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2,4-Trichlorobenzene ND 0.20 EPA 8260C	1,2,3-Trichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
4-Chlorotoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,3,5-Trimethylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 tert-Butylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2,4-Trimethylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 sec-Butylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,3-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 p-Isopropyltoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,4-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 n-Butylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Dibromo-3-chloropropane ND 1.0 EPA 8260C 5-16-17 5-16-17 1,2,4-Trichlorobenzene ND 0.20 EPA 8260C	n-Propylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
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tert-Butylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2,4-Trimethylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 sec-Butylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,3-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 p-Isopropyltoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,4-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 n-Butylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Dibromo-3-chloropropane ND 1.0 EPA 8260C 5-16-17 5-16-17 1,2,4-Trichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Hexachlorobutadiene ND 0.20 EPA 8260C 5-16-17 5-16-17 ND 1.3 EPA 8260C 5-16-17 5-1	4-Chlorotoluene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2,4-Trimethylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 sec-Butylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,3-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 p-Isopropyltoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,4-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 n-Butylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Dibromo-3-chloropropane ND 1.0 EPA 8260C 5-16-17 5-16-17 1,2,4-Trichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Naphthalene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2,3-Trichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17	1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
sec-Butylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,3-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 p-Isopropyltoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,4-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 n-Butylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Dibromo-3-chloropropane ND 1.0 EPA 8260C 5-16-17 5-16-17 1,2,4-Trichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Hexachlorobutadiene ND 0.20 EPA 8260C 5-16-17 5-16-17 Naphthalene ND 1.3 EPA 8260C 5-16-17 5-16-17 1,2,3-Trichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17	tert-Butylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,3-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 p-Isopropyltoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,4-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 n-Butylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Dibromo-3-chloropropane ND 1.0 EPA 8260C 5-16-17 5-16-17 1,2,4-Trichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Hexachlorobutadiene ND 0.20 EPA 8260C 5-16-17 5-16-17 Naphthalene ND 1.3 EPA 8260C 5-16-17 5-16-17 1,2,3-Trichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17	1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
p-Isopropyltoluene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,4-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 n-Butylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Dibromo-3-chloropropane ND 1.0 EPA 8260C 5-16-17 5-16-17 1,2,4-Trichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Hexachlorobutadiene ND 0.20 EPA 8260C 5-16-17 5-16-17 Naphthalene ND 1.3 EPA 8260C 5-16-17 5-16-17 1,2,3-Trichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17	sec-Butylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,4-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 n-Butylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Dibromo-3-chloropropane ND 1.0 EPA 8260C 5-16-17 5-16-17 1,2,4-Trichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Hexachlorobutadiene ND 0.20 EPA 8260C 5-16-17 5-16-17 Naphthalene ND 1.3 EPA 8260C 5-16-17 5-16-17 1,2,3-Trichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17	1,3-Dichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 n-Butylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Dibromo-3-chloropropane ND 1.0 EPA 8260C 5-16-17 5-16-17 1,2,4-Trichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Hexachlorobutadiene ND 0.20 EPA 8260C 5-16-17 5-16-17 Naphthalene ND 1.3 EPA 8260C 5-16-17 5-16-17 1,2,3-Trichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17	p-Isopropyltoluene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
n-Butylbenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 1,2-Dibromo-3-chloropropane ND 1.0 EPA 8260C 5-16-17 5-16-17 1,2,4-Trichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Hexachlorobutadiene ND 0.20 EPA 8260C 5-16-17 5-16-17 Naphthalene ND 1.3 EPA 8260C 5-16-17 5-16-17 1,2,3-Trichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17	1,4-Dichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dibromo-3-chloropropane ND 1.0 EPA 8260C 5-16-17 5-16-17 1,2,4-Trichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Hexachlorobutadiene ND 0.20 EPA 8260C 5-16-17 5-16-17 Naphthalene ND 1.3 EPA 8260C 5-16-17 5-16-17 1,2,3-Trichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17	1,2-Dichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2,4-Trichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17 Hexachlorobutadiene ND 0.20 EPA 8260C 5-16-17 5-16-17 Naphthalene ND 1.3 EPA 8260C 5-16-17 5-16-17 1,2,3-Trichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17	n-Butylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Hexachlorobutadiene ND 0.20 EPA 8260C 5-16-17 5-16-17 Naphthalene ND 1.3 EPA 8260C 5-16-17 5-16-17 1,2,3-Trichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17	1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	5-16-17	5-16-17	
Naphthalene ND 1.3 EPA 8260C 5-16-17 5-16-17 1,2,3-Trichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17	1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2,3-Trichlorobenzene ND 0.20 EPA 8260C 5-16-17 5-16-17	Hexachlorobutadiene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
• •	Naphthalene	ND	1.3	EPA 8260C	5-16-17	5-16-17	
Currentel Paraut Paraut Paraut I imite	1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Surrogate. Percent Recovery Control Limits	Surrogate:	Percent Recovery	Control Limits				_
Dibromofluoromethane 114 77-129	Dibromofluoromethane	114	77-129				
Toluene-d8 104 80-127	Toluene-d8	104	80-127				

4-Bromofluorobenzene

80-125

95

Project: 15-05986-040

VOLATILES by EPA 8260C METHOD BLANK QUALITY CONTROL

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0516W1					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Chloromethane	ND	1.0	EPA 8260C	5-16-17	5-16-17	
Vinyl Chloride	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromomethane	ND	0.40	EPA 8260C	5-16-17	5-16-17	
Chloroethane	ND	1.0	EPA 8260C	5-16-17	5-16-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Acetone	ND	5.0	EPA 8260C	5-16-17	5-16-17	
lodomethane	ND	2.1	EPA 8260C	5-16-17	5-16-17	
Carbon Disulfide	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Methylene Chloride	ND	1.0	EPA 8260C	5-16-17	5-16-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Vinyl Acetate	ND	1.0	EPA 8260C	5-16-17	5-16-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Butanone	ND	5.0	EPA 8260C	5-16-17	5-16-17	
Bromochloromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Chloroform	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Benzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Trichloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Dibromomethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromodichloromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Chloroethyl Vinyl Ether	ND	11	EPA 8260C	5-16-17	5-16-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	5-16-17	5-16-17	
Toluene	ND	1.0	EPA 8260C	5-16-17	5-16-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-16-17	5-16-17	

Project: 15-05986-040

VOLATILES by EPA 8260C METHOD BLANK QUALITY CONTROL

page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Analyto	rtoouit	. 42	mourou	rioparoa	Analyzou	1 lugo
Laboratory ID:	MB0516W1					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Tetrachloroethene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Hexanone	ND	2.0	EPA 8260C	5-16-17	5-16-17	
Dibromochloromethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Chlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Ethylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
m,p-Xylene	ND	0.40	EPA 8260C	5-16-17	5-16-17	
o-Xylene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Styrene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromoform	ND	1.0	EPA 8260C	5-16-17	5-16-17	
Isopropylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Bromobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,1,2,2-Tetrachloroethane	ND	0.27	EPA 8260C	5-16-17	5-16-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	5-16-17	5-16-17	
n-Propylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
tert-Butylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
sec-Butylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
n-Butylbenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	5-16-17	5-16-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Naphthalene	ND	1.3	EPA 8260C	5-16-17	5-16-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	5-16-17	5-16-17	
Surrogate:	Percent Recovery	Control Limits				

Surrogate: Percent Recovery Control Limit
Dibromofluoromethane 116 77-129
Toluene-d8 105 80-127
4-Bromofluorobenzene 89 80-125



Project: 15-05986-040

VOLATILES by EPA 8260C MS/MSD QUALITY CONTROL

					Source	Per	cent	Recovery		RPD	
Analyte	Result		Spike	Level	Result	Recovery		Limits	RPD	Limit	Flags
MATRIX SPIKES											
Laboratory ID:	05-17	76-01									
	MS	MSD	MS	MSD		MS	MSD				
1,1-Dichloroethene	9.76	9.54	10.0	10.0	ND	98	95	65-119	2	15	
Benzene	10.1	10.4	10.0	10.0	ND	101	104	75-117	3	15	
Trichloroethene	8.95	9.04	10.0	10.0	ND	90	90	66-120	1	15	
Toluene	10.2	10.5	10.0	10.0	ND	102	105	79-120	3	15	
Chlorobenzene	9.48	9.47	10.0	10.0	ND	95	95	76-120	0	15	
Surrogate:											
Dibromofluoromethane						121	121	77-129			
Toluene-d8						102	106	80-127			
4-Bromofluorobenzene						96	92	80-125			

Project: 15-05986-040

PAHs EPA 8270D/SIM

·				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW1					
Laboratory ID:	05-184-01					
Naphthalene	ND	0.095	EPA 8270D/SIM	5-16-17	5-17-17	
2-Methylnaphthalene	ND	0.095	EPA 8270D/SIM	5-16-17	5-17-17	
1-Methylnaphthalene	ND	0.095	EPA 8270D/SIM	5-16-17	5-17-17	
Acenaphthylene	ND	0.095	EPA 8270D/SIM	5-16-17	5-17-17	
Acenaphthene	ND	0.095	EPA 8270D/SIM	5-16-17	5-17-17	
Fluorene	ND	0.095	EPA 8270D/SIM	5-16-17	5-17-17	
Phenanthrene	ND	0.095	EPA 8270D/SIM	5-16-17	5-17-17	
Anthracene	ND	0.095	EPA 8270D/SIM	5-16-17	5-17-17	
Fluoranthene	ND	0.095	EPA 8270D/SIM	5-16-17	5-17-17	
Pyrene	ND	0.095	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo[a]anthracene	ND	0.0095	EPA 8270D/SIM	5-16-17	5-17-17	
Chrysene	ND	0.0095	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo[b]fluoranthene	ND	0.0095	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo(j,k)fluoranthene	ND	0.0095	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo[a]pyrene	ND	0.0095	EPA 8270D/SIM	5-16-17	5-17-17	
Indeno(1,2,3-c,d)pyrene	ND	0.0095	EPA 8270D/SIM	5-16-17	5-17-17	
Dibenz[a,h]anthracene	ND	0.0095	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo[g,h,i]perylene	ND	0.0095	EPA 8270D/SIM	5-16-17	5-17-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	46	30 - 124				
Pyrene-d10	45	40 - 143				

Project: 15-05986-040

PAHs EPA 8270D/SIM

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW2	·		•	•	
Laboratory ID:	05-184-02					
Naphthalene	ND	0.099	EPA 8270D/SIM	5-16-17	5-17-17	
2-Methylnaphthalene	ND	0.099	EPA 8270D/SIM	5-16-17	5-17-17	
1-Methylnaphthalene	ND	0.099	EPA 8270D/SIM	5-16-17	5-17-17	
Acenaphthylene	ND	0.099	EPA 8270D/SIM	5-16-17	5-17-17	
Acenaphthene	ND	0.099	EPA 8270D/SIM	5-16-17	5-17-17	
Fluorene	ND	0.099	EPA 8270D/SIM	5-16-17	5-17-17	
Phenanthrene	ND	0.099	EPA 8270D/SIM	5-16-17	5-17-17	
Anthracene	ND	0.099	EPA 8270D/SIM	5-16-17	5-17-17	
Fluoranthene	ND	0.099	EPA 8270D/SIM	5-16-17	5-17-17	
Pyrene	ND	0.099	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo[a]anthracene	ND	0.0099	EPA 8270D/SIM	5-16-17	5-17-17	
Chrysene	ND	0.0099	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo[b]fluoranthene	ND	0.0099	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo(j,k)fluoranthene	ND	0.0099	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo[a]pyrene	ND	0.0099	EPA 8270D/SIM	5-16-17	5-17-17	
Indeno(1,2,3-c,d)pyrene	ND	0.0099	EPA 8270D/SIM	5-16-17	5-17-17	
Dibenz[a,h]anthracene	ND	0.0099	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo[g,h,i]perylene	ND	0.0099	EPA 8270D/SIM	5-16-17	5-17-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	88	30 - 124				
Pyrene-d10	87	40 - 143				
Tamahana dada	405	07 407				

Project: 15-05986-040

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW3					
Laboratory ID:	05-184-03					
Naphthalene	ND	0.095	EPA 8270D/SIM	5-16-17	5-18-17	
2-Methylnaphthalene	ND	0.095	EPA 8270D/SIM	5-16-17	5-18-17	
1-Methylnaphthalene	ND	0.095	EPA 8270D/SIM	5-16-17	5-18-17	
Acenaphthylene	ND	0.095	EPA 8270D/SIM	5-16-17	5-18-17	
Acenaphthene	ND	0.095	EPA 8270D/SIM	5-16-17	5-18-17	
Fluorene	ND	0.095	EPA 8270D/SIM	5-16-17	5-18-17	
Phenanthrene	ND	0.095	EPA 8270D/SIM	5-16-17	5-18-17	
Anthracene	ND	0.095	EPA 8270D/SIM	5-16-17	5-18-17	
Fluoranthene	ND	0.095	EPA 8270D/SIM	5-16-17	5-18-17	
Pyrene	ND	0.095	EPA 8270D/SIM	5-16-17	5-18-17	
Benzo[a]anthracene	ND	0.0095	EPA 8270D/SIM	5-16-17	5-18-17	
Chrysene	ND	0.0095	EPA 8270D/SIM	5-16-17	5-18-17	
Benzo[b]fluoranthene	ND	0.0095	EPA 8270D/SIM	5-16-17	5-18-17	
Benzo(j,k)fluoranthene	ND	0.0095	EPA 8270D/SIM	5-16-17	5-18-17	
Benzo[a]pyrene	ND	0.0095	EPA 8270D/SIM	5-16-17	5-18-17	
Indeno(1,2,3-c,d)pyrene	ND	0.0095	EPA 8270D/SIM	5-16-17	5-18-17	
Dibenz[a,h]anthracene	ND	0.0095	EPA 8270D/SIM	5-16-17	5-18-17	
Benzo[g,h,i]perylene	ND	0.0095	EPA 8270D/SIM	5-16-17	5-18-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	78	30 - 124				
Pyrene-d10	77	40 - 143				
Torphonyl d11	വാ	27 127				

Date

Date

Date of Report: May 23, 2017 Samples Submitted: May 12, 2017 Laboratory Reference: 1705-184

Project: 15-05986-040

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW4					
Laboratory ID:	05-184-04					
Naphthalene	ND	0.096	EPA 8270D/SIM	5-16-17	5-18-17	
2-Methylnaphthalene	ND	0.096	EPA 8270D/SIM	5-16-17	5-18-17	
1-Methylnaphthalene	ND	0.096	EPA 8270D/SIM	5-16-17	5-18-17	
Acenaphthylene	ND	0.096	EPA 8270D/SIM	5-16-17	5-18-17	
Acenaphthene	ND	0.096	EPA 8270D/SIM	5-16-17	5-18-17	
Fluorene	ND	0.096	EPA 8270D/SIM	5-16-17	5-18-17	
Phenanthrene	ND	0.096	EPA 8270D/SIM	5-16-17	5-18-17	
Anthracene	ND	0.096	EPA 8270D/SIM	5-16-17	5-18-17	
Fluoranthene	ND	0.096	EPA 8270D/SIM	5-16-17	5-18-17	
Pyrene	ND	0.096	EPA 8270D/SIM	5-16-17	5-18-17	
Benzo[a]anthracene	ND	0.0096	EPA 8270D/SIM	5-16-17	5-18-17	
Chrysene	ND	0.0096	EPA 8270D/SIM	5-16-17	5-18-17	
Benzo[b]fluoranthene	ND	0.0096	EPA 8270D/SIM	5-16-17	5-18-17	
Benzo(j,k)fluoranthene	ND	0.0096	EPA 8270D/SIM	5-16-17	5-18-17	
Benzo[a]pyrene	ND	0.0096	EPA 8270D/SIM	5-16-17	5-18-17	
Indeno(1,2,3-c,d)pyrene	ND	0.0096	EPA 8270D/SIM	5-16-17	5-18-17	
Dibenz[a,h]anthracene	ND	0.0096	EPA 8270D/SIM	5-16-17	5-18-17	
Benzo[g,h,i]perylene	ND	0.0096	EPA 8270D/SIM	5-16-17	5-18-17	
Surrogate:	Percent Recovery	Control Limits			_	
2-Fluorobiphenyl	104	30 - 124				
Pyrene-d10	84	40 - 143				

Project: 15-05986-040

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW5					
Laboratory ID:	05-184-05					
Naphthalene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
2-Methylnaphthalene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
1-Methylnaphthalene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Acenaphthylene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Acenaphthene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Fluorene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Phenanthrene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Anthracene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Fluoranthene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Pyrene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Chrysene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	58	30 - 124				
Pyrene-d10	60	40 - 143				

Date

Date

Date of Report: May 23, 2017 Samples Submitted: May 12, 2017 Laboratory Reference: 1705-184

Project: 15-05986-040

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW6					
Laboratory ID:	05-184-06					
Naphthalene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
2-Methylnaphthalene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
1-Methylnaphthalene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Acenaphthylene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Acenaphthene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Fluorene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Phenanthrene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Anthracene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Fluoranthene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Pyrene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Chrysene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Surrogate:	Percent Recovery	Control Limits	_		_	
2-Fluorobiphenyl	80	30 - 124				
Pyrene-d10	79	40 - 143				

Project: 15-05986-040

PAHs EPA 8270D/SIM

Matrix: Water Units: ug/L

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW DUP					
Laboratory ID:	05-184-07					
Naphthalene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
2-Methylnaphthalene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
1-Methylnaphthalene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Acenaphthylene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Acenaphthene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Fluorene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Phenanthrene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Anthracene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Fluoranthene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Pyrene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Chrysene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	94	30 - 124				
Pyrene-d10	90	40 - 143				

Terphenyl-d14 27 - 127 111

Project: 15-05986-040

PAHS EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

Matrix: Water Units: ug/L

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0516W2					
Naphthalene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
2-Methylnaphthalene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
1-Methylnaphthalene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Acenaphthylene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Acenaphthene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Fluorene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Phenanthrene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Anthracene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Fluoranthene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Pyrene	ND	0.10	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Chrysene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270D/SIM	5-16-17	5-17-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	64	30 - 124				
Pyrene-d10	84	40 - 143				

27 - 127

99

Terphenyl-d14

Project: 15-05986-040

PAHS EPA 8270D/SIM SB/SBD QUALITY CONTROL

					P	ercent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Re	covery	Limits	RPD	Limit	Flags
SPIKE BLANKS										
Laboratory ID:	SB05	16W2								
	SB	SBD	SB	SBD	SE	SBD				
Naphthalene	0.402	0.392	0.500	0.500	80	78	29 - 101	3	47	
Acenaphthylene	0.435	0.454	0.500	0.500	87	91	20 - 117	4	50	
Acenaphthene	0.479	0.461	0.500	0.500	96	92	37 - 109	4	43	
Fluorene	0.453	0.486	0.500	0.500	91	97	47 - 108	7	34	
Phenanthrene	0.451	0.476	0.500	0.500	90	95	49 - 109	5	28	
Anthracene	0.436	0.458	0.500	0.500	87	92	34 - 140	5	32	
Fluoranthene	0.477	0.488	0.500	0.500	95	98	45 - 120	2	39	
Pyrene	0.575	0.493	0.500	0.500	11	5 99	42 - 133	15	39	
Benzo[a]anthracene	0.521	0.523	0.500	0.500	10-	105	71 - 117	0	28	
Chrysene	0.495	0.493	0.500	0.500	99	99	53 - 110	0	25	
Benzo[b]fluoranthene	0.504	0.500	0.500	0.500	10	100	53 - 123	1	37	
Benzo(j,k)fluoranthene	0.504	0.509	0.500	0.500	10	102	52 - 119	1	41	
Benzo[a]pyrene	0.462	0.466	0.500	0.500	92	93	37 - 129	1	33	
Indeno(1,2,3-c,d)pyrene	0.507	0.500	0.500	0.500	10	100	45 - 128	1	31	
Dibenz[a,h]anthracene	0.511	0.498	0.500	0.500	10:	2 100	54 - 120	3	30	
Benzo[g,h,i]perylene	0.498	0.494	0.500	0.500	100	99	49 - 117	1	29	
Surrogate:										
2-Fluorobiphenyl					86	81	30 - 124			
Pyrene-d10					91	91	40 - 143			
Terphenyl-d14					11.	3 100	27 - 127			

Project: 15-05986-040

TOTAL METALS EPA 200.8/7470A

	- " ' '			Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-184-01 MW1					
Antimony	ND	5.6	200.8	5-17-17	5-17-17	
Arsenic	ND	3.3	200.8	5-17-17	5-17-17	
Beryllium	ND	11	200.8	5-17-17	5-17-17	
Cadmium	ND	4.4	200.8	5-17-17	5-17-17	
Chromium	ND	11	200.8	5-17-17	5-17-17	
Copper	ND	11	200.8	5-17-17	5-17-17	
Lead	ND	1.1	200.8	5-17-17	5-17-17	
Mercury	ND	0.50	7470A	5-16-17	5-16-17	
Nickel	ND	22	200.8	5-17-17	5-17-17	
Selenium	ND	5.6	200.8	5-17-17	5-17-17	
Silver	ND	11	200.8	5-17-17	5-17-17	
Thallium	ND	5.6	200.8	5-17-17	5-17-17	
Zinc	160	28	200.8	5-17-17	5-17-17	

Project: 15-05986-040

TOTAL METALS EPA 200.8/7470A

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-184-02					
Client ID:	MW2					
Antimony	ND	5.6	200.8	5-17-17	5-17-17	
Arsenic	ND	3.3	200.8	5-17-17	5-17-17	
Beryllium	ND	11	200.8	5-17-17	5-17-17	
Cadmium	ND	4.4	200.8	5-17-17	5-17-17	
Chromium	ND	11	200.8	5-17-17	5-17-17	
Copper	ND	11	200.8	5-17-17	5-17-17	
Lead	ND	1.1	200.8	5-17-17	5-17-17	
Mercury	ND	0.50	7470A	5-16-17	5-16-17	
Nickel	ND	22	200.8	5-17-17	5-17-17	
Selenium	ND	5.6	200.8	5-17-17	5-17-17	
Silver	ND	11	200.8	5-17-17	5-17-17	
Thallium	ND	5.6	200.8	5-17-17	5-17-17	
Zinc	42	28	200.8	5-17-17	5-17-17	

Project: 15-05986-040

TOTAL METALS EPA 200.8/7470A

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-184-03					
Client ID:	MW3					
Antimony	ND	5.6	200.8	5-17-17	5-17-17	
Arsenic	ND	3.3	200.8	5-17-17	5-17-17	
Beryllium	ND	11	200.8	5-17-17	5-17-17	
Cadmium	ND	4.4	200.8	5-17-17	5-17-17	
Chromium	ND	11	200.8	5-17-17	5-17-17	
Copper	ND	11	200.8	5-17-17	5-17-17	
Lead	ND	1.1	200.8	5-17-17	5-17-17	
Mercury	ND	0.50	7470A	5-16-17	5-16-17	
Nickel	ND	22	200.8	5-17-17	5-17-17	
Selenium	ND	5.6	200.8	5-17-17	5-17-17	
Silver	ND	11	200.8	5-17-17	5-17-17	
Thallium	ND	5.6	200.8	5-17-17	5-17-17	
Zinc	170	28	200.8	5-17-17	5-17-17	

Project: 15-05986-040

TOTAL METALS EPA 200.8/7470A

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-184-04					
Client ID:	MW4					
Antimony	ND	5.6	200.8	5-17-17	5-17-17	
Arsenic	8.3	3.3	200.8	5-17-17	5-17-17	
Beryllium	ND	11	200.8	5-17-17	5-17-17	
Cadmium	ND	4.4	200.8	5-17-17	5-17-17	
Chromium	ND	11	200.8	5-17-17	5-17-17	
Copper	ND	11	200.8	5-17-17	5-17-17	
Lead	ND	1.1	200.8	5-17-17	5-17-17	
Mercury	ND	0.50	7470A	5-16-17	5-16-17	
Nickel	ND	22	200.8	5-17-17	5-17-17	
Selenium	ND	5.6	200.8	5-17-17	5-17-17	
Silver	ND	11	200.8	5-17-17	5-17-17	
Thallium	ND	5.6	200.8	5-17-17	5-17-17	
Zinc	80	28	200.8	5-17-17	5-17-17	

Project: 15-05986-040

TOTAL METALS EPA 200.8/7470A

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-184-05					
Client ID:	MW5					
Antimony	ND	5.6	200.8	5-17-17	5-17-17	
Arsenic	ND	3.3	200.8	5-17-17	5-17-17	
Beryllium	ND	11	200.8	5-17-17	5-17-17	
Cadmium	ND	4.4	200.8	5-17-17	5-17-17	
Chromium	ND	11	200.8	5-17-17	5-17-17	
Copper	ND	11	200.8	5-17-17	5-17-17	
Lead	4.0	1.1	200.8	5-17-17	5-17-17	
Mercury	ND	0.50	7470A	5-16-17	5-16-17	
Nickel	ND	22	200.8	5-17-17	5-17-17	
Selenium	ND	5.6	200.8	5-17-17	5-17-17	
Silver	ND	11	200.8	5-17-17	5-17-17	
Thallium	ND	5.6	200.8	5-17-17	5-17-17	
Zinc	37	28	200.8	5-17-17	5-17-17	

Project: 15-05986-040

TOTAL METALS EPA 200.8/7470A

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-184-06					
Client ID:	MW6					
Antimony	ND	5.6	200.8	5-17-17	5-17-17	
Arsenic	ND	3.3	200.8	5-17-17	5-17-17	
Beryllium	ND	11	200.8	5-17-17	5-17-17	
Cadmium	ND	4.4	200.8	5-17-17	5-17-17	
Chromium	ND	11	200.8	5-17-17	5-17-17	
Copper	ND	11	200.8	5-17-17	5-17-17	
Lead	ND	1.1	200.8	5-17-17	5-17-17	
Mercury	ND	0.50	7470A	5-16-17	5-16-17	
Nickel	ND	22	200.8	5-17-17	5-17-17	
Selenium	ND	5.6	200.8	5-17-17	5-17-17	
Silver	ND	11	200.8	5-17-17	5-17-17	
Thallium	ND	5.6	200.8	5-17-17	5-17-17	
Zinc	ND	28	200.8	5-17-17	5-17-17	

Project: 15-05986-040

TOTAL METALS EPA 200.8/7470A

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-184-07					
Client ID:	MW DUP					
Antimony	ND	5.6	200.8	5-17-17	5-17-17	
Arsenic	ND	3.3	200.8	5-17-17	5-17-17	
Beryllium	ND	11	200.8	5-17-17	5-17-17	
Cadmium	ND	4.4	200.8	5-17-17	5-17-17	
Chromium	ND	11	200.8	5-17-17	5-17-17	
Copper	ND	11	200.8	5-17-17	5-17-17	
Lead	ND	1.1	200.8	5-17-17	5-17-17	
Mercury	ND	0.50	7470A	5-16-17	5-16-17	
Nickel	ND	22	200.8	5-17-17	5-17-17	
Selenium	ND	5.6	200.8	5-17-17	5-17-17	
Silver	ND	11	200.8	5-17-17	5-17-17	
Thallium	ND	5.6	200.8	5-17-17	5-17-17	
Zinc	35	28	200.8	5-17-17	5-17-17	

Project: 15-05986-040

TOTAL METALS EPA 200.8/7470A METHOD BLANK QUALITY CONTROL

Date Extracted: 5-16&17-17

Date Analyzed: 5-16&17-17

Matrix: Water
Units: ug/L (ppb)

Lab ID: MB0516W1&MB0517WM1

Analyte	Method	Result	PQL
Antimony	200.8	ND	5.6
Arsenic	200.8	ND	3.3
Beryllium	200.8	ND	11
Cadmium	200.8	ND	4.4
Chromium	200.8	ND	11
Copper	200.8	ND	11
Lead	200.8	ND	1.1
Mercury	7470A	ND	0.50
Nickel	200.8	ND	22
Selenium	200.8	ND	5.6
Silver	200.8	ND	11
Thallium	200.8	ND	5.6
Zinc	200.8	ND	28

Project: 15-05986-040

TOTAL METALS EPA 200.8/7470A DUPLICATE QUALITY CONTROL

Date Extracted: 5-16&17-17

Date Analyzed: 5-16&17-17

Matrix: Water
Units: ug/L (ppb)

Lab ID: 05-184-01

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Antimony	ND	ND	NA	5.6	
Arsenic	ND	ND	NA	3.3	
Beryllium	ND	ND	NA	11	
Cadmium	ND	ND	NA	4.4	
Chromium	ND	ND	NA	11	
Copper	ND	ND	NA	11	
Lead	ND	ND	NA	1.1	
Mercury	ND	ND	NA	0.50	
Nickel	ND	ND	NA	22	
Selenium	ND	ND	NA	5.6	
Silver	ND	ND	NA	11	
Thallium	ND	ND	NA	5.6	
Zinc	159	155	3	28	

Project: 15-05986-040

TOTAL METALS EPA 200.8/7470A MS/MSD QUALITY CONTROL

Date Extracted: 5-16&17-17

Date Analyzed: 5-16&17-17

Matrix: Water
Units: ug/L (ppb)

Lab ID: 05-184-01

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Antimony	222	225	101	222	100	1	
Arsenic	222	215	97	216	97	0	
Beryllium	222	227	102	223	100	2	
Cadmium	222	223	100	214	97	4	
Chromium	222	208	94	207	93	0	
Copper	222	216	97	216	97	0	
Lead	222	213	96	213	96	0	
Mercury	12.5	12.5	100	11.4	91	9	
Nickel	222	209	94	204	92	2	
Selenium	222	216	97	215	97	1	
Silver	222	211	95	208	94	2	
Thallium	222	207	93	206	93	1	
Zinc	222	367	94	371	95	1	

Project: 15-05986-040

DISSOLVED METALS EPA 200.8/7470A

	,			Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-184-01 MW1					
Antimony	ND	5.0	200.8		5-16-17	
Arsenic	ND	3.0	200.8		5-16-17	
Beryllium	ND	10	200.8		5-16-17	
Cadmium	ND	4.0	200.8		5-16-17	
Chromium	ND	10	200.8		5-16-17	
Copper	ND	10	200.8		5-16-17	
Lead	ND	1.0	200.8		5-16-17	
Mercury	ND	0.50	7470A		5-16-17	
Nickel	ND	20	200.8		5-16-17	
Selenium	ND	5.0	200.8		5-16-17	
Silver	ND	10	200.8		5-16-17	
Thallium	ND	5.0	200.8		5-16-17	
Zinc	170	25	200.8		5-16-17	

Project: 15-05986-040

DISSOLVED METALS EPA 200.8/7470A

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-184-02					
Client ID:	MW2					
Antimony	ND	5.0	200.8		5-16-17	
Arsenic	ND	3.0	200.8		5-16-17	
Beryllium	ND	10	200.8		5-16-17	
Cadmium	ND	4.0	200.8		5-16-17	
Chromium	ND	10	200.8		5-16-17	
Copper	ND	10	200.8		5-16-17	
Lead	ND	1.0	200.8		5-16-17	
Mercury	ND	0.50	7470A		5-16-17	
Nickel	ND	20	200.8		5-16-17	
Selenium	ND	5.0	200.8		5-16-17	
Silver	ND	10	200.8		5-16-17	
Thallium	ND	5.0	200.8		5-16-17	
Zinc	ND	25	200.8		5-16-17	

Project: 15-05986-040

DISSOLVED METALS EPA 200.8/7470A

	- " ' '			Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-184-03					
Client ID:	MW3					
Antimony	ND	5.0	200.8		5-16-17	
Arsenic	ND	3.0	200.8		5-16-17	
Beryllium	ND	10	200.8		5-16-17	
Cadmium	ND	4.0	200.8		5-16-17	
Chromium	ND	10	200.8		5-16-17	
Copper	ND	10	200.8		5-16-17	
Lead	ND	1.0	200.8		5-16-17	
Mercury	ND	0.50	7470A		5-16-17	
Nickel	ND	20	200.8		5-16-17	
Selenium	ND	5.0	200.8		5-16-17	
Silver	ND	10	200.8		5-16-17	
Thallium	ND	5.0	200.8		5-16-17	
Zinc	140	25	200.8		5-16-17	

Project: 15-05986-040

DISSOLVED METALS EPA 200.8/7470A

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-184-04					
Client ID:	MW4					
Antimony	ND	5.0	200.8		5-16-17	
Arsenic	8.3	3.0	200.8		5-16-17	
Beryllium	ND	10	200.8		5-16-17	
Cadmium	ND	4.0	200.8		5-16-17	
Chromium	ND	10	200.8		5-16-17	
Copper	ND	10	200.8		5-16-17	
Lead	ND	1.0	200.8		5-16-17	
Mercury	ND	0.50	7470A		5-16-17	
Nickel	ND	20	200.8		5-16-17	
Selenium	ND	5.0	200.8		5-16-17	
Silver	ND	10	200.8		5-16-17	
Thallium	ND	5.0	200.8		5-16-17	
Zinc	37	25	200.8		5-16-17	

Project: 15-05986-040

DISSOLVED METALS EPA 200.8/7470A

	- " ' '			Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-184-05					
Client ID:	MW5					
Antimony	ND	5.0	200.8		5-16-17	
Arsenic	ND	3.0	200.8		5-16-17	
Beryllium	ND	10	200.8		5-16-17	
Cadmium	ND	4.0	200.8		5-16-17	
Chromium	ND	10	200.8		5-16-17	
Copper	ND	10	200.8		5-16-17	
Lead	3.0	1.0	200.8		5-16-17	
Mercury	ND	0.50	7470A		5-16-17	
Nickel	ND	20	200.8		5-16-17	
Selenium	ND	5.0	200.8		5-16-17	
Silver	ND	10	200.8		5-16-17	
Thallium	ND	5.0	200.8		5-16-17	
Zinc	35	25	200.8		5-16-17	

Project: 15-05986-040

DISSOLVED METALS EPA 200.8/7470A

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-184-06					
Client ID:	MW6					
Antimony	ND	5.0	200.8		5-16-17	
Arsenic	ND	3.0	200.8		5-16-17	
Beryllium	ND	10	200.8		5-16-17	
Cadmium	ND	4.0	200.8		5-16-17	
Chromium	ND	10	200.8		5-16-17	
Copper	ND	10	200.8		5-16-17	
Lead	ND	1.0	200.8		5-16-17	
Mercury	ND	0.50	7470A		5-16-17	
Nickel	ND	20	200.8		5-16-17	
Selenium	ND	5.0	200.8		5-16-17	
Silver	ND	10	200.8		5-16-17	
Thallium	ND	5.0	200.8		5-16-17	
Zinc	ND	25	200.8		5-16-17	

Project: 15-05986-040

DISSOLVED METALS EPA 200.8/7470A

	- " ' '			Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-184-07					
Client ID:	MW DUP					
Antimony	ND	5.0	200.8		5-16-17	
Arsenic	ND	3.0	200.8		5-16-17	
Beryllium	ND	10	200.8		5-16-17	
Cadmium	ND	4.0	200.8		5-16-17	
Chromium	ND	10	200.8		5-16-17	
Copper	ND	10	200.8		5-16-17	
Lead	ND	1.0	200.8		5-16-17	
Mercury	ND	0.50	7470A		5-16-17	
Nickel	ND	20	200.8		5-16-17	
Selenium	ND	5.0	200.8		5-16-17	
Silver	ND	10	200.8		5-16-17	
Thallium	ND	5.0	200.8		5-16-17	
Zinc	ND	25	200.8		5-16-17	

Project: 15-05986-040

DISSOLVED METALS EPA 200.8/7470A

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-184-08					
Client ID:	Filter Blank					
Antimony	ND	5.0	200.8		5-16-17	
Arsenic	ND	3.0	200.8		5-16-17	
Beryllium	ND	10	200.8		5-16-17	
Cadmium	ND	4.0	200.8		5-16-17	
Chromium	ND	10	200.8		5-16-17	
Copper	ND	10	200.8		5-16-17	
Lead	ND	1.0	200.8		5-16-17	
Mercury	ND	0.50	7470A		5-16-17	
Nickel	ND	20	200.8		5-16-17	
Selenium	ND	5.0	200.8		5-16-17	
Silver	ND	10	200.8		5-16-17	
Thallium	ND	5.0	200.8		5-16-17	
Zinc	ND	25	200.8		5-16-17	

Project: 15-05986-040

DISSOLVED METALS EPA 200.8/7470A METHOD BLANK QUALITY CONTROL

Date Analyzed: 5-16-17

Matrix: Water Units: ug/L (ppb)

Lab ID: MB0516D1&MB0516D2

Analyte	Method	Result	PQL
Antimony	200.8	ND	5.0
Arsenic	200.8	ND	3.0
Beryllium	200.8	ND	10
Cadmium	200.8	ND	4.0
Chromium	200.8	ND	10
Copper	200.8	ND	10
Lead	200.8	ND	1.0
Mercury	7470A	ND	0.50
Nickel	200.8	ND	20
Selenium	200.8	ND	5.0
Silver	200.8	ND	10
Thallium	200.8	ND	5.0
Zinc	200.8	ND	25

Project: 15-05986-040

DISSOLVED METALS EPA 200.8/7470A DUPLICATE QUALITY CONTROL

Date Analyzed: 5-16-17

Matrix: Water
Units: ug/L (ppb)

Lab ID: 05-184-01

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Antimony	ND	ND	NA	5.0	
Arsenic	ND	ND	NA	3.0	
Beryllium	ND	ND	NA	10	
Cadmium	ND	ND	NA	4.0	
Chromium	ND	ND	NA	10	
Copper	ND	ND	NA	10	
Lead	ND	ND	NA	1.0	
Mercury	ND	ND	NA	0.5	
Nickel	ND	ND	NA	20	
Selenium	ND	ND	NA	5.0	
Silver	ND	ND	NA	10	
Thallium	ND	ND	NA	5.0	
Zinc	172	167	3	25	

Project: 15-05986-040

DISSOLVED METALS EPA 200.8/7470A MS/MSD QUALITY CONTROL

Date Analyzed: 5-16-17

Matrix: Water
Units: ug/L (ppb)

Lab ID: 05-184-01

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Antimony	200	197	99	209	105	6	
Arsenic	200	195	98	203	101	4	
Beryllium	200	193	96	207	103	7	
Cadmium	200	198	99	208	104	5	
Chromium	200	180	90	191	95	6	
Copper	200	197	99	205	102	4	
Lead	200	186	93	199	100	7	
Mercury	12.5	10.9	87	10.8	86	1	
Nickel	200	184	92	195	97	6	
Selenium	200	204	102	217	109	6	
Silver	200	185	92	196	98	6	
Thallium	200	191	95	205	103	7	
Zinc	200	365	97	376	102	3	



Data Qualifiers and Abbreviations

- A Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
- B The analyte indicated was also found in the blank sample.
- C The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
- E The value reported exceeds the quantitation range and is an estimate.
- F Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
- H The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
- I Compound recovery is outside of the control limits.
- J The value reported was below the practical quantitation limit. The value is an estimate.
- K Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
- L The RPD is outside of the control limits.
- M Hydrocarbons in the gasoline range are impacting the diesel range result.
- M1 Hydrocarbons in the gasoline range (toluene-napthalene) are present in the sample.
- N Hydrocarbons in the lube oil range are impacting the diesel range result.
- N1 Hydrocarbons in diesel range are impacting lube oil range results.
- O Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
- P The RPD of the detected concentrations between the two columns is greater than 40.
- Q Surrogate recovery is outside of the control limits.
- S Surrogate recovery data is not available due to the necessary dilution of the sample.
- T The sample chromatogram is not similar to a typical .
- U The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- U1 The practical quantitation limit is elevated due to interferences present in the sample.
- V Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
- W Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
- X Sample extract treated with a mercury cleanup procedure.
- X1- Sample extract treated with a Sulfuric acid/Silica gel cleanup procedure.
- Y The calibration verification for this analyte exceeded the 20% drift specified in method 8260C, and therefore the reported result should be considered an estimate. The overall performance of the calibration verification standard met the acceptance criteria of the method.

7 -

ND - Not Detected at PQL

PQL - Practical Quantitation Limit

RPD - Relative Percent Difference



OnSite Environmental Inc.

Chain of Custody

	9	1
Page	of	

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Frojec	Pacific Patt Manager:	x/L	Sta	ndard (7 Days) H analysis 5 Days	ave)	ço				SGCI		8260C	s Only)	MIS	-level)		ides 8(Pesticides	icides				664A	Metals	PPWHORIS			
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Sampl	ed by: Corge 1 Ane	r Brianna Blaud		(other)	_	er of Con	HCID	H-Gx/BTEX	1-Gx	NWTPH-Dx (PAcid / SG Clean-up)	Volatiles 8260C	nated Vo	EDB EPA 8011 (Waters Only)	ivolatiles 8270D/SIM I low-level PAHs)	PAHs 8270D/SIM (low-level)	082A	Organochlorine	Organophosphorus	Chlorinated Acid Herbicides 81	RCRA Metals	Total MTCA Metals	letals	HEM (oil and grease) 1664A	of PP	Dissolved		1	nre
Lab ID		ample Identification	Date Sampled	Time Sampled	Matrix	Numbe	NWTPH-HCID	NWTPH-(NWTPH-Gx	NWTPH	Volatile	Halogenated	EDB EF	Semivolatiles (with low-level	PAHs 8	PCBs 8082A	Organo	Organo	Chlorina	Total RC	Total M	TCLP Metals	HEM (oi	Total	Diss		tojevy 70	% Moisture
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2	MWZ		9121=	11:10	water	11			X	X	X				X									X	X			
3	MW3		5/12/17	13:43	water	u			X	X	X				X									X	X		T	
4	MW4		5/12/13		water				X	X	X				X									X	X		1	
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14648 NE 95th Street, Redmond, WA 98052 • (425) 883-3881

June 13, 2017

Bruce Carpenter Herrera Environmental Consultants, Inc. 2200 6th Avenue, Suite 1100 Seattle, WA 98121

Re: Analytical Data for Project 15-05986-040

Laboratory Reference No. 1705-299

Dear Bruce:

Enclosed are the analytical results and associated quality control data for samples submitted on May 24, 2017.

The standard policy of OnSite Environmental, Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

David Baumeister Project Manager

Enclosures

Project: 15-05986-040

Case Narrative

Samples were collected on May 23 and 24, 2017 and received by the laboratory on May 24, 2017. They were maintained at the laboratory at a temperature of 2°C to 6°C.

Please note that any and all soil sample results are reported on a dry-weight basis, unless otherwise noted below.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

NWTPH Gx and Volatiles EPA 8260C (soil) Analysis

Per EPA Method 5035A, samples were received by the laboratory in pre-weighed 40 mL VOA vials within 48 hours of sample collection. They were stored in a freezer at between -7°C and -20°C until extraction or analysis.

PAHs EPA 8270D/SIM Analysis

The Spike Blank had one recovery slightly above control limits. Samples PP5-W and PP7-W each had one surrogate recovery out of control limits. This is within allowance of our standard operating procedure as long as the recovery is above 10%.

Please note that any other QA/QC issues associated with these extractions and analyses will be indicated with a footnote reference and discussed in detail on the Data Qualifier page.



Project: 15-05986-040

NWTPH-Gx

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP1-2					
Laboratory ID:	05-299-01					
Gasoline	ND	6.9	NWTPH-Gx	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	94	63-124				
Client ID:	PP1-10					
Laboratory ID:	05-299-03					
Gasoline	ND	8.0	NWTPH-Gx	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	99	63-124				
Client ID:	PP2-2.5					
Laboratory ID:	05-299-04					
Gasoline	ND	7.8	NWTPH-Gx	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	96	63-124				
Client ID:	PP2-10					
Laboratory ID:	05-299-05					
Gasoline	ND	9.0	NWTPH-Gx	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	102	63-124				
Client ID:	PP3-2.5					
Laboratory ID:	05-299-07					
Gasoline	ND	6.6	NWTPH-Gx	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	97	63-124				
Client ID:	PP3-10					
Laboratory ID:	05-299-08					
Gasoline	ND	7.0	NWTPH-Gx	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	92	63-124				

Project: 15-05986-040

NWTPH-Gx

Matrix: Soil

onite. mg/ng (ppm)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP4-3					
Laboratory ID:	05-299-10					
Gasoline	ND	7.4	NWTPH-Gx	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	97	63-124				
Client ID:	PP4-10					
Laboratory ID:	05-299-11					
Gasoline	ND	7.1	NWTPH-Gx	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	92	63-124				
Client ID:	PP5-2.5					
Laboratory ID:	05-299-13					
Gasoline	ND	7.0	NWTPH-Gx	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	100	63-124				
Client ID:	PP5-10					
Laboratory ID:	05-299-14					
Gasoline	ND	8.1	NWTPH-Gx	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	101	63-124				
Client ID:	PP6-2					
Laboratory ID:	05-299-16					
Gasoline	ND	5.7	NWTPH-Gx	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	93	63-124				
Client ID:	PP6-10					
Laboratory ID:	05-299-17					
Gasoline	ND	9.7	NWTPH-Gx	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	111	63-124				
Client ID:	PP7-0					
Laboratory ID:	05-299-19					
Gasoline	ND	6.6	NWTPH-Gx	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	86	63-124				

Project: 15-05986-040

NWTPH-Gx

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP7-10					
Laboratory ID:	05-299-20					
Gasoline	ND	7.8	NWTPH-Gx	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	104	63-124				
Client ID:	PP8-0.5					
Laboratory ID:	05-299-22					
Gasoline	ND	6.1	NWTPH-Gx	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	107	63-124				
Client ID:	PP8-10					
Laboratory ID:	05-299-23					
Gasoline	ND	9.6	NWTPH-Gx	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	113	63-124				
Client ID:	PP9-0					
Laboratory ID:	05-299-25					
Gasoline	ND	5.6	NWTPH-Gx	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	100	63-124				
Client ID:	PP9-10					
Laboratory ID:	05-299-26					
Gasoline	ND	7.0	NWTPH-Gx	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	102	63-124				
Client ID:	PP10-0					
Laboratory ID:	05-299-28					
Gasoline	ND	6.8	NWTPH-Gx	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	87	63-124				

Project: 15-05986-040

NWTPH-Gx QUALITY CONTROL

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0531S1					
Gasoline	ND	5.0	NWTPH-Gx	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	96	63-124				
Laboratory ID:	MB0531S2					
Gasoline	ND	5.0	NWTPH-Gx	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits		·		
Fluorobenzene	96	63-124				

					Source	Perc	cent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Reco	very	Limits	RPD	Limit	Flags
DUPLICATE											
Laboratory ID:	05-29	99-01									
	ORIG	DUP									
Gasoline	ND	ND	NA	NA		N	Α	NA	NA	30	
Surrogate:											_
Fluorobenzene						94	92	63-124			
Laboratory ID:	05-29	99-08									
	ORIG	DUP									
Gasoline	ND	ND	NA	NA		N	Α	NA	NA	30	
Surrogate:		•	•						•		
Fluorobenzene						92	92	63-124			

Project: 15-05986-040

NWTPH-Gx

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP1-W					
Laboratory ID:	05-299-02					
Gasoline	ND	100	NWTPH-Gx	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	95	61-118				
Client ID:	PP2-W					
Laboratory ID:	05-299-06					
Gasoline	ND	100	NWTPH-Gx	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	94	61-118				
Client ID:	PP3-W					
Laboratory ID:	05-299-09					
Gasoline	ND	100	NWTPH-Gx	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	96	61-118				
Client ID:	PP4-W					
Laboratory ID:	05-299-12					
Gasoline	ND	100	NWTPH-Gx	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	96	61-118				
Client ID:	PP5-W					
Laboratory ID:	05-299-15					
Gasoline	210	100	NWTPH-Gx	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	95	61-118				
Client ID:	PP6-W					
Laboratory ID:	05-299-18					
Gasoline	ND	100	NWTPH-Gx	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	97	61-118				

Project: 15-05986-040

NWTPH-Gx

Ginto: 49,2 (pps)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP7-W					
Laboratory ID:	05-299-21					
Gasoline	ND	100	NWTPH-Gx	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	107	61-118				
Client ID:	PP8-W					
Laboratory ID:	05-299-24					
Gasoline	ND	100	NWTPH-Gx	5-26-17	5-26-17	-
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	104	61-118				
Client ID:	PP9-W					
Laboratory ID:	05-299-27					
Gasoline	ND	100	NWTPH-Gx	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	97	61-118				
Client ID:	PP10-W					
Laboratory ID:	05-299-29					
Gasoline	ND	100	NWTPH-Gx	5-26-17	5-26-17	_
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	98	61-118				

Project: 15-05986-040

NWTPH-Gx QUALITY CONTROL

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0526W2					
Gasoline	ND	100	NWTPH-Gx	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	91	61-118				

Analyte	Res	sult	Spike	Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
DUPLICATE										
Laboratory ID:	05-29	92-02								
-	ORIG	DUP								
Gasoline	ND	ND	NA	NA		NA	NA	NA	30	
Surrogate:										
Fluorobenzene						105 95	61-118			

Project: 15-05986-040

NWTPH-Dx

Matrix: Soil

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP1-2				-	
Laboratory ID:	05-299-01					
Diesel Range Organics	ND	31	NWTPH-Dx	5-30-17	5-31-17	X1
Lube Oil Range Organics	ND	62	NWTPH-Dx	5-30-17	5-31-17	X1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	88	50-150				
Client ID:	PP1-10					
Laboratory ID:	05-299-03					
Diesel Range Organics	ND	33	NWTPH-Dx	5-30-17	5-31-17	X1
Lube Oil Range Organics	ND	67	NWTPH-Dx	5-30-17	5-31-17	X1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	105	50-150				
Client ID:	PP2-2.5					
Laboratory ID:	05-299-04					
Diesel Range Organics	ND	33	NWTPH-Dx	5-30-17	5-31-17	X1
Lube Oil Range Organics	ND	66	NWTPH-Dx	5-30-17	5-31-17	X1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	95	50-150				
Client ID:	PP2-10					
Laboratory ID:	05-299-05					
Diesel Range Organics	ND	36	NWTPH-Dx	5-30-17	5-31-17	X1
Lube Oil Range Organics	ND	73	NWTPH-Dx	5-30-17	5-31-17	X1
Surrogate:	Percent Recovery	Control Limits	TWO THE DA	0 00 11	0 0 1 17	7(1
o-Terphenyl	81	50-150				
o respinenty.	0.	00 700				
Client ID:	PP3-2.5					
Laboratory ID:	05-299-07	04	NIM/TOLL D	F 00 47	E 04 47	V/4
Diesel Range Organics Lube Oil	ND 140	31 63	NWTPH-Dx NWTPH-Dx	5-30-17 5-30-17	5-31-17 5-31-17	X1 X1
Surrogate:	Percent Recovery	Control Limits	INVVIFII-DX	3-30-17	3-31-17	Λ1
o-Terphenyl	97	50-150				
Client ID:	PP3-10					
Laboratory ID:	05-299-08					
Diesel Range Organics	ND	31	NWTPH-Dx	5-30-17	5-31-17	X1
Lube Oil Range Organics	ND	62	NWTPH-Dx	5-30-17	5-31-17	X1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	71	50-150				

Project: 15-05986-040

NWTPH-Dx

Matrix: Soil

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP4-3			•	-	-
Laboratory ID:	05-299-10					
Diesel Range Organics	ND	54	NWTPH-Dx	5-30-17	5-31-17	U1,X1
Lube Oil	500	63	NWTPH-Dx	5-30-17	5-31-17	X1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	105	50-150				
Client ID:	PP4-10					
Laboratory ID:	05-299-11					
Diesel Range Organics	ND	55	NWTPH-Dx	5-30-17	5-31-17	U1,X1
Lube Oil	570	62	NWTPH-Dx	5-30-17	5-31-17	X1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	104	50-150				
Client ID:	PP5-2.5					
Laboratory ID:	05-299-13					
Diesel Range Organics	ND	68	NWTPH-Dx	5-30-17	5-31-17	U1,X1
Lube Oil	620	66	NWTPH-Dx	5-30-17	5-31-17	X1
Surrogate:	Percent Recovery	Control Limits	IIII DX	0 00 11	00111	7.1
o-Terphenyl	112	50-150				
Client ID:	PP5-10					
Laboratory ID:	05-299-14					
Diesel Range Organics	ND	34	NWTPH-Dx	5-30-17	5-31-17	X1
Lube Oil Range Organics	ND ND	67	NWTPH-Dx	5-30-17 5-30-17	5-31-17 5-31-17	X1 X1
Surrogate:	Percent Recovery	Control Limits	INVVIFII-DX	3-30-17	3-31-17	Λ1
o-Terphenyl	100	50-150				
о-тегрпенуі	100	30-130				
Client ID:	PP6-2					
Laboratory ID:	05-299-16					
Diesel Range Organics Lube Oil	ND 62	29 58	NWTPH-Dx NWTPH-Dx	5-30-17 5-30-17	5-31-17 5-31-17	X1 X1
Surrogate:	Percent Recovery	Control Limits	INVVII-II-DX	J-JU-11	J-J 1-1 <i>1</i>	A1
o-Terphenyl	110	50-150				
Стагрнануі	110	00-100				
Client ID:	PP6-10					
Laboratory ID:	05-299-17					
Diesel Range Organics	ND	37	NWTPH-Dx	6-2-17	6-2-17	X1
Lube Oil Range Organics	ND	75	NWTPH-Dx	6-2-17	6-2-17	X1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	98	50-150				

Project: 15-05986-040

NWTPH-Dx

Matrix: Soil

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP7-0			-	_	
Laboratory ID:	05-299-19					
Diesel Range Organics	ND	31	NWTPH-Dx	5-30-17	5-31-17	X1
Lube Oil	63	61	NWTPH-Dx	5-30-17	5-31-17	X1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	101	50-150				
Client ID:	PP7-10					
Laboratory ID:	05-299-20					
Diesel Range Organics	ND	31	NWTPH-Dx	5-30-17	5-31-17	X1
Lube Oil Range Organics	ND	63	NWTPH-Dx	5-30-17	5-31-17	X1 X1
Surrogate:	Percent Recovery	Control Limits	INVVII II-DX	3-30-17	3-31-17	Λī
o-Terphenyl	86	50-150				
o respiration	00	00 700				
Client ID:	PP8-0.5					
Laboratory ID:	05-299-22					
Diesel Range Organics	ND	28	NWTPH-Dx	5-30-17	5-31-17	X1
Lube Oil Range Organics	ND	57	NWTPH-Dx	5-30-17	5-31-17	X1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	120	50-150				
Oli 4 ID-	DD0 40					
Client ID:	PP8-10					
Laboratory ID:	05-299-23	200	NIM/TOLL Dec	F 00 47	F 04 47	V/4
Diesel Range Organics	ND	36 71	NWTPH-Dx	5-30-17	5-31-17	X1
Lube Oil Range Organics	ND		NWTPH-Dx	5-30-17	5-31-17	X1
Surrogate:	Percent Recovery 81	Control Limits 50-150				
o-Terphenyl	01	50-150				
Client ID:	PP9-0					
Laboratory ID:	05-299-25					
Diesel Range Organics	ND	26	NWTPH-Dx	5-30-17	5-31-17	X1
Lube Oil Range Organics	ND	52	NWTPH-Dx	5-30-17	5-31-17	X1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	102	50-150				
0" . 1 D	BBC 10					
Client ID:	PP9-10					
Laboratory ID:	05-299-26			:-		
Diesel Range Organics	ND	31	NWTPH-Dx	5-30-17	5-31-17	X1
Lube Oil Range Organics	ND 1.D	62	NWTPH-Dx	5-30-17	5-31-17	X1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	92	50-150				

Project: 15-05986-040

NWTPH-Dx

Matrix: Soil

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP10-0			•	•	
Laboratory ID:	05-299-28					
Diesel Range Organics	ND	31	NWTPH-Dx	5-30-17	5-31-17	X1
Lube Oil Range Organics	ND	61	NWTPH-Dx	5-30-17	5-31-17	X1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	85	50-150				

Project: 15-05986-040

NWTPH-Dx **QUALITY CONTROL**

Matrix: Soil

Units: mg/Kg (ppm)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0530S3					
Diesel Range Organics	ND	25	NWTPH-Dx	5-30-17	5-31-17	X1
Lube Oil Range Organics	ND	50	NWTPH-Dx	5-30-17	5-31-17	X1
Surrogate:	Percent Recovery	Control Limits				•
o-Terphenyl	70	50-150				
Laboratory ID:	MB0602S1					
Diesel Range Organics	ND	25	NWTPH-Dx	6-2-17	6-2-17	X1
Lube Oil Range Organics	ND	50	NWTPH-Dx	6-2-17	6-2-17	X1
Surrogate:	Percent Recovery	Control Limits				•
o-Terphenyl	9.3	50-150				

o-Terphenyl 50-150

					Source	Perce	ent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Recov	ery	Limits	RPD	Limit	Flags
DUPLICATE											
Laboratory ID:	05-29	99-03									
	ORIG	DUP									
Diesel Range	ND	ND	NA	NA		NA		NA	NA	NA	X1
Lube Oil Range	ND	ND	NA	NA		NA		NA	NA	NA	X1
Surrogate:											
o-Terphenyl						105	89	50-150			
Laboratory ID:	05-29	99-17									
	ORIG	DUP									
Diesel Range	ND	ND	NA	NA		NA		NA	NA	NA	X1
Lube Oil Range	ND	ND	NA	NA		NA		NA	NA	NA	X1
Surrogate:											
o-Terphenyl						98	92	50-150			
Laboratory ID:	05-29	99-19									
	ORIG	DUP									
Diesel Range	ND	ND	NA	NA		NA		NA	NA	NA	X1
Lube Oil	51.5	ND	NA	NA		NA	L	NA	NA	NA	X1
Surrogate:											
o-Terphenyl						101	98	50-150			

Project: 15-05986-040

NWTPH-Dx

5 (11 /				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP1-W					
Laboratory ID:	05-299-02					
Diesel Range Organics	ND	0.27	NWTPH-Dx	5-26-17	5-26-17	X1
Lube Oil Range Organics	ND	0.43	NWTPH-Dx	5-26-17	5-26-17	X1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	83	50-150				
Client ID:	PP2-W					
Laboratory ID:	05-299-06					
Diesel Range Organics	ND	0.37	NWTPH-Dx	5-26-17	5-26-17	X1
Lube Oil Range Organics	ND	0.60	NWTPH-Dx	5-26-17 5-26-17	5-26-17 5-26-17	X1 X1
Surrogate:	Percent Recovery	Control Limits	INVVII II-DX	3-20-17	3-20-17	ΧΙ
o-Terphenyl	81	50-150				
o . orphonyr	51	00 100				
Client ID:	PP3-W					
Laboratory ID:	05-299-09					
Diesel Range Organics	ND	0.26	NWTPH-Dx	5-26-17	5-30-17	X1
Lube Oil Range Organics	ND	0.42	NWTPH-Dx	5-26-17	5-30-17	X1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	105	50-150				
011 4 ID	DD 4 144					
Client ID:	PP4-W					
Laboratory ID:	05-299-12	0.00	NIA/TOUR	5.00.47	5.00.47	27.4
Diesel Range Organics	ND	0.26	NWTPH-Dx	5-26-17	5-26-17	X1
Lube Oil Range Organics	ND	0.41	NWTPH-Dx	5-26-17	5-26-17	X1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	96	50-150				
Client ID:	PP5-W					
Laboratory ID:	05-299-15					
Diesel Range Organics	ND	0.27	NWTPH-Dx	5-26-17	5-30-17	X1
Lube Oil Range Organics	ND	0.43	NWTPH-Dx	5-26-17	5-30-17	X1 X1
Surrogate:	Percent Recovery	Control Limits		0 =0 11	0 00 11	
o-Terphenyl	100	50-150				
Client ID:	PP6-W					
Laboratory ID:	05-299-18					
Diesel Range Organics	ND	0.26	NWTPH-Dx	5-26-17	5-26-17	X1
Lube Oil Range Organics	ND	0.42	NWTPH-Dx	5-26-17	5-26-17	X1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	113	50-150				

Project: 15-05986-040

NWTPH-Dx

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP7-W					
Laboratory ID:	05-299-21					
Diesel Range Organics	ND	0.28	NWTPH-Dx	5-26-17	5-26-17	X1
Lube Oil Range Organics	ND	0.44	NWTPH-Dx	5-26-17	5-26-17	X1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	90	50-150				
Client ID:	PP8-W					
Laboratory ID:	05-299-24					
Diesel Range Organics	ND	0.26	NWTPH-Dx	5-26-17	5-26-17	X1
Lube Oil Range Organics	ND	0.41	NWTPH-Dx	5-26-17	5-26-17	X1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	113	50-150				
Client ID:	PP9-W					
Laboratory ID:	05-299-27					
Diesel Range Organics	ND	0.28	NWTPH-Dx	5-26-17	5-26-17	X1
Lube Oil Range Organics	ND	0.45	NWTPH-Dx	5-26-17	5-26-17	X1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	104	50-150				
, ,						
Client ID:	DD40 W					
Client ID:	PP10-W					
Laboratory ID:	05-299-29					
Diesel Range Organics	ND	0.27	NWTPH-Dx	5-26-17	5-26-17	X1
Lube Oil Range Organics	ND	0.43	NWTPH-Dx	5-26-17	5-26-17	X1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	92	50-150				

Project: 15-05986-040

NWTPH-Dx QUALITY CONTROL

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0526W1					
Diesel Range Organics	ND	0.25	NWTPH-Dx	5-26-17	5-26-17	X1
Lube Oil Range Organics	ND	0.40	NWTPH-Dx	5-26-17	5-26-17	X1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	92	50-150				

					Source	Perce	ent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Recov	ery	Limits	RPD	Limit	Flags
DUPLICATE											
Laboratory ID:	05-29	99-02									
	ORIG	DUP									
Diesel Range	ND	ND	NA	NA		NA		NA	NA	NA	X1
Lube Oil Range	ND	ND	NA	NA		NA	ı	NA	NA	NA	X1
Surrogate:											
o-Terphenyl						83	96	50-150			
Laboratory ID:	05-29	99-27									
	ORIG	DUP									
Diesel Range	ND	ND	NA	NA		NA		NA	NA	NA	X1
Lube Oil Range	ND	ND	NA	NA		NA		NA	NA	NA	X1
Surrogate: o-Terphenyl						104	88	50-150			

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

Matrix: Soil Units: mg/kg

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP1-2					
Laboratory ID:	05-299-01					
Dichlorodifluoromethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Chloromethane	ND	0.0057	EPA 8260C	5-25-17	5-25-17	
Vinyl Chloride	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Bromomethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Chloroethane	ND	0.0057	EPA 8260C	5-25-17	5-25-17	
Trichlorofluoromethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,1-Dichloroethene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Acetone	ND	0.011	EPA 8260C	5-25-17	5-25-17	
lodomethane	ND	0.0072	EPA 8260C	5-25-17	5-25-17	
Carbon Disulfide	ND	0.0016	EPA 8260C	5-25-17	5-25-17	
Methylene Chloride	ND	0.011	EPA 8260C	5-25-17	5-25-17	
(trans) 1,2-Dichloroethene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Methyl t-Butyl Ether	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,1-Dichloroethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Vinyl Acetate	ND	0.0057	EPA 8260C	5-25-17	5-25-17	
2,2-Dichloropropane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
(cis) 1,2-Dichloroethene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
2-Butanone	ND	0.0057	EPA 8260C	5-25-17	5-25-17	
Bromochloromethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Chloroform	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,1,1-Trichloroethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Carbon Tetrachloride	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,1-Dichloropropene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Benzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,2-Dichloroethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Trichloroethene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,2-Dichloropropane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Dibromomethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Bromodichloromethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
2-Chloroethyl Vinyl Ether	ND	0.0057	EPA 8260C	5-25-17	5-25-17	
(cis) 1,3-Dichloropropene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Methyl Isobutyl Ketone	ND	0.0057	EPA 8260C	5-25-17	5-25-17	
Toluene	0.019	0.0057	EPA 8260C	5-25-17	5-25-17	
(trans) 1,3-Dichloropropene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	

Project: 15-05986-040

VOLATILES EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP1-2					
Laboratory ID:	05-299-01					
1,1,2-Trichloroethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Γetrachloroethene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,3-Dichloropropane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
2-Hexanone	ND	0.0057	EPA 8260C	5-25-17	5-25-17	
Dibromochloromethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,2-Dibromoethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Chlorobenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,1,1,2-Tetrachloroethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Ethylbenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
n,p-Xylene	ND	0.0023	EPA 8260C	5-25-17	5-25-17	
o-Xylene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Styrene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Bromoform	ND	0.0057	EPA 8260C	5-25-17	5-25-17	
sopropylbenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Bromobenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
I,1,2,2-Tetrachloroethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
I,2,3-Trichloropropane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
n-Propylbenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
2-Chlorotoluene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1-Chlorotoluene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,3,5-Trimethylbenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
ert-Butylbenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,2,4-Trimethylbenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
sec-Butylbenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,3-Dichlorobenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
o-Isopropyltoluene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,4-Dichlorobenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,2-Dichlorobenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
n-Butylbenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,2-Dibromo-3-chloropropane		0.0057	EPA 8260C	5-25-17	5-25-17	
I,2,4-Trichlorobenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Hexachlorobutadiene	ND	0.0057	EPA 8260C	5-25-17	5-25-17	
Naphthalene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,2,3-Trichlorobenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	119	73-134				
Toluene-d8	109	81-124				
1-Bromofluorobenzene	95	80-131				
+ DIGITIONAGIONGINZGING	30	00-131				

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

Matrix: Soil Units: mg/kg

• • •				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP1-10					
Laboratory ID:	05-299-03	0.0040		5.05.47	5.05.47	
Dichlorodifluoromethane	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
Chloromethane	ND	0.0064	EPA 8260C	5-25-17	5-25-17	
Vinyl Chloride	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
Bromomethane	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
Chloroethane	ND	0.0064	EPA 8260C	5-25-17	5-25-17	
Trichlorofluoromethane	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
1,1-Dichloroethene	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
Acetone	0.051	0.013	EPA 8260C	5-25-17	5-25-17	
lodomethane	ND	0.0081	EPA 8260C	5-25-17	5-25-17	
Carbon Disulfide	ND	0.0018	EPA 8260C	5-25-17	5-25-17	
Methylene Chloride	ND	0.013	EPA 8260C	5-25-17	5-25-17	
(trans) 1,2-Dichloroethene	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
Methyl t-Butyl Ether	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
1,1-Dichloroethane	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
Vinyl Acetate	ND	0.0064	EPA 8260C	5-25-17	5-25-17	
2,2-Dichloropropane	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
(cis) 1,2-Dichloroethene	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
2-Butanone	0.012	0.0064	EPA 8260C	5-25-17	5-25-17	
Bromochloromethane	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
Chloroform	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
1,1,1-Trichloroethane	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
Carbon Tetrachloride	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
1,1-Dichloropropene	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
Benzene	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
1,2-Dichloroethane	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
Trichloroethene	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
1,2-Dichloropropane	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
Dibromomethane	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
Bromodichloromethane	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
2-Chloroethyl Vinyl Ether	ND	0.0064	EPA 8260C	5-25-17	5-25-17	
(cis) 1,3-Dichloropropene	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
Methyl Isobutyl Ketone	ND	0.0064	EPA 8260C	5-25-17	5-25-17	
Toluene	0.020	0.0064	EPA 8260C	5-25-17	5-25-17	
(trans) 1,3-Dichloropropene	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
, , ,	-					

Project: 15-05986-040

VOLATILES EPA 8260C page 2 of 2

Analyte Result PQL Method Prepared Analyzed Flags					Date	Date	
Laboratory ID: 05-299-03	Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
1,1,2-Trichloroethane	Client ID:	PP1-10					
Tetrachloroethene ND 0.0013 EPA 8260C 5-25-17 5-25-17 [3,3-Dichloropropane ND 0.0013 EPA 8260C 5-25-17 5-25-17 5-25-17 [5] [5] [5] [5] [5] [5] [5] [5] [5] [5]	Laboratory ID:	05-299-03					
1,3-Dichloropropane ND 0.0013 EPA 8260C 5-25-17 5-25-17 E-Hexanone ND 0.0064 EPA 8260C 5-25-17 5-25-17 E-Hexanone ND 0.0064 EPA 8260C 5-25-17 5-25-17 E-Hexanone ND 0.0013 EPA 8260C 5-25-17 5-25-17 E-Hexanone ND 0.	1,1,2-Trichloroethane	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
Part	Tetrachloroethene	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
Dibromochloromethane	1,3-Dichloropropane	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
1,2-Dibromoethane	2-Hexanone	ND	0.0064	EPA 8260C	5-25-17	5-25-17	
Chlorobenzene	Dibromochloromethane	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
1,1,1,2-Tetrachloroethane	1,2-Dibromoethane	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
Ethylbenzene	Chlorobenzene	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
N.D. N.D. N.D. 0.0026 EPA 8260C 5-25-17 5-	1,1,1,2-Tetrachloroethane	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
ND 0.0013 EPA 8260C 5-25-17 5-25-17	Ethylbenzene	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
Styrene ND 0.0013 EPA 8260C 5-25-17 5-25-17 September ND 0.0064 EPA 8260C 5-25-17 5-25-17 September ND 0.0013 EPA 8260C 5-25-17 September September ND 0.0013 EPA 8260C 5-25-17 September ND 0.0013 EPA 8260C 5-25-17 September ND 0.0013 EPA 8260C 5-25-17 September September ND 0.0013 EPA 8260C 5-25-17 September September ND 0.0013 EPA 8260C September September September ND 0.0013 EPA 8260C September September September ND 0.0013 EPA 8260C September Septembe	n,p-Xylene	ND	0.0026	EPA 8260C	5-25-17	5-25-17	
Second Form ND 0.0064 EPA 8260C 5-25-17 5-25	o-Xylene	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
Sopropy benzene ND	Styrene	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
Sammobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,1,2,2-Tetrachloroethane ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichloropropane ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichloropropane ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2,5-Trimethylbenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2,4-Trimethylbenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,3-Dichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,3-Dichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,4-Dichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,4-Dichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-Dichlorobenzene ND 0.0064 EPA 8260C 5-25-17 5-25-17 1,2-Dichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-2	Bromoform	ND	0.0064	EPA 8260C	5-25-17	5-25-17	
1,1,2,2-Tetrachloroethane	sopropylbenzene	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
1,2,3-Trichloropropane ND 0.0013 EPA 8260C 5-25-17 5-25-17 1-20-10-10-10-10-10-10-10-10-10-10-10-10-10	Bromobenzene	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
ND 0.0013 EPA 8260C 5-25-17 5-25-17 5-25-17 1-2-Chlorotoluene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1-2-Chlorotoluene ND 0.0064 EPA 8260C 5-25-17 5-25-17 1-2-Chlorotoluene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1-2-Chlo	I,1,2,2-Tetrachloroethane	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
C-Chlorotoluene	,2,3-Trichloropropane	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
A-Chlorotoluene ND 0.0013 EPA 8260C 5-25-17 5-25-17 13,3-5-Trimethylbenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 14,2,4-Trimethylbenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 14,2,4-Trimethylbenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 15,2-5-17 15,2	n-Propylbenzene	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
1,3,5-Trimethylbenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2,4-Trimethylbenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2,4-Trimethylbenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,3-Dichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,3-Dichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,4-Dichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,4-Dichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2-Dichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2-Dibromo-3-chloropropane ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene 116 73-134 Toluene-d8 109 81-124 Toluene-d8 109	2-Chlorotoluene	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
ert-Butylbenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 sec-Butylbenzene ND 0.0064 EPA 8260C 5-25-17 5-25-17 sec-Butylbenzene ND 0.0013 EPA 8	1-Chlorotoluene	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
1,2,4-Trimethylbenzene	1,3,5-Trimethylbenzene	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
Sec-Butylbenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,3-Dichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,4-Dichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,4-Dichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2-Dichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2-Dibromo-3-chloropropane ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0064 EPA 8260C 5-25-17 5-25-17 Naphthalene ND 0.0013 EPA 8260	ert-Butylbenzene	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
1,3-Dichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 b-Isopropyltoluene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,4-Dichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2-Dichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 n-Butylbenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2-Dibromo-3-chloropropane ND 0.0064 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 Hexachlorobutadiene ND 0.0064 EPA 8260C 5-25-17 5-25-17 Naphthalene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 116 73-134 Toluene-d8 109	1,2,4-Trimethylbenzene	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
Delsopropyltoluene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,4-Dichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2-Dichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2-Dichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2-Dibromo-3-chloropropane ND 0.0064 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0064 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0064 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0013 EPA 8260C 5-25-	sec-Butylbenzene	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
I,4-Dichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 I,2-Dichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 n-Butylbenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 I,2-Dibromo-3-chloropropane ND 0.0064 EPA 8260C 5-25-17 5-25-17 I,2,4-Trichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 Hexachlorobutadiene ND 0.0064 EPA 8260C 5-25-17 5-25-17 Naphthalene ND 0.0013 EPA 8260C 5-25-17 5-25-17 I,2,3-Trichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 116 73-134 Toluene-d8 109 81-124	1,3-Dichlorobenzene	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
1,2-Dichlorobenzene	o-Isopropyltoluene	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2-Dibromo-3-chloropropane ND 0.0064 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0064 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene Percent Recovery Control Limits Dibromofluoromethane 116 73-134 Toluene-d8 109 81-124	I,4-Dichlorobenzene	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
1,2-Dibromo-3-chloropropane ND 0.0064 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 Hexachlorobutadiene ND 0.0064 EPA 8260C 5-25-17 5-25-17 Naphthalene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 116 73-134 Toluene-d8 109 81-124	I,2-Dichlorobenzene	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
1,2,4-Trichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 Hexachlorobutadiene ND 0.0064 EPA 8260C 5-25-17 5-25-17 Naphthalene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 116 73-134 Toluene-d8 109 81-124	n-Butylbenzene	ND	0.0013	EPA 8260C	5-25-17	5-25-17	
I,2,4-Trichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 Hexachlorobutadiene ND 0.0064 EPA 8260C 5-25-17 5-25-17 Naphthalene ND 0.0013 EPA 8260C 5-25-17 5-25-17 I,2,3-Trichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 116 73-134 Toluene-d8 109 81-124	,2-Dibromo-3-chloropropane	ND	0.0064	EPA 8260C	5-25-17	5-25-17	
Naphthalene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 116 73-134 Toluene-d8 109 81-124			0.0013	EPA 8260C	5-25-17	5-25-17	
Naphthalene ND 0.0013 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0013 EPA 8260C 5-25-17 5-25-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 116 73-134 Toluene-d8 109 81-124	Hexachlorobutadiene	ND	0.0064	EPA 8260C	5-25-17	5-25-17	
ND 0.0013 EPA 8260C 5-25-17 5-25-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 116 73-134 Toluene-d8 109 81-124							
Surrogate: Percent Recovery Control Limits Dibromofluoromethane 116 73-134 Toluene-d8 109 81-124	•	ND	0.0013	EPA 8260C	5-25-17		
Dibromofluoromethane 116 73-134 Toluene-d8 109 81-124							
Toluene-d8 109 81-124							

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

Matrix: Soil Units: mg/kg

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP2-2.5					
Laboratory ID:	05-299-04					
Dichlorodifluoromethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Chloromethane	ND	0.0062	EPA 8260C	5-26-17	5-26-17	
Vinyl Chloride	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Bromomethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Chloroethane	ND	0.0062	EPA 8260C	5-26-17	5-26-17	
Trichlorofluoromethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Acetone	0.071	0.012	EPA 8260C	5-26-17	5-26-17	
lodomethane	ND	0.0062	EPA 8260C	5-26-17	5-26-17	
Carbon Disulfide	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Methylene Chloride	ND	0.012	EPA 8260C	5-26-17	5-26-17	
(trans) 1,2-Dichloroethene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Methyl t-Butyl Ether	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Vinyl Acetate	ND	0.0062	EPA 8260C	5-26-17	5-26-17	
2,2-Dichloropropane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
(cis) 1,2-Dichloroethene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
2-Butanone	0.019	0.0062	EPA 8260C	5-26-17	5-26-17	
Bromochloromethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Chloroform	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,1,1-Trichloroethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Carbon Tetrachloride	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloropropene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Benzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloroethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Trichloroethene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloropropane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Dibromomethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Bromodichloromethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
2-Chloroethyl Vinyl Ether	ND	0.0079	EPA 8260C	5-26-17	5-26-17	
(cis) 1,3-Dichloropropene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Methyl Isobutyl Ketone	ND	0.0062	EPA 8260C	5-26-17	5-26-17	
Toluene	0.025	0.0062	EPA 8260C	5-26-17	5-26-17	
(trans) 1,3-Dichloropropene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	

Project: 15-05986-040

VOLATILES EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP2-2.5					
Laboratory ID:	05-299-04					
1,1,2-Trichloroethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Tetrachloroethene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,3-Dichloropropane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
2-Hexanone	ND	0.0062	EPA 8260C	5-26-17	5-26-17	
Dibromochloromethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromoethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Chlorobenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,1,1,2-Tetrachloroethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Ethylbenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
m,p-Xylene	ND	0.0025	EPA 8260C	5-26-17	5-26-17	
o-Xylene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Styrene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Bromoform	ND	0.0062	EPA 8260C	5-26-17	5-26-17	
Isopropylbenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Bromobenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,1,2,2-Tetrachloroethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichloropropane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
n-Propylbenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
2-Chlorotoluene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
4-Chlorotoluene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,3,5-Trimethylbenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
tert-Butylbenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trimethylbenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
sec-Butylbenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,3-Dichlorobenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
p-Isopropyltoluene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,4-Dichlorobenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,2-Dichlorobenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
n-Butylbenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromo-3-chloropropane		0.0062	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trichlorobenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Hexachlorobutadiene	ND	0.0062	EPA 8260C	5-26-17	5-26-17	
Naphthalene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichlorobenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits		0 20 11	0 20 11	
Dibromofluoromethane	111	73-134				
Toluene-d8	101	73-13 4 81-124				
1 Oluene-uo 4-Bromofluorobenzene	96	80-131				
+-DIOITIOIIUOIODEIIZEITE	90	00-131				

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

Matrix: Soil Units: mg/kg

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP2-10					
Laboratory ID:	05-299-05					
Dichlorodifluoromethane	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
Chloromethane	ND	0.0069	EPA 8260C	5-25-17	5-25-17	
Vinyl Chloride	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
Bromomethane	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
Chloroethane	ND	0.0069	EPA 8260C	5-25-17	5-25-17	
Trichlorofluoromethane	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
1,1-Dichloroethene	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
Acetone	0.088	0.014	EPA 8260C	5-25-17	5-25-17	
lodomethane	ND	0.0087	EPA 8260C	5-25-17	5-25-17	
Carbon Disulfide	ND	0.0019	EPA 8260C	5-25-17	5-25-17	
Methylene Chloride	ND	0.014	EPA 8260C	5-25-17	5-25-17	
(trans) 1,2-Dichloroethene	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
Methyl t-Butyl Ether	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
1,1-Dichloroethane	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
Vinyl Acetate	ND	0.0069	EPA 8260C	5-25-17	5-25-17	
2,2-Dichloropropane	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
(cis) 1,2-Dichloroethene	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
2-Butanone	0.023	0.0069	EPA 8260C	5-25-17	5-25-17	
Bromochloromethane	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
Chloroform	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
1,1,1-Trichloroethane	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
Carbon Tetrachloride	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
1,1-Dichloropropene	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
Benzene	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
1,2-Dichloroethane	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
Trichloroethene	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
1,2-Dichloropropane	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
Dibromomethane	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
Bromodichloromethane	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
2-Chloroethyl Vinyl Ether	ND	0.0069	EPA 8260C	5-25-17	5-25-17	
(cis) 1,3-Dichloropropene	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
Methyl Isobutyl Ketone	ND	0.0069	EPA 8260C	5-25-17	5-25-17	
Toluene	0.019	0.0069	EPA 8260C	5-25-17	5-25-17	
(trans) 1,3-Dichloropropene	ND	0.0014	EPA 8260C	5-25-17	5-25-17	

Project: 15-05986-040

VOLATILES EPA 8260C

page 2 of 2

Cilent ID:					Date	Date	
Laboratory ID: 05-299-05 1,1,2-Trichloroethane ND		Result	PQL	Method	Prepared	Analyzed	Flags
1,1,2-Trichloroethane							
Tetrachloroethene ND 0.0014 EPA 8260C 5-25-17 5-25-17 [3,3-Dichloropropane] ND 0.0014 EPA 8260C 5-25-17 5-25-17 5-25-17 [2,3-Dichloropropane] ND 0.0069 EPA 8260C 5-25-17 5-25-17 [3,2-Dichloropropane] ND 0.0014 EPA 8260C 5-25-17 5-25-17 [3,1,1,2-Tetrachloroethane] ND 0.0014 EPA 8260C 5-25-17 5-25-17 [3,1,1,2-Tetrachloroethane] ND 0.0014 EPA 8260C 5-25-17 5-25-17 [3,1,1,2-Tetrachloroethane] ND 0.0014 EPA 8260C 5-25-17 5-25-17 [3,2-Tetrachloroethane] ND 0.0014 EPA 8260C 5-25-17 5-25-17 [3,3-Trichloropropane] ND 0.0014 EPA 8260C 5-25-17 5-25-17 [3,3-Trichloropenzene] ND 0.0014 EPA 8260C 5-25-17 5-25-17 [3,2-Trichloropenzene] ND 0.0014 EPA 8260C 5-25-17 5-25-17 [3,2-Trichloropenzene] ND 0.0014 EPA 8260C 5-25-17 5-25-17 [3,2-Trichloropenzene] ND 0							
1,3-Dichloropropane							
2-Hexanone							
Dibromochloromethane	1,3-Dichloropropane		0.0014	EPA 8260C		5-25-17	
1,2-Dibromoethane	2-Hexanone	ND	0.0069	EPA 8260C	5-25-17	5-25-17	
Chlorobenzene	Dibromochloromethane	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
1,1,1,2-Tetrachloroethane	1,2-Dibromoethane	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
Ethylbenzene	Chlorobenzene	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
N.D. Xylene N.D. 0.0028 EPA 8260C 5-25-17	1,1,1,2-Tetrachloroethane	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
ND 0.0014 EPA 8260C 5-25-17	Ethylbenzene	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
Styrene ND 0.0014 EPA 8260C 5-25-17 5-25-17 Septimor ND 0.0069 EPA 8260C 5-25-17 5-25-17 Septimor ND 0.0014 EPA 8260C 5-25-17 Septimor Septimor ND 0.0014 EPA 8260C 5-25-17 Septimor Septimor ND 0.0014 EPA 8260C S-25-17 Septimor Septimor ND 0.0014 EPA 8260C S-25-17 Septimor Septimor ND 0.0014 EPA 8260C S-25-17 S-25-17 S-25-17 Septimor ND 0.0014 EPA 8260C S-25-17 S-25-17 S-25-17 Septimor ND 0.0014 EPA 8260C S-25-17	m,p-Xylene	ND	0.0028	EPA 8260C	5-25-17	5-25-17	
Second Form ND 0.0069 EPA 8260C 5-25-17 5-25	o-Xylene	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
Sepropy S	Styrene	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
Second Decomposition	Bromoform	ND	0.0069	EPA 8260C	5-25-17	5-25-17	
Second Denomber ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,1,2,2-Tetrachloroethane ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichloropropane ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichloropropane ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-Dibromo-3-chloropropane ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2,3-Trimethylbenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,3,5-Trimethylbenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2,4-Trimethylbenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-Dibromo-3-chloropropane ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-Dibromo-3-chloropropane ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-Dibromo-3-chloropropane ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-Dibromo-3-chloropropane ND 0.0014 EPA 8260C 5-25-17 5-25-17	Isopropylbenzene	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
1,2,3-Trichloropropane	Bromobenzene	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
n-Propylbenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,3-5-17 1,3-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,4-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-Dichlorobenzene ND 0.	1,1,2,2-Tetrachloroethane	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
n-Propylbenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,3-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,4-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-Dibromo-3-chloropropane ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-Dibromo-3-chloropropane ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-3-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-3-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-Dibromo-3-chloropropane ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-	1,2,3-Trichloropropane	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
C-Chlorotoluene	n-Propylbenzene	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
1,3,5-Trimethylbenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 tert-Butylbenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2,4-Trimethylbenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 sec-Butylbenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,3-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,4-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,4-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-Dibromo-3-chloropropane ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0069 EPA 8260C 5-25-17 5-25-17 Hexachlorobutadiene ND 0.0014 EPA 8260C 5-25-17 5-25-17 ND 0.0014 EPA 8260C	2-Chlorotoluene	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
Itert-Butylbenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2,4-Trimethylbenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 sec-Butylbenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,3-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,4-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,4-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-Dibromo-3-chloropropane ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-Dibromo-3-chloropropane ND 0.0069 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 Hexachlorobutadiene ND 0.0014 EPA 8260C 5-25-17 5-25-17 ND 0.0014 EPA 8260C	4-Chlorotoluene	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
Sect	1,3,5-Trimethylbenzene	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
1,2,4-Trimethylbenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 sec-Butylbenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,3-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,4-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 n-Butylbenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-Dibromo-3-chloropropane ND 0.0069 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 Hexachlorobutadiene ND 0.0069 EPA 8260C 5-25-17 5-25-17 Naphthalene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 121 73-134		ND	0.0014	EPA 8260C	5-25-17	5-25-17	
Seec-Butylbenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,3-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,3-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,4-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-Dibromo-3-chloropropane ND 0.0069 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 Naphthalene ND 0.0014 EPA 8260C		ND	0.0014	EPA 8260C	5-25-17		
1,3-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 p-Isopropyltoluene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,4-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 n-Butylbenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-Dibromo-3-chloropropane ND 0.0069 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 Hexachlorobutadiene ND 0.0069 EPA 8260C 5-25-17 5-25-17 Naphthalene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 121 73-134 Toluene-d8 114 81-124		ND					
Description		ND	0.0014		5-25-17	5-25-17	
1,4-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 n-Butylbenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-Dibromo-3-chloropropane ND 0.0069 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 Hexachlorobutadiene ND 0.0069 EPA 8260C 5-25-17 5-25-17 Naphthalene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 121 73-134 Toluene-d8 114 81-124		ND	0.0014		5-25-17		
1,2-Dichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 n-Butylbenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-Dibromo-3-chloropropane ND 0.0069 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 Hexachlorobutadiene ND 0.0069 EPA 8260C 5-25-17 5-25-17 Naphthalene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 121 73-134 Toluene-d8 114 81-124							
ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2-Dibromo-3-chloropropane ND 0.0069 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 5-25-17 Hexachlorobutadiene ND 0.0069 EPA 8260C 5-25-17 5-25-17 5-25-17 Naphthalene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 121 73-134 Toluene-d8 114 81-124 81-124 114 81-124 114 81-124 114 115 11							
1,2-Dibromo-3-chloropropane ND 0.0069 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 Hexachlorobutadiene ND 0.0069 EPA 8260C 5-25-17 5-25-17 Naphthalene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 121 73-134 Toluene-d8 114 81-124							
1,2,4-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 Hexachlorobutadiene ND 0.0069 EPA 8260C 5-25-17 5-25-17 Naphthalene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 121 73-134 Toluene-d8 114 81-124	-						
Hexachlorobutadiene ND 0.0069 EPA 8260C 5-25-17 5-25-17 Naphthalene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 121 73-134 Toluene-d8 114 81-124							
Naphthalene ND 0.0014 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 121 73-134 Toluene-d8 114 81-124							
1,2,3-Trichlorobenzene ND 0.0014 EPA 8260C 5-25-17 5-25-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 121 73-134 Toluene-d8 114 81-124							
Surrogate: Percent Recovery Control Limits Dibromofluoromethane 121 73-134 Toluene-d8 114 81-124							
Dibromofluoromethane 121 73-134 Toluene-d8 114 81-124					0 _0	0 =0	
Toluene-d8 114 81-124		=					
	4-Bromofluorobenzene	108	80-131				

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

Matrix: Soil Units: mg/kg

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP3-2.5					
Laboratory ID:	05-299-07					
Dichlorodifluoromethane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Chloromethane	ND	0.0060	EPA 8260C	5-25-17	5-25-17	
Vinyl Chloride	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Bromomethane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Chloroethane	ND	0.0060	EPA 8260C	5-25-17	5-25-17	
Trichlorofluoromethane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
1,1-Dichloroethene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Acetone	0.058	0.012	EPA 8260C	5-25-17	5-25-17	
Iodomethane	ND	0.0076	EPA 8260C	5-25-17	5-25-17	
Carbon Disulfide	ND	0.0017	EPA 8260C	5-25-17	5-25-17	
Methylene Chloride	ND	0.012	EPA 8260C	5-25-17	5-25-17	
(trans) 1,2-Dichloroethene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Methyl t-Butyl Ether	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
1,1-Dichloroethane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Vinyl Acetate	ND	0.0060	EPA 8260C	5-25-17	5-25-17	
2,2-Dichloropropane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
(cis) 1,2-Dichloroethene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
2-Butanone	0.013	0.0060	EPA 8260C	5-25-17	5-25-17	
Bromochloromethane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Chloroform	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
1,1,1-Trichloroethane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Carbon Tetrachloride	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
1,1-Dichloropropene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Benzene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
1,2-Dichloroethane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Trichloroethene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
1,2-Dichloropropane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Dibromomethane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Bromodichloromethane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
2-Chloroethyl Vinyl Ether	ND	0.0060	EPA 8260C	5-25-17	5-25-17	
(cis) 1,3-Dichloropropene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Methyl Isobutyl Ketone	ND	0.0060	EPA 8260C	5-25-17	5-25-17	
Toluene	ND	0.0060	EPA 8260C	5-25-17	5-25-17	
(trans) 1,3-Dichloropropene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	

Project: 15-05986-040

VOLATILES EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP3-2.5					
Laboratory ID:	05-299-07					
1,1,2-Trichloroethane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Tetrachloroethene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
1,3-Dichloropropane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
2-Hexanone	ND	0.0060	EPA 8260C	5-25-17	5-25-17	
Dibromochloromethane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
1,2-Dibromoethane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Chlorobenzene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
1,1,1,2-Tetrachloroethane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Ethylbenzene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
m,p-Xylene	ND	0.0024	EPA 8260C	5-25-17	5-25-17	
o-Xylene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Styrene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Bromoform	ND	0.0060	EPA 8260C	5-25-17	5-25-17	
Isopropylbenzene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Bromobenzene	ND	0.069	EPA 8260C	5-26-17	5-26-17	
1,1,2,2-Tetrachloroethane	ND	0.069	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichloropropane	ND	0.069	EPA 8260C	5-26-17	5-26-17	
n-Propylbenzene	ND	0.069	EPA 8260C	5-26-17	5-26-17	
2-Chlorotoluene	ND	0.069	EPA 8260C	5-26-17	5-26-17	
4-Chlorotoluene	ND	0.069	EPA 8260C	5-26-17	5-26-17	
1,3,5-Trimethylbenzene	ND	0.069	EPA 8260C	5-26-17	5-26-17	
tert-Butylbenzene	ND	0.069	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trimethylbenzene	ND	0.069	EPA 8260C	5-26-17	5-26-17	
sec-Butylbenzene	ND	0.069	EPA 8260C	5-26-17	5-26-17	
1,3-Dichlorobenzene	ND	0.069	EPA 8260C	5-26-17	5-26-17	
p-Isopropyltoluene	ND	0.069	EPA 8260C	5-26-17	5-26-17	
1,4-Dichlorobenzene	ND	0.069	EPA 8260C	5-26-17	5-26-17	
1,2-Dichlorobenzene	ND	0.069	EPA 8260C	5-26-17	5-26-17	
n-Butylbenzene	ND	0.069	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromo-3-chloropropane	ND	0.34	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trichlorobenzene	ND	0.069	EPA 8260C	5-26-17	5-26-17	
Hexachlorobutadiene	ND	0.34	EPA 8260C	5-26-17	5-26-17	
Naphthalene	ND	0.069	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichlorobenzene	ND	0.069	EPA 8260C	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	105	73-134				
Toluene-d8	96	81-124				

 Dibromofluoromethane
 105
 73-134

 Toluene-d8
 96
 81-124

 4-Bromofluorobenzene
 80
 80-131



Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP3-10					
Laboratory ID:	05-299-08					
Dichlorodifluoromethane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Chloromethane	ND	0.0061	EPA 8260C	5-25-17	5-25-17	
Vinyl Chloride	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Bromomethane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Chloroethane	ND	0.0061	EPA 8260C	5-25-17	5-25-17	
Trichlorofluoromethane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
1,1-Dichloroethene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Acetone	0.028	0.012	EPA 8260C	5-25-17	5-25-17	
lodomethane	ND	0.0077	EPA 8260C	5-25-17	5-25-17	
Carbon Disulfide	ND	0.0017	EPA 8260C	5-25-17	5-25-17	
Methylene Chloride	ND	0.012	EPA 8260C	5-25-17	5-25-17	
(trans) 1,2-Dichloroethene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Methyl t-Butyl Ether	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
1,1-Dichloroethane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Vinyl Acetate	ND	0.0061	EPA 8260C	5-25-17	5-25-17	
2,2-Dichloropropane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
(cis) 1,2-Dichloroethene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
2-Butanone	ND	0.0061	EPA 8260C	5-25-17	5-25-17	
Bromochloromethane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Chloroform	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
1,1,1-Trichloroethane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Carbon Tetrachloride	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
1,1-Dichloropropene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Benzene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
1,2-Dichloroethane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Trichloroethene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
1,2-Dichloropropane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Dibromomethane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Bromodichloromethane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
2-Chloroethyl Vinyl Ether	ND	0.0061	EPA 8260C	5-25-17	5-25-17	
(cis) 1,3-Dichloropropene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Methyl Isobutyl Ketone	ND	0.0061	EPA 8260C	5-25-17	5-25-17	
Toluene	0.016	0.0061	EPA 8260C	5-25-17	5-25-17	
(trans) 1,3-Dichloropropene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	

Project: 15-05986-040

VOLATILES EPA 8260C

Client ID:					Date	Date	
Laboratory ID:	Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
1,1,2-Trichloroethane							
Tetrachloroethene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,3-Dichloropropane ND 0.0012 EPA 8260C 5-25-17		05-299-08					
1,3-Dichloropropane ND 0.0012 EPA 8260C 5-25-17 5-25-17	1,1,2-Trichloroethane		0.0012	EPA 8260C	5-25-17	5-25-17	
Part	Tetrachloroethene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Dibromochloromethane	1,3-Dichloropropane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
1,2-Dibromoethane ND 0.0012 EPA 8260C 5-25-17	2-Hexanone	ND	0.0061	EPA 8260C	5-25-17	5-25-17	
Chlorobenzene	Dibromochloromethane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
1,1,1,2-Tetrachloroethane	1,2-Dibromoethane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Ethylbenzene	Chlorobenzene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
m,p-Xylene ND 0.0025 EPA 8260C 5-25-17 5-25-17 Styrene ND 0.0012 EPA 8260C 5-25-17 5-25-17 Styrene ND 0.0012 EPA 8260C 5-25-17 5-25-17 Styrene ND 0.0012 EPA 8260C 5-25-17 5-25-17 Styrene ND 0.0061 EPA 8260C 5-25-17 5-25-17 Styrene ND 0.0061 EPA 8260C 5-25-17 5-25-17 Styrene ND 0.0012 EPA 8260C 5-25-17 5-25-17 Styrene ND 0.0012 EPA 8260C 5-25-17 5-25-17 Styrene ND 0.0012 EPA 8260C 5-25-17 5-25-17 Styrene ND 0.0012 EPA 8260C 5-25-17 5-25-17 Styrene ND 0.0012 EPA 8260C 5-25-17 5-25-1	1,1,1,2-Tetrachloroethane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
ND ND ND ND ND ND ND ND	Ethylbenzene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Styrene ND 0.0012 EPA 8260C 5-25-17 5-25-17 Styrene ND 0.0061 EPA 8260C 5-25-17 5-25-17 Styrene ND 0.0061 EPA 8260C 5-25-17 5-25-17 Styrene ND 0.0012 EPA 8260C 5-25-17 5-25-17 Styrene Styrene Styrene ND 0.0012 EPA 8260C 5-25-17 5-25-17 Styrene Styr	n,p-Xylene	ND	0.0025	EPA 8260C	5-25-17	5-25-17	
Second Composition ND 0.0061 EPA 8260C 5-25-17	o-Xylene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Sopropy benzene ND 0.0012 EPA 8260C 5-25-17	Styrene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Sammobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,1,2,2-Tetrachloroethane ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichloropropane ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichloropropane ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobluene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2,5-Trimethylbenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,3,5-Trimethylbenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2,4-Trimethylbenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2,4-Trimethylbenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,3-Dichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,3-Dichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,3-Dichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,4-Dichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,4-Dichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2-Dichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17	Bromoform	ND	0.0061	EPA 8260C	5-25-17	5-25-17	
1,1,2,2-Tetrachloroethane	sopropylbenzene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
1,2,3-Trichloropropane	Bromobenzene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
A-Propylbenzene	I,1,2,2-Tetrachloroethane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
ND	I,2,3-Trichloropropane	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
A-Chlorotoluene ND 0.0012 EPA 8260C 5-25-17 5-25-17 [1,3,5-Trimethylbenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 [1,2,4-Trimethylbenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 [1,2,4-Trimethylbenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 [1,2,4-Trimethylbenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 [1,3-Dichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 [1,3-Dichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 [1,4-Dichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 [1,4-Dichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 [1,2-Dichlorobenzene ND 0.0061 EPA 8260C 5-25-17 5-25-17 [1,2,4-Trichlorobenzene ND 0.0061 EPA 8260C 5-25-17 5-25-17 [1,2,4-Trichlorobenzene ND 0.0061 EPA 8260C 5-25-17 5-25-17 [1,2,3-Trichlorobenzene ND 0.0012 EPA 826	n-Propylbenzene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
1,3,5-Trimethylbenzene	2-Chlorotoluene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
ert-Butylbenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2,4-Trimethylbenzene ND 0.0012 EPA 8260C 5-25-17 5-25-1	1-Chlorotoluene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
1,2,4-Trimethylbenzene	1,3,5-Trimethylbenzene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
1,2,4-Trimethylbenzene		ND	0.0012	EPA 8260C	5-25-17	5-25-17	
Sec-Butylbenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,3-Dichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,4-Dichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,4-Dichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2-Dichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2-Dichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2-Dibromo-3-chloropropane ND 0.0061 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 Naphthalene ND 0.0012 EPA 8260C 5-25-17 5-25-17 ND 0.0012		ND	0.0012	EPA 8260C	5-25-17	5-25-17	
I,3-Dichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 p-Isopropyltoluene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,4-Dichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2-Dichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 n-Butylbenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2-Dibromo-3-chloropropane ND 0.0061 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 Nexachlorobutadiene ND 0.0061 EPA 8260C 5-25-17 5-25-17 Naphthalene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 104 73-134 Toluene-d8 106		ND	0.0012	EPA 8260C	5-25-17	5-25-17	
1,4-Dichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2-Dichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 n-Butylbenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2-Dibromo-3-chloropropane ND 0.0061 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 Hexachlorobutadiene ND 0.0061 EPA 8260C 5-25-17 5-25-17 Naphthalene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 104 73-134 Toluene-d8 106 81-124		ND	0.0012	EPA 8260C	5-25-17	5-25-17	
1,4-Dichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2-Dichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 n-Butylbenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2-Dibromo-3-chloropropane ND 0.0061 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 Hexachlorobutadiene ND 0.0061 EPA 8260C 5-25-17 5-25-17 Naphthalene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 104 73-134 Toluene-d8 106 81-124	o-Isopropyltoluene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2-Dibromo-3-chloropropane ND 0.0061 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobutadiene ND 0.0061 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2		ND	0.0012	EPA 8260C	5-25-17	5-25-17	
1,2-Dibromo-3-chloropropane ND 0.0061 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 Hexachlorobutadiene ND 0.0061 EPA 8260C 5-25-17 5-25-17 Naphthalene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 104 73-134 Toluene-d8 106 81-124	1,2-Dichlorobenzene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
1,2-Dibromo-3-chloropropane ND 0.0061 EPA 8260C 5-25-17 5-25-17 1,2,4-Trichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 Hexachlorobutadiene ND 0.0061 EPA 8260C 5-25-17 5-25-17 Naphthalene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 104 73-134 Toluene-d8 106 81-124	n-Butylbenzene	ND	0.0012	EPA 8260C	5-25-17	5-25-17	
1,2,4-Trichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 Hexachlorobutadiene ND 0.0061 EPA 8260C 5-25-17 5-25-17 Naphthalene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 104 73-134 Toluene-d8 106 81-124		ND	0.0061	EPA 8260C	5-25-17	5-25-17	
Hexachlorobutadiene ND 0.0061 EPA 8260C 5-25-17 5-25-17 Naphthalene ND 0.0012 EPA 8260C 5-25-17 5-25-17 1,2,3-Trichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 104 73-134 Toluene-d8 106 81-124			0.0012	EPA 8260C	5-25-17	5-25-17	
Naphthalene ND 0.0012 EPA 8260C 5-25-17 5-25-17 I,2,3-Trichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 104 73-134 Toluene-d8 106 81-124	·						
I,2,3-Trichlorobenzene ND 0.0012 EPA 8260C 5-25-17 5-25-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 104 73-134 Toluene-d8 106 81-124							
Surrogate: Percent Recovery Control Limits Dibromofluoromethane 104 73-134 Toluene-d8 106 81-124	•						
Dibromofluoromethane 104 73-134 Toluene-d8 106 81-124					-	•	
Toluene-d8 106 81-124		-					
		100	80-131				

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP4-3					
Laboratory ID:	05-299-10					
Dichlorodifluoromethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Chloromethane	ND	0.0057	EPA 8260C	5-25-17	5-25-17	
Vinyl Chloride	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Bromomethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Chloroethane	ND	0.0057	EPA 8260C	5-25-17	5-25-17	
Trichlorofluoromethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,1-Dichloroethene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Acetone	ND	0.011	EPA 8260C	5-25-17	5-25-17	
lodomethane	ND	0.0071	EPA 8260C	5-25-17	5-25-17	
Carbon Disulfide	ND	0.0016	EPA 8260C	5-25-17	5-25-17	
Methylene Chloride	ND	0.011	EPA 8260C	5-25-17	5-25-17	
(trans) 1,2-Dichloroethene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Methyl t-Butyl Ether	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,1-Dichloroethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Vinyl Acetate	ND	0.0057	EPA 8260C	5-25-17	5-25-17	
2,2-Dichloropropane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
(cis) 1,2-Dichloroethene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
2-Butanone	ND	0.0057	EPA 8260C	5-25-17	5-25-17	
Bromochloromethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Chloroform	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,1,1-Trichloroethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Carbon Tetrachloride	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,1-Dichloropropene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Benzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,2-Dichloroethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Trichloroethene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,2-Dichloropropane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Dibromomethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Bromodichloromethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
2-Chloroethyl Vinyl Ether	ND	0.0057	EPA 8260C	5-25-17	5-25-17	
(cis) 1,3-Dichloropropene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Methyl Isobutyl Ketone	ND	0.0057	EPA 8260C	5-25-17	5-25-17	
Toluene	0.016	0.0057	EPA 8260C	5-25-17	5-25-17	
(trans) 1,3-Dichloropropene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	

Project: 15-05986-040

VOLATILES EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP4-3					
Laboratory ID:	05-299-10					
1,1,2-Trichloroethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Tetrachloroethene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,3-Dichloropropane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
2-Hexanone	ND	0.0057	EPA 8260C	5-25-17	5-25-17	
Dibromochloromethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,2-Dibromoethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Chlorobenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,1,1,2-Tetrachloroethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Ethylbenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
m,p-Xylene	ND	0.0023	EPA 8260C	5-25-17	5-25-17	
o-Xylene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Styrene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Bromoform	ND	0.0057	EPA 8260C	5-25-17	5-25-17	
Isopropylbenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Bromobenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,1,2,2-Tetrachloroethane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,2,3-Trichloropropane	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
n-Propylbenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
2-Chlorotoluene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
4-Chlorotoluene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,3,5-Trimethylbenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
tert-Butylbenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,2,4-Trimethylbenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
sec-Butylbenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,3-Dichlorobenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
p-Isopropyltoluene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,4-Dichlorobenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,2-Dichlorobenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
n-Butylbenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,2-Dibromo-3-chloropropane	ND	0.0057	EPA 8260C	5-25-17	5-25-17	
1,2,4-Trichlorobenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Hexachlorobutadiene	ND	0.0057	EPA 8260C	5-25-17	5-25-17	
Naphthalene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
1,2,3-Trichlorobenzene	ND	0.0011	EPA 8260C	5-25-17	5-25-17	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	107	73-134				
Toluene-d8	105	81-124				

4-Bromofluorobenzene

80-131

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP4-10					
Laboratory ID:	05-299-11					
Dichlorodifluoromethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Chloromethane	ND	0.0064	EPA 8260C	5-26-17	5-26-17	
Vinyl Chloride	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Bromomethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Chloroethane	ND	0.0064	EPA 8260C	5-26-17	5-26-17	
Trichlorofluoromethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Acetone	0.051	0.013	EPA 8260C	5-26-17	5-26-17	
lodomethane	ND	0.0064	EPA 8260C	5-26-17	5-26-17	
Carbon Disulfide	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Methylene Chloride	ND	0.013	EPA 8260C	5-26-17	5-26-17	
(trans) 1,2-Dichloroethene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Methyl t-Butyl Ether	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Vinyl Acetate	ND	0.0064	EPA 8260C	5-26-17	5-26-17	
2,2-Dichloropropane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
(cis) 1,2-Dichloroethene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
2-Butanone	0.0068	0.0064	EPA 8260C	5-26-17	5-26-17	
Bromochloromethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Chloroform	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1,1-Trichloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Carbon Tetrachloride	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloropropene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Benzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Trichloroethene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloropropane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Dibromomethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Bromodichloromethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
2-Chloroethyl Vinyl Ether	ND	0.0083	EPA 8260C	5-26-17	5-26-17	
(cis) 1,3-Dichloropropene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Methyl Isobutyl Ketone	ND	0.0064	EPA 8260C	5-26-17	5-26-17	
Toluene	0.011	0.0064	EPA 8260C	5-26-17	5-26-17	
(trans) 1,3-Dichloropropene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	

Project: 15-05986-040

VOLATILES EPA 8260C

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP4-10					
Laboratory ID:	05-299-11					
1,1,2-Trichloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Tetrachloroethene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,3-Dichloropropane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
2-Hexanone	ND	0.0064	EPA 8260C	5-26-17	5-26-17	
Dibromochloromethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromoethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Chlorobenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1,1,2-Tetrachloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Ethylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
m,p-Xylene	ND	0.0026	EPA 8260C	5-26-17	5-26-17	
o-Xylene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Styrene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Bromoform	ND	0.0064	EPA 8260C	5-26-17	5-26-17	
Isopropylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Bromobenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1,2,2-Tetrachloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichloropropane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
n-Propylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
2-Chlorotoluene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
4-Chlorotoluene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,3,5-Trimethylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
tert-Butylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trimethylbenzene	0.0024	0.0013	EPA 8260C	5-26-17	5-26-17	
sec-Butylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,3-Dichlorobenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
o-Isopropyltoluene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,4-Dichlorobenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2-Dichlorobenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
n-Butylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromo-3-chloropropane		0.0064	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trichlorobenzene	ND	0.0013	EPA 8260C	5-26-17 5-26-17	5-26-17	
Hexachlorobutadiene	ND	0.0064	EPA 8260C	5-26-17 5-26-17	5-26-17 5-26-17	
Naphthalene	0.0019	0.0004	EPA 8260C	5-26-17 5-26-17	5-26-17 5-26-17	
1,2,3-Trichlorobenzene	0.0019 ND	0.0013	EPA 8260C	5-26-17 5-26-17	5-26-17 5-26-17	
			EFA 02000	J-ZU-11	J-ZU-1 <i>1</i>	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	102	73-134				
Toluene-d8	99	81-124				
4-Bromofluorobenzene	94	80-131				

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP5-2.5					
Laboratory ID:	05-299-13					
Dichlorodifluoromethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Chloromethane	ND	0.0058	EPA 8260C	5-26-17	5-26-17	
Vinyl Chloride	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Bromomethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Chloroethane	ND	0.0058	EPA 8260C	5-26-17	5-26-17	
Trichlorofluoromethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Acetone	ND	0.012	EPA 8260C	5-26-17	5-26-17	
lodomethane	ND	0.0058	EPA 8260C	5-26-17	5-26-17	
Carbon Disulfide	ND	0.0058	EPA 8260C	5-26-17	5-26-17	
Methylene Chloride	ND	0.012	EPA 8260C	5-26-17	5-26-17	
(trans) 1,2-Dichloroethene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Methyl t-Butyl Ether	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Vinyl Acetate	ND	0.0058	EPA 8260C	5-26-17	5-26-17	
2,2-Dichloropropane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
(cis) 1,2-Dichloroethene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
2-Butanone	ND	0.0058	EPA 8260C	5-26-17	5-26-17	
Bromochloromethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Chloroform	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,1,1-Trichloroethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Carbon Tetrachloride	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloropropene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Benzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloroethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Trichloroethene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloropropane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Dibromomethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Bromodichloromethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
2-Chloroethyl Vinyl Ether	ND	0.0074	EPA 8260C	5-26-17	5-26-17	
(cis) 1,3-Dichloropropene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Methyl Isobutyl Ketone	ND	0.0058	EPA 8260C	5-26-17	5-26-17	
Toluene	ND	0.0058	EPA 8260C	5-26-17	5-26-17	
(trans) 1,3-Dichloropropene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	

Project: 15-05986-040

VOLATILES EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP5-2.5					
Laboratory ID:	05-299-13					
1,1,2-Trichloroethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Tetrachloroethene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,3-Dichloropropane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
2-Hexanone	ND	0.0058	EPA 8260C	5-26-17	5-26-17	
Dibromochloromethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromoethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Chlorobenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,1,1,2-Tetrachloroethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Ethylbenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
m,p-Xylene	ND	0.0023	EPA 8260C	5-26-17	5-26-17	
o-Xylene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Styrene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Bromoform	ND	0.0058	EPA 8260C	5-26-17	5-26-17	
Isopropylbenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Bromobenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,1,2,2-Tetrachloroethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichloropropane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
n-Propylbenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
2-Chlorotoluene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
4-Chlorotoluene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,3,5-Trimethylbenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
tert-Butylbenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trimethylbenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
sec-Butylbenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,3-Dichlorobenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
p-Isopropyltoluene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,4-Dichlorobenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,2-Dichlorobenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
n-Butylbenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromo-3-chloropropane	ND	0.0058	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trichlorobenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Hexachlorobutadiene	ND	0.0058	EPA 8260C	5-26-17	5-26-17	
Naphthalene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichlorobenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	105	73-134				
Toluene-d8	101	81-124				

4-Bromofluorobenzene

80-131

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP5-10					
Laboratory ID:	05-299-14					
Dichlorodifluoromethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Chloromethane	ND	0.0065	EPA 8260C	5-26-17	5-26-17	
Vinyl Chloride	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Bromomethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Chloroethane	ND	0.0065	EPA 8260C	5-26-17	5-26-17	
Trichlorofluoromethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Acetone	0.046	0.013	EPA 8260C	5-26-17	5-26-17	
lodomethane	ND	0.0065	EPA 8260C	5-26-17	5-26-17	
Carbon Disulfide	0.0024	0.0013	EPA 8260C	5-26-17	5-26-17	
Methylene Chloride	ND	0.013	EPA 8260C	5-26-17	5-26-17	
(trans) 1,2-Dichloroethene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Methyl t-Butyl Ether	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Vinyl Acetate	ND	0.0065	EPA 8260C	5-26-17	5-26-17	
2,2-Dichloropropane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
(cis) 1,2-Dichloroethene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
2-Butanone	0.013	0.0065	EPA 8260C	5-26-17	5-26-17	
Bromochloromethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Chloroform	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1,1-Trichloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Carbon Tetrachloride	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloropropene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Benzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Trichloroethene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloropropane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Dibromomethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Bromodichloromethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
2-Chloroethyl Vinyl Ether	ND	0.0083	EPA 8260C	5-26-17	5-26-17	
(cis) 1,3-Dichloropropene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Methyl Isobutyl Ketone	ND	0.0065	EPA 8260C	5-26-17	5-26-17	
Toluene	0.016	0.0065	EPA 8260C	5-26-17	5-26-17	
(trans) 1,3-Dichloropropene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	

Project: 15-05986-040

VOLATILES EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP5-10					
Laboratory ID:	05-299-14					
1,1,2-Trichloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Tetrachloroethene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,3-Dichloropropane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
2-Hexanone	ND	0.0065	EPA 8260C	5-26-17	5-26-17	
Dibromochloromethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromoethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Chlorobenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1,1,2-Tetrachloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Ethylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
m,p-Xylene	ND	0.0026	EPA 8260C	5-26-17	5-26-17	
o-Xylene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Styrene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Bromoform	ND	0.0065	EPA 8260C	5-26-17	5-26-17	
Isopropylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Bromobenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1,2,2-Tetrachloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichloropropane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
n-Propylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
2-Chlorotoluene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
4-Chlorotoluene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,3,5-Trimethylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
tert-Butylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trimethylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
sec-Butylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,3-Dichlorobenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
p-Isopropyltoluene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,4-Dichlorobenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2-Dichlorobenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
n-Butylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromo-3-chloropropane		0.0065	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trichlorobenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Hexachlorobutadiene	ND	0.0065	EPA 8260C	5-26-17	5-26-17	
Naphthalene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichlorobenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits		-	-	
Dibromofluoromethane	102	73-134				
Toluene-d8	101	81-124				
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4-Bromofluorobenzene

80-131

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP6-2					
Laboratory ID:	05-299-16					
Dichlorodifluoromethane	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
Chloromethane	ND	0.0045	EPA 8260C	5-26-17	5-26-17	
Vinyl Chloride	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
Bromomethane	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
Chloroethane	ND	0.0045	EPA 8260C	5-26-17	5-26-17	
Trichlorofluoromethane	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
Acetone	0.011	0.0090	EPA 8260C	5-26-17	5-26-17	
lodomethane	ND	0.0045	EPA 8260C	5-26-17	5-26-17	
Carbon Disulfide	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
Methylene Chloride	ND	0.0090	EPA 8260C	5-26-17	5-26-17	
(trans) 1,2-Dichloroethene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
Methyl t-Butyl Ether	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethane	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
Vinyl Acetate	ND	0.0045	EPA 8260C	5-26-17	5-26-17	
2,2-Dichloropropane	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
(cis) 1,2-Dichloroethene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
2-Butanone	ND	0.0045	EPA 8260C	5-26-17	5-26-17	
Bromochloromethane	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
Chloroform	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
1,1,1-Trichloroethane	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
Carbon Tetrachloride	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloropropene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
Benzene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloroethane	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
Trichloroethene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloropropane	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
Dibromomethane	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
Bromodichloromethane	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
2-Chloroethyl Vinyl Ether	ND	0.0058	EPA 8260C	5-26-17	5-26-17	
(cis) 1,3-Dichloropropene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
Methyl Isobutyl Ketone	ND	0.0045	EPA 8260C	5-26-17	5-26-17	
Toluene	0.0096	0.0045	EPA 8260C	5-26-17	5-26-17	
(trans) 1,3-Dichloropropene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	

Project: 15-05986-040

VOLATILES EPA 8260C

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP6-2					
_aboratory ID:	05-299-16					
1,1,2-Trichloroethane	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
Γetrachloroethene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
1,3-Dichloropropane	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
2-Hexanone	ND	0.0045	EPA 8260C	5-26-17	5-26-17	
Dibromochloromethane	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromoethane	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
Chlorobenzene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
1,1,1,2-Tetrachloroethane	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
Ethylbenzene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
n,p-Xylene	ND	0.0018	EPA 8260C	5-26-17	5-26-17	
o-Xylene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
Styrene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
Bromoform	ND	0.0045	EPA 8260C	5-26-17	5-26-17	
sopropylbenzene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
Bromobenzene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
I,1,2,2-Tetrachloroethane	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichloropropane	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
n-Propylbenzene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
2-Chlorotoluene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
1-Chlorotoluene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
1,3,5-Trimethylbenzene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
ert-Butylbenzene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trimethylbenzene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
sec-Butylbenzene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
1,3-Dichlorobenzene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
o-Isopropyltoluene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
1,4-Dichlorobenzene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
1,2-Dichlorobenzene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
n-Butylbenzene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromo-3-chloropropane	ND	0.0045	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trichlorobenzene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
	ND	0.0045	EPA 8260C	5-26-17	5-26-17	
Naphthalene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichlorobenzene	ND	0.00090	EPA 8260C	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	108	73-134				
Toluene-d8						
	101	81-124				

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

• • •		-01		Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP6-10					
Laboratory ID:	05-299-17	0.0045		5.00.47	5.00.17	
Dichlorodifluoromethane	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
Chloromethane	ND	0.0076	EPA 8260C	5-26-17	5-26-17	
Vinyl Chloride	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
Bromomethane	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
Chloroethane	ND	0.0076	EPA 8260C	5-26-17	5-26-17	
Trichlorofluoromethane	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
Acetone	0.10	0.015	EPA 8260C	5-26-17	5-26-17	
lodomethane	ND	0.0076	EPA 8260C	5-26-17	5-26-17	
Carbon Disulfide	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
Methylene Chloride	ND	0.015	EPA 8260C	5-26-17	5-26-17	
(trans) 1,2-Dichloroethene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
Methyl t-Butyl Ether	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethane	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
Vinyl Acetate	ND	0.0076	EPA 8260C	5-26-17	5-26-17	
2,2-Dichloropropane	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
(cis) 1,2-Dichloroethene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
2-Butanone	0.025	0.0076	EPA 8260C	5-26-17	5-26-17	
Bromochloromethane	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
Chloroform	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
1,1,1-Trichloroethane	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
Carbon Tetrachloride	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloropropene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
Benzene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloroethane	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
Trichloroethene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloropropane	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
Dibromomethane	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
Bromodichloromethane	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
2-Chloroethyl Vinyl Ether	ND	0.0097	EPA 8260C	5-26-17	5-26-17	
(cis) 1,3-Dichloropropene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
Methyl Isobutyl Ketone	ND	0.0076	EPA 8260C	5-26-17	5-26-17	
Toluene	0.0083	0.0076	EPA 8260C	5-26-17	5-26-17	
(trans) 1,3-Dichloropropene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	

Project: 15-05986-040

VOLATILES EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP6-10					
Laboratory ID:	05-299-17					
1,1,2-Trichloroethane	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
Tetrachloroethene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
1,3-Dichloropropane	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
2-Hexanone	ND	0.0076	EPA 8260C	5-26-17	5-26-17	
Dibromochloromethane	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromoethane	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
Chlorobenzene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
1,1,1,2-Tetrachloroethane	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
Ethylbenzene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
m,p-Xylene	ND	0.0030	EPA 8260C	5-26-17	5-26-17	
o-Xylene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
Styrene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
Bromoform	ND	0.0076	EPA 8260C	5-26-17	5-26-17	
Isopropylbenzene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
Bromobenzene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
1,1,2,2-Tetrachloroethane	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichloropropane	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
n-Propylbenzene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
2-Chlorotoluene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
4-Chlorotoluene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
1,3,5-Trimethylbenzene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
tert-Butylbenzene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trimethylbenzene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
sec-Butylbenzene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
1,3-Dichlorobenzene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
p-Isopropyltoluene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
1,4-Dichlorobenzene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
1,2-Dichlorobenzene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
n-Butylbenzene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromo-3-chloropropane	ND	0.0076	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trichlorobenzene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
Hexachlorobutadiene	ND	0.0076	EPA 8260C	5-26-17	5-26-17	
Naphthalene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichlorobenzene	ND	0.0015	EPA 8260C	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	105	73-134				
Toluene-d8	111	81-124				

4-Bromofluorobenzene

80-131

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP7-0					
Laboratory ID:	05-299-19					
Dichlorodifluoromethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Chloromethane	ND	0.0071	EPA 8260C	5-26-17	5-26-17	
Vinyl Chloride	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Bromomethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Chloroethane	ND	0.0071	EPA 8260C	5-26-17	5-26-17	
Trichlorofluoromethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Acetone	ND	0.014	EPA 8260C	5-26-17	5-26-17	
lodomethane	ND	0.0071	EPA 8260C	5-26-17	5-26-17	
Carbon Disulfide	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Methylene Chloride	ND	0.014	EPA 8260C	5-26-17	5-26-17	
(trans) 1,2-Dichloroethene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Methyl t-Butyl Ether	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Vinyl Acetate	ND	0.0071	EPA 8260C	5-26-17	5-26-17	
2,2-Dichloropropane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
(cis) 1,2-Dichloroethene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
2-Butanone	ND	0.0071	EPA 8260C	5-26-17	5-26-17	
Bromochloromethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Chloroform	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,1,1-Trichloroethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Carbon Tetrachloride	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloropropene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Benzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloroethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Trichloroethene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloropropane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Dibromomethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Bromodichloromethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
2-Chloroethyl Vinyl Ether	ND	0.0091	EPA 8260C	5-26-17	5-26-17	
(cis) 1,3-Dichloropropene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Methyl Isobutyl Ketone	ND	0.0071	EPA 8260C	5-26-17	5-26-17	
Toluene	ND	0.0071	EPA 8260C	5-26-17	5-26-17	
(trans) 1,3-Dichloropropene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	

Project: 15-05986-040

VOLATILES EPA 8260C

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP7-0					
Laboratory ID:	05-299-19					
1,1,2-Trichloroethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Tetrachloroethene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,3-Dichloropropane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
2-Hexanone	ND	0.0071	EPA 8260C	5-26-17	5-26-17	
Dibromochloromethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromoethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Chlorobenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,1,1,2-Tetrachloroethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Ethylbenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
m,p-Xylene	ND	0.0028	EPA 8260C	5-26-17	5-26-17	
o-Xylene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Styrene	0.012	0.0014	EPA 8260C	5-26-17	5-26-17	
Bromoform	ND	0.0071	EPA 8260C	5-26-17	5-26-17	
Isopropylbenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Bromobenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,1,2,2-Tetrachloroethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichloropropane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
n-Propylbenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
2-Chlorotoluene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
4-Chlorotoluene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,3,5-Trimethylbenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
tert-Butylbenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trimethylbenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
sec-Butylbenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,3-Dichlorobenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
p-Isopropyltoluene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,4-Dichlorobenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,2-Dichlorobenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
n-Butylbenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromo-3-chloropropane		0.0071	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trichlorobenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Hexachlorobutadiene	ND	0.0074	EPA 8260C	5-26-17 5-26-17	5-26-17	
Naphthalene	ND	0.0014	EPA 8260C	5-26-17 5-26-17	5-26-17	
1,2,3-Trichlorobenzene	ND ND	0.0014	EPA 8260C	5-26-17 5-26-17	5-26-17 5-26-17	
Surrogate:	Percent Recovery	Control Limits	Li A 02000	J-2U-11	0-20-11	
Surrogate. Dibromofluoromethane	103	73-134				
Toluene-d8	100	81-124				
4-Bromofluorobenzene	90	80-131				

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP7-10					
Laboratory ID:	05-299-20					
Dichlorodifluoromethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Chloromethane	ND	0.0065	EPA 8260C	5-26-17	5-26-17	
Vinyl Chloride	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Bromomethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Chloroethane	ND	0.0065	EPA 8260C	5-26-17	5-26-17	
Trichlorofluoromethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Acetone	0.019	0.013	EPA 8260C	5-26-17	5-26-17	
lodomethane	ND	0.0065	EPA 8260C	5-26-17	5-26-17	
Carbon Disulfide	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Methylene Chloride	ND	0.013	EPA 8260C	5-26-17	5-26-17	
(trans) 1,2-Dichloroethene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Methyl t-Butyl Ether	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Vinyl Acetate	ND	0.0065	EPA 8260C	5-26-17	5-26-17	
2,2-Dichloropropane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
(cis) 1,2-Dichloroethene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
2-Butanone	ND	0.0065	EPA 8260C	5-26-17	5-26-17	
Bromochloromethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Chloroform	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1,1-Trichloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Carbon Tetrachloride	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloropropene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Benzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Trichloroethene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloropropane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Dibromomethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Bromodichloromethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
2-Chloroethyl Vinyl Ether	ND	0.0083	EPA 8260C	5-26-17	5-26-17	
(cis) 1,3-Dichloropropene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Methyl Isobutyl Ketone	ND	0.0065	EPA 8260C	5-26-17	5-26-17	
Toluene	0.022	0.0065	EPA 8260C	5-26-17	5-26-17	
(trans) 1,3-Dichloropropene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	

Project: 15-05986-040

VOLATILES EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP7-10					
Laboratory ID:	05-299-20					
1,1,2-Trichloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Tetrachloroethene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,3-Dichloropropane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
2-Hexanone	ND	0.0065	EPA 8260C	5-26-17	5-26-17	
Dibromochloromethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromoethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Chlorobenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1,1,2-Tetrachloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Ethylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
m,p-Xylene	ND	0.0026	EPA 8260C	5-26-17	5-26-17	
o-Xylene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Styrene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Bromoform	ND	0.0065	EPA 8260C	5-26-17	5-26-17	
Isopropylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Bromobenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1,2,2-Tetrachloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichloropropane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
n-Propylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
2-Chlorotoluene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
4-Chlorotoluene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,3,5-Trimethylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
tert-Butylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trimethylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
sec-Butylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,3-Dichlorobenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
p-Isopropyltoluene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,4-Dichlorobenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2-Dichlorobenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
n-Butylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromo-3-chloropropane	ND	0.0065	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trichlorobenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Hexachlorobutadiene	ND	0.0065	EPA 8260C	5-26-17	5-26-17	
Naphthalene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichlorobenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	105	73-134				
Toluene-d8	103	81-124				

4-Bromofluorobenzene

80-131

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

• • •		-01		Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP8-0.5					
Laboratory ID:	05-299-22	0.0040		5.00.47	5.00.17	
Dichlorodifluoromethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Chloromethane	ND	0.0065	EPA 8260C	5-26-17	5-26-17	
Vinyl Chloride	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Bromomethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Chloroethane	ND	0.0065	EPA 8260C	5-26-17	5-26-17	
Trichlorofluoromethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Acetone	0.027	0.013	EPA 8260C	5-26-17	5-26-17	
lodomethane	ND	0.0065	EPA 8260C	5-26-17	5-26-17	
Carbon Disulfide	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Methylene Chloride	ND	0.013	EPA 8260C	5-26-17	5-26-17	
(trans) 1,2-Dichloroethene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Methyl t-Butyl Ether	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Vinyl Acetate	ND	0.0065	EPA 8260C	5-26-17	5-26-17	
2,2-Dichloropropane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
(cis) 1,2-Dichloroethene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
2-Butanone	ND	0.0065	EPA 8260C	5-26-17	5-26-17	
Bromochloromethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Chloroform	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1,1-Trichloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Carbon Tetrachloride	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloropropene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Benzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Trichloroethene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloropropane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Dibromomethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Bromodichloromethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
2-Chloroethyl Vinyl Ether	ND	0.0083	EPA 8260C	5-26-17	5-26-17	
(cis) 1,3-Dichloropropene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Methyl Isobutyl Ketone	ND	0.0065	EPA 8260C	5-26-17	5-26-17	
Toluene	0.014	0.0065	EPA 8260C	5-26-17	5-26-17	
(trans) 1,3-Dichloropropene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	

Project: 15-05986-040

VOLATILES EPA 8260C

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP8-0.5					
Laboratory ID:	05-299-22					
1,1,2-Trichloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Tetrachloroethene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,3-Dichloropropane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
2-Hexanone	ND	0.0065	EPA 8260C	5-26-17	5-26-17	
Dibromochloromethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromoethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Chlorobenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1,1,2-Tetrachloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Ethylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
n,p-Xylene	ND	0.0026	EPA 8260C	5-26-17	5-26-17	
o-Xylene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Styrene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Bromoform	ND	0.0065	EPA 8260C	5-26-17	5-26-17	
sopropylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Bromobenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
I,1,2,2-Tetrachloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichloropropane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
n-Propylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
2-Chlorotoluene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1-Chlorotoluene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,3,5-Trimethylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
ert-Butylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trimethylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
sec-Butylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,3-Dichlorobenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
o-Isopropyltoluene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,4-Dichlorobenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2-Dichlorobenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
n-Butylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromo-3-chloropropane		0.0065	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trichlorobenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Hexachlorobutadiene	ND	0.0065	EPA 8260C	5-26-17	5-26-17	
Naphthalene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichlorobenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits		0 =0	<u> </u>	
Dibromofluoromethane	115	73-134				
Toluene-d8	108	81-124				
4-Bromofluorobenzene	102	80-131				
+ DI OHIOHAOHODGHZGHG	102	00-131				

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP8-10					
Laboratory ID:	05-299-23					
Dichlorodifluoromethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Chloromethane	ND	0.0070	EPA 8260C	5-26-17	5-26-17	
Vinyl Chloride	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Bromomethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Chloroethane	ND	0.0070	EPA 8260C	5-26-17	5-26-17	
Trichlorofluoromethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Acetone	0.016	0.014	EPA 8260C	5-26-17	5-26-17	
lodomethane	ND	0.0070	EPA 8260C	5-26-17	5-26-17	
Carbon Disulfide	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Methylene Chloride	ND	0.014	EPA 8260C	5-26-17	5-26-17	
(trans) 1,2-Dichloroethene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Methyl t-Butyl Ether	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Vinyl Acetate	ND	0.0070	EPA 8260C	5-26-17	5-26-17	
2,2-Dichloropropane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
(cis) 1,2-Dichloroethene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
2-Butanone	ND	0.0070	EPA 8260C	5-26-17	5-26-17	
Bromochloromethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Chloroform	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,1,1-Trichloroethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Carbon Tetrachloride	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloropropene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Benzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloroethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Trichloroethene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloropropane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Dibromomethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Bromodichloromethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
2-Chloroethyl Vinyl Ether	ND	0.0089	EPA 8260C	5-26-17	5-26-17	
(cis) 1,3-Dichloropropene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Methyl Isobutyl Ketone	ND	0.0070	EPA 8260C	5-26-17	5-26-17	
Toluene	0.021	0.0070	EPA 8260C	5-26-17	5-26-17	
(trans) 1,3-Dichloropropene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	

Project: 15-05986-040

VOLATILES EPA 8260C

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP8-10					
Laboratory ID:	05-299-23					
1,1,2-Trichloroethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Tetrachloroethene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,3-Dichloropropane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
2-Hexanone	ND	0.0070	EPA 8260C	5-26-17	5-26-17	
Dibromochloromethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromoethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Chlorobenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,1,1,2-Tetrachloroethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Ethylbenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
m,p-Xylene	ND	0.0028	EPA 8260C	5-26-17	5-26-17	
o-Xylene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Styrene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Bromoform	ND	0.0070	EPA 8260C	5-26-17	5-26-17	
Isopropylbenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Bromobenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,1,2,2-Tetrachloroethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichloropropane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
n-Propylbenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
2-Chlorotoluene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
4-Chlorotoluene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,3,5-Trimethylbenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
tert-Butylbenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trimethylbenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
sec-Butylbenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,3-Dichlorobenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
p-Isopropyltoluene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,4-Dichlorobenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,2-Dichlorobenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
n-Butylbenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromo-3-chloropropane		0.0070	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trichlorobenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Hexachlorobutadiene	ND	0.0070	EPA 8260C	5-26-17	5-26-17	
Naphthalene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichlorobenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits		0 20 11	0 20 11	
Dibromofluoromethane	103	73-134				
Toluene-d8	102	81-124				
1 Oluene-do 4-Bromofluorobenzene	95	80-131				
4-DI OHTOHUOLODENZENE	90	00-131				

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP9-0					
Laboratory ID:	05-299-25					
Dichlorodifluoromethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Chloromethane	ND	0.0063	EPA 8260C	5-26-17	5-26-17	
Vinyl Chloride	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Bromomethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Chloroethane	ND	0.0063	EPA 8260C	5-26-17	5-26-17	
Trichlorofluoromethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Acetone	ND	0.013	EPA 8260C	5-26-17	5-26-17	
lodomethane	ND	0.0063	EPA 8260C	5-26-17	5-26-17	
Carbon Disulfide	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Methylene Chloride	ND	0.013	EPA 8260C	5-26-17	5-26-17	
(trans) 1,2-Dichloroethene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Methyl t-Butyl Ether	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Vinyl Acetate	ND	0.0063	EPA 8260C	5-26-17	5-26-17	
2,2-Dichloropropane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
(cis) 1,2-Dichloroethene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
2-Butanone	ND	0.0063	EPA 8260C	5-26-17	5-26-17	
Bromochloromethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Chloroform	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1,1-Trichloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Carbon Tetrachloride	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloropropene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Benzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Trichloroethene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloropropane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Dibromomethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Bromodichloromethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
2-Chloroethyl Vinyl Ether	ND	0.0080	EPA 8260C	5-26-17	5-26-17	
(cis) 1,3-Dichloropropene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Methyl Isobutyl Ketone	ND	0.0063	EPA 8260C	5-26-17	5-26-17	
Toluene	0.014	0.0063	EPA 8260C	5-26-17	5-26-17	
(trans) 1,3-Dichloropropene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	

Project: 15-05986-040

VOLATILES EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP9-0					
Laboratory ID:	05-299-25					
1,1,2-Trichloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Tetrachloroethene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,3-Dichloropropane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
2-Hexanone	ND	0.0063	EPA 8260C	5-26-17	5-26-17	
Dibromochloromethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromoethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Chlorobenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
1,1,1,2-Tetrachloroethane	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Ethylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
m,p-Xylene	ND	0.0025	EPA 8260C	5-26-17	5-26-17	
o-Xylene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Styrene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Bromoform	ND	0.0063	EPA 8260C	5-26-17	5-26-17	
Isopropylbenzene	ND	0.0013	EPA 8260C	5-26-17	5-26-17	
Bromobenzene	ND	0.057	EPA 8260C	5-30-17	5-30-17	
1,1,2,2-Tetrachloroethane	ND	0.057	EPA 8260C	5-30-17	5-30-17	
1,2,3-Trichloropropane	ND	0.057	EPA 8260C	5-30-17	5-30-17	
n-Propylbenzene	ND	0.057	EPA 8260C	5-30-17	5-30-17	
2-Chlorotoluene	ND	0.057	EPA 8260C	5-30-17	5-30-17	
4-Chlorotoluene	ND	0.057	EPA 8260C	5-30-17	5-30-17	
1,3,5-Trimethylbenzene	ND	0.057	EPA 8260C	5-30-17	5-30-17	
tert-Butylbenzene	ND	0.057	EPA 8260C	5-30-17	5-30-17	
1,2,4-Trimethylbenzene	ND	0.057	EPA 8260C	5-30-17	5-30-17	
sec-Butylbenzene	ND	0.057	EPA 8260C	5-30-17	5-30-17	
1,3-Dichlorobenzene	ND	0.057	EPA 8260C	5-30-17	5-30-17	
p-Isopropyltoluene	ND	0.057	EPA 8260C	5-30-17	5-30-17	
1,4-Dichlorobenzene	ND	0.057	EPA 8260C	5-30-17	5-30-17	
1,2-Dichlorobenzene	ND	0.057	EPA 8260C	5-30-17	5-30-17	
n-Butylbenzene	ND	0.057	EPA 8260C	5-30-17	5-30-17	
1,2-Dibromo-3-chloropropane		0.29	EPA 8260C	5-30-17	5-30-17	
1,2,4-Trichlorobenzene	ND	0.057	EPA 8260C	5-30-17	5-30-17	
Hexachlorobutadiene	ND	0.29	EPA 8260C	5-30-17	5-30-17	
Naphthalene	ND	0.057	EPA 8260C	5-30-17	5-30-17	
1,2,3-Trichlorobenzene	ND	0.057	EPA 8260C	5-30-17	5-30-17	
Surrogate:	Percent Recovery	Control Limits		-	-	
Dibromofluoromethane	105	73-134				
Toluene-d8	97	81-124				
		·- ·				

4-Bromofluorobenzene

80-131

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP9-10					
Laboratory ID:	05-299-26					
Dichlorodifluoromethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Chloromethane	ND	0.0062	EPA 8260C	5-26-17	5-26-17	
Vinyl Chloride	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Bromomethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Chloroethane	ND	0.0062	EPA 8260C	5-26-17	5-26-17	
Trichlorofluoromethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Acetone	0.015	0.012	EPA 8260C	5-26-17	5-26-17	
lodomethane	ND	0.0062	EPA 8260C	5-26-17	5-26-17	
Carbon Disulfide	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Methylene Chloride	ND	0.012	EPA 8260C	5-26-17	5-26-17	
(trans) 1,2-Dichloroethene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Methyl t-Butyl Ether	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Vinyl Acetate	ND	0.0062	EPA 8260C	5-26-17	5-26-17	
2,2-Dichloropropane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
(cis) 1,2-Dichloroethene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
2-Butanone	ND	0.0062	EPA 8260C	5-26-17	5-26-17	
Bromochloromethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Chloroform	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,1,1-Trichloroethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Carbon Tetrachloride	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloropropene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Benzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloroethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Trichloroethene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloropropane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Dibromomethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Bromodichloromethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
2-Chloroethyl Vinyl Ether	ND	0.0079	EPA 8260C	5-26-17	5-26-17	
(cis) 1,3-Dichloropropene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Methyl Isobutyl Ketone	ND	0.0062	EPA 8260C	5-26-17	5-26-17	
Toluene	0.015	0.0062	EPA 8260C	5-26-17	5-26-17	
(trans) 1,3-Dichloropropene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	

Project: 15-05986-040

VOLATILES EPA 8260C

Client ID:					Date	Date	
Laboratory ID: 05-299-26 1,1,2-Trichloroethane ND 0.0012 EPA 8260C 5-26-17 5-26-17 Tetrachloroethane ND 0.0012 EPA 8260C 5-26-17 5-26-17 Tetrachloroethane ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,3-Dichloropropane ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,3-Dichloropropane ND 0.0062 EPA 8260C 5-26-17 5-26-17 1,2-Dibromoethane ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2-Dibromoethane ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,1,1-2-Tetrachloroethane ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,1,1,1-2-Tetrachloroethane ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,1,1,1-1-10-Tetrachloroethane ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,1,1,1-10-Tetrachloroethane ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,1,1-1-10-Tetrachloroethane ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,1,1-1-10-Tetrachloroethane ND 0.0012			PQL	Method	Prepared	Analyzed	Flags
1,1,2-Trichloroethane							
Tetrachloroethene ND 0.0012 EPA 8260C 5-26-17 5-26-17 [1,3-Dichloropropane] ND 0.0012 EPA 8260C 5-26-17 5-26-17 5-26-17 [2-Dibromochloromethane] ND 0.0062 EPA 8260C 5-26-17 5-26-17 [2-Dibromochloromethane] ND 0.0012 EPA 8260C 5-26-17 5-26-17 [1,1,2-Tetrachloroethane] ND 0.0012 EPA 8260C 5-26-17 5-26-17 [1,1,1,2-Tetrachloroethane] ND 0.0012 EPA 8260C 5-26-17 5-26-17 [2-Dibromochloromethane] ND 0.0							
1,3-Dichloropropane ND 0.0012 EPA 8260C 5-26-17 5-26-17 2-Hexanone ND 0.0062 EPA 8260C 5-26-17 5-26-17 Dibromochloromethane ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2-Dibromoethane ND 0.0012 EPA 8260C 5-26-17 5-26-17 Chlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 Chlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 Ethylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 Ethylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 Syylene ND 0.0012 EPA 8260C 5-26-17 5-26-17 Bromoform ND 0.0062 EPA 8260C 5-26-17 5-26-17 Bromoform ND 0.0012 EPA 8260C 5-26-17 5-26-17 Sopropylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17							
2-Hexanone							
Dibromochloromethane	1,3-Dichloropropane		0.0012	EPA 8260C		5-26-17	
1,2-Dibromoethane	2-Hexanone	ND	0.0062	EPA 8260C	5-26-17	5-26-17	
Chlorobenzene	Dibromochloromethane		0.0012	EPA 8260C	5-26-17	5-26-17	
1,1,1,2-Tetrachloroethane	1,2-Dibromoethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Ethylbenzene	Chlorobenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
N.D. N.D. N.D. 0.0025 EPA 8260C 5-26-17 5-	1,1,1,2-Tetrachloroethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
December ND December September ND December December September ND December Decem	Ethylbenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Styrene ND 0.0012 EPA 8260C 5-26-17 5-26-17 Sepromoform ND 0.0062 EPA 8260C 5-26-17 5-26-17 Sepromoform ND 0.0012 EPA 8260C 5-26-17 5-26-17 Sepromobenzene ND 0.0012 EPA 8260C	m,p-Xylene	ND	0.0025	EPA 8260C	5-26-17	5-26-17	
Second Form ND 0.0062 EPA 8260C 5-26-17 5-26	o-Xylene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Sepropy S	Styrene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Second beautiful Second Processes ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,1,2,2-Tetrachloroethane ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichloropropane ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichloropropane ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichloropropane ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,3,5-Trimethylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,3,5-Trimethylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,3,5-Trimethylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2,4-Trimethylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2,4-Trimethylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,3-Dichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,3-Dichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,3-Dichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,4-Dichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2-Dichlorobenzene ND	Bromoform	ND	0.0062	EPA 8260C	5-26-17	5-26-17	
1,1,2,2-Tetrachloroethane ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichloropropane ND 0.0012 EPA 8260C 5-26-17 5-26-17 n-Propylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 2-Chlorotoluene ND 0.0012 EPA 8260C 5-26-17 5-26-17 4-Chlorotoluene ND 0.0012 EPA 8260C 5-26-17 5-26-17 4-Chlorotoluene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,3,5-Trimethylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1-1,2,4-Trimethylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1-2,4-Trimethylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1-3,5-Dichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,3-Dichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,4-Dichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2-Dichlorobenzene	Isopropylbenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichloropropane	Bromobenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
n-Propylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,3-5-Trimethylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,3-5-Trimethylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,3-5-Trimethylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,3-4-Trimethylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,3-Dichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,4-Dichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,4-Dichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2-Dichlorobenzene ND	1,1,2,2-Tetrachloroethane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
C-Chlorotoluene	1,2,3-Trichloropropane	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
C-Chlorotoluene	n-Propylbenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,3,5-Trimethylbenzene	2-Chlorotoluene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
tert-Butylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 12,4-Trimethylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-1	4-Chlorotoluene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
Itert-Butylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2,4-Trimethylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 sec-Butylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,3-Dichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,4-Dichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,4-Dichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2-Dichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 n-Butylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 n-Butylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 n-Butylbenzene ND 0.0062 EPA 8260C 5-26-17 5-26-17 1,2-Dibromo-3-chloropropane ND 0.0012 EPA 8260C 5-26-17 5-26-17 Hexachlorobutadiene ND 0.0062 EPA	1,3,5-Trimethylbenzene	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trimethylbenzene		ND	0.0012	EPA 8260C	5-26-17	5-26-17	
See-Butylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,3-Dichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,3-Dichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,4-Dichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2-Dichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2-Dichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2-Dibromo-3-chloropropane ND 0.0062 EPA 8260C 5-26-17 5-26-17 1,2,4-Trichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 Hexachlorobutadiene ND 0.0062 EPA 8260C 5-26-17 5-26-17 Naphthalene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 106 73-134 Toluene-d8 105 81-124	·	ND	0.0012	EPA 8260C	5-26-17	5-26-17	
1,3-Dichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 p-Isopropyltoluene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,4-Dichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2-Dichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 n-Butylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2-Dibromo-3-chloropropane ND 0.0062 EPA 8260C 5-26-17 5-26-17 1,2,4-Trichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 Hexachlorobutadiene ND 0.0062 EPA 8260C 5-26-17 5-26-17 Naphthalene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 106 73-134 Toluene-d8 105 81-124		ND					
Description ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,4-Dichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2-Dichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 5-26-17 1,2-Dibromo-3-chloropropane ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2,4-Trichlorobenzene ND 0.0062 EPA 8260C 5-26-17 5-26-17 1,2,4-Trichlorobenzene ND 0.0062 EPA 8260C 5-26-17 5-26-17 1,2,4-Trichlorobenzene ND 0.0062 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 106 73-134 Toluene-d8 105 81-124 105 81-124 105 105 100		ND	0.0012		5-26-17	5-26-17	
1,4-Dichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2-Dichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 n-Butylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2-Dibromo-3-chloropropane ND 0.0062 EPA 8260C 5-26-17 5-26-17 1,2,4-Trichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 Hexachlorobutadiene ND 0.0062 EPA 8260C 5-26-17 5-26-17 Naphthalene ND 0.0062 EPA 8260C 5-26-17 5-26-17 Naphthalene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 106 73-134 Toluene-d8 105 81-124		ND	0.0012		5-26-17		
1,2-Dichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 n-Butylbenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2-Dibromo-3-chloropropane ND 0.0062 EPA 8260C 5-26-17 5-26-17 1,2,4-Trichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 Hexachlorobutadiene ND 0.0062 EPA 8260C 5-26-17 5-26-17 Naphthalene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 106 73-134 Toluene-d8 105 81-124							
ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2-Dibromo-3-chloropropane ND 0.0062 EPA 8260C 5-26-17 5-26-17 1,2,4-Trichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 5-26-17 Hexachlorobutadiene ND 0.0062 EPA 8260C 5-26-17 5-26-17 5-26-17 Naphthalene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 106 73-134 Toluene-d8 105 81-124 ND NO NO NO NO NO NO NO							
1,2-Dibromo-3-chloropropane ND 0.0062 EPA 8260C 5-26-17 5-26-17 1,2,4-Trichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 Hexachlorobutadiene ND 0.0062 EPA 8260C 5-26-17 5-26-17 Naphthalene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 106 73-134 Toluene-d8 105 81-124							
1,2,4-Trichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 Hexachlorobutadiene ND 0.0062 EPA 8260C 5-26-17 5-26-17 Naphthalene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 106 73-134 Toluene-d8 105 81-124	-						
Hexachlorobutadiene ND 0.0062 EPA 8260C 5-26-17 5-26-17 Naphthalene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 106 73-134 Toluene-d8 105 81-124							
Naphthalene ND 0.0012 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 106 73-134 Toluene-d8 105 81-124							
1,2,3-Trichlorobenzene ND 0.0012 EPA 8260C 5-26-17 5-26-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 106 73-134 Toluene-d8 105 81-124							
Surrogate: Percent Recovery Control Limits Dibromofluoromethane 106 73-134 Toluene-d8 105 81-124	•						
Dibromofluoromethane 106 73-134 Toluene-d8 105 81-124					0 _0	0 = 0	
Toluene-d8 105 81-124							
	4-Bromofluorobenzene	102	80-131				

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP10-0					
Laboratory ID:	05-299-28					
Dichlorodifluoromethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Chloromethane	ND	0.0070	EPA 8260C	5-26-17	5-26-17	
Vinyl Chloride	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Bromomethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Chloroethane	ND	0.0070	EPA 8260C	5-26-17	5-26-17	
Trichlorofluoromethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Acetone	ND	0.014	EPA 8260C	5-26-17	5-26-17	
lodomethane	ND	0.0070	EPA 8260C	5-26-17	5-26-17	
Carbon Disulfide	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Methylene Chloride	ND	0.014	EPA 8260C	5-26-17	5-26-17	
(trans) 1,2-Dichloroethene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Methyl t-Butyl Ether	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Vinyl Acetate	ND	0.0070	EPA 8260C	5-26-17	5-26-17	
2,2-Dichloropropane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
(cis) 1,2-Dichloroethene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
2-Butanone	ND	0.0070	EPA 8260C	5-26-17	5-26-17	
Bromochloromethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Chloroform	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,1,1-Trichloroethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Carbon Tetrachloride	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloropropene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Benzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloroethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Trichloroethene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloropropane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Dibromomethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Bromodichloromethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
2-Chloroethyl Vinyl Ether	ND	0.0090	EPA 8260C	5-26-17	5-26-17	
(cis) 1,3-Dichloropropene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Methyl Isobutyl Ketone	ND	0.0070	EPA 8260C	5-26-17	5-26-17	
Toluene	ND	0.0070	EPA 8260C	5-26-17	5-26-17	
(trans) 1,3-Dichloropropene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	

Project: 15-05986-040

VOLATILES EPA 8260C

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP10-0					
Laboratory ID:	05-299-28					
1,1,2-Trichloroethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Γetrachloroethene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,3-Dichloropropane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
2-Hexanone	ND	0.0070	EPA 8260C	5-26-17	5-26-17	
Dibromochloromethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromoethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Chlorobenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,1,1,2-Tetrachloroethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Ethylbenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
n,p-Xylene	ND	0.0028	EPA 8260C	5-26-17	5-26-17	
o-Xylene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Styrene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Bromoform	ND	0.0070	EPA 8260C	5-26-17	5-26-17	
sopropylbenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Bromobenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
I,1,2,2-Tetrachloroethane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichloropropane	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
n-Propylbenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
2-Chlorotoluene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1-Chlorotoluene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,3,5-Trimethylbenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
ert-Butylbenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trimethylbenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
sec-Butylbenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,3-Dichlorobenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
o-Isopropyltoluene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
,4-Dichlorobenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,2-Dichlorobenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
n-Butylbenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromo-3-chloropropane		0.0070	EPA 8260C	5-26-17	5-26-17	
I,2,4-Trichlorobenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Hexachlorobutadiene	ND	0.0070	EPA 8260C	5-26-17	5-26-17	
Naphthalene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichlorobenzene	ND	0.0014	EPA 8260C	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	103	73-134				
Toluene-d8	98	81-124				
4-Bromofluorobenzene	90	80-131				
+ DIGITIONIUOIODGITZGITG	30	00-131				

Project: 15-05986-040

VOLATILES by EPA 8260C METHOD BLANK QUALITY CONTROL

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0525S1					
Dichlorodifluoromethane	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
Chloromethane	ND	0.0050	EPA 8260C	5-25-17	5-25-17	
Vinyl Chloride	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
Bromomethane	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
Chloroethane	ND	0.0050	EPA 8260C	5-25-17	5-25-17	
Trichlorofluoromethane	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
1,1-Dichloroethene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
Acetone	ND	0.010	EPA 8260C	5-25-17	5-25-17	
lodomethane	ND	0.0063	EPA 8260C	5-25-17	5-25-17	
Carbon Disulfide	ND	0.0014	EPA 8260C	5-25-17	5-25-17	
Methylene Chloride	ND	0.010	EPA 8260C	5-25-17	5-25-17	
(trans) 1,2-Dichloroethene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
Methyl t-Butyl Ether	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
1,1-Dichloroethane	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
Vinyl Acetate	ND	0.0050	EPA 8260C	5-25-17	5-25-17	
2,2-Dichloropropane	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
(cis) 1,2-Dichloroethene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
2-Butanone	ND	0.0050	EPA 8260C	5-25-17	5-25-17	
Bromochloromethane	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
Chloroform	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
1,1,1-Trichloroethane	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
Carbon Tetrachloride	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
1,1-Dichloropropene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
Benzene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
1,2-Dichloroethane	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
Trichloroethene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
1,2-Dichloropropane	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
Dibromomethane	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
Bromodichloromethane	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
2-Chloroethyl Vinyl Ether	ND	0.0050	EPA 8260C	5-25-17	5-25-17	
(cis) 1,3-Dichloropropene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
Methyl Isobutyl Ketone	ND	0.0050	EPA 8260C	5-25-17	5-25-17	
Toluene	ND	0.0050	EPA 8260C	5-25-17	5-25-17	
(trans) 1,3-Dichloropropene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	

Project: 15-05986-040

VOLATILES by EPA 8260C METHOD BLANK QUALITY CONTROL

page 2 of 2

Amalada	Decell	DOI	B# a4b a al	Date	Date	-
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0525S1					
1,1,2-Trichloroethane	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
Tetrachloroethene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
1,3-Dichloropropane	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
2-Hexanone	ND	0.0050	EPA 8260C	5-25-17	5-25-17	
Dibromochloromethane	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
1,2-Dibromoethane	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
Chlorobenzene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
1,1,1,2-Tetrachloroethane	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
Ethylbenzene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
m,p-Xylene	ND	0.0020	EPA 8260C	5-25-17	5-25-17	
o-Xylene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
Styrene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
Bromoform	ND	0.0050	EPA 8260C	5-25-17	5-25-17	
Isopropylbenzene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
Bromobenzene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
1,1,2,2-Tetrachloroethane	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
1,2,3-Trichloropropane	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
n-Propylbenzene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
2-Chlorotoluene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
4-Chlorotoluene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
1,3,5-Trimethylbenzene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
tert-Butylbenzene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
1,2,4-Trimethylbenzene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
sec-Butylbenzene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
1,3-Dichlorobenzene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
p-Isopropyltoluene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
1,4-Dichlorobenzene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
1,2-Dichlorobenzene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
n-Butylbenzene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
1,2-Dibromo-3-chloropropane		0.0050	EPA 8260C	5-25-17	5-25-17	
1,2,4-Trichlorobenzene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
Hexachlorobutadiene	ND	0.0050	EPA 8260C	5-25-17	5-25-17	
Naphthalene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
1,2,3-Trichlorobenzene	ND	0.0010	EPA 8260C	5-25-17	5-25-17	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	109	73-134				
Toluene-d8	108	81-124				
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4-Bromofluorobenzene

80-131

Project: 15-05986-040

VOLATILES by EPA 8260C METHOD BLANK QUALITY CONTROL

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0526S1					
Dichlorodifluoromethane	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
Chloromethane	ND	0.0050	EPA 8260C	5-26-17	5-26-17	
Vinyl Chloride	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
Bromomethane	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
Chloroethane	ND	0.0050	EPA 8260C	5-26-17	5-26-17	
Trichlorofluoromethane	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
Acetone	ND	0.010	EPA 8260C	5-26-17	5-26-17	
lodomethane	ND	0.0050	EPA 8260C	5-26-17	5-26-17	
Carbon Disulfide	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
Methylene Chloride	ND	0.010	EPA 8260C	5-26-17	5-26-17	
(trans) 1,2-Dichloroethene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
Methyl t-Butyl Ether	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethane	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
Vinyl Acetate	ND	0.0050	EPA 8260C	5-26-17	5-26-17	
2,2-Dichloropropane	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
(cis) 1,2-Dichloroethene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
2-Butanone	ND	0.0050	EPA 8260C	5-26-17	5-26-17	
Bromochloromethane	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
Chloroform	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
1,1,1-Trichloroethane	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
Carbon Tetrachloride	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloropropene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
Benzene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloroethane	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
Trichloroethene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloropropane	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
Dibromomethane	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
Bromodichloromethane	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
2-Chloroethyl Vinyl Ether	ND	0.0064	EPA 8260C	5-26-17	5-26-17	
(cis) 1,3-Dichloropropene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
Methyl Isobutyl Ketone	ND	0.0050	EPA 8260C	5-26-17	5-26-17	
Toluene	ND	0.0050	EPA 8260C	5-26-17	5-26-17	
(trans) 1,3-Dichloropropene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	

Project: 15-05986-040

VOLATILES by EPA 8260C METHOD BLANK QUALITY CONTROL

page 2 of 2

Amalada	Decel	DOI	Madh a d	Date	Date	FI
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0526S1					
1,1,2-Trichloroethane	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
Tetrachloroethene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
1,3-Dichloropropane	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
2-Hexanone	ND	0.0050	EPA 8260C	5-26-17	5-26-17	
Dibromochloromethane	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromoethane	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
Chlorobenzene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
1,1,1,2-Tetrachloroethane	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
Ethylbenzene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
m,p-Xylene	ND	0.0020	EPA 8260C	5-26-17	5-26-17	
o-Xylene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
Styrene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
Bromoform	ND	0.0050	EPA 8260C	5-26-17	5-26-17	
Isopropylbenzene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
Bromobenzene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
1,1,2,2-Tetrachloroethane	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichloropropane	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
n-Propylbenzene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
2-Chlorotoluene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
4-Chlorotoluene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
1,3,5-Trimethylbenzene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
tert-Butylbenzene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trimethylbenzene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
sec-Butylbenzene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
1,3-Dichlorobenzene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
p-Isopropyltoluene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
1,4-Dichlorobenzene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
1,2-Dichlorobenzene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
n-Butylbenzene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromo-3-chloropropane		0.0050	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trichlorobenzene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
Hexachlorobutadiene	ND	0.0050	EPA 8260C	5-26-17	5-26-17	
Naphthalene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichlorobenzene	ND	0.0010	EPA 8260C	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	113	73-134				
Toluene-d8	111	81-124				
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4-Bromofluorobenzene

80-131

Project: 15-05986-040

VOLATILES by EPA 8260C METHOD BLANK QUALITY CONTROL

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0530S1					
Dichlorodifluoromethane	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
Chloromethane	ND	0.0050	EPA 8260C	5-30-17	5-30-17	
Vinyl Chloride	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
Bromomethane	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
Chloroethane	ND	0.0050	EPA 8260C	5-30-17	5-30-17	
Trichlorofluoromethane	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
1,1-Dichloroethene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
Acetone	ND	0.010	EPA 8260C	5-30-17	5-30-17	
lodomethane	ND	0.0050	EPA 8260C	5-30-17	5-30-17	
Carbon Disulfide	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
Methylene Chloride	ND	0.010	EPA 8260C	5-30-17	5-30-17	
(trans) 1,2-Dichloroethene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
Methyl t-Butyl Ether	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
1,1-Dichloroethane	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
Vinyl Acetate	ND	0.0050	EPA 8260C	5-30-17	5-30-17	
2,2-Dichloropropane	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
(cis) 1,2-Dichloroethene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
2-Butanone	ND	0.0050	EPA 8260C	5-30-17	5-30-17	
Bromochloromethane	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
Chloroform	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
1,1,1-Trichloroethane	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
Carbon Tetrachloride	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
1,1-Dichloropropene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
Benzene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
1,2-Dichloroethane	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
Trichloroethene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
1,2-Dichloropropane	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
Dibromomethane	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
Bromodichloromethane	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
2-Chloroethyl Vinyl Ether	ND	0.0050	EPA 8260C	5-30-17	5-30-17	
(cis) 1,3-Dichloropropene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
Methyl Isobutyl Ketone	ND	0.0050	EPA 8260C	5-30-17	5-30-17	
Toluene	ND	0.0050	EPA 8260C	5-30-17	5-30-17	
(trans) 1,3-Dichloropropene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	

Project: 15-05986-040

VOLATILES by EPA 8260C METHOD BLANK QUALITY CONTROL

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0530S1					
1,1,2-Trichloroethane	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
Tetrachloroethene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
1,3-Dichloropropane	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
2-Hexanone	ND	0.0050	EPA 8260C	5-30-17	5-30-17	
Dibromochloromethane	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
1,2-Dibromoethane	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
Chlorobenzene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
1,1,1,2-Tetrachloroethane	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
Ethylbenzene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
m,p-Xylene	ND	0.0020	EPA 8260C	5-30-17	5-30-17	
o-Xylene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
Styrene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
Bromoform	ND	0.0050	EPA 8260C	5-30-17	5-30-17	
Isopropylbenzene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
Bromobenzene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
1,1,2,2-Tetrachloroethane	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
1,2,3-Trichloropropane	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
n-Propylbenzene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
2-Chlorotoluene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
4-Chlorotoluene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
1,3,5-Trimethylbenzene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
tert-Butylbenzene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
1,2,4-Trimethylbenzene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
sec-Butylbenzene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
1,3-Dichlorobenzene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
p-Isopropyltoluene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
1,4-Dichlorobenzene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
1,2-Dichlorobenzene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
n-Butylbenzene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
1,2-Dibromo-3-chloropropane	ND	0.0050	EPA 8260C	5-30-17	5-30-17	
1,2,4-Trichlorobenzene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
Hexachlorobutadiene	ND	0.0050	EPA 8260C	5-30-17	5-30-17	
Naphthalene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
1,2,3-Trichlorobenzene	ND	0.0010	EPA 8260C	5-30-17	5-30-17	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	107	73-134				
Toluene-d8	110	81-124				
. 5.50110 40		0. 121				

4-Bromofluorobenzene

80-131

Project: 15-05986-040

VOLATILES by EPA 8260C SB/SBD QUALITY CONTROL

					Per	cent	Recovery		RPD	
Analyte	Result		Spike	Spike Level		overy	Limits	RPD	Limit	Flags
SPIKE BLANKS										
Laboratory ID:	SB05	25S1								
	SB	SBD	SB	SBD	SB	SBD				
1,1-Dichloroethene	0.0613	0.0591	0.0500	0.0500	123	118	66-127	4	15	
Benzene	0.0547	0.0554	0.0500	0.0500	109	111	76-122	1	15	
Trichloroethene	0.0542	0.0536	0.0500	0.0500	108	107	78-120	1	15	
Toluene	0.0544	0.0552	0.0500	0.0500	109	110	83-120	1	15	
Chlorobenzene	0.0509	0.0511	0.0500	0.0500	102	102	81-120	0	15	
Surrogate:										
Dibromofluoromethane					100	102	73-134			
Toluene-d8					101	103	81-124			
4-Bromofluorobenzene					97	99	80-131			

Project: 15-05986-040

VOLATILES by EPA 8260C SB/SBD QUALITY CONTROL

Matrix: Soil Units: mg/kg

Analyte	Res	sult	Spike	Level		cent overy	Recovery Limits	RPD	RPD Limit	Flags
SPIKE BLANKS										
Laboratory ID:	SB05	26S1								
	SB	SBD	SB	SBD	SB	SBD				
1,1-Dichloroethene	0.0554	0.0593	0.0500	0.0500	111	119	66-127	7	15	
Benzene	0.0527	0.0554	0.0500	0.0500	105	111	76-122	5	15	
Trichloroethene	0.0482	0.0511	0.0500	0.0500	96	102	78-120	6	15	
Toluene	0.0498	0.0534	0.0500	0.0500	100	107	83-120	7	15	
Chlorobenzene	0.0478	0.0483	0.0500	0.0500	96	97	81-120	1	15	
Surrogate:										
Dibromofluoromethane					103	104	73-134			
Toluene-d8					99	99	81-124			
4-Bromofluorobenzene					97	97	80-131			

Project: 15-05986-040

VOLATILES by EPA 8260C SB/SBD QUALITY CONTROL

Matrix: Soil Units: mg/kg

					Per	cent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Reco	overy	Limits	RPD	Limit	Flags
SPIKE BLANKS										
Laboratory ID:	SB05	30S1								
	SB	SBD	SB	SBD	SB	SBD				
1,1-Dichloroethene	0.0613	0.0542	0.0500	0.0500	123	108	66-127	12	15	
Benzene	0.0567	0.0513	0.0500	0.0500	113	103	76-122	10	15	
Trichloroethene	0.0551	0.0498	0.0500	0.0500	110	100	78-120	10	15	
Toluene	0.0560	0.0511	0.0500	0.0500	112	102	83-120	9	15	
Chlorobenzene	0.0510	0.0466	0.0500	0.0500	102	93	81-120	9	15	
Surrogate:										
Dibromofluoromethane					98	94	73-134			
Toluene-d8					100	98	81-124			
4-Bromofluorobenzene					94	85	80-131			

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP1-W					
Laboratory ID:	05-299-02					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Chloromethane	ND	1.0	EPA 8260C	5-30-17	5-30-17	
Vinyl Chloride	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Bromomethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Chloroethane	ND	1.0	EPA 8260C	5-30-17	5-30-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Acetone	6.7	5.0	EPA 8260C	5-30-17	5-30-17	
Iodomethane	ND	1.0	EPA 8260C	5-30-17	5-30-17	
Carbon Disulfide	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Methylene Chloride	ND	1.0	EPA 8260C	5-30-17	5-30-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Vinyl Acetate	ND	1.0	EPA 8260C	5-30-17	5-30-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
2-Butanone	ND	5.0	EPA 8260C	5-30-17	5-30-17	
Bromochloromethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Chloroform	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Benzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Trichloroethene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Dibromomethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Bromodichloromethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	5-30-17	5-30-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	5-30-17	5-30-17	
Toluene	ND	1.0	EPA 8260C	5-30-17	5-30-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-30-17	5-30-17	

Project: 15-05986-040

VOLATILES EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP1-W					
Laboratory ID:	05-299-02					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Tetrachloroethene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
2-Hexanone	ND	2.0	EPA 8260C	5-30-17	5-30-17	
Dibromochloromethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Chlorobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Ethylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
m,p-Xylene	ND	0.40	EPA 8260C	5-30-17	5-30-17	
o-Xylene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Styrene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Bromoform	ND	1.0	EPA 8260C	5-30-17	5-30-17	
Isopropylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Bromobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
n-Propylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
tert-Butylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
sec-Butylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
n-Butylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	5-30-17	5-30-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Naphthalene	ND	1.0	EPA 8260C	5-30-17	5-30-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	113	77-129				
-						

 Dibromofluoromethane
 113
 77-129

 Toluene-d8
 113
 80-127

 4-Bromofluorobenzene
 84
 80-125



Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP2-W					
Laboratory ID:	05-299-06					
Dichlorodifluoromethane	ND	0.26	EPA 8260C	5-26-17	5-26-17	
Chloromethane	ND	1.4	EPA 8260C	5-26-17	5-26-17	
Vinyl Chloride	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromomethane	ND	0.33	EPA 8260C	5-26-17	5-26-17	
Chloroethane	ND	1.0	EPA 8260C	5-26-17	5-26-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Acetone	75	5.0	EPA 8260C	5-26-17	5-26-17	
Iodomethane	ND	1.7	EPA 8260C	5-26-17	5-26-17	
Carbon Disulfide	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Methylene Chloride	ND	1.0	EPA 8260C	5-26-17	5-26-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Vinyl Acetate	ND	1.0	EPA 8260C	5-26-17	5-26-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Butanone	ND	5.0	EPA 8260C	5-26-17	5-26-17	
Bromochloromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Chloroform	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Benzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Trichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Dibromomethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromodichloromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Chloroethyl Vinyl Ether	ND	6.6	EPA 8260C	5-26-17	5-26-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	5-26-17	5-26-17	
Toluene	ND	1.0	EPA 8260C	5-26-17	5-26-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-26-17	5-26-17	

Project: 15-05986-040

VOLATILES EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP2-W					
Laboratory ID:	05-299-06					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Tetrachloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Hexanone	ND	2.0	EPA 8260C	5-26-17	5-26-17	
Dibromochloromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Chlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Ethylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
m,p-Xylene	ND	0.40	EPA 8260C	5-26-17	5-26-17	
o-Xylene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Styrene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromoform	ND	1.0	EPA 8260C	5-26-17	5-26-17	
Isopropylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
n-Propylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
tert-Butylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
sec-Butylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
n-Butylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Naphthalene	ND	1.4	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	102	77-129				
-						

 Dibromofluoromethane
 102
 77-129

 Toluene-d8
 84
 80-127

 4-Bromofluorobenzene
 123
 80-125



Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP3-W					
Laboratory ID:	05-299-09					
Dichlorodifluoromethane	ND	0.26	EPA 8260C	5-26-17	5-26-17	
Chloromethane	ND	1.4	EPA 8260C	5-26-17	5-26-17	
Vinyl Chloride	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromomethane	ND	0.33	EPA 8260C	5-26-17	5-26-17	
Chloroethane	ND	1.0	EPA 8260C	5-26-17	5-26-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Acetone	5.3	5.0	EPA 8260C	5-26-17	5-26-17	
Iodomethane	ND	1.7	EPA 8260C	5-26-17	5-26-17	
Carbon Disulfide	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Methylene Chloride	ND	1.0	EPA 8260C	5-26-17	5-26-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Vinyl Acetate	ND	1.0	EPA 8260C	5-26-17	5-26-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Butanone	ND	5.0	EPA 8260C	5-26-17	5-26-17	
Bromochloromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Chloroform	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Benzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Trichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Dibromomethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromodichloromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Chloroethyl Vinyl Ether	ND	6.6	EPA 8260C	5-26-17	5-26-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	5-26-17	5-26-17	
Toluene	ND	1.0	EPA 8260C	5-26-17	5-26-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-26-17	5-26-17	

Project: 15-05986-040

VOLATILES EPA 8260C

page 2 of 2

Amalista	Dec. 16	DC:	M -41 - 1	Date	Date	-
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP3-W					
Laboratory ID:	05-299-09					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Tetrachloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Hexanone	ND	2.0	EPA 8260C	5-26-17	5-26-17	
Dibromochloromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Chlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Ethylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
m,p-Xylene	ND	0.40	EPA 8260C	5-26-17	5-26-17	
o-Xylene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Styrene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromoform	ND	1.0	EPA 8260C	5-26-17	5-26-17	
Isopropylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
n-Propylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
tert-Butylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
sec-Butylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
n-Butylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromo-3-chloropropane		1.0	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Naphthalene	ND	1.4	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits		0 _ 0	0 = 0 11	
Dibromofluoromethane	101	77-129				
Toluene-d8	83	80-127				
i Gladi Id-ad	00	00-121				

4-Bromofluorobenzene

80-125

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP4-W					
Laboratory ID:	05-299-12					
Dichlorodifluoromethane	ND	0.26	EPA 8260C	5-26-17	5-26-17	
Chloromethane	ND	1.4	EPA 8260C	5-26-17	5-26-17	
Vinyl Chloride	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromomethane	ND	0.33	EPA 8260C	5-26-17	5-26-17	
Chloroethane	ND	1.0	EPA 8260C	5-26-17	5-26-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Acetone	5.6	5.0	EPA 8260C	5-26-17	5-26-17	
lodomethane	ND	1.7	EPA 8260C	5-26-17	5-26-17	
Carbon Disulfide	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Methylene Chloride	ND	1.0	EPA 8260C	5-26-17	5-26-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Vinyl Acetate	ND	1.0	EPA 8260C	5-26-17	5-26-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Butanone	ND	5.0	EPA 8260C	5-26-17	5-26-17	
Bromochloromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Chloroform	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Benzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Trichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Dibromomethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromodichloromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Chloroethyl Vinyl Ether	ND	6.6	EPA 8260C	5-26-17	5-26-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	5-26-17	5-26-17	
Toluene	ND	1.0	EPA 8260C	5-26-17	5-26-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-26-17	5-26-17	

Project: 15-05986-040

VOLATILES EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP4-W					
Laboratory ID:	05-299-12					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Tetrachloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Hexanone	ND	2.0	EPA 8260C	5-26-17	5-26-17	
Dibromochloromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Chlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Ethylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
m,p-Xylene	ND	0.40	EPA 8260C	5-26-17	5-26-17	
o-Xylene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Styrene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromoform	ND	1.0	EPA 8260C	5-26-17	5-26-17	
Isopropylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
n-Propylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
tert-Butylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
sec-Butylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
n-Butylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Naphthalene	ND	1.4	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	99	77-129				
-						

4-Bromofluorobenzene

Toluene-d8

80-127

80-125

81

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP5-W					
Laboratory ID:	05-299-15					
Dichlorodifluoromethane	ND	0.26	EPA 8260C	5-26-17	5-26-17	
Chloromethane	ND	1.4	EPA 8260C	5-26-17	5-26-17	
Vinyl Chloride	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromomethane	ND	0.33	EPA 8260C	5-26-17	5-26-17	
Chloroethane	ND	1.0	EPA 8260C	5-26-17	5-26-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Acetone	5.5	5.0	EPA 8260C	5-26-17	5-26-17	
lodomethane	ND	1.7	EPA 8260C	5-26-17	5-26-17	
Carbon Disulfide	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Methylene Chloride	ND	1.0	EPA 8260C	5-26-17	5-26-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Vinyl Acetate	ND	1.0	EPA 8260C	5-26-17	5-26-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Butanone	ND	5.0	EPA 8260C	5-26-17	5-26-17	
Bromochloromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Chloroform	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Benzene	6.4	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Trichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Dibromomethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromodichloromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Chloroethyl Vinyl Ether	ND	6.6	EPA 8260C	5-26-17	5-26-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	5-26-17	5-26-17	
Toluene	ND	1.0	EPA 8260C	5-26-17	5-26-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-26-17	5-26-17	

Project: 15-05986-040

VOLATILES EPA 8260C page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP5-W					
Laboratory ID:	05-299-15					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Tetrachloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Hexanone	ND	2.0	EPA 8260C	5-26-17	5-26-17	
Dibromochloromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Chlorobenzene	0.43	0.20	EPA 8260C	5-26-17	5-26-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Ethylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
m,p-Xylene	32	0.40	EPA 8260C	5-26-17	5-26-17	
o-Xylene	0.57	0.20	EPA 8260C	5-26-17	5-26-17	
Styrene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromoform	ND	1.0	EPA 8260C	5-26-17	5-26-17	
Isopropylbenzene	0.45	0.20	EPA 8260C	5-26-17	5-26-17	
Bromobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
n-Propylbenzene	0.95	0.20	EPA 8260C	5-26-17	5-26-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,3,5-Trimethylbenzene	3.1	0.20	EPA 8260C	5-26-17	5-26-17	
tert-Butylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trimethylbenzene	12	0.20	EPA 8260C	5-26-17	5-26-17	
sec-Butylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
n-Butylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Naphthalene	ND	1.4	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	100	77-129				
Toluene-d8	90	80-127				

4-Bromofluorobenzene

80-125

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP6-W					
Laboratory ID:	05-299-18					
Dichlorodifluoromethane	ND	0.26	EPA 8260C	5-26-17	5-26-17	
Chloromethane	ND	1.4	EPA 8260C	5-26-17	5-26-17	
Vinyl Chloride	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromomethane	ND	0.33	EPA 8260C	5-26-17	5-26-17	
Chloroethane	ND	1.0	EPA 8260C	5-26-17	5-26-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Acetone	7.9	5.0	EPA 8260C	5-26-17	5-26-17	
Iodomethane	ND	1.7	EPA 8260C	5-26-17	5-26-17	
Carbon Disulfide	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Methylene Chloride	ND	1.0	EPA 8260C	5-26-17	5-26-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Vinyl Acetate	ND	1.0	EPA 8260C	5-26-17	5-26-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Butanone	ND	5.0	EPA 8260C	5-26-17	5-26-17	
Bromochloromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Chloroform	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Benzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Trichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Dibromomethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromodichloromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Chloroethyl Vinyl Ether	ND	6.6	EPA 8260C	5-26-17	5-26-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	5-26-17	5-26-17	
Toluene	ND	1.0	EPA 8260C	5-26-17	5-26-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-26-17	5-26-17	

Project: 15-05986-040

VOLATILES EPA 8260C page 2 of 2

Analyte					Date	Date	
Laboratory ID: 05-299-18 1,1,2 Trichiloroethane ND 0.20 EPA 8260C 5-26-17 5-26-17 Tetrachloroethane ND 0.20 EPA 8260C 5-26-17 5-26-17 1,3-Dichloropropane ND 0.20 EPA 8260C 5-26-17 5-26-17 1,3-Dichloropropane ND 0.20 EPA 8260C 5-26-17 5-26-17 1,3-Dichloropropane ND 0.20 EPA 8260C 5-26-17 5-26-17 1,3-Dichloromothane ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-Dibromoethane ND 0.20 EPA 8260C 5-26-17 5-26-17 1,1-1,2-Tetrachloroethane ND 0.20 EPA 8260C 5-26-17 5-26-17 1,1,1,2-Tetrachloroethane ND 0.20 EPA 8260C 5-26-17 5-26-17 1,1,1,2-Tetrachloropropane ND 0.20 EPA 8260C 5-26-17 5-26-17 1,1,1,2-Tetrachloropropane ND 0.20 EPA 8260C 5-26-17 5-26-17 1,1,2-Tetrachloropropane ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-26-17 1,2-26-17 1,2-26-17 1,2-26-17 1,2-26-17 1,2-26-17 1,2-26-17 1,2-26-17 1,2-26-17 1,2-26-17	Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
1,1,2-Trichloroethane	Client ID:	PP6-W					
Tetrachloroethene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,3-Dichloropropane ND 0.20 EPA 8260C 5-26-17 5-26-17 2-Hexanone ND 2.0 EPA 8260C 5-26-17 5-26-17 Dibromochloromethane ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-Dibromoethane ND 0.20 EPA 8260C 5-26-17 5-26-17 Chlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Ethylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Styrene ND 0.20 EPA 8260C 5-26-17 5-26-17 Styrene ND 0.20 EPA 8260C 5-26-17 5-26-17	Laboratory ID:	05-299-18					
1,3-Dichloropropane ND 0.20 EPA 8260C 5-26-17 5-26-17 2-Hexanone ND 2.0 EPA 8260C 5-26-17 5-26-17 Dibromochtoromethane ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-Dibromoethane ND 0.20 EPA 8260C 5-26-17 5-26-17 1,1,1,2-Tetrachloroethane ND 0.20 EPA 8260C 5-26-17 5-26-17 Ethylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Ethylbenzene ND 0.40 EPA 8260C 5-26-17 5-26-17 Ethylene ND 0.40 EPA 8260C 5-26-17 5-26-17 Ethylene ND 0.40 EPA 8260C 5-26-17 5-26-17 Eyrene ND 0.20 EPA 8260C 5-26-17 5-26-17 Brylene ND 0.20 EPA 8260C 5-26-17 5-26-17 Brylene ND 0.20 EPA 8260C 5-26-17 5-26-17	1,1,2-Trichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Hexanone ND 2.0 EPA 8260C 5-26-17 5-26-17 Dibromochloromethane ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-Dibromochlane ND 0.20 EPA 8260C 5-26-17 5-26-17 Chlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,1,1,2-Tetrachloroethane ND 0.20 EPA 8260C 5-26-17 5-26-17 Ethylbenzene ND 0.40 EPA 8260C 5-26-17 5-26-17 mp-Xylene ND 0.40 EPA 8260C 5-26-17 5-26-17 o-Xylene ND 0.20 EPA 8260C 5-26-17 5-26-17 Styrene ND 0.20 EPA 8260C 5-26-17 5-26-17 Styrene ND 0.20 EPA 8260C 5-26-17 5-26-17 Isopropylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Isopropylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17	Tetrachloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Dibromochloromethane	1,3-Dichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromoethane ND 0.20 EPA 8260C 5-26-17 5-26-17 Chlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,1,1,2-Tetrachloroethane ND 0.20 EPA 8260C 5-26-17 5-26-17 Ethylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 m,p-Xylene ND 0.40 EPA 8260C 5-26-17 5-26-17 o-Xylene ND 0.20 EPA 8260C 5-26-17 5-26-17 Styrene ND 0.20 EPA 8260C 5-26-17 5-26-17 Bromoform ND 1.0 EPA 8260C 5-26-17 5-26-17 Isopropylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Bromobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Isopropylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2,3-Tribolropropane ND 0.20 EPA 8260C 5-26-17 5-26-17 <tr< td=""><td>2-Hexanone</td><td>ND</td><td>2.0</td><td>EPA 8260C</td><td>5-26-17</td><td>5-26-17</td><td></td></tr<>	2-Hexanone	ND	2.0	EPA 8260C	5-26-17	5-26-17	
Chlorobenzene	Dibromochloromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1,1,2-Tetrachloroethane ND 0.20 EPA 8260C 5-26-17 5-26-17 Ethylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 m,p-Xylene ND 0.40 EPA 8260C 5-26-17 5-26-17 O-Xylene ND 0.20 EPA 8260C 5-26-17 5-26-17 Styrene ND 0.20 EPA 8260C 5-26-17 5-26-17 Bromoform ND 1.0 EPA 8260C 5-26-17 5-26-17 Bromobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 J.2.3-Trichloroptoane ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2.3-Trichlorobulane ND 0.20 EPA 8260C 5-26-17 5-26-17	1,2-Dibromoethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Ethylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 m,p-Xylene ND 0.40 EPA 8260C 5-26-17 5-26-17 o-Xylene ND 0.20 EPA 8260C 5-26-17 5-26-17 Styrene ND 0.20 EPA 8260C 5-26-17 5-26-17 Bromoform ND 1.0 EPA 8260C 5-26-17 5-26-17 Bromobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Bromobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,1,2,2-Tetrachloroethane ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichloropropane ND 0.20 EPA 8260C 5-26-17 5-26-17 n-Propylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 4-Chlorotoluene ND 0.20 EPA 8260C 5-26-17 5-26-17 4-Chlorotoluene ND 0.20 EPA 8260C 5-26-17 5-26-17	Chlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
m.p-Xylene ND 0.40 EPA 8260C 5-26-17 5-26-17 c-Xylene ND 0.20 EPA 8260C 5-26-17 5-26-17 Styrene ND 0.20 EPA 8260C 5-26-17 5-26-17 Bromoform ND 1.0 EPA 8260C 5-26-17 5-26-17 Isopropylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Bromobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Bromobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,1,2,2-Tetrachloroethane ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichloropropane ND 0.20 EPA 8260C 5-26-17 5-26-17 -Propylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 -Propylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 -Chlorotoluene ND 0.20 EPA 8260C 5-26-17 5-26-17	1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
o-Xylene ND 0.20 EPA 8260C 5-26-17 5-26-17 Styrene ND 0.20 EPA 8260C 5-26-17 5-26-17 Bromoform ND 1.0 EPA 8260C 5-26-17 5-26-17 Isopropylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Bromobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,1,2,2-Tetrachloroethane ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichloropropane ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichloropropane ND 0.20 EPA 8260C 5-26-17 5-26-17 -P-Popylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 -Chlorotoluene ND 0.20 EPA 8260C 5-26-17 5-26-17 4-Chlorotoluene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,3,5-Trimethylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 </td <td>Ethylbenzene</td> <td>ND</td> <td>0.20</td> <td>EPA 8260C</td> <td>5-26-17</td> <td>5-26-17</td> <td></td>	Ethylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Styrene ND 0.20 EPA 8260C 5-26-17 5-26-17 Bromoform ND 1.0 EPA 8260C 5-26-17 5-26-17 Isopropylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Bromobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 J.2,3-Trichloropropane ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichloropropane ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichloropropane ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichloropropane ND 0.20 EPA 8260C 5-26-17 5-26-17 2-Chlorotoluene ND 0.20 EPA 8260C 5-26-17 5-26-17 4-Chlorotoluene ND 0.20 EPA 8260C 5-26-17 5-26-17 4-Chlorotoluene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,3,5-Trimethylbenzene ND 0.20 EPA 8260C 5-26-17 <	m,p-Xylene	ND	0.40	EPA 8260C	5-26-17	5-26-17	
Bromoform ND 1.0 EPA 8260C 5-26-17 5-26-17 Isopropylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Bromobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,1,2,3-Trichlorobrane ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichlorobropropane ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichlorobropropane ND 0.20 EPA 8260C 5-26-17 5-26-17 2-Chlorotoluene ND 0.20 EPA 8260C 5-26-17 5-26-17 4-Chlorotoluene ND 0.20 EPA 8260C 5-26-17 5-26-17 4-Chlorotoluene ND 0.20 EPA 8260C 5-26-17 5-26-17 4-Chlorotoluene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,3,5-Trimethylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2,4-Trimethylbenzene ND 0.20 EPA 8260C 5-26-17	o-Xylene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Isopropylbenzene	Styrene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromobenzene ND	Bromoform	ND	1.0	EPA 8260C	5-26-17	5-26-17	
1,1,2,2-Tetrachloroethane ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichloropropane ND 0.20 EPA 8260C 5-26-17 5-26-17 n-Propylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 2-Chlorotoluene ND 0.20 EPA 8260C 5-26-17 5-26-17 4-Chlorotoluene ND 0.20 EPA 8260C 5-26-17 5-26-17 4-Chlorotoluene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,3,5-Trimethylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2,4-Trimethylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-A-Trimethylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-A-Trichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,4-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-Dichlorobenzene ND 0.20 EPA 8260C	Isopropylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichloropropane ND 0.20 EPA 8260C 5-26-17 5-26-17 n-Propylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 2-Chlorotoluene ND 0.20 EPA 8260C 5-26-17 5-26-17 4-Chlorotoluene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,3,5-Trimethylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2,4-Trimethylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,3-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,4-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-Dichlorobenzene ND 0.20 EPA 8260C <td>Bromobenzene</td> <td>ND</td> <td>0.20</td> <td>EPA 8260C</td> <td>5-26-17</td> <td>5-26-17</td> <td></td>	Bromobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
n-Propylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 2-Chlorotoluene ND 0.20 EPA 8260C 5-26-17 5-26-17 4-Chlorotoluene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,3,5-Trimethylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2,4-Trimethylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,3-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,3-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,4-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-Dichlorobenzene ND 0.20 EPA 8260C	1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Chlorotoluene ND 0.20 EPA 8260C 5-26-17 5-26-17 4-Chlorotoluene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,3,5-Trimethylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 tert-Butylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2,4-Trimethylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 sec-Butylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,3-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,3-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,4-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 n-Butylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-Dibromo-3-chloropropane ND 0.20 EPA 8260C	1,2,3-Trichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
4-Chlorotoluene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,3,5-Trimethylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 tert-Butylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2,4-Trimethylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 sec-Butylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,3-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,3-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 p-Isopropyltoluene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,4-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 n-Butylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-Brichlorobenzene ND 0.20 EPA 8260C	n-Propylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,3,5-Trimethylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 tert-Butylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2,4-Trimethylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 sec-Butylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,3-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 p-Isopropyltoluene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,4-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 n-Butylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-Dibromo-3-chloropropane ND 1.0 EPA 8260C 5-26-17 5-26-17 1,2,4-Trichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 ND 0.20 EPA 8260C 5-26-17 <td< td=""><td>2-Chlorotoluene</td><td>ND</td><td>0.20</td><td>EPA 8260C</td><td>5-26-17</td><td>5-26-17</td><td></td></td<>	2-Chlorotoluene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
tert-Butylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2,4-Trimethylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 sec-Butylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,3-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 p-Isopropyltoluene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,4-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 n-Butylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-Dibromo-3-chloropropane ND 1.0 EPA 8260C 5-26-17 5-26-17 1,2,4-Trichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Hexachlorobutadiene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichlorobenzene ND 0.20 EPA 8260C	4-Chlorotoluene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trimethylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 sec-Butylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,3-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 p-Isopropyltoluene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,4-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 n-Butylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-Dibromo-3-chloropropane ND 1.0 EPA 8260C 5-26-17 5-26-17 1,2,4-Trichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Naphthalene ND 1.4 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Surrogate: Percent Recovery Control Limits	1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
sec-Butylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,3-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 p-Isopropyltoluene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,4-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 n-Butylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-Dibromo-3-chloropropane ND 1.0 EPA 8260C 5-26-17 5-26-17 1,2,4-Trichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Hexachlorobutadiene ND 0.20 EPA 8260C 5-26-17 5-26-17 Naphthalene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Surrogate: Percent Recovery Control Limits </td <td>tert-Butylbenzene</td> <td>ND</td> <td>0.20</td> <td>EPA 8260C</td> <td>5-26-17</td> <td>5-26-17</td> <td></td>	tert-Butylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,3-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 p-Isopropyltoluene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,4-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 n-Butylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-Dibromo-3-chloropropane ND 1.0 EPA 8260C 5-26-17 5-26-17 1,2,4-Trichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Hexachlorobutadiene ND 0.20 EPA 8260C 5-26-17 5-26-17 Naphthalene ND 1.4 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 97 77-129	1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
p-Isopropyltoluene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,4-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 n-Butylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-Dibromo-3-chloropropane ND 1.0 EPA 8260C 5-26-17 5-26-17 1,2,4-Trichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Hexachlorobutadiene ND 0.20 EPA 8260C 5-26-17 5-26-17 Naphthalene ND 1.4 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 97 77-129 77-129	sec-Butylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,4-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 n-Butylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-Dibromo-3-chloropropane ND 1.0 EPA 8260C 5-26-17 5-26-17 1,2,4-Trichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Hexachlorobutadiene ND 0.20 EPA 8260C 5-26-17 5-26-17 Naphthalene ND 1.4 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 97 77-129	1,3-Dichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 n-Butylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-Dibromo-3-chloropropane ND 1.0 EPA 8260C 5-26-17 5-26-17 1,2,4-Trichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Hexachlorobutadiene ND 0.20 EPA 8260C 5-26-17 5-26-17 Naphthalene ND 1.4 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 97 77-129	p-Isopropyltoluene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
n-Butylbenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 1,2-Dibromo-3-chloropropane ND 1.0 EPA 8260C 5-26-17 5-26-17 1,2,4-Trichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Hexachlorobutadiene ND 0.20 EPA 8260C 5-26-17 5-26-17 Naphthalene ND 1.4 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 97 77-129 Control Limits	1,4-Dichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromo-3-chloropropane ND 1.0 EPA 8260C 5-26-17 5-26-17 1,2,4-Trichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Hexachlorobutadiene ND 0.20 EPA 8260C 5-26-17 5-26-17 Naphthalene ND 1.4 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 97 77-129 77-129	1,2-Dichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Hexachlorobutadiene ND 0.20 EPA 8260C 5-26-17 5-26-17 Naphthalene ND 1.4 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 97 77-129	n-Butylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Hexachlorobutadiene ND 0.20 EPA 8260C 5-26-17 5-26-17 Naphthalene ND 1.4 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 97 77-129	1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	5-26-17	5-26-17	
Naphthalene ND 1.4 EPA 8260C 5-26-17 5-26-17 1,2,3-Trichlorobenzene ND 0.20 EPA 8260C 5-26-17 5-26-17 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 97 77-129			0.20	EPA 8260C	5-26-17	5-26-17	
1,2,3-TrichlorobenzeneND0.20EPA 8260C5-26-175-26-17Surrogate:Percent Recovery Control LimitsDibromofluoromethane9777-129	Hexachlorobutadiene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Surrogate: Percent Recovery Control Limits Dibromofluoromethane 97 77-129	Naphthalene	ND	1.4	EPA 8260C	5-26-17	5-26-17	
Dibromofluoromethane 97 77-129	1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
	Surrogate:	Percent Recovery	Control Limits				
Toluene-d8 98 80-127	Dibromofluoromethane	97	77-129				
	Toluene-d8	98	80-127				

4-Bromofluorobenzene

80-125

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP7-W					
Laboratory ID:	05-299-21					
Dichlorodifluoromethane	ND	0.26	EPA 8260C	5-26-17	5-26-17	
Chloromethane	ND	1.4	EPA 8260C	5-26-17	5-26-17	
Vinyl Chloride	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromomethane	ND	0.33	EPA 8260C	5-26-17	5-26-17	
Chloroethane	ND	1.0	EPA 8260C	5-26-17	5-26-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Acetone	6.1	5.0	EPA 8260C	5-26-17	5-26-17	
Iodomethane	ND	1.7	EPA 8260C	5-26-17	5-26-17	
Carbon Disulfide	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Methylene Chloride	ND	1.0	EPA 8260C	5-26-17	5-26-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Vinyl Acetate	ND	1.0	EPA 8260C	5-26-17	5-26-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Butanone	ND	5.0	EPA 8260C	5-26-17	5-26-17	
Bromochloromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Chloroform	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Benzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Trichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Dibromomethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromodichloromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Chloroethyl Vinyl Ether	ND	6.6	EPA 8260C	5-26-17	5-26-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	5-26-17	5-26-17	
Toluene	5.2	1.0	EPA 8260C	5-26-17	5-26-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-26-17	5-26-17	

Project: 15-05986-040

VOLATILES EPA 8260C page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP7-W					
Laboratory ID:	05-299-21					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Tetrachloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Hexanone	ND	2.0	EPA 8260C	5-26-17	5-26-17	
Dibromochloromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Chlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Ethylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
m,p-Xylene	ND	0.40	EPA 8260C	5-26-17	5-26-17	
o-Xylene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Styrene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromoform	ND	1.0	EPA 8260C	5-26-17	5-26-17	
Isopropylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
n-Propylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
tert-Butylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
sec-Butylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
n-Butylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Naphthalene	ND	1.4	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	96	77-129				
Toluene-d8	83	80-127				

4-Bromofluorobenzene

80-125

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

Analysis	Pagult	DOL	Mathad	Date	Date	Elema
Analyte Client ID:	Result PP8-W	PQL	Method	Prepared	Analyzed	Flags
	_					
Laboratory ID:	05-299-24 ND	0.26	EPA 8260C	5-26-17	5-26-17	
Dichlorodifluoromethane						
Chloromethane	ND	1.4	EPA 8260C	5-26-17	5-26-17	
Vinyl Chloride	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromomethane	ND	0.33	EPA 8260C	5-26-17	5-26-17	
Chloroethane	ND	1.0	EPA 8260C	5-26-17	5-26-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Acetone	5.2	5.0	EPA 8260C	5-26-17	5-26-17	
lodomethane	ND	1.7	EPA 8260C	5-26-17	5-26-17	
Carbon Disulfide	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Methylene Chloride	ND	1.0	EPA 8260C	5-26-17	5-26-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Vinyl Acetate	ND	1.0	EPA 8260C	5-26-17	5-26-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Butanone	ND	5.0	EPA 8260C	5-26-17	5-26-17	
Bromochloromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Chloroform	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Benzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Trichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Dibromomethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromodichloromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Chloroethyl Vinyl Ether	ND	6.6	EPA 8260C	5-26-17	5-26-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	5-26-17	5-26-17	
Toluene	ND	1.0	EPA 8260C	5-26-17	5-26-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-26-17	5-26-17	

Project: 15-05986-040

VOLATILES EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP8-W					
Laboratory ID:	05-299-24					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Tetrachloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Hexanone	ND	2.0	EPA 8260C	5-26-17	5-26-17	
Dibromochloromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Chlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Ethylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
m,p-Xylene	ND	0.40	EPA 8260C	5-26-17	5-26-17	
o-Xylene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Styrene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromoform	ND	1.0	EPA 8260C	5-26-17	5-26-17	
Isopropylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
n-Propylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
tert-Butylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
sec-Butylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
n-Butylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Naphthalene	ND	1.4	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	97	77-129				

Surrogate: Percent Recovery Control Limi
Dibromofluoromethane 97 77-129
Toluene-d8 102 80-127
4-Bromofluorobenzene 103 80-125



Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

Client ID:					Date	Date	
Laboratory ID: 05-299-27	Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Dichlorodifluoromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Chloromethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Vinyl Chloride ND 0.20 EPA 8260C 5-30-17 5-30-17 Bromomethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Chloroethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Trichloroffluoromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 IDichloroethene ND 0.20 EPA 8260C 5-30-17 5-30-17 Acetone 7.2 5.0 EPA 8260C 5-30-17 5-30-17 Iodomethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Acetone 7.2 5.0 EPA 8260C 5-30-17 5-30-17 Iodomethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Methylene Chloride ND 1.0 EPA 8260C 5-30-17 5-30-17	Client ID:	PP9-W					
Chloromethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Vinyl Chloride ND 0.20 EPA 8260C 5-30-17 5-30-17 Bromomethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Chloroethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Trichlorofluoromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Trichloroethene ND 0.20 EPA 8260C 5-30-17 5-30-17 Acetone 7.2 5.0 EPA 8260C 5-30-17 5-30-17 Iodomethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Carbon Disulfide 0.22 0.20 EPA 8260C 5-30-17 5-30-17 Carbon Disulfide 0.22 0.20 EPA 8260C 5-30-17 5-30-17 Vamblylene Chloride ND 1.0 EPA 8260C 5-30-17 5-30-17 Vinyl Acetate ND 0.20 EPA 8260C 5-30-17 5-30-17	Laboratory ID:	05-299-27					
Vinyl Chloride ND 0.20 EPA 8260C 5-30-17 5-30-17 Bromomethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Chloroethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Trichloroftuoromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 1,1-Dichloroethene ND 0.20 EPA 8260C 5-30-17 5-30-17 Acetone 7.2 5.0 EPA 8260C 5-30-17 5-30-17 Acetone 7.2 5.0 EPA 8260C 5-30-17 5-30-17 Iodomethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Methylene Chloride ND 1.0 EPA 8260C 5-30-17 5-30-17 Methyl Ebutyl Ether ND 0.20 EPA 8260C 5-30-17 5-30-17 Vinyl Acetate ND 0.20 EPA 8260C 5-30-17 5-30-17 2,2-Dichloropropane ND 0.20 EPA 8260C 5-30-17 5-30-17	Dichlorodifluoromethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Bromomethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Chloroethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Trichlorofluoromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 1,1-Dichloroethene ND 0.20 EPA 8260C 5-30-17 5-30-17 Acetone 7.2 5.0 EPA 8260C 5-30-17 5-30-17 Acetone ND 1.0 EPA 8260C 5-30-17 5-30-17 Iodomethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Carbon Disulfide 0.22 0.20 EPA 8260C 5-30-17 5-30-17 Wethylene Chloride ND 1.0 EPA 8260C 5-30-17 5-30-17 (trans) 1,2-Dichloroethene ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyl t- Butyl Ether ND 0.20 EPA 8260C 5-30-17 5-30-17 1,1-Dichloroethane ND 0.20 EPA 8260C 5-30-17 5-30-17 </td <td>Chloromethane</td> <td>ND</td> <td>1.0</td> <td>EPA 8260C</td> <td>5-30-17</td> <td>5-30-17</td> <td></td>	Chloromethane	ND	1.0	EPA 8260C	5-30-17	5-30-17	
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Trichlorofluoromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 1,1-Dichloroethene ND 0.20 EPA 8260C 5-30-17 5-30-17 Acetone 7.2 5.0 EPA 8260C 5-30-17 5-30-17 Iodomethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Carbon Disulfide 0.22 0.20 EPA 8260C 5-30-17 5-30-17 Y Methylene Chloride ND 1.0 EPA 8260C 5-30-17 5-30-17 Y Methyl t-Butyl Ether ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyl t-Butyl Ether ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyl t-Butyl Ether ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyl t-Butyl Ether ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyl t-Butyl Ether ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyl t-Butyl Ether ND 0.20 EPA 8260C 5-30-17 5-30-17 5-30-17	Bromomethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,1-Dichloroethene ND 0.20 EPA 8260C 5-30-17 5-30-17 Acetone 7.2 5.0 EPA 8260C 5-30-17 5-30-17 Iodomethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Carbon Disulfide 0.22 0.20 EPA 8260C 5-30-17 5-30-17 Y Methylene Chloride ND 1.0 EPA 8260C 5-30-17 5-30-17 Y Methylene Chloride ND 0.20 EPA 8260C 5-30-17 5-30-17 Y Methyl t-Butyl Ether ND 0.20 EPA 8260C 5-30-17 5-30-17 Y Methyl t-Butyl Ether ND 0.20 EPA 8260C 5-30-17 5-30-17 Y Methyl t-Butyl Ether ND 0.20 EPA 8260C 5-30-17 5-30-17 Y Methyl t-Butyl Ether ND 0.20 EPA 8260C 5-30-17 5-30-17 Y Y Y Y Y Y Y Y Y Y Y	Chloroethane	ND	1.0	EPA 8260C	5-30-17	5-30-17	
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lodomethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Carbon Disulfide 0.22 0.20 EPA 8260C 5-30-17 5-30-17 Y Methylene Chloride ND 1.0 EPA 8260C 5-30-17 5-30-17 Y Methyl Ebutyl Ether ND 0.20 EPA 8260C 5-30-17 5-30-17 T-30-17 1,1-Dichloroethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Vinyl Acetate ND 1.0 EPA 8260C 5-30-17 5-30-17 2,2-Dichloropropane ND 0.20 EPA 8260C 5-30-17 5-30-17 2,2-Dichloroethene ND 0.20 EPA 8260C 5-30-17 5-30-17 2,2-Dichloroethene ND 0.20 EPA 8260C 5-30-17 5-30-17 2-Butanone ND 0.20 EPA 8260C 5-30-17 5-30-17 2-Butanone ND 0.20 EPA 8260C 5-30-17 5-30-17 Pomochloromethane ND 0.20 <t< td=""><td>1,1-Dichloroethene</td><td>ND</td><td>0.20</td><td>EPA 8260C</td><td>5-30-17</td><td>5-30-17</td><td></td></t<>	1,1-Dichloroethene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
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1,1-Dichloropropene ND 0.20 EPA 8260C 5-30-17 5-30-17 Benzene ND 0.20 EPA 8260C 5-30-17 5-30-17 1,2-Dichloroethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Trichloroethene ND 0.20 EPA 8260C 5-30-17 5-30-17 1,2-Dichloropropane ND 0.20 EPA 8260C 5-30-17 5-30-17 Dibromomethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Bromodichloromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 5-30-17 5-30-17 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 5-30-17 5-30-17 Toluene ND 1.0 EPA 8260C 5-30-17 5-30-17	1,1,1-Trichloroethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Benzene ND 0.20 EPA 8260C 5-30-17 5-30-17 1,2-Dichloroethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Trichloroethene ND 0.20 EPA 8260C 5-30-17 5-30-17 1,2-Dichloropropane ND 0.20 EPA 8260C 5-30-17 5-30-17 Dibromomethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Bromodichloromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 5-30-17 5-30-17 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 5-30-17 5-30-17 Toluene ND 1.0 EPA 8260C 5-30-17 5-30-17	Carbon Tetrachloride	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,2-Dichloroethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Trichloroethene ND 0.20 EPA 8260C 5-30-17 5-30-17 1,2-Dichloropropane ND 0.20 EPA 8260C 5-30-17 5-30-17 Dibromomethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Bromodichloromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 5-30-17 5-30-17 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 5-30-17 5-30-17 Toluene ND 1.0 EPA 8260C 5-30-17 5-30-17	1,1-Dichloropropene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Trichloroethene ND 0.20 EPA 8260C 5-30-17 5-30-17 1,2-Dichloropropane ND 0.20 EPA 8260C 5-30-17 5-30-17 Dibromomethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Bromodichloromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 5-30-17 5-30-17 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 5-30-17 5-30-17 Toluene ND 1.0 EPA 8260C 5-30-17 5-30-17	Benzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,2-Dichloropropane ND 0.20 EPA 8260C 5-30-17 5-30-17 Dibromomethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Bromodichloromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 5-30-17 5-30-17 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 5-30-17 5-30-17 Toluene ND 1.0 EPA 8260C 5-30-17 5-30-17	1,2-Dichloroethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Dibromomethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Bromodichloromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 5-30-17 5-30-17 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 5-30-17 5-30-17 Toluene ND 1.0 EPA 8260C 5-30-17 5-30-17	Trichloroethene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Bromodichloromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 5-30-17 5-30-17 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 5-30-17 5-30-17 Toluene ND 1.0 EPA 8260C 5-30-17 5-30-17	1,2-Dichloropropane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 5-30-17 5-30-17 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 5-30-17 5-30-17 Toluene ND 1.0 EPA 8260C 5-30-17 5-30-17	Dibromomethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
(cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 5-30-17 5-30-17 Toluene ND 1.0 EPA 8260C 5-30-17 5-30-17	Bromodichloromethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Methyl Isobutyl Ketone ND 2.0 EPA 8260C 5-30-17 5-30-17 Toluene ND 1.0 EPA 8260C 5-30-17 5-30-17	2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	5-30-17	5-30-17	
Toluene ND 1.0 EPA 8260C 5-30-17 5-30-17	(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Toluene ND 1.0 EPA 8260C 5-30-17 5-30-17	Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	5-30-17	5-30-17	
(trans) 1,3-Dichloropropene ND 0.20 EPA 8260C 5-30-17 5-30-17		ND	1.0	EPA 8260C	5-30-17	5-30-17	
	(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-30-17	5-30-17	

Project: 15-05986-040

VOLATILES EPA 8260C page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP9-W					
Laboratory ID:	05-299-27					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Tetrachloroethene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
2-Hexanone	ND	2.0	EPA 8260C	5-30-17	5-30-17	
Dibromochloromethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Chlorobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Ethylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
m,p-Xylene	ND	0.40	EPA 8260C	5-30-17	5-30-17	
o-Xylene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Styrene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Bromoform	ND	1.0	EPA 8260C	5-30-17	5-30-17	
Isopropylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Bromobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
n-Propylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
tert-Butylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
sec-Butylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
n-Butylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,2-Dibromo-3-chloropropane		1.0	EPA 8260C	5-30-17	5-30-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Naphthalene	ND	1.0	EPA 8260C	5-30-17	5-30-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Surrogate:	Percent Recovery	Control Limits	_			
Dibromofluoromethane	119	77-129				
Toluene-d8	116	80-127				
. 5.50110 40		00 121				

4-Bromofluorobenzene

80-125

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP10-W					
Laboratory ID:	05-299-29					
Dichlorodifluoromethane	ND	0.26	EPA 8260C	5-26-17	5-26-17	
Chloromethane	ND	1.4	EPA 8260C	5-26-17	5-26-17	
Vinyl Chloride	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromomethane	ND	0.33	EPA 8260C	5-26-17	5-26-17	
Chloroethane	ND	1.0	EPA 8260C	5-26-17	5-26-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Acetone	ND	5.0	EPA 8260C	5-26-17	5-26-17	
lodomethane	ND	1.7	EPA 8260C	5-26-17	5-26-17	
Carbon Disulfide	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Methylene Chloride	ND	1.0	EPA 8260C	5-26-17	5-26-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Vinyl Acetate	ND	1.0	EPA 8260C	5-26-17	5-26-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Butanone	ND	5.0	EPA 8260C	5-26-17	5-26-17	
Bromochloromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Chloroform	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Benzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Trichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Dibromomethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromodichloromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Chloroethyl Vinyl Ether	ND	6.6	EPA 8260C	5-26-17	5-26-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	5-26-17	5-26-17	
Toluene	ND	1.0	EPA 8260C	5-26-17	5-26-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-26-17	5-26-17	

Project: 15-05986-040

VOLATILES EPA 8260C page 2 of 2

Analyte				Date	Date	
	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP10-W					
Laboratory ID:	05-299-29					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Tetrachloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Hexanone	ND	2.0	EPA 8260C	5-26-17	5-26-17	
Dibromochloromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Chlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Ethylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
m,p-Xylene	ND	0.40	EPA 8260C	5-26-17	5-26-17	
o-Xylene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Styrene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromoform	ND	1.0	EPA 8260C	5-26-17	5-26-17	
Isopropylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
n-Propylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
tert-Butylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
sec-Butylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
n-Butylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Naphthalene	ND	1.4	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	104	77-129				
Toluene-d8	98	80-127				

4-Bromofluorobenzene

80-125

Project: 15-05986-040

VOLATILES EPA 8260C page 1 of 2

Client ID: Trip Blank Laboratory ID: 05-299-30 Dichlorodifluoromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Chloromethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Vinyl Chloride ND 0.20 EPA 8260C 5-30-17 5-30-17 Bromomethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Chloroethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Trichlorofluoromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Trichlorofluoromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Indomethane ND 5.0 EPA 8260C 5-30-17 5-30-17 Iodomethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Carbon Disulfide ND 0.20 EPA 8260C 5-30-17 5-30-17 Wethylene Chloride ND 1.0 EPA 8260C 5-30-17 5-30-17 <t< th=""><th></th><th></th><th></th><th></th><th>Date</th><th>Date</th><th></th></t<>					Date	Date	
Laboratory ID: 05-299-30 Dichlorodifluoromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Chloromethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Vinyl Chloride ND 0.20 EPA 8260C 5-30-17 5-30-17 Bromomethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Chloroethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Tricklorofluoromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Tricklorofluoromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Acetone ND 5.0 EPA 8260C 5-30-17 5-30-17 Acetone ND 1.0 EPA 8260C 5-30-17 5-30-17 Carbon Disulfide ND 1.0 EPA 8260C 5-30-17 5-30-17 Methylene Chloride ND 1.0 EPA 8260C 5-30-17 5-30-17 Methyle Bulyl Ether ND 0.20	Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Dichlorodiffluoromethane	Client ID:	Trip Blank					
Chloromethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Vinyl Chloride ND 0.20 EPA 8260C 5-30-17 5-30-17 Bromomethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Chloroethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Trichlorofluoromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 T,1-Dichloroethene ND 0.20 EPA 8260C 5-30-17 5-30-17 Acetone ND 5.0 EPA 8260C 5-30-17 5-30-17 Iodomethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Carbon Disulfide ND 1.0 EPA 8260C 5-30-17 5-30-17 Methylene Chloride ND 1.0 EPA 8260C 5-30-17 5-30-17 Methylene Chloride ND 0.20 EPA 8260C 5-30-17 5-30-17 Methylene Chloride ND 0.20 EPA 8260C 5-30-17 5-30-17	Laboratory ID:	05-299-30					
Vinyl Chloride ND 0.20 EPA 8260C 5-30-17 5-30-17 Bromomethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Chloroethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Trichlorofluoromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 1,1-Dichloroethene ND 0.20 EPA 8260C 5-30-17 5-30-17 Acetone ND 5.0 EPA 8260C 5-30-17 5-30-17 Acetone ND 1.0 EPA 8260C 5-30-17 5-30-17 Iodomethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Methylene Chloride ND 1.0 EPA 8260C 5-30-17 5-30-17 Methyl t-Butyl Ether ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyl t-Butyl Ether ND 0.20 EPA 8260C 5-30-17 5-30-17 Vinyl Acetate ND 0.20 EPA 8260C 5-30-17 5-30-17	Dichlorodifluoromethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Bromomethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Chloroethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Trichlorofluoromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Int-Dichloroethene ND 0.20 EPA 8260C 5-30-17 5-30-17 Acetone ND 5.0 EPA 8260C 5-30-17 5-30-17 Iodomethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Carbon Disulfide ND 1.0 EPA 8260C 5-30-17 5-30-17 Methylene Chloride ND 1.0 EPA 8260C 5-30-17 5-30-17 Methyl L-Butyl Ether ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyl L-Butyl Ether ND 0.20 EPA 8260C 5-30-17 5-30-17 1,1-Dichloroethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Vinyl Acetate ND 0.20 EPA 8260C 5-30-17 5-30-17	Chloromethane	ND	1.0	EPA 8260C	5-30-17	5-30-17	
Chloroethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Trichlorofluoromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 1,1-Dichloroethene ND 0.20 EPA 8260C 5-30-17 5-30-17 Acetone ND 5.0 EPA 8260C 5-30-17 5-30-17 Iodomethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Carbon Disulfide ND 1.0 EPA 8260C 5-30-17 5-30-17 Methylene Chloride ND 1.0 EPA 8260C 5-30-17 5-30-17 Methylene Chloride ND 1.0 EPA 8260C 5-30-17 5-30-17 Methylene Chloride ND 0.20 EPA 8260C 5-30-17 5-30-17 Methylene Chloride ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyle Ether ND 0.20 EPA 8260C 5-30-17 5-30-17 1,1-Dichloroethane ND 0.20 EPA 8260C 5-30-17 5-30-17	Vinyl Chloride	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Trichlorofluoromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 1,1-Dichloroethene ND 0.20 EPA 8260C 5-30-17 5-30-17 Acetone ND 5.0 EPA 8260C 5-30-17 5-30-17 Iodomethane ND 1.0 EPA 8260C 5-30-17 5-30-17 Carbon Disulfide ND 0.20 EPA 8260C 5-30-17 5-30-17 Methylene Chloride ND 1.0 EPA 8260C 5-30-17 5-30-17 Methyl EButyl Ether ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyl L-Butyl Ether ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyl L-Butyl Ether ND 0.20 EPA 8260C 5-30-17 5-30-17 Vinyl Acetate ND 0.20 EPA 8260C 5-30-17 5-30-17 2,2-Dichloropropane ND 0.20 EPA 8260C 5-30-17 5-30-17 2,2-Dichloropropane ND 0.20 EPA 8260C 5-30-17	Bromomethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
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(trans) 1,2-Dichloroethene ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyl t-Butyl Ether ND 0.20 EPA 8260C 5-30-17 5-30-17 1,1-Dichloroethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Vinyl Acetate ND 1.0 EPA 8260C 5-30-17 5-30-17 2,2-Dichloropropane ND 0.20 EPA 8260C 5-30-17 5-30-17 (cis) 1,2-Dichloroethene ND 0.20 EPA 8260C 5-30-17 5-30-17 2-Butanone ND 5.0 EPA 8260C 5-30-17 5-30-17 2-Butanone ND 0.20 EPA 8260C 5-30-17 5-30-17 Bromochloromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Chloroform ND 0.20 EPA 8260C 5-30-17 5-30-17 Chloroethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Carbon Tetrachloride ND 0.20 EPA 8260C 5-30-17 <	Carbon Disulfide	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Methyl t-Butyl Ether ND 0.20 EPA 8260C 5-30-17 5-30-17 1,1-Dichloroethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Vinyl Acetate ND 1.0 EPA 8260C 5-30-17 5-30-17 2,2-Dichloropropane ND 0.20 EPA 8260C 5-30-17 5-30-17 (cis) 1,2-Dichloroethene ND 0.20 EPA 8260C 5-30-17 5-30-17 2-Butanone ND 5.0 EPA 8260C 5-30-17 5-30-17 Bromochloromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Chloroform ND 0.20 EPA 8260C 5-30-17 5-30-17 Chloroform ND 0.20 EPA 8260C 5-30-17 5-30-17 Carbon Tetrachloride ND 0.20 EPA 8260C 5-30-17 5-30-17 1,1-Dichloropropene ND 0.20 EPA 8260C 5-30-17 5-30-17 1,2-Dichloroethane ND 0.20 EPA 8260C 5-30-17 <t< td=""><td>Methylene Chloride</td><td>ND</td><td>1.0</td><td>EPA 8260C</td><td>5-30-17</td><td>5-30-17</td><td></td></t<>	Methylene Chloride	ND	1.0	EPA 8260C	5-30-17	5-30-17	
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(cis) 1,2-Dichloroethene ND 0.20 EPA 8260C 5-30-17 5-30-17 2-Butanone ND 5.0 EPA 8260C 5-30-17 5-30-17 Bromochloromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Chloroform ND 0.20 EPA 8260C 5-30-17 5-30-17 1,1,1-Trichloroethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Carbon Tetrachloride ND 0.20 EPA 8260C 5-30-17 5-30-17 Carbon Tetrachloride ND 0.20 EPA 8260C 5-30-17 5-30-17 Carbon Tetrachloride ND 0.20 EPA 8260C 5-30-17 5-30-17 1,1-Dichloropropene ND 0.20 EPA 8260C 5-30-17 5-30-17 Benzene ND 0.20 EPA 8260C 5-30-17 5-30-17 1,2-Dichloroethane ND 0.20 EPA 8260C 5-30-17 5-30-17 1,2-Dichloropropane ND 0.20 EPA 8260C 5-30-17	Vinyl Acetate	ND	1.0	EPA 8260C	5-30-17	5-30-17	
2-Butanone ND 5.0 EPA 8260C 5-30-17 5-30-17 Bromochloromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Chloroform ND 0.20 EPA 8260C 5-30-17 5-30-17 1,1,1-Trichloroethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Carbon Tetrachloride ND 0.20 EPA 8260C 5-30-17 5-30-17 1,1-Dichloropropene ND 0.20 EPA 8260C 5-30-17 5-30-17 Benzene ND 0.20 EPA 8260C 5-30-17 5-30-17 1,2-Dichloroethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Trichloroethene ND 0.20 EPA 8260C 5-30-17 5-30-17 1,2-Dichloropropane ND 0.20 EPA 8260C 5-30-17 5-30-17 Dibromomethane ND 0.20 EPA 8260C 5-30-17 5-30-17 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 5-30-17 5	2,2-Dichloropropane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
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1,1,1-Trichloroethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Carbon Tetrachloride ND 0.20 EPA 8260C 5-30-17 5-30-17 1,1-Dichloropropene ND 0.20 EPA 8260C 5-30-17 5-30-17 Benzene ND 0.20 EPA 8260C 5-30-17 5-30-17 1,2-Dichloroethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Trichloroethene ND 0.20 EPA 8260C 5-30-17 5-30-17 1,2-Dichloropropane ND 0.20 EPA 8260C 5-30-17 5-30-17 1,2-Dichloropropane ND 0.20 EPA 8260C 5-30-17 5-30-17 Dibromomethane ND 0.20 EPA 8260C 5-30-17 5-30-17 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 5-30-17 5-30-17 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyl Isobutyl Ketone ND 1.0 EPA 8260C <t< td=""><td>Bromochloromethane</td><td>ND</td><td>0.20</td><td>EPA 8260C</td><td>5-30-17</td><td>5-30-17</td><td></td></t<>	Bromochloromethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
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Benzene ND 0.20 EPA 8260C 5-30-17 5-30-17 1,2-Dichloroethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Trichloroethene ND 0.20 EPA 8260C 5-30-17 5-30-17 1,2-Dichloropropane ND 0.20 EPA 8260C 5-30-17 5-30-17 1,2-Dichloropropane ND 0.20 EPA 8260C 5-30-17 5-30-17 Dibromomethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Bromodichloromethane ND 1.0 EPA 8260C 5-30-17 5-30-17 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 5-30-17 5-30-17 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 5-30-17 5-30-17 Toluene ND 1.0 EPA 8260C 5-30-17 5-30-17	Carbon Tetrachloride	ND	0.20	EPA 8260C	5-30-17	5-30-17	
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1,2-Dichloropropane ND 0.20 EPA 8260C 5-30-17 5-30-17 Dibromomethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Bromodichloromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 5-30-17 5-30-17 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 5-30-17 5-30-17 Toluene ND 1.0 EPA 8260C 5-30-17 5-30-17	1,2-Dichloroethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Dibromomethane ND 0.20 EPA 8260C 5-30-17 5-30-17 Bromodichloromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 5-30-17 5-30-17 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 5-30-17 5-30-17 Toluene ND 1.0 EPA 8260C 5-30-17 5-30-17	Trichloroethene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Bromodichloromethane ND 0.20 EPA 8260C 5-30-17 5-30-17 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 5-30-17 5-30-17 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 5-30-17 5-30-17 Toluene ND 1.0 EPA 8260C 5-30-17 5-30-17	1,2-Dichloropropane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 5-30-17 5-30-17 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 5-30-17 5-30-17 Toluene ND 1.0 EPA 8260C 5-30-17 5-30-17	Dibromomethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
(cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 5-30-17 5-30-17 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 5-30-17 5-30-17 Toluene ND 1.0 EPA 8260C 5-30-17 5-30-17	Bromodichloromethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Methyl Isobutyl Ketone ND 2.0 EPA 8260C 5-30-17 5-30-17 Toluene ND 1.0 EPA 8260C 5-30-17 5-30-17	2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	5-30-17	5-30-17	
Toluene ND 1.0 EPA 8260C 5-30-17 5-30-17	(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Toluene ND 1.0 EPA 8260C 5-30-17 5-30-17	Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	5-30-17	5-30-17	
(trans) 1,3-Dichloropropene ND 0.20 EPA 8260C 5-30-17 5-30-17	Toluene	ND	1.0	EPA 8260C	5-30-17	5-30-17	
	(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-30-17	5-30-17	

Project: 15-05986-040

VOLATILES EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	Trip Blank					
Laboratory ID:	05-299-30					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Tetrachloroethene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
2-Hexanone	ND	2.0	EPA 8260C	5-30-17	5-30-17	
Dibromochloromethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Chlorobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Ethylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
m,p-Xylene	ND	0.40	EPA 8260C	5-30-17	5-30-17	
o-Xylene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Styrene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Bromoform	ND	1.0	EPA 8260C	5-30-17	5-30-17	
Isopropylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Bromobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
n-Propylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
tert-Butylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
sec-Butylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
n-Butylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,2-Dibromo-3-chloropropane		1.0	EPA 8260C	5-30-17	5-30-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Naphthalene	ND	1.0	EPA 8260C	5-30-17	5-30-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Surrogate:	Percent Recovery	Control Limits	2.7102000	0 00 11	0 00 17	
Dibromofluoromethane	126	77-129				
Dibioinolluololliellialie	120	11-129				

4-Bromofluorobenzene

Toluene-d8

80-127

80-125

118

Project: 15-05986-040

VOLATILES by EPA 8260C METHOD BLANK QUALITY CONTROL

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0526W1					
Dichlorodifluoromethane	ND	0.26	EPA 8260C	5-26-17	5-26-17	
Chloromethane	ND	1.4	EPA 8260C	5-26-17	5-26-17	
Vinyl Chloride	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromomethane	ND	0.33	EPA 8260C	5-26-17	5-26-17	
Chloroethane	ND	1.0	EPA 8260C	5-26-17	5-26-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Acetone	ND	5.0	EPA 8260C	5-26-17	5-26-17	
Iodomethane	ND	1.7	EPA 8260C	5-26-17	5-26-17	
Carbon Disulfide	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Methylene Chloride	ND	1.0	EPA 8260C	5-26-17	5-26-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Vinyl Acetate	ND	1.0	EPA 8260C	5-26-17	5-26-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Butanone	ND	5.0	EPA 8260C	5-26-17	5-26-17	
Bromochloromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Chloroform	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Benzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Trichloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Dibromomethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromodichloromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Chloroethyl Vinyl Ether	ND	6.6	EPA 8260C	5-26-17	5-26-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	5-26-17	5-26-17	
Toluene	ND	1.0	EPA 8260C	5-26-17	5-26-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-26-17	5-26-17	

Project: 15-05986-040

VOLATILES by EPA 8260C METHOD BLANK QUALITY CONTROL

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0526W1					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Tetrachloroethene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Hexanone	ND	2.0	EPA 8260C	5-26-17	5-26-17	
Dibromochloromethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Chlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Ethylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
m,p-Xylene	ND	0.40	EPA 8260C	5-26-17	5-26-17	
o-Xylene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Styrene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromoform	ND	1.0	EPA 8260C	5-26-17	5-26-17	
Isopropylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Bromobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	5-26-17	5-26-17	
n-Propylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
tert-Butylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
sec-Butylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
n-Butylbenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
1,2-Dibromo-3-chloropropane		1.0	EPA 8260C	5-26-17	5-26-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Naphthalene	ND	1.4	EPA 8260C	5-26-17	5-26-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	5-26-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits	, , 52000	0 20 11	0 20 11	
Dibromofly oromothers	ne	77 120				

 Dibromofluoromethane
 96
 77-129

 Toluene-d8
 90
 80-127

 4-Bromofluorobenzene
 102
 80-125



Project: 15-05986-040

VOLATILES by EPA 8260C METHOD BLANK QUALITY CONTROL

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0530W1					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Chloromethane	ND	1.0	EPA 8260C	5-30-17	5-30-17	
Vinyl Chloride	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Bromomethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Chloroethane	ND	1.0	EPA 8260C	5-30-17	5-30-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Acetone	ND	5.0	EPA 8260C	5-30-17	5-30-17	
lodomethane	ND	1.0	EPA 8260C	5-30-17	5-30-17	
Carbon Disulfide	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Methylene Chloride	ND	1.0	EPA 8260C	5-30-17	5-30-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Vinyl Acetate	ND	1.0	EPA 8260C	5-30-17	5-30-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
2-Butanone	ND	5.0	EPA 8260C	5-30-17	5-30-17	
Bromochloromethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Chloroform	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Benzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Trichloroethene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Dibromomethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Bromodichloromethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	5-30-17	5-30-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	5-30-17	5-30-17	
Toluene	ND	1.0	EPA 8260C	5-30-17	5-30-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	5-30-17	5-30-17	

Project: 15-05986-040

VOLATILES by EPA 8260C METHOD BLANK QUALITY CONTROL

page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
7 interior to	rtoouit		mourou		7 <u>7</u>	1 1490
Laboratory ID:	MB0530W1					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Tetrachloroethene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
2-Hexanone	ND	2.0	EPA 8260C	5-30-17	5-30-17	
Dibromochloromethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Chlorobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Ethylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
m,p-Xylene	ND	0.40	EPA 8260C	5-30-17	5-30-17	
o-Xylene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Styrene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Bromoform	ND	1.0	EPA 8260C	5-30-17	5-30-17	
Isopropylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Bromobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	5-30-17	5-30-17	
n-Propylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
tert-Butylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
sec-Butylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
n-Butylbenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	5-30-17	5-30-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Naphthalene	ND	1.0	EPA 8260C	5-30-17	5-30-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	5-30-17	5-30-17	
Surrogate:	Percent Recovery	Control Limits				

Surrogate: Percent Recovery Control Limit
Dibromofluoromethane 120 77-129
Toluene-d8 121 80-127
4-Bromofluorobenzene 89 80-125



Project: 15-05986-040

VOLATILES by EPA 8260C MS/MSD QUALITY CONTROL

					Source	Per	cent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Rec	overy	Limits	RPD	Limit	Flags
MATRIX SPIKES											
Laboratory ID:	05-27	74-07									
	MS	MSD	MS	MSD		MS	MSD				
1,1-Dichloroethene	10.7	10.5	10.0	10.0	ND	107	105	65-119	2	15	
Benzene	10.6	10.7	10.0	10.0	ND	106	107	75-117	1	15	
Trichloroethene	9.44	9.58	10.0	10.0	ND	94	96	66-120	1	15	
Toluene	10.7	10.5	10.0	10.0	ND	107	105	79-120	2	15	
Chlorobenzene	9.15	9.48	10.0	10.0	ND	92	95	76-120	4	15	
Surrogate:											
Dibromofluoromethane						96	97	77-129			
Toluene-d8						93	91	80-127			
4-Bromofluorobenzene						105	106	80-125			

Project: 15-05986-040

VOLATILES by EPA 8260C SB/SBD QUALITY CONTROL

					Per	cent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Reco	overy	Limits	RPD	Limit	Flags
SPIKE BLANKS										
Laboratory ID:	SB05	30W1								
	SB	SBD	SB	SBD	SB	SBD				
1,1-Dichloroethene	9.80	9.93	10.0	10.0	98	99	63-127	1	17	
Benzene	10.7	11.0	10.0	10.0	107	110	76-121	3	12	
Trichloroethene	9.13	9.38	10.0	10.0	91	94	64-120	3	15	
Toluene	10.1	10.6	10.0	10.0	101	106	82-120	5	13	
Chlorobenzene	10.1	10.7	10.0	10.0	101	107	80-120	6	14	
Surrogate:										
Dibromofluoromethane					118	124	77-129			
Toluene-d8					112	113	80-127			
4-Bromofluorobenzene					89	88	80-125			

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

page 1 of 2

Matrix: Soil Units: mg/Kg

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP1-2					
Laboratory ID:	05-299-01					
n-Nitrosodimethylamine	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Pyridine	ND	0.41	EPA 8270D	5-31-17	6-1-17	
Phenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Aniline	ND	0.21	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethyl)ether	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Chlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,3-Dichlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,4-Dichlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Benzyl alcohol	ND	0.21	EPA 8270D	5-31-17	6-1-17	
1,2-Dichlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Methylphenol (o-Cresol)	ND	0.041	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroisopropyl)ether	ND	0.041	EPA 8270D	5-31-17	6-1-17	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.041	EPA 8270D	5-31-17	6-1-17	
n-Nitroso-di-n-propylamine	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Hexachloroethane	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Nitrobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Isophorone	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Nitrophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4-Dimethylphenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethoxy)methane	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4-Dichlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,2,4-Trichlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Naphthalene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
4-Chloroaniline	ND	0.21	EPA 8270D	5-31-17	6-1-17	
Hexachlorobutadiene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
4-Chloro-3-methylphenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Methylnaphthalene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
1-Methylnaphthalene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Hexachlorocyclopentadiene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4,6-Trichlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,3-Dichloroaniline	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4,5-Trichlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Chloronaphthalene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Nitroaniline	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,4-Dinitrobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Dimethylphthalate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,3-Dinitrobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,6-Dinitrotoluene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,2-Dinitrobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Acenaphthylene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
3-Nitroaniline	ND	0.041	EPA 8270D	5-31-17	6-1-17	

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP1-2					
Laboratory ID:	05-299-01					
2,4-Dinitrophenol	ND	0.21	EPA 8270D	5-31-17	6-1-17	
Acenaphthene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
4-Nitrophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4-Dinitrotoluene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Dibenzofuran	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,3,5,6-Tetrachlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,3,4,6-Tetrachlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Diethylphthalate	ND	0.21	EPA 8270D	5-31-17	6-1-17	
4-Chlorophenyl-phenylether	ND	0.041	EPA 8270D	5-31-17	6-1-17	
4-Nitroaniline	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Fluorene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
4,6-Dinitro-2-methylphenol	ND	0.21	EPA 8270D	5-31-17	6-1-17	
n-Nitrosodiphenylamine	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,2-Diphenylhydrazine	ND	0.041	EPA 8270D	5-31-17	6-1-17	
4-Bromophenyl-phenylether	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Hexachlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Pentachlorophenol	ND	0.21	EPA 8270D	5-31-17	6-1-17	
Phenanthrene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Anthracene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Carbazole	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Di-n-butylphthalate	ND	0.21	EPA 8270D	5-31-17	6-1-17	
Fluoranthene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Benzidine	ND	0.41	EPA 8270D	5-31-17	6-1-17	
Pyrene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Butylbenzylphthalate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
bis-2-Ethylhexyladipate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
3,3'-Dichlorobenzidine	ND	0.21	EPA 8270D	5-31-17	6-1-17	
Benzo[a]anthracene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Chrysene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
bis(2-Ethylhexyl)phthalate	0.052	0.041	EPA 8270D	5-31-17	6-1-17	
Di-n-octylphthalate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Benzo[b]fluoranthene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo(j,k)fluoranthene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[a]pyrene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Indeno[1,2,3-cd]pyrene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Dibenz[a,h]anthracene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[g,h,i]perylene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits			0 0 1 11	
2-Fluorophenol	63	18 - 109				
Phenol-d6	66	25 - 111				
Nitrobenzene-d5	66	22 - 113				
2-Fluorobiphenyl	56	30 - 114				
2,4,6-Tribromophenol	59	22 - 116				
Terphenyl-d14	<i>57</i>	33 - 114				
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Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

page 1 of 2

Matrix: Soil Units: mg/Kg

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP1-10					
Laboratory ID:	05-299-03					
n-Nitrosodimethylamine	ND	0.045	EPA 8270D	5-31-17	5-31-17	
Pyridine	ND	0.45	EPA 8270D	5-31-17	5-31-17	
Phenol	ND	0.045	EPA 8270D	5-31-17	5-31-17	
Aniline	ND	0.22	EPA 8270D	5-31-17	5-31-17	
bis(2-Chloroethyl)ether	ND	0.045	EPA 8270D	5-31-17	5-31-17	
2-Chlorophenol	ND	0.045	EPA 8270D	5-31-17	5-31-17	
1,3-Dichlorobenzene	ND	0.045	EPA 8270D	5-31-17	5-31-17	
1,4-Dichlorobenzene	ND	0.045	EPA 8270D	5-31-17	5-31-17	
Benzyl alcohol	ND	0.22	EPA 8270D	5-31-17	5-31-17	
1,2-Dichlorobenzene	ND	0.045	EPA 8270D	5-31-17	5-31-17	
2-Methylphenol (o-Cresol)	ND	0.045	EPA 8270D	5-31-17	5-31-17	
bis(2-Chloroisopropyl)ether	ND	0.045	EPA 8270D	5-31-17	5-31-17	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.045	EPA 8270D	5-31-17	5-31-17	
n-Nitroso-di-n-propylamine	ND	0.045	EPA 8270D	5-31-17	5-31-17	
Hexachloroethane	ND	0.045	EPA 8270D	5-31-17	5-31-17	
Nitrobenzene	ND	0.045	EPA 8270D	5-31-17	5-31-17	
Isophorone	ND	0.045	EPA 8270D	5-31-17	5-31-17	
2-Nitrophenol	ND	0.045	EPA 8270D	5-31-17	5-31-17	
2,4-Dimethylphenol	ND	0.045	EPA 8270D	5-31-17	5-31-17	
bis(2-Chloroethoxy)methane	ND	0.045	EPA 8270D	5-31-17	5-31-17	
2,4-Dichlorophenol	ND	0.045	EPA 8270D	5-31-17	5-31-17	
1,2,4-Trichlorobenzene	ND	0.045	EPA 8270D	5-31-17	5-31-17	
Naphthalene	ND	0.0089	EPA 8270D/SIM	5-31-17	5-31-17	
4-Chloroaniline	ND	0.22	EPA 8270D	5-31-17	5-31-17	
Hexachlorobutadiene	ND	0.045	EPA 8270D	5-31-17	5-31-17	
4-Chloro-3-methylphenol	ND	0.045	EPA 8270D	5-31-17	5-31-17	
2-Methylnaphthalene	ND	0.0089	EPA 8270D/SIM	5-31-17	5-31-17	
1-Methylnaphthalene	ND	0.0089	EPA 8270D/SIM	5-31-17	5-31-17	
Hexachlorocyclopentadiene	ND	0.045	EPA 8270D	5-31-17	5-31-17	
2,4,6-Trichlorophenol	ND	0.045	EPA 8270D	5-31-17	5-31-17	
2,3-Dichloroaniline	ND	0.045	EPA 8270D	5-31-17	5-31-17	
2,4,5-Trichlorophenol	ND	0.045	EPA 8270D	5-31-17	5-31-17	
2-Chloronaphthalene	ND	0.045	EPA 8270D	5-31-17	5-31-17	
2-Nitroaniline	ND	0.045	EPA 8270D	5-31-17	5-31-17	
1,4-Dinitrobenzene	ND	0.045	EPA 8270D	5-31-17	5-31-17	
Dimethylphthalate	ND	0.045	EPA 8270D	5-31-17	5-31-17	
1,3-Dinitrobenzene	ND	0.045	EPA 8270D	5-31-17	5-31-17	
2,6-Dinitrotoluene	ND	0.045	EPA 8270D	5-31-17	5-31-17	
1,2-Dinitrobenzene	ND	0.045	EPA 8270D	5-31-17	5-31-17	
Acenaphthylene	ND	0.0089	EPA 8270D/SIM	5-31-17	5-31-17	
3-Nitroaniline	ND	0.045	EPA 8270D	5-31-17	5-31-17	

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP1-10					
Laboratory ID:	05-299-03					
2,4-Dinitrophenol	ND	0.22	EPA 8270D	5-31-17	5-31-17	
Acenaphthene	ND	0.0089	EPA 8270D/SIM	5-31-17	5-31-17	
4-Nitrophenol	ND	0.045	EPA 8270D	5-31-17	5-31-17	
2,4-Dinitrotoluene	ND	0.045	EPA 8270D	5-31-17	5-31-17	
Dibenzofuran	ND	0.045	EPA 8270D	5-31-17	5-31-17	
2,3,5,6-Tetrachlorophenol	ND	0.045	EPA 8270D	5-31-17	5-31-17	
2,3,4,6-Tetrachlorophenol	ND	0.045	EPA 8270D	5-31-17	5-31-17	
Diethylphthalate	ND	0.22	EPA 8270D	5-31-17	5-31-17	
4-Chlorophenyl-phenylether	ND	0.045	EPA 8270D	5-31-17	5-31-17	
4-Nitroaniline	ND	0.045	EPA 8270D	5-31-17	5-31-17	
Fluorene	ND	0.0089	EPA 8270D/SIM	5-31-17	5-31-17	
4,6-Dinitro-2-methylphenol	ND	0.22	EPA 8270D	5-31-17	5-31-17	
n-Nitrosodiphenylamine	ND	0.045	EPA 8270D	5-31-17	5-31-17	
1,2-Diphenylhydrazine	ND	0.045	EPA 8270D	5-31-17	5-31-17	
4-Bromophenyl-phenylether	ND	0.045	EPA 8270D	5-31-17	5-31-17	
Hexachlorobenzene	ND	0.045	EPA 8270D	5-31-17	5-31-17	
Pentachlorophenol	ND	0.22	EPA 8270D	5-31-17	5-31-17	
Phenanthrene	ND	0.0089	EPA 8270D/SIM	5-31-17	5-31-17	
Anthracene	ND	0.0089	EPA 8270D/SIM	5-31-17	5-31-17	
Carbazole	ND	0.045	EPA 8270D	5-31-17	5-31-17	
Di-n-butylphthalate	ND	0.22	EPA 8270D	5-31-17	5-31-17	
Fluoranthene	ND	0.0089	EPA 8270D/SIM	5-31-17	5-31-17	
Benzidine	ND	0.45	EPA 8270D	5-31-17	5-31-17	
Pyrene	ND	0.0089	EPA 8270D/SIM	5-31-17	5-31-17	
Butylbenzylphthalate	ND	0.045	EPA 8270D	5-31-17	5-31-17	
bis-2-Ethylhexyladipate	ND	0.045	EPA 8270D	5-31-17	5-31-17	
3,3'-Dichlorobenzidine	ND	0.22	EPA 8270D	5-31-17	5-31-17	
Benzo[a]anthracene	ND	0.0089	EPA 8270D/SIM	5-31-17	5-31-17	
Chrysene	ND	0.0089	EPA 8270D/SIM	5-31-17	5-31-17	
bis(2-Ethylhexyl)phthalate	ND	0.045	EPA 8270D	5-31-17	5-31-17	
Di-n-octylphthalate	ND	0.045	EPA 8270D	5-31-17	5-31-17	
Benzo[b]fluoranthene	ND	0.0089	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo(j,k)fluoranthene	ND	0.0089	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[a]pyrene	ND	0.0089	EPA 8270D/SIM	5-31-17	5-31-17	
Indeno[1,2,3-cd]pyrene	ND	0.0089	EPA 8270D/SIM	5-31-17	5-31-17	
Dibenz[a,h]anthracene	ND	0.0089	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[g,h,i]perylene	ND	0.0089	EPA 8270D/SIM	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorophenol	83	18 - 109				
Phenol-d6	81	25 - 111				
Nitrobenzene-d5	87	22 - 113				
2-Fluorobiphenyl	84	30 - 114				
2,4,6-Tribromophenol						
=, :, c ::c: c::.cp::.c:	85	22 - 116				

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

page 1 of 2

Matrix: Soil Units: mg/Kg

Analyte Result PQL Method Prepared Analyzed Flags Client ID: PP2-2.5 Section 10 05-299-04 Section 10 Section 17 6-1-17 n-Nitrosodimethylamine ND 0.044 EPA 8270D 5-31-17 6-1-17 Pyridine ND 0.044 EPA 8270D 5-31-17 6-1-17 Phenol ND 0.044 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.022 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.044 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Benzyl alcohol ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.044 EPA					Date	Date	
Laboratory ID: 05-299-04 n-Nitrosodimethylamine ND 0.044 EPA 8270D 5-31-17 6-1-17 Pyridine ND 0.44 EPA 8270D 5-31-17 6-1-17 Phenol ND 0.044 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.022 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.044 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 4,4-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Benzyl alcohol ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND <th< th=""><th>Analyte</th><th>Result</th><th>PQL</th><th>Method</th><th>Prepared</th><th>Analyzed</th><th>Flags</th></th<>	Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
n-Nitrosodimethylamine ND 0.044 EPA 8270D 5-31-17 6-1-17 Pyridine ND 0.44 EPA 8270D 5-31-17 6-1-17 Phenol ND 0.044 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.044 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.044 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.044 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17	Client ID:	PP2-2.5					
Pyridine ND 0.44 EPA 8270D 5-31-17 6-1-17 Phenol ND 0.044 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.022 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-	Laboratory ID:	05-299-04					
Pyridine ND 0.44 EPA 8270D 5-31-17 6-1-17 Phenol ND 0.044 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.022 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-		ND	0.044	EPA 8270D	5-31-17	6-1-17	
Aniline ND 0.22 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Benzyl alcohol ND 0.22 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (m,p-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 N-Aughtylphenol (m,p-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 N-Bexachloroethane ND 0.044 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.044 EPA 8270D		ND	0.44	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethyl)ether ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Benzyl alcohol ND 0.22 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroisopropyl)ether ND 0.044 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 n-Nitroso-di-n-propylamine ND 0.044 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.044 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.044 EPA	-	ND	0.044	EPA 8270D	5-31-17	6-1-17	
2-Chlorophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Benzyl alcohol ND 0.22 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroisopropyl)ether ND 0.044 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 n-Nitroso-di-n-propylamine ND 0.044 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.044 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.044 EPA 8270D	Aniline	ND	0.22	EPA 8270D	5-31-17	6-1-17	
1,3-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Benzyl alcohol ND 0.22 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroisopropyl)ether ND 0.044 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 n-Nitroso-di-n-propylamine ND 0.044 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.044 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Nitrophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 2,4-Dimethylphenol ND 0.044 EPA 8270	bis(2-Chloroethyl)ether	ND	0.044	EPA 8270D	5-31-17	6-1-17	
1,4-DichlorobenzeneND0.044EPA 8270D5-31-176-1-17Benzyl alcoholND0.22EPA 8270D5-31-176-1-171,2-DichlorobenzeneND0.044EPA 8270D5-31-176-1-172-Methylphenol (o-Cresol)ND0.044EPA 8270D5-31-176-1-17bis(2-Chloroisopropyl)etherND0.044EPA 8270D5-31-176-1-17(3+4)-Methylphenol (m,p-Cresol)ND0.044EPA 8270D5-31-176-1-17n-Nitroso-di-n-propylamineND0.044EPA 8270D5-31-176-1-17HexachloroethaneND0.044EPA 8270D5-31-176-1-17NitrobenzeneND0.044EPA 8270D5-31-176-1-17IsophoroneND0.044EPA 8270D5-31-176-1-172-NitrophenolND0.044EPA 8270D5-31-176-1-172,4-DimethylphenolND0.044EPA 8270D5-31-176-1-17bis(2-Chloroethoxy)methaneND0.044EPA 8270D5-31-176-1-17	2-Chlorophenol	ND	0.044	EPA 8270D	5-31-17	6-1-17	
Benzyl alcohol ND 0.22 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroisopropyl)ether ND 0.044 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 n-Nitroso-di-n-propylamine ND 0.044 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.044 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Nitrophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 2,4-Dimethylphenol ND 0.044 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethoxy)methane ND 0.044 EPA 8270D<	1,3-Dichlorobenzene	ND	0.044	EPA 8270D	5-31-17	6-1-17	
1,2-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroisopropyl)ether ND 0.044 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 n-Nitroso-di-n-propylamine ND 0.044 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.044 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Nitrophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 2,4-Dimethylphenol ND 0.044 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethoxy)methane ND 0.044 EPA 8270D 5-31-17 6-1-17	1,4-Dichlorobenzene	ND	0.044	EPA 8270D	5-31-17	6-1-17	
1,2-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroisopropyl)ether ND 0.044 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 n-Nitroso-di-n-propylamine ND 0.044 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.044 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Nitrophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 2,4-Dimethylphenol ND 0.044 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethoxy)methane ND 0.044 EPA 8270D 5-31-17 6-1-17	Benzyl alcohol	ND	0.22	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroisopropyl)ether ND 0.044 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 n-Nitroso-di-n-propylamine ND 0.044 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.044 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Nitrophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 2,4-Dimethylphenol ND 0.044 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethoxy)methane ND 0.044 EPA 8270D 5-31-17 6-1-17		ND	0.044	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroisopropyl)ether ND 0.044 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 n-Nitroso-di-n-propylamine ND 0.044 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.044 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Nitrophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 2,4-Dimethylphenol ND 0.044 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethoxy)methane ND 0.044 EPA 8270D 5-31-17 6-1-17	2-Methylphenol (o-Cresol)	ND	0.044	EPA 8270D	5-31-17	6-1-17	
n-Nitroso-di-n-propylamine ND 0.044 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.044 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Nitrophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 2,4-Dimethylphenol ND 0.044 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethoxy)methane ND 0.044 EPA 8270D 5-31-17 6-1-17		ND	0.044	EPA 8270D	5-31-17	6-1-17	
n-Nitroso-di-n-propylamine ND 0.044 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.044 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Nitrophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 2,4-Dimethylphenol ND 0.044 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethoxy)methane ND 0.044 EPA 8270D 5-31-17 6-1-17	(3+4)-Methylphenol (m,p-Cresol)	ND	0.044	EPA 8270D	5-31-17	6-1-17	
Nitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Nitrophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 2,4-Dimethylphenol ND 0.044 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethoxy)methane ND 0.044 EPA 8270D 5-31-17 6-1-17		ND	0.044	EPA 8270D	5-31-17	6-1-17	
Isophorone ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Nitrophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 2,4-Dimethylphenol ND 0.044 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethoxy)methane ND 0.044 EPA 8270D 5-31-17 6-1-17	Hexachloroethane	ND	0.044	EPA 8270D	5-31-17	6-1-17	
2-Nitrophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 2,4-Dimethylphenol ND 0.044 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethoxy)methane ND 0.044 EPA 8270D 5-31-17 6-1-17	Nitrobenzene	ND	0.044	EPA 8270D	5-31-17	6-1-17	
2-Nitrophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 2,4-Dimethylphenol ND 0.044 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethoxy)methane ND 0.044 EPA 8270D 5-31-17 6-1-17	Isophorone	ND	0.044	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethoxy)methane ND 0.044 EPA 8270D 5-31-17 6-1-17		ND	0.044	EPA 8270D	5-31-17	6-1-17	
	2,4-Dimethylphenol	ND	0.044	EPA 8270D	5-31-17	6-1-17	
•	bis(2-Chloroethoxy)methane	ND	0.044	EPA 8270D	5-31-17	6-1-17	
2,4-Dichlorophenol ND 0.044 EPA 8270D 5-31-17 6-1-17	2,4-Dichlorophenol	ND	0.044	EPA 8270D	5-31-17	6-1-17	
1,2,4-Trichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17	1,2,4-Trichlorobenzene	ND	0.044	EPA 8270D	5-31-17	6-1-17	
Naphthalene ND 0.0088 EPA 8270D/SIM 5-31-17 5-31-17	Naphthalene	ND	0.0088	EPA 8270D/SIM	5-31-17	5-31-17	
4-Chloroaniline ND 0.22 EPA 8270D 5-31-17 6-1-17	4-Chloroaniline	ND	0.22	EPA 8270D	5-31-17	6-1-17	
Hexachlorobutadiene ND 0.044 EPA 8270D 5-31-17 6-1-17	Hexachlorobutadiene	ND	0.044	EPA 8270D	5-31-17	6-1-17	
4-Chloro-3-methylphenol ND 0.044 EPA 8270D 5-31-17 6-1-17	4-Chloro-3-methylphenol	ND	0.044	EPA 8270D	5-31-17	6-1-17	
2-Methylnaphthalene ND 0.0088 EPA 8270D/SIM 5-31-17 5-31-17	2-Methylnaphthalene	ND	0.0088	EPA 8270D/SIM	5-31-17	5-31-17	
1-Methylnaphthalene ND 0.0088 EPA 8270D/SIM 5-31-17 5-31-17	1-Methylnaphthalene	ND	0.0088	EPA 8270D/SIM	5-31-17	5-31-17	
Hexachlorocyclopentadiene ND 0.044 EPA 8270D 5-31-17 6-1-17		ND	0.044	EPA 8270D	5-31-17	6-1-17	
2,4,6-Trichlorophenol ND 0.044 EPA 8270D 5-31-17 6-1-17	2,4,6-Trichlorophenol	ND	0.044	EPA 8270D	5-31-17	6-1-17	
2,3-Dichloroaniline ND 0.044 EPA 8270D 5-31-17 6-1-17	2,3-Dichloroaniline	ND	0.044	EPA 8270D	5-31-17	6-1-17	
2,4,5-Trichlorophenol ND 0.044 EPA 8270D 5-31-17 6-1-17	2,4,5-Trichlorophenol	ND	0.044	EPA 8270D	5-31-17	6-1-17	
2-Chloronaphthalene ND 0.044 EPA 8270D 5-31-17 6-1-17	2-Chloronaphthalene	ND	0.044	EPA 8270D	5-31-17	6-1-17	
2-Nitroaniline ND 0.044 EPA 8270D 5-31-17 6-1-17	2-Nitroaniline	ND	0.044	EPA 8270D	5-31-17	6-1-17	
1,4-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17	1,4-Dinitrobenzene	ND	0.044	EPA 8270D	5-31-17	6-1-17	
Dimethylphthalate ND 0.044 EPA 8270D 5-31-17 6-1-17	Dimethylphthalate	ND	0.044	EPA 8270D	5-31-17	6-1-17	
1,3-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17	1,3-Dinitrobenzene	ND	0.044	EPA 8270D	5-31-17	6-1-17	
2,6-Dinitrotoluene ND 0.044 EPA 8270D 5-31-17 6-1-17	2,6-Dinitrotoluene	ND	0.044	EPA 8270D	5-31-17	6-1-17	
1,2-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17	1,2-Dinitrobenzene	ND	0.044	EPA 8270D	5-31-17	6-1-17	
Acenaphthylene ND 0.0088 EPA 8270D/SIM 5-31-17 5-31-17	Acenaphthylene	ND	0.0088	EPA 8270D/SIM	5-31-17		
3-Nitroaniline ND 0.044 EPA 8270D 5-31-17 6-1-17	3-Nitroaniline	ND	0.044	EPA 8270D	5-31-17	6-1-17	

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP2-2.5	. ~-	ou		7 iii ii i j	ugo
Laboratory ID:	05-299-04					
2,4-Dinitrophenol	ND	0.22	EPA 8270D	5-31-17	6-1-17	
Acenaphthene	ND	0.0088	EPA 8270D/SIM	5-31-17	5-31-17	
4-Nitrophenol	ND	0.044	EPA 8270D	5-31-17	6-1-17	
2,4-Dinitrotoluene	ND	0.044	EPA 8270D	5-31-17	6-1-17	
Dibenzofuran	ND	0.044	EPA 8270D	5-31-17	6-1-17	
2,3,5,6-Tetrachlorophenol	ND	0.044	EPA 8270D	5-31-17	6-1-17	
2,3,4,6-Tetrachlorophenol	ND	0.044	EPA 8270D	5-31-17	6-1-17	
Diethylphthalate	ND	0.22	EPA 8270D	5-31-17	6-1-17	
4-Chlorophenyl-phenylether	ND	0.044	EPA 8270D	5-31-17	6-1-17	
4-Nitroaniline	ND	0.044	EPA 8270D	5-31-17	6-1-17	
Fluorene	ND	0.0088	EPA 8270D/SIM	5-31-17	5-31-17	
4,6-Dinitro-2-methylphenol	ND	0.22	EPA 8270D	5-31-17	6-1-17	
n-Nitrosodiphenylamine	ND	0.044	EPA 8270D	5-31-17	6-1-17	
1,2-Diphenylhydrazine	ND	0.044	EPA 8270D	5-31-17	6-1-17	
4-Bromophenyl-phenylether	ND	0.044	EPA 8270D	5-31-17	6-1-17	
Hexachlorobenzene	ND	0.044	EPA 8270D	5-31-17	6-1-17	
Pentachlorophenol	ND	0.22	EPA 8270D	5-31-17	6-1-17	
Phenanthrene	ND	0.0088	EPA 8270D/SIM	5-31-17	5-31-17	
Anthracene	ND	0.0088	EPA 8270D/SIM	5-31-17	5-31-17	
Carbazole	ND	0.044	EPA 8270D	5-31-17	6-1-17	
Di-n-butylphthalate	ND	0.22	EPA 8270D	5-31-17	6-1-17	
Fluoranthene	ND	0.0088	EPA 8270D/SIM	5-31-17	5-31-17	
Benzidine	ND	0.44	EPA 8270D	5-31-17	6-1-17	
Pyrene	ND	0.0088	EPA 8270D/SIM	5-31-17	5-31-17	
Butylbenzylphthalate	ND	0.044	EPA 8270D	5-31-17	6-1-17	
bis-2-Ethylhexyladipate	ND	0.044	EPA 8270D	5-31-17	6-1-17	
3,3'-Dichlorobenzidine	ND	0.22	EPA 8270D	5-31-17	6-1-17	
Benzo[a]anthracene	ND	0.0088	EPA 8270D/SIM	5-31-17	5-31-17	
Chrysene	ND	0.0088	EPA 8270D/SIM	5-31-17	5-31-17	
bis(2-Ethylhexyl)phthalate	ND	0.044	EPA 8270D	5-31-17	6-1-17	
Di-n-octylphthalate	ND	0.044	EPA 8270D	5-31-17	6-1-17	
Benzo[b]fluoranthene	ND	0.0088	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo(j,k)fluoranthene	ND	0.0088	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[a]pyrene	ND	0.0088	EPA 8270D/SIM	5-31-17	5-31-17	
Indeno[1,2,3-cd]pyrene	ND	0.0088	EPA 8270D/SIM	5-31-17	5-31-17	
Dibenz[a,h]anthracene	ND	0.0088	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[g,h,i]perylene	ND	0.0088	EPA 8270D/SIM	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorophenol	86	18 - 109				
Phenol-d6	86	25 - 111				
Nitrobenzene-d5	85	22 - 113				
2-Fluorobiphenyl	69	30 - 114				
2,4,6-Tribromophenol	73	22 - 116				
Terphenyl-d14	66	33 - 114				

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

page 1 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP2-10			•	-	
Laboratory ID:	05-299-05					
n-Nitrosodimethylamine	ND	0.048	EPA 8270D	5-31-17	6-1-17	
Pyridine	ND	0.48	EPA 8270D	5-31-17	6-1-17	
Phenol	ND	0.048	EPA 8270D	5-31-17	6-1-17	
Aniline	ND	0.24	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethyl)ether	ND	0.048	EPA 8270D	5-31-17	6-1-17	
2-Chlorophenol	ND	0.048	EPA 8270D	5-31-17	6-1-17	
1,3-Dichlorobenzene	ND	0.048	EPA 8270D	5-31-17	6-1-17	
1,4-Dichlorobenzene	ND	0.048	EPA 8270D	5-31-17	6-1-17	
Benzyl alcohol	ND	0.24	EPA 8270D	5-31-17	6-1-17	
1,2-Dichlorobenzene	ND	0.048	EPA 8270D	5-31-17	6-1-17	
2-Methylphenol (o-Cresol)	ND	0.048	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroisopropyl)ether	ND	0.048	EPA 8270D	5-31-17	6-1-17	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.048	EPA 8270D	5-31-17	6-1-17	
n-Nitroso-di-n-propylamine	ND	0.048	EPA 8270D	5-31-17	6-1-17	
Hexachloroethane	ND	0.048	EPA 8270D	5-31-17	6-1-17	
Nitrobenzene	ND	0.048	EPA 8270D	5-31-17	6-1-17	
Isophorone	ND	0.048	EPA 8270D	5-31-17	6-1-17	
2-Nitrophenol	ND	0.048	EPA 8270D	5-31-17	6-1-17	
2,4-Dimethylphenol	ND	0.048	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethoxy)methane	ND	0.048	EPA 8270D	5-31-17	6-1-17	
2,4-Dichlorophenol	ND	0.048	EPA 8270D	5-31-17	6-1-17	
1,2,4-Trichlorobenzene	ND	0.048	EPA 8270D	5-31-17	6-1-17	
Naphthalene	ND	0.0097	EPA 8270D/SIM	5-31-17	5-31-17	
4-Chloroaniline	ND	0.24	EPA 8270D	5-31-17	6-1-17	
Hexachlorobutadiene	ND	0.048	EPA 8270D	5-31-17	6-1-17	
4-Chloro-3-methylphenol	ND	0.048	EPA 8270D	5-31-17	6-1-17	
2-Methylnaphthalene	ND	0.0097	EPA 8270D/SIM	5-31-17	5-31-17	
1-Methylnaphthalene	ND	0.0097	EPA 8270D/SIM	5-31-17	5-31-17	
Hexachlorocyclopentadiene	ND	0.048	EPA 8270D	5-31-17	6-1-17	
2,4,6-Trichlorophenol	ND	0.048	EPA 8270D	5-31-17	6-1-17	
2,3-Dichloroaniline	ND	0.048	EPA 8270D	5-31-17	6-1-17	
2,4,5-Trichlorophenol	ND	0.048	EPA 8270D	5-31-17	6-1-17	
2-Chloronaphthalene	ND	0.048	EPA 8270D	5-31-17	6-1-17	
2-Nitroaniline	ND	0.048	EPA 8270D	5-31-17	6-1-17	
1,4-Dinitrobenzene	ND	0.048	EPA 8270D	5-31-17	6-1-17	
Dimethylphthalate	ND	0.048	EPA 8270D	5-31-17	6-1-17	
1,3-Dinitrobenzene	ND	0.048	EPA 8270D	5-31-17	6-1-17	
2,6-Dinitrotoluene	ND	0.048	EPA 8270D	5-31-17	6-1-17	
1,2-Dinitrobenzene	ND	0.048	EPA 8270D	5-31-17	6-1-17	
Acenaphthylene	ND	0.0097	EPA 8270D/SIM	5-31-17	5-31-17	
3-Nitroaniline	ND	0.048	EPA 8270D	5-31-17	6-1-17	

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP2-10					
Laboratory ID:	05-299-05					
2,4-Dinitrophenol	ND	0.24	EPA 8270D	5-31-17	6-1-17	
Acenaphthene	ND	0.0097	EPA 8270D/SIM	5-31-17	5-31-17	
4-Nitrophenol	ND	0.048	EPA 8270D	5-31-17	6-1-17	
2,4-Dinitrotoluene	ND	0.048	EPA 8270D	5-31-17	6-1-17	
Dibenzofuran	ND	0.048	EPA 8270D	5-31-17	6-1-17	
2,3,5,6-Tetrachlorophenol	ND	0.048	EPA 8270D	5-31-17	6-1-17	
2,3,4,6-Tetrachlorophenol	ND	0.048	EPA 8270D	5-31-17	6-1-17	
Diethylphthalate	ND	0.24	EPA 8270D	5-31-17	6-1-17	
4-Chlorophenyl-phenylether	ND	0.048	EPA 8270D	5-31-17	6-1-17	
4-Nitroaniline	ND	0.048	EPA 8270D	5-31-17	6-1-17	
Fluorene	ND	0.0097	EPA 8270D/SIM	5-31-17	5-31-17	
4,6-Dinitro-2-methylphenol	ND	0.24	EPA 8270D	5-31-17	6-1-17	
n-Nitrosodiphenylamine	ND	0.048	EPA 8270D	5-31-17	6-1-17	
1,2-Diphenylhydrazine	ND	0.048	EPA 8270D	5-31-17	6-1-17	
4-Bromophenyl-phenylether	ND	0.048	EPA 8270D	5-31-17	6-1-17	
Hexachlorobenzene	ND	0.048	EPA 8270D	5-31-17	6-1-17	
Pentachlorophenol	ND	0.24	EPA 8270D	5-31-17	6-1-17	
Phenanthrene	ND	0.0097	EPA 8270D/SIM	5-31-17	5-31-17	
Anthracene	ND	0.0097	EPA 8270D/SIM	5-31-17	5-31-17	
Carbazole	ND	0.048	EPA 8270D	5-31-17	6-1-17	
Di-n-butylphthalate	ND	0.24	EPA 8270D	5-31-17	6-1-17	
Fluoranthene	ND	0.0097	EPA 8270D/SIM	5-31-17	5-31-17	
Benzidine	ND	0.48	EPA 8270D	5-31-17	6-1-17	
Pyrene	ND	0.0097	EPA 8270D/SIM	5-31-17	5-31-17	
Butylbenzylphthalate	ND	0.048	EPA 8270D	5-31-17	6-1-17	
bis-2-Ethylhexyladipate	ND	0.048	EPA 8270D	5-31-17	6-1-17	
3,3'-Dichlorobenzidine	ND	0.24	EPA 8270D	5-31-17	6-1-17	
Benzo[a]anthracene	ND	0.0097	EPA 8270D/SIM	5-31-17	5-31-17	
Chrysene	ND	0.0097	EPA 8270D/SIM	5-31-17	5-31-17	
bis(2-Ethylhexyl)phthalate	ND	0.048	EPA 8270D	5-31-17	6-1-17	
Di-n-octylphthalate	ND	0.048	EPA 8270D	5-31-17	6-1-17	
Benzo[b]fluoranthene	ND	0.0097	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo(j,k)fluoranthene	ND	0.0097	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[a]pyrene	ND	0.0097	EPA 8270D/SIM	5-31-17	5-31-17	
Indeno[1,2,3-cd]pyrene	ND	0.0097	EPA 8270D/SIM	5-31-17	5-31-17	
Dibenz[a,h]anthracene	ND	0.0097	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[g,h,i]perylene	ND	0.0097	EPA 8270D/SIM	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorophenol	75	18 - 109				
Phenol-d6	72	25 - 111				
Nitrobenzene-d5	73	22 - 113				
2-Fluorobiphenyl	61	30 - 114				
2,4,6-Tribromophenol	61	22 - 116				
Terphenyl-d14	62	33 - 114				
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Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

page 1 of 2

Analyte Result PQL Method Prepared Analyzed Flags Client ID: PP3-2.5 Laboratory ID: 05-299-07					Date	Date	
Laboratory ID: 05-299-07 n-Nitrosodimethylamine ND 0.042 EPA 8270D 5-31-17 6-1-17 Pyridine ND 0.42 EPA 8270D 5-31-17 6-1-17 Phenol ND 0.042 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.042 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.042 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Chlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND	Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
n-Nitrosodimethylamine ND 0.042 EPA 8270D 5-31-17 6-1-17 Pyridine ND 0.42 EPA 8270D 5-31-17 6-1-17 Phenol ND 0.042 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.042 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.042 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.042 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17	Client ID:	PP3-2.5					
Pyridine ND 0.42 EPA 8270D 5-31-17 6-1-17 Phenol ND 0.042 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.21 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (m,p-Cresol) ND 0.042 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.042 EPA 8270D 5-31	Laboratory ID:	05-299-07					
Pyridine ND 0.42 EPA 8270D 5-31-17 6-1-17 Phenol ND 0.042 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.21 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (m,p-Cresol) ND 0.042 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.042 EPA 8270D 5-31		ND	0.042	EPA 8270D	5-31-17	6-1-17	
Aniline ND 0.21 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 1,2-Dich		ND	0.42	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethyl)ether ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 Benzyl alcohol ND 0.21 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.042 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroisopropyl)ether ND 0.042 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.042 EPA 8270D 5-31-17 6-1-17 n-Nitroso-di-n-propylamine ND 0.042 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.042 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.042 EPA	Phenol	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2-Chlorophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 Benzyl alcohol ND 0.21 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.042 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroisopropyl)ether ND 0.042 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.042 EPA 8270D 5-31-17 6-1-17 n-Nitroso-di-n-propylamine ND 0.042 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.042 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Nitrophenol ND 0.042 EPA 8270D <td>Aniline</td> <td>ND</td> <td>0.21</td> <td>EPA 8270D</td> <td>5-31-17</td> <td>6-1-17</td> <td></td>	Aniline	ND	0.21	EPA 8270D	5-31-17	6-1-17	
1,3-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 Benzyl alcohol ND 0.21 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.042 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroisopropyl)ether ND 0.042 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.042 EPA 8270D 5-31-17 6-1-17 n-Nitroso-di-n-propylamine ND 0.042 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.042 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Nitrophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Nitrophenol ND 0.042 EPA 8270D <td>bis(2-Chloroethyl)ether</td> <td>ND</td> <td>0.042</td> <td>EPA 8270D</td> <td>5-31-17</td> <td>6-1-17</td> <td></td>	bis(2-Chloroethyl)ether	ND	0.042	EPA 8270D	5-31-17	6-1-17	
1,3-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 Benzyl alcohol ND 0.21 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.042 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroisopropyl)ether ND 0.042 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.042 EPA 8270D 5-31-17 6-1-17 n-Nitroso-di-n-propylamine ND 0.042 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.042 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Nitrophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Nitrophenol ND 0.042 EPA 8270D <td>2-Chlorophenol</td> <td>ND</td> <td>0.042</td> <td>EPA 8270D</td> <td>5-31-17</td> <td>6-1-17</td> <td></td>	2-Chlorophenol	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Benzyl alcohol ND 0.21 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.042 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroisopropyl)ether ND 0.042 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.042 EPA 8270D 5-31-17 6-1-17 n-Nitroso-di-n-propylamine ND 0.042 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.042 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Nitrophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 2,4-Dimethylphenol ND 0.042 EPA 8270D 5-31-17 6-1-17	1,3-Dichlorobenzene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
1,2-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.042 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroisopropyl)ether ND 0.042 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.042 EPA 8270D 5-31-17 6-1-17 n-Nitroso-di-n-propylamine ND 0.042 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.042 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Nitrophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 2,4-Dimethylphenol ND 0.042 EPA 8270D 5-31-17 6-1-17	1,4-Dichlorobenzene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
1,2-Dichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.042 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroisopropyl)ether ND 0.042 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.042 EPA 8270D 5-31-17 6-1-17 n-Nitroso-di-n-propylamine ND 0.042 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.042 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Nitrophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 2,4-Dimethylphenol ND 0.042 EPA 8270D 5-31-17 6-1-17	Benzyl alcohol	ND	0.21	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroisopropyl)ether ND 0.042 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.042 EPA 8270D 5-31-17 6-1-17 n-Nitroso-di-n-propylamine ND 0.042 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.042 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Nitrophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 2,4-Dimethylphenol ND 0.042 EPA 8270D 5-31-17 6-1-17		ND	0.042	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroisopropyl)ether ND 0.042 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.042 EPA 8270D 5-31-17 6-1-17 n-Nitroso-di-n-propylamine ND 0.042 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.042 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Nitrophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 2,4-Dimethylphenol ND 0.042 EPA 8270D 5-31-17 6-1-17	2-Methylphenol (o-Cresol)	ND	0.042	EPA 8270D	5-31-17	6-1-17	
n-Nitroso-di-n-propylamine ND 0.042 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.042 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Nitrophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 2,4-Dimethylphenol ND 0.042 EPA 8270D 5-31-17 6-1-17		ND	0.042	EPA 8270D	5-31-17	6-1-17	
n-Nitroso-di-n-propylamine ND 0.042 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.042 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Nitrophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 2,4-Dimethylphenol ND 0.042 EPA 8270D 5-31-17 6-1-17	(3+4)-Methylphenol (m,p-Cresol)	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Nitrobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Nitrophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 2,4-Dimethylphenol ND 0.042 EPA 8270D 5-31-17 6-1-17		ND	0.042	EPA 8270D	5-31-17	6-1-17	
Isophorone ND 0.042 EPA 8270D 5-31-17 6-1-17 2-Nitrophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 2,4-Dimethylphenol ND 0.042 EPA 8270D 5-31-17 6-1-17	Hexachloroethane	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Value ND 0.042 EPA 8270D 5-31-17 6-1-17 2,4-Dimethylphenol ND 0.042 EPA 8270D 5-31-17 6-1-17	Nitrobenzene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2-Nitrophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 2,4-Dimethylphenol ND 0.042 EPA 8270D 5-31-17 6-1-17	Isophorone	ND	0.042	EPA 8270D	5-31-17	6-1-17	
		ND	0.042	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethoxy)methane ND 0.042 EPA 8270D 5-31-17 6-1-17	2,4-Dimethylphenol	ND	0.042	EPA 8270D	5-31-17	6-1-17	
	bis(2-Chloroethoxy)methane	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2,4-Dichlorophenol ND 0.042 EPA 8270D 5-31-17 6-1-17		ND	0.042	EPA 8270D	5-31-17	6-1-17	
1,2,4-Trichlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17	1,2,4-Trichlorobenzene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Naphthalene ND 0.0084 EPA 8270D/SIM 5-31-17 5-31-17	Naphthalene	ND	0.0084	EPA 8270D/SIM	5-31-17	5-31-17	
4-Chloroaniline ND 0.21 EPA 8270D 5-31-17 6-1-17	4-Chloroaniline	ND	0.21	EPA 8270D	5-31-17	6-1-17	
Hexachlorobutadiene ND 0.042 EPA 8270D 5-31-17 6-1-17	Hexachlorobutadiene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
4-Chloro-3-methylphenol ND 0.042 EPA 8270D 5-31-17 6-1-17	4-Chloro-3-methylphenol	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2-Methylnaphthalene ND 0.0084 EPA 8270D/SIM 5-31-17 5-31-17	2-Methylnaphthalene	ND	0.0084	EPA 8270D/SIM	5-31-17	5-31-17	
1-Methylnaphthalene ND 0.0084 EPA 8270D/SIM 5-31-17 5-31-17	1-Methylnaphthalene	ND	0.0084	EPA 8270D/SIM	5-31-17	5-31-17	
Hexachlorocyclopentadiene ND 0.042 EPA 8270D 5-31-17 6-1-17		ND	0.042	EPA 8270D	5-31-17	6-1-17	
2,4,6-Trichlorophenol ND 0.042 EPA 8270D 5-31-17 6-1-17	2,4,6-Trichlorophenol	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2,3-Dichloroaniline ND 0.042 EPA 8270D 5-31-17 6-1-17	2,3-Dichloroaniline	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2,4,5-Trichlorophenol ND 0.042 EPA 8270D 5-31-17 6-1-17	2,4,5-Trichlorophenol	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2-Chloronaphthalene ND 0.042 EPA 8270D 5-31-17 6-1-17	2-Chloronaphthalene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2-Nitroaniline ND 0.042 EPA 8270D 5-31-17 6-1-17	2-Nitroaniline	ND	0.042	EPA 8270D	5-31-17	6-1-17	
1,4-Dinitrobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17	1,4-Dinitrobenzene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Dimethylphthalate ND 0.042 EPA 8270D 5-31-17 6-1-17	Dimethylphthalate	ND	0.042	EPA 8270D	5-31-17	6-1-17	
1,3-Dinitrobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17	1,3-Dinitrobenzene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2,6-Dinitrotoluene ND 0.042 EPA 8270D 5-31-17 6-1-17	2,6-Dinitrotoluene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
1,2-Dinitrobenzene ND 0.042 EPA 8270D 5-31-17 6-1-17	1,2-Dinitrobenzene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Acenaphthylene ND 0.0084 EPA 8270D/SIM 5-31-17 5-31-17	Acenaphthylene	ND	0.0084	EPA 8270D/SIM	5-31-17		
3-Nitroaniline ND 0.042 EPA 8270D 5-31-17 6-1-17	3-Nitroaniline	ND	0.042	EPA 8270D	5-31-17	6-1-17	

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP3-2.5					110.90
Laboratory ID:	05-299-07					
2,4-Dinitrophenol	ND	0.21	EPA 8270D	5-31-17	6-1-17	
Acenaphthene	ND	0.0084	EPA 8270D/SIM	5-31-17	5-31-17	
4-Nitrophenol	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2,4-Dinitrotoluene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Dibenzofuran	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2,3,5,6-Tetrachlorophenol	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2,3,4,6-Tetrachlorophenol	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Diethylphthalate	ND	0.21	EPA 8270D	5-31-17	6-1-17	
4-Chlorophenyl-phenylether	ND	0.042	EPA 8270D	5-31-17	6-1-17	
4-Nitroaniline	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Fluorene	ND	0.0084	EPA 8270D/SIM	5-31-17	5-31-17	
4,6-Dinitro-2-methylphenol	ND	0.21	EPA 8270D	5-31-17	6-1-17	
n-Nitrosodiphenylamine	ND	0.042	EPA 8270D	5-31-17	6-1-17	
1,2-Diphenylhydrazine	ND	0.042	EPA 8270D	5-31-17	6-1-17	
4-Bromophenyl-phenylether	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Hexachlorobenzene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Pentachlorophenol	ND	0.21	EPA 8270D	5-31-17	6-1-17	
Phenanthrene	ND	0.0084	EPA 8270D/SIM	5-31-17	5-31-17	
Anthracene	ND	0.0084	EPA 8270D/SIM	5-31-17	5-31-17	
Carbazole	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Di-n-butylphthalate	ND	0.21	EPA 8270D	5-31-17	6-1-17	
Fluoranthene	ND	0.0084	EPA 8270D/SIM	5-31-17	5-31-17	
Benzidine	ND	0.42	EPA 8270D	5-31-17	6-1-17	
Pyrene	ND	0.0084	EPA 8270D/SIM	5-31-17	5-31-17	
Butylbenzylphthalate	ND	0.042	EPA 8270D	5-31-17	6-1-17	
bis-2-Ethylhexyladipate	ND	0.042	EPA 8270D	5-31-17	6-1-17	
3,3'-Dichlorobenzidine	ND	0.21	EPA 8270D	5-31-17	6-1-17	
Benzo[a]anthracene	ND	0.0084	EPA 8270D/SIM	5-31-17	5-31-17	
Chrysene	ND	0.0084	EPA 8270D/SIM	5-31-17	5-31-17	
bis(2-Ethylhexyl)phthalate	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Di-n-octylphthalate	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Benzo[b]fluoranthene	ND	0.0084	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo(j,k)fluoranthene	ND	0.0084	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[a]pyrene	ND	0.0084	EPA 8270D/SIM	5-31-17	5-31-17	
Indeno[1,2,3-cd]pyrene	ND	0.0084	EPA 8270D/SIM	5-31-17	5-31-17	
Dibenz[a,h]anthracene	ND	0.0084	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[g,h,i]perylene	ND	0.0084	EPA 8270D/SIM	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorophenol	62	18 - 109				
Phenol-d6	64	25 - 111				
Nitrobenzene-d5	63	22 - 113				
2-Fluorobiphenyl						
	53	30 - 114				
2,4,6-Tribromophenol	53 57	30 - 114 22 - 116				

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP3-10					
Laboratory ID:	05-299-08					
n-Nitrosodimethylamine	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Pyridine	ND	0.41	EPA 8270D	5-31-17	6-1-17	
Phenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Aniline	ND	0.21	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethyl)ether	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Chlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,3-Dichlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,4-Dichlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Benzyl alcohol	ND	0.21	EPA 8270D	5-31-17	6-1-17	
1,2-Dichlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Methylphenol (o-Cresol)	ND	0.041	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroisopropyl)ether	ND	0.041	EPA 8270D	5-31-17	6-1-17	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.041	EPA 8270D	5-31-17	6-1-17	
n-Nitroso-di-n-propylamine	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Hexachloroethane	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Nitrobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Isophorone	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Nitrophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4-Dimethylphenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethoxy)methane	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4-Dichlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,2,4-Trichlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Naphthalene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
4-Chloroaniline	ND	0.21	EPA 8270D	5-31-17	6-1-17	
Hexachlorobutadiene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
4-Chloro-3-methylphenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Methylnaphthalene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
1-Methylnaphthalene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Hexachlorocyclopentadiene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4,6-Trichlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,3-Dichloroaniline	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4,5-Trichlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Chloronaphthalene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Nitroaniline	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,4-Dinitrobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Dimethylphthalate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,3-Dinitrobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,6-Dinitrotoluene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,2-Dinitrobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Acenaphthylene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
3-Nitroaniline	ND	0.041	EPA 8270D	5-31-17	6-1-17	

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP3-10					
Laboratory ID:	05-299-08					
2,4-Dinitrophenol	ND	0.21	EPA 8270D	5-31-17	6-1-17	
Acenaphthene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
4-Nitrophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4-Dinitrotoluene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Dibenzofuran	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,3,5,6-Tetrachlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,3,4,6-Tetrachlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Diethylphthalate	ND	0.21	EPA 8270D	5-31-17	6-1-17	
4-Chlorophenyl-phenylether	ND	0.041	EPA 8270D	5-31-17	6-1-17	
4-Nitroaniline	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Fluorene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
4,6-Dinitro-2-methylphenol	ND	0.21	EPA 8270D	5-31-17	6-1-17	
n-Nitrosodiphenylamine	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,2-Diphenylhydrazine	ND	0.041	EPA 8270D	5-31-17	6-1-17	
4-Bromophenyl-phenylether	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Hexachlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Pentachlorophenol	ND	0.21	EPA 8270D	5-31-17	6-1-17	
Phenanthrene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Anthracene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Carbazole	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Di-n-butylphthalate	ND	0.21	EPA 8270D	5-31-17	6-1-17	
Fluoranthene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Benzidine	ND	0.41	EPA 8270D	5-31-17	6-1-17	
Pyrene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Butylbenzylphthalate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
bis-2-Ethylhexyladipate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
3,3'-Dichlorobenzidine	ND	0.21	EPA 8270D	5-31-17	6-1-17	
Benzo[a]anthracene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Chrysene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
bis(2-Ethylhexyl)phthalate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Di-n-octylphthalate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Benzo[b]fluoranthene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo(j,k)fluoranthene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[a]pyrene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Indeno[1,2,3-cd]pyrene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Dibenz[a,h]anthracene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[g,h,i]perylene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorophenol	74	18 - 109				
Phenol-d6	71	25 - 111				
Nitrobenzene-d5	75	22 - 113				
2-Fluorobiphenyl	70	30 - 114				
2,4,6-Tribromophenol						
	66	22 - 116				

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP4-3					
Laboratory ID:	05-299-10					
n-Nitrosodimethylamine	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Pyridine	ND	0.42	EPA 8270D	5-31-17	6-1-17	
Phenol	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Aniline	ND	0.21	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethyl)ether	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2-Chlorophenol	ND	0.042	EPA 8270D	5-31-17	6-1-17	
1,3-Dichlorobenzene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
1,4-Dichlorobenzene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Benzyl alcohol	ND	0.21	EPA 8270D	5-31-17	6-1-17	
1,2-Dichlorobenzene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2-Methylphenol (o-Cresol)	ND	0.042	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroisopropyl)ether	ND	0.042	EPA 8270D	5-31-17	6-1-17	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.042	EPA 8270D	5-31-17	6-1-17	
n-Nitroso-di-n-propylamine	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Hexachloroethane	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Nitrobenzene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Isophorone	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2-Nitrophenol	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2,4-Dimethylphenol	ND	0.042	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethoxy)methane	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2,4-Dichlorophenol	ND	0.042	EPA 8270D	5-31-17	6-1-17	
1,2,4-Trichlorobenzene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Naphthalene	ND	0.0084	EPA 8270D/SIM	5-31-17	5-31-17	
4-Chloroaniline	ND	0.21	EPA 8270D	5-31-17	6-1-17	
Hexachlorobutadiene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
4-Chloro-3-methylphenol	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2-Methylnaphthalene	0.017	0.0084	EPA 8270D/SIM	5-31-17	5-31-17	
1-Methylnaphthalene	ND	0.0084	EPA 8270D/SIM	5-31-17	5-31-17	
Hexachlorocyclopentadiene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2,4,6-Trichlorophenol	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2,3-Dichloroaniline	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2,4,5-Trichlorophenol	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2-Chloronaphthalene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2-Nitroaniline	ND	0.042	EPA 8270D	5-31-17	6-1-17	
1,4-Dinitrobenzene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Dimethylphthalate	ND	0.042	EPA 8270D	5-31-17	6-1-17	
1,3-Dinitrobenzene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2,6-Dinitrotoluene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
1,2-Dinitrobenzene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Acenaphthylene	ND	0.0084	EPA 8270D/SIM	5-31-17	5-31-17	
3-Nitroaniline	ND	0.042	EPA 8270D	5-31-17	6-1-17	

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

Analyte Result PQL Method Prepared Analyzed Client ID: PP4-3 Beaboratory ID: 05-299-10 05-299-10 5-31-17 6-1-17 2,4-Dinitrophenol ND 0.0084 EPA 8270D 5-31-17 6-1-17 4-Nitrophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 2,4-Dinitrotoluene ND 0.042 EPA 8270D 5-31-17 6-1-17 2,4-Dinitrotoluene ND 0.042 EPA 8270D 5-31-17 6-1-17 2,4-Dinitrotoluene ND 0.042 EPA 8270D 5-31-17 6-1-17 Dibenzofuran ND 0.042 EPA 8270D 5-31-17 6-1-17 2,3,5,6-Tetrachlorophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 2,3,4,6-Tetrachlorophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 4-Chlorophenyl-phenylether ND 0.042 EPA 8270D 5-31-17 6-1-17 4-Nitroaniline ND 0.042	Flags
Laboratory ID: 05-299-10 2,4-Dinitrophenol ND	
2,4-Dinitrophenol ND 0.21 EPA 8270D 5-31-17 6-1-17 Acenaphthene ND 0.0084 EPA 8270D/SIM 5-31-17 5-31-17 4-Nitrophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 2,4-Dinitrotoluene ND 0.042 EPA 8270D 5-31-17 6-1-17 2,4-Dinitrotoluene ND 0.042 EPA 8270D 5-31-17 6-1-17 Dibehylphtogram ND 0.042 EPA 8270D 5-31-17 6-1-17 2,3,5,6-Tetrachlorophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 2,3,4,6-Tetrachlorophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 2,3,4,6-Tetrachlorophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 4-Chlorophenyl-phenylether ND 0.042 EPA 8270D 5-31-17 6-1-17 4-Nitroaniline ND 0.042 EPA 8270D 5-31-17 6-1-17 4,6-Dinitro-2-methylphenol ND 0.042 EPA 827	
Acenaphthene ND 0.0084 EPA 8270D/SIM 5-31-17 5-31-17 4-Nitrophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 2,4-Dinitrotoluene ND 0.042 EPA 8270D 5-31-17 6-1-17 Dibenzofuran ND 0.042 EPA 8270D 5-31-17 6-1-17 2,3,5,6-Tetrachlorophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 2,3,4,6-Tetrachlorophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 2,3,4,6-Tetrachlorophenol ND 0.042 EPA 8270D 5-31-17 6-1-17 4-Chlorophenyl-phenylether ND 0.21 EPA 8270D 5-31-17 6-1-17 4-Nitroaniline ND 0.042 EPA 8270D 5-31-17 6-1-17 4-Nitroaniline ND 0.0042 EPA 8270D 5-31-17 6-1-17 4,6-Dinitro-2-methylphenol ND 0.042 EPA 8270D 5-31-17 6-1-17 4,6-Dinitro-2-methylphdrazine ND 0.042 E	
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Di-n-butylphthalate ND 0.21 EPA 8270D 5-31-17 6-1-17	
Fluoranthene 0.042 0.0084 EPA 8270D/SIM 5-31-17 5-31-17	
Benzidine ND 0.42 EPA 8270D 5-31-17 6-1-17	
Pyrene 0.046 0.042 EPA 8270D 5-31-17 6-1-17	
Butylbenzylphthalate ND 0.042 EPA 8270D 5-31-17 6-1-17	
bis-2-Ethylhexyladipate ND 0.042 EPA 8270D 5-31-17 6-1-17	
3,3'-Dichlorobenzidine ND 0.21 EPA 8270D 5-31-17 6-1-17	
Benzo[a]anthracene 0.029 0.0084 EPA 8270D/SIM 5-31-17 5-31-17	
Chrysene 0.037 0.0084 EPA 8270D/SIM 5-31-17 5-31-17	
bis(2-Ethylhexyl)phthalate ND 0.042 EPA 8270D 5-31-17 6-1-17	
Di-n-octylphthalate ND 0.042 EPA 8270D 5-31-17 6-1-17	
Benzo[b]fluoranthene 0.042 0.0084 EPA 8270D/SIM 5-31-17 5-31-17	
Benzo(j,k)fluoranthene 0.014 0.0084 EPA 8270D/SIM 5-31-17 5-31-17	
Benzo[a]pyrene 0.034 0.0084 EPA 8270D/SIM 5-31-17 5-31-17	
Indeno[1,2,3-cd]pyrene 0.027 0.0084 EPA 8270D/SIM 5-31-17 5-31-17	
Dibenz[a,h]anthracene ND 0.0084 EPA 8270D/SIM 5-31-17 5-31-17	
Benzo[g,h,i]perylene 0.032 0.0084 EPA 8270D/SIM 5-31-17 5-31-17	
Surrogate: Percent Recovery Control Limits	
2-Fluorophenol 65 18 - 109	
Phenol-d6 66 25 - 111	
Nitrobenzene-d5 67 22 - 113	
2-Fluorobiphenyl 58 30 - 114	
2,4,6-Tribromophenol 61 22 - 116	
Terphenyl-d14 57 33 - 114	

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP4-10					
Laboratory ID:	05-299-11					
n-Nitrosodimethylamine	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Pyridine	ND	0.41	EPA 8270D	5-31-17	6-1-17	
Phenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Aniline	ND	0.21	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethyl)ether	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Chlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,3-Dichlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,4-Dichlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Benzyl alcohol	ND	0.21	EPA 8270D	5-31-17	6-1-17	
1,2-Dichlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Methylphenol (o-Cresol)	ND	0.041	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroisopropyl)ether	ND	0.041	EPA 8270D	5-31-17	6-1-17	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.041	EPA 8270D	5-31-17	6-1-17	
n-Nitroso-di-n-propylamine	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Hexachloroethane	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Nitrobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Isophorone	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Nitrophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4-Dimethylphenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethoxy)methane	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4-Dichlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,2,4-Trichlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Naphthalene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
4-Chloroaniline	ND	0.21	EPA 8270D	5-31-17	6-1-17	
Hexachlorobutadiene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
4-Chloro-3-methylphenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Methylnaphthalene	0.010	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
1-Methylnaphthalene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Hexachlorocyclopentadiene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4,6-Trichlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,3-Dichloroaniline	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4,5-Trichlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Chloronaphthalene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Nitroaniline	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,4-Dinitrobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Dimethylphthalate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,3-Dinitrobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,6-Dinitrotoluene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,2-Dinitrobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Acenaphthylene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
3-Nitroaniline	ND	0.041	EPA 8270D	5-31-17	6-1-17	

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP4-10	. ~-	ou		7 iii ii i j	. iugo
Laboratory ID:	05-299-11					
2,4-Dinitrophenol	ND	0.21	EPA 8270D	5-31-17	6-1-17	
Acenaphthene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
4-Nitrophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4-Dinitrotoluene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Dibenzofuran	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,3,5,6-Tetrachlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,3,4,6-Tetrachlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Diethylphthalate	ND	0.21	EPA 8270D	5-31-17	6-1-17	
4-Chlorophenyl-phenylether	ND	0.041	EPA 8270D	5-31-17	6-1-17	
4-Nitroaniline	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Fluorene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
4,6-Dinitro-2-methylphenol	ND	0.21	EPA 8270D	5-31-17	6-1-17	
n-Nitrosodiphenylamine	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,2-Diphenylhydrazine	ND	0.041	EPA 8270D	5-31-17	6-1-17	
4-Bromophenyl-phenylether	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Hexachlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Pentachlorophenol	ND	0.21	EPA 8270D	5-31-17	6-1-17	
Phenanthrene	0.011	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Anthracene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Carbazole	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Di-n-butylphthalate	ND	0.21	EPA 8270D	5-31-17	6-1-17	
Fluoranthene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Benzidine	ND	0.41	EPA 8270D	5-31-17	6-1-17	
Pyrene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Butylbenzylphthalate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
bis-2-Ethylhexyladipate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
3,3'-Dichlorobenzidine	ND	0.21	EPA 8270D	5-31-17	6-1-17	
Benzo[a]anthracene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Chrysene	0.014	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
bis(2-Ethylhexyl)phthalate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Di-n-octylphthalate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Benzo[b]fluoranthene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo(j,k)fluoranthene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[a]pyrene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Indeno[1,2,3-cd]pyrene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Dibenz[a,h]anthracene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[g,h,i]perylene	ND	0.0082	EPA 8270D/SIM	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorophenol	72	18 - 109				
Phenol-d6	72	25 - 111				
Nitrobenzene-d5	68	22 - 113				
2-Fluorobiphenyl	65	30 - 114				
2,4,6-Tribromophenol	59	22 - 116				
Terphenyl-d14	67	33 - 114				

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

page 1 of 2

Analyte					Date	Date	
Laboratory ID:	Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Nitrosodimethylamine ND	Client ID:	PP5-2.5					
Pyridine ND	Laboratory ID:	05-299-13					
Phenol ND 0.044 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.22 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Misco-Chloroisopropylether ND 0.044 EPA 8270D 5-31-17 6-1-17 4-Hylphenol (o-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 -Misco-Chlorostopropylether ND 0.044 EPA 8270D 5-31-17 6-1-17 -Na 0.044 EPA 8270D 5-31-17 6-	n-Nitrosodimethylamine	ND	0.044	EPA 8270D	5-31-17	6-1-17	
Aniline ND 0.22 EPA 8270D 5-31-17 6-1-7 bis(2-Chloroethyl)ether ND 0.044 EPA 8270D 5-31-17 6-1-7 2-Chlorophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Hexachlorobetane ND 0.044 EPA 8270D 5-31-17 <td>Pyridine</td> <td>ND</td> <td>0.44</td> <td>EPA 8270D</td> <td>5-31-17</td> <td>6-1-17</td> <td></td>	Pyridine	ND	0.44	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethyl)ether ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Mitrophylphenol (o-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 4-Sig-2-Chloroisopropylpether ND 0.044 EPA 8270D 5-31-17 6-1-17 4-Sig-2-Chloroisopropylpether ND 0.044 EPA 8270D 5-31-17 6-1-17 Hexachloroctethane ND 0.044 EPA 8270D 5-31-17 6-1-17 Isoporone ND 0.044	Phenol	ND	0.044	EPA 8270D	5-31-17	6-1-17	
2-Chlorophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorophenol ND 0.044 EPA 8270D 5-31-1	Aniline	ND	0.22	EPA 8270D	5-31-17	6-1-17	
2-Chlorophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorophenol ND 0.044 EPA 8270D 5-31-1	bis(2-Chloroethyl)ether	ND	0.044	EPA 8270D	5-31-17	6-1-17	
1,4-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Benzy alcohol ND 0.22 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroisopropyl)ether ND 0.044 EPA 8270D 5-31-17 6-1-17 si3(-4)-Methylphenol (m,P-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 n-Nitroso-di-n-propylamine ND 0.044 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.044 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 ND (1)-Chioroethoxy)methane ND 0.044 EPA 8270D 5-31-17 6-1-17 2,4-Dichlorophenol ND 0.044		ND	0.044	EPA 8270D	5-31-17	6-1-17	
Benzyl alcohol ND 0.22 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dis(2-Chloroisopropyl)ether ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-d)-Methylphenol (m,p-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-d)-Methylphenol (m,p-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-d)-Methylphenol (m,p-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-d)-Methylphenol ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-d-Nitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,4-Chloroa-methylphenol ND 0.044 EPA 8270D 5-31-17 6-1-17 1-4-Chloroa-methylphenol ND 0.044 EPA 8270D 5-31-17 6-1-17 1-4-Dinitrobenzene ND 0.044	1,3-Dichlorobenzene	ND	0.044	EPA 8270D	5-31-17	6-1-17	
1,2-Dichlorobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 isi(2-Chloroisopropyl)ether ND 0.044 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m.p-Cresol) ND 0.044 EPA 8270D 5-31-17 6-1-17 N-Nitroso-di-n-propylamine ND 0.044 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.044 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.044 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.044 EPA 8270D 5-31-17 6-1-17 2,4-Dimethylphenol ND 0.044 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethoxy)methane ND 0.044 EPA 8270D 5-31-17 6-1-17 2,4-Dimethylphenol ND 0.044 EPA 8270	1,4-Dichlorobenzene	ND	0.044	EPA 8270D	5-31-17	6-1-17	
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Hexachloroethane		ND	0.044	EPA 8270D	5-31-17	6-1-17	
Nitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Nitrophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 2,4-Dimethylphenol ND 0.044 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethoxy)methane ND 0.044 EPA 8270D 5-31-17 6-1-17 2,4-Dichlorophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 2,4-Dichlorophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 2,4-Dichlorophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2,4-Trichlorophenol ND 0.0087 EPA 8270D/SIM 5-31-17 6-1-17 ND 0.0087 EPA 8270D 5-31-17 6-1-17 4-Chloroa-smethylphenol ND 0.044 EPA 8270D 5-31-17 6-1-17 4-Methylnaphthalene ND 0.0087 EPA 8270D/SIM 5-31-17 6-1-1		ND			5-31-17	6-1-17	
Isophorone	Nitrobenzene	ND		EPA 8270D	5-31-17	6-1-17	
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4-Chloroaniline ND 0.22 EPA 8270D 5-31-17 6-1-17 Hexachlorobutadiene ND 0.044 EPA 8270D 5-31-17 6-1-17 4-Chloro-3-methylphenol ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Methylnaphthalene ND 0.0087 EPA 8270D/SIM 5-31-17 5-31-17 1-Methylnaphthalene ND 0.0087 EPA 8270D/SIM 5-31-17 5-31-17 1-Methylnaphthalene ND 0.0087 EPA 8270D/SIM 5-31-17 5-31-17 1-Methylnaphthalene ND 0.0044 EPA 8270D 5-31-17 6-1-17 2-Methylnaphthalene ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Methylnaphthalene ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Methylnaphthalene ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Chloronaphthalene ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Nitroaniline ND 0.044 EPA 8		ND	0.0087		5-31-17	5-31-17	
4-Chloro-3-methylphenol ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Methylnaphthalene ND 0.0087 EPA 8270D/SIM 5-31-17 5-31-17 1-Methylnaphthalene ND 0.0087 EPA 8270D/SIM 5-31-17 5-31-17 Hexachlorocyclopentadiene ND 0.044 EPA 8270D 5-31-17 6-1-17 2,4,6-Trichlorophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 2,3-Dichloroaniline ND 0.044 EPA 8270D 5-31-17 6-1-17 2,4,5-Trichlorophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Chloronaphthalene ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Nitroaniline ND 0.044 EPA 8270D 5-31-17 6-1-17 1,4-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 2,6-Dinitrotoluene ND 0.044		ND	0.22	EPA 8270D	5-31-17	6-1-17	
2-MethylnaphthaleneND0.0087EPA 8270D/SIM5-31-175-31-171-MethylnaphthaleneND0.0087EPA 8270D/SIM5-31-175-31-17HexachlorocyclopentadieneND0.044EPA 8270D5-31-176-1-172,4,6-TrichlorophenolND0.044EPA 8270D5-31-176-1-172,3-DichloroanilineND0.044EPA 8270D5-31-176-1-172,4,5-TrichlorophenolND0.044EPA 8270D5-31-176-1-172-ChloronaphthaleneND0.044EPA 8270D5-31-176-1-172-NitroanilineND0.044EPA 8270D5-31-176-1-171,4-DinitrobenzeneND0.044EPA 8270D5-31-176-1-171,3-DinitrobenzeneND0.044EPA 8270D5-31-176-1-171,3-DinitrobenzeneND0.044EPA 8270D5-31-176-1-172,6-DinitrotolueneND0.044EPA 8270D5-31-176-1-171,2-DinitrobenzeneND0.044EPA 8270D5-31-176-1-17AcenaphthyleneND0.044EPA 8270D/SIM5-31-175-31-17	Hexachlorobutadiene	ND	0.044	EPA 8270D	5-31-17	6-1-17	
2-MethylnaphthaleneND0.0087EPA 8270D/SIM5-31-175-31-171-MethylnaphthaleneND0.0087EPA 8270D/SIM5-31-175-31-17HexachlorocyclopentadieneND0.044EPA 8270D5-31-176-1-172,4,6-TrichlorophenolND0.044EPA 8270D5-31-176-1-172,3-DichloroanilineND0.044EPA 8270D5-31-176-1-172,4,5-TrichlorophenolND0.044EPA 8270D5-31-176-1-172-ChloronaphthaleneND0.044EPA 8270D5-31-176-1-172-NitroanilineND0.044EPA 8270D5-31-176-1-171,4-DinitrobenzeneND0.044EPA 8270D5-31-176-1-17DimethylphthalateND0.044EPA 8270D5-31-176-1-171,3-DinitrobenzeneND0.044EPA 8270D5-31-176-1-172,6-DinitrotolueneND0.044EPA 8270D5-31-176-1-171,2-DinitrobenzeneND0.044EPA 8270D5-31-176-1-17AcenaphthyleneND0.044EPA 8270D/SIM5-31-175-31-17	4-Chloro-3-methylphenol	ND	0.044	EPA 8270D	5-31-17	6-1-17	
1-Methylnaphthalene ND 0.0087 EPA 8270D/SIM 5-31-17 5-31-17 Hexachlorocyclopentadiene ND 0.044 EPA 8270D 5-31-17 6-1-17 2,4,6-Trichlorophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 2,3-Dichloroaniline ND 0.044 EPA 8270D 5-31-17 6-1-17 2,4,5-Trichlorophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Chloronaphthalene ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Nitroaniline ND 0.044 EPA 8270D 5-31-17 6-1-17 1,4-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 2,6-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2-Dinitrobenzene ND 0.044 EPA 8270D/SI		ND	0.0087	EPA 8270D/SIM	5-31-17	5-31-17	
HexachlorocyclopentadieneND0.044EPA 8270D5-31-176-1-172,4,6-TrichlorophenolND0.044EPA 8270D5-31-176-1-172,3-DichloroanilineND0.044EPA 8270D5-31-176-1-172,4,5-TrichlorophenolND0.044EPA 8270D5-31-176-1-172-ChloronaphthaleneND0.044EPA 8270D5-31-176-1-172-NitroanilineND0.044EPA 8270D5-31-176-1-171,4-DinitrobenzeneND0.044EPA 8270D5-31-176-1-17DimethylphthalateND0.044EPA 8270D5-31-176-1-171,3-DinitrobenzeneND0.044EPA 8270D5-31-176-1-172,6-DinitrotolueneND0.044EPA 8270D5-31-176-1-171,2-DinitrobenzeneND0.044EPA 8270D5-31-176-1-17AcenaphthyleneND0.0087EPA 8270D/SIM5-31-175-31-17		ND	0.0087		5-31-17		
2,4,6-Trichlorophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 2,3-Dichloroaniline ND 0.044 EPA 8270D 5-31-17 6-1-17 2,4,5-Trichlorophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Chloronaphthalene ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Nitroaniline ND 0.044 EPA 8270D 5-31-17 6-1-17 1,4-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 2,6-Dinitrotoluene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Acenaphthylene ND 0.0087 EPA 8270D/SIM 5-31-17 5-31-17	Hexachlorocyclopentadiene	ND	0.044	EPA 8270D	5-31-17	6-1-17	
2,4,5-Trichlorophenol ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Chloronaphthalene ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Nitroaniline ND 0.044 EPA 8270D 5-31-17 6-1-17 1,4-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Dimethylphthalate ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 2,6-Dinitrotoluene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Acenaphthylene ND 0.0087 EPA 8270D/SIM 5-31-17 5-31-17		ND	0.044	EPA 8270D	5-31-17	6-1-17	
2-Chloronaphthalene ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Nitroaniline ND 0.044 EPA 8270D 5-31-17 6-1-17 1,4-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Dimethylphthalate ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 2,6-Dinitrotoluene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Acenaphthylene ND 0.0087 EPA 8270D/SIM 5-31-17 5-31-17	2,3-Dichloroaniline	ND	0.044	EPA 8270D	5-31-17	6-1-17	
2-Chloronaphthalene ND 0.044 EPA 8270D 5-31-17 6-1-17 2-Nitroaniline ND 0.044 EPA 8270D 5-31-17 6-1-17 1,4-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Dimethylphthalate ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 2,6-Dinitrotoluene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Acenaphthylene ND 0.0087 EPA 8270D/SIM 5-31-17 5-31-17	2,4,5-Trichlorophenol	ND	0.044	EPA 8270D	5-31-17	6-1-17	
2-Nitroaniline ND 0.044 EPA 8270D 5-31-17 6-1-17 1,4-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Dimethylphthalate ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 2,6-Dinitrotoluene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Acenaphthylene ND 0.0087 EPA 8270D/SIM 5-31-17 5-31-17		ND	0.044	EPA 8270D	5-31-17	6-1-17	
Dimethylphthalate ND 0.044 EPA 8270D 5-31-17 6-1-17 1,3-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 2,6-Dinitrotoluene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Acenaphthylene ND 0.0087 EPA 8270D/SIM 5-31-17 5-31-17	*	ND	0.044	EPA 8270D	5-31-17	6-1-17	
1,3-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 2,6-Dinitrotoluene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Acenaphthylene ND 0.0087 EPA 8270D/SIM 5-31-17 5-31-17	1,4-Dinitrobenzene	ND	0.044	EPA 8270D	5-31-17	6-1-17	
2,6-Dinitrotoluene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Acenaphthylene ND 0.0087 EPA 8270D/SIM 5-31-17 5-31-17		ND	0.044		5-31-17	6-1-17	
2,6-Dinitrotoluene ND 0.044 EPA 8270D 5-31-17 6-1-17 1,2-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Acenaphthylene ND 0.0087 EPA 8270D/SIM 5-31-17 5-31-17	1,3-Dinitrobenzene	ND	0.044	EPA 8270D	5-31-17	6-1-17	
1,2-Dinitrobenzene ND 0.044 EPA 8270D 5-31-17 6-1-17 Acenaphthylene ND 0.0087 EPA 8270D/SIM 5-31-17 5-31-17	2,6-Dinitrotoluene	ND	0.044	EPA 8270D	5-31-17	6-1-17	
		ND	0.044			6-1-17	
	Acenaphthylene	ND	0.0087	EPA 8270D/SIM	5-31-17	5-31-17	
		ND	0.044	EPA 8270D	5-31-17	6-1-17	

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

	Result	PQL	Method	Prepared	Analyzed	Flags
Analyte Client ID:	PP5-2.5					
Laboratory ID:	05-299-13					
2,4-Dinitrophenol	ND	0.22	EPA 8270D	5-31-17	6-1-17	
Acenaphthene	ND	0.0087	EPA 8270D/SIM	5-31-17	5-31-17	
4-Nitrophenol	ND	0.044	EPA 8270D	5-31-17	6-1-17	
2,4-Dinitrotoluene	ND	0.044	EPA 8270D	5-31-17	6-1-17	
Dibenzofuran	ND	0.044	EPA 8270D	5-31-17	6-1-17	
2,3,5,6-Tetrachlorophenol	ND	0.044	EPA 8270D	5-31-17	6-1-17	
2,3,4,6-Tetrachlorophenol	ND	0.044	EPA 8270D	5-31-17	6-1-17	
Diethylphthalate	ND	0.22	EPA 8270D	5-31-17	6-1-17	
4-Chlorophenyl-phenylether	ND	0.044	EPA 8270D	5-31-17	6-1-17	
4-Nitroaniline	ND	0.044	EPA 8270D	5-31-17	6-1-17	
Fluorene	ND	0.0087	EPA 8270D/SIM	5-31-17	5-31-17	
4,6-Dinitro-2-methylphenol	ND	0.22	EPA 8270D	5-31-17	6-1-17	
n-Nitrosodiphenylamine	ND	0.044	EPA 8270D	5-31-17	6-1-17	
1,2-Diphenylhydrazine	ND	0.044	EPA 8270D	5-31-17	6-1-17	
4-Bromophenyl-phenylether	ND	0.044	EPA 8270D	5-31-17	6-1-17	
Hexachlorobenzene	ND	0.044	EPA 8270D	5-31-17	6-1-17	
Pentachlorophenol	ND	0.22	EPA 8270D	5-31-17	6-1-17	
Phenanthrene	ND	0.0087	EPA 8270D/SIM	5-31-17	5-31-17	
Anthracene	ND	0.0087	EPA 8270D/SIM	5-31-17	5-31-17	
Carbazole	ND	0.044	EPA 8270D	5-31-17	6-1-17	
Di-n-butylphthalate	ND	0.22	EPA 8270D	5-31-17	6-1-17	
Fluoranthene	ND	0.0087	EPA 8270D/SIM	5-31-17	5-31-17	
Benzidine	ND	0.44	EPA 8270D	5-31-17	6-1-17	
Pyrene	ND	0.0087	EPA 8270D/SIM	5-31-17	5-31-17	
Butylbenzylphthalate	ND	0.044	EPA 8270D	5-31-17	6-1-17	
bis-2-Ethylhexyladipate	ND	0.044	EPA 8270D	5-31-17	6-1-17	
3,3'-Dichlorobenzidine	ND	0.22	EPA 8270D	5-31-17	6-1-17	
Benzo[a]anthracene	ND	0.0087	EPA 8270D/SIM	5-31-17	5-31-17	
Chrysene	ND	0.0087	EPA 8270D/SIM	5-31-17	5-31-17	
bis(2-Ethylhexyl)phthalate	ND	0.044	EPA 8270D	5-31-17	6-1-17	
Di-n-octylphthalate	ND	0.044	EPA 8270D	5-31-17	6-1-17	
Benzo[b]fluoranthene	ND	0.0087	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo(j,k)fluoranthene	ND	0.0087	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[a]pyrene	ND	0.0087	EPA 8270D/SIM	5-31-17	5-31-17	
Indeno[1,2,3-cd]pyrene	ND	0.0087	EPA 8270D/SIM	5-31-17	5-31-17	
Dibenz[a,h]anthracene	ND	0.0087	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[g,h,i]perylene	ND	0.0087	EPA 8270D/SIM	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorophenol	60	18 - 109				
Phenol-d6	60	25 - 111				
Nitrobenzene-d5	57	22 - 113				
2-Fluorobiphenyl	51	30 - 114				
2,4,6-Tribromophenol	46	22 - 116				
Terphenyl-d14	45	33 - 114				

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP5-10					
Laboratory ID:	05-299-14					
n-Nitrosodimethylamine	ND	0.045	EPA 8270D	5-31-17	6-1-17	
Pyridine	ND	0.45	EPA 8270D	5-31-17	6-1-17	
Phenol	ND	0.045	EPA 8270D	5-31-17	6-1-17	
Aniline	ND	0.22	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethyl)ether	ND	0.045	EPA 8270D	5-31-17	6-1-17	
2-Chlorophenol	ND	0.045	EPA 8270D	5-31-17	6-1-17	
1,3-Dichlorobenzene	ND	0.045	EPA 8270D	5-31-17	6-1-17	
1,4-Dichlorobenzene	ND	0.045	EPA 8270D	5-31-17	6-1-17	
Benzyl alcohol	ND	0.22	EPA 8270D	5-31-17	6-1-17	
1,2-Dichlorobenzene	ND	0.045	EPA 8270D	5-31-17	6-1-17	
2-Methylphenol (o-Cresol)	ND	0.045	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroisopropyl)ether	ND	0.045	EPA 8270D	5-31-17	6-1-17	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.045	EPA 8270D	5-31-17	6-1-17	
n-Nitroso-di-n-propylamine	ND	0.045	EPA 8270D	5-31-17	6-1-17	
Hexachloroethane	ND	0.045	EPA 8270D	5-31-17	6-1-17	
Nitrobenzene	ND	0.045	EPA 8270D	5-31-17	6-1-17	
Isophorone	ND	0.045	EPA 8270D	5-31-17	6-1-17	
2-Nitrophenol	ND	0.045	EPA 8270D	5-31-17	6-1-17	
2,4-Dimethylphenol	ND	0.045	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethoxy)methane	ND	0.045	EPA 8270D	5-31-17	6-1-17	
2,4-Dichlorophenol	ND	0.045	EPA 8270D	5-31-17	6-1-17	
1,2,4-Trichlorobenzene	ND	0.045	EPA 8270D	5-31-17	6-1-17	
Naphthalene	ND	0.0090	EPA 8270D/SIM	5-31-17	5-31-17	
4-Chloroaniline	ND	0.22	EPA 8270D	5-31-17	6-1-17	
Hexachlorobutadiene	ND	0.045	EPA 8270D	5-31-17	6-1-17	
4-Chloro-3-methylphenol	ND	0.045	EPA 8270D	5-31-17	6-1-17	
2-Methylnaphthalene	ND	0.0090	EPA 8270D/SIM	5-31-17	5-31-17	
1-Methylnaphthalene	ND	0.0090	EPA 8270D/SIM	5-31-17	5-31-17	
Hexachlorocyclopentadiene	ND	0.045	EPA 8270D	5-31-17	6-1-17	
2,4,6-Trichlorophenol	ND	0.045	EPA 8270D	5-31-17	6-1-17	
2,3-Dichloroaniline	ND	0.045	EPA 8270D	5-31-17	6-1-17	
2,4,5-Trichlorophenol	ND	0.045	EPA 8270D	5-31-17	6-1-17	
2-Chloronaphthalene	ND	0.045	EPA 8270D	5-31-17	6-1-17	
2-Nitroaniline	ND	0.045	EPA 8270D	5-31-17	6-1-17	
1,4-Dinitrobenzene	ND	0.045	EPA 8270D	5-31-17	6-1-17	
Dimethylphthalate	ND	0.045	EPA 8270D	5-31-17	6-1-17	
1,3-Dinitrobenzene	ND	0.045	EPA 8270D	5-31-17	6-1-17	
2,6-Dinitrotoluene	ND	0.045	EPA 8270D	5-31-17	6-1-17	
1,2-Dinitrobenzene	ND	0.045	EPA 8270D	5-31-17	6-1-17	
Acenaphthylene	ND	0.0090	EPA 8270D/SIM	5-31-17	5-31-17	
3-Nitroaniline	ND	0.045	EPA 8270D	5-31-17	6-1-17	

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP5-10					
Laboratory ID:	05-299-14					
2,4-Dinitrophenol	ND	0.22	EPA 8270D	5-31-17	6-1-17	
Acenaphthene	ND	0.0090	EPA 8270D/SIM	5-31-17	5-31-17	
4-Nitrophenol	ND	0.045	EPA 8270D	5-31-17	6-1-17	
2,4-Dinitrotoluene	ND	0.045	EPA 8270D	5-31-17	6-1-17	
Dibenzofuran	ND	0.045	EPA 8270D	5-31-17	6-1-17	
2,3,5,6-Tetrachlorophenol	ND	0.045	EPA 8270D	5-31-17	6-1-17	
2,3,4,6-Tetrachlorophenol	ND	0.045	EPA 8270D	5-31-17	6-1-17	
Diethylphthalate	ND	0.22	EPA 8270D	5-31-17	6-1-17	
4-Chlorophenyl-phenylether	ND	0.045	EPA 8270D	5-31-17	6-1-17	
4-Nitroaniline	ND	0.045	EPA 8270D	5-31-17	6-1-17	
Fluorene	ND	0.0090	EPA 8270D/SIM	5-31-17	5-31-17	
4,6-Dinitro-2-methylphenol	ND	0.22	EPA 8270D	5-31-17	6-1-17	
n-Nitrosodiphenylamine	ND	0.045	EPA 8270D	5-31-17	6-1-17	
1,2-Diphenylhydrazine	ND	0.045	EPA 8270D	5-31-17	6-1-17	
4-Bromophenyl-phenylether	ND	0.045	EPA 8270D	5-31-17	6-1-17	
Hexachlorobenzene	ND	0.045	EPA 8270D	5-31-17	6-1-17	
Pentachlorophenol	ND	0.22	EPA 8270D	5-31-17	6-1-17	
Phenanthrene	ND	0.0090	EPA 8270D/SIM	5-31-17	5-31-17	
Anthracene	ND	0.0090	EPA 8270D/SIM	5-31-17	5-31-17	
Carbazole	ND	0.045	EPA 8270D	5-31-17	6-1-17	
Di-n-butylphthalate	ND	0.22	EPA 8270D	5-31-17	6-1-17	
Fluoranthene	ND	0.0090	EPA 8270D/SIM	5-31-17	5-31-17	
Benzidine	ND	0.45	EPA 8270D	5-31-17	6-1-17	
Pyrene	ND	0.0090	EPA 8270D/SIM	5-31-17	5-31-17	
Butylbenzylphthalate	ND	0.045	EPA 8270D	5-31-17	6-1-17	
bis-2-Ethylhexyladipate	ND	0.045	EPA 8270D	5-31-17	6-1-17	
3,3'-Dichlorobenzidine	ND	0.22	EPA 8270D	5-31-17	6-1-17	
Benzo[a]anthracene	ND	0.0090	EPA 8270D/SIM	5-31-17	5-31-17	
Chrysene	ND	0.0090	EPA 8270D/SIM	5-31-17	5-31-17	
bis(2-Ethylhexyl)phthalate	ND	0.045	EPA 8270D	5-31-17	6-1-17	
Di-n-octylphthalate	ND	0.045	EPA 8270D	5-31-17	6-1-17	
Benzo[b]fluoranthene	ND	0.0090	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo(j,k)fluoranthene	ND	0.0090	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[a]pyrene	ND	0.0090	EPA 8270D/SIM	5-31-17	5-31-17	
Indeno[1,2,3-cd]pyrene	ND	0.0090	EPA 8270D/SIM	5-31-17	5-31-17	
Dibenz[a,h]anthracene	ND	0.0090	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[g,h,i]perylene	ND	0.0090	EPA 8270D/SIM	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorophenol	67	18 - 109				
Phenol-d6	67	25 - 111				
Nitrobenzene-d5	52	22 - 113				
2-Fluorobiphenyl	56	30 - 114				
	<i>5</i> 5	22 - 116				
	62					
2,4,6-Tribromophenol Terphenyl-d14	55					

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

page 1 of 2

Analyte Result PQL Method Prepared Analyzed Flags Client ID: PP6-2 Laboratory ID: 05-299-16 n-Nitrosodimethylamine ND 0.039 EPA 8270D 5-31-17 6-1-17 Pyridine ND 0.39 EPA 8270D 5-31-17 6-1-17 Phenol ND 0.039 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.19 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.039 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.039 EPA 8270D 5-31-17 6-1-17
Laboratory ID: 05-299-16 n-Nitrosodimethylamine ND 0.039 EPA 8270D 5-31-17 6-1-17 Pyridine ND 0.39 EPA 8270D 5-31-17 6-1-17 Phenol ND 0.039 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.19 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.039 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.039 EPA 8270D 5-31-17 6-1-17
n-Nitrosodimethylamine ND 0.039 EPA 8270D 5-31-17 6-1-17 Pyridine ND 0.39 EPA 8270D 5-31-17 6-1-17 Phenol ND 0.039 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.19 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.039 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.039 EPA 8270D 5-31-17 6-1-17
Pyridine ND 0.39 EPA 8270D 5-31-17 6-1-17 Phenol ND 0.039 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.19 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.039 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.039 EPA 8270D 5-31-17 6-1-17
Pyridine ND 0.39 EPA 8270D 5-31-17 6-1-17 Phenol ND 0.039 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.19 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.039 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.039 EPA 8270D 5-31-17 6-1-17
Aniline ND 0.19 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.039 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.039 EPA 8270D 5-31-17 6-1-17
bis(2-Chloroethyl)ether ND 0.039 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.039 EPA 8270D 5-31-17 6-1-17
2-Chlorophenol ND 0.039 EPA 8270D 5-31-17 6-1-17
1,3-Dichlorobenzene ND 0.039 EPA 8270D 5-31-17 6-1-17
1,4-Dichlorobenzene ND 0.039 EPA 8270D 5-31-17 6-1-17
Benzyl alcohol ND 0.19 EPA 8270D 5-31-17 6-1-17
1,2-Dichlorobenzene ND 0.039 EPA 8270D 5-31-17 6-1-17
2-Methylphenol (o-Cresol) ND 0.039 EPA 8270D 5-31-17 6-1-17
bis(2-Chloroisopropyl)ether ND 0.039 EPA 8270D 5-31-17 6-1-17
(3+4)-Methylphenol (m,p-Cresol) ND 0.039 EPA 8270D 5-31-17 6-1-17
n-Nitroso-di-n-propylamine ND 0.039 EPA 8270D 5-31-17 6-1-17
Hexachloroethane ND 0.039 EPA 8270D 5-31-17 6-1-17
Nitrobenzene ND 0.039 EPA 8270D 5-31-17 6-1-17
Isophorone ND 0.039 EPA 8270D 5-31-17 6-1-17
2-Nitrophenol ND 0.039 EPA 8270D 5-31-17 6-1-17
2,4-Dimethylphenol ND 0.039 EPA 8270D 5-31-17 6-1-17
bis(2-Chloroethoxy)methane ND 0.039 EPA 8270D 5-31-17 6-1-17
2,4-Dichlorophenol ND 0.039 EPA 8270D 5-31-17 6-1-17
1,2,4-Trichlorobenzene ND 0.039 EPA 8270D 5-31-17 6-1-17
Naphthalene ND 0.0077 EPA 8270D/SIM 5-31-17 6-1-17
4-Chloroaniline ND 0.19 EPA 8270D 5-31-17 6-1-17
Hexachlorobutadiene ND 0.039 EPA 8270D 5-31-17 6-1-17
4-Chloro-3-methylphenol ND 0.039 EPA 8270D 5-31-17 6-1-17
2-Methylnaphthalene ND 0.0077 EPA 8270D/SIM 5-31-17 6-1-17
1-Methylnaphthalene ND 0.0077 EPA 8270D/SIM 5-31-17 6-1-17
Hexachlorocyclopentadiene ND 0.039 EPA 8270D 5-31-17 6-1-17
2,4,6-Trichlorophenol ND 0.039 EPA 8270D 5-31-17 6-1-17
2,3-Dichloroaniline ND 0.039 EPA 8270D 5-31-17 6-1-17
2,4,5-Trichlorophenol ND 0.039 EPA 8270D 5-31-17 6-1-17
2-Chloronaphthalene ND 0.039 EPA 8270D 5-31-17 6-1-17
2-Nitroaniline ND 0.039 EPA 8270D 5-31-17 6-1-17
1,4-Dinitrobenzene ND 0.039 EPA 8270D 5-31-17 6-1-17
Dimethylphthalate ND 0.039 EPA 8270D 5-31-17 6-1-17
1,3-Dinitrobenzene ND 0.039 EPA 8270D 5-31-17 6-1-17
2,6-Dinitrotoluene ND 0.039 EPA 8270D 5-31-17 6-1-17
1,2-Dinitrobenzene ND 0.039 EPA 8270D 5-31-17 6-1-17
Acenaphthylene ND 0.0077 EPA 8270D/SIM 5-31-17 6-1-17
3-Nitroaniline ND 0.039 EPA 8270D 5-31-17 6-1-17

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP6-2					
Laboratory ID:	05-299-16					
2,4-Dinitrophenol	ND	0.19	EPA 8270D	5-31-17	6-1-17	
Acenaphthene	ND	0.0077	EPA 8270D/SIM	5-31-17	6-1-17	
4-Nitrophenol	ND	0.039	EPA 8270D	5-31-17	6-1-17	
2,4-Dinitrotoluene	ND	0.039	EPA 8270D	5-31-17	6-1-17	
Dibenzofuran	ND	0.039	EPA 8270D	5-31-17	6-1-17	
2,3,5,6-Tetrachlorophenol	ND	0.039	EPA 8270D	5-31-17	6-1-17	
2,3,4,6-Tetrachlorophenol	ND	0.039	EPA 8270D	5-31-17	6-1-17	
Diethylphthalate	ND	0.19	EPA 8270D	5-31-17	6-1-17	
4-Chlorophenyl-phenylether	ND	0.039	EPA 8270D	5-31-17	6-1-17	
4-Nitroaniline	ND	0.039	EPA 8270D	5-31-17	6-1-17	
Fluorene	ND	0.0077	EPA 8270D/SIM	5-31-17	6-1-17	
4,6-Dinitro-2-methylphenol	ND	0.19	EPA 8270D	5-31-17	6-1-17	
n-Nitrosodiphenylamine	ND	0.039	EPA 8270D	5-31-17	6-1-17	
1,2-Diphenylhydrazine	ND	0.039	EPA 8270D	5-31-17	6-1-17	
4-Bromophenyl-phenylether	ND	0.039	EPA 8270D	5-31-17	6-1-17	
Hexachlorobenzene	ND	0.039	EPA 8270D	5-31-17	6-1-17	
Pentachlorophenol	ND	0.19	EPA 8270D	5-31-17	6-1-17	
Phenanthrene	ND	0.0077	EPA 8270D/SIM	5-31-17	6-1-17	
Anthracene	ND	0.0077	EPA 8270D/SIM	5-31-17	6-1-17	
Carbazole	ND	0.039	EPA 8270D	5-31-17	6-1-17	
Di-n-butylphthalate	ND	0.19	EPA 8270D	5-31-17	6-1-17	
Fluoranthene	0.013	0.0077	EPA 8270D/SIM	5-31-17	6-1-17	
Benzidine	ND	0.39	EPA 8270D	5-31-17	6-1-17	
Pyrene	0.015	0.0077	EPA 8270D/SIM	5-31-17	6-1-17	
Butylbenzylphthalate	ND	0.039	EPA 8270D	5-31-17	6-1-17	
bis-2-Ethylhexyladipate	ND	0.039	EPA 8270D	5-31-17	6-1-17	
3,3'-Dichlorobenzidine	ND	0.19	EPA 8270D	5-31-17	6-1-17	
Benzo[a]anthracene	0.0084	0.0077	EPA 8270D/SIM	5-31-17	6-1-17	
Chrysene	0.012	0.0077	EPA 8270D/SIM	5-31-17	6-1-17	
bis(2-Ethylhexyl)phthalate	ND	0.039	EPA 8270D	5-31-17	6-1-17	
Di-n-octylphthalate	ND	0.039	EPA 8270D	5-31-17	6-1-17	
Benzo[b]fluoranthene	0.016	0.0077	EPA 8270D/SIM	5-31-17	6-1-17	
Benzo(j,k)fluoranthene	ND	0.0077	EPA 8270D/SIM	5-31-17	6-1-17	
Benzo[a]pyrene	0.012	0.0077	EPA 8270D/SIM	5-31-17	6-1-17	
Indeno[1,2,3-cd]pyrene	0.0092	0.0077	EPA 8270D/SIM	5-31-17	6-1-17	
Dibenz[a,h]anthracene	ND	0.0077	EPA 8270D/SIM	5-31-17	6-1-17	
Benzo[g,h,i]perylene	0.011	0.0077	EPA 8270D/SIM	5-31-17	6-1-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorophenol	79	18 - 109				
Phenol-d6	82	25 - 111				
Nitrobenzene-d5	80	22 - 113				
2-Fluorobiphenyl	63	30 - 114				
2,4,6-Tribromophenol	74	22 - 116				
Terphenyl-d14	63	33 - 114				
ı erpnenyı-a14	63	33 - 114				

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

page 1 of 2

Analyte	Populé					
	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP6-10					
Laboratory ID:	05-299-17					
n-Nitrosodimethylamine	ND	0.050	EPA 8270D	5-31-17	6-1-17	
Pyridine	ND	0.50	EPA 8270D	5-31-17	6-1-17	
Phenol	ND	0.050	EPA 8270D	5-31-17	6-1-17	
Aniline	ND	0.25	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethyl)ether	ND	0.050	EPA 8270D	5-31-17	6-1-17	
2-Chlorophenol	ND	0.050	EPA 8270D	5-31-17	6-1-17	
1,3-Dichlorobenzene	ND	0.050	EPA 8270D	5-31-17	6-1-17	
1,4-Dichlorobenzene	ND	0.050	EPA 8270D	5-31-17	6-1-17	
Benzyl alcohol	ND	0.25	EPA 8270D	5-31-17	6-1-17	
1,2-Dichlorobenzene	ND	0.050	EPA 8270D	5-31-17	6-1-17	
2-Methylphenol (o-Cresol)	ND	0.050	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroisopropyl)ether	ND	0.050	EPA 8270D	5-31-17	6-1-17	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.050	EPA 8270D	5-31-17	6-1-17	
n-Nitroso-di-n-propylamine	ND	0.050	EPA 8270D	5-31-17	6-1-17	
Hexachloroethane	ND	0.050	EPA 8270D	5-31-17	6-1-17	
Nitrobenzene	ND	0.050	EPA 8270D	5-31-17	6-1-17	
Isophorone	ND	0.050	EPA 8270D	5-31-17	6-1-17	
2-Nitrophenol	ND	0.050	EPA 8270D	5-31-17	6-1-17	
2,4-Dimethylphenol	ND	0.050	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethoxy)methane	ND	0.050	EPA 8270D	5-31-17	6-1-17	
2,4-Dichlorophenol	ND	0.050	EPA 8270D	5-31-17	6-1-17	
1,2,4-Trichlorobenzene	ND	0.050	EPA 8270D	5-31-17	6-1-17	
Naphthalene	ND	0.0099	EPA 8270D/SIM	5-31-17	5-31-17	
4-Chloroaniline	ND	0.25	EPA 8270D	5-31-17	6-1-17	
Hexachlorobutadiene	ND	0.050	EPA 8270D	5-31-17	6-1-17	
4-Chloro-3-methylphenol	ND	0.050	EPA 8270D	5-31-17	6-1-17	
2-Methylnaphthalene	ND	0.0099	EPA 8270D/SIM	5-31-17	5-31-17	
1-Methylnaphthalene	ND	0.0099	EPA 8270D/SIM	5-31-17	5-31-17	
Hexachlorocyclopentadiene	ND	0.050	EPA 8270D	5-31-17	6-1-17	
2,4,6-Trichlorophenol	ND	0.050	EPA 8270D	5-31-17	6-1-17	
2,3-Dichloroaniline	ND	0.050	EPA 8270D	5-31-17	6-1-17	
2,4,5-Trichlorophenol	ND	0.050	EPA 8270D	5-31-17	6-1-17	
2-Chloronaphthalene	ND	0.050	EPA 8270D	5-31-17	6-1-17	
2-Nitroaniline	ND	0.050	EPA 8270D	5-31-17	6-1-17	
1,4-Dinitrobenzene	ND	0.050	EPA 8270D	5-31-17	6-1-17	
Dimethylphthalate	ND	0.050	EPA 8270D	5-31-17	6-1-17	
1,3-Dinitrobenzene	ND	0.050	EPA 8270D	5-31-17	6-1-17	
2,6-Dinitrotoluene	ND	0.050	EPA 8270D	5-31-17	6-1-17	
1,2-Dinitrobenzene	ND	0.050	EPA 8270D	5-31-17	6-1-17	
Acenaphthylene	ND	0.0099	EPA 8270D/SIM	5-31-17	5-31-17	
3-Nitroaniline	ND	0.050	EPA 8270D	5-31-17	6-1-17	

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP6-10					110.90
Laboratory ID:	05-299-17					
2,4-Dinitrophenol	ND	0.25	EPA 8270D	5-31-17	6-1-17	
Acenaphthene	ND	0.0099	EPA 8270D/SIM	5-31-17	5-31-17	
4-Nitrophenol	ND	0.050	EPA 8270D	5-31-17	6-1-17	
2,4-Dinitrotoluene	ND	0.050	EPA 8270D	5-31-17	6-1-17	
Dibenzofuran	ND	0.050	EPA 8270D	5-31-17	6-1-17	
2,3,5,6-Tetrachlorophenol	ND	0.050	EPA 8270D	5-31-17	6-1-17	
2,3,4,6-Tetrachlorophenol	ND	0.050	EPA 8270D	5-31-17	6-1-17	
Diethylphthalate	ND	0.25	EPA 8270D	5-31-17	6-1-17	
4-Chlorophenyl-phenylether	ND	0.050	EPA 8270D	5-31-17	6-1-17	
4-Nitroaniline	ND	0.050	EPA 8270D	5-31-17	6-1-17	
Fluorene	ND	0.0099	EPA 8270D/SIM	5-31-17	5-31-17	
4,6-Dinitro-2-methylphenol	ND	0.25	EPA 8270D	5-31-17	6-1-17	
n-Nitrosodiphenylamine	ND	0.050	EPA 8270D	5-31-17	6-1-17	
1,2-Diphenylhydrazine	ND	0.050	EPA 8270D	5-31-17	6-1-17	
4-Bromophenyl-phenylether	ND	0.050	EPA 8270D	5-31-17	6-1-17	
Hexachlorobenzene	ND	0.050	EPA 8270D	5-31-17	6-1-17	
Pentachlorophenol	ND	0.25	EPA 8270D	5-31-17	6-1-17	
Phenanthrene	ND	0.0099	EPA 8270D/SIM	5-31-17	5-31-17	
Anthracene	ND	0.0099	EPA 8270D/SIM	5-31-17	5-31-17	
Carbazole	ND	0.050	EPA 8270D	5-31-17	6-1-17	
Di-n-butylphthalate	ND	0.25	EPA 8270D	5-31-17	6-1-17	
Fluoranthene	ND	0.0099	EPA 8270D/SIM	5-31-17	5-31-17	
Benzidine	ND	0.50	EPA 8270D	5-31-17	6-1-17	
Pyrene	ND	0.0099	EPA 8270D/SIM	5-31-17	5-31-17	
Butylbenzylphthalate	ND	0.050	EPA 8270D	5-31-17	6-1-17	
bis-2-Ethylhexyladipate	ND	0.050	EPA 8270D	5-31-17	6-1-17	
3,3'-Dichlorobenzidine	ND	0.25	EPA 8270D	5-31-17	6-1-17	
Benzo[a]anthracene	ND	0.0099	EPA 8270D/SIM	5-31-17	5-31-17	
Chrysene	ND	0.0099	EPA 8270D/SIM	5-31-17	5-31-17	
bis(2-Ethylhexyl)phthalate	ND	0.050	EPA 8270D	5-31-17	6-1-17	
Di-n-octylphthalate	ND	0.050	EPA 8270D	5-31-17	6-1-17	
Benzo[b]fluoranthene	ND	0.0099	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo(j,k)fluoranthene	ND	0.0099	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[a]pyrene	ND	0.0099	EPA 8270D/SIM	5-31-17	5-31-17	
Indeno[1,2,3-cd]pyrene	ND	0.0099	EPA 8270D/SIM	5-31-17	5-31-17	
Dibenz[a,h]anthracene	ND	0.0099	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[g,h,i]perylene	ND	0.0099	EPA 8270D/SIM	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorophenol	31	18 - 109				
Phenol-d6	34	25 - 111				
Nitrobenzene-d5	31	22 - 113				
2-Fluorobiphenyl	31					
2,4,6-Tribromophenol	35 44	30 - 114 22 - 116				

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP7-0					
Laboratory ID:	05-299-19					
n-Nitrosodimethylamine	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Pyridine	ND	0.41	EPA 8270D	5-31-17	6-1-17	
Phenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Aniline	ND	0.20	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethyl)ether	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Chlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,3-Dichlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,4-Dichlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Benzyl alcohol	ND	0.20	EPA 8270D	5-31-17	6-1-17	
1,2-Dichlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Methylphenol (o-Cresol)	ND	0.041	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroisopropyl)ether	ND	0.041	EPA 8270D	5-31-17	6-1-17	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.041	EPA 8270D	5-31-17	6-1-17	
n-Nitroso-di-n-propylamine	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Hexachloroethane	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Nitrobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Isophorone	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Nitrophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4-Dimethylphenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethoxy)methane	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4-Dichlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,2,4-Trichlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Naphthalene	ND	0.0082	EPA 8270D/SIM	5-31-17	6-1-17	
4-Chloroaniline	ND	0.20	EPA 8270D	5-31-17	6-1-17	
Hexachlorobutadiene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
4-Chloro-3-methylphenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Methylnaphthalene	ND	0.0082	EPA 8270D/SIM	5-31-17	6-1-17	
1-Methylnaphthalene	ND	0.0082	EPA 8270D/SIM	5-31-17	6-1-17	
Hexachlorocyclopentadiene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4,6-Trichlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,3-Dichloroaniline	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4,5-Trichlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Chloronaphthalene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Nitroaniline	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,4-Dinitrobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Dimethylphthalate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,3-Dinitrobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,6-Dinitrotoluene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,2-Dinitrobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Acenaphthylene	ND	0.0082	EPA 8270D/SIM	5-31-17	6-1-17	
3-Nitroaniline	ND	0.041	EPA 8270D	5-31-17	6-1-17	

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP7-0					
Laboratory ID:	05-299-19					
2,4-Dinitrophenol	ND	0.20	EPA 8270D	5-31-17	6-1-17	
Acenaphthene	ND	0.0082	EPA 8270D/SIM	5-31-17	6-1-17	
4-Nitrophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4-Dinitrotoluene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Dibenzofuran	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,3,5,6-Tetrachlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,3,4,6-Tetrachlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Diethylphthalate	ND	0.20	EPA 8270D	5-31-17	6-1-17	
4-Chlorophenyl-phenylether	ND	0.041	EPA 8270D	5-31-17	6-1-17	
4-Nitroaniline	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Fluorene	ND	0.0082	EPA 8270D/SIM	5-31-17	6-1-17	
4,6-Dinitro-2-methylphenol	ND	0.20	EPA 8270D	5-31-17	6-1-17	
n-Nitrosodiphenylamine	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,2-Diphenylhydrazine	ND	0.041	EPA 8270D	5-31-17	6-1-17	
4-Bromophenyl-phenylether	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Hexachlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Pentachlorophenol	ND	0.20	EPA 8270D	5-31-17	6-1-17	
Phenanthrene	ND	0.0082	EPA 8270D/SIM	5-31-17	6-1-17	
Anthracene	ND	0.0082	EPA 8270D/SIM	5-31-17	6-1-17	
Carbazole	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Di-n-butylphthalate	ND	0.20	EPA 8270D	5-31-17	6-1-17	
Fluoranthene	ND	0.0082	EPA 8270D/SIM	5-31-17	6-1-17	
Benzidine	ND	0.41	EPA 8270D	5-31-17	6-1-17	
Pyrene	ND	0.0082	EPA 8270D/SIM	5-31-17	6-1-17	
Butylbenzylphthalate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
bis-2-Ethylhexyladipate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
3,3'-Dichlorobenzidine	ND	0.20	EPA 8270D	5-31-17	6-1-17	
Benzo[a]anthracene	ND	0.0082	EPA 8270D/SIM	5-31-17	6-1-17	
Chrysene	ND	0.0082	EPA 8270D/SIM	5-31-17	6-1-17	
bis(2-Ethylhexyl)phthalate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Di-n-octylphthalate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Benzo[b]fluoranthene	ND	0.0082	EPA 8270D/SIM	5-31-17	6-1-17	
Benzo(j,k)fluoranthene	ND	0.0082	EPA 8270D/SIM	5-31-17	6-1-17	
Benzo[a]pyrene	ND	0.0082	EPA 8270D/SIM	5-31-17	6-1-17	
Indeno[1,2,3-cd]pyrene	ND	0.0082	EPA 8270D/SIM	5-31-17	6-1-17	
Dibenz[a,h]anthracene	ND	0.0082	EPA 8270D/SIM	5-31-17	6-1-17	
Benzo[g,h,i]perylene	ND	0.0082	EPA 8270D/SIM	5-31-17	6-1-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorophenol	62	18 - 109				
Phenol-d6	69	25 - 111				
Nitrobenzene-d5	69	22 - 113				
2-Fluorobiphenyl	71	30 - 114				
2,4,6-Tribromophenol	 78	22 - 116				
Terphenyl-d14	76	33 - 114				
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Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP7-10					
Laboratory ID:	05-299-20					
n-Nitrosodimethylamine	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Pyridine	ND	0.42	EPA 8270D	5-31-17	6-1-17	
Phenol	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Aniline	ND	0.21	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethyl)ether	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2-Chlorophenol	ND	0.042	EPA 8270D	5-31-17	6-1-17	
1,3-Dichlorobenzene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
1,4-Dichlorobenzene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Benzyl alcohol	ND	0.21	EPA 8270D	5-31-17	6-1-17	
1,2-Dichlorobenzene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2-Methylphenol (o-Cresol)	ND	0.042	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroisopropyl)ether	ND	0.042	EPA 8270D	5-31-17	6-1-17	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.042	EPA 8270D	5-31-17	6-1-17	
n-Nitroso-di-n-propylamine	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Hexachloroethane	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Nitrobenzene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Isophorone	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2-Nitrophenol	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2,4-Dimethylphenol	ND	0.042	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethoxy)methane	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2,4-Dichlorophenol	ND	0.042	EPA 8270D	5-31-17	6-1-17	
1,2,4-Trichlorobenzene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Naphthalene	ND	0.0084	EPA 8270D/SIM	5-31-17	5-31-17	
4-Chloroaniline	ND	0.21	EPA 8270D	5-31-17	6-1-17	
Hexachlorobutadiene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
4-Chloro-3-methylphenol	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2-Methylnaphthalene	ND	0.0084	EPA 8270D/SIM	5-31-17	5-31-17	
1-Methylnaphthalene	ND	0.0084	EPA 8270D/SIM	5-31-17	5-31-17	
Hexachlorocyclopentadiene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2,4,6-Trichlorophenol	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2,3-Dichloroaniline	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2,4,5-Trichlorophenol	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2-Chloronaphthalene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2-Nitroaniline	ND	0.042	EPA 8270D	5-31-17	6-1-17	
1,4-Dinitrobenzene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Dimethylphthalate	ND	0.042	EPA 8270D	5-31-17	6-1-17	
1,3-Dinitrobenzene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
2,6-Dinitrotoluene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
1,2-Dinitrobenzene	ND	0.042	EPA 8270D	5-31-17	6-1-17	
Acenaphthylene	ND	0.0084	EPA 8270D/SIM	5-31-17	5-31-17	
3-Nitroaniline	ND	0.042	EPA 8270D	5-31-17	6-1-17	

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

Analyte Result PQL Method Prepared Analyte Client ID: PP7-10	zed Flags
VIIGILIU. FF/*IV	
Laboratory ID: 05-299-20	
2,4-Dinitrophenol ND 0.21 EPA 8270D 5-31-17 6-1-1	17
Acenaphthene ND 0.0084 EPA 8270D/SIM 5-31-17 5-31-	
4-Nitrophenol ND 0.042 EPA 8270D 5-31-17 6-1-1	
2,4-Dinitrotoluene ND 0.042 EPA 8270D 5-31-17 6-1-1	17
Dibenzofuran ND 0.042 EPA 8270D 5-31-17 6-1-1	17
2,3,5,6-Tetrachlorophenol ND 0.042 EPA 8270D 5-31-17 6-1-1	17
2,3,4,6-Tetrachlorophenol ND 0.042 EPA 8270D 5-31-17 6-1-1	17
Diethylphthalate ND 0.21 EPA 8270D 5-31-17 6-1-1	17
4-Chlorophenyl-phenylether ND 0.042 EPA 8270D 5-31-17 6-1-1	
4-Nitroaniline ND 0.042 EPA 8270D 5-31-17 6-1-1	17
Fluorene ND 0.0084 EPA 8270D/SIM 5-31-17 5-31-	17
4,6-Dinitro-2-methylphenol ND 0.21 EPA 8270D 5-31-17 6-1-1	17
n-Nitrosodiphenylamine ND 0.042 EPA 8270D 5-31-17 6-1-1	
1,2-Diphenylhydrazine ND 0.042 EPA 8270D 5-31-17 6-1-1	
4-Bromophenyl-phenylether ND 0.042 EPA 8270D 5-31-17 6-1-1	17
Hexachlorobenzene ND 0.042 EPA 8270D 5-31-17 6-1-1	
Pentachlorophenol ND 0.21 EPA 8270D 5-31-17 6-1-1	
Phenanthrene ND 0.0084 EPA 8270D/SIM 5-31-17 5-31-	
Anthracene ND 0.0084 EPA 8270D/SIM 5-31-17 5-31-	
Carbazole ND 0.042 EPA 8270D 5-31-17 6-1-1	
Di-n-butylphthalate ND 0.21 EPA 8270D 5-31-17 6-1-1	
Fluoranthene ND 0.0084 EPA 8270D/SIM 5-31-17 5-31-	
Benzidine ND 0.42 EPA 8270D 5-31-17 6-1-1	
Pyrene ND 0.0084 EPA 8270D/SIM 5-31-17 5-31-	
Butylbenzylphthalate ND 0.042 EPA 8270D 5-31-17 6-1-1	
bis-2-Ethylhexyladipate ND 0.042 EPA 8270D 5-31-17 6-1-1	
3,3'-Dichlorobenzidine ND 0.21 EPA 8270D 5-31-17 6-1-1	
Benzo[a]anthracene ND 0.0084 EPA 8270D/SIM 5-31-17 5-31-	17
Chrysene ND 0.0084 EPA 8270D/SIM 5-31-17 5-31-	
bis(2-Ethylhexyl)phthalate ND 0.042 EPA 8270D 5-31-17 6-1-1	17
Di-n-octylphthalate ND 0.042 EPA 8270D 5-31-17 6-1-1	17
Benzo[b]fluoranthene ND 0.0084 EPA 8270D/SIM 5-31-17 5-31-	
Benzo(j,k)fluoranthene ND 0.0084 EPA 8270D/SIM 5-31-17 5-31-	
Benzo[a]pyrene ND 0.0084 EPA 8270D/SIM 5-31-17 5-31-	17
Indeno[1,2,3-cd]pyrene ND 0.0084 EPA 8270D/SIM 5-31-17 5-31-	
Dibenz[a,h]anthracene ND 0.0084 EPA 8270D/SIM 5-31-17 5-31-	
Benzo[g,h,i]perylene ND 0.0084 EPA 8270D/SIM 5-31-17 5-31-	17
Surrogate: Percent Recovery Control Limits	
2-Fluorophenol 71 18 - 109	
Phenol-d6 74 25 - 111	
Nitrobenzene-d5 72 22 - 113	
2-Fluorobiphenyl 70 30 - 114	
2,4,6-Tribromophenol 60 22 - 116	
Terphenyl-d14 64 33 - 114	

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

page 1 of 2

Analyte	Dogult	PQL	Mathad	Date	Date	Elogo
Analyte Client ID:	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP8-0.5 05-299-22					
Laboratory ID:		0.020	EDA 0270D	E 24 47	6 1 17	
n-Nitrosodimethylamine	ND ND	0.038	EPA 8270D	5-31-17	6-1-17	
Pyridine	ND	0.38	EPA 8270D	5-31-17	6-1-17	
Phenol	ND	0.038	EPA 8270D	5-31-17	6-1-17	
Aniline	ND	0.19	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethyl)ether	ND	0.038	EPA 8270D	5-31-17	6-1-17	
2-Chlorophenol	ND	0.038	EPA 8270D	5-31-17	6-1-17	
1,3-Dichlorobenzene	ND	0.038	EPA 8270D	5-31-17	6-1-17	
1,4-Dichlorobenzene	ND	0.038	EPA 8270D	5-31-17	6-1-17	
Benzyl alcohol	0.70	0.19	EPA 8270D	5-31-17	6-1-17	
1,2-Dichlorobenzene	ND	0.038	EPA 8270D	5-31-17	6-1-17	
2-Methylphenol (o-Cresol)	ND	0.038	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroisopropyl)ether	ND	0.038	EPA 8270D	5-31-17	6-1-17	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.038	EPA 8270D	5-31-17	6-1-17	
n-Nitroso-di-n-propylamine	ND	0.038	EPA 8270D	5-31-17	6-1-17	
Hexachloroethane	ND	0.038	EPA 8270D	5-31-17	6-1-17	
Nitrobenzene	ND	0.038	EPA 8270D	5-31-17	6-1-17	
Isophorone	ND	0.038	EPA 8270D	5-31-17	6-1-17	
2-Nitrophenol	ND	0.038	EPA 8270D	5-31-17	6-1-17	
2,4-Dimethylphenol	ND	0.038	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethoxy)methane	ND	0.038	EPA 8270D	5-31-17	6-1-17	
2,4-Dichlorophenol	ND	0.038	EPA 8270D	5-31-17	6-1-17	
1,2,4-Trichlorobenzene	ND	0.038	EPA 8270D	5-31-17	6-1-17	
Naphthalene	ND	0.0076	EPA 8270D/SIM	5-31-17	5-31-17	
4-Chloroaniline	ND	0.19	EPA 8270D	5-31-17	6-1-17	
Hexachlorobutadiene	ND	0.038	EPA 8270D	5-31-17	6-1-17	
4-Chloro-3-methylphenol	ND	0.038	EPA 8270D	5-31-17	6-1-17	
2-Methylnaphthalene	ND	0.0076	EPA 8270D/SIM	5-31-17	5-31-17	
1-Methylnaphthalene	ND	0.0076	EPA 8270D/SIM	5-31-17	5-31-17	
Hexachlorocyclopentadiene	ND	0.038	EPA 8270D	5-31-17	6-1-17	
2,4,6-Trichlorophenol	ND	0.038	EPA 8270D	5-31-17	6-1-17	
2,3-Dichloroaniline	ND	0.038	EPA 8270D	5-31-17	6-1-17	
2,4,5-Trichlorophenol	ND	0.038	EPA 8270D	5-31-17	6-1-17	
2-Chloronaphthalene	ND	0.038	EPA 8270D	5-31-17	6-1-17	
2-Nitroaniline	ND	0.038	EPA 8270D	5-31-17	6-1-17	
1,4-Dinitrobenzene	ND	0.038	EPA 8270D	5-31-17	6-1-17	
Dimethylphthalate	ND	0.038	EPA 8270D	5-31-17	6-1-17	
1,3-Dinitrobenzene	ND	0.038	EPA 8270D	5-31-17	6-1-17	
2,6-Dinitrotoluene	ND	0.038	EPA 8270D	5-31-17	6-1-17	
1,2-Dinitrobenzene	ND	0.038	EPA 8270D	5-31-17	6-1-17	
Acenaphthylene	ND	0.0076	EPA 8270D/SIM	5-31-17	5-31-17	
3-Nitroaniline	ND	0.038	EPA 8270D	5-31-17	6-1-17	

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP8-0.5					110.90
Laboratory ID:	05-299-22					
2,4-Dinitrophenol	ND	0.19	EPA 8270D	5-31-17	6-1-17	
Acenaphthene	ND	0.0076	EPA 8270D/SIM	5-31-17	5-31-17	
4-Nitrophenol	ND	0.038	EPA 8270D	5-31-17	6-1-17	
2,4-Dinitrotoluene	ND	0.038	EPA 8270D	5-31-17	6-1-17	
Dibenzofuran	ND	0.038	EPA 8270D	5-31-17	6-1-17	
2,3,5,6-Tetrachlorophenol	ND	0.038	EPA 8270D	5-31-17	6-1-17	
2,3,4,6-Tetrachlorophenol	ND	0.038	EPA 8270D	5-31-17	6-1-17	
Diethylphthalate	ND	0.19	EPA 8270D	5-31-17	6-1-17	
4-Chlorophenyl-phenylether	ND	0.038	EPA 8270D	5-31-17	6-1-17	
4-Nitroaniline	ND	0.038	EPA 8270D	5-31-17	6-1-17	
Fluorene	ND	0.0076	EPA 8270D/SIM	5-31-17	5-31-17	
4,6-Dinitro-2-methylphenol	ND	0.19	EPA 8270D	5-31-17	6-1-17	
n-Nitrosodiphenylamine	ND	0.038	EPA 8270D	5-31-17	6-1-17	
1,2-Diphenylhydrazine	ND	0.038	EPA 8270D	5-31-17	6-1-17	
4-Bromophenyl-phenylether	ND	0.038	EPA 8270D	5-31-17	6-1-17	
Hexachlorobenzene	ND	0.038	EPA 8270D	5-31-17	6-1-17	
Pentachlorophenol	ND	0.19	EPA 8270D	5-31-17	6-1-17	
Phenanthrene	ND	0.0076	EPA 8270D/SIM	5-31-17	5-31-17	
Anthracene	ND	0.0076	EPA 8270D/SIM	5-31-17	5-31-17	
Carbazole	ND	0.038	EPA 8270D	5-31-17	6-1-17	
Di-n-butylphthalate	ND	0.19	EPA 8270D	5-31-17	6-1-17	
Fluoranthene	ND	0.0076	EPA 8270D/SIM	5-31-17	5-31-17	
Benzidine	ND	0.38	EPA 8270D	5-31-17	6-1-17	
Pyrene	ND	0.0076	EPA 8270D/SIM	5-31-17	5-31-17	
Butylbenzylphthalate	ND	0.038	EPA 8270D	5-31-17	6-1-17	
bis-2-Ethylhexyladipate	ND	0.038	EPA 8270D	5-31-17	6-1-17	
3,3'-Dichlorobenzidine	ND	0.19	EPA 8270D	5-31-17	6-1-17	
Benzo[a]anthracene	ND	0.0076	EPA 8270D/SIM	5-31-17	5-31-17	
Chrysene	ND	0.0076	EPA 8270D/SIM	5-31-17	5-31-17	
bis(2-Ethylhexyl)phthalate	ND	0.038	EPA 8270D	5-31-17	6-1-17	
Di-n-octylphthalate	ND	0.038	EPA 8270D	5-31-17	6-1-17	
Benzo[b]fluoranthene	ND	0.0076	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo(j,k)fluoranthene	ND	0.0076	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[a]pyrene	ND	0.0076	EPA 8270D/SIM	5-31-17	5-31-17	
Indeno[1,2,3-cd]pyrene	ND	0.0076	EPA 8270D/SIM	5-31-17	5-31-17	
Dibenz[a,h]anthracene	ND	0.0076	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[g,h,i]perylene	ND	0.0076	EPA 8270D/SIM	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorophenol	80	18 - 109				
Phenol-d6	85	25 - 111				
Nitrobenzene-d5	85	22 - 113				
2-Fluorobiphenyl	00					
	87	30 - 114				
2,4,6-Tribromophenol						

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

page 1 of 2

Analyte Result PQL Method Prepared Analyzed Flags Client ID: PP8-10 Laboratory ID: 05-299-23 n-Nitrosodimethylamine ND 0.047 EPA 8270D 5-31-17 6-1-17 Pyridine ND 0.47 EPA 8270D 5-31-17 6-1-17 Phenol ND 0.047 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.24 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.047 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.047 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.047 EPA 8270D 5-31-17 6-1-17					Date	Date	
Laboratory ID: 05-299-23 n-Nitrosodimethylamine ND 0.047 EPA 8270D 5-31-17 6-1-17 Pyridine ND 0.47 EPA 8270D 5-31-17 6-1-17 Phenol ND 0.047 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.24 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.047 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.047 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.047 EPA 8270D 5-31-17 6-1-17	Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
n-Nitrosodimethylamine ND 0.047 EPA 8270D 5-31-17 6-1-17 Pyridine ND 0.47 EPA 8270D 5-31-17 6-1-17 Phenol ND 0.047 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.24 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.047 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.047 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.047 EPA 8270D 5-31-17 6-1-17	Client ID:	PP8-10					
Pyridine ND 0.47 EPA 8270D 5-31-17 6-1-17 Phenol ND 0.047 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.24 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.047 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.047 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.047 EPA 8270D 5-31-17 6-1-17	Laboratory ID:	05-299-23					
Pyridine ND 0.47 EPA 8270D 5-31-17 6-1-17 Phenol ND 0.047 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.24 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.047 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.047 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.047 EPA 8270D 5-31-17 6-1-17		ND	0.047	EPA 8270D	5-31-17	6-1-17	
Aniline ND 0.24 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.047 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.047 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.047 EPA 8270D 5-31-17 6-1-17		ND	0.47	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethyl)ether ND 0.047 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.047 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.047 EPA 8270D 5-31-17 6-1-17	_	ND	0.047	EPA 8270D	5-31-17	6-1-17	
2-Chlorophenol ND 0.047 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.047 EPA 8270D 5-31-17 6-1-17	Aniline	ND	0.24	EPA 8270D	5-31-17	6-1-17	
1,3-Dichlorobenzene ND 0.047 EPA 8270D 5-31-17 6-1-17	bis(2-Chloroethyl)ether	ND	0.047	EPA 8270D	5-31-17	6-1-17	
1,3-Dichlorobenzene ND 0.047 EPA 8270D 5-31-17 6-1-17	2-Chlorophenol	ND	0.047	EPA 8270D	5-31-17	6-1-17	
· · · · · · · · · · · · · · · · · · ·	1,3-Dichlorobenzene	ND	0.047	EPA 8270D	5-31-17	6-1-17	
1,4-Dichlorobenzene ND 0.047 EPA 8270D 5-31-17 6-1-17	1,4-Dichlorobenzene	ND	0.047	EPA 8270D	5-31-17	6-1-17	
Benzyl alcohol ND 0.24 EPA 8270D 5-31-17 6-1-17	Benzyl alcohol	ND	0.24	EPA 8270D	5-31-17	6-1-17	
1,2-Dichlorobenzene ND 0.047 EPA 8270D 5-31-17 6-1-17		ND	0.047	EPA 8270D	5-31-17	6-1-17	
2-Methylphenol (o-Cresol) ND 0.047 EPA 8270D 5-31-17 6-1-17	2-Methylphenol (o-Cresol)	ND	0.047	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroisopropyl)ether ND 0.047 EPA 8270D 5-31-17 6-1-17		ND		EPA 8270D	5-31-17	6-1-17	
(3+4)-Methylphenol (m,p-Cresol) ND 0.047 EPA 8270D 5-31-17 6-1-17	(3+4)-Methylphenol (m,p-Cresol)	ND	0.047	EPA 8270D	5-31-17	6-1-17	
n-Nitroso-di-n-propylamine ND 0.047 EPA 8270D 5-31-17 6-1-17		ND	0.047	EPA 8270D	5-31-17	6-1-17	
Hexachloroethane ND 0.047 EPA 8270D 5-31-17 6-1-17	Hexachloroethane	ND	0.047	EPA 8270D	5-31-17	6-1-17	
Nitrobenzene ND 0.047 EPA 8270D 5-31-17 6-1-17	Nitrobenzene	ND	0.047	EPA 8270D	5-31-17	6-1-17	
Isophorone ND 0.047 EPA 8270D 5-31-17 6-1-17	Isophorone	ND	0.047	EPA 8270D	5-31-17	6-1-17	
2-Nitrophenol ND 0.047 EPA 8270D 5-31-17 6-1-17		ND	0.047	EPA 8270D	5-31-17	6-1-17	
2,4-Dimethylphenol ND 0.047 EPA 8270D 5-31-17 6-1-17	2,4-Dimethylphenol	ND	0.047	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethoxy)methane ND 0.047 EPA 8270D 5-31-17 6-1-17	bis(2-Chloroethoxy)methane	ND	0.047	EPA 8270D	5-31-17	6-1-17	
2,4-Dichlorophenol ND 0.047 EPA 8270D 5-31-17 6-1-17		ND	0.047	EPA 8270D	5-31-17	6-1-17	
1,2,4-Trichlorobenzene ND 0.047 EPA 8270D 5-31-17 6-1-17	1,2,4-Trichlorobenzene	ND	0.047	EPA 8270D	5-31-17	6-1-17	
Naphthalene ND 0.0095 EPA 8270D/SIM 5-31-17 5-31-17	Naphthalene	ND	0.0095	EPA 8270D/SIM	5-31-17	5-31-17	
4-Chloroaniline ND 0.24 EPA 8270D 5-31-17 6-1-17	4-Chloroaniline	ND	0.24	EPA 8270D	5-31-17	6-1-17	
Hexachlorobutadiene ND 0.047 EPA 8270D 5-31-17 6-1-17	Hexachlorobutadiene	ND	0.047	EPA 8270D	5-31-17	6-1-17	
4-Chloro-3-methylphenol ND 0.047 EPA 8270D 5-31-17 6-1-17	4-Chloro-3-methylphenol	ND	0.047	EPA 8270D	5-31-17	6-1-17	
2-Methylnaphthalene ND 0.0095 EPA 8270D/SIM 5-31-17 5-31-17	2-Methylnaphthalene	ND	0.0095	EPA 8270D/SIM	5-31-17	5-31-17	
1-Methylnaphthalene ND 0.0095 EPA 8270D/SIM 5-31-17 5-31-17	1-Methylnaphthalene	ND	0.0095	EPA 8270D/SIM	5-31-17	5-31-17	
Hexachlorocyclopentadiene ND 0.047 EPA 8270D 5-31-17 6-1-17		ND	0.047	EPA 8270D	5-31-17	6-1-17	
2,4,6-Trichlorophenol ND 0.047 EPA 8270D 5-31-17 6-1-17	2,4,6-Trichlorophenol	ND	0.047	EPA 8270D	5-31-17	6-1-17	
2,3-Dichloroaniline ND 0.047 EPA 8270D 5-31-17 6-1-17	2,3-Dichloroaniline	ND	0.047	EPA 8270D	5-31-17	6-1-17	
2,4,5-Trichlorophenol ND 0.047 EPA 8270D 5-31-17 6-1-17	2,4,5-Trichlorophenol	ND	0.047	EPA 8270D	5-31-17	6-1-17	
2-Chloronaphthalene ND 0.047 EPA 8270D 5-31-17 6-1-17	2-Chloronaphthalene	ND	0.047	EPA 8270D	5-31-17	6-1-17	
2-Nitroaniline ND 0.047 EPA 8270D 5-31-17 6-1-17	2-Nitroaniline	ND	0.047	EPA 8270D	5-31-17	6-1-17	
1,4-Dinitrobenzene ND 0.047 EPA 8270D 5-31-17 6-1-17	1,4-Dinitrobenzene	ND	0.047	EPA 8270D	5-31-17	6-1-17	
Dimethylphthalate ND 0.047 EPA 8270D 5-31-17 6-1-17	Dimethylphthalate	ND	0.047	EPA 8270D	5-31-17	6-1-17	
1,3-Dinitrobenzene ND 0.047 EPA 8270D 5-31-17 6-1-17	1,3-Dinitrobenzene	ND	0.047	EPA 8270D	5-31-17	6-1-17	
2,6-Dinitrotoluene ND 0.047 EPA 8270D 5-31-17 6-1-17	2,6-Dinitrotoluene	ND	0.047	EPA 8270D	5-31-17	6-1-17	
1,2-Dinitrobenzene ND 0.047 EPA 8270D 5-31-17 6-1-17	1,2-Dinitrobenzene	ND	0.047	EPA 8270D	5-31-17	6-1-17	
Acenaphthylene ND 0.0095 EPA 8270D/SIM 5-31-17 5-31-17	Acenaphthylene	ND	0.0095	EPA 8270D/SIM	5-31-17	5-31-17	
3-Nitroaniline ND 0.047 EPA 8270D 5-31-17 6-1-17	3-Nitroaniline	ND	0.047	EPA 8270D	5-31-17	6-1-17	

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP8-10					110.90
Laboratory ID:	05-299-23					
2,4-Dinitrophenol	ND	0.24	EPA 8270D	5-31-17	6-1-17	
Acenaphthene	ND	0.0095	EPA 8270D/SIM	5-31-17	5-31-17	
4-Nitrophenol	ND	0.047	EPA 8270D	5-31-17	6-1-17	
2,4-Dinitrotoluene	ND	0.047	EPA 8270D	5-31-17	6-1-17	
Dibenzofuran	ND	0.047	EPA 8270D	5-31-17	6-1-17	
2,3,5,6-Tetrachlorophenol	ND	0.047	EPA 8270D	5-31-17	6-1-17	
2,3,4,6-Tetrachlorophenol	ND	0.047	EPA 8270D	5-31-17	6-1-17	
Diethylphthalate	ND	0.24	EPA 8270D	5-31-17	6-1-17	
4-Chlorophenyl-phenylether	ND	0.047	EPA 8270D	5-31-17	6-1-17	
4-Nitroaniline	ND	0.047	EPA 8270D	5-31-17	6-1-17	
Fluorene	ND	0.0095	EPA 8270D/SIM	5-31-17	5-31-17	
4,6-Dinitro-2-methylphenol	ND	0.24	EPA 8270D	5-31-17	6-1-17	
n-Nitrosodiphenylamine	ND	0.047	EPA 8270D	5-31-17	6-1-17	
1,2-Diphenylhydrazine	ND	0.047	EPA 8270D	5-31-17	6-1-17	
4-Bromophenyl-phenylether	ND	0.047	EPA 8270D	5-31-17	6-1-17	
Hexachlorobenzene	ND	0.047	EPA 8270D	5-31-17	6-1-17	
Pentachlorophenol	ND	0.24	EPA 8270D	5-31-17	6-1-17	
Phenanthrene	ND	0.0095	EPA 8270D/SIM	5-31-17	5-31-17	
Anthracene	ND	0.0095	EPA 8270D/SIM	5-31-17	5-31-17	
Carbazole	ND	0.047	EPA 8270D	5-31-17	6-1-17	
Di-n-butylphthalate	ND	0.24	EPA 8270D	5-31-17	6-1-17	
Fluoranthene	ND	0.0095	EPA 8270D/SIM	5-31-17	5-31-17	
Benzidine	ND	0.47	EPA 8270D	5-31-17	6-1-17	
Pyrene	ND	0.0095	EPA 8270D/SIM	5-31-17	5-31-17	
Butylbenzylphthalate	ND	0.047	EPA 8270D	5-31-17	6-1-17	
bis-2-Ethylhexyladipate	ND	0.047	EPA 8270D	5-31-17	6-1-17	
3,3'-Dichlorobenzidine	ND	0.24	EPA 8270D	5-31-17	6-1-17	
Benzo[a]anthracene	ND	0.0095	EPA 8270D/SIM	5-31-17	5-31-17	
Chrysene	ND	0.0095	EPA 8270D/SIM	5-31-17	5-31-17	
bis(2-Ethylhexyl)phthalate	ND	0.047	EPA 8270D	5-31-17	6-1-17	
Di-n-octylphthalate	ND	0.047	EPA 8270D	5-31-17	6-1-17	
Benzo[b]fluoranthene	ND	0.0095	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo(j,k)fluoranthene	ND	0.0095	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[a]pyrene	ND	0.0095	EPA 8270D/SIM	5-31-17	5-31-17	
Indeno[1,2,3-cd]pyrene	ND	0.0095	EPA 8270D/SIM	5-31-17	5-31-17	
Dibenz[a,h]anthracene	ND	0.0095	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[g,h,i]perylene	ND	0.0095	EPA 8270D/SIM	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorophenol	76	18 - 109				
Phenol-d6	73	25 - 111				
Nitrobenzene-d5	82	22 - 113				
2-Fluorobiphenyl						
		30 - 114				
2,4,6-Tribromophenol	78 70	30 - 114 22 - 116				

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

page 1 of 2

	-		•• 41 .	Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP9-0					
Laboratory ID:	05-299-25	0.005	EDA 0070D	E 04 47	0.4.47	
n-Nitrosodimethylamine	ND	0.035	EPA 8270D	5-31-17	6-1-17	
Pyridine	ND	0.35	EPA 8270D	5-31-17	6-1-17	
Phenol	ND	0.035	EPA 8270D	5-31-17	6-1-17	
Aniline	ND	0.17	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethyl)ether	ND	0.035	EPA 8270D	5-31-17	6-1-17	
2-Chlorophenol	ND	0.035	EPA 8270D	5-31-17	6-1-17	
1,3-Dichlorobenzene	ND	0.035	EPA 8270D	5-31-17	6-1-17	
1,4-Dichlorobenzene	ND	0.035	EPA 8270D	5-31-17	6-1-17	
Benzyl alcohol	ND	0.17	EPA 8270D	5-31-17	6-1-17	
1,2-Dichlorobenzene	ND	0.035	EPA 8270D	5-31-17	6-1-17	
2-Methylphenol (o-Cresol)	ND	0.035	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroisopropyl)ether	ND	0.035	EPA 8270D	5-31-17	6-1-17	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.035	EPA 8270D	5-31-17	6-1-17	
n-Nitroso-di-n-propylamine	ND	0.035	EPA 8270D	5-31-17	6-1-17	
Hexachloroethane	ND	0.035	EPA 8270D	5-31-17	6-1-17	
Nitrobenzene	ND	0.035	EPA 8270D	5-31-17	6-1-17	
Isophorone	ND	0.035	EPA 8270D	5-31-17	6-1-17	
2-Nitrophenol	ND	0.035	EPA 8270D	5-31-17	6-1-17	
2,4-Dimethylphenol	ND	0.035	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethoxy)methane	ND	0.035	EPA 8270D	5-31-17	6-1-17	
2,4-Dichlorophenol	ND	0.035	EPA 8270D	5-31-17	6-1-17	
1,2,4-Trichlorobenzene	ND	0.035	EPA 8270D	5-31-17	6-1-17	
Naphthalene	ND	0.0070	EPA 8270D/SIM	5-31-17	5-31-17	
4-Chloroaniline	ND	0.17	EPA 8270D	5-31-17	6-1-17	
Hexachlorobutadiene	ND	0.035	EPA 8270D	5-31-17	6-1-17	
4-Chloro-3-methylphenol	ND	0.035	EPA 8270D	5-31-17	6-1-17	
2-Methylnaphthalene	ND	0.0070	EPA 8270D/SIM	5-31-17	5-31-17	
1-Methylnaphthalene	ND	0.0070	EPA 8270D/SIM	5-31-17	5-31-17	
Hexachlorocyclopentadiene	ND	0.035	EPA 8270D	5-31-17	6-1-17	
2,4,6-Trichlorophenol	ND	0.035	EPA 8270D	5-31-17	6-1-17	
2,3-Dichloroaniline	ND	0.035	EPA 8270D	5-31-17	6-1-17	
2,4,5-Trichlorophenol	ND	0.035	EPA 8270D	5-31-17	6-1-17	
2-Chloronaphthalene	ND	0.035	EPA 8270D	5-31-17	6-1-17	
2-Nitroaniline	ND	0.035	EPA 8270D	5-31-17	6-1-17	
1,4-Dinitrobenzene	ND	0.035	EPA 8270D	5-31-17	6-1-17	
Dimethylphthalate	ND	0.035	EPA 8270D	5-31-17	6-1-17	
1,3-Dinitrobenzene	ND	0.035	EPA 8270D	5-31-17	6-1-17	
2,6-Dinitrotoluene	ND	0.035	EPA 8270D	5-31-17	6-1-17	
1,2-Dinitrobenzene	ND	0.035	EPA 8270D	5-31-17	6-1-17	
Acenaphthylene	ND	0.0070	EPA 8270D/SIM	5-31-17	5-31-17	
3-Nitroaniline	ND	0.035	EPA 8270D	5-31-17	6-1-17	

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP9-0					- - -
Laboratory ID:	05-299-25					
2,4-Dinitrophenol	ND	0.17	EPA 8270D	5-31-17	6-1-17	
Acenaphthene	ND	0.0070	EPA 8270D/SIM	5-31-17	5-31-17	
4-Nitrophenol	ND	0.035	EPA 8270D	5-31-17	6-1-17	
2,4-Dinitrotoluene	ND	0.035	EPA 8270D	5-31-17	6-1-17	
Dibenzofuran	ND	0.035	EPA 8270D	5-31-17	6-1-17	
2,3,5,6-Tetrachlorophenol	ND	0.035	EPA 8270D	5-31-17	6-1-17	
2,3,4,6-Tetrachlorophenol	ND	0.035	EPA 8270D	5-31-17	6-1-17	
Diethylphthalate	ND	0.17	EPA 8270D	5-31-17	6-1-17	
4-Chlorophenyl-phenylether	ND	0.035	EPA 8270D	5-31-17	6-1-17	
4-Nitroaniline	ND	0.035	EPA 8270D	5-31-17	6-1-17	
Fluorene	ND	0.0070	EPA 8270D/SIM	5-31-17	5-31-17	
4,6-Dinitro-2-methylphenol	ND	0.17	EPA 8270D	5-31-17	6-1-17	
n-Nitrosodiphenylamine	ND	0.035	EPA 8270D	5-31-17	6-1-17	
1,2-Diphenylhydrazine	ND	0.035	EPA 8270D	5-31-17	6-1-17	
4-Bromophenyl-phenylether	ND	0.035	EPA 8270D	5-31-17	6-1-17	
Hexachlorobenzene	ND	0.035	EPA 8270D	5-31-17	6-1-17	
Pentachlorophenol	ND	0.17	EPA 8270D	5-31-17	6-1-17	
Phenanthrene	ND	0.0070	EPA 8270D/SIM	5-31-17	5-31-17	
Anthracene	ND	0.0070	EPA 8270D/SIM	5-31-17	5-31-17	
Carbazole	ND	0.035	EPA 8270D	5-31-17	6-1-17	
Di-n-butylphthalate	ND	0.17	EPA 8270D	5-31-17	6-1-17	
Fluoranthene	ND	0.0070	EPA 8270D/SIM	5-31-17	5-31-17	
Benzidine	ND	0.35	EPA 8270D	5-31-17	6-1-17	
Pyrene	ND	0.0070	EPA 8270D/SIM	5-31-17	5-31-17	
Butylbenzylphthalate	ND	0.035	EPA 8270D	5-31-17	6-1-17	
bis-2-Ethylhexyladipate	ND	0.035	EPA 8270D	5-31-17	6-1-17	
3,3'-Dichlorobenzidine	ND	0.17	EPA 8270D	5-31-17	6-1-17	
Benzo[a]anthracene	ND	0.0070	EPA 8270D/SIM	5-31-17	5-31-17	
Chrysene	ND	0.0070	EPA 8270D/SIM	5-31-17	5-31-17	
bis(2-Ethylhexyl)phthalate	ND	0.035	EPA 8270D	5-31-17	6-1-17	
Di-n-octylphthalate	ND	0.13	EPA 8270D	5-31-17	6-1-17	U1
Benzo[b]fluoranthene	ND	0.0070	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo(j,k)fluoranthene	ND	0.0070	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[a]pyrene	ND	0.0070	EPA 8270D/SIM	5-31-17	5-31-17	
Indeno[1,2,3-cd]pyrene	ND	0.0070	EPA 8270D/SIM	5-31-17	5-31-17	
Dibenz[a,h]anthracene	ND	0.0070	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[g,h,i]perylene	ND	0.0070	EPA 8270D/SIM	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits			0 0 1 11	
2-Fluorophenol	56	18 - 109				
Phenol-d6	63	25 - 111				
Nitrobenzene-d5	62	22 - 113				
2-Fluorobiphenyl	66	30 - 114				
2,4,6-Tribromophenol	69	22 - 116				
Terphenyl-d14	69	33 - 114				
Torphonyr a 1+	00	00 117				

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

page 1 of 2

Analyte Result PQL Method Prepared Analyzed Flags Client ID: PP9-10 Laboratory ID: 05-299-26					Date	Date	
Laboratory ID: 05-299-26	Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
n-Nitrosodimethylamine ND 0.041 EPA 8270D 5-31-17 6-1-17 Pyridine ND 0.41 EPA 8270D 5-31-17 6-1-17 Phenol ND 0.041 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.21 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.041 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.041 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.041 EPA 8270D 5-31-17 6-1-17 c-Chlorophenol ND 0.041 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.041 EPA 8270D 5-31-17 6-1-	Client ID:	PP9-10					
Pyridine ND 0.41 EPA 8270D 5-31-17 6-1-17 Phenol ND 0.041 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.21 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.041 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.041 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 Benzyl alcohol ND 0.041 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.041 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroisopropyl)ether ND 0.041 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.041 EPA 8270D 5-31-17 </td <td>Laboratory ID:</td> <td>05-299-26</td> <td></td> <td></td> <td></td> <td></td> <td></td>	Laboratory ID:	05-299-26					
Pyridine ND 0.41 EPA 8270D 5-31-17 6-1-17 Phenol ND 0.041 EPA 8270D 5-31-17 6-1-17 Aniline ND 0.21 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.041 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.041 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 Benzyl alcohol ND 0.041 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.041 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroisopropyl)ether ND 0.041 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.041 EPA 8270D 5-31-17 </td <td>n-Nitrosodimethylamine</td> <td>ND</td> <td>0.041</td> <td>EPA 8270D</td> <td>5-31-17</td> <td>6-1-17</td> <td></td>	n-Nitrosodimethylamine	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Aniline ND 0.21 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroethyl)ether ND 0.041 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.041 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 Benzyl alcohol ND 0.21 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.041 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroisopropyl)ether ND 0.041 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.041 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.041 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.041 EPA 8270D		ND	0.41	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethyl)ether ND 0.041 EPA 8270D 5-31-17 6-1-17 2-Chlorophenol ND 0.041 EPA 8270D 5-31-17 6-1-17 1,3-Dichlorobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 Benzyl alcohol ND 0.21 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.041 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroisopropyl)ether ND 0.041 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.041 EPA 8270D 5-31-17 6-1-17 n-Nitroso-di-n-propylamine ND 0.041 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.041 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.041 EP	Phenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-ChlorophenolND0.041EPA 8270D5-31-176-1-171,3-DichlorobenzeneND0.041EPA 8270D5-31-176-1-171,4-DichlorobenzeneND0.041EPA 8270D5-31-176-1-17Benzyl alcoholND0.21EPA 8270D5-31-176-1-171,2-DichlorobenzeneND0.041EPA 8270D5-31-176-1-172-Methylphenol (o-Cresol)ND0.041EPA 8270D5-31-176-1-17bis(2-Chloroisopropyl)etherND0.041EPA 8270D5-31-176-1-17(3+4)-Methylphenol (m,p-Cresol)ND0.041EPA 8270D5-31-176-1-17n-Nitroso-di-n-propylamineND0.041EPA 8270D5-31-176-1-17HexachloroethaneND0.041EPA 8270D5-31-176-1-17NitrobenzeneND0.041EPA 8270D5-31-176-1-17IsophoroneND0.041EPA 8270D5-31-176-1-17	Aniline	ND	0.21	EPA 8270D	5-31-17	6-1-17	
1,3-Dichlorobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 1,4-Dichlorobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 Benzyl alcohol ND 0.21 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.041 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroisopropyl)ether ND 0.041 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.041 EPA 8270D 5-31-17 6-1-17 n-Nitroso-di-n-propylamine ND 0.041 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.041 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.041 EPA 8270D 5-31-17 6-1-17	bis(2-Chloroethyl)ether	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,4-Dichlorobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 Benzyl alcohol ND 0.21 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.041 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroisopropyl)ether ND 0.041 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.041 EPA 8270D 5-31-17 6-1-17 n-Nitroso-di-n-propylamine ND 0.041 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.041 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.041 EPA 8270D 5-31-17 6-1-17	2-Chlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Benzyl alcohol ND 0.21 EPA 8270D 5-31-17 6-1-17 1,2-Dichlorobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.041 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroisopropyl)ether ND 0.041 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.041 EPA 8270D 5-31-17 6-1-17 n-Nitroso-di-n-propylamine ND 0.041 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.041 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.041 EPA 8270D 5-31-17 6-1-17	1,3-Dichlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,2-Dichlorobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 2-Methylphenol (o-Cresol) ND 0.041 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroisopropyl)ether ND 0.041 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.041 EPA 8270D 5-31-17 6-1-17 n-Nitroso-di-n-propylamine ND 0.041 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.041 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.041 EPA 8270D 5-31-17 6-1-17	1,4-Dichlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Methylphenol (o-Cresol) ND 0.041 EPA 8270D 5-31-17 6-1-17 bis(2-Chloroisopropyl)ether ND 0.041 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.041 EPA 8270D 5-31-17 6-1-17 n-Nitroso-di-n-propylamine ND 0.041 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.041 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.041 EPA 8270D 5-31-17 6-1-17	Benzyl alcohol	ND	0.21	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroisopropyl)ether ND 0.041 EPA 8270D 5-31-17 6-1-17 (3+4)-Methylphenol (m,p-Cresol) ND 0.041 EPA 8270D 5-31-17 6-1-17 n-Nitroso-di-n-propylamine ND 0.041 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.041 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.041 EPA 8270D 5-31-17 6-1-17	1,2-Dichlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
(3+4)-Methylphenol (m,p-Cresol) ND 0.041 EPA 8270D 5-31-17 6-1-17 n-Nitroso-di-n-propylamine ND 0.041 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.041 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.041 EPA 8270D 5-31-17 6-1-17	2-Methylphenol (o-Cresol)	ND	0.041	EPA 8270D	5-31-17	6-1-17	
n-Nitroso-di-n-propylamine ND 0.041 EPA 8270D 5-31-17 6-1-17 Hexachloroethane ND 0.041 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.041 EPA 8270D 5-31-17 6-1-17	bis(2-Chloroisopropyl)ether	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Hexachloroethane ND 0.041 EPA 8270D 5-31-17 6-1-17 Nitrobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.041 EPA 8270D 5-31-17 6-1-17	(3+4)-Methylphenol (m,p-Cresol)	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Nitrobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17 Isophorone ND 0.041 EPA 8270D 5-31-17 6-1-17	n-Nitroso-di-n-propylamine	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Isophorone ND 0.041 EPA 8270D 5-31-17 6-1-17	Hexachloroethane	ND	0.041	EPA 8270D	5-31-17	6-1-17	
·	Nitrobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2 Nitrophenal ND 0.041 EDA 8270D 5-31.17 6.1-17	Isophorone	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Nill Ophichol ND 0.041 LIA 0270D 3-31-17 0-1-17	2-Nitrophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4-Dimethylphenol ND 0.041 EPA 8270D 5-31-17 6-1-17	2,4-Dimethylphenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethoxy)methane ND 0.041 EPA 8270D 5-31-17 6-1-17	bis(2-Chloroethoxy)methane	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4-Dichlorophenol ND 0.041 EPA 8270D 5-31-17 6-1-17	2,4-Dichlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,2,4-Trichlorobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17	1,2,4-Trichlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Naphthalene ND 0.0083 EPA 8270D/SIM 5-31-17 5-31-17	Naphthalene	ND	0.0083	EPA 8270D/SIM	5-31-17	5-31-17	
4-Chloroaniline ND 0.21 EPA 8270D 5-31-17 6-1-17	4-Chloroaniline	ND	0.21	EPA 8270D	5-31-17	6-1-17	
Hexachlorobutadiene ND 0.041 EPA 8270D 5-31-17 6-1-17	Hexachlorobutadiene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
4-Chloro-3-methylphenol ND 0.041 EPA 8270D 5-31-17 6-1-17	4-Chloro-3-methylphenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Methylnaphthalene ND 0.0083 EPA 8270D/SIM 5-31-17 5-31-17	2-Methylnaphthalene	ND	0.0083	EPA 8270D/SIM	5-31-17	5-31-17	
1-Methylnaphthalene ND 0.0083 EPA 8270D/SIM 5-31-17 5-31-17	1-Methylnaphthalene	ND	0.0083	EPA 8270D/SIM	5-31-17	5-31-17	
Hexachlorocyclopentadiene ND 0.041 EPA 8270D 5-31-17 6-1-17		ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4,6-Trichlorophenol ND 0.041 EPA 8270D 5-31-17 6-1-17	2,4,6-Trichlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,3-Dichloroaniline ND 0.041 EPA 8270D 5-31-17 6-1-17	2,3-Dichloroaniline	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4,5-Trichlorophenol ND 0.041 EPA 8270D 5-31-17 6-1-17	2,4,5-Trichlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Chloronaphthalene ND 0.041 EPA 8270D 5-31-17 6-1-17	2-Chloronaphthalene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Nitroaniline ND 0.041 EPA 8270D 5-31-17 6-1-17	2-Nitroaniline	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,4-Dinitrobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17	1,4-Dinitrobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Dimethylphthalate ND 0.041 EPA 8270D 5-31-17 6-1-17	Dimethylphthalate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,3-Dinitrobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17	1,3-Dinitrobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,6-Dinitrotoluene ND 0.041 EPA 8270D 5-31-17 6-1-17	2,6-Dinitrotoluene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,2-Dinitrobenzene ND 0.041 EPA 8270D 5-31-17 6-1-17	1,2-Dinitrobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Acenaphthylene ND 0.0083 EPA 8270D/SIM 5-31-17 5-31-17	Acenaphthylene	ND	0.0083	EPA 8270D/SIM	5-31-17	5-31-17	
3-Nitroaniline ND 0.041 EPA 8270D 5-31-17 6-1-17	3-Nitroaniline	ND	0.041	EPA 8270D	5-31-17	6-1-17	

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP9-10					
Laboratory ID:	05-299-26					
2,4-Dinitrophenol	ND	0.21	EPA 8270D	5-31-17	6-1-17	
Acenaphthene	ND	0.0083	EPA 8270D/SIM	5-31-17	5-31-17	
4-Nitrophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4-Dinitrotoluene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Dibenzofuran	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,3,5,6-Tetrachlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,3,4,6-Tetrachlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Diethylphthalate	ND	0.21	EPA 8270D	5-31-17	6-1-17	
4-Chlorophenyl-phenylether	ND	0.041	EPA 8270D	5-31-17	6-1-17	
4-Nitroaniline	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Fluorene	ND	0.0083	EPA 8270D/SIM	5-31-17	5-31-17	
4,6-Dinitro-2-methylphenol	ND	0.21	EPA 8270D	5-31-17	6-1-17	
n-Nitrosodiphenylamine	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,2-Diphenylhydrazine	ND	0.041	EPA 8270D	5-31-17	6-1-17	
4-Bromophenyl-phenylether	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Hexachlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Pentachlorophenol	ND	0.21	EPA 8270D	5-31-17	6-1-17	
Phenanthrene	ND	0.0083	EPA 8270D/SIM	5-31-17	5-31-17	
Anthracene	ND	0.0083	EPA 8270D/SIM	5-31-17	5-31-17	
Carbazole	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Di-n-butylphthalate	ND	0.21	EPA 8270D	5-31-17	6-1-17	
Fluoranthene	ND	0.0083	EPA 8270D/SIM	5-31-17	5-31-17	
Benzidine	ND	0.41	EPA 8270D	5-31-17	6-1-17	
Pyrene	ND	0.0083	EPA 8270D/SIM	5-31-17	5-31-17	
Butylbenzylphthalate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
bis-2-Ethylhexyladipate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
3,3'-Dichlorobenzidine	ND	0.21	EPA 8270D	5-31-17	6-1-17	
Benzo[a]anthracene	ND	0.0083	EPA 8270D/SIM	5-31-17	5-31-17	
Chrysene	ND	0.0083	EPA 8270D/SIM	5-31-17	5-31-17	
bis(2-Ethylhexyl)phthalate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Di-n-octylphthalate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Benzo[b]fluoranthene	ND	0.0083	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo(j,k)fluoranthene	ND	0.0083	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[a]pyrene	ND	0.0083	EPA 8270D/SIM	5-31-17	5-31-17	
Indeno[1,2,3-cd]pyrene	ND	0.0083	EPA 8270D/SIM	5-31-17	5-31-17	
Dibenz[a,h]anthracene	ND	0.0083	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[g,h,i]perylene	ND	0.0083	EPA 8270D/SIM	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits		· · · ·	0 0 1 11	
2-Fluorophenol	79	18 - 109				
Phenol-d6	78	25 - 111				
Nitrobenzene-d5	80	22 - 113				
2-Fluorobiphenyl	71	30 - 114				
2,4,6-Tribromophenol	67	22 - 116				
Terphenyl-d14	65	33 - 114				
· 1	30					

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP10-0					
Laboratory ID:	05-299-28					
n-Nitrosodimethylamine	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Pyridine	ND	0.41	EPA 8270D	5-31-17	6-1-17	
Phenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Aniline	ND	0.20	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethyl)ether	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Chlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,3-Dichlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,4-Dichlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Benzyl alcohol	ND	0.20	EPA 8270D	5-31-17	6-1-17	
1,2-Dichlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Methylphenol (o-Cresol)	ND	0.041	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroisopropyl)ether	ND	0.041	EPA 8270D	5-31-17	6-1-17	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.041	EPA 8270D	5-31-17	6-1-17	
n-Nitroso-di-n-propylamine	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Hexachloroethane	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Nitrobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Isophorone	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Nitrophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4-Dimethylphenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
bis(2-Chloroethoxy)methane	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4-Dichlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,2,4-Trichlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Naphthalene	ND	0.0081	EPA 8270D/SIM	5-31-17	6-1-17	
4-Chloroaniline	ND	0.20	EPA 8270D	5-31-17	6-1-17	
Hexachlorobutadiene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
4-Chloro-3-methylphenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Methylnaphthalene	ND	0.0081	EPA 8270D/SIM	5-31-17	6-1-17	
1-Methylnaphthalene	ND	0.0081	EPA 8270D/SIM	5-31-17	6-1-17	
Hexachlorocyclopentadiene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4,6-Trichlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,3-Dichloroaniline	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4,5-Trichlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Chloronaphthalene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2-Nitroaniline	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,4-Dinitrobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Dimethylphthalate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,3-Dinitrobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,6-Dinitrotoluene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,2-Dinitrobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Acenaphthylene	ND	0.0081	EPA 8270D/SIM	5-31-17	6-1-17	
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Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP10-0					
Laboratory ID:	05-299-28					
2,4-Dinitrophenol	ND	0.20	EPA 8270D	5-31-17	6-1-17	
Acenaphthene	ND	0.0081	EPA 8270D/SIM	5-31-17	6-1-17	
4-Nitrophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,4-Dinitrotoluene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Dibenzofuran	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,3,5,6-Tetrachlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
2,3,4,6-Tetrachlorophenol	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Diethylphthalate	ND	0.20	EPA 8270D	5-31-17	6-1-17	
4-Chlorophenyl-phenylether	ND	0.041	EPA 8270D	5-31-17	6-1-17	
4-Nitroaniline	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Fluorene	ND	0.0081	EPA 8270D/SIM	5-31-17	6-1-17	
4,6-Dinitro-2-methylphenol	ND	0.20	EPA 8270D	5-31-17	6-1-17	
n-Nitrosodiphenylamine	ND	0.041	EPA 8270D	5-31-17	6-1-17	
1,2-Diphenylhydrazine	ND	0.041	EPA 8270D	5-31-17	6-1-17	
4-Bromophenyl-phenylether	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Hexachlorobenzene	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Pentachlorophenol	ND	0.20	EPA 8270D	5-31-17	6-1-17	
Phenanthrene	ND	0.0081	EPA 8270D/SIM	5-31-17	6-1-17	
Anthracene	ND	0.0081	EPA 8270D/SIM	5-31-17	6-1-17	
Carbazole	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Di-n-butylphthalate	ND	0.20	EPA 8270D	5-31-17	6-1-17	
Fluoranthene	ND	0.0081	EPA 8270D/SIM	5-31-17	6-1-17	
Benzidine	ND	0.41	EPA 8270D	5-31-17	6-1-17	
Pyrene	ND	0.0081	EPA 8270D/SIM	5-31-17	6-1-17	
Butylbenzylphthalate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
bis-2-Ethylhexyladipate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
3,3'-Dichlorobenzidine	ND	0.20	EPA 8270D	5-31-17	6-1-17	
Benzo[a]anthracene	ND	0.0081	EPA 8270D/SIM	5-31-17	6-1-17	
Chrysene	ND	0.0081	EPA 8270D/SIM	5-31-17	6-1-17	
bis(2-Ethylhexyl)phthalate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Di-n-octylphthalate	ND	0.041	EPA 8270D	5-31-17	6-1-17	
Benzo[b]fluoranthene	ND	0.0081	EPA 8270D/SIM	5-31-17	6-1-17	
Benzo(j,k)fluoranthene	ND	0.0081	EPA 8270D/SIM	5-31-17	6-1-17	
Benzo[a]pyrene	ND	0.0081	EPA 8270D/SIM	5-31-17	6-1-17	
Indeno[1,2,3-cd]pyrene	ND	0.0081	EPA 8270D/SIM	5-31-17	6-1-17	
Dibenz[a,h]anthracene	ND	0.0081	EPA 8270D/SIM	5-31-17	6-1-17	
Benzo[g,h,i]perylene	ND	0.0081	EPA 8270D/SIM	5-31-17	6-1-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorophenol	75	18 - 109				
Phenol-d6	79	25 - 111				
Nitrobenzene-d5	84	22 - 113				
2-Fluorobiphenyl	77	30 - 114				
	11	JU 117				
2,4,6-Tribromophenol	81	22 - 116				

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

page 1 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0531S1					
n-Nitrosodimethylamine	ND	0.033	EPA 8270D	5-31-17	5-31-17	
Pyridine	ND	0.33	EPA 8270D	5-31-17	5-31-17	
Phenol	ND	0.033	EPA 8270D	5-31-17	5-31-17	
Aniline	ND	0.17	EPA 8270D	5-31-17	5-31-17	
bis(2-Chloroethyl)ether	ND	0.033	EPA 8270D	5-31-17	5-31-17	
2-Chlorophenol	ND	0.033	EPA 8270D	5-31-17	5-31-17	
1,3-Dichlorobenzene	ND	0.033	EPA 8270D	5-31-17	5-31-17	
1,4-Dichlorobenzene	ND	0.033	EPA 8270D	5-31-17	5-31-17	
Benzyl alcohol	ND	0.17	EPA 8270D	5-31-17	5-31-17	
1,2-Dichlorobenzene	ND	0.033	EPA 8270D	5-31-17	5-31-17	
2-Methylphenol (o-Cresol)	ND	0.033	EPA 8270D	5-31-17	5-31-17	
bis(2-Chloroisopropyl)ether	ND	0.033	EPA 8270D	5-31-17	5-31-17	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.033	EPA 8270D	5-31-17	5-31-17	
n-Nitroso-di-n-propylamine	ND	0.033	EPA 8270D	5-31-17	5-31-17	
Hexachloroethane	ND	0.033	EPA 8270D	5-31-17	5-31-17	
Nitrobenzene	ND	0.033	EPA 8270D	5-31-17	5-31-17	
Isophorone	ND	0.033	EPA 8270D	5-31-17	5-31-17	
2-Nitrophenol	ND	0.033	EPA 8270D	5-31-17	5-31-17	
2,4-Dimethylphenol	ND	0.033	EPA 8270D	5-31-17	5-31-17	
bis(2-Chloroethoxy)methane	ND	0.033	EPA 8270D	5-31-17	5-31-17	
2,4-Dichlorophenol	ND	0.033	EPA 8270D	5-31-17	5-31-17	
1,2,4-Trichlorobenzene	ND	0.033	EPA 8270D	5-31-17	5-31-17	
Naphthalene	ND	0.0067	EPA 8270D/SIM	5-31-17	5-31-17	
4-Chloroaniline	ND	0.17	EPA 8270D	5-31-17	5-31-17	
Hexachlorobutadiene	ND	0.033	EPA 8270D	5-31-17	5-31-17	
4-Chloro-3-methylphenol	ND	0.033	EPA 8270D	5-31-17	5-31-17	
2-Methylnaphthalene	ND	0.0067	EPA 8270D/SIM	5-31-17	5-31-17	
1-Methylnaphthalene	ND	0.0067	EPA 8270D/SIM	5-31-17	5-31-17	
Hexachlorocyclopentadiene	ND	0.033	EPA 8270D	5-31-17	5-31-17	
2,4,6-Trichlorophenol	ND	0.033	EPA 8270D	5-31-17	5-31-17	
2,3-Dichloroaniline	ND	0.033	EPA 8270D	5-31-17	5-31-17	
2,4,5-Trichlorophenol	ND	0.033	EPA 8270D	5-31-17	5-31-17	
2-Chloronaphthalene	ND	0.033	EPA 8270D	5-31-17	5-31-17	
2-Nitroaniline	ND	0.033	EPA 8270D	5-31-17	5-31-17	
1,4-Dinitrobenzene	ND	0.033	EPA 8270D	5-31-17	5-31-17	
Dimethylphthalate	ND	0.033	EPA 8270D	5-31-17	5-31-17	
1,3-Dinitrobenzene	ND	0.033	EPA 8270D	5-31-17	5-31-17	
2,6-Dinitrotoluene	ND	0.033	EPA 8270D	5-31-17	5-31-17	
1,2-Dinitrobenzene	ND	0.033	EPA 8270D	5-31-17	5-31-17	
Acenaphthylene	ND	0.0067	EPA 8270D/SIM	5-31-17	5-31-17	
3-Nitroaniline	ND	0.033	EPA 8270D	5-31-17	5-31-17	

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0531S1					
2,4-Dinitrophenol	ND	0.17	EPA 8270D	5-31-17	5-31-17	
Acenaphthene	ND	0.0067	EPA 8270D/SIM	5-31-17	5-31-17	
4-Nitrophenol	ND	0.033	EPA 8270D	5-31-17	5-31-17	
2,4-Dinitrotoluene	ND	0.033	EPA 8270D	5-31-17	5-31-17	
Dibenzofuran	ND	0.033	EPA 8270D	5-31-17	5-31-17	
2,3,5,6-Tetrachlorophenol	ND	0.033	EPA 8270D	5-31-17	5-31-17	
2,3,4,6-Tetrachlorophenol	ND	0.033	EPA 8270D	5-31-17	5-31-17	
Diethylphthalate	ND	0.17	EPA 8270D	5-31-17	5-31-17	
4-Chlorophenyl-phenylether		0.033	EPA 8270D	5-31-17	5-31-17	
4-Nitroaniline	ND	0.033	EPA 8270D	5-31-17	5-31-17	
Fluorene	ND	0.0067	EPA 8270D/SIM	5-31-17	5-31-17	
4,6-Dinitro-2-methylphenol	ND	0.17	EPA 8270D	5-31-17	5-31-17	
n-Nitrosodiphenylamine	ND	0.033	EPA 8270D	5-31-17	5-31-17	
1,2-Diphenylhydrazine	ND	0.033	EPA 8270D	5-31-17	5-31-17	
4-Bromophenyl-phenylether	ND	0.033	EPA 8270D	5-31-17	5-31-17	
Hexachlorobenzene	ND	0.033	EPA 8270D	5-31-17	5-31-17	
Pentachlorophenol	ND	0.17	EPA 8270D	5-31-17	5-31-17	
Phenanthrene	ND	0.0067	EPA 8270D/SIM	5-31-17	5-31-17	
Anthracene	ND	0.0067	EPA 8270D/SIM	5-31-17	5-31-17	
Carbazole	ND	0.033	EPA 8270D	5-31-17	5-31-17	
Di-n-butylphthalate	ND	0.17	EPA 8270D	5-31-17	5-31-17	
Fluoranthene	ND	0.0067	EPA 8270D/SIM	5-31-17	5-31-17	
Benzidine	ND	0.33	EPA 8270D	5-31-17	5-31-17	
Pyrene	ND	0.0067	EPA 8270D/SIM	5-31-17	5-31-17	
Butylbenzylphthalate	ND	0.033	EPA 8270D	5-31-17	5-31-17	
bis-2-Ethylhexyladipate	ND	0.033	EPA 8270D	5-31-17	5-31-17	
3,3'-Dichlorobenzidine	ND	0.17	EPA 8270D	5-31-17	5-31-17	
Benzo[a]anthracene	ND	0.0067	EPA 8270D/SIM	5-31-17	5-31-17	
Chrysene	ND	0.0067	EPA 8270D/SIM	5-31-17	5-31-17	
bis(2-Ethylhexyl)phthalate	ND	0.033	EPA 8270D	5-31-17	5-31-17	
Di-n-octylphthalate	ND	0.033	EPA 8270D	5-31-17	5-31-17	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[a]pyrene	ND	0.0067	EPA 8270D/SIM	5-31-17	5-31-17	
Indeno[1,2,3-cd]pyrene	ND	0.0067	EPA 8270D/SIM	5-31-17	5-31-17	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270D/SIM	5-31-17	5-31-17	
Benzo[g,h,i]perylene	ND	0.0067	EPA 8270D/SIM	5-31-17	5-31-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorophenol	76	18 - 109				
Phenol-d6	76	25 - 111				
Nitrobenzene-d5	78	22 - 113				
2-Fluorobiphenyl	78	30 - 114				
2,4,6-Tribromophenol	<i>7</i> 5	22 - 116				
Terphenyl-d14	79	33 - 114				

Project: 15-05986-040

SEMIVOLATILES EPA 8270D/SIM MS/MSD QUALITY CONTROL

					Source	Per	cent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Rec	overy	Limits	RPD	Limit	Flags
MATRIX SPIKES											
Laboratory ID:	05-29	99-03									
	MS	MSD	MS	MSD		MS	MSD				
Phenol	0.904	0.784	1.33	1.33	ND	68	59	25 - 99	14	36	
2-Chlorophenol	0.937	0.814	1.33	1.33	ND	70	61	21 - 104	14	38	
1,4-Dichlorobenzene	0.479	0.408	0.667	0.667	ND	72	61	20 - 110	16	40	
n-Nitroso-di-n-propylamine	0.436	0.371	0.667	0.667	ND	65	56	24 - 100	16	38	
1,2,4-Trichlorobenzene	0.469	0.401	0.667	0.667	ND	70	60	21 - 110	16	40	
4-Chloro-3-methylphenol	0.915	0.836	1.33	1.33	ND	69	63	26 - 109	9	29	
Acenaphthene	0.422	0.359	0.667	0.667	ND	63	54	33 - 99	16	30	
4-Nitrophenol	0.980	0.840	1.33	1.33	ND	74	63	21 - 107	15	29	
2,4-Dinitrotoluene	0.436	0.383	0.667	0.667	ND	65	57	20 - 106	13	30	
Pentachlorophenol	0.987	0.818	1.33	1.33	ND	74	62	20 - 113	19	31	
Pyrene	0.462	0.409	0.667	0.667	ND	69	61	24 - 115	12	28	
Surrogate:											
2-Fluorophenol						76	65	18 - 109			
Phenol-d6						78	67	25 - 111			
Nitrobenzene-d5						74	66	22 - 113			
2-Fluorobiphenyl						77	65	30 - 114			
2,4,6-Tribromophenol						72	65	22 - 116			
Terphenyl-d14						74	67	33 - 114			

Project: 15-05986-040

PAHs EPA 8270D/SIM

Matrix: Water Units: ug/L

	-			Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP1-W					
Laboratory ID:	05-299-02					
Naphthalene	ND	0.096	EPA 8270D/SIM	5-25-17	5-25-17	
2-Methylnaphthalene	ND	0.096	EPA 8270D/SIM	5-25-17	5-25-17	
1-Methylnaphthalene	ND	0.096	EPA 8270D/SIM	5-25-17	5-25-17	
Acenaphthylene	ND	0.096	EPA 8270D/SIM	5-25-17	5-25-17	
Acenaphthene	ND	0.096	EPA 8270D/SIM	5-25-17	5-25-17	
Fluorene	ND	0.096	EPA 8270D/SIM	5-25-17	5-25-17	
Phenanthrene	ND	0.096	EPA 8270D/SIM	5-25-17	5-25-17	
Anthracene	ND	0.096	EPA 8270D/SIM	5-25-17	5-25-17	
Fluoranthene	ND	0.096	EPA 8270D/SIM	5-25-17	5-25-17	
Pyrene	ND	0.096	EPA 8270D/SIM	5-25-17	5-25-17	
Benzo[a]anthracene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-25-17	
Chrysene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-25-17	
Benzo[b]fluoranthene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-25-17	
Benzo(j,k)fluoranthene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-25-17	
Benzo[a]pyrene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-25-17	
Indeno(1,2,3-c,d)pyrene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-25-17	
Dibenz[a,h]anthracene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-25-17	
Benzo[g,h,i]perylene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-25-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	46	30 - 124				
Pyrene-d10	42	40 - 143				
Torphonyl d11	<i>1E</i>	27 427				

Project: 15-05986-040

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP2-W					
Laboratory ID:	05-299-06					
Naphthalene	ND	0.12	EPA 8270D/SIM	5-25-17	5-25-17	
2-Methylnaphthalene	ND	0.12	EPA 8270D/SIM	5-25-17	5-25-17	
1-Methylnaphthalene	ND	0.12	EPA 8270D/SIM	5-25-17	5-25-17	
Acenaphthylene	ND	0.12	EPA 8270D/SIM	5-25-17	5-25-17	
Acenaphthene	ND	0.12	EPA 8270D/SIM	5-25-17	5-25-17	
Fluorene	ND	0.12	EPA 8270D/SIM	5-25-17	5-25-17	
Phenanthrene	ND	0.12	EPA 8270D/SIM	5-25-17	5-25-17	
Anthracene	ND	0.12	EPA 8270D/SIM	5-25-17	5-25-17	
Fluoranthene	ND	0.12	EPA 8270D/SIM	5-25-17	5-25-17	
Pyrene	ND	0.12	EPA 8270D/SIM	5-25-17	5-25-17	
Benzo[a]anthracene	ND	0.012	EPA 8270D/SIM	5-25-17	5-25-17	
Chrysene	0.012	0.012	EPA 8270D/SIM	5-25-17	5-25-17	
Benzo[b]fluoranthene	0.020	0.012	EPA 8270D/SIM	5-25-17	5-25-17	
Benzo(j,k)fluoranthene	ND	0.012	EPA 8270D/SIM	5-25-17	5-25-17	
Benzo[a]pyrene	ND	0.012	EPA 8270D/SIM	5-25-17	5-25-17	
Indeno(1,2,3-c,d)pyrene	0.014	0.012	EPA 8270D/SIM	5-25-17	5-25-17	
Dibenz[a,h]anthracene	ND	0.012	EPA 8270D/SIM	5-25-17	5-25-17	
Benzo[g,h,i]perylene	0.020	0.012	EPA 8270D/SIM	5-25-17	5-25-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	72	30 - 124				
Pyrene-d10	80	40 - 143				

Project: 15-05986-040

PAHs EPA 8270D/SIM

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP3-W	ı «L	Metriou	i repared	Allalyzed	i iags
Laboratory ID:	05-299-09					
Naphthalene	ND	0.096	EPA 8270D/SIM	5-25-17	5-25-17	
2-Methylnaphthalene	ND	0.096	EPA 8270D/SIM	5-25-17	5-25-17	
1-Methylnaphthalene	ND	0.096	EPA 8270D/SIM	5-25-17	5-25-17	
Acenaphthylene	ND	0.096	EPA 8270D/SIM	5-25-17	5-25-17	
Acenaphthene	ND	0.096	EPA 8270D/SIM	5-25-17	5-25-17	
Fluorene	ND	0.096	EPA 8270D/SIM	5-25-17	5-25-17	
Phenanthrene	ND	0.096	EPA 8270D/SIM	5-25-17	5-25-17	
Anthracene	ND	0.096	EPA 8270D/SIM	5-25-17	5-25-17	
Fluoranthene	ND	0.096	EPA 8270D/SIM	5-25-17	5-25-17	
Pyrene	ND	0.096	EPA 8270D/SIM	5-25-17	5-25-17	
Benzo[a]anthracene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-25-17	
Chrysene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-25-17	
Benzo[b]fluoranthene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-25-17	
Benzo(j,k)fluoranthene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-25-17	
Benzo[a]pyrene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-25-17	
ndeno(1,2,3-c,d)pyrene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-25-17	
Dibenz[a,h]anthracene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-25-17	
Benzo[g,h,i]perylene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-25-17	
Surrogate:	Percent Recovery	Control Limits	_			•
2-Fluorobiphenyl	88	30 - 124				
Pyrene-d10	83	40 - 143				
Tamahana dada	400	07 407				

Project: 15-05986-040

PAHs EPA 8270D/SIM

Analysia	Beault .	DOL	Mathad	Date	Date	Elect
Analyte	Result PP4-W	PQL	Method	Prepared	Analyzed	Flags
Client ID:						
Laboratory ID:	05-299-12					
Naphthalene	ND	0.10	EPA 8270D/SIM	5-25-17	5-26-17	
2-Methylnaphthalene	ND	0.10	EPA 8270D/SIM	5-25-17	5-26-17	
1-Methylnaphthalene	ND	0.10	EPA 8270D/SIM	5-25-17	5-26-17	
Acenaphthylene	ND	0.10	EPA 8270D/SIM	5-25-17	5-26-17	
Acenaphthene	ND	0.10	EPA 8270D/SIM	5-25-17	5-26-17	
Fluorene	ND	0.10	EPA 8270D/SIM	5-25-17	5-26-17	
Phenanthrene	ND	0.10	EPA 8270D/SIM	5-25-17	5-26-17	
Anthracene	ND	0.10	EPA 8270D/SIM	5-25-17	5-26-17	
Fluoranthene	ND	0.10	EPA 8270D/SIM	5-25-17	5-26-17	
Pyrene	ND	0.10	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[a]anthracene	0.013	0.010	EPA 8270D/SIM	5-25-17	5-26-17	
Chrysene	0.021	0.010	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[b]fluoranthene	0.036	0.010	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo(j,k)fluoranthene	0.011	0.010	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[a]pyrene	0.024	0.010	EPA 8270D/SIM	5-25-17	5-26-17	
Indeno(1,2,3-c,d)pyrene	0.040	0.010	EPA 8270D/SIM	5-25-17	5-26-17	
Dibenz[a,h]anthracene	0.011	0.010	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[g,h,i]perylene	0.046	0.010	EPA 8270D/SIM	5-25-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	49	30 - 124				
Pyrene-d10	47	40 - 143				
Townshamud ald 4	50	07 407				

Project: 15-05986-040

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP5-W					
Laboratory ID:	05-299-15					
Naphthalene	0.18	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
2-Methylnaphthalene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
1-Methylnaphthalene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
Acenaphthylene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
Acenaphthene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
Fluorene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
Phenanthrene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
Anthracene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
Fluoranthene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
Pyrene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[a]anthracene	ND	0.0097	EPA 8270D/SIM	5-25-17	5-26-17	
Chrysene	ND	0.0097	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[b]fluoranthene	ND	0.0097	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo(j,k)fluoranthene	ND	0.0097	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[a]pyrene	ND	0.0097	EPA 8270D/SIM	5-25-17	5-26-17	
Indeno(1,2,3-c,d)pyrene	ND	0.0097	EPA 8270D/SIM	5-25-17	5-26-17	
Dibenz[a,h]anthracene	ND	0.0097	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[g,h,i]perylene	ND	0.0097	EPA 8270D/SIM	5-25-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	35	30 - 124				
Pyrene-d10	31	40 - 143				Q
Terphenyl-d14	40	27 - 127				

Project: 15-05986-040

PAHs EPA 8270D/SIM

Matrix: Water Units: ug/L

G				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP6-W					
Laboratory ID:	05-299-18					
Naphthalene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
2-Methylnaphthalene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
1-Methylnaphthalene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
Acenaphthylene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
Acenaphthene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
Fluorene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
Phenanthrene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
Anthracene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
Fluoranthene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
Pyrene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[a]anthracene	0.052	0.0097	EPA 8270D/SIM	5-25-17	5-26-17	
Chrysene	0.080	0.0097	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[b]fluoranthene	0.14	0.0097	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo(j,k)fluoranthene	0.042	0.0097	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[a]pyrene	0.051	0.0097	EPA 8270D/SIM	5-25-17	5-26-17	
Indeno(1,2,3-c,d)pyrene	0.066	0.0097	EPA 8270D/SIM	5-25-17	5-26-17	
Dibenz[a,h]anthracene	0.018	0.0097	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[g,h,i]perylene	0.055	0.0097	EPA 8270D/SIM	5-25-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	66	30 - 124				
Pyrene-d10	64	40 - 143				
Terphenyl-d14	78	27 - 127				

Project: 15-05986-040

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP7-W					
Laboratory ID:	05-299-21					
Naphthalene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
2-Methylnaphthalene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
1-Methylnaphthalene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
Acenaphthylene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
Acenaphthene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
Fluorene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
Phenanthrene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
Anthracene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
Fluoranthene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
Pyrene	ND	0.097	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[a]anthracene	ND	0.0097	EPA 8270D/SIM	5-25-17	5-26-17	
Chrysene	ND	0.0097	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[b]fluoranthene	ND	0.0097	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo(j,k)fluoranthene	ND	0.0097	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[a]pyrene	ND	0.0097	EPA 8270D/SIM	5-25-17	5-26-17	
Indeno(1,2,3-c,d)pyrene	ND	0.0097	EPA 8270D/SIM	5-25-17	5-26-17	
Dibenz[a,h]anthracene	ND	0.0097	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[g,h,i]perylene	ND	0.0097	EPA 8270D/SIM	5-25-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	63	30 - 124				
Pyrene-d10	23	40 - 143				Q
Terphenyl-d14	54	27 - 127				

Project: 15-05986-040

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP8-W					
Laboratory ID:	05-299-24					
Naphthalene	ND	0.095	EPA 8270D/SIM	5-25-17	5-26-17	
2-Methylnaphthalene	ND	0.095	EPA 8270D/SIM	5-25-17	5-26-17	
1-Methylnaphthalene	ND	0.095	EPA 8270D/SIM	5-25-17	5-26-17	
Acenaphthylene	ND	0.095	EPA 8270D/SIM	5-25-17	5-26-17	
Acenaphthene	ND	0.095	EPA 8270D/SIM	5-25-17	5-26-17	
Fluorene	ND	0.095	EPA 8270D/SIM	5-25-17	5-26-17	
Phenanthrene	ND	0.095	EPA 8270D/SIM	5-25-17	5-26-17	
Anthracene	ND	0.095	EPA 8270D/SIM	5-25-17	5-26-17	
Fluoranthene	ND	0.095	EPA 8270D/SIM	5-25-17	5-26-17	
Pyrene	ND	0.095	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[a]anthracene	ND	0.0095	EPA 8270D/SIM	5-25-17	5-26-17	
Chrysene	ND	0.0095	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[b]fluoranthene	ND	0.0095	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo(j,k)fluoranthene	ND	0.0095	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[a]pyrene	ND	0.0095	EPA 8270D/SIM	5-25-17	5-26-17	
Indeno(1,2,3-c,d)pyrene	ND	0.0095	EPA 8270D/SIM	5-25-17	5-26-17	
Dibenz[a,h]anthracene	ND	0.0095	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[g,h,i]perylene	ND	0.0095	EPA 8270D/SIM	5-25-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	75	30 - 124				
Pyrene-d10	79	40 - 143				
Torphonyl d11	100	27 427				

Project: 15-05986-040

PAHs EPA 8270D/SIM

naluta					Date	
nalyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP9-W					
aboratory ID:	05-299-27					
laphthalene	ND	0.096	EPA 8270D/SIM	5-25-17	5-26-17	
-Methylnaphthalene	ND	0.096	EPA 8270D/SIM	5-25-17	5-26-17	
1-Methylnaphthalene ND		0.096	EPA 8270D/SIM	5-25-17	5-26-17	
cenaphthylene	ND	0.096	EPA 8270D/SIM	5-25-17	5-26-17	
cenaphthene	ND	0.096	EPA 8270D/SIM	5-25-17	5-26-17	
·luorene	ND	0.096	EPA 8270D/SIM	5-25-17	5-26-17	
henanthrene	ND	0.096	EPA 8270D/SIM	5-25-17	5-26-17	
nthracene	ND	0.096	EPA 8270D/SIM	5-25-17	5-26-17	
·luoranthene	ND	0.096	EPA 8270D/SIM	5-25-17	5-26-17	
yrene	ND	0.096	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[a]anthracene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-26-17	
Chrysene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[b]fluoranthene	0.019	0.0096	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo(j,k)fluoranthene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[a]pyrene	0.0098	0.0096	EPA 8270D/SIM	5-25-17	5-26-17	
ndeno(1,2,3-c,d)pyrene	0.011	0.0096	EPA 8270D/SIM	5-25-17	5-26-17	
Dibenz[a,h]anthracene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[g,h,i]perylene	0.012	0.0096	EPA 8270D/SIM	5-25-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits			_	
-Fluorobiphenyl	62	30 - 124				
Pyrene-d10	44	40 - 143				
erphenyl-d14	67	27 - 127				

Project: 15-05986-040

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP10-W					
Laboratory ID:	05-299-29					
Naphthalene	ND	0.096	EPA 8270D/SIM	5-25-17	5-26-17	
2-Methylnaphthalene	ND	0.096	EPA 8270D/SIM	5-25-17	5-26-17	
1-Methylnaphthalene	ND	0.096	EPA 8270D/SIM	5-25-17	5-26-17	
Acenaphthylene	ND	0.096	EPA 8270D/SIM	5-25-17	5-26-17	
Acenaphthene	ND	0.096	EPA 8270D/SIM	5-25-17	5-26-17	
Fluorene	ND	0.096	EPA 8270D/SIM	5-25-17	5-26-17	
Phenanthrene	ND	0.096	EPA 8270D/SIM	5-25-17	5-26-17	
Anthracene	ND	0.096	EPA 8270D/SIM	5-25-17	5-26-17	
Fluoranthene	ND	0.096	EPA 8270D/SIM	5-25-17	5-26-17	
Pyrene	ND	0.096	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[a]anthracene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-26-17	
Chrysene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[b]fluoranthene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo(j,k)fluoranthene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[a]pyrene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-26-17	
Indeno(1,2,3-c,d)pyrene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-26-17	
Dibenz[a,h]anthracene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-26-17	
Benzo[g,h,i]perylene	ND	0.0096	EPA 8270D/SIM	5-25-17	5-26-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	44	30 - 124				
Pyrene-d10	51	40 - 143				
Torphopul d11	61	27 127				

Project: 15-05986-040

PAHs EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

Matrix: Water Units: ug/L

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0525W1					
Naphthalene	ND	0.10	EPA 8270D/SIM	5-25-17	5-25-17	
2-Methylnaphthalene	ND	0.10	EPA 8270D/SIM	5-25-17	5-25-17	
1-Methylnaphthalene	ND	0.10	EPA 8270D/SIM	5-25-17	5-25-17	
Acenaphthylene	ND	0.10	EPA 8270D/SIM	5-25-17	5-25-17	
Acenaphthene	ND	0.10	EPA 8270D/SIM	5-25-17	5-25-17	
Fluorene	ND	0.10	EPA 8270D/SIM	5-25-17	5-25-17	
Phenanthrene	ND	0.10	EPA 8270D/SIM	5-25-17	5-25-17	
Anthracene	ND	0.10	EPA 8270D/SIM	5-25-17	5-25-17	
Fluoranthene	ND	0.10	EPA 8270D/SIM	5-25-17	5-25-17	
Pyrene	ND	0.10	EPA 8270D/SIM	5-25-17	5-25-17	
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	5-25-17	5-25-17	
Chrysene	ND	0.010	EPA 8270D/SIM	5-25-17	5-25-17	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	5-25-17	5-25-17	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	5-25-17	5-25-17	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	5-25-17	5-25-17	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	5-25-17	5-25-17	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	5-25-17	5-25-17	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270D/SIM	5-25-17	5-25-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	81	30 - 124				
Pyrene-d10	87	40 - 143				

Terphenyl-d14 27 - 127 100



Project: 15-05986-040

PAHS EPA 8270D/SIM SB/SBD QUALITY CONTROL

					Per	cent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Rec	overy	Limits	RPD	Limit	Flags
SPIKE BLANKS										
Laboratory ID:	SB05	25W1								
	SB	SBD	SB	SBD	SB	SBD				
Naphthalene	0.462	0.450	0.500	0.500	92	90	29 - 101	3	47	
Acenaphthylene	0.465	0.442	0.500	0.500	93	88	20 - 117	5	50	
Acenaphthene	0.478	0.469	0.500	0.500	96	94	37 - 109	2	43	
Fluorene	0.493	0.479	0.500	0.500	99	96	47 - 108	3	34	
Phenanthrene	0.468	0.435	0.500	0.500	94	87	49 - 109	7	28	
Anthracene	0.526	0.509	0.500	0.500	105	102	34 - 140	3	32	
Fluoranthene	0.528	0.513	0.500	0.500	106	103	45 - 120	3	39	
Pyrene	0.533	0.544	0.500	0.500	107	109	42 - 133	2	39	
Benzo[a]anthracene	0.550	0.523	0.500	0.500	110	105	71 - 117	5	28	
Chrysene	0.559	0.543	0.500	0.500	112	109	53 - 110	3	25	I
Benzo[b]fluoranthene	0.536	0.512	0.500	0.500	107	102	53 - 123	5	37	
Benzo(j,k)fluoranthene	0.550	0.554	0.500	0.500	110	111	52 - 119	1	41	
Benzo[a]pyrene	0.512	0.508	0.500	0.500	102	102	37 - 129	1	33	
Indeno(1,2,3-c,d)pyrene	0.638	0.608	0.500	0.500	128	122	45 - 128	5	31	
Dibenz[a,h]anthracene	0.513	0.486	0.500	0.500	103	97	54 - 120	5	30	
Benzo[g,h,i]perylene	0.494	0.486	0.500	0.500	99	97	49 - 117	2	29	
Surrogate:										
2-Fluorobiphenyl					85	81	30 - 124			
Pyrene-d10					93	86	40 - 143			
Terphenyl-d14					106	102	27 - 127			

Project: 15-05986-040

TOTAL METALS EPA 6010C/7471B

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID: Client ID:	05-299-01 PP1-2					
Arsenic	ND	12	6010C	6-2-17	6-2-17	
Barium	62	3.1	6010C	6-2-17	6-2-17	
Cadmium	ND	0.62	6010C	6-2-17	6-2-17	
Chromium	20	0.62	6010C	6-2-17	6-2-17	
Lead	8.9	6.2	6010C	6-2-17	6-2-17	
Mercury	ND	0.31	7471B	5-31-17	5-31-17	
Selenium	ND	12	6010C	6-2-17	6-2-17	
Silver	ND	1.2	6010C	6-2-17	6-2-17	
Lab ID: Client ID:	05-299-03 PP1-10					
Arsenic	ND	13	6010C	6-2-17	6-5-17	
Barium	30	3.3	6010C	6-2-17	6-5-17	
Cadmium	ND	0.67	6010C	6-2-17	6-5-17	
Chromium	13	0.67	6010C	6-2-17	6-5-17	
Lead	ND	6.7	6010C	6-2-17	6-5-17	
Mercury	ND	0.33	7471B	5-31-17	5-31-17	
Selenium	ND	13	6010C	6-2-17	6-5-17	
Silver	ND	1.3	6010C	6-2-17	6-5-17	

Project: 15-05986-040

TOTAL METALS EPA 6010C/7471B

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-299-04					
Client ID:	PP2-2.5					
Arsenic	ND	13	6010C	6-2-17	6-2-17	
Barium	31	3.3	6010C	6-2-17	6-2-17	
Cadmium	ND	0.66	6010C	6-2-17	6-2-17	
Chromium	11	0.66	6010C	6-2-17	6-2-17	
Lead	ND	6.6	6010C	6-2-17	6-2-17	
Mercury	ND	0.33	7471B	5-31-17	5-31-17	
Selenium	ND	13	6010C	6-2-17	6-2-17	
Silver	ND	1.3	6010C	6-2-17	6-2-17	

Lab ID: Client ID:	05-299-05 PP2-10					
Arsenic	ND	15	6010C	6-2-17	6-2-17	
Barium	21	3.6	6010C	6-2-17	6-2-17	
Cadmium	ND	0.73	6010C	6-2-17	6-2-17	
Chromium	11	0.73	6010C	6-2-17	6-2-17	
Lead	ND	7.3	6010C	6-2-17	6-2-17	
Mercury	ND	0.36	7471B	5-31-17	5-31-17	
Selenium	ND	15	6010C	6-2-17	6-2-17	
Silver	ND	1.5	6010C	6-2-17	6-2-17	

Project: 15-05986-040

TOTAL METALS EPA 6010C/7471B

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-299-07					
Client ID:	PP3-2.5					
Arsenic	ND	13	6010C	6-2-17	6-2-17	
Barium	80	3.1	6010C	6-2-17	6-2-17	
Cadmium	ND	0.63	6010C	6-2-17	6-2-17	
Chromium	27	0.63	6010C	6-2-17	6-2-17	
Lead	25	6.3	6010C	6-2-17	6-2-17	
Mercury	ND	0.31	7471B	5-31-17	5-31-17	
Selenium	ND	13	6010C	6-2-17	6-2-17	
Silver	ND	1.3	6010C	6-2-17	6-2-17	

Lab ID:	05-299-08					
Client ID:	PP3-10					
Arsenic	ND	12	6010C	6-2-17	6-2-17	
Barium	26	3.1	6010C	6-2-17	6-2-17	
Cadmium	ND	0.62	6010C	6-2-17	6-2-17	
Chromium	9.5	0.62	6010C	6-2-17	6-2-17	
Lead	ND	6.2	6010C	6-2-17	6-2-17	
Mercury	ND	0.31	7471B	5-31-17	5-31-17	
Selenium	ND	12	6010C	6-2-17	6-2-17	
Silver	ND	1.2	6010C	6-2-17	6-2-17	

Project: 15-05986-040

TOTAL METALS EPA 6010C/7471B

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-299-10					
Client ID:	PP4-3					
Arsenic	ND	13	6010C	6-2-17	6-2-17	
Barium	78	3.2	6010C	6-2-17	6-2-17	
Cadmium	ND	0.63	6010C	6-2-17	6-2-17	
Chromium	29	0.63	6010C	6-2-17	6-2-17	
Lead	84	6.3	6010C	6-2-17	6-2-17	
Mercury	ND	0.32	7471B	5-31-17	5-31-17	
Selenium	ND	13	6010C	6-2-17	6-2-17	
Silver	ND	1.3	6010C	6-2-17	6-2-17	

Lab ID: Client ID:	05-299-11 PP4-10					
Arsenic	ND	12	6010C	6-2-17	6-2-17	
Barium	38	3.1	6010C	6-2-17	6-2-17	
Cadmium	ND	0.62	6010C	6-2-17	6-2-17	
Chromium	13	0.62	6010C	6-2-17	6-2-17	
_ead	ND	6.2	6010C	6-2-17	6-2-17	
Mercury	ND	0.31	7471B	5-31-17	5-31-17	
Selenium	ND	12	6010C	6-2-17	6-2-17	
Silver	ND	1.2	6010C	6-2-17	6-2-17	

Project: 15-05986-040

TOTAL METALS EPA 6010C/7471B

Matrix: Soil

		PQL		Date Prepared	Date Analyzed	Flags
Analyte	Result		EPA Method			
Lab ID:	05-299-13					
Client ID:	PP5-2.5					
Arsenic	ND	13	6010C	6-2-17	6-2-17	
Barium	47	3.3	6010C	6-2-17	6-2-17	
Cadmium	ND	0.65	6010C	6-2-17	6-2-17	
Chromium	23	0.65	6010C	6-2-17	6-2-17	
Lead	27	6.5	6010C	6-2-17	6-2-17	
Mercury	ND	0.33	7471B	5-31-17	5-31-17	
Selenium	ND	13	6010C	6-2-17	6-2-17	
Silver	ND	1.3	6010C	6-2-17	6-2-17	

Lab ID: Client ID:	05-299-14 PP5-10					
Arsenic	ND	13	6010C	6-2-17	6-2-17	
Barium	28	3.4	6010C	6-2-17	6-2-17	
Cadmium	ND	0.67	6010C	6-2-17	6-2-17	
Chromium	13	0.67	6010C	6-2-17	6-2-17	
Lead	ND	6.7	6010C	6-2-17	6-2-17	
Mercury	ND	0.34	7471B	5-31-17	5-31-17	
Selenium	ND	13	6010C	6-2-17	6-2-17	
Silver	ND	1.3	6010C	6-2-17	6-2-17	

Project: 15-05986-040

TOTAL METALS EPA 6010C/7471B

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-299-16					
Client ID:	PP6-2					
Arsenic	ND	12	6010C	6-2-17	6-2-17	
Barium	53	2.9	6010C	6-2-17	6-2-17	
Cadmium	ND	0.58	6010C	6-2-17	6-2-17	
Chromium	29	0.58	6010C	6-2-17	6-2-17	
Lead	9.7	5.8	6010C	6-2-17	6-2-17	
Mercury	ND	0.29	7471B	5-31-17	5-31-17	
Selenium	ND	12	6010C	6-2-17	6-2-17	
Silver	ND	1.2	6010C	6-2-17	6-2-17	

Lab ID:	05-299-17					
Client ID:	PP6-10					
Arsenic	ND	15	6010C	6-2-17	6-2-17	
Barium	34	3.7	6010C	6-2-17	6-2-17	
Cadmium	ND	0.74	6010C	6-2-17	6-2-17	
Chromium	14	0.74	6010C	6-2-17	6-2-17	
Lead	ND	7.4	6010C	6-2-17	6-2-17	
Mercury	ND	0.37	7471B	5-31-17	5-31-17	
Selenium	ND	15	6010C	6-2-17	6-2-17	
Silver	ND	1.5	6010C	6-2-17	6-2-17	

Project: 15-05986-040

TOTAL METALS EPA 6010C/7471B

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-299-19					
Client ID:	PP7-0					
Arsenic	ND	12	6010C	6-2-17	6-2-17	
Barium	31	3.1	6010C	6-2-17	6-2-17	
Cadmium	ND	0.61	6010C	6-2-17	6-2-17	
Chromium	12	0.61	6010C	6-2-17	6-2-17	
Lead	ND	6.1	6010C	6-2-17	6-2-17	
Mercury	ND	0.31	7471B	5-31-17	5-31-17	
Selenium	ND	12	6010C	6-2-17	6-2-17	
Silver	ND	1.2	6010C	6-2-17	6-2-17	
	_		_		•	

Lab ID: Client ID:	05-299-20 PP7-10					
Arsenic	ND	13	6010C	6-2-17	6-2-17	
Barium	22	3.1	6010C	6-2-17	6-2-17	
Cadmium	ND	0.63	6010C	6-2-17	6-2-17	
Chromium	11	0.63	6010C	6-2-17	6-2-17	
Lead	ND	6.3	6010C	6-2-17	6-2-17	
Mercury	ND	0.31	7471B	5-31-17	5-31-17	
Selenium	ND	13	6010C	6-2-17	6-2-17	
Silver	ND	1.3	6010C	6-2-17	6-2-17	

Project: 15-05986-040

TOTAL METALS EPA 6010C/7471B

Matrix: Soil

				Date	Date Analyzed	Flags
Analyte	Result	PQL	EPA Method	Prepared		
Lab ID: Client ID:	05-299-22 PP8-0.5					
Arsenic	ND	11	6010C	6-2-17	6-2-17	
Barium	22	2.8	6010C	6-2-17	6-2-17	
Cadmium	ND	0.57	6010C	6-2-17	6-2-17	
Chromium	13	0.57	6010C	6-2-17	6-2-17	
Lead	ND	5.7	6010C	6-2-17	6-2-17	
Mercury	ND	0.28	7471B	5-31-17	5-31-17	
Selenium	ND	11	6010C	6-2-17	6-2-17	
Silver	ND	1.1	6010C	6-2-17	6-2-17	

Project: 15-05986-040

TOTAL METALS EPA 6010C/7471B

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-299-23 PP8-10					
Arsenic	ND	14	6010C	6-2-17	6-2-17	
Barium	37	3.6	6010C	6-2-17	6-2-17	
Cadmium	ND	0.71	6010C	6-2-17	6-2-17	
Chromium	13	0.71	6010C	6-2-17	6-2-17	
Lead	ND	7.1	6010C	6-2-17	6-2-17	
Mercury	ND	0.36	7471B	5-31-17	5-31-17	
Selenium	ND	14	6010C	6-2-17	6-2-17	
Silver	ND	1.4	6010C	6-2-17	6-2-17	
Lab ID: Client ID:	05-299-25 PP9-0					
Arsenic	ND	10	6010C	6-2-17	6-2-17	
Barium	62	2.6	6010C	6-2-17	6-2-17	
Cadmium	ND	0.52	6010C	6-2-17	6-2-17	
Chromium	25	0.52	6010C	6-2-17	6-2-17	
Lead	ND	5.2	6010C	6-2-17	6-2-17	
Mercury	ND	0.26	7471B	5-31-17	5-31-17	
Selenium	ND	10	6010C	6-2-17	6-2-17	
Silver	ND	1.0	6010C	6-2-17	6-2-17	

Project: 15-05986-040

TOTAL METALS EPA 6010C/7471B

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-299-26					
Client ID:	PP9-10					
Arsenic	ND	12	6010C	6-2-17	6-2-17	
Barium	28	3.1	6010C	6-2-17	6-2-17	
Cadmium	ND	0.62	6010C	6-2-17	6-2-17	
Chromium	14	0.62	6010C	6-2-17	6-2-17	
Lead	ND	6.2	6010C	6-2-17	6-2-17	
Mercury	ND	0.31	7471B	5-31-17	5-31-17	
Selenium	ND	12	6010C	6-2-17	6-2-17	
Silver	ND	1.2	6010C	6-2-17	6-2-17	

Lab ID: Client ID:	05-299-28 PP10-0					
Arsenic	ND	12	6010C	6-2-17	6-2-17	
Barium	30	3.0	6010C	6-2-17	6-2-17	
Cadmium	ND	0.61	6010C	6-2-17	6-2-17	
Chromium	13	0.61	6010C	6-2-17	6-2-17	
Lead	ND	6.1	6010C	6-2-17	6-2-17	
Mercury	ND	0.30	7471B	5-31-17	5-31-17	
Selenium	ND	12	6010C	6-2-17	6-2-17	
Silver	ND	1.2	6010C	6-2-17	6-2-17	

Project: 15-05986-040

TOTAL METALS EPA 6010C METHOD BLANK QUALITY CONTROL

Date Extracted: 6-2-17
Date Analyzed: 6-2-17

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: MB0602SM2

Analyte	Method	Result	PQL
Arsenic	6010C	ND	10
Barium	6010C	ND	2.5
Cadmium	6010C	ND	0.50
Chromium	6010C	ND	0.50
Lead	6010C	ND	5.0
Selenium	6010C	ND	10
Silver	6010C	ND	1.0

Project: 15-05986-040

TOTAL MERCURY EPA 7471B METHOD BLANK QUALITY CONTROL

Date Extracted: 5-31-17
Date Analyzed: 5-31-17

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: MB0531S2

Analyte Method Result PQL

Mercury 7471B **ND** 0.25

Project: 15-05986-040

TOTAL METALS EPA 6010C DUPLICATE QUALITY CONTROL

Date Extracted: 6-2-17
Date Analyzed: 6-2-17

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 05-299-01

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Arsenic	ND	ND	NA	10	
Barium	50.2	41.8	18	2.5	
Cadmium	ND	ND	NA	0.50	
Chromium	15.9	19.1	18	0.50	
Lead	7.20	5.45	28	5.0	С
Selenium	ND	ND	NA	10	
Silver	ND	ND	NA	1.0	

Project: 15-05986-040

TOTAL MERCURY EPA 7471B DUPLICATE QUALITY CONTROL

Date Extracted: 5-31-17
Date Analyzed: 5-31-17

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 05-299-03

Analyte Sample Duplicate
Result Result RPD PQL Flags

Mercury ND ND NA 0.25

Project: 15-05986-040

TOTAL METALS EPA 6010C MS/MSD QUALITY CONTROL

Date Extracted: 6-2-17
Date Analyzed: 6-2-17

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 05-299-01

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Arsenic	100	98.8	99	99.1	99	0	
Barium	100	151	101	151	101	0	
Cadmium	50.0	49.4	99	49.6	99	0	
Chromium	100	114	98	114	98	0	
Lead	250	248	96	249	97	1	
Selenium	100	96.6	97	95.2	95	2	
Silver	25.0	23.3	93	23.5	94	1	

Project: 15-05986-040

TOTAL MERCURY EPA 7471B MS/MSD QUALITY CONTROL

Date Extracted: 5-31-17
Date Analyzed: 5-31-17

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 05-299-03

	Spike		Percent		Percent		
Analyte	Level	MS	Recovery	MSD	Recovery	RPD	Flags
Mercury	0.500	0.537	107	0.543	109	1	

Project: 15-05986-040

TOTAL METALS EPA 200.8/7470A

Matrix: Water Units: ug/L (ppb)

	- " '			Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-299-02 PP1-W					
Antimony	ND	5.6	200.8	5-26-17	5-26-17	
Arsenic	36	3.3	200.8	5-26-17	5-26-17	
Beryllium	ND	11	200.8	5-26-17	5-26-17	
Cadmium	ND	4.4	200.8	5-26-17	5-26-17	
Chromium	73	11	200.8	5-26-17	5-26-17	
Copper	80	11	200.8	5-26-17	5-26-17	
Lead	250	1.1	200.8	5-26-17	5-26-17	
Mercury	ND	0.50	7470A	5-25-17	5-25-17	
Nickel	33	22	200.8	5-26-17	5-26-17	
Selenium	ND	5.6	200.8	5-26-17	5-26-17	
Silver	ND	11	200.8	5-26-17	5-26-17	
Thallium	ND	5.6	200.8	5-26-17	5-26-17	
Zinc	700	56	200.8	5-26-17	6-1-17	

Project: 15-05986-040

TOTAL METALS EPA 200.8/7470A

Matrix: Water
Units: ug/L (ppb)

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-299-06					
Client ID:	PP2-W					
Antimony	ND	5.6	200.8	5-26-17	5-26-17	
Arsenic	910	6.7	200.8	5-26-17	5-26-17	
Beryllium	34	22	200.8	5-26-17	5-26-17	
Cadmium	24	8.9	200.8	5-26-17	5-26-17	
Chromium	2000	110	200.8	5-26-17	6-1-17	
Copper	4200	110	200.8	5-26-17	6-1-17	
Lead	2100	22	200.8	5-26-17	6-1-17	
Mercury	4.3	0.50	7470A	5-25-17	5-25-17	
Nickel	1800	220	200.8	5-26-17	6-1-17	
Selenium	53	11	200.8	5-26-17	5-26-17	
Silver	ND	11	200.8	5-26-17	5-26-17	
Thallium	9.6	5.6	200.8	5-26-17	6-1-17	
Zinc	5500	560	200.8	5-26-17	6-1-17	

Project: 15-05986-040

TOTAL METALS EPA 200.8/7470A

Matrix: Water Units: ug/L (ppb)

	- " ' '			Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-299-09					
Client ID:	PP3-W					
Antimony	ND	5.6	200.8	5-26-17	5-26-17	
Arsenic	18	3.3	200.8	5-26-17	5-26-17	
Beryllium	ND	11	200.8	5-26-17	5-26-17	
Cadmium	ND	4.4	200.8	5-26-17	5-26-17	
Chromium	63	11	200.8	5-26-17	5-26-17	
Copper	120	11	200.8	5-26-17	5-26-17	
Lead	54	1.1	200.8	5-26-17	5-26-17	
Mercury	ND	0.50	7470A	5-25-17	5-25-17	
Nickel	32	22	200.8	5-26-17	5-26-17	
Selenium	ND	5.6	200.8	5-26-17	5-26-17	
Silver	ND	11	200.8	5-26-17	5-26-17	
Thallium	ND	5.6	200.8	5-26-17	5-26-17	
Zinc	130	28	200.8	5-26-17	5-26-17	

Project: 15-05986-040

TOTAL METALS EPA 200.8/7470A

Matrix: Water
Units: ug/L (ppb)

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-299-12					
Client ID:	PP4-W					
Antimony	47	5.6	200.8	5-26-17	5-26-17	
Arsenic	110	3.3	200.8	5-26-17	5-26-17	
Beryllium	ND	11	200.8	5-26-17	5-26-17	
Cadmium	ND	4.4	200.8	5-26-17	5-26-17	
Chromium	240	11	200.8	5-26-17	5-26-17	
Copper	460	11	200.8	5-26-17	5-26-17	
Lead	2800	11	200.8	5-26-17	6-1-17	
Mercury	1.2	0.50	7470A	5-25-17	5-25-17	
Nickel	190	22	200.8	5-26-17	5-26-17	
Selenium	ND	5.6	200.8	5-26-17	5-26-17	
Silver	ND	11	200.8	5-26-17	5-26-17	
Thallium	ND	5.6	200.8	5-26-17	5-26-17	
Zinc	1000	280	200.8	5-26-17	6-1-17	

Project: 15-05986-040

TOTAL METALS EPA 200.8/7470A

Matrix: Water Units: ug/L (ppb)

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-299-15					
Client ID:	PP5-W					
Antimony	ND	5.6	200.8	5-26-17	5-26-17	
Arsenic	14	3.3	200.8	5-26-17	5-26-17	
Beryllium	ND	11	200.8	5-26-17	5-26-17	
Cadmium	ND	4.4	200.8	5-26-17	5-26-17	
Chromium	29	11	200.8	5-26-17	5-26-17	
Copper	50	11	200.8	5-26-17	5-26-17	
Lead	32	1.1	200.8	5-26-17	5-26-17	
Mercury	ND	0.50	7470A	5-25-17	5-25-17	
Nickel	ND	22	200.8	5-26-17	5-26-17	
Selenium	ND	5.6	200.8	5-26-17	5-26-17	
Silver	ND	11	200.8	5-26-17	5-26-17	
Thallium	ND	5.6	200.8	5-26-17	5-26-17	
Zinc	49	28	200.8	5-26-17	5-26-17	

Project: 15-05986-040

TOTAL METALS EPA 200.8/7470A

Matrix: Water Units: ug/L (ppb)

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-299-18					
Client ID:	PP6-W					
Antimony	ND	5.6	200.8	5-26-17	5-26-17	
Arsenic	14	3.3	200.8	5-26-17	5-26-17	
Beryllium	ND	11	200.8	5-26-17	5-26-17	
Cadmium	ND	4.4	200.8	5-26-17	5-26-17	
Chromium	45	11	200.8	5-26-17	5-26-17	
Copper	54	11	200.8	5-26-17	5-26-17	
Lead	55	1.1	200.8	5-26-17	5-26-17	
Mercury	ND	0.50	7470A	5-25-17	5-25-17	
Nickel	46	22	200.8	5-26-17	5-26-17	
Selenium	ND	5.6	200.8	5-26-17	5-26-17	
Silver	ND	11	200.8	5-26-17	5-26-17	
Thallium	ND	5.6	200.8	5-26-17	5-26-17	
Zinc	100	28	200.8	5-26-17	5-26-17	

Project: 15-05986-040

TOTAL METALS EPA 200.8/7470A

Matrix: Water
Units: ug/L (ppb)

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-299-21					
Client ID:	PP7-W					
Antimony	ND	5.6	200.8	5-26-17	5-26-17	
Arsenic	110	3.3	200.8	5-26-17	5-26-17	
Beryllium	ND	11	200.8	5-26-17	5-26-17	
Cadmium	ND	4.4	200.8	5-26-17	5-26-17	
Chromium	210	11	200.8	5-26-17	5-26-17	
Copper	570	22	200.8	5-26-17	6-1-17	
Lead	250	1.1	200.8	5-26-17	5-26-17	
Mercury	0.68	0.50	7470A	5-25-17	5-25-17	
Nickel	210	22	200.8	5-26-17	5-26-17	
Selenium	6.0	5.6	200.8	5-26-17	5-26-17	
Silver	ND	11	200.8	5-26-17	5-26-17	
Thallium	ND	5.6	200.8	5-26-17	5-26-17	
Zinc	520	28	200.8	5-26-17	5-26-17	

Project: 15-05986-040

TOTAL METALS EPA 200.8/7470A

Matrix: Water Units: ug/L (ppb)

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-299-24					
Client ID:	PP8-W					
Antimony	ND	5.6	200.8	5-26-17	5-26-17	
Arsenic	ND	3.3	200.8	5-26-17	5-26-17	
Beryllium	ND	11	200.8	5-26-17	5-26-17	
Cadmium	ND	4.4	200.8	5-26-17	5-26-17	
Chromium	ND	11	200.8	5-26-17	5-26-17	
Copper	ND	11	200.8	5-26-17	5-26-17	
Lead	ND	1.1	200.8	5-26-17	5-26-17	
Mercury	ND	0.50	7470A	5-25-17	5-25-17	
Nickel	ND	22	200.8	5-26-17	5-26-17	
Selenium	ND	5.6	200.8	5-26-17	5-26-17	
Silver	ND	11	200.8	5-26-17	5-26-17	
Thallium	ND	5.6	200.8	5-26-17	5-26-17	
Zinc	71	28	200.8	5-26-17	5-26-17	

Project: 15-05986-040

TOTAL METALS EPA 200.8/7470A

Matrix: Water Units: ug/L (ppb)

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-299-27					
Client ID:	PP9-W					
Antimony	ND	5.6	200.8	5-26-17	5-26-17	
Arsenic	30	3.3	200.8	5-26-17	5-26-17	
Beryllium	ND	11	200.8	5-26-17	5-26-17	
Cadmium	ND	4.4	200.8	5-26-17	5-26-17	
Chromium	49	11	200.8	5-26-17	5-26-17	
Copper	110	11	200.8	5-26-17	5-26-17	
Lead	61	1.1	200.8	5-26-17	5-26-17	
Mercury	ND	0.50	7470A	5-25-17	5-25-17	
Nickel	44	22	200.8	5-26-17	5-26-17	
Selenium	ND	5.6	200.8	5-26-17	5-26-17	
Silver	ND	11	200.8	5-26-17	5-26-17	
Thallium	ND	5.6	200.8	5-26-17	5-26-17	
Zinc	210	28	200.8	5-26-17	5-26-17	

Project: 15-05986-040

TOTAL METALS EPA 200.8/7470A

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-299-29					
Client ID:	PP10-W					
Antimony	ND	5.6	200.8	5-26-17	5-26-17	
Arsenic	ND	3.3	200.8	5-26-17	5-26-17	
Beryllium	ND	11	200.8	5-26-17	5-26-17	
Cadmium	ND	4.4	200.8	5-26-17	5-26-17	
Chromium	ND	11	200.8	5-26-17	5-26-17	
Copper	ND	11	200.8	5-26-17	5-26-17	
Lead	ND	1.1	200.8	5-26-17	5-26-17	
Mercury	ND	0.50	7470A	5-25-17	5-25-17	
Nickel	ND	22	200.8	5-26-17	5-26-17	
Selenium	ND	5.6	200.8	5-26-17	5-26-17	
Silver	ND	11	200.8	5-26-17	5-26-17	
Thallium	ND	5.6	200.8	5-26-17	5-26-17	
Zinc	78	28	200.8	5-26-17	5-26-17	

Project: 15-05986-040

TOTAL METALS EPA 200.8 METHOD BLANK QUALITY CONTROL

Date Extracted: 5-26-17 Date Analyzed: 5-26-17

Matrix: Water
Units: ug/L (ppb)

Lab ID: MB0526WM1

Analyte	Method	Result	PQL
Antimony	200.8	ND	5.6
Arsenic	200.8	ND	3.3
Beryllium	200.8	ND	11
Cadmium	200.8	ND	4.4
Chromium	200.8	ND	11
Copper	200.8	ND	11
Lead	200.8	ND	1.1
Nickel	200.8	ND	22
Selenium	200.8	ND	5.6
Silver	200.8	ND	11
Thallium	200.8	ND	5.6
Zinc	200.8	ND	28

Project: 15-05986-040

TOTAL MERCURY EPA 7470A METHOD BLANK QUALITY CONTROL

Date Extracted: 5-25-17
Date Analyzed: 5-25-17

Matrix: Water
Units: ug/L (ppb)

Lab ID: MB0525W1

Analyte Method Result PQL

Mercury 7470A **ND** 0.50

Project: 15-05986-040

TOTAL METALS EPA 200.8 DUPLICATE QUALITY CONTROL

Date Extracted: 5-26-17
Date Analyzed: 5-26-17

Matrix: Water
Units: ug/L (ppb)

Lab ID: 05-299-29

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Antimony	ND	ND	NA	5.6	
Arsenic	ND	ND	NA	3.3	
Beryllium	ND	ND	NA	11	
Cadmium	ND	ND	NA	4.4	
Chromium	ND	ND	NA	11	
Copper	ND	ND	NA	11	
Lead	ND	ND	NA	1.1	
Nickel	ND	ND	NA	22	
Selenium	ND	ND	NA	5.6	
Silver	ND	ND	NA	11	
Thallium	ND	ND	NA	5.6	
Zinc	78.3	76.7	2	28	

Project: 15-05986-040

TOTAL MERCURY EPA 7470A DUPLICATE QUALITY CONTROL

Date Extracted: 5-25-17
Date Analyzed: 5-25-17

Matrix: Water
Units: ug/L (ppb)

Lab ID: 05-299-02

Analyte Sample Duplicate
Result Result RPD PQL Flags

Mercury ND ND NA 0.50

Project: 15-05986-040

TOTAL METALS EPA 200.8 MS/MSD QUALITY CONTROL

Date Extracted: 5-26-17
Date Analyzed: 5-26-17

Matrix: Water
Units: ug/L (ppb)

Lab ID: 05-299-29

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Antimony	222	246	111	246	111	0	
Arsenic	222	247	111	247	111	0	
Beryllium	222	233	105	235	106	1	
Cadmium	222	240	108	239	108	0	
Chromium	222	219	99	224	101	3	
Copper	222	218	98	222	100	2	
Lead	222	242	109	238	107	1	
Nickel	222	221	99	225	101	2	
Selenium	222	271	122	266	120	2	
Silver	222	223	100	223	100	0	
Thallium	222	246	111	244	110	1	
Zinc	222	317	108	329	113	4	

Project: 15-05986-040

TOTAL MERCURY EPA 7470A MS/MSD QUALITY CONTROL

Date Extracted: 5-25-17
Date Analyzed: 5-25-17

Matrix: Water
Units: ug/L (ppb)

Lab ID: 05-299-02

	Spike		Percent		Percent		
Analyte	Level	MS	Recovery	MSD	Recovery	RPD	Flags
Mercury	12.5	13.0	104	13.0	104	0	

Project: 15-05986-040

DISSOLVED METALS EPA 200.8/7470A

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-299-02 PP1-W					
Antimony	ND	5.0	200.8		5-26-17	
Arsenic	18	3.0	200.8		5-26-17	
Beryllium	ND	10	200.8		5-26-17	
Cadmium	ND	4.0	200.8		5-26-17	
Chromium	ND	10	200.8		5-26-17	
Copper	ND	10	200.8		5-26-17	
Lead	ND	1.0	200.8		5-26-17	
Mercury	ND	0.50	7470A		5-25-17	
Nickel	ND	20	200.8		5-26-17	
Selenium	ND	5.0	200.8		5-26-17	
Silver	ND	10	200.8		5-26-17	
Thallium	ND	5.0	200.8		5-26-17	
Zinc	39	25	200.8		5-26-17	

Project: 15-05986-040

DISSOLVED METALS EPA 200.8/7470A

	- " ' '			Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-299-06					
Client ID:	PP2-W					
Antimony	ND	5.0	200.8		5-26-17	
Arsenic	28	3.0	200.8		5-26-17	
Beryllium	ND	10	200.8		5-26-17	
Cadmium	ND	4.0	200.8		5-26-17	
Chromium	ND	10	200.8		5-26-17	
Copper	ND	10	200.8		5-26-17	
Lead	1.2	1.0	200.8		5-26-17	
Mercury	ND	0.50	7470A		5-25-17	
Nickel	ND	20	200.8		5-26-17	
Selenium	ND	5.0	200.8		5-26-17	
Silver	ND	10	200.8		5-26-17	
Thallium	ND	5.0	200.8		5-26-17	
Zinc	ND	25	200.8		5-26-17	

Project: 15-05986-040

DISSOLVED METALS EPA 200.8/7470A

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-299-09					
Client ID:	PP3-W					
Antimony	ND	5.0	200.8		5-26-17	
Arsenic	3.5	3.0	200.8		5-26-17	
Beryllium	ND	10	200.8		5-26-17	
Cadmium	ND	4.0	200.8		5-26-17	
Chromium	ND	10	200.8		5-26-17	
Copper	ND	10	200.8		5-26-17	
Lead	ND	1.0	200.8		5-26-17	
Mercury	ND	0.50	7470A		5-25-17	
Nickel	ND	20	200.8		5-26-17	
Selenium	ND	5.0	200.8		5-26-17	
Silver	ND	10	200.8		5-26-17	
Thallium	ND	5.0	200.8		5-26-17	
Zinc	ND	25	200.8		5-26-17	

Project: 15-05986-040

DISSOLVED METALS EPA 200.8/7470A

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-299-12					
Client ID:	PP4-W					
Antimony	7.1	5.0	200.8		5-26-17	
Arsenic	7.1	3.0	200.8		5-26-17	
Beryllium	ND	10	200.8		5-26-17	
Cadmium	ND	4.0	200.8		5-26-17	
Chromium	ND	10	200.8		5-26-17	
Copper	ND	10	200.8		5-26-17	
Lead	3.1	1.0	200.8		5-26-17	
Mercury	ND	0.50	7470A		5-25-17	
Nickel	ND	20	200.8		5-26-17	
Selenium	ND	5.0	200.8		5-26-17	
Silver	ND	10	200.8		5-26-17	
Thallium	ND	1.0	200.8		5-26-17	
Zinc	ND	25	200.8		5-26-17	

Project: 15-05986-040

DISSOLVED METALS EPA 200.8/7470A

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-299-15					
Client ID:	PP5-W					
Antimony	ND	5.0	200.8		5-26-17	
Arsenic	3.3	3.0	200.8		5-26-17	
Beryllium	ND	10	200.8		5-26-17	
Cadmium	ND	4.0	200.8		5-26-17	
Chromium	ND	10	200.8		5-26-17	
Copper	ND	10	200.8		5-26-17	
Lead	ND	1.0	200.8		5-26-17	
Mercury	ND	0.50	7470A		5-25-17	
Nickel	ND	20	200.8		5-26-17	
Selenium	ND	5.0	200.8		5-26-17	
Silver	ND	10	200.8		5-26-17	
Thallium	ND	1.0	200.8		5-26-17	
Zinc	ND	25	200.8		5-26-17	

Project: 15-05986-040

DISSOLVED METALS EPA 200.8/7470A

	- " ' '			Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-299-18					
Client ID:	PP6-W					
Antimony	ND	5.0	200.8		5-26-17	
Arsenic	3.6	3.0	200.8		5-26-17	
Beryllium	ND	10	200.8		5-26-17	
Cadmium	ND	4.0	200.8		5-26-17	
Chromium	ND	10	200.8		5-26-17	
Copper	ND	10	200.8		5-26-17	
Lead	ND	1.0	200.8		5-26-17	
Mercury	ND	0.50	7470A		5-25-17	
Nickel	ND	20	200.8		5-26-17	
Selenium	ND	5.0	200.8		5-26-17	
Silver	ND	10	200.8		5-26-17	
Thallium	ND	1.0	200.8		5-26-17	
Zinc	ND	25	200.8		5-26-17	

Project: 15-05986-040

DISSOLVED METALS EPA 200.8/7470A

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-299-21					
Client ID:	PP7-W					
Antimony	ND	5.0	200.8		5-26-17	
Arsenic	ND	3.0	200.8		5-26-17	
Beryllium	ND	10	200.8		5-26-17	
Cadmium	ND	4.0	200.8		5-26-17	
Chromium	ND	10	200.8		5-26-17	
Copper	ND	10	200.8		5-26-17	
Lead	ND	1.0	200.8		5-26-17	
Mercury	ND	0.50	7470A		5-25-17	
Nickel	ND	20	200.8		5-26-17	
Selenium	ND	5.0	200.8		5-26-17	
Silver	ND	10	200.8		5-26-17	
Thallium	ND	5.0	200.8		5-26-17	
Zinc	ND	25	200.8		5-26-17	

Project: 15-05986-040

DISSOLVED METALS EPA 200.8/7470A

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-299-24					
Client ID:	PP8-W					
Antimony	ND	5.0	200.8		5-26-17	
Arsenic	ND	3.0	200.8		5-26-17	
Beryllium	ND	10	200.8		5-26-17	
Cadmium	ND	4.0	200.8		5-26-17	
Chromium	ND	10	200.8		5-26-17	
Copper	ND	10	200.8		5-26-17	
Lead	ND	1.0	200.8		5-26-17	
Mercury	ND	0.50	7470A		5-25-17	
Nickel	ND	20	200.8		5-26-17	
Selenium	ND	5.0	200.8		5-26-17	
Silver	ND	10	200.8		5-26-17	
Thallium	ND	5.0	200.8		5-26-17	
Zinc	62	25	200.8		5-26-17	

Project: 15-05986-040

DISSOLVED METALS EPA 200.8/7470A

				Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-299-27					
Client ID:	PP9-W					
Antimony	ND	5.0	200.8		5-26-17	
Arsenic	ND	3.0	200.8		5-26-17	
Beryllium	ND	10	200.8		5-26-17	
Cadmium	ND	4.0	200.8		5-26-17	
Chromium	ND	10	200.8		5-26-17	
Copper	ND	10	200.8		5-26-17	
Lead	ND	1.0	200.8		5-26-17	
Mercury	ND	0.50	7470A		5-25-17	
Nickel	ND	20	200.8		5-26-17	
Selenium	ND	5.0	200.8		5-26-17	
Silver	ND	10	200.8		5-26-17	
Thallium	ND	1.0	200.8		5-26-17	
Zinc	ND	25	200.8		5-26-17	

Project: 15-05986-040

DISSOLVED METALS EPA 200.8/7470A

	- " ' '			Date	Date	
Analyte	Result	PQL	EPA Method	Prepared	Analyzed	Flags
Lab ID:	05-299-29					
Client ID:	PP10-W					
Antimony	ND	5.0	200.8		5-26-17	
Arsenic	ND	3.0	200.8		5-26-17	
Beryllium	ND	10	200.8		5-26-17	
Cadmium	ND	4.0	200.8		5-26-17	
Chromium	ND	10	200.8		5-26-17	
Copper	ND	10	200.8		5-26-17	
Lead	ND	1.0	200.8		5-26-17	
Mercury	ND	0.50	7470A		5-25-17	
Nickel	ND	20	200.8		5-26-17	
Selenium	ND	5.0	200.8		5-26-17	
Silver	ND	10	200.8		5-26-17	
Thallium	ND	1.0	200.8		5-26-17	
Zinc	64	25	200.8		5-26-17	

Project: 15-05986-040

DISSOLVED METALS EPA 200.8 METHOD BLANK QUALITY CONTROL

Date Analyzed: 5-26-17

Matrix: Water Units: ug/L (ppb)

Lab ID: MB0524F1

Analyte	Method	Result	PQL
Antimony	200.8	ND	5.0
Arsenic	200.8	ND	3.0
Beryllium	200.8	ND	10
Cadmium	200.8	ND	4.0
Chromium	200.8	ND	10
Copper	200.8	ND	10
Lead	200.8	ND	1.0
Nickel	200.8	ND	20
Selenium	200.8	ND	5.0
Silver	200.8	ND	10
Thallium	200.8	ND	5.0
Zinc	200.8	ND	25

Project: 15-05986-040

DISSOLVED MERCURY EPA 7470A METHOD BLANK QUALITY CONTROL

Date Analyzed: 5-25-17

Matrix: Water
Units: ug/L (ppb)

Lab ID: MB0524F1

Analyte Method Result PQL

Mercury 7470A **ND** 0.50

Project: 15-05986-040

DISSOLVED METALS EPA 200.8 DUPLICATE QUALITY CONTROL

Date Analyzed: 5-26-17

Matrix: Water
Units: ug/L (ppb)

Lab ID: 05-299-29

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Antimony	ND	ND	NA	5.0	
Arsenic	ND	ND	NA	3.0	
Beryllium	ND	ND	NA	10	
Cadmium	ND	ND	NA	4.0	
Chromium	ND	ND	NA	10	
Copper	ND	ND	NA	10	
Lead	ND	ND	NA	1.0	
Nickel	ND	ND	NA	20	
Selenium	ND	ND	NA	5.0	
Silver	ND	ND	NA	10	
Thallium	ND	ND	NA	5.0	
Zinc	64.0	64.5	1	25	

Project: 15-05986-040

DISSOLVED MERCURY EPA 7470A DUPLICATE QUALITY CONTROL

Date Analyzed: 5-25-17

Matrix: Water
Units: ug/L (ppb)

Lab ID: 05-299-02

Sample Duplicate

Analyte Result Result RPD PQL Flags

Mercury ND ND NA 0.5

Project: 15-05986-040

DISSOLVED METALS EPA 200.8 MS/MSD QUALITY CONTROL

Date Analyzed: 5-26-17

Matrix: Water
Units: ug/L (ppb)

Lab ID: 05-299-29

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Antimony	200	199	99	208	104	5	
Arsenic	200	201	101	208	104	3	
Beryllium	200	191	95	200	100	5	
Cadmium	200	199	100	205	102	3	
Chromium	200	183	92	188	94	2	
Copper	200	182	91	184	92	1	
Lead	200	194	97	200	100	3	
Nickel	200	184	92	189	94	3	
Selenium	200	214	107	224	112	5	
Silver	200	183	91	196	98	7	
Thallium	200	185	93	204	102	10	
Zinc	200	271	103	274	105	1	

Project: 15-05986-040

DISSOLVED MERCURY EPA 7470A MS/MSD QUALITY CONTROL

Date Analyzed: 5-25-17

Matrix: Water
Units: ug/L (ppb)

Lab ID: 05-299-02

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Mercury	12.5	12.6	100	12.4	99	1	

Project: 15-05986-040

PCBs EPA 8082A

Matrix: Soil

Units: mg/Kg (ppm)

• • •				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP3-2.5					
Laboratory ID:	05-299-07					
Aroclor 1016	ND	0.063	EPA 8082A	6-6-17	6-6-17	
Aroclor 1221	ND	0.063	EPA 8082A	6-6-17	6-6-17	
Aroclor 1232	ND	0.063	EPA 8082A	6-6-17	6-6-17	
Aroclor 1242	ND	0.063	EPA 8082A	6-6-17	6-6-17	
Aroclor 1248	ND	0.063	EPA 8082A	6-6-17	6-6-17	
Aroclor 1254	ND	0.063	EPA 8082A	6-6-17	6-6-17	
Aroclor 1260	ND	0.063	EPA 8082A	6-6-17	6-6-17	
Surrogate:	Percent Recovery	Control Limits				
DCB	90	42-139				
Client ID:	PP4-3					
Laboratory ID:	05-299-10					
Aroclor 1016	ND	0.063	EPA 8082A	6-6-17	6-6-17	
Aroclor 1221	ND	0.063	EPA 8082A	6-6-17	6-6-17	
Aroclor 1232	ND	0.063	EPA 8082A	6-6-17	6-6-17	
Aroclor 1242	ND	0.063	EPA 8082A	6-6-17	6-6-17	
Aroclor 1248	ND	0.063	EPA 8082A	6-6-17	6-6-17	
Aroclor 1254	ND	0.063	EPA 8082A	6-6-17	6-6-17	
Aroclor 1260	0.18	0.063	EPA 8082A	6-6-17	6-6-17	
Surrogate:	Percent Recovery	Control Limits				
DCB	88	42-139				
Client ID:	PP4-10					
Laboratory ID:	05-299-11					
Aroclor 1016	ND	0.062	EPA 8082A	6-6-17	6-6-17	
Aroclor 1221	ND	0.062	EPA 8082A	6-6-17	6-6-17	
Aroclor 1232	ND	0.062	EPA 8082A	6-6-17	6-6-17	
Aroclor 1242	ND	0.062	EPA 8082A	6-6-17	6-6-17	
Aroclor 1248	ND	0.062	EPA 8082A	6-6-17	6-6-17	
Aroclor 1254	ND	0.062	EPA 8082A	6-6-17	6-6-17	
Aroclor 1260	ND	0.062	EPA 8082A	6-6-17	6-6-17	
Surrogate:	Percent Recovery	Control Limits				
DCB	91	42-139				

Project: 15-05986-040

PCBs EPA 8082A

Matrix: Soil

Units: mg/Kg (ppm)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP5-2.5					
Laboratory ID:	05-299-13					
Aroclor 1016	ND	0.065	EPA 8082A	6-6-17	6-6-17	
Aroclor 1221	ND	0.065	EPA 8082A	6-6-17	6-6-17	
Aroclor 1232	ND	0.065	EPA 8082A	6-6-17	6-6-17	
Aroclor 1242	ND	0.065	EPA 8082A	6-6-17	6-6-17	
Aroclor 1248	ND	0.065	EPA 8082A	6-6-17	6-6-17	
Aroclor 1254	ND	0.065	EPA 8082A	6-6-17	6-6-17	
Aroclor 1260	ND	0.065	EPA 8082A	6-6-17	6-6-17	
Surrogate:	Percent Recovery	Control Limits				
DCB	80	42-139				
Client ID:	PP6-2					
Laboratory ID:	05-299-16					
Aroclor 1016	ND	0.058	EPA 8082A	6-6-17	6-6-17	
Aroclor 1221	ND	0.058	EPA 8082A	6-6-17	6-6-17	
Aroclor 1232	ND	0.058	EPA 8082A	6-6-17	6-6-17	
Aroclor 1242	ND	0.058	EPA 8082A	6-6-17	6-6-17	
Aroclor 1248	ND	0.058	EPA 8082A	6-6-17	6-6-17	
Aroclor 1254	0.12	0.058	EPA 8082A	6-6-17	6-6-17	
Aroclor 1260	ND	0.058	EPA 8082A	6-6-17	6-6-17	
Surrogate:	Percent Recovery	Control Limits				
DCB	84	42-139				
Client ID:	PP7-0					
Laboratory ID:	05-299-19					
Aroclor 1016	ND	0.061	EPA 8082A	6-6-17	6-6-17	
Aroclor 1221	ND	0.061	EPA 8082A	6-6-17	6-6-17	
Aroclor 1232	ND	0.061	EPA 8082A	6-6-17	6-6-17	
Aroclor 1242	ND	0.061	EPA 8082A	6-6-17	6-6-17	
Aroclor 1248	ND	0.061	EPA 8082A	6-6-17	6-6-17	
Aroclor 1254	ND	0.061	EPA 8082A	6-6-17	6-6-17	
Aroclor 1260	ND	0.061	EPA 8082A	6-6-17	6-6-17	
Surrogate:	Percent Recovery	Control Limits				
DCB	77	42-139				

Project: 15-05986-040

PCBs EPA 8082A QUALITY CONTROL

Matrix: Soil

Units: mg/Kg (ppm)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0606S1					
Aroclor 1016	ND	0.050	EPA 8082A	6-6-17	6-6-17	
Aroclor 1221	ND	0.050	EPA 8082A	6-6-17	6-6-17	
Aroclor 1232	ND	0.050	EPA 8082A	6-6-17	6-6-17	
Aroclor 1242	ND	0.050	EPA 8082A	6-6-17	6-6-17	
Aroclor 1248	ND	0.050	EPA 8082A	6-6-17	6-6-17	
Aroclor 1254	ND	0.050	EPA 8082A	6-6-17	6-6-17	
Aroclor 1260	ND	0.050	EPA 8082A	6-6-17	6-6-17	
<u> </u>	D 15	0 , 11: "	•	•		

Surrogate: Percent Recovery Control Limits DCB 97 42-139

Analyte	Re	sult	Spike	Level	Source Result		rcent covery	Recovery Limits	RPD	RPD Limit	Flags
MATRIX SPIKES											
Laboratory ID:	06-0	20-14									
	MS	MSD	MS	MSD		MS	MSD				
Aroclor 1260	0.340	0.307	0.500	0.500	ND	68	61	26-127	10	22	
Surrogate:											
DCB						73	70	42-139			

Project: 15-05986-040

% MOISTURE

Date Analyzed: 5-25-17

Client ID	Lab ID	% Moisture
PP1-2	05-299-01	19
PP1-10	05-299-03	25
PP2-2.5	05-299-04	24
PP2-10	05-299-05	31
PP3-2.5	05-299-07	20
PP3-10	05-299-08	19
PP4-3	05-299-10	21
PP4-10	05-299-11	19
PP5-2.5	05-299-13	24
PP5-10	05-299-14	26
PP6-2	05-299-16	14
PP6-10	05-299-17	33
PP7-0	05-299-19	18
PP7-10	05-299-20	20
PP8-0.5	05-299-22	12
PP8-10	05-299-23	30
PP9-0	05-299-25	4
PP9-10	05-299-26	19
PP10-0	05-299-28	18



Data Qualifiers and Abbreviations

- A Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
- B The analyte indicated was also found in the blank sample.
- C The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
- E The value reported exceeds the quantitation range and is an estimate.
- F Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
- H The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
- I Compound recovery is outside of the control limits.
- J The value reported was below the practical quantitation limit. The value is an estimate.
- K Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
- L The RPD is outside of the control limits.
- M Hydrocarbons in the gasoline range are impacting the diesel range result.
- M1 Hydrocarbons in the gasoline range (toluene-napthalene) are present in the sample.
- N Hydrocarbons in the lube oil range are impacting the diesel range result.
- N1 Hydrocarbons in diesel range are impacting lube oil range results.
- O Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
- P The RPD of the detected concentrations between the two columns is greater than 40.
- Q Surrogate recovery is outside of the control limits.
- S Surrogate recovery data is not available due to the necessary dilution of the sample.
- T The sample chromatogram is not similar to a typical .
- U The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- U1 The practical quantitation limit is elevated due to interferences present in the sample.
- V Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
- W Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
- X Sample extract treated with a mercury cleanup procedure.
- X1- Sample extract treated with a Sulfuric acid/Silica gel cleanup procedure.
- Y The calibration verification for this analyte exceeded the 20% drift specified in method 8260C, and therefore the reported result should be considered an estimate. The overall performance of the calibration verification standard met the acceptance criteria of the method.

7 -

ND - Not Detected at PQL

PQL - Practical Quantitation Limit

RPD - Relative Percent Difference





Chain of Custody

Analytical Laboratory Testing Services 14648 NE 95th Street • Redmond, WA 98052	Turnaround Req (in working da		Ì	Labo	orato	ory	Nun	nbe	r:	0 5	5 -	25	9									
Phone: (425) 883-3881 · www.onsite-env.com Company: Herrera Environ mental Project Number: 15-05986-040 Project Name: Pacific Park Project Manager: Broce Carpenter Sampled by: Broce Carpenter Brianna Bland		1 Day 3 Days	5	TEX		NWTPH-DX VI Acid / SG Clean-up)	3260C	260C	(Waters Only)	PAHS)	SIM (low-level)	ne Pesticides 80818	arus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	etals	HEM (oil and grease) 1664A	I Priority Pollestant Mades	ssolva P.P. Metals		ure
Lab ID Sample Identification	Date Time Sampled Sampled	Matrix 2	Namue	NWTPH-HCID NWTPH-Gx/B	NWTPH-Gx	NWTP	Volatiles 8	Halogenated	Semitor	(with low-level	PAHs 8	PCBs 8082A	Organo	Chloring	Total RC	Total M	TCLP Metals	HEM (o	Total	D SS		% Moisture
1 PPI-2	5/23/17 9:20	5 6	5		X	X	X								X							X
2 PPI-W	5/23/17 10:00	WI	1		X	X	X				X								X	X		
3 PP1-10	5/23/17 9:45	5 6			X	X	X								X							X
4 PP2-2.5	5/23/17 10:35	5 6	,		X	X	X			X					X							X
5 PP2-10	5/23/17 10:50	S 6	,		X	X	X								X							X
6 PP2-W	5/23/17 11:00	W. C	7		X	X	X				X			H					X	X		X
7 PP3-2.5	5/24/17 10:05	5 4	5		X	X	X				6	()		1	X						1	X
8 PP3-10	5/24/17 10:15		2		X	X	X			X					X							X
9 PP3-W	5/24/17 10:00	1			X	X	X		7		X			1	X	ik			X	X		
10 1 -	5/24/17 8:30				X	X	X			X	6	3)			X							X
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Page 2 of 3

Analytical Laboratory Testing Services 14648 NE 95th Street • Redmond, WA 98052	Turnaround Request (in working days) Laboratory Number					er:	0	5	-2	9	9											
Phone: (425) 883-3881 · www.onsite-env.com Company: Herrera Env. ron mental Project Number: 15-05986 - o 40 Project Name: Project Manager: Brue Carpenter Sampled by: Brue Carpenter Brue Carpenter Brianna Hard	Date Time	Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX	NWTPH-GX	5	Halogenated Volatiles 8260C	EPA 801	Semivolatiles 8270D/SIM (with low-level PAHs)	PAHs 8270D/SIM (low-level)	PCBs 8082A	Organochlorine Pesticides 8081B	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	fotal MTCA Metals	TCLP Metals	HEM (oil and grease) 1664A	otal P.P. Metals	wed P.P. Meta		Moisture
Lab ID Sample Identification PP4-18	Sampled Sampled Matrix 5/24/17 8:46 5	6	Z	Z	2 2	* * * * * * * * * * * * * * * * * * *	Ĩ	Ш	X	4		0	0	0	X	12	F	エ		l		× ×
12 PPH-W	5/24/17 9:00 W	11			X	X				X				1					X	X		
13 PP5-2.5	5/23/17 15:40 S	6			$\stackrel{*}{\bigwedge}$	$\langle \! \rangle$			X	(X)			X							X
14 PPS- 10	5/23/17 15:50 5	6			X	X		1	X						X							X
15 PP5-W	5/33/17 16:00 W	11			$\langle \rangle$	X				X									X	X		
16 PP6-2	5/23/17 14:35 S	6			\bigvee	X			X	(8				X							X
17 PP6-10	5/23/17 14.45 5	6			$\langle \rangle$	\otimes			X		R				X							X
18 PP6-W	5/23/17 15:00 W	11			${\otimes}$	\otimes				X									X	X		
19 PP7-0	5/24/17 14:30 5	6			\Diamond	\propto			X	($\langle \hat{\mathbf{x}} \rangle$				X							X
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Page _3_ of _3_

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Company: Herrera F Project Number: 15-059 Project Name: Pacific Project Manager: Sampled by: Brue Care	26-040 Parte Carpenter Briance Bland ample Identification	☐ 2 E	Days [andard (7 Days) PH analysis 5 Day (other)	1 Day 3 Days ys) Matrix	Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX	NWTPH-Gx	NWTPH-Dx (X Acid / SG Clean-up)	voiatiles 8260C Halogenated Volatiles 8260C	EDB EPA 8011 (Waters Only)	Semivolatiles 8270D/SIM (with low-level PAHs)	PAHs 8270D/SIM (low-level)	PCBs 8082A	Organochlorine Pesticides 8081B	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	TCLP Metals	HEM (oil and grease) 1664A	Total Priort, Pollotatualis	resolved TP Metals		6 Moisture
21 PP7-W	ample lucininication	5/24/1		W	III		_	X	X	1	1	Fr.	X			0	0	Е	F		_	X	X		%
22 PP8 · 0	.5	5/24/	17 13:20	5	6			X	X	1		X		1				X			1				X
23 PP8-1		5/24/1	7 13:30	5	6			X	X	1		X						X				П			X
24 PP8-0		5/24/1	1 1 1 1 1 1	W	11			X	X			X	X									X	X		
25 TP9-0		5/24/1	711:50	5	6			X		X		X						X							X
26 PP9-1	0	5/24/1	7 121.00	5.	6			X	X	\langle		X						X							X
27 PP9-1	J	5/24/1	7 12:10	W.	11			\bigvee					X									X	X		
28 PP 10-	0	5/24/1	7 15:10	5.	6			$\stackrel{\checkmark}{}$	$\langle \rangle$			X						X							X
29 PP10-		5/24/1	7 13:45	W	11			\swarrow	\Diamond				\times									\times	X		
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APPENDIX E

Data Quality Assurance Review



Herrera Environmental Consultants, Inc.

Memorandum

To Project File 15-05986-040

From Gina Catarra, Herrera Environmental Consultants

Date July 18, 2017

Subject Data Quality Assurance Review of the Pacific Park/Dumpsite Data

This memorandum presents a review of data quality for 19 soil samples (including one field duplicate), 17 groundwater samples (including one field duplicate), one filter blank sample, and two trip blank samples collected for the Pacific Park/Dumpsite property between May 12 and 24, 2017. OnSite Environmental, Inc., of Redmond, Washington analyzed the samples for:

- Gasoline-range petroleum hydrocarbons (TPH-G) by Ecology's NWPTH-Gx method
- Diesel- and lube oil-range petroleum hydrocarbons by Ecology's NWTPH-Dx method
- Volatile organic compounds (VOCs) by EPA Method 8260C
- Semivolatile organic compounds (SVOCs) by EPA Method 8270D/SIM (only soil samples)
- Polycyclic aromatic hydrocarbons (PAHs) by EPA Method 8270D/SIM (only groundwater samples)
- Polychlorinated biphenyls (PCBs) by EPA Method 8082A
- Total Resource Conservation and Recovery Act (RCRA) metals (arsenic, barium, cadmium, chromium, lead, mercury, selenium, and silver) by EPA 6010C/7471B (soil samples)
- Total and dissolved Priority Pollutant metals (antimony, arsenic, beryllium, cadmium, chromium, copper, lead, mercury, nickel, selenium, silver, thallium, and zinc) by EPA Method 200.8/7470A.

Results for the following samples were validated.

Sample ID	Lab SDG	Date Collected	Matrix	Analyses
MW1	1705-184	5/12/2017	Groundwater	TPH-G, TPH-D, VOCs, PAHs, total/dissolved metals
MW2	1705-184	5/12/2017	Groundwater	TPH-G, TPH-D, VOCs, PAHs, total/dissolved metals
MW3	1705-184	5/12/2017	Groundwater	TPH-G, TPH-D, VOCs, PAHs, total/dissolved metals
MW4	1705-184	5/12/2017	Groundwater	TPH-G, TPH-D, VOCs, PAHs, total/dissolved metals
MW5	1705-184	5/12/2017	Groundwater	TPH-G, TPH-D, VOCs, PAHs, total/dissolved metals
MW6	1705-184	5/12/2017	Groundwater	TPH-G, TPH-D, VOCs, PAHs, total/dissolved metals
$MWDUP^a$	1705-184	5/12/2017	Groundwater	TPH-G, TPH-D, VOCs, PAHs, total/dissolved metals
Filter Blank ^b	1705-184	5/12/2017	Water	Metals
Trip Blank ^c	1705-184	5/12/2017	Water	TPH-G, VOCs
PP1-2	17-05-299	5/23/2017	Soil	TPH-G, TPH-D, VOCs, SVOCs, total metals
PP1-W	17-05-299	5/23/2017	Groundwater	TPH-G, TPH-D, VOCs, PAHs, total/dissolved metals
PP1-10	17-05-299	5/23/2017	Soil	TPH-G, TPH-D, VOCs, SVOCs, total metals
PP2-2.5	17-05-299	5/23/2017	Soil	TPH-G, TPH-D, VOCs, SVOCs, total metals
PP2-10	17-05-299	5/23/2017	Soil	TPH-G, TPH-D, VOCs, SVOCs, total metals
PP2-W	17-05-299	5/23/2017	Groundwater	TPH-G, TPH-D, VOCs, PAHs, total/dissolved metals
PP3-2.5	17-05-299	5/23/2017	Soil	TPH-G, TPH-D, VOCs, SVOCs, PCBs, total metals
PP3-10	17-05-299	5/23/2017	Soil	TPH-G, TPH-D, VOCs, SVOCs, total metals
PP3-W	17-05-299	5/23/2017	Groundwater	TPH-G, TPH-D, VOCs, PAHs, total/dissolved metals
PP4-3	17-05-299	5/24/2017	Soil	TPH-G, TPH-D, VOCs, SVOCs, PCBs, total metals
PP4-10	17-05-299	5/24/2017	Soil	TPH-G, TPH-D, VOCs, SVOCs, PCBs, total metals
PP4-W	17-05-299	5/24/2017	Groundwater	TPH-G, TPH-D, VOCs, PAHs, total/dissolved metals
PP5-2.5	17-05-299	5/23/2017	Soil	TPH-G, TPH-D, VOCs, SVOCs, PCBs, total metals
PP5-10	17-05-299	5/23/2017	Soil	TPH-G, TPH-D, VOCs, SVOCs, total metals
PP5-W	17-05-299	5/23/2017	Groundwater	TPH-G, TPH-D, VOCs, PAHs, total/dissolved metals
PP6-2	17-05-299	5/23/2017	Soil	TPH-G, TPH-D, VOCs, SVOCs, PCBs, total metals
PP6-10	17-05-299	5/23/2017	Soil	TPH-G, TPH-D, VOCs, SVOCs, total metals
PP6-W	17-05-299	5/23/2017	Groundwater	TPH-G, TPH-D, VOCs, PAHs, total/dissolved metals
PP7-0	17-05-299	5/24/2017	Soil	TPH-G, TPH-D, VOCs, SVOCs, PCBs, total metals
PP7-10	17-05-299	5/24/2017	Soil	TPH-G, TPH-D, VOCs, SVOCs, total metals
PP7-W	17-05-299	5/24/2017	Groundwater	TPH-G, TPH-D, VOCs, PAHs, total/dissolved metals
PP8-0.5	17-05-299	5/24/2017	Soil	TPH-G, TPH-D, VOCs, SVOCs, total metals
PP8-10	17-05-299	5/24/2017	Soil	TPH-G, TPH-D, VOCs, SVOCs, total metals
PP8-W	17-05-299	5/24/2017	Groundwater	TPH-G, TPH-D, VOCs, PAHs, total/dissolved metals
PP9-0	17-05-299	5/24/2017	Soil	TPH-G, TPH-D, VOCs, SVOCs, total metals
PP9-10	17-05-299	5/24/2017	Soil	TPH-G, TPH-D, VOCs, SVOCs, total metals
PP9-W	17-05-299	5/24/2017	Groundwater	TPH-G, TPH-D, VOCs, PAHs, total/dissolved metals
PP10-0 ^d	17-05-299	5/24/2017	Soil	TPH-G, TPH-D, VOCs, SVOCs, total metals
PP10-W ^e	17-05-299	5/24/2017	Groundwater	TPH-G, TPH-D, VOCs, PAHs, total/dissolved metals
Trip Blank ^c	17-05-299	5/24/2017	Water	TPG-G, VOCs

Field duplicate of MW2
Field filter blank
Trip blank
Field duplicate of PP7-10
Field duplicate of PP7-W

The data were reviewed using guidance and quality control criteria documented in the *Pacific Park/Dumpsite Environmental and Economic Assessment Sampling and Analysis Plan* (SAP) (Herrera 2017), and in accordance with quality control (QC) criteria established by the laboratory and in the specified methods.

Quality control data summaries submitted by the laboratories were reviewed; raw data were not submitted by the laboratories. Data qualifiers (flags) were added to the sample results in the laboratory reports. Data validation results are summarized below, followed by definitions of data qualifiers.

Custody, Preservation, Holding Times, and Completeness—Acceptable

The samples were properly preserved and sample custody was maintained from sample collection to receipt at the laboratories. All samples were analyzed within the required method holding time. The laboratory report was complete and contained results for all samples and tests requested on the chain-of-custody (COC) forms.

Laboratory Reporting Limits—Acceptable

The laboratory reporting limits met those specified in the SAP.

Method Blank Analysis—Acceptable

Method blanks were analyzed at the required frequency for all analyses. Method blanks did not contain levels of target analytes above the laboratory reporting limits.

Trip Blank Analysis—Acceptable

Trip blanks were analyzed for gasoline and VOCs at the required frequency. Trip blanks did not contain levels of target analytes above the laboratory reporting limits.

Filter Blank Analysis—Acceptable

A filter blank was analyzed with the dissolved metals as required by the SAP. The filter blank did not contain levels of target analytes above the laboratory reporting limits

Laboratory Control Sample Analysis—Acceptable with Discussion

Blank spike (BS) or blank spike/blank spike duplicate (BS/BSD) samples were analyzed with project samples for VOCs and PAHs at the required frequency. With the exception noted below, the percent recovery values met the control limits established by the SAP.

The BS percent recovery for chrysene (112 percent) exceeded the upper control limit of 110 percent. Chrysene was detected above the reporting limits in three associated samples (PP2-W, PP4-W, and PP6-W). However, no data were qualified because the exceedance was marginal (2 percent) and the percent recovery in the BSD (109 percent) was within control limits.

Surrogate Analysis—Acceptable

Surrogate compounds were added to all samples as required by the specified methods. The percent recovery values for all surrogate compounds met the criteria established by the laboratory or specified method.

Matrix Spike Analysis—Acceptable

Matrix spike/matrix spike duplicate (MS/MSD) samples were analyzed with groundwater samples for VOCs (batch samples) and total/dissolved metals (sample MW1), and soil samples for SVOCs (sample PP1-10) and metals (sample PP1-0). All percent recovery values met those established in the SAP.

Laboratory Duplicate Analysis—Acceptable

Project-sample laboratory duplicates were analyzed for with groundwater samples for TPH-G, TPH-D, and total/dissolved metals, and with soil samples for TPH-G, TPH-D, and metals. The relative percent difference (RPD) was calculated for each analyte where both duplicate values were greater than five times the reporting limit (RL). The RPD values met the control limits established in the SAP.

Field Duplicate Analysis—Acceptable

Field duplicates were analyzed for groundwater samples MW2 and PP7-W, and soil sample PP7-10. The RPD was calculated for each analyte where both the original sample and duplicate values were greater than five times the reporting limit (RL). The RPD values met the control limits established in the SAP.

Definition of Data Qualifiers

The following data qualifier definitions are taken from *USEPA National Functional Guidelines* for *Inorganic Superfund Data Review* (USEPA 2014).

- U The analyte was analyzed for, but was not detected above the level of the reported sample quantitation limit.
- J The result is an estimated quantity. The associated numerical value is the approximate concentration of the analyte in the sample.
- **UJ** The material was analyzed for, but was not detected. The reported quantitation limit is approximate and may be inaccurate or imprecise.
- **R** The data are unusable. (Note: analyte may or may not be present.)

References

Herrera. 2017. Pacific Park/Dumpsite Environmental and Economic Assessment, Pacific, Washington, Sampling and Analysis Plan. Prepared by Herrera Environmental Consultants for River and Floodplain Management Section, King County Water and Land Resources Division. April.

USEPA. 2014. National functional guidelines for inorganic superfund data review. US Environmental Protection Agency, Office of Superfund Remediation and Technology Innovation (OSRTI), Washington, D.C. (EPA-540-R-013-001).

APPENDIX F

Geophysical Investigation Report, Pacific Right Bank Project



GEOPHYSICAL INVESTIGATION REPORT

PACIFIC RIGHT BANK PROJECT PACIFIC, WASHINGTON Contract E00495E17

FOR

HERRERA ENVIRONMENTAL CONSULTANTS, INC. SEATTLE, WASHINGTON

SEPTEMBER, 2018

PHILIP H. DUOOS
GEOPHYSICAL CONSULTANT

September 5, 2018 Our Ref.: 1261-18

Mr. George Iftner Herrera Environmental Consulting, Inc. 2200 Sixth Ave., Suite 1100 Seattle, WA 98121

Report: Geophysical Investigation

Pacific Right Bank Project Pacific, Washington Contract E00495E17

Dear Mr. Iftner:

This report describes the methodology and results of the geophysical investigation that I performed at the site as a subcontractor to Herrera in support of your Contract E00495E17 for the King County Department of Natural Resources and Parks, Water and Land Resources Division. The contract is to prepare an Environmental Impact Statement (EIS) and provide design services for the Pacific Right Bank project. The ultimate goal at the site is to design and construct a permanent flood facility, which requires a knowledge of the environmental and geotechnical conditions at the site.

The purpose of the geophysical investigation was to preliminarily characterize the lateral and vertical extent of buried waste to aid in targeting geotechnical and environmental explorations. The geophysical results also provide additional information between the numerous intrusive explorations (monitoring wells, push probes, etc.) and will help to refine waste volume estimates.

INTRODUCTION

The Pacific Right Bank project is located on the right (west) bank of the White River in the City of Pacific between the A Street and BNSF bridges (RM 6.3) and the King-Pierce county boundary line (RM 5.5). Downstream of the BNSF bridge abutment to about RM 5.9, the existing right bank consists of a levee and concrete revetment constructed in 1919. Filling of the former channel landward of the levee with municipal waste and dredge spoils occurred from the early-1920s through the late-1960s, when the area was developed into a park. (As stated in the Contract Request for Proposal).

Based on the history of the site, it was determined that the use of geophysical methods may be helpful in determining the lateral and vertical extent of buried waste at the site. The presence of municipal waste typically contains ferrous and non-ferrous metal debris. Dredge spoils often contain a greater amount of finer-grained materials (such as silt) which can increase the electrical conductivity of the subsurface. Several types of contamination such as total metals, salts and nitrates will also increase the electrical conductivity.

The Phase II Environmental Site Assessment, Pacific Right Bank Levee Setback Project, prepared for King County by Shannon & Wilson, Inc., May 2016 was reviewed prior to the geophysical field work. The report identified areas with debris that includes concrete and brick, burnt wood, metal debris, wire, porcelain, glass and plastic. Most of the debris was observed at depths less than about 20 feet. The presence of elevated levels of metals (such as lead, chromium and cadmium) and other contaminants were also detected.

The geophysical survey was performed during the period of January 30 - 31, 2018. Draft results were provided to Herrera on February 20 prior to your additional explorations at the site using push probes. The area of investigation was the open area of the park, and did not extend into the wooded area to the south. The site is relatively flat and contained large areas of ponded water during the survey. Several buildings, playground structures, metal benches, reinforced concrete slabs and other metal features are concentrated in the northern portion of the area. A baseball field near the west side of the park is bordered by a metal fence and bleachers to the west.

METHODOLOGY

Electromagnetic (EM) and magnetic geophysical methods were used to primarily determine the general lateral extent of buried debris. An Electromagnetic (EM) system was used to measure the electrical properties of the subsurface to an effective depth of about 18 feet. The electrical conductivity of the subsurface can be affected by various properties such as soil type, moisture content, and the presence of certain contaminants such as total metals, nitrates, etc. The data is also affected by the presence of buried ferrous metal and non-ferrous metal objects.

A magnetometer was also use to search for buried ferrous metal – the magnetometer is more sensitive to smaller amounts of ferrous material than the EM system. A brief description of the methods are provided in **Appendix A**.

My proposal recommended that electrical resistivity imaging (ERI) be used along one long line across the site to help determine the vertical extent of buried debris. However, upon obtaining more information regarding the site conditions, the ERI method as proposed would not have been able to provide the required resolution of the shallow subsurface conditions (less than 20 feet), and therefore was not used.

Reference Baselines

The geophysical surveys were referenced to numerous baselines that were marked using 300-foot tape measures, wood stakes and wire pin flags. The initial reference line runs along the levee and is labelled 100W and runs grid south to grid north (henceforth called south and north). True north is approximately 37 degrees west of grid north. Line 100W was marked with wood stakes placed at 100-foot intervals. Reference baselines were then established heading to the west, and spaced approximately 300 feet apart, and marked with wire pin flags at 50-foot intervals.

Additional north-south baselines were established along the west side of the site. A primary reference baseline was located along Line 650W between the Hesco barrier wall and the baseball field. Numerous shorter baselines oriented north-south were established in the northern portion of the site due to the angle of the property boundary.

Numerous reference baseline locations and visible reference features were located using a submeter GPS system for fairly accurate plotting of the baselines on the air photo used for Figure 5. Field measurements were also made to buildings and other features.

Electromagnetic Survey

The electromagnetic survey was performed using a DualEM Conductivity meter. Data were digitally recorded at one-second intervals while walking each line, with data recorded at about 3 to 4 foot intervals along each line. Transect lines were oriented east-west and typically spaced 100 feet apart as per the original scope of work. However, in the southern portion of the site a closer

line spacing of 50 feet was used as time allowed, and to provide more detail in an area with large changes in the EM conductivity data observed in the field. One short line was oriented north-south along Line 650W.

Magnetometer Survey

Magnetometer data were recorded with an EG&G Geometrics G-858 Cesium Magnetometer primarily along lines oriented east-west and spaced 100 feet apart as per the scope of work. While additional EM data were being recorded in the southern portion of the site, a second operator gathered additional magnetic data in the northern portion of the site at 50-foot line spacing. One long magnetic line was run north-south along Line 200W. The magnetic data was recorded at a faster data-collection rate resulting in data sampling at about 1 to 2 foot intervals along each line.

Both total field and vertical gradient magnetic data were obtained. The total magnetic field is affected by large amounts of ferrous material to a greater depth. The vertical gradient data is more sensitive to shallow buried ferrous material.

INTERPRETATION RESULTS

The interpretation considered all of the geophysical data. The data were also affected by the various structures and the wire mesh in the HESCO barrier wall around the western perimeter of the site. The locations of these various sources of interference were noted in our field notes and taken into consideration during the data analysis. The locations of major cultural interferences are shown on the data contour maps.

Electromagnetic Data

Both EM conductivity data (measured in milli-Siemens/meter – mS/m) and the inphase component of the primary field (measured in parts/thousand – ppt) were recorded. **Figure 1** is a contour map of the conductivity data which responds both to buried metal and changes in the electrical properties of the soil to a depth of about 18 feet. The southwestern portion of the site has elevated subsurface conductivities over fairly broad areas. These elevated values may indicate increased amounts, and/or shallower depths of conductive materials such as scattered amounts of small metal debris, possible conductive soil contamination such as elevated levels of total metals contamination, salts, nitrates, etc., an increase in finer-grained materials such as clay and silt. The smooth nature of the contours would indicate that the high conductivities are not caused by large buried metal objects.

The negative EM conductivity values near coordinate 350W, 700E is classified as a high metal anomaly. These values are typical of larger amounts of buried metal, and may indicate a denser mass of smaller objects or a cluster of larger metal objects (both ferrous and non-ferrous).

Several other locations along the eastern portion of the site also have conductivity values that are slightly negative. One area near 160W, 1300N is classified as a moderate metal anomalous zone, and indicates a buried metal objects with less mass than in the high metal anomalous zone. Negative values were also recorded along the western edge of the existing dike, and may indicate a linear metal feature such as a buried wire fence or a small-diameter pipe.

Figure 2 is a contour map of the EM inphase data which responds primarily to large concentrations of buried metal to a depth of about 18 feet. However, in areas with highly conductive soils the inphase values can also be affected. Elevated inphase values (above 4 ppt) are associated with the two southern areas with increased conductivity. The very low inphase values near 350W, 700N indicate large amounts of buried metal. Smaller and more scattered inphase anomalies are

observed over much of the eastern portion of the site and are interpreted to indicate scattered buried metal debris.

Magnetic Data

Both total field and vertical gradient magnetic data were obtained. **Figure 3** is a contour map of the total magnetic field data which is affected by ferrous material, and is more effective at detecting large amounts at a greater depth. Typical total field anomalies are indicated by high values to the south of the ferrous mass and low values to the north.

The interpreted buried metal near 350W, 700N shows such a response. A large magnitude total field anomaly is also located near 520W, 600N but is limited in areal extent. A smaller anomaly is located near 160W, 700N. The linear feature along the west side of the dike (near Line 100W) is very pronounced in the total field data.

The magnetic data were affected by the structures and playground equipment in a similar manner as the EM data. The magnetic data was also affected by the baseball backstop and nearby metal bleachers to a greater extent than the EM data.

Figure 4 is a contour map of the vertical gradient data which is more sensitive to smaller amounts of ferrous material and at shallower depths. The high values in the central portion of the site near 350W, 700N are very prominent and correlate well with the very low EM data values. The anomaly near 520W, 600N near the baseball field fence is also apparent. The vertical gradient data does not show a large anomaly near 160W, 700N where the total field data does. This may indicate that the object is deeper than some of the other features. The total field data is also more affected by permanent and/or remnant magnetization of some metal objects such as pipes.

The linear feature along the west side of the dike is also apparent. The scattered anomalies in the vertical gradient magnetic data are similar in distribution as the EM inphase anomalies. The vertical gradient data is much more sensitive to small amounts of buried ferrous material than the EM inphase data.

Combined Electromagnetic and Magnetic Interpretation

Figure 5 shows the interpretation results of the geophysical data. Areas interpreted to have buried debris with metal are categorized into three types of anomalous zones based on both the EM and magnetic data. The rankings (High, Moderate and Low Anomalous Zones) are based on the magnitude and character of the EM and magnetic data. The High anomalous zones are interpreted to have large amounts of buried metal. The Moderate zones are interpreted to have a lesser concentration of metal objects, or they may not extend as deep (so less amount of buried metal). The Low Anomalous zones are interpreted to contain low amounts of metal. Areas outside the anomalous zones may contain minor amounts of scattered buried metal that was not interpreted from the data.

The magnitude of an anomaly can also be affected by the depth of burial of the feature. However, the depth is difficult to determine at sites such as this with general waste and numerous sources of anomalies.

The two questionable high metal anomalies indicate a strong total field magnetic anomaly (see Figure 3), but there is no corresponding EM response (Figures 1 and 2). One is located near 170W, 700N, the other near 520W, 600N. These types of anomalies may occur over a vertical pipe (such as an old steel well casing, or a vertical metal post). It may also indicate a deeper depth of burial. The EM tool has an effective depth of 18 feet, and the magnetometer can detect large amounts of metal at greater depths.

The linear feature that is near the western edge of the levee, and runs along about Line 100W is observed on both the EM and magnetic data. It may be an old metal or wire fence now buried, a small diameter buried pipe, or perhaps some metal structure used in the construction of the levee.

Selected EM conductivity contour lines are also shown on Figure 5 and indicate areas with higher electrical conductivity. The elevated values may be caused by scattered metal debris as the magnetic data is also elevated in the general areas where the EM values are greater. The EM contours may also show fill material with higher electrical conductivity due to finer-grained materials such as silt, or various types of possible contamination including metals, salts and nitrates. The increase in conductivity values may be related to the concentration of the higher conductivity material and/or a greater thickness of the material.

These results were interpreted based on the geophysical data. Upon completion of the numerous push probes advanced at the site after the geophysical survey, the exploration logs were reviewed. Additionally, the earlier results from previous test pits, wells and geoprobe investigations were also reviewed in greater detail.

The various logs of the intrusive investigations were summarized into seven categories based on subsurface conditions that may affect the EM conductivity and/or magnetic data. The locations of the explorations (test pits, probes, etc.) are shown on Figure 5. I have added the following colored symbols indicating the general categories based on my notes of the logs as follows:

Large Red Hexagon: Buried metal debris observed

Brown Circle: Elevated Shallow Metals Contamination (S&W, 2016 results only)

Dark Green Circle:
Light Green Circle:
Yellow Circle:
Debris and silty fill material
Debris with little silt fill material
Silt fill material with no debris
Fill material with little silt

Dark Blue Circle: Little fill material

Explorations with observed debris also have the depth range of the debris noted. Some explorations detected traces of small pieces of glass at depths that I did not include in my notes as indicating primary layers of debris for the purposes of correlating to the geophysical data.

The results of the test pits are probably more reliable for categorizing the fill material for debris and the type of debris. Several probes that did not encounter metal debris were located in areas interpreted to contain buried metal based on very strong anomalies in the geophysical data. However, a small diameter probe could easily miss a discrete buried metal object. At most waste sites there is always some percentage of metal debris mixed in with the non-metal waste.

The explorations generally agree with the geophysical interpretation. Most of the explorations with observed debris are within anomalous zones. Several explorations outside the anomalous zones detected debris including buried metal. However, other nearby explorations indicated no debris, so the detected debris is not interpreted to be wide spread.

Areas with higher EM conductivity values (shown by the data contours) were interpreted to be related to areas with thicker and/or deeper zones of debris, or increased amounts of metal. The summarized exploration results do not seem to relate to either of these initial possible interpretations. The increased conductivities also do not seem to be related to higher silt content. Water content can increase the conductivity, but most of the site was water-saturated at the time of the survey. Additionally, the higher conductivity values (20 mS/m) near 400W, 600N are in a low area with standing water, while the higher values (20 mS/m) near 300W, 400N are in an area with higher ground elevations. This would seem to indicate that the depth to water (and percentage of the material that is saturated) is not a significant factor.

Review of the existing lab results from earlier explorations seems to indicate a possible correlation of the total metals contamination of the shallow soils to the geophysical results. At the time of this report, the results from the Shannon and Wilson Phase II Environmental Site Assessment (May, 2016) and results for push probes PP-1 through PP-10 from June, 2017 (Herrera) were available. The Shannon and Wilson report indicated numerous locations with total metals contamination at shallow soil depths (generally between 0 and 5 feet deep) that exceed the Washington state Model Toxics Control Act (MTCA) levels. These locations are shown on Figure 5 with a brown circle. The majority of these locations are within an anomalous zone indicating metal debris or an area with elevated EM conductivity. The Herrera push probes PP-1 through PP-10 did not indicate high levels of total metals contamination of the soil. The lab results from the most recent round of push probe explorations was not available.

The higher conductivity zones may also be related to other factors not observed in the boring logs such as the mineralogy of the sediments. Iron-rich sediments from mafic materials may cause elevated conductivity values.

CONCLUSIONS

The use of these geophysical techniques provided a rapid and non-intrusive means of investigating the area of interest for large zones of buried debris containing metal and possible soil contamination to provide a broad overview of the site conditions. However, because of the numerous variables involved in geophysical investigations, there is a possibility that some features may not have been detected.

Please contact me if you have any questions or comments regarding this information, or if you require further assistance. I appreciated the opportunity to work with you on this project and look forward to providing you with geophysical services in the future.

Sincerely,

Philip H. Duoos

Geophysical Consultant

Attachments:

Description of Methods

1. Aum

Figure 1: EM Conductivity Data Contour Map Figure 2: EM Inphase Data Contour Map Figure 3: Total Magnetic Field Contour Map

Figure 4: Vertical Gradient Magnetic Field Contour Map

Figure 5: Geophysical Results Map



ATTACHMENT A

DESCRIPTION OF METHODS

ELECTROMAGNETICS (DualEM)

The DualEM instrument measures subsurface conductance using the principles of electromagnetic induction to depths of about 18 feet, and can detect large amounts of metal at greater depths. The DualEM is portable, rapid and non-destructive. It has a fixed boom containing the transmitter and receiver coils so that handling and data gathering is easily achieved by one operator.

Factors which may increase subsurface conductivities (measured in milli-Siemens/meter – mS/m) include higher moisture content, greater amounts of finer materials, increased clay and/or silt content, soil contamination and/or ground water contamination. The presence of buried metal can also affect the conductivity data. The detectability of metal objects (buried pipes, drums, etc.) can be enhanced by measuring the change in the magnitude of the primary field (in-phase component, measured in parts/thousand – ppt) of the induced magnetic field.

Several factors can limit the effectiveness of the EM method including the proximity of cultural interferences (such as buildings, fences and reinforced concrete) the presence of highly conductive materials (such as clays and water), and the size, depth and conductivity contrast of the target.

EG&G 858 CESIUM MAGNETOMETER/GRADIOMETER

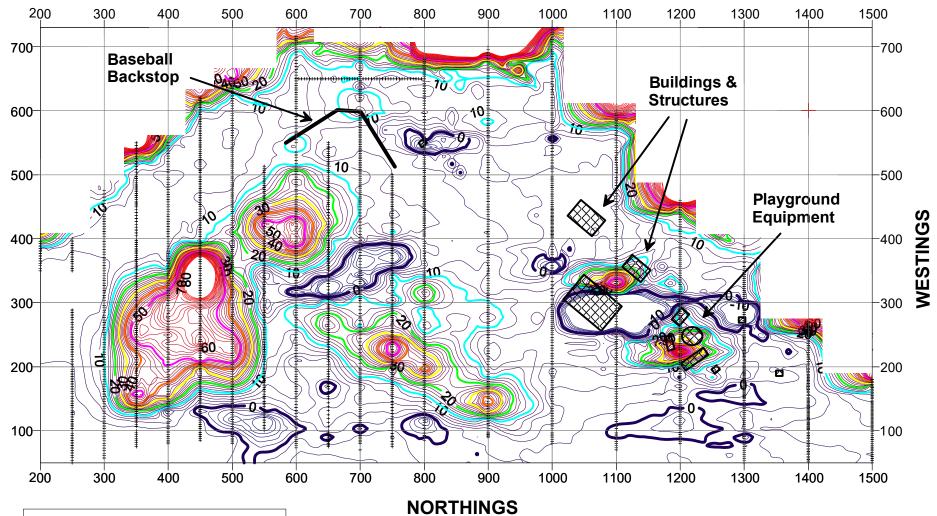
The EG&G 858 magnetometer/gradiometer is a rapid, effective and non-destructive instrument used to locate buried ferrous material (drums, pipes, mineral deposits, archaeological objects, etc.). The gradiometer consists of two sensors (one at a height of six feet, the other at a height of three feet) and a digital recording unit carried in a harness by one operator. Data are recorded and later downloaded to a computer.

Two types of measurements are recorded during a gradiometer survey: the total field and vertical gradient. The total field measurement (measured in nanoTeslas – nT) recorded using the top sensor as it is farther from the ground and less susceptible to small amounts of scattered ferrous material). The total field measurement can detect large ferrous metal objects to a great depth. The total field data is also affected by regional geology and naturally occurring changes in the earth's magnetic field during the day.

The vertical gradient data is the difference in the earth's magnetic field measured between the two magnetometer sensors. The vertical gradient data (nT/meter) are more affected by near-surface sources and provides better resolution of shallow buried objects. The gradient data is less affected by cultural interferences such as fences and vehicles, and is also not affected by the naturally occurring changes in the earth's total field throughout the day.

The magnetic data may also detect disturbances in the native soils due to trenching and grading activities. These activities disturb the original remnant magnetic orientation of the soil particles when they were deposited. Natural changes in the soils due to different materials (zones of gravels or cobbles) or natural erosion may also be the source of these minor disturbances in the gradient data.

Several factors can limit the effectiveness of the magnetometry method including the proximity of cultural interferences (such as buildings, fences and reinforced concrete), and the size, depth and magnetic susceptibility of the target.



Electromagnetic Conductivity Data

-10 to 80 mS/m at 2 mS/m intervals

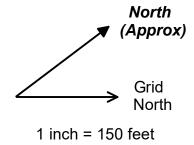
0: dark blue 10: light blue

20: green 30: yellow

40: orange

50: magenta

Philip H. Duoos, Geophysical Consultant for Herrera, Inc. Duoos Project No. 1261-18

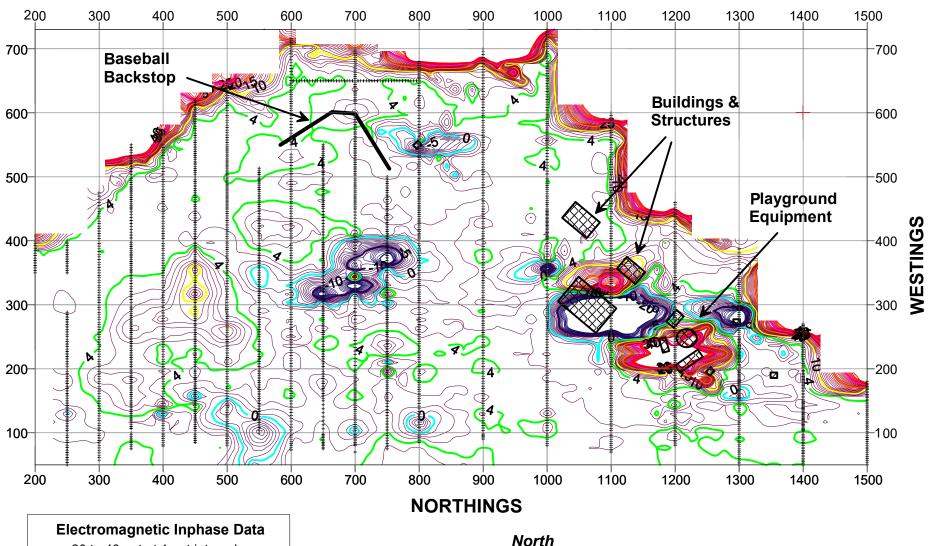


Electromagnetic Conductivity Data

Pacific Right Bank Project Pacific, Washington

Contract E00495E17

9/5/2018 **FIG. 1**



-20 to 40 ppt at 1 ppt intervals

- dark blue
 - light blue 0:
- green
- yellow
- orange
- magenta

(Approx) 1 inch = 150 feet

Grid

North

Philip H. Duoos, Geophysical Consultant for Herrera, Inc. Duoos Project No. 1261-18

Electromagnetic Inphase Data

Pacific Right Bank Project Pacific, Washington

Contract E00495E17

9/5/2018

FIG. 2

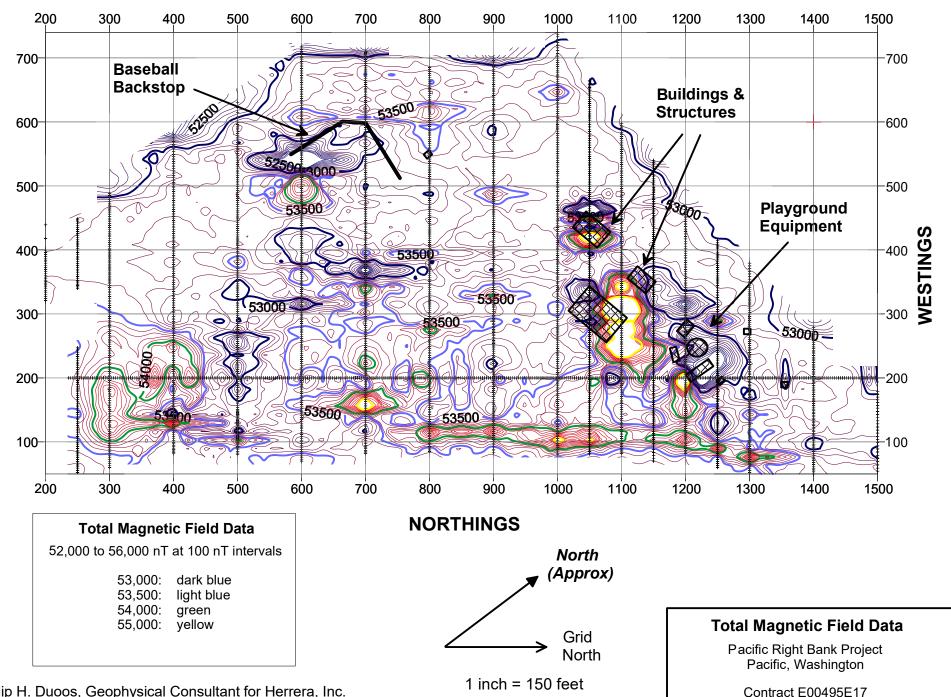
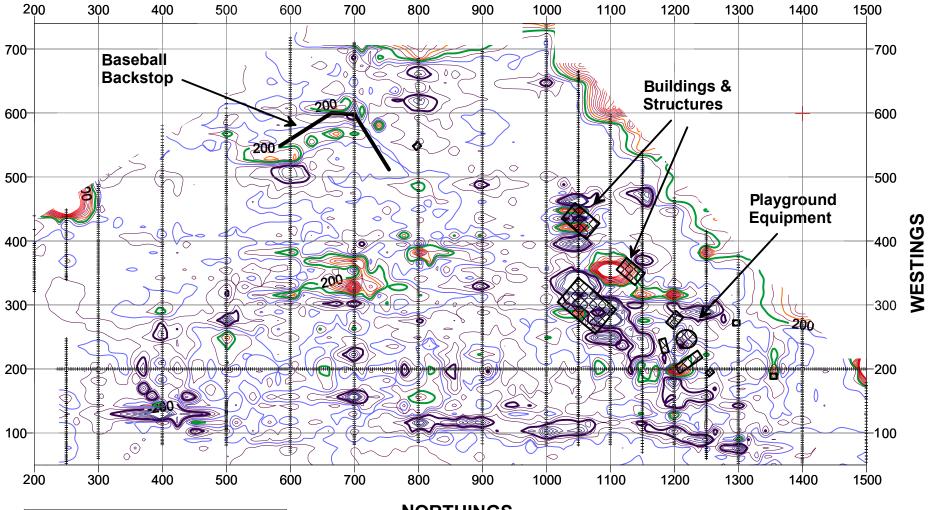


FIG. 3

9/5/2018

Philip H. Duoos, Geophysical Consultant for Herrera, Inc. Duoos Project No. 1261-18



Vertical Gradient Magnetic Data

-600 to 1160 nT/m at 80 nT/m intervals

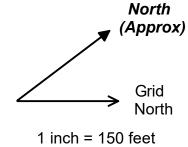
-200: dark blue 40: light blue (thin)

200: green

360: orange (thin)

Philip H. Duoos, Geophysical Consultant for Herrera, Inc. Duoos Project No. 1261-18

NORTHINGS



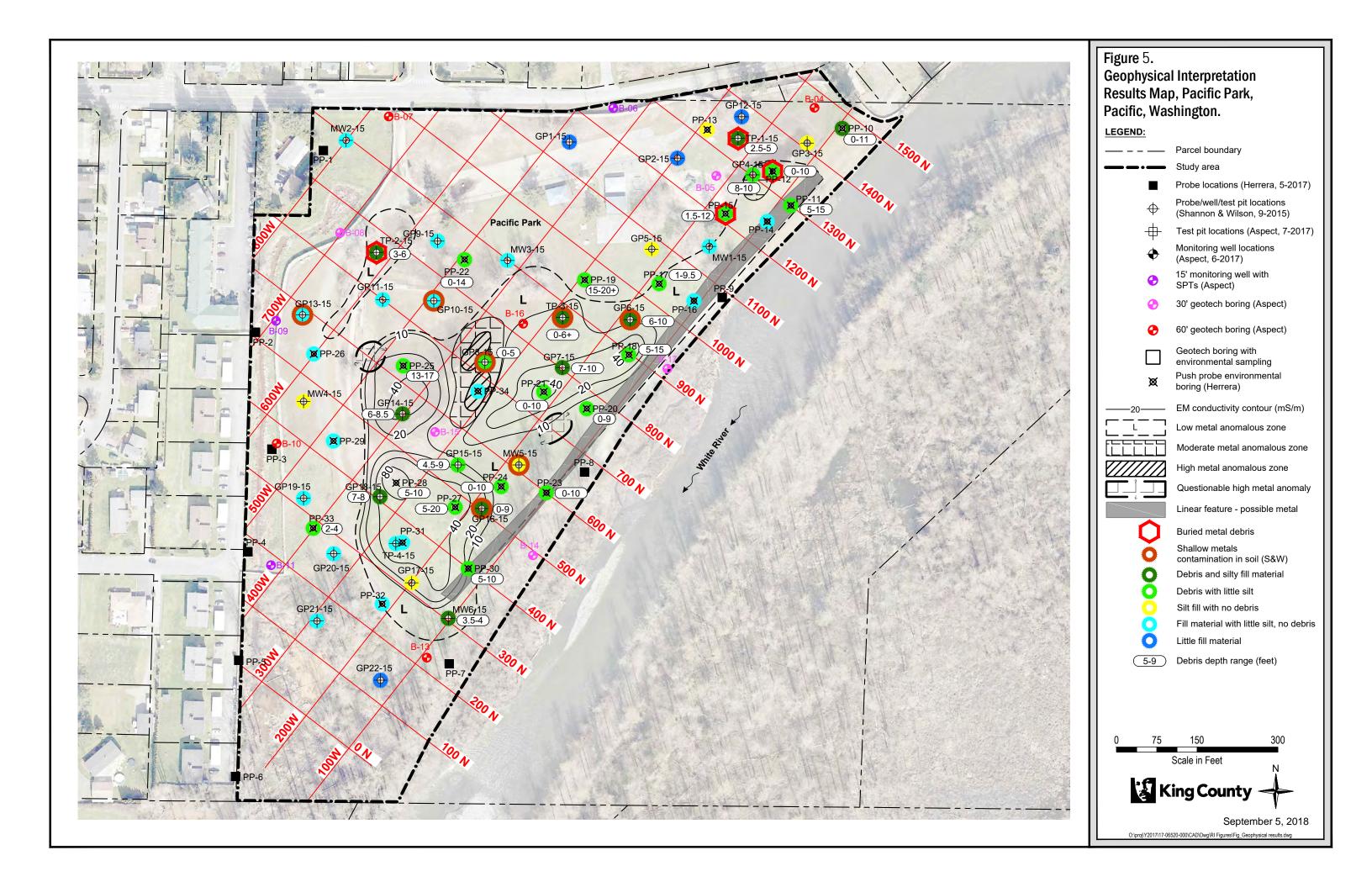
Vertical Gradient Magnetic Data

Pacific Right Bank Project Pacific, Washington

Contract E00495E17

9/5/2018

FIG. 4



APPENDIX G

Laboratory Analytical Reports





14648 NE 95th Street, Redmond, WA 98052 • (425) 883-3881

March 16, 2018

Bruce Carpenter Herrera Environmental Consultants, Inc. 2200 6th Avenue, Suite 1100 Seattle, WA 98121

Re: Analytical Data for Project 17-06520-000

Laboratory Reference No. 1802-233

Dear Bruce:

Enclosed are the analytical results and associated quality control data for samples submitted on February 22, 2018.

The standard policy of OnSite Environmental, Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

David Baumeister Project Manager

Enclosures



Project: 17-06520-000

Case Narrative

Samples were collected on February 20 and 21, 2018 and received by the laboratory on February 22, 2018. They were maintained at the laboratory at a temperature of 2°C to 6°C.

Please note that any and all soil sample results are reported on a dry-weight basis, unless otherwise noted below.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

Total Metals EPA 6010D/7471B Analysis

The Matrix Spike/ Matrix Spike Duplicate recoveries for Chromium are outside control limits due to matrix inhomogeneity. The samples were re-extracted and re-analyzed with similar results. The Spike Blank recovery was 99%.

The Matrix Spike/Matrix Spike Duplicate RPD for Chromium is outside control limits due to matrix inhomogeneity. The samples were re-extracted and re-analyzed with similar results.

Any other QA/QC issues associated with this extraction and analysis will be indicated with a footnote reference and discussed in detail on the Data Qualifier page.

Project: 17-06520-000

NWTPH-HCID

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-15_5					
Laboratory ID:	02-233-13					
Gasoline Range Organics	ND	35	NWTPH-HCID	2-23-18	2-23-18	
Diesel Range Organics	ND	87	NWTPH-HCID	2-23-18	2-23-18	
Lube Oil Range Organics	ND	170	NWTPH-HCID	2-23-18	2-23-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	96	50-150				
Client ID:	B-15_7.5					
Laboratory ID:	02-233-14					
Gasoline Range Organics	ND	30	NWTPH-HCID	2-23-18	2-23-18	
Diesel Range Organics	ND	76	NWTPH-HCID	2-23-18	2-23-18	
Lube Oil Range Organics	ND	150	NWTPH-HCID	2-23-18	2-23-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	90	50-150				
Client ID:	D 45 45					
	B-15_15 02-233-15					
Laboratory ID:		07	NATELLIOID	0.00.10	0.00.40	
Gasoline Range Organics	ND	27	NWTPH-HCID	2-23-18	2-23-18	
Diesel Range Organics	ND	67	NWTPH-HCID	2-23-18	2-23-18	
Lube Oil Range Organics	ND	140	NWTPH-HCID	2-23-18	2-23-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	86	50-150				

Project: 17-06520-000

NWTPH-HCID QUALITY CONTROL

Matrix: Soil

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0223S3					
Gasoline Range Organics	ND	20	NWTPH-HCID	2-23-18	2-23-18	
Diesel Range Organics	ND	50	NWTPH-HCID	2-23-18	2-23-18	
Lube Oil Range Organics	ND	100	NWTPH-HCID	2-23-18	2-23-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	95	50-150				

Project: 17-06520-000

NWTPH-HCID

Matrix: Soil

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	B-04_12.5	I QL	Metriou	rrepared	Allalyzeu	i iags
Laboratory ID:	02-233-03					
Gasoline Range Organics	ND	25	NWTPH-HCID	2-26-18	2-26-18	
Diesel Range Organics	ND	64	NWTPH-HCID	2-26-18	2-26-18	
Lube Oil Range Organics	ND	130	NWTPH-HCID	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	103	50-150				
011 4 15						
Client ID:	B-05_7.5					
Laboratory ID:	02-233-06	0.4	NIMEDIALIO	0.00.40	0.00.40	
Gasoline Range Organics	ND Data ata d	24	NWTPH-HCID	2-26-18	2-26-18	
Diesel Range Organics	Detected	60	NWTPH-HCID	2-26-18	2-26-18	
Lube Oil Range Organics	Detected Percent Recovery	120	NWTPH-HCID	2-26-18	2-26-18	
Surrogate: o-Terphenyl	Percent Recovery 108	Control Limits 50-150				
0-Terprienyi	108	<i>50-150</i>				
Client ID:	B-05_12.5					
Laboratory ID:	02-233-07					
Gasoline Range Organics	ND	25	NWTPH-HCID	2-26-18	2-26-18	
Diesel Range Organics	ND	63	NWTPH-HCID	2-26-18	2-26-18	
Lube Oil Range Organics	ND	130	NWTPH-HCID	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	100	50-150				
Client ID:	B-14_5					
Laboratory ID:	02-233-09					
Gasoline Range Organics	ND	27	NWTPH-HCID	2-26-18	2-26-18	
Diesel Range Organics	ND	68	NWTPH-HCID	2-26-18	2-26-18	
Lube Oil Range Organics	ND	140	NWTPH-HCID	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	97	50-150				
Client ID:	B-14_10					
Laboratory ID:	02-233-10					
Gasoline Range Organics	ND	25	NWTPH-HCID	2-26-18	2-26-18	
Diesel Range Organics	ND	62	NWTPH-HCID	2-26-18	2-26-18	
Lube Oil Range Organics	ND	120	NWTPH-HCID	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	100	50-150				

Project: 17-06520-000

NWTPH-HCID

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-17_10					
Laboratory ID:	02-233-18					
Gasoline Range Organics	ND	25	NWTPH-HCID	2-26-18	2-26-18	
Diesel Range Organics	ND	61	NWTPH-HCID	2-26-18	2-26-18	
Lube Oil Range Organics	ND	120	NWTPH-HCID	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	100	50-150				
Client ID:	B-17_15					
Laboratory ID:	02-233-19					
Gasoline Range Organics	ND	22	NWTPH-HCID	2-26-18	2-26-18	
Diesel Range Organics	ND	54	NWTPH-HCID	2-26-18	2-26-18	
Lube Oil Range Organics	ND	110	NWTPH-HCID	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	99	50-150				

Project: 17-06520-000

NWTPH-HCID QUALITY CONTROL

Matrix: Soil

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0226S2					
Gasoline Range Organics	ND	20	NWTPH-HCID	2-26-18	2-26-18	
Diesel Range Organics	ND	50	NWTPH-HCID	2-26-18	2-26-18	
Lube Oil Range Organics	ND	100	NWTPH-HCID	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	100	50-150				

Project: 17-06520-000

NWTPH-Dx

Matrix: Soil

Analyte	Result	PQL	Method	Date Prepared	Date Analvzed	Flags
Client ID:	B-05_7.5	1 0,2	Mictiloa	Troparca	Anaryzea	i iugo
Laboratory ID:	02-233-06					
Diesel Range Organics	440	30	NWTPH-Dx	3-1-18	3-1-18	_
Lube Oil Range Organics	ND	220	NWTPH-Dx	3-1-18	3-1-18	U1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	101	50-150				

Project: 17-06520-000

NWTPH-Dx QUALITY CONTROL

Matrix: Soil

Analyte	Result	PQL	Method	Date Prepared	Date Analvzed	Flage
	Nesuit	FUL	METHOR	riepaieu	Allalyzeu	Flags
METHOD BLANK						
Laboratory ID:	MB0301S2					
Diesel Range Organics	ND	25	NWTPH-Dx	3-1-18	3-1-18	
Lube Oil Range Organics	ND	50	NWTPH-Dx	3-1-18	3-1-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	93	50-150				

Amaluta	Por	sult	Cnika	Lovel	Source Result	Perc		Recovery Limits	RPD	RPD Limit	Elogo
Analyte	Res	Suit	Бріке	Level	Resuit	Reco	very	Limits	RPU	Limit	Flags
DUPLICATE											
Laboratory ID:	03-00	03-06									
	ORIG	DUP									
Mineral Oil	ND	ND	NA	NA		N.	A	NA	NA	NA	
Surrogate:											
o-Terphenyl						90	70	50-150			

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-15_5					
Laboratory ID:	02-233-13					
Benzo[a]anthracene	ND	0.012	EPA 8270D/SIM	3-1-18	3-1-18	
Chrysene	ND	0.012	EPA 8270D/SIM	3-1-18	3-1-18	
Benzo[b]fluoranthene	0.014	0.012	EPA 8270D/SIM	3-1-18	3-1-18	
Benzo(j,k)fluoranthene	ND	0.012	EPA 8270D/SIM	3-1-18	3-1-18	
Benzo[a]pyrene	ND	0.012	EPA 8270D/SIM	3-1-18	3-1-18	
Indeno(1,2,3-c,d)pyrene	ND	0.012	EPA 8270D/SIM	3-1-18	3-1-18	
Dibenz[a,h]anthracene	ND	0.012	EPA 8270D/SIM	3-1-18	3-1-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	74	32 - 115				
Pyrene-d10	89	35 - 129				
Terphenyl-d14	93	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

			Date	Date	
Result	PQL	Method	Prepared	Analyzed	Flags
B-15_7.5					
02-233-14					
ND	0.010	EPA 8270D/SIM	3-1-18	3-1-18	
ND	0.010	EPA 8270D/SIM	3-1-18	3-1-18	
ND	0.010	EPA 8270D/SIM	3-1-18	3-1-18	
ND	0.010	EPA 8270D/SIM	3-1-18	3-1-18	
ND	0.010	EPA 8270D/SIM	3-1-18	3-1-18	
ND	0.010	EPA 8270D/SIM	3-1-18	3-1-18	
ND	0.010	EPA 8270D/SIM	3-1-18	3-1-18	
Percent Recovery	Control Limits				
77	32 - 115				
94	35 - 129				
96	33 - 114				
	B-15_7.5 02-233-14 ND ND ND ND ND ND ND ND ND Percent Recovery 77 94	B-15_7.5 02-233-14 ND 0.010 ND 0.010 ND 0.010 ND 0.010 ND 0.010 ND 0.010 Percent Recovery Control Limits 77 32 - 115 94 35 - 129	B-15_7.5 02-233-14 ND 0.010 EPA 8270D/SIM Percent Recovery Control Limits 77 32 - 115 94 35 - 129	Result PQL Method Prepared B-15_7.5 02-233-14 ND 0.010 EPA 8270D/SIM 3-1-18 Percent Recovery Control Limits 77 32 - 115 32 - 115 94 35 - 129 35 - 129	Result PQL Method Prepared Analyzed B-15_7.5 02-233-14 Secondary 100 Best School 100 Best School 100 3-1-18 3-1-18 ND 0.010 Best School 100 3-1-18 3-1-18 3-1-18 ND 0.010 Best Best School 100 3-1-18 3-1-18 3-1-18 ND 0.010 Best Best School 100 3-1-18 3-1-18 3-1-18 ND 0.010 Best Best School 100 3-1-18 3-1-18 3-1-18 ND 0.010 Best Best School 100 3-1-18 3-1-18 3-1-18 Percent Recovery Control Limits 77 32 - 115 32 - 115 3-1-18 3-1-18

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-15_15					
Laboratory ID:	02-233-15					
Benzo[a]anthracene	ND	0.0090	EPA 8270D/SIM	3-1-18	3-1-18	
Chrysene	ND	0.0090	EPA 8270D/SIM	3-1-18	3-1-18	
Benzo[b]fluoranthene	ND	0.0090	EPA 8270D/SIM	3-1-18	3-1-18	
Benzo(j,k)fluoranthene	ND	0.0090	EPA 8270D/SIM	3-1-18	3-1-18	
Benzo[a]pyrene	ND	0.0090	EPA 8270D/SIM	3-1-18	3-1-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0090	EPA 8270D/SIM	3-1-18	3-1-18	
Dibenz[a,h]anthracene	ND	0.0090	EPA 8270D/SIM	3-1-18	3-1-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	80	32 - 115				
Pyrene-d10	85	35 - 129				
Terphenyl-d14	91	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-04_12.5					
Laboratory ID:	02-233-03					
Benzo[a]anthracene	ND	0.0085	EPA 8270D/SIM	3-1-18	3-4-18	
Chrysene	ND	0.0085	EPA 8270D/SIM	3-1-18	3-4-18	
Benzo[b]fluoranthene	ND	0.0085	EPA 8270D/SIM	3-1-18	3-4-18	
Benzo(j,k)fluoranthene	ND	0.0085	EPA 8270D/SIM	3-1-18	3-4-18	
Benzo[a]pyrene	ND	0.0085	EPA 8270D/SIM	3-1-18	3-4-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0085	EPA 8270D/SIM	3-1-18	3-4-18	
Dibenz[a,h]anthracene	ND	0.0085	EPA 8270D/SIM	3-1-18	3-4-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	85	32 - 115				
Pyrene-d10	97	35 - 129				
Terphenyl-d14	105	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-05_7.5					
Laboratory ID:	02-233-06					
Benzo[a]anthracene	ND	0.0080	EPA 8270D/SIM	3-1-18	3-5-18	
Chrysene	ND	0.0080	EPA 8270D/SIM	3-1-18	3-5-18	
Benzo[b]fluoranthene	ND	0.0080	EPA 8270D/SIM	3-1-18	3-5-18	
Benzo(j,k)fluoranthene	ND	0.0080	EPA 8270D/SIM	3-1-18	3-5-18	
Benzo[a]pyrene	ND	0.0080	EPA 8270D/SIM	3-1-18	3-5-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0080	EPA 8270D/SIM	3-1-18	3-5-18	
Dibenz[a,h]anthracene	ND	0.0080	EPA 8270D/SIM	3-1-18	3-5-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	84	32 - 115				
Pyrene-d10	109	35 - 129				
Terphenyl-d14	110	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-05_12.5					
Laboratory ID:	02-233-07					
Benzo[a]anthracene	ND	0.0084	EPA 8270D/SIM	3-1-18	3-4-18	
Chrysene	ND	0.0084	EPA 8270D/SIM	3-1-18	3-4-18	
Benzo[b]fluoranthene	ND	0.0084	EPA 8270D/SIM	3-1-18	3-4-18	
Benzo(j,k)fluoranthene	ND	0.0084	EPA 8270D/SIM	3-1-18	3-4-18	
Benzo[a]pyrene	ND	0.0084	EPA 8270D/SIM	3-1-18	3-4-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0084	EPA 8270D/SIM	3-1-18	3-4-18	
Dibenz[a,h]anthracene	ND	0.0084	EPA 8270D/SIM	3-1-18	3-4-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	87	32 - 115				
Pyrene-d10	93	35 - 129				
Terphenyl-d14	100	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-14_5					
Laboratory ID:	02-233-09					
Benzo[a]anthracene	ND	0.0090	EPA 8270D/SIM	3-1-18	3-5-18	
Chrysene	ND	0.0090	EPA 8270D/SIM	3-1-18	3-5-18	
Benzo[b]fluoranthene	ND	0.0090	EPA 8270D/SIM	3-1-18	3-5-18	
Benzo(j,k)fluoranthene	ND	0.0090	EPA 8270D/SIM	3-1-18	3-5-18	
Benzo[a]pyrene	ND	0.0090	EPA 8270D/SIM	3-1-18	3-5-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0090	EPA 8270D/SIM	3-1-18	3-5-18	
Dibenz[a,h]anthracene	ND	0.0090	EPA 8270D/SIM	3-1-18	3-5-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	84	32 - 115				
Pyrene-d10	96	35 - 129				
Terphenyl-d14	102	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-14_10					
Laboratory ID:	02-233-10					
Benzo[a]anthracene	ND	0.0082	EPA 8270D/SIM	3-1-18	3-4-18	
Chrysene	ND	0.0082	EPA 8270D/SIM	3-1-18	3-4-18	
Benzo[b]fluoranthene	ND	0.0082	EPA 8270D/SIM	3-1-18	3-4-18	
Benzo(j,k)fluoranthene	ND	0.0082	EPA 8270D/SIM	3-1-18	3-4-18	
Benzo[a]pyrene	ND	0.0082	EPA 8270D/SIM	3-1-18	3-4-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0082	EPA 8270D/SIM	3-1-18	3-4-18	
Dibenz[a,h]anthracene	ND	0.0082	EPA 8270D/SIM	3-1-18	3-4-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	91	32 - 115				
Pyrene-d10	91	35 - 129				
Terphenyl-d14	98	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-17_10					
Laboratory ID:	02-233-18					
Benzo[a]anthracene	ND	0.0082	EPA 8270D/SIM	3-1-18	3-4-18	
Chrysene	ND	0.0082	EPA 8270D/SIM	3-1-18	3-4-18	
Benzo[b]fluoranthene	ND	0.0082	EPA 8270D/SIM	3-1-18	3-4-18	
Benzo(j,k)fluoranthene	ND	0.0082	EPA 8270D/SIM	3-1-18	3-4-18	
Benzo[a]pyrene	ND	0.0082	EPA 8270D/SIM	3-1-18	3-4-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0082	EPA 8270D/SIM	3-1-18	3-4-18	
Dibenz[a,h]anthracene	ND	0.0082	EPA 8270D/SIM	3-1-18	3-4-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	87	32 - 115				
Pyrene-d10	93	35 - 129				
Terphenyl-d14	103	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-17_15					
Laboratory ID:	02-233-19					
Benzo[a]anthracene	ND	0.0072	EPA 8270D/SIM	3-1-18	3-4-18	
Chrysene	ND	0.0072	EPA 8270D/SIM	3-1-18	3-4-18	
Benzo[b]fluoranthene	ND	0.0072	EPA 8270D/SIM	3-1-18	3-4-18	
Benzo(j,k)fluoranthene	ND	0.0072	EPA 8270D/SIM	3-1-18	3-4-18	
Benzo[a]pyrene	ND	0.0072	EPA 8270D/SIM	3-1-18	3-4-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0072	EPA 8270D/SIM	3-1-18	3-4-18	
Dibenz[a,h]anthracene	ND	0.0072	EPA 8270D/SIM	3-1-18	3-4-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	90	32 - 115				
Pyrene-d10	100	35 - 129				
Terphenyl-d14	108	33 - 114				

Date of Report: March 16, 2018 Samples Submitted: February 22, 2018

Laboratory Reference: 1802-233

Project: 17-06520-000

CPAHS EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0301S2					
Benzo[a]anthracene	ND	0.0067	EPA 8270D/SIM	3-1-18	3-1-18	
Chrysene	ND	0.0067	EPA 8270D/SIM	3-1-18	3-1-18	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270D/SIM	3-1-18	3-1-18	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270D/SIM	3-1-18	3-1-18	
Benzo[a]pyrene	ND	0.0067	EPA 8270D/SIM	3-1-18	3-1-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0067	EPA 8270D/SIM	3-1-18	3-1-18	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270D/SIM	3-1-18	3-1-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	92	32 - 115				
Pyrene-d10	97	35 - 129				
Terphenyl-d14	104	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM MS/MSD QUALITY CONTROL

					Source	Per	cent	Recovery		RPD	
Analyte	Re	sult	Spike	Level	Result	Rec	overy	Limits	RPD	Limit	Flags
MATRIX SPIKES											
Laboratory ID:	02-23	34-04									
	MS	MSD	MS	MSD		MS	MSD				
Benzo[a]anthracene	0.172	0.183	0.167	0.167	ND	103	110	27 - 143	6	23	
Chrysene	0.161	0.170	0.167	0.167	ND	96	102	22 - 130	5	24	
Benzo[b]fluoranthene	0.160	0.169	0.167	0.167	ND	96	101	15 - 141	5	26	
Benzo(j,k)fluoranthene	0.161	0.168	0.167	0.167	ND	96	101	42 - 112	4	24	
Benzo[a]pyrene	0.161	0.168	0.167	0.167	ND	96	101	33 - 126	4	26	
Indeno(1,2,3-c,d)pyrene	0.154	0.157	0.167	0.167	ND	92	94	30 - 125	2	25	
Dibenz[a,h]anthracene	0.156	0.159	0.167	0.167	ND	93	95	31 - 124	2	22	
Surrogate:											
2-Fluorobiphenyl						81	88	32 - 115			
Pyrene-d10						93	95	35 - 129			
Terphenyl-d14						97	97	33 - 114			

Project: 17-06520-000

PCBs EPA 8082A

Matrix: Soil

Units: mg/Kg (ppm)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-05_7.5					
Laboratory ID:	02-233-06					
Aroclor 1016	ND	0.060	EPA 8082A	3-1-18	3-1-18	
Aroclor 1221	ND	0.060	EPA 8082A	3-1-18	3-1-18	
Aroclor 1232	ND	0.060	EPA 8082A	3-1-18	3-1-18	
Aroclor 1242	ND	0.060	EPA 8082A	3-1-18	3-1-18	
Aroclor 1248	ND	0.060	EPA 8082A	3-1-18	3-1-18	
Aroclor 1254	ND	0.060	EPA 8082A	3-1-18	3-1-18	
Aroclor 1260	ND	0.060	EPA 8082A	3-1-18	3-1-18	

Surrogate: Percent Recovery Control Limits DCB 76 40-134

Project: 17-06520-000

PCBs EPA 8082A QUALITY CONTROL

Matrix: Soil

Units: mg/Kg (ppm)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0301S1					
Aroclor 1016	ND	0.050	EPA 8082A	3-1-18	3-1-18	
Aroclor 1221	ND	0.050	EPA 8082A	3-1-18	3-1-18	
Aroclor 1232	ND	0.050	EPA 8082A	3-1-18	3-1-18	
Aroclor 1242	ND	0.050	EPA 8082A	3-1-18	3-1-18	
Aroclor 1248	ND	0.050	EPA 8082A	3-1-18	3-1-18	
Aroclor 1254	ND	0.050	EPA 8082A	3-1-18	3-1-18	
Aroclor 1260	ND	0.050	EPA 8082A	3-1-18	3-1-18	

Surrogate: Percent Recovery Control Limits DCB 84 40-134

					Source	Pe	rcent	Recovery		RPD	
Analyte	Re	sult	Spike	Level	Result	Red	covery	Limits	RPD	Limit	Flags
MATRIX SPIKES											
Laboratory ID:	02-2	30-02									
	MS	MSD	MS	MSD		MS	MSD				
Aroclor 1260	0.408	0.443	0.500	0.500	ND	82	89	34-126	8	16	
Surrogate:	•		•			•	•		•		•
DCB						74	81	40-134			

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B

Matrix: Soil

Ssg,. 15 (pp)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-15_5					
Laboratory ID:	02-233-13					
Arsenic	ND	17	EPA 6010D	2-26-18	2-27-18	
Cadmium	1.3	0.87	EPA 6010D	2-26-18	2-27-18	
Chromium	35	0.87	EPA 6010D	2-26-18	2-27-18	
Lead	75	8.7	EPA 6010D	2-26-18	2-27-18	
Mercury	ND	0.43	EPA 7471B	2-23-18	2-23-18	
- 11						
Client ID:	B-15_7.5					
Laboratory ID:	02-233-14					
Arsenic	ND	15	EPA 6010D	2-26-18	2-27-18	
Cadmium	ND	0.76	EPA 6010D	2-26-18	2-27-18	
Chromium	31	0.76	EPA 6010D	2-26-18	2-27-18	
Lead	49	7.6	EPA 6010D	2-26-18	2-27-18	
Mercury	ND	0.38	EPA 7471B	2-23-18	2-23-18	
Client ID:	B-15_15					
Laboratory ID:	02-233-15					
Arsenic	ND	13	EPA 6010D	2-26-18	2-27-18	
Cadmium	ND	0.67	EPA 6010D	2-26-18	2-27-18	
Chromium	17	0.67	EPA 6010D	2-26-18	2-27-18	
Lead	ND	6.7	EPA 6010D	2-26-18	2-27-18	
Mercury	ND	0.34	EPA 7471B	2-23-18	2-23-18	

Date of Report: March 16, 2018 Samples Submitted: February 22, 2018

Laboratory Reference: 1802-233

Project: 17-06520-000

TOTAL METAL EPA 6010D/7471B QUALITY CONTROL

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0226SM1					
Arsenic	ND	10	EPA 6010D	2-26-18	2-27-18	
Cadmium	ND	0.50	EPA 6010D	2-26-18	2-27-18	
Chromium	ND	0.50	EPA 6010D	2-26-18	2-27-18	
Lead	ND	5.0	EPA 6010D	2-26-18	2-27-18	
Laboratory ID:	MB0223S1					
Mercury	ND	0.25	EPA 7471B	2-23-18	2-23-18	

	_				Source	_	rcent	Recovery		RPD	
Analyte	Re	sult	Spike	Level	Result	Rec	overy	Limits	RPD	Limit	Flags
DUPLICATE											
Laboratory ID:	02-2	55-01									
	ORIG	DUP									
Arsenic	ND	ND	NA	NA			NA	NA	NA	20	
Cadmium	ND	ND	NA	NA		ı	NA	NA	NA	20	
Chromium	262	241	NA	NA		ı	NA	NA	8	20	
Lead	9.60	9.45	NA	NA		ı	NA	NA	2	20	
Laboratory ID:	02-2	17-01									
Mercury	ND	ND	NA	NA		-	NA	NA	NA	20	
Weredry	ND	ND	INA	IVA		<u>'</u>	INA	IVA	IVA	20	
MATRIX SPIKES											
Laboratory ID:	02-2	55-01									
	MS	MSD	MS	MSD		MS	MSD				
Arsenic	96.4	95.2	100	100	ND	96	95	75-125	1	20	
Cadmium	45.5	46.3	50.0	50.0	ND	91	93	75-125	2	20	
Chromium	186	479	100	100	262	-76	218	75-125	88	20	
Lead	231	235	250	250	9.60	89	90	75-125	2	20	
Laboratory ID:	02-2	17-01									
Mercury	0.540	0.571	0.500	0.500	ND	108	114	80-120	6	20	

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B

Matrix: Soil

Offits. Hig/Kg (ppHI)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-04_12.5					
Laboratory ID:	02-233-03					
Arsenic	ND	13	EPA 6010D	2-28-18	2-28-18	
Cadmium	ND	0.64	EPA 6010D	2-28-18	2-28-18	
Chromium	16	0.64	EPA 6010D	2-28-18	2-28-18	
Lead	ND	6.4	EPA 6010D	2-28-18	2-28-18	
Mercury	ND	0.32	EPA 7471B	2-28-18	2-28-18	
OII						
Client ID:	B-05_7.5					
Laboratory ID:	02-233-06					
Arsenic	ND	12	EPA 6010D	2-28-18	2-28-18	
Cadmium	ND	0.60	EPA 6010D	2-28-18	2-28-18	
Chromium	430	3.0	EPA 6010D	2-28-18	2-28-18	
Lead	ND	6.0	EPA 6010D	2-28-18	2-28-18	
Mercury	ND	0.30	EPA 7471B	2-28-18	2-28-18	
Client ID:	B-05_12.5					
Laboratory ID:	02-233-07					
Arsenic	ND	13	EPA 6010D	2-28-18	2-28-18	
Cadmium	ND	0.63	EPA 6010D	2-28-18	2-28-18	
Chromium	11	0.63	EPA 6010D	2-28-18	2-28-18	
Lead	ND	6.3	EPA 6010D	2-28-18	2-28-18	
Mercury	ND	0.32	EPA 7471B	2-28-18	2-28-18	
Client ID:	B-14_5					
Laboratory ID:	02-233-09					
Arsenic	ND	13	EPA 6010D	2-28-18	2-28-18	
Cadmium	ND	0.67	EPA 6010D	2-28-18	2-28-18	
Chromium	23	0.67	EPA 6010D	2-28-18	2-28-18	
Lead	31	6.7	EPA 6010D	2-28-18	2-28-18	
Mercury	ND	0.34	EPA 7471B	2-28-18	2-28-18	

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B

Matrix: Soil

omis. mg/rtg (ppm)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-14_10					
Laboratory ID:	02-233-10					
Arsenic	ND	12	EPA 6010D	2-28-18	2-28-18	
Cadmium	ND	0.62	EPA 6010D	2-28-18	2-28-18	
Chromium	17	0.62	EPA 6010D	2-28-18	2-28-18	
Lead	33	6.2	EPA 6010D	2-28-18	2-28-18	
Mercury	ND	0.31	EPA 7471B	2-28-18	2-28-18	
Client ID:	B-17_10					
Laboratory ID:	02-233-18					
Arsenic	ND	12	EPA 6010D	2-28-18	2-28-18	
Cadmium	ND	0.61	EPA 6010D	2-28-18	2-28-18	
Chromium	23	0.61	EPA 6010D	2-28-18	2-28-18	
Lead	ND	6.1	EPA 6010D	2-28-18	2-28-18	
Mercury	ND	0.31	EPA 7471B	2-28-18	2-28-18	
Client ID:	B-17_15					
Laboratory ID:	02-233-19					
Arsenic	ND	11	EPA 6010D	2-28-18	2-28-18	
Cadmium	ND	0.54	EPA 6010D	2-28-18	2-28-18	
Chromium	9.6	0.54	EPA 6010D	2-28-18	2-28-18	
Lead	ND	5.4	EPA 6010D	2-28-18	2-28-18	
Mercury	ND	0.27	EPA 7471B	2-28-18	2-28-18	

Date of Report: March 16, 2018 Samples Submitted: February 22, 2018

Laboratory Reference: 1802-233

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B QUALITY CONTROL

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0228SM1					
Arsenic	ND	10	EPA 6010D	2-28-18	2-28-18	
Cadmium	ND	0.50	EPA 6010D	2-28-18	2-28-18	
Chromium	ND	0.50	EPA 6010D	2-28-18	2-28-18	
Lead	ND	5.0	EPA 6010D	2-28-18	2-28-18	
Laboratory ID:	MB0228S1					
Mercury	ND	0.25	EPA 7471B	2-28-18	2-28-18	

					Source	Pei	rcent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Rec	overy	Limits	RPD	Limit	Flags
DUPLICATE											
Laboratory ID:	02-23	30-03									
	ORIG	DUP									
Arsenic	ND	ND	NA	NA		1	NA	NA	NA	20	
Cadmium	ND	ND	NA	NA		1	NΑ	NA	NA	20	
Chromium	56.1	67.8	NA	NA		1	NΑ	NA	19	20	
Lead	ND	ND	NA	NA		1	NA	NA	NA	20	
Laboratory ID:	02-2	49-03									
Mercury	ND	ND	NA	NA		1	NA	NA	NA	20	
MATRIX SPIKES											
Laboratory ID:	02-23	30-03									
	MS	MSD	MS	MSD		MS	MSD				
Arsenic	95.9	94.6	100	100	ND	96	95	75-125	1	20	
Cadmium	44.7	44.6	50.0	50.0	ND	89	89	75-125	0	20	
Chromium	143	137	100	100	56.1	87	81	75-125	4	20	
Lead	215	217	250	250	ND	86	87	75-125	1	20	
Laboratory ID:	02-24	49-03									
Mercury	0.551	0.567	0.500	0.500	0.00670	109	112	80-120	3	20	

Project: 17-06520-000

TCLP CHROMIUM EPA 1311/6010D

Matrix: TCLP Extract Units: mg/L (ppm)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-05_7.5					
Laboratory ID:	02-233-06					
Chromium	ND	0.020	EPA 6010D	3-14-18	3-14-18	

Project: 17-06520-000

TCLP CHROMIUM EPA 1311/6010D QUALITY CONTROL

Matrix: TCLP Extract Units: mg/L (ppm)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0314TM1					
Chromium	ND	0.020	EPA 6010D	3-14-18	3-14-18	

					Source	Pe	rcent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Rec	overy	Limits	RPD	Limit	Flags
DUPLICATE											
Laboratory ID:	02-1	51-15									
	ORIG	DUP									
Chromium	ND	ND	NA	NA			NA	NA	NA	20	
MATRIX SPIKES											
Laboratory ID:	02-1	51-15									
	MS	MSD	MS	MSD		MS	MSD				
Chromium	3.82	3.80	4.00	4.00	ND	96	95	75-125	1	20	

Project: 17-06520-000

% MOISTURE

Date Analyzed: 2-23,26&27-18

Client ID	Lab ID	% Moisture
B-04_12.5	02-233-03	21
B-05_7.5	02-233-06	16
B-05_12.5	02-233-07	21
B-14_5	02-233-09	26
B-14_10	02-233-10	19
B-15_5	02-233-13	42
B-15_7.5	02-233-14	34
B-15_15	02-233-15	26
B-17_10	02-233-18	18
B-17_15	02-233-19	7



Data Qualifiers and Abbreviations

- A Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
- B The analyte indicated was also found in the blank sample.
- C The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
- E The value reported exceeds the quantitation range and is an estimate.
- F Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
- H The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
- I Compound recovery is outside of the control limits.
- J The value reported was below the practical quantitation limit. The value is an estimate.
- K Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
- L The RPD is outside of the control limits.
- M Hydrocarbons in the gasoline range are impacting the diesel range result.
- M1 Hydrocarbons in the gasoline range (toluene-naphthalene) are present in the sample.
- N Hydrocarbons in the lube oil range are impacting the diesel range result.
- N1 Hydrocarbons in diesel range are impacting lube oil range results.
- O Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
- P The RPD of the detected concentrations between the two columns is greater than 40.
- Q Surrogate recovery is outside of the control limits.
- S Surrogate recovery data is not available due to the necessary dilution of the sample.
- T The sample chromatogram is not similar to a typical .
- U The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- U1 The practical quantitation limit is elevated due to interferences present in the sample.
- V Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
- W Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
- X Sample extract treated with a mercury cleanup procedure.
- X1- Sample extract treated with a sulfuric acid/silica gel cleanup procedure.
- Y The calibration verification for this analyte exceeded the 20% drift specified in method 8260C, and therefore the reported result should be considered an estimate. The overall performance of the calibration verification standard met the acceptance criteria of the method.

7 -

ND - Not Detected at PQL

PQL - Practical Quantitation Limit

RPD - Relative Percent Difference



OnSite Environmental Inc.

Chain of Custody

Page of 2

	Analytical Laboratory Testing Services 14648 NE 95th Street • Redmond, WA 98052		urnaround Req (in working da			L	.abd	orate	ory	Nur	nbe	er:			02	2 -	2	3	3							
Project	Number: 1706620-000 Name: Pacific Right Bank Manager: Bruce Carpenter	☐ 2 E		1 Day	Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX		NWTPH-Dx (☐ Acid / SG Clean-up)	Volatiles 8260C	genated Volatiles 8	EPA 8011 (Waters Only)	Semivolatiles 8270D/SIM (with low-level PAHs)	s 8270D/SIM (low-level) PAHS		Organochlorine Pesticides 8081B	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	TCLP Metals	HEM (oil and grease) 1664A		ELP Gr		% Moisture
Lab ID	Sample Identification	Sample	d Sampled	Matrix	Nun	NW	NN	NN	NN	Vola	Halo	EDB	Serr (with	PAH		Orga	Orga	SHO	Tota	-	,	HEN		+	1	% W
1	B-04_2.5	2/30/16	8 +300	soil	1	A	A		A				98	¥	A					*	08					X
2	B-04_ 7.5	2/20/18	1305		2	A	A		A					A	A					A				t		X
3	13-04_12.5	2/21/18	1430		2	9	A		A						A									-		
T	13-04_17.5	2/2/18	1450		2	A	A		A					A	A					A						X
5	B-05_ 2.5	2/20/18	1300		2	M	A		A					A	A					A						X
6	B-05_87.5	2/20/18	1305		2		A		100						8									X		
7	B-05_12.5	2/20/18	1310		2	6	A		A					•	A					6						•
8	B-05_17.5	2/20/18			a	A	A		A					A	A					A						X
9	13-14=5	2/21/200		-	22	-			A		1				A					6						
10	B-14_10	2/21/18		1	2	9	1		A						Δ									1		
	Signature		Company	-			Date			Time			Com	men	ts/Sp	ecial	Instr	uction	ns	7						1
Relinqu	dished Hammy MAT	0	Herrera	a			2	22/1	8	15	05			1		^	2	2.1	/	-	1		~	- A	0	
Receiv	ed		OSF				21	22	118	- 1	50	5	- 1	H	-	H	K	-M.	, V.	2	5	٥.	11	14	>	
Relinq	uished												6	00	100	led	2	23	118	3.	02	6	T	1		
Receiv	ed												6	20	die	lec	12	15	28/	18.	03	10	-0	3		
Relinqu	uished												>	(-1	Ada	led	31	181	18.	D3	(s	(S)	2	/		
Receiv	ed											1										Level				
Review	/ed/Date		Reviewed/Dat	te									Chro	matc	gram	s wit	h fina	al rep	ort [Ele	ctroni	c Data	ı Deliv	erable	s (EDDs) 🗆

OnSite Environmental Inc.

Chain of Custody

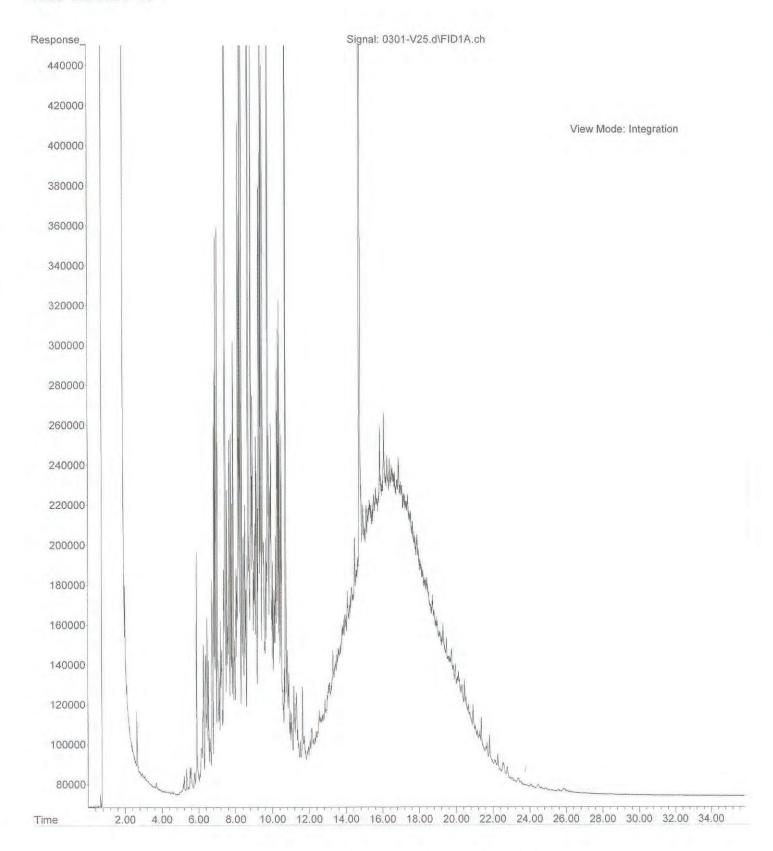
Page 2 of Z

	Analytical Laboratory Testing Services 14648 NE 95th Street • Redmond, WA 98052		rnaround Red n working da			L	abo	orato	ry l	Nun	ber		C	12	- 2	23	3								
Project Number Project Name: Project Manag Sampled by:	1706520-000 Pacific Right Bank			1 Day	Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX		NWTPH-Dx (Acid / SG Clean-up)	Volatiles 8260C	EDB EPA 8011 (Waters Only)	Semivolatiles 8270D/SIM (with low-level PAHs)	PAHS 8270D/SIM (low-level) _ DALLA	PCBs 8082A	Organochlorine Pesticides 8081B	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	TCLP Metals	HEM (oil and grease) 1664A				% Moisture
11	8-14_15	2/21/18	9:15	Soil	2	A			A				A						A						X
12	B-14_20	212/18	9:20		2	A	h		A				A	A					A						X
13	B-15_5	2/21/18	13:30		2	X	A		A				R						8)					0
114	B-15_7.5	2/21/18	12:35		2		A		A				(8)	A				((V)						8
15	B-15 + 15	2/21/18	1366		2	X	A		A				X	A					(1)						0
No	B-15_18.5	2/21/18	1305		2	A	A		A				A	A					A						X
17	B-17_7.5	2/20/18	₹·14:35		2	A	A		A				A	A					6						X
18	B-17_ 10	2/20/18	14:40		2	•	A		A					A											
19	B-17_15	2/20/18	14:45		2	9	A		A					A											1
20	B-17-20	2/20/18	M:50	V	2	A	A		A				A	A					A						X
Relinquished	Signature ALAMAU MIT		l proce	· M	400		Date			Time	~	Cor	nmer	its/Sp	ecial	Instr	uction	15							
Received	HALWING WICE		Herre	901				22/		150															
Relinquished	00		OSE				213	2211	0	15	08														
Received									+			1													
Relinquished									1																
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Reviewed/Da	te		Reviewed/Da	te								Chro	omato	ogran	ns wit	h fina	al rep	ort 🗆	Elec	ctroni	c Data	Delive	erables ((EDDs)	

File :X:\DIESELS\VIGO\DATA\V180301\0301-V25.d
Operator : JT
Acquired : 1 Mar 2018 22:59 using AcqMethod V
Instrument : Vigo using AcqMethod V171020F.M

Sample Name: 02-233-06

Misc Info : Vial Number: 25





14648 NE 95th Street, Redmond, WA 98052 • (425) 883-3881

March 8, 2018

Bruce Carpenter Herrera Environmental Consultants, Inc. 2200 6th Avenue, Suite 1100 Seattle, WA 98121

Re: Analytical Data for Project 17-06520-000

Laboratory Reference No. 1802-249

Dear Bruce:

Enclosed are the analytical results and associated quality control data for samples submitted on February 26, 2018.

The standard policy of OnSite Environmental, Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

David Baumeister Project Manager

Enclosures

Date of Report: March 8, 2018

Samples Submitted: February 26, 2018 Laboratory Reference: 1802-249

Project: 17-06520-000

Case Narrative

Samples were collected on February 22 and 23, 2018 and received by the laboratory on February 26, 2018. They were maintained at the laboratory at a temperature of 2°C to 6°C.

Please note that any and all soil sample results are reported on a dry-weight basis, unless otherwise noted below.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

Project: 17-06520-000

NWTPH-HCID

Matrix: Soil

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	B-06_2.5	. 42	Motriou	Tioparoa	Allaryzou	i lago
Laboratory ID:	02-249-01					
Gasoline Range Organics	ND	25	NWTPH-HCID	2-26-18	2-26-18	
Diesel Range Organics	ND	62	NWTPH-HCID	2-26-18	2-26-18	
Lube Oil	Detected	120	NWTPH-HCID	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits	111111111111111111111111111111111111111	2 20 .0	2 20 10	
o-Terphenyl	101	50-150				
	-					
Client ID:	B-06_12.5					
Laboratory ID:	02-249-03					
Gasoline Range Organics	ND	26	NWTPH-HCID	2-26-18	2-26-18	
Diesel Range Organics	ND	78	NWTPH-HCID	2-26-18	2-26-18	U1
Lube Oil	Detected	130	NWTPH-HCID	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	103	50-150				
Client ID:	B-09_5					
Laboratory ID:	02-249-04					
Gasoline Range Organics	ND	31	NWTPH-HCID	2-26-18	2-26-18	
Diesel Range Organics	ND	78	NWTPH-HCID	2-26-18	2-26-18	
Lube Oil Range Organics	ND	160	NWTPH-HCID	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	98	50-150				
Client ID:	B-09_15					
Laboratory ID:	02-249-06					
Gasoline Range Organics	ND	31	NWTPH-HCID	2-26-18	2-26-18	
Diesel Range Organics	ND	78	NWTPH-HCID	2-26-18	2-26-18	
Lube Oil Range Organics	ND	160	NWTPH-HCID	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	109	50-150				
Client ID:	B-13_7.5					
Laboratory ID:	02-249-07					
Gasoline Range Organics	ND	34	NWTPH-HCID	2-26-18	2-26-18	
Diesel Range Organics	ND	84	NWTPH-HCID	2-26-18	2-26-18	
Lube Oil Range Organics	ND	170	NWTPH-HCID	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	112	50-150				

Project: 17-06520-000

NWTPH-HCID

Matrix: Soil

3 3 (11)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-13_15					
Laboratory ID:	02-249-09					
Gasoline Range Organics	ND	26	NWTPH-HCID	2-26-18	2-26-18	
Diesel Range Organics	ND	66	NWTPH-HCID	2-26-18	2-26-18	
Lube Oil Range Organics	ND	130	NWTPH-HCID	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	96	50-150				
Client ID:	B-16_2.5					
Laboratory ID:	02-249-11					
Gasoline Range Organics	ND	29	NWTPH-HCID	2-26-18	2-26-18	
Diesel Range Organics	ND	71	NWTPH-HCID	2-26-18	2-26-18	
Lube Oil Range Organics	ND	140	NWTPH-HCID	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	96	50-150				
Client ID:	B-16_7.5					
Laboratory ID:	02-249-12					
Gasoline Range Organics	ND	23	NWTPH-HCID	2-26-18	2-26-18	
Diesel Range Organics	ND	58	NWTPH-HCID	2-26-18	2-26-18	
Lube Oil Range Organics	ND	120	NWTPH-HCID	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	99	50-150				
Client ID:	B-16_17.5					
Laboratory ID:	02-249-14					
Gasoline Range Organics	ND	27	NWTPH-HCID	2-26-18	2-26-18	
Diesel Range Organics	ND	66	NWTPH-HCID	2-26-18	2-26-18	
Lube Oil Range Organics	ND	130	NWTPH-HCID	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits			2 20 10	
o-Terphenyl	92	50-150				
	<u></u>					

Date of Report: March 8, 2018

Samples Submitted: February 26, 2018 Laboratory Reference: 1802-249

Project: 17-06520-000

NWTPH-HCID QUALITY CONTROL

Matrix: Soil

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0226S2					
Gasoline Range Organics	ND	20	NWTPH-HCID	2-26-18	2-26-18	_
Diesel Range Organics	ND	50	NWTPH-HCID	2-26-18	2-26-18	
Lube Oil Range Organics	ND	100	NWTPH-HCID	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	100	50-150				

Date of Report: March 8, 2018 Samples Submitted: February 26, 2018

Laboratory Reference: 1802-249

Project: 17-06520-000

NWTPH-Dx

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-06_2.5					
Laboratory ID:	02-249-01					
Diesel Range Organics	ND	31	NWTPH-Dx	3-2-18	3-2-18	
Lube Oil	150	62	NWTPH-Dx	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	86	50-150				
Client ID:	B-06_10					
Laboratory ID:	02-249-02					
Diesel Range Organics	ND	31	NWTPH-Dx	3-2-18	3-2-18	
Lube Oil Range Organics	ND	61	NWTPH-Dx	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	85	50-150				
Client ID:	B-06_12.5					
Laboratory ID:	02-249-03					
Diesel Range Organics	ND	32	NWTPH-Dx	3-2-18	3-2-18	
Lube Oil	130	64	NWTPH-Dx	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenvl	89	50-150				

Date of Report: March 8, 2018

Samples Submitted: February 26, 2018 Laboratory Reference: 1802-249

Project: 17-06520-000

NWTPH-Dx QUALITY CONTROL

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0302S1					
Diesel Range Organics	ND	25	NWTPH-Dx	3-2-18	3-2-18	
Lube Oil Range Organics	ND	50	NWTPH-Dx	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	97	50-150				

					Source	Percent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Recovery	Limits	RPD	Limit	Flags
DUPLICATE										
Laboratory ID:	02-24	19-02								
	ORIG	DUP								
Diesel Range	ND	ND	NA	NA		NA	NA	NA	NA	
Lube Oil Range	ND	ND	NA	NA		NA	NA	NA	NA	
Surrogate:										
o-Terphenyl						85 86	50-150			

Project: 17-06520-000

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-06_2.5					
Laboratory ID:	02-249-01					
Benzo[a]anthracene	ND	0.0083	EPA 8270D/SIM	2-28-18	2-28-18	
Chrysene	ND	0.0083	EPA 8270D/SIM	2-28-18	2-28-18	
Benzo[b]fluoranthene	ND	0.0083	EPA 8270D/SIM	2-28-18	2-28-18	
Benzo(j,k)fluoranthene	ND	0.0083	EPA 8270D/SIM	2-28-18	2-28-18	
Benzo[a]pyrene	ND	0.0083	EPA 8270D/SIM	2-28-18	2-28-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0083	EPA 8270D/SIM	2-28-18	2-28-18	
Dibenz[a,h]anthracene	ND	0.0083	EPA 8270D/SIM	2-28-18	2-28-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	57	32 - 115				
Pyrene-d10	68	35 - 129				
Terphenyl-d14	64	33 - 114				

Project: 17-06520-000

PAHs EPA 8270D/SIM

			Date	Date	
Result	PQL	Method	Prepared	Analyzed	Flags
B-06_12.5					
02-249-03					
ND	0.0086	EPA 8270D/SIM	2-28-18	3-5-18	
ND	0.0086	EPA 8270D/SIM	2-28-18	3-5-18	
ND	0.0086	EPA 8270D/SIM	2-28-18	3-5-18	
ND	0.0086	EPA 8270D/SIM	2-28-18	3-5-18	
ND	0.0086	EPA 8270D/SIM	2-28-18	3-5-18	
ND	0.0086	EPA 8270D/SIM	2-28-18	3-5-18	
ND	0.0086	EPA 8270D/SIM	2-28-18	3-5-18	
Percent Recovery	Control Limits				
78	32 - 115				
90	35 - 129				
91	33 - 114				
	B-06_12.5 02-249-03 ND ND ND ND ND ND ND ND Percent Recovery 78 90	B-06_12.5 02-249-03 0.0086 ND 0.0086 ND 0.0086 ND 0.0086 ND 0.0086 ND 0.0086 ND 0.0086 Percent Recovery Control Limits 78 32 - 115 90 35 - 129	B-06_12.5 02-249-03 ND 0.0086 EPA 8270D/SIM Percent Recovery Control Limits 78 32 - 115 90 35 - 129	Result PQL Method Prepared B-06_12.5 02-249-03 ND 0.0086 EPA 8270D/SIM 2-28-18 Percent Recovery Control Limits 78 32 - 115 32 - 115 90 35 - 129 35 - 129	Result PQL Method Prepared Analyzed B-06_12.5 02-249-03 Secondary 12-24-18 3-5-18 ND 0.0086 EPA 8270D/SIM 2-28-18 3-5-18 Percent Recovery Control Limits 78 32 - 115 90 35 - 129 35 - 129 35 - 129

Project: 17-06520-000

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-09_5					
Laboratory ID:	02-249-04					
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	2-28-18	2-28-18	
Chrysene	0.012	0.010	EPA 8270D/SIM	2-28-18	2-28-18	
Benzo[b]fluoranthene	0.011	0.010	EPA 8270D/SIM	2-28-18	2-28-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	2-28-18	2-28-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	2-28-18	2-28-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	2-28-18	2-28-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	2-28-18	2-28-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	39	32 - 115				
Pyrene-d10	45	35 - 129				
Terphenyl-d14	42	33 - 114				

Date of Report: March 8, 2018 Samples Submitted: February 26, 2018

Laboratory Reference: 1802-249

Project: 17-06520-000

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-09_15					
Laboratory ID:	02-249-06					
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	2-28-18	3-4-18	
Chrysene	ND	0.010	EPA 8270D/SIM	2-28-18	3-4-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	2-28-18	3-4-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	2-28-18	3-4-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	2-28-18	3-4-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	2-28-18	3-4-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	2-28-18	3-4-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	94	32 - 115				
Pyrene-d10	97	35 - 129				
Terphenyl-d14	106	33 - 114				

Date of Report: March 8, 2018

Samples Submitted: February 26, 2018 Laboratory Reference: 1802-249

Project: 17-06520-000

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-13_7.5					
Laboratory ID:	02-249-07					
Benzo[a]anthracene	ND	0.011	EPA 8270D/SIM	2-28-18	2-28-18	
Chrysene	0.016	0.011	EPA 8270D/SIM	2-28-18	2-28-18	
Benzo[b]fluoranthene	0.014	0.011	EPA 8270D/SIM	2-28-18	2-28-18	
Benzo(j,k)fluoranthene	ND	0.011	EPA 8270D/SIM	2-28-18	2-28-18	
Benzo[a]pyrene	ND	0.011	EPA 8270D/SIM	2-28-18	2-28-18	
Indeno(1,2,3-c,d)pyrene	ND	0.011	EPA 8270D/SIM	2-28-18	2-28-18	
Dibenz[a,h]anthracene	ND	0.011	EPA 8270D/SIM	2-28-18	2-28-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	47	32 - 115				
Pyrene-d10	57	35 - 129				
Terphenyl-d14	55	33 - 114				

Date of Report: March 8, 2018 Samples Submitted: February 26, 2018

Laboratory Reference: 1802-249

Project: 17-06520-000

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-13_15					
Laboratory ID:	02-249-09					
Benzo[a]anthracene	ND	0.0088	EPA 8270D/SIM	2-28-18	3-4-18	
Chrysene	ND	0.0088	EPA 8270D/SIM	2-28-18	3-4-18	
Benzo[b]fluoranthene	ND	0.0088	EPA 8270D/SIM	2-28-18	3-4-18	
Benzo(j,k)fluoranthene	ND	0.0088	EPA 8270D/SIM	2-28-18	3-4-18	
Benzo[a]pyrene	ND	0.0088	EPA 8270D/SIM	2-28-18	3-4-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0088	EPA 8270D/SIM	2-28-18	3-4-18	
Dibenz[a,h]anthracene	ND	0.0088	EPA 8270D/SIM	2-28-18	3-4-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	79	32 - 115				
Pyrene-d10	85	35 - 129				
Terphenyl-d14	91	33 - 114				

Date of Report: March 8, 2018

Samples Submitted: February 26, 2018 Laboratory Reference: 1802-249

Project: 17-06520-000

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-16_2.5					
Laboratory ID:	02-249-11					
Benzo[a]anthracene	ND	0.0095	EPA 8270D/SIM	2-28-18	3-5-18	
Chrysene	ND	0.0095	EPA 8270D/SIM	2-28-18	3-5-18	
Benzo[b]fluoranthene	ND	0.0095	EPA 8270D/SIM	2-28-18	3-5-18	
Benzo(j,k)fluoranthene	ND	0.0095	EPA 8270D/SIM	2-28-18	3-5-18	
Benzo[a]pyrene	ND	0.0095	EPA 8270D/SIM	2-28-18	3-5-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0095	EPA 8270D/SIM	2-28-18	3-5-18	
Dibenz[a,h]anthracene	ND	0.0095	EPA 8270D/SIM	2-28-18	3-5-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	64	32 - 115				
Pyrene-d10	74	35 - 129				
Terphenyl-d14	77	33 - 114				
reipnenyi-u 14	//	33 - 11 4				

Samples Submitted: February 26, 2018 Laboratory Reference: 1802-249

Project: 17-06520-000

PAHs EPA 8270D/SIM

			Date	Date	
Result	PQL	Method	Prepared	Analyzed	Flags
B-16_7.5					
02-249-12					
ND	0.0077	EPA 8270D/SIM	2-28-18	3-4-18	
ND	0.0077	EPA 8270D/SIM	2-28-18	3-4-18	
ND	0.0077	EPA 8270D/SIM	2-28-18	3-4-18	
ND	0.0077	EPA 8270D/SIM	2-28-18	3-4-18	
ND	0.0077	EPA 8270D/SIM	2-28-18	3-4-18	
ND	0.0077	EPA 8270D/SIM	2-28-18	3-4-18	
ND	0.0077	EPA 8270D/SIM	2-28-18	3-4-18	
Percent Recovery	Control Limits				
86	32 - 115				
90	35 - 129				
93	33 - 114				
	B-16_7.5 02-249-12 ND ND ND ND ND ND ND ND Percent Recovery 86 90	B-16_7.5 02-249-12 0.0077 ND 0.0077 ND 0.0077 ND 0.0077 ND 0.0077 ND 0.0077 ND 0.0077 Percent Recovery Control Limits 86 32 - 115 90 35 - 129	B-16_7.5 02-249-12 ND 0.0077 EPA 8270D/SIM Percent Recovery Control Limits 86 32 - 115 90 35 - 129	Result PQL Method Prepared B-16_7.5 02-249-12	Result PQL Method Prepared Analyzed B-16_7.5 02-249-12 Separate 3-4-18 ND 0.0077 EPA 8270D/SIM 2-28-18 3-4-18 Percent Recovery Control Limits 86 32 - 115 32 - 115 90 35 - 129 35 - 129

Samples Submitted: February 26, 2018 Laboratory Reference: 1802-249

Project: 17-06520-000

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-16_17.5					
Laboratory ID:	02-249-14					
Benzo[a]anthracene	ND	0.0089	EPA 8270D/SIM	2-28-18	3-5-18	
Chrysene	ND	0.0089	EPA 8270D/SIM	2-28-18	3-5-18	
Benzo[b]fluoranthene	ND	0.0089	EPA 8270D/SIM	2-28-18	3-5-18	
Benzo(j,k)fluoranthene	ND	0.0089	EPA 8270D/SIM	2-28-18	3-5-18	
Benzo[a]pyrene	ND	0.0089	EPA 8270D/SIM	2-28-18	3-5-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0089	EPA 8270D/SIM	2-28-18	3-5-18	
Dibenz[a,h]anthracene	ND	0.0089	EPA 8270D/SIM	2-28-18	3-5-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	88	32 - 115				
Pyrene-d10	95	35 - 129				
Terphenyl-d14	98	33 - 114				

Samples Submitted: February 26, 2018

Laboratory Reference: 1802-249

Project: 17-06520-000

PAHS EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0228S2					
Benzo[a]anthracene	ND	0.0067	EPA 8270D/SIM	2-28-18	2-28-18	
Chrysene	ND	0.0067	EPA 8270D/SIM	2-28-18	2-28-18	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270D/SIM	2-28-18	2-28-18	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270D/SIM	2-28-18	2-28-18	
Benzo[a]pyrene	ND	0.0067	EPA 8270D/SIM	2-28-18	2-28-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0067	EPA 8270D/SIM	2-28-18	2-28-18	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270D/SIM	2-28-18	2-28-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	57	32 - 115				
Pyrene-d10	76	35 - 129				
Terphenyl-d14	75	33 - 114				

Samples Submitted: February 26, 2018 Laboratory Reference: 1802-249

Project: 17-06520-000

PAHS EPA 8270D/SIM MS/MSD QUALITY CONTROL

					Source	Per	cent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Rec	overy	Limits	RPD	Limit	Flags
MATRIX SPIKES											
Laboratory ID:	02-24	49-11									
	MS	MSD	MS	MSD		MS	MSD				
Benzo[a]anthracene	0.156	0.161	0.167	0.167	ND	93	96	27 - 143	3	23	
Chrysene	0.146	0.151	0.167	0.167	ND	87	90	22 - 130	3	24	
Benzo[b]fluoranthene	0.163	0.169	0.167	0.167	ND	98	101	15 - 141	4	26	
Benzo(j,k)fluoranthene	0.142	0.148	0.167	0.167	ND	85	89	42 - 112	4	24	
Benzo[a]pyrene	0.141	0.145	0.167	0.167	ND	84	87	33 - 126	3	26	
Indeno(1,2,3-c,d)pyrene	0.133	0.139	0.167	0.167	ND	80	83	30 - 125	4	25	
Dibenz[a,h]anthracene	0.134	0.139	0.167	0.167	ND	80	83	31 - 124	4	22	
Surrogate:											
2-Fluorobiphenyl						67	73	32 - 115			
Pyrene-d10						84	87	35 - 129			
Terphenyl-d14						86	89	33 - 114			

Samples Submitted: February 26, 2018 Laboratory Reference: 1802-249

Project: 17-06520-000

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-06_10					
Laboratory ID:	02-249-02					
Benzo[a]anthracene	ND	0.0082	EPA 8270D/SIM	3-5-18	3-5-18	
Chrysene	ND	0.0082	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[b]fluoranthene	ND	0.0082	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo(j,k)fluoranthene	ND	0.0082	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[a]pyrene	ND	0.0082	EPA 8270D/SIM	3-5-18	3-5-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0082	EPA 8270D/SIM	3-5-18	3-5-18	
Dibenz[a,h]anthracene	ND	0.0082	EPA 8270D/SIM	3-5-18	3-5-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	71	32 - 115				
Pyrene-d10	78	35 - 129				
Terphenyl-d14	82	33 - 114				

Samples Submitted: February 26, 2018 Laboratory Reference: 1802-249

Project: 17-06520-000

PAHS EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0305S1					
Benzo[a]anthracene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Chrysene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[a]pyrene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	84	32 - 115				
Pyrene-d10	94	35 - 129				
Terphenyl-d14	102	33 - 114				

Date of Report: March 8, 2018 Samples Submitted: February 26, 2018

Laboratory Reference: 1802-249

Project: 17-06520-000

PAHS EPA 8270D/SIM SB/SBD QUALITY CONTROL

					Per	cent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Reco	overy	Limits	RPD	Limit	Flags
SPIKE BLANKS										
Laboratory ID:	SB03	05S1								
	SB	SBD	SB	SBD	SB	SBD				
Benzo[a]anthracene	0.0878	0.0851	0.0833	0.0833	105	102	64 - 135	3	15	
Chrysene	0.0869	0.0844	0.0833	0.0833	104	101	70 - 119	3	15	
Benzo[b]fluoranthene	0.0826	0.0807	0.0833	0.0833	99	97	54 - 135	2	15	
Benzo(j,k)fluoranthene	0.0872	0.0839	0.0833	0.0833	105	101	66 - 122	4	15	
Benzo[a]pyrene	0.0816	0.0789	0.0833	0.0833	98	95	62 - 125	3	15	
Indeno(1,2,3-c,d)pyrene	0.0752	0.0724	0.0833	0.0833	90	87	55 - 129	4	15	
Dibenz[a,h]anthracene	0.0769	0.0756	0.0833	0.0833	92	91	58 - 125	2	15	
Surrogate:										
2-Fluorobiphenyl					88	87	32 - 115			
Pyrene-d10					99	96	35 - 129			
Terphenyl-d14					105	102	33 - 114			

Samples Submitted: February 26, 2018 Laboratory Reference: 1802-249

Project: 17-06520-000

PCBs EPA 8082A

Matrix: Soil

			Date	Date	
Result	PQL	Method	Prepared	Analyzed	Flags
B-06_2.5					
02-249-01					
ND	0.062	EPA 8082A	3-6-18	3-6-18	
ND	0.062	EPA 8082A	3-6-18	3-6-18	
ND	0.062	EPA 8082A	3-6-18	3-6-18	
ND	0.062	EPA 8082A	3-6-18	3-6-18	
ND	0.062	EPA 8082A	3-6-18	3-6-18	
ND	0.062	EPA 8082A	3-6-18	3-6-18	
ND	0.062	EPA 8082A	3-6-18	3-6-18	
Percent Recovery	Control Limits				
79	40-134				
B-06_12.5					
02-249-03					
ND	0.064	EPA 8082A	3-6-18	3-6-18	
ND	0.064	EPA 8082A	3-6-18	3-6-18	
ND	0.064	EPA 8082A	3-6-18	3-6-18	
ND	0.064	EPA 8082A	3-6-18	3-6-18	
ND	0.064	EPA 8082A	3-6-18	3-6-18	
ND	0.064	EPA 8082A	3-6-18	3-6-18	
ND	0.064	EPA 8082A	3-6-18	3-6-18	
Percent Recovery	Control Limits				
77	40-134				
	B-06_2.5 02-249-01 ND ND ND ND ND ND ND Percent Recovery 79 B-06_12.5 02-249-03 ND	B-06_2.5 02-249-01 ND	B-06_2.5	Result PQL Method Prepared B-06_2.5 02-249-01 3-6-18 ND 0.062 EPA 8082A 3-6-18 Percent Recovery Control Limits 79 40-134 B-06_12.5 02-249-03 0.064 EPA 8082A 3-6-18 ND 0.064 EPA 8082A <t< td=""><td>Result PQL Method Prepared Analyzed B-06_2.5 02-249-01 B-06_2.5 02-249-01 B-06_2.5 02-249-01 B-06_2.5 02-249-01 B-06_2.5 02-249-01 B-06_2.5 02-249-03 B-06_12.5 02-249-03 B-06_12.5 00-064 B-06_4.25 00-064 B-06_6.26 00-064 B-06_6.26 00-064 B-06_6.26 00-064 B-06_6.26 00-064 B-06_6.26 00-064 B-06_6.26 00-064 B-06_6.26 00-064 B-06_6.26 00-064 B-06_6.26 00-064 B-06_6.26 00-064 00-064 B-06_6.26 00-064 00-064 B-06_6.26 00-064 00-06</td></t<>	Result PQL Method Prepared Analyzed B-06_2.5 02-249-01 B-06_2.5 02-249-01 B-06_2.5 02-249-01 B-06_2.5 02-249-01 B-06_2.5 02-249-01 B-06_2.5 02-249-03 B-06_12.5 02-249-03 B-06_12.5 00-064 B-06_4.25 00-064 B-06_6.26 00-064 B-06_6.26 00-064 B-06_6.26 00-064 B-06_6.26 00-064 B-06_6.26 00-064 B-06_6.26 00-064 B-06_6.26 00-064 B-06_6.26 00-064 B-06_6.26 00-064 B-06_6.26 00-064 00-064 B-06_6.26 00-064 00-064 B-06_6.26 00-064 00-06

Date of Report: March 8, 2018 Samples Submitted: February 26, 2018

Laboratory Reference: 1802-249

Project: 17-06520-000

PCBs EPA 8082A QUALITY CONTROL

Matrix: Soil

Units: mg/Kg (ppm)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0306S1					
Aroclor 1016	ND	0.050	EPA 8082A	3-6-18	3-6-18	
Aroclor 1221	ND	0.050	EPA 8082A	3-6-18	3-6-18	
Aroclor 1232	ND	0.050	EPA 8082A	3-6-18	3-6-18	
Aroclor 1242	ND	0.050	EPA 8082A	3-6-18	3-6-18	
Aroclor 1248	ND	0.050	EPA 8082A	3-6-18	3-6-18	
Aroclor 1254	ND	0.050	EPA 8082A	3-6-18	3-6-18	
Aroclor 1260	ND	0.050	EPA 8082A	3-6-18	3-6-18	

Surrogate: Percent Recovery Control Limits DCB 93 40-134

					Source	Pe	rcent	Recovery		RPD	
Analyte	Re	sult	Spike	Level	Result	Rec	overy	Limits	RPD	Limit	Flags
MATRIX SPIKES											
Laboratory ID:	02-2	49-01									
	MS	MSD	MS	MSD		MS	MSD				
Aroclor 1260	0.384	0.388	0.500	0.500	ND	77	78	34-126	1	16	
Surrogate:											
DCB						73	74	40-134			

Date of Report: March 8, 2018 Samples Submitted: February 26, 2018 Laboratory Reference: 1802-249

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B

Matrix: Soil

Units: mg/kg (ppm)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-06_2.5			•	-	
Laboratory ID:	02-249-01					
Arsenic	ND	12	EPA 6010D	3-6-18	3-6-18	
Cadmium	0.75	0.62	EPA 6010D	3-6-18	3-6-18	
Chromium	13	0.62	EPA 6010D	3-6-18	3-6-18	
Lead	20	6.2	EPA 6010D	3-6-18	3-6-18	
Mercury	ND	0.31	EPA 7471B	2-28-18	2-28-18	
Client ID:	B-06_12.5					
Laboratory ID:	02-249-03					
Arsenic	ND	13	EPA 6010D	3-6-18	3-6-18	
Cadmium	ND	0.64	EPA 6010D	3-6-18	3-6-18	
Chromium	11	0.64	EPA 6010D	3-6-18	3-6-18	
Lead	ND	6.4	EPA 6010D	3-6-18	3-6-18	
Mercury	ND	0.32	EPA 7471B	2-28-18	2-28-18	
Client ID:	B-09_5					
Laboratory ID:	02-249-04					
Arsenic	ND	15	EPA 6010D	3-6-18	3-6-18	
Cadmium	ND	0.77	EPA 6010D	3-6-18	3-6-18	
Chromium	16	0.77	EPA 6010D	3-6-18	3-6-18	
Lead	20	7.7	EPA 6010D	3-6-18	3-6-18	
Mercury	ND	0.39	EPA 7471B	2-28-18	2-28-18	
Client ID:	B-09_15					
Laboratory ID:	02-249-06					
Arsenic	ND	16	EPA 6010D	3-6-18	3-6-18	
Cadmium	ND	0.78	EPA 6010D	3-6-18	3-6-18	
Chromium	24	0.78	EPA 6010D	3-6-18	3-6-18	
Lead	ND	7.8	EPA 6010D	3-6-18	3-6-18	
Mercury	ND	0.39	EPA 7471B	2-28-18	2-28-18	

Date of Report: March 8, 2018 Samples Submitted: February 26, 2018 Laboratory Reference: 1802-249

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B

Matrix: Soil

ome. mg/rtg (ppm)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-13_7.5					
Laboratory ID:	02-249-07					
Arsenic	ND	17	EPA 6010D	3-6-18	3-6-18	
Cadmium	ND	0.84	EPA 6010D	3-6-18	3-6-18	
Chromium	24	0.84	EPA 6010D	3-6-18	3-6-18	
Lead	12	8.4	EPA 6010D	3-6-18	3-6-18	
Mercury	ND	0.42	EPA 7471B	2-28-18	2-28-18	
Client ID:	B-13_15					
Laboratory ID:	02-249-09					
Arsenic	ND	13	EPA 6010D	3-6-18	3-6-18	
Cadmium	ND	0.66	EPA 6010D	3-6-18	3-6-18	
Chromium	8.9	0.66	EPA 6010D	3-6-18	3-6-18	
Lead	ND	6.6	EPA 6010D	3-6-18	3-6-18	
Mercury	ND	0.33	EPA 7471B	2-28-18	2-28-18	
Client ID:	B-16_2.5					
Laboratory ID:	02-249-11					
Arsenic	ND	14	EPA 6010D	3-6-18	3-6-18	
Cadmium	ND	0.71	EPA 6010D	3-6-18	3-6-18	
Chromium	18	0.71	EPA 6010D	3-6-18	3-6-18	
Lead	9.2	7.1	EPA 6010D	3-6-18	3-6-18	
Mercury	ND	0.36	EPA 7471B	2-28-18	2-28-18	
Client ID:	B-16_7.5					
Laboratory ID:	02-249-12					
Arsenic	ND	12	EPA 6010D	3-6-18	3-6-18	
Cadmium	1.6	0.58	EPA 6010D	3-6-18	3-6-18	
Chromium	28	0.58	EPA 6010D	3-6-18	3-6-18	
Lead	68	5.8	EPA 6010D	3-6-18	3-6-18	
Mercury	ND	0.29	EPA 7471B	2-28-18	2-28-18	

Samples Submitted: February 26, 2018 Laboratory Reference: 1802-249

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-16_17.5					
Laboratory ID:	02-249-14					
Arsenic	ND	13	EPA 6010D	3-6-18	3-6-18	
Cadmium	ND	0.66	EPA 6010D	3-6-18	3-6-18	
Chromium	16	0.66	EPA 6010D	3-6-18	3-6-18	
Lead	ND	6.6	EPA 6010D	3-6-18	3-6-18	
Mercury	ND	0.33	FPA 7471B	2-28-18	2-28-18	

Date of Report: March 8, 2018 Samples Submitted: February 26, 2018

Laboratory Reference: 1802-249

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B QUALITY CONTROL

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0306SM1					
Arsenic	ND	10	EPA 6010D	3-6-18	3-6-18	
Cadmium	ND	0.50	EPA 6010D	3-6-18	3-6-18	
Chromium	ND	0.50	EPA 6010D	3-6-18	3-6-18	
Lead	ND	5.0	EPA 6010D	3-6-18	3-6-18	
Laboratory ID:	MB0228S1					
Mercury	ND	0.25	EPA 7471B	2-28-18	2-28-18	<u> </u>

					Source	Pe	rcent	Recovery		RPD	
Analyte	Re	sult	Spike	Level	Result	Rec	overy	Limits	RPD	Limit	Flags
DUPLICATE											
Laboratory ID:	02-2	64-89									
	ORIG	DUP									
Arsenic	ND	ND	NA	NA			NA AV	NA	NA	20	
Cadmium	ND	ND	NA	NA		ı	NA	NA	NA	20	
Chromium	27.7	26.3	NA	NA		ı	NA	NA	5	20	
Lead	11.2	7.15	NA	NA			NA	NA	44	20	
Laboratory ID:	02-2	49-03									
Mercury	ND	ND	NA	NA			NA	NA	NA	20	
MATRIX SPIKES											
Laboratory ID:	02-2	64-89									
	MS	MSD	MS	MSD		MS	MSD				
Arsenic	95.3	97.2	100	100	ND	95	97	75-125	2	20	
Cadmium	44.6	45.3	50.0	50.0	ND	89	91	75-125	1	20	
Chromium	124	114	100	100	27.7	97	86	75-125	9	20	
Lead	226	230	250	250	11.2	86	87	75-125	1	20	
Laboratory ID:	02-2	49-03									
Mercury	0.551	0.567	0.500	0.500	0.00670	109	112	80-120	3	20	

Samples Submitted: February 26, 2018 Laboratory Reference: 1802-249

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-06_10					_
Laboratory ID:	02-249-02					
Arsenic	ND	12	EPA 6010D	3-6-18	3-6-18	_
Cadmium	ND	0.61	EPA 6010D	3-6-18	3-6-18	
Chromium	13	0.61	EPA 6010D	3-6-18	3-6-18	
Lead	ND	6.1	EPA 6010D	3-6-18	3-6-18	
Mercury	ND	0.31	EPA 7471B	3-7-18	3-7-18	

Date of Report: March 8, 2018 Samples Submitted: February 26, 2018 Laboratory Reference: 1802-249

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B QUALITY CONTROL

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0306SM1					
Arsenic	ND	10	EPA 6010D	3-5-18	3-5-18	
Cadmium	ND	0.50	EPA 6010D	3-5-18	3-5-18	
Chromium	ND	0.50	EPA 6010D	3-5-18	3-5-18	
Lead	ND	5.0	EPA 6010D	3-5-18	3-5-18	
Laboratory ID:	MB0307S2					
Mercury	ND	0.25	EPA 7471B	3-7-18	3-7-18	

Amalusta	Dav	14	Cmiles	Laval	Source	_	rcent	Recovery	DDD	RPD	Flores
Analyte	Kes	sult	Бріке	Level	Result	Rec	overy	Limits	RPD	Limit	Flags
DUPLICATE											
Laboratory ID:	02-26	64-89									
	ORIG	DUP									
Arsenic	ND	ND	NA	NA		1	NA	NA	NA	20	,
Cadmium	ND	ND	NA	NA		1	NΑ	NA	NA	20	
Chromium	27.7	26.3	NA	NA		1	NΑ	NA	5	20	
Lead	11.2	7.15	NA	NA		1	NA	NA	44	20	
Laboratory ID:	02-27	71-21									
Mercury	ND	ND	NA	NA		ľ	NA	NA	NA	20	
MATRIX SPIKES											
Laboratory ID:	02-26	64-89									
Education y 12.	MS	MSD	MS	MSD		MS	MSD				
Arsenic	95.5	97.0	100	100	ND	96	97	75-125	2	20	
Cadmium	44.6	45.3	50.0	50.0	ND	89	91	75-125	1	20	
Chromium	125	114	100	100	27.7	97	86	75-125	9	20	
Lead	226	230	250	250	11.2	86	87	75-125	2	20	
Laboratory ID:	02-27	71-21									
Mercury	0.605	0.590	0.500	0.500	0.0364	114	111	80-120	3	20	

Samples Submitted: February 26, 2018 Laboratory Reference: 1802-249

Project: 17-06520-000

% MOISTURE

Date Analyzed: 2-26&3-2-18

Client ID	Lab ID	% Moisture
B-06_2.5	02-249-01	19
B-06_10	02-249-02	19
B-06_12.5	02-249-03	22
B-09_5	02-249-04	35
B-09_15	02-249-06	36
B-13_7.5	02-249-07	40
B-13_15	02-249-09	24
B-16_2.5	02-249-11	30
B-16_7.5	02-249-12	14
B-16_17.5	02-249-14	25



Data Qualifiers and Abbreviations

- A Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
- B The analyte indicated was also found in the blank sample.
- C The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
- E The value reported exceeds the quantitation range and is an estimate.
- F Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
- H The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
- I Compound recovery is outside of the control limits.
- J The value reported was below the practical quantitation limit. The value is an estimate.
- K Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
- L The RPD is outside of the control limits.
- M Hydrocarbons in the gasoline range are impacting the diesel range result.
- M1 Hydrocarbons in the gasoline range (toluene-naphthalene) are present in the sample.
- N Hydrocarbons in the lube oil range are impacting the diesel range result.
- N1 Hydrocarbons in diesel range are impacting lube oil range results.
- O Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
- P The RPD of the detected concentrations between the two columns is greater than 40.
- Q Surrogate recovery is outside of the control limits.
- S Surrogate recovery data is not available due to the necessary dilution of the sample.
- T The sample chromatogram is not similar to a typical
- U The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- U1 The practical quantitation limit is elevated due to interferences present in the sample.
- V Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
- W Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
- X Sample extract treated with a mercury cleanup procedure.
- X1- Sample extract treated with a sulfuric acid/silica gel cleanup procedure.
- Y The calibration verification for this analyte exceeded the 20% drift specified in method 8260C, and therefore the reported result should be considered an estimate. The overall performance of the calibration verification standard met the acceptance criteria of the method.

7 -

ND - Not Detected at PQL

PQL - Practical Quantitation Limit

RPD - Relative Percent Difference





Chain of Custody

Page 1 of 2

Analytical Laboratory Testing Services 14648 NE 95th Street • Redmond, WA 98052		naround Req n working day			Lá	abo	ratory	Num	ber:	()2	- 1	2 4	9									
Phone: (425) 883-3881 · www.onsite-env.com Company: Aspect Consulting Project Number: 1706520-000 Project Name: Pacific Right Bank Project Manager: Brace Carpenter Sampled by: T4F /AAF Lab ID Sample Identification			1 Day 3 Days ys)	Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX	NWTPH-Gx NWTPH-bx (☐ Acid / SG Clean-up)	Volatiles 8260C	EDB EPA 8011 (Waters Only)	Semivolatiles 8270D/SIM with low-level PAHs)	PAHS 8270D/SIM (low-level) C PAHS		Organochlorine Pesticides 8081B	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	TCLP Metals	HEM (oil and grease) 1664A				% Moisture
1 B-06_2.5	2/22/19		5	2			X	5				8					•						
2 13-06_110	2122/18		5	2			8				8						(3)					(8
3 13-06-12.5	2122/16		5	2	•		X				•	X											•
4 3-09-5	2/22/18	14:05	5	2	•												•						
5 B-09_10	2/22/18	14:15	5	2																			
6 B-09-15	2/22/16	14:30.	5	2	•						•						•						•
Signature	Co	ompany				Date		Time		Cor	mmei	nts/Sp	ecial	Instr	uction	ns							
Relinquished Received Relinquished Received Relinquished Relinquished		ASPECT (OM	Cons	SULT	ive	2			40	-				-		_	6	1,8	1. 3	DB ST	(s-A)	rA /	>
Received										Dat	a Pa	ckage	: Sta	andar	rd 🗆	Le	vel III		Level	IV 🗆			
Reviewed/Date		Reviewed/Da	te							Chr	omat	ogran	ns wit	th fina	al rep	oort [Ele	ectron	ic Data	Deliver	ables (E	DDs)	1



Chain of Custody

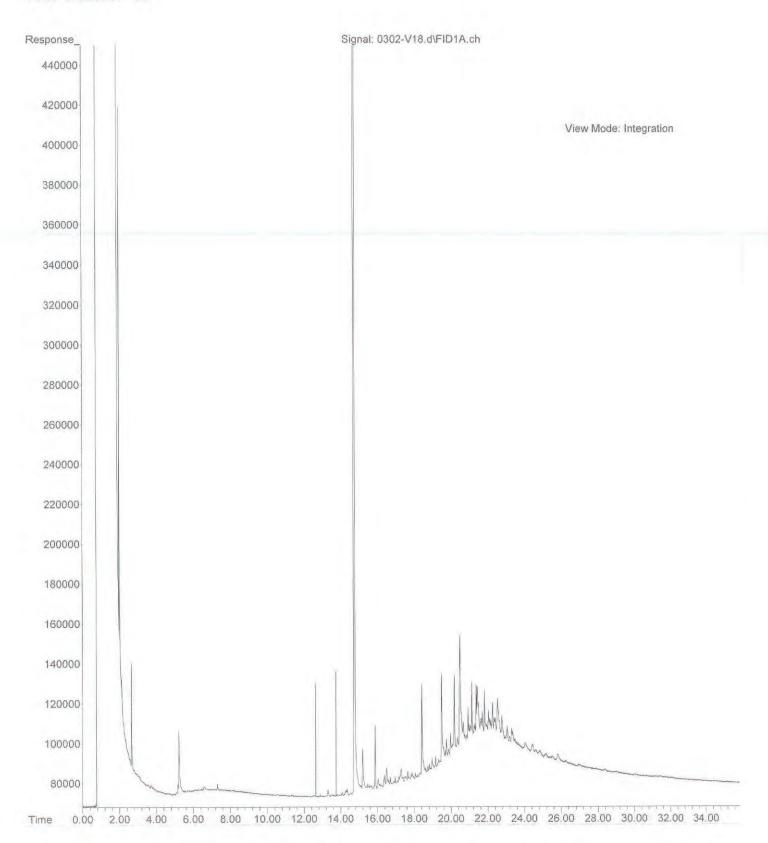
Page _ 2 of _ 2

Analytical Laboratory Testing Services 14648 NE 95th Street • Redmond, WA 98052		naround Requ n working day			La	abo	rato	ry l	Num	ber:	()2	-	2	19									
Phone: (425) 883-3881 · www.onsite-env.com Company: ASPECT CONSULTING Project Number: 1706520-000 Project Name: PACIFIC RIGHT BANK Project Manager: BRUCE CARPENTER Sampled by: AAF Lab ID Sample Identification		-	1 Day 3 Days ys)	Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX	NWTPH-Gx	Notified Boston	voiatiles 8200C Halogenated Volatiles 8260C	EDB EPA 8011 (Waters Only)	Semivolatiles 8270D/SIM (with low-level PAHs)	PAHs 8270D/SIM (low-level)	PCBs 8082A	Organochlorine Pesticides 8081B	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	TCLP Metals	HEM (oil and grease) 1664A				% Moisture
7 B-13-7.5	2/23	9:15	5	2	•							•						•						•
8 B-13_12,5	2/23	9:30	5	2																				
9 B-13_15	2/23	9:36	5	2	0							•						0						•
	2/23	9:43	5	2																				
11 B-16_2.5	2/23	1:00	5	2	•							•						•						•
	2/23	1:10	5	2	9													0						
13 B-16_12,5	2/23	1:25	5	2																				
14 B-16_17.5	2/23	1:33	5	2	0			-		1		•						0				1		•
								+	+	+														H
Signature		ompany				Date	_		Time		Cor	nmer	nts/Sp	ecial	Instru	uction	ns							
Relinquished	4	SPECT C	CONSUL	TI	NG	21	123		16:0	40														- 1
Received		0	8C			21	23/1	8	16	40														
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Reviewed/Date		Reviewed/Dat	te								-		_		_							erables	(EDDs)) 🗆

File :X:\DIESELS\VIGO\DATA\V180302\0302-V18.d
Operator : JT
Acquired : 2 Mar 2018 19:17 using AcqMethod V171020F.M
Instrument : Vigo

Sample Name: 03-249-01

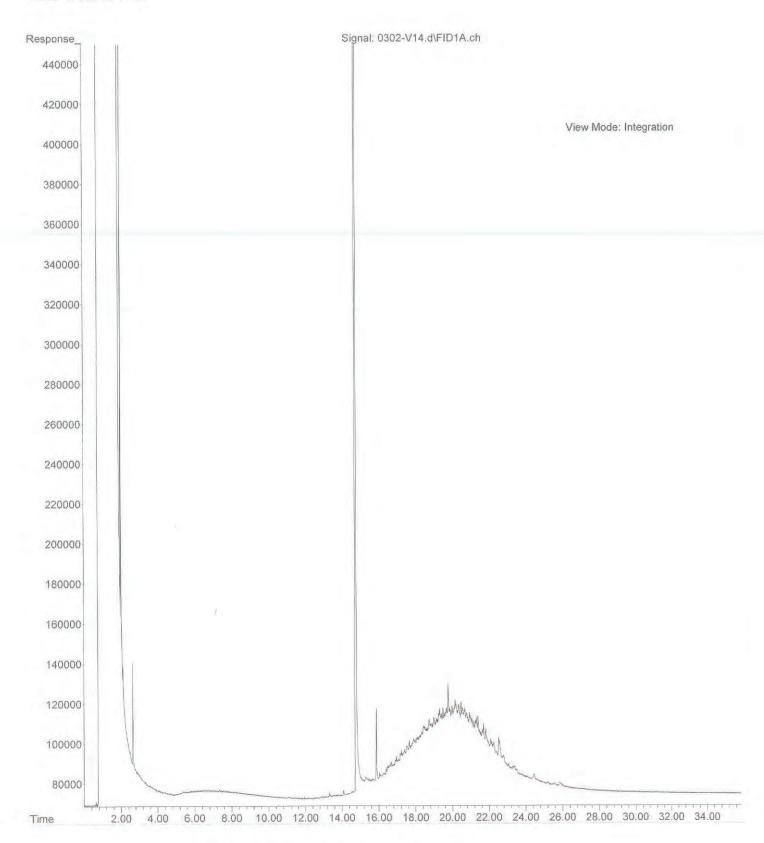
Misc Info : Vial Number: 18



File :X:\DIESELS\VIGO\DATA\V180302\0302-V14.d
Operator : JT
Acquired : 2 Mar 2018 16:37 using AcqMethod V171020F.M

Instrument: Vigo Sample Name: 02-249-03

Misc Info : Vial Number: 14





14648 NE 95th Street, Redmond, WA 98052 • (425) 883-3881

March 16, 2018

Bruce Carpenter Herrera Environmental Consultants, Inc. 2200 6th Avenue, Suite 1100 Seattle, WA 98121

Re: Analytical Data for Project 17-06520-000

Laboratory Reference No. 1802-234

Dear Bruce:

Enclosed are the analytical results and associated quality control data for samples submitted on February 22, 2018.

The standard policy of OnSite Environmental, Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

David Baumeister Project Manager

Enclosures

Date of Report: March 16, 2018 Samples Submitted: February 22, 2018 Laboratory Reference: 1802-234

Project: 17-06520-000

Case Narrative

Samples were collected on February 21, 2018 and received by the laboratory on February 22, 2018. They were maintained at the laboratory at a temperature of 2°C to 6°C.

Please note that any and all soil sample results are reported on a dry-weight basis, unless otherwise noted below.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

NWTPH-HCID

Matrix: Soil

Units: mg/kg (ppm)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP12-2			-		
Laboratory ID:	02-234-02					
Gasoline Range Organics	ND	25	NWTPH-HCID	2-23-18	2-23-18	
Diesel Range Organics	ND	65	NWTPH-HCID	2-23-18	2-23-18	U1
Lube Oil	Detected	120	NWTPH-HCID	2-23-18	2-23-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	100	50-150				
Client ID:	PP12-7					
Laboratory ID:	02-234-03					
Gasoline Range Organics	ND	29	NWTPH-HCID	2-23-18	2-23-18	
Diesel Range Organics	ND	120	NWTPH-HCID	2-23-18	2-23-18	U1
Lube Oil	Detected	150	NWTPH-HCID	2-23-18	2-23-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	81	50-150				
Client ID:	PP13-10					
Laboratory ID:	02-234-08					
Gasoline Range Organics	ND	25	NWTPH-HCID	2-23-18	2-23-18	
Diesel Range Organics	ND	62	NWTPH-HCID	2-23-18	2-23-18	
Lube Oil Range Organics	ND	120	NWTPH-HCID	2-23-18	2-23-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	89	50-150				
, ,						
Client ID:	PP15-6					
Laboratory ID:	02-234-11					
Gasoline Range Organics	ND	31	NWTPH-HCID	2-23-18	2-23-18	
Diesel Range Organics	Detected	76	NWTPH-HCID	2-23-18	2-23-18	
Lube Oil Range Organics	Detected	150	NWTPH-HCID	2-23-18	2-23-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	113	50-150				
Client ID:	PP15-12					
Laboratory ID:	02-234-12					
Gasoline Range Organics	ND	24	NWTPH-HCID	2-23-18	2-23-18	
Diesel Range Organics	ND	59	NWTPH-HCID	2-23-18	2-23-18	
Lube Oil Range Organics	ND	120	NWTPH-HCID	2-23-18	2-23-18	
Surrogate:	Percent Recovery	Control Limits		2 20 10	2 23 10	
o-Terphenyl	89	50-150				
C. Orphony:	55	00 100				

NWTPH-HCID

Matrix: Soil

• • •		201		Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP19-7					
Laboratory ID:	02-234-15					
Gasoline Range Organics	ND	25	NWTPH-HCID	2-23-18	2-23-18	
Diesel Range Organics	Detected	63	NWTPH-HCID	2-23-18	2-23-18	
Lube Oil Range Organics	ND	130	NWTPH-HCID	2-23-18	2-23-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	99	50-150				
Client ID:	PP19-10					
Laboratory ID:	02-234-16					
Gasoline Range Organics	ND	24	NWTPH-HCID	2-23-18	2-23-18	
Diesel Range Organics	ND	59	NWTPH-HCID	2-23-18	2-23-18	
Lube Oil Range Organics	ND	120	NWTPH-HCID	2-23-18	2-23-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	96	50-150				
Client ID:	PP19-15					
Laboratory ID:	02-234-17					
Gasoline Range Organics	ND	24	NWTPH-HCID	2-23-18	2-23-18	
Diesel Range Organics	ND	60	NWTPH-HCID	2-23-18	2-23-18	
Lube Oil Range Organics	ND	120	NWTPH-HCID	2-23-18	2-23-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	91	50-150				
, ,						
Client ID:	PP22-4					
Laboratory ID:	02-234-18					
Gasoline Range Organics	ND	26	NWTPH-HCID	2-23-18	2-23-18	
Diesel Range Organics	ND	64	NWTPH-HCID	2-23-18	2-23-18	
Lube Oil Range Organics	ND	130	NWTPH-HCID	2-23-18	2-23-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	95	50-150				
Client ID:	PP22-9					
Laboratory ID:	02-234-19					
Gasoline Range Organics	ND	25	NWTPH-HCID	2-23-18	2-23-18	
Diesel Range Organics	ND	61	NWTPH-HCID	2-23-18	2-23-18	
Lube Oil Range Organics	ND ND	120	NWTPH-HCID	2-23-18	2-23-18	
Surrogate:	Percent Recovery	Control Limits	1444 II II-IIOID	2 20-10	Z ZU-10	
o-Terphenyl	94	50-150				
o respirestlys	J *1	JU-1JU				

NWTPH-HCID

Matrix: Soil

Units: mg/kg (ppm)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP22-13			-	-	
Laboratory ID:	02-234-20					
Gasoline Range Organics	ND	23	NWTPH-HCID	2-23-18	2-23-18	
Diesel Range Organics	ND	58	NWTPH-HCID	2-23-18	2-23-18	
Lube Oil Range Organics	ND	120	NWTPH-HCID	2-23-18	2-23-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	95	50-150				
Client ID:	PP25-7					
Laboratory ID:	02-234-23					
Gasoline Range Organics	ND	27	NWTPH-HCID	2-23-18	2-23-18	
Diesel Range Organics	ND	69	NWTPH-HCID	2-23-18	2-23-18	
Lube Oil Range Organics	ND	140	NWTPH-HCID	2-23-18	2-23-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	117	50-150				
Client ID:	PP25-13					
Laboratory ID:	02-234-24					
Gasoline Range Organics	Detected	34	NWTPH-HCID	2-23-18	2-23-18	
Diesel Range Organics	Detected	85	NWTPH-HCID	2-23-18	2-23-18	
Lube Oil Range Organics	Detected	170	NWTPH-HCID	2-23-18	2-23-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	108	50-150				
o respirent	700	00 100				
Client ID:	PP26-11					
Laboratory ID:	02-234-28					
Gasoline Range Organics	ND	23	NWTPH-HCID	2-23-18	2-23-18	
Diesel Range Organics	ND	58	NWTPH-HCID	2-23-18	2-23-18	
Lube Oil Range Organics	ND	120	NWTPH-HCID	2-23-18	2-23-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	92	50-150				
, ,						
Client ID:	PP26-17					
Laboratory ID:	02-234-29					
Gasoline Range Organics	ND	31	NWTPH-HCID	2-23-18	2-23-18	
Diesel Range Organics	ND	77	NWTPH-HCID	2-23-18	2-23-18	
Lube Oil Range Organics	ND	150	NWTPH-HCID	2-23-18	2-23-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	87	50-150				
, ,						

NWTPH-HCID

Matrix: Soil

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP34-8					111.90
Laboratory ID:	02-234-31					
Gasoline Range Organics	ND	24	NWTPH-HCID	2-23-18	2-23-18	
Diesel Range Organics	ND	61	NWTPH-HCID	2-23-18	2-23-18	
Lube Oil Range Organics	ND	120	NWTPH-HCID	2-23-18	2-23-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	102	50-150				

NWTPH-HCID QUALITY CONTROL

Matrix: Soil

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0223S3					
Gasoline Range Organics	ND	20	NWTPH-HCID	2-23-18	2-23-18	
Diesel Range Organics	ND	50	NWTPH-HCID	2-23-18	2-23-18	
Lube Oil Range Organics	ND	100	NWTPH-HCID	2-23-18	2-23-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	95	50-150				

Date of Report: March 16, 2018 Samples Submitted: February 22, 2018 Laboratory Reference: 1802-234

Project: 17-06520-000

NWTPH-Gx/BTEX

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP25-13					
Laboratory ID:	02-234-24					
Benzene	ND	0.026	EPA 8021B	2-28-18	3-1-18	
Toluene	ND	0.13	EPA 8021B	2-28-18	3-1-18	
Ethyl Benzene	ND	0.13	EPA 8021B	2-28-18	3-1-18	
m,p-Xylene	ND	0.13	EPA 8021B	2-28-18	3-1-18	
o-Xylene	ND	0.13	EPA 8021B	2-28-18	3-1-18	
Gasoline	400	13	NWTPH-Gx	2-28-18	3-1-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	79	66-130				
Client ID:	PP25-17					
Laboratory ID:	02-234-25					
Benzene	ND	0.030	EPA 8021B	2-28-18	3-1-18	
Toluene	ND	0.15	EPA 8021B	2-28-18	3-1-18	
Ethyl Benzene	ND	0.15	EPA 8021B	2-28-18	3-1-18	
m,p-Xylene	ND	0.15	EPA 8021B	2-28-18	3-1-18	
o-Xylene	ND	0.15	EPA 8021B	2-28-18	3-1-18	
Gasoline	ND	15	NWTPH-Gx	2-28-18	3-1-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	82	66-130				

Date of Report: March 16, 2018 Samples Submitted: February 22, 2018 Laboratory Reference: 1802-234

Project: 17-06520-000

NWTPH-Gx/BTEX QUALITY CONTROL

Matrix: Soil

Units: mg/kg (ppm)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0228S2					
Benzene	ND	0.020	EPA 8021B	2-28-18	2-28-18	
Toluene	ND	0.050	EPA 8021B	2-28-18	2-28-18	
Ethyl Benzene	ND	0.050	EPA 8021B	2-28-18	2-28-18	
m,p-Xylene	ND	0.050	EPA 8021B	2-28-18	2-28-18	
o-Xylene	ND	0.050	EPA 8021B	2-28-18	2-28-18	
Gasoline	ND	5.0	NWTPH-Gx	2-28-18	2-28-18	
_			·		·	·

Percent Recovery Control Limits Surrogate: Fluorobenzene 86 66-130

					Source	Per	cent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Rec	overy	Limits	RPD	Limit	Flags
DUPLICATE											
Laboratory ID:	02-26	65-02									
	ORIG	DUP									
Benzene	ND	ND	NA	NA		١	1A	NA	NA	30	
Toluene	ND	ND	NA	NA		N	١A	NA	NA	30	
Ethyl Benzene	ND	ND	NA	NA		١	1A	NA	NA	30	
m,p-Xylene	ND	ND	NA	NA		١	1A	NA	NA	30	
o-Xylene	ND	ND	NA	NA		١	1A	NA	NA	30	
Gasoline	ND	ND	NA	NA		١	IA.	NA	NA	30	
Surrogate:											
Fluorobenzene						103	96	66-130			
SPIKE BLANKS											
Laboratory ID:	SB02	28S1									
	SB	SBD	SB	SBD		SB	SBD				
Benzene	0.859	0.879	1.00	1.00		86	88	70-120	2	11	
Toluene	0.871	0.894	1.00	1.00		87	89	73-121	3	14	
Ethyl Benzene	0.881	0.902	1.00	1.00		88	90	74-121	2	11	
m,p-Xylene	0.881	0.903	1.00	1.00		88	90	75-124	2	13	
o-Xylene	0.899	0.924	1.00	1.00		90	92	75-121	3	12	
Surrogate:											
Fluorobenzene						85	86	66-130			

NWTPH-Dx

Matrix: Soil

5 5 (1)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP12-2					
Laboratory ID:	02-234-02					
Diesel Range Organics	ND	31	NWTPH-Dx	2-28-18	2-28-18	
Lube Oil	69	62	NWTPH-Dx	2-28-18	2-28-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	50	50-150				
Client ID:	PP12-7					
Laboratory ID:	02-234-03					
Diesel Range Organics	ND	130	NWTPH-Dx	2-28-18	2-28-18	U1
Lube Oil	800	74	NWTPH-Dx	2-28-18	2-28-18	01
Surrogate:	Percent Recovery	Control Limits	TWO THE DA	2 20 10	2 20 10	
o-Terphenyl	101	50-150				
Client ID:	PP12-10					
Laboratory ID:	02-234-04					
Diesel Range Organics	ND	32	NWTPH-Dx	2-28-18	2-28-18	
Lube Oil	82	64	NWTPH-Dx	2-28-18	2-28-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	87	50-150				
Client ID:	PP15-6					
Laboratory ID:	02-234-11					
Diesel Range Organics	81	38	NWTPH-Dx	2-28-18	2-28-18	
Lube Oil Range Organics	550	76	NWTPH-Dx	2-28-18	2-28-18	
Surrogate:	Percent Recovery	Control Limits	INVVIIII-DX	2-20-10	2-20-10	
o-Terphenyl	97	50-150				
o respirativi	37	00 100				
Client ID:	PP19-7					
Laboratory ID:	02-234-15					
Diesel Range Organics	400	32	NWTPH-Dx	2-28-18	2-28-18	
Lube Oil Range Organics	370	63	NWTPH-Dx	2-28-18	2-28-18	N1
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	101	50-150				
Olion (ID)	DD05 40					
Client ID:	PP25-13					
Laboratory ID:	02-234-24	40	NIM/TOUR	0.00.10	0.00.10	
Diesel Range Organics	1400	43	NWTPH-Dx	2-28-18	2-28-18	
Lube Oil Range Organics	1200	85	NWTPH-Dx	2-28-18	2-28-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	85	50-150				

NWTPH-Dx

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP25-17					
Laboratory ID:	02-234-25					
Diesel Range Organics	49	47	NWTPH-Dx	2-28-18	2-28-18	_
Lube Oil Range Organics	130	93	NWTPH-Dx	2-28-18	2-28-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	71	50-150				

NWTPH-Dx QUALITY CONTROL

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0228S1					
Diesel Range Organics	ND	25	NWTPH-Dx	2-28-18	2-28-18	
Lube Oil Range Organics	ND	50	NWTPH-Dx	2-28-18	2-28-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	109	50-150				

Analyte	Res	sult	Spike	Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
DUPLICATE										
Laboratory ID:	02-23	34-04								
	ORIG	DUP								
Diesel Range	ND	ND	NA	NA		NA	NA	NA	NA	
Lube Oil	63.8	61.5	NA	NA		NA	NA	4	NA	
Surrogate:										
o-Terphenyl						87 88	50-150			

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP12-2					
Laboratory ID:	02-234-02					
Benzo[a]anthracene	0.0084	0.0082	EPA 8270D/SIM	2-26-18	2-26-18	
Chrysene	0.014	0.0082	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[b]fluoranthene	0.016	0.0082	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo(j,k)fluoranthene	ND	0.0082	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[a]pyrene	0.015	0.0082	EPA 8270D/SIM	2-26-18	2-26-18	
Indeno(1,2,3-c,d)pyrene	0.012	0.0082	EPA 8270D/SIM	2-26-18	2-26-18	
Dibenz[a,h]anthracene	ND	0.0082	EPA 8270D/SIM	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	74	32 - 115				
Pyrene-d10	80	35 - 129				
Terphenyl-d14	77	33 - 114				

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP12-7					
Laboratory ID:	02-234-03					
Benzo[a]anthracene	0.010	0.0098	EPA 8270D/SIM	2-26-18	2-26-18	
Chrysene	0.013	0.0098	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[b]fluoranthene	0.014	0.0098	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo(j,k)fluoranthene	ND	0.0098	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[a]pyrene	0.011	0.0098	EPA 8270D/SIM	2-26-18	2-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0098	EPA 8270D/SIM	2-26-18	2-26-18	
Dibenz[a,h]anthracene	ND	0.0098	EPA 8270D/SIM	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	85	32 - 115				
Pyrene-d10	83	35 - 129				
Terphenyl-d14	81	33 - 114				

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP13-10					
Laboratory ID:	02-234-08					
Benzo[a]anthracene	ND	0.0083	EPA 8270D/SIM	2-26-18	2-26-18	
Chrysene	ND	0.0083	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[b]fluoranthene	ND	0.0083	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo(j,k)fluoranthene	ND	0.0083	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[a]pyrene	ND	0.0083	EPA 8270D/SIM	2-26-18	2-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0083	EPA 8270D/SIM	2-26-18	2-26-18	
Dibenz[a,h]anthracene	ND	0.0083	EPA 8270D/SIM	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	82	32 - 115				
Pyrene-d10	87	35 - 129				
Terphenyl-d14	85	33 - 114				

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP15-6					
Laboratory ID:	02-234-11					
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	2-26-18	2-26-18	
Chrysene	ND	0.010	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	2-26-18	2-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	2-26-18	2-26-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	83	32 - 115				
Pyrene-d10	84	35 - 129				
Terphenyl-d14	83	33 - 114				

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP15-12					
Laboratory ID:	02-234-12					
Benzo[a]anthracene	ND	0.0079	EPA 8270D/SIM	2-26-18	2-26-18	
Chrysene	ND	0.0079	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[b]fluoranthene	ND	0.0079	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo(j,k)fluoranthene	ND	0.0079	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[a]pyrene	ND	0.0079	EPA 8270D/SIM	2-26-18	2-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0079	EPA 8270D/SIM	2-26-18	2-26-18	
Dibenz[a,h]anthracene	ND	0.0079	EPA 8270D/SIM	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	91	32 - 115				
Pyrene-d10	92	35 - 129				
Terphenyl-d14	92	33 - 114				

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP19-7					
Laboratory ID:	02-234-15					
Benzo[a]anthracene	ND	0.0084	EPA 8270D/SIM	2-26-18	2-26-18	
Chrysene	ND	0.0084	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[b]fluoranthene	ND	0.0084	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo(j,k)fluoranthene	ND	0.0084	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[a]pyrene	ND	0.0084	EPA 8270D/SIM	2-26-18	2-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0084	EPA 8270D/SIM	2-26-18	2-26-18	
Dibenz[a,h]anthracene	ND	0.0084	EPA 8270D/SIM	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	82	32 - 115				
Pyrene-d10	98	35 - 129				
Terphenyl-d14	88	33 - 114				

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP19-10					
Laboratory ID:	02-234-16					
Benzo[a]anthracene	ND	0.0078	EPA 8270D/SIM	2-26-18	2-26-18	
Chrysene	ND	0.0078	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[b]fluoranthene	ND	0.0078	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo(j,k)fluoranthene	ND	0.0078	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[a]pyrene	ND	0.0078	EPA 8270D/SIM	2-26-18	2-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0078	EPA 8270D/SIM	2-26-18	2-26-18	
Dibenz[a,h]anthracene	ND	0.0078	EPA 8270D/SIM	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	77	32 - 115				
Pyrene-d10	88	35 - 129				
Terphenyl-d14	88	33 - 114				

Laboratory Reference: 1802-234 Project: 17-06520-000

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP19-15					
Laboratory ID:	02-234-17					
Benzo[a]anthracene	ND	0.0080	EPA 8270D/SIM	2-26-18	2-26-18	
Chrysene	ND	0.0080	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[b]fluoranthene	ND	0.0080	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo(j,k)fluoranthene	ND	0.0080	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[a]pyrene	ND	0.0080	EPA 8270D/SIM	2-26-18	2-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0080	EPA 8270D/SIM	2-26-18	2-26-18	
Dibenz[a,h]anthracene	ND	0.0080	EPA 8270D/SIM	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	81	32 - 115				
Pyrene-d10	93	35 - 129				
Terphenyl-d14	89	33 - 114				

Laboratory Reference: 1802-234 Project: 17-06520-000

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP22-4					
Laboratory ID:	02-234-18					
Benzo[a]anthracene	0.056	0.0086	EPA 8270D/SIM	2-26-18	2-26-18	
Chrysene	0.086	0.0086	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[b]fluoranthene	0.088	0.0086	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo(j,k)fluoranthene	0.029	0.0086	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[a]pyrene	0.064	0.0086	EPA 8270D/SIM	2-26-18	2-26-18	
Indeno(1,2,3-c,d)pyrene	0.051	0.0086	EPA 8270D/SIM	2-26-18	2-26-18	
Dibenz[a,h]anthracene	0.017	0.0086	EPA 8270D/SIM	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	80	32 - 115				
Pyrene-d10	91	35 - 129				
Terphenyl-d14	78	33 - 114				

Laboratory Reference: 1802-234 Project: 17-06520-000

PAHs EPA 8270D/SIM

0 0				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP22-9					
Laboratory ID:	02-234-19					
Benzo[a]anthracene	ND	0.0082	EPA 8270D/SIM	2-26-18	2-26-18	
Chrysene	ND	0.0082	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[b]fluoranthene	ND	0.0082	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo(j,k)fluoranthene	ND	0.0082	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[a]pyrene	ND	0.0082	EPA 8270D/SIM	2-26-18	2-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0082	EPA 8270D/SIM	2-26-18	2-26-18	
Dibenz[a,h]anthracene	ND	0.0082	EPA 8270D/SIM	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	85	32 - 115				
Pyrene-d10	84	35 - 129				
Terphenyl-d14	86	33 - 114				

Laboratory Reference: 1802-234 Project: 17-06520-000

PAHs EPA 8270D/SIM

0 0				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP22-13					·
Laboratory ID:	02-234-20					
Benzo[a]anthracene	ND	0.0077	EPA 8270D/SIM	2-26-18	2-26-18	
Chrysene	ND	0.0077	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[b]fluoranthene	ND	0.0077	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo(j,k)fluoranthene	ND	0.0077	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[a]pyrene	ND	0.0077	EPA 8270D/SIM	2-26-18	2-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0077	EPA 8270D/SIM	2-26-18	2-26-18	
Dibenz[a,h]anthracene	ND	0.0077	EPA 8270D/SIM	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	80	32 - 115				
Pyrene-d10	86	35 - 129				
Terphenyl-d14	89	33 - 114				

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP25-7					
Laboratory ID:	02-234-23					
Benzo[a]anthracene	ND	0.0091	EPA 8270D/SIM	2-26-18	2-26-18	
Chrysene	ND	0.0091	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[b]fluoranthene	ND	0.0091	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo(j,k)fluoranthene	ND	0.0091	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[a]pyrene	ND	0.0091	EPA 8270D/SIM	2-26-18	2-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0091	EPA 8270D/SIM	2-26-18	2-26-18	
Dibenz[a,h]anthracene	ND	0.0091	EPA 8270D/SIM	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	91	32 - 115				
Pyrene-d10	86	35 - 129				
Terphenyl-d14	86	33 - 114				

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP25-13					
Laboratory ID:	02-234-24					
Benzo[a]anthracene	ND	0.011	EPA 8270D/SIM	2-26-18	2-26-18	
Chrysene	0.012	0.011	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[b]fluoranthene	ND	0.011	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo(j,k)fluoranthene	ND	0.011	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[a]pyrene	ND	0.011	EPA 8270D/SIM	2-26-18	2-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.011	EPA 8270D/SIM	2-26-18	2-26-18	
Dibenz[a,h]anthracene	ND	0.011	EPA 8270D/SIM	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	77	32 - 115				
Pyrene-d10	87	35 - 129				
Terphenyl-d14	81	33 - 114				

Laboratory Reference: 1802-234 Project: 17-06520-000

PAHs EPA 8270D/SIM

3 3				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP26-11					·
Laboratory ID:	02-234-28					
Benzo[a]anthracene	ND	0.0077	EPA 8270D/SIM	2-26-18	2-26-18	
Chrysene	ND	0.0077	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[b]fluoranthene	ND	0.0077	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo(j,k)fluoranthene	ND	0.0077	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[a]pyrene	ND	0.0077	EPA 8270D/SIM	2-26-18	2-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0077	EPA 8270D/SIM	2-26-18	2-26-18	
Dibenz[a,h]anthracene	ND	0.0077	EPA 8270D/SIM	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	86	32 - 115				
Pyrene-d10	87	35 - 129				
Terphenyl-d14	88	33 - 114				

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP26-17					
Laboratory ID:	02-234-29					
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	2-26-18	2-26-18	
Chrysene	ND	0.010	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	2-26-18	2-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	2-26-18	2-26-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	88	32 - 115				
Pyrene-d10	82	35 - 129				
Terphenyl-d14	84	33 - 114				

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP34-8					
Laboratory ID:	02-234-31					
Benzo[a]anthracene	ND	0.0081	EPA 8270D/SIM	2-26-18	2-26-18	
Chrysene	ND	0.0081	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[b]fluoranthene	ND	0.0081	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo(j,k)fluoranthene	ND	0.0081	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[a]pyrene	ND	0.0081	EPA 8270D/SIM	2-26-18	2-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0081	EPA 8270D/SIM	2-26-18	2-26-18	
Dibenz[a,h]anthracene	ND	0.0081	EPA 8270D/SIM	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	80	32 - 115				
Pyrene-d10	88	35 - 129				
Terphenyl-d14	87	33 - 114				

Laboratory Reference: 1802-234

Project: 17-06520-000

PAHS EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
						_
Laboratory ID:	MB0226S1					
Benzo[a]anthracene	ND	0.0067	EPA 8270D/SIM	2-26-18	2-26-18	_
Chrysene	ND	0.0067	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270D/SIM	2-26-18	2-26-18	
Benzo[a]pyrene	ND	0.0067	EPA 8270D/SIM	2-26-18	2-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0067	EPA 8270D/SIM	2-26-18	2-26-18	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270D/SIM	2-26-18	2-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	96	32 - 115				
Pyrene-d10	91	35 - 129				
Terphenyl-d14	93	33 - 114				

Laboratory Reference: 1802-234 Project: 17-06520-000

PAHs EPA 8270D/SIM MS/MSD QUALITY CONTROL

					Source	Per	cent	Recovery		RPD	
Analyte	Re	sult	Spike	Level	Result	Rec	overy	Limits	RPD	Limit	Flags
MATRIX SPIKES											
Laboratory ID:	02-2	34-19									
	MS	MSD	MS	MSD		MS	MSD				
Benzo[a]anthracene	0.0813	0.0814	0.0833	0.0833	ND	98	98	27 - 143	0	23	
Chrysene	0.0831	0.0829	0.0833	0.0833	ND	100	100	22 - 130	0	24	
Benzo[b]fluoranthene	0.0727	0.0737	0.0833	0.0833	ND	87	88	15 - 141	1	26	
Benzo(j,k)fluoranthene	0.0856	0.0857	0.0833	0.0833	ND	103	103	42 - 112	0	24	
Benzo[a]pyrene	0.0761	0.0761	0.0833	0.0833	ND	91	91	33 - 126	0	26	
Indeno(1,2,3-c,d)pyrene	0.0917	0.0912	0.0833	0.0833	ND	110	109	30 - 125	1	25	
Dibenz[a,h]anthracene	0.0747	0.0756	0.0833	0.0833	ND	90	91	31 - 124	1	22	
Surrogate:											
2-Fluorobiphenyl						93	90	32 - 115			
Pyrene-d10						88	88	35 - 129			
Terphenyl-d14						90	89	33 - 114			

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP12-10					
Laboratory ID:	02-234-04					
Benzo[a]anthracene	ND	0.0086	EPA 8270D/SIM	3-1-18	3-1-18	
Chrysene	ND	0.0086	EPA 8270D/SIM	3-1-18	3-1-18	
Benzo[b]fluoranthene	ND	0.0086	EPA 8270D/SIM	3-1-18	3-1-18	
Benzo(j,k)fluoranthene	ND	0.0086	EPA 8270D/SIM	3-1-18	3-1-18	
Benzo[a]pyrene	ND	0.0086	EPA 8270D/SIM	3-1-18	3-1-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0086	EPA 8270D/SIM	3-1-18	3-1-18	
Dibenz[a,h]anthracene	ND	0.0086	EPA 8270D/SIM	3-1-18	3-1-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	86	32 - 115				
Pyrene-d10	97	35 - 129				
Terphenyl-d14	103	33 - 114				

Date of Report: March 16, 2018 Samples Submitted: February 22, 2018 Laboratory Reference: 1802-234

Project: 17-06520-000

PAHs EPA 8270D/SIM

0 0				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP25-17					
Laboratory ID:	02-234-25					
Benzo[a]anthracene	ND	0.012	EPA 8270D/SIM	3-1-18	3-4-18	
Chrysene	ND	0.012	EPA 8270D/SIM	3-1-18	3-4-18	
Benzo[b]fluoranthene	ND	0.012	EPA 8270D/SIM	3-1-18	3-4-18	
Benzo(j,k)fluoranthene	ND	0.012	EPA 8270D/SIM	3-1-18	3-4-18	
Benzo[a]pyrene	ND	0.012	EPA 8270D/SIM	3-1-18	3-4-18	
Indeno(1,2,3-c,d)pyrene	ND	0.012	EPA 8270D/SIM	3-1-18	3-4-18	
Dibenz[a,h]anthracene	ND	0.012	EPA 8270D/SIM	3-1-18	3-4-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	77	32 - 115				
Pyrene-d10	81	35 - 129				
Terphenyl-d14	87	33 - 114				

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP34-15					
Laboratory ID:	02-234-33					
Benzo[a]anthracene	ND	0.0093	EPA 8270D/SIM	3-1-18	3-4-18	
Chrysene	ND	0.0093	EPA 8270D/SIM	3-1-18	3-4-18	
Benzo[b]fluoranthene	ND	0.0093	EPA 8270D/SIM	3-1-18	3-4-18	
Benzo(j,k)fluoranthene	ND	0.0093	EPA 8270D/SIM	3-1-18	3-4-18	
Benzo[a]pyrene	ND	0.0093	EPA 8270D/SIM	3-1-18	3-4-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0093	EPA 8270D/SIM	3-1-18	3-4-18	
Dibenz[a,h]anthracene	ND	0.0093	EPA 8270D/SIM	3-1-18	3-4-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	72	32 - 115				
Pyrene-d10	74	35 - 129				
Terphenyl-d14	80	33 - 114				

Laboratory Reference: 1802-234

Project: 17-06520-000

PAHS EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0301S2					
Benzo[a]anthracene	ND	0.0067	EPA 8270D/SIM	3-1-18	3-1-18	
Chrysene	ND	0.0067	EPA 8270D/SIM	3-1-18	3-1-18	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270D/SIM	3-1-18	3-1-18	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270D/SIM	3-1-18	3-1-18	
Benzo[a]pyrene	ND	0.0067	EPA 8270D/SIM	3-1-18	3-1-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0067	EPA 8270D/SIM	3-1-18	3-1-18	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270D/SIM	3-1-18	3-1-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	92	32 - 115				
Pyrene-d10	97	35 - 129				
Terphenyl-d14	104	33 - 114				

Laboratory Reference: 1802-234 Project: 17-06520-000

PAHs EPA 8270D/SIM MS/MSD QUALITY CONTROL

					Source	Per	cent	Recovery		RPD	
Analyte	Re	sult	Spike	Level	Result	Rec	overy	Limits	RPD	Limit	Flags
MATRIX SPIKES											
Laboratory ID:	02-2	34-04									
	MS	MSD	MS	MSD		MS	MSD				
Benzo[a]anthracene	0.172	0.183	0.167	0.167	ND	103	110	27 - 143	6	23	
Chrysene	0.161	0.170	0.167	0.167	ND	96	102	22 - 130	5	24	
Benzo[b]fluoranthene	0.160	0.169	0.167	0.167	ND	96	101	15 - 141	5	26	
Benzo(j,k)fluoranthene	0.161	0.168	0.167	0.167	ND	96	101	42 - 112	4	24	
Benzo[a]pyrene	0.161	0.168	0.167	0.167	ND	96	101	33 - 126	4	26	
Indeno(1,2,3-c,d)pyrene	0.154	0.157	0.167	0.167	ND	92	94	30 - 125	2	25	
Dibenz[a,h]anthracene	0.156	0.159	0.167	0.167	ND	93	95	31 - 124	2	22	
Surrogate:											
2-Fluorobiphenyl						81	88	32 - 115			
Pyrene-d10						93	95	35 - 129			
Terphenyl-d14						97	97	33 - 114			

PCBs EPA 8082A

Matrix: Soil

onits. Hig/kg (ppin)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP12-2					
Laboratory ID:	02-234-02					
Aroclor 1016	ND	0.062	EPA 8082A	3-1-18	3-1-18	
Aroclor 1221	ND	0.062	EPA 8082A	3-1-18	3-1-18	
Aroclor 1232	ND	0.062	EPA 8082A	3-1-18	3-1-18	
Aroclor 1242	ND	0.062	EPA 8082A	3-1-18	3-1-18	
Aroclor 1248	ND	0.062	EPA 8082A	3-1-18	3-1-18	
Aroclor 1254	ND	0.062	EPA 8082A	3-1-18	3-1-18	
Aroclor 1260	ND	0.062	EPA 8082A	3-1-18	3-1-18	
Surrogate:	Percent Recovery	Control Limits				
DCB	80	40-134				
Client ID:	PP12-7					
Laboratory ID:	02-234-03					
Aroclor 1016	ND	0.073	EPA 8082A	3-1-18	3-1-18	
Aroclor 1221	ND	0.073	EPA 8082A	3-1-18	3-1-18	
Aroclor 1232	ND	0.073	EPA 8082A	3-1-18	3-1-18	
Aroclor 1242	ND	0.073	EPA 8082A	3-1-18	3-1-18	
Aroclor 1248	ND	0.073	EPA 8082A	3-1-18	3-1-18	
Aroclor 1254	0.29	0.073	EPA 8082A	3-1-18	3-1-18	
Aroclor 1260	ND	0.073	EPA 8082A	3-1-18	3-1-18	
Surrogate:	Percent Recovery	Control Limits				
DCB	72	40-134				
Client ID:	PP12-10					
Laboratory ID:	02-234-04					
Aroclor 1016	ND	0.064	EPA 8082A	3-1-18	3-1-18	
Aroclor 1221	ND	0.064	EPA 8082A	3-1-18	3-1-18	
Aroclor 1232	ND	0.064	EPA 8082A	3-1-18	3-1-18	
Aroclor 1242	ND	0.064	EPA 8082A	3-1-18	3-1-18	
Aroclor 1248	ND	0.064	EPA 8082A	3-1-18	3-1-18	
Aroclor 1254	ND	0.064	EPA 8082A	3-1-18	3-1-18	
Aroclor 1260	ND	0.064	EPA 8082A	3-1-18	3-1-18	
Surrogate:	Percent Recovery	Control Limits				
DCB	77	40-134				

PCBs EPA 8082A

Matrix: Soil

Units: mg/Kg (ppm)

Analysis	Result	PQL	Method	Date Prepared	Date Analyzed	Elogo
Analyte Client ID:	PP15-6	PQL	Wethod	Prepared	Analyzeu	Flags
Laboratory ID:	02-234-11	0.070	EDA 0000A	0.4.40	0.4.40	
Aroclor 1016	ND	0.076	EPA 8082A	3-1-18	3-1-18	
Aroclor 1221	ND	0.076	EPA 8082A	3-1-18	3-1-18	
Aroclor 1232	ND	0.076	EPA 8082A	3-1-18	3-1-18	
Aroclor 1242	ND	0.076	EPA 8082A	3-1-18	3-1-18	
Aroclor 1248	ND	0.076	EPA 8082A	3-1-18	3-1-18	
Aroclor 1254	ND	0.076	EPA 8082A	3-1-18	3-1-18	
Aroclor 1260	ND 	0.076	EPA 8082A	3-1-18	3-1-18	
Surrogate:	Percent Recovery	Control Limits				
DCB	77	40-134				
Client ID:	PP25-13					
Laboratory ID:	02-234-24					
Aroclor 1016	ND	0.085	EPA 8082A	3-1-18	3-7-18	Χ
Aroclor 1221	ND	0.085	EPA 8082A	3-1-18	3-7-18	Χ
Aroclor 1232	ND	0.085	EPA 8082A	3-1-18	3-7-18	Χ
Aroclor 1242	0.58	0.085	EPA 8082A	3-1-18	3-7-18	Χ
Aroclor 1248	ND	0.085	EPA 8082A	3-1-18	3-7-18	Χ
Aroclor 1254	ND	0.085	EPA 8082A	3-1-18	3-7-18	Χ
Aroclor 1260	0.78	0.085	EPA 8082A	3-1-18	3-7-18	Χ
Surrogate:	Percent Recovery	Control Limits				
DCB	82	40-134				
Client ID:	PP25-17					
Laboratory ID:	02-234-25					
Aroclor 1016	ND	0.093	EPA 8082A	3-1-18	3-7-18	
Aroclor 1221	ND	0.093	EPA 8082A	3-1-18	3-7-18	
Aroclor 1232	ND	0.093	EPA 8082A	3-1-18	3-7-18	
Aroclor 1242	ND	0.093	EPA 8082A	3-1-18	3-7-18	
Aroclor 1248	ND	0.093	EPA 8082A	3-1-18	3-7-18	
Aroclor 1254	ND	0.093	EPA 8082A	3-1-18	3-7-18	
Aroclor 1260	ND	0.093	EPA 8082A	3-1-18	3-7-18	
Surrogate:	Percent Recovery	Control Limits				

DCB 74 40-134

PCBs EPA 8082A METHOD BLANK QUALITY CONTROL

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0301S1					
Aroclor 1016	ND	0.050	EPA 8082A	3-1-18	3-7-18	Х
Aroclor 1221	ND	0.050	EPA 8082A	3-1-18	3-7-18	X
Aroclor 1232	ND	0.050	EPA 8082A	3-1-18	3-7-18	X
Aroclor 1242	ND	0.050	EPA 8082A	3-1-18	3-7-18	X
Aroclor 1248	ND	0.050	EPA 8082A	3-1-18	3-7-18	X
Aroclor 1254	ND	0.050	EPA 8082A	3-1-18	3-7-18	X
Aroclor 1260	ND	0.050	EPA 8082A	3-1-18	3-7-18	X
Surrogate:	Percent Recovery	Control Limits				
DCB	92	40-134				
Laboratory ID:	MB0301S1					
Aroclor 1016	ND	0.050	EPA 8082A	3-1-18	3-1-18	
Aroclor 1221	ND	0.050	EPA 8082A	3-1-18	3-1-18	
Aroclor 1232	ND	0.050	EPA 8082A	3-1-18	3-1-18	
Aroclor 1242	ND	0.050	EPA 8082A	3-1-18	3-1-18	
Aroclor 1248	ND	0.050	EPA 8082A	3-1-18	3-1-18	
Aroclor 1254	ND	0.050	EPA 8082A	3-1-18	3-1-18	
Aroclor 1260	ND	0.050	EPA 8082A	3-1-18	3-1-18	
Surrogate:	Percent Recovery	Control Limits				
DCB	84	40-134				

Analyte	Re	sult	Spike	Level	Source Result		rcent covery	Recovery Limits	RPD	RPD Limit	Flags
MATRIX SPIKES											
Laboratory ID:	02-2	30-02									
	MS	MSD	MS	MSD		MS	MSD				
Aroclor 1260	0.408	0.443	0.500	0.500	ND	82	89	34-126	8	16	
Surrogate:											
DCB						74	81	40-134			

Laboratory Reference: 1802-234 Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B

Matrix: Soil

·····9/··19 (PP····)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP12-2					
Laboratory ID:	02-234-02					
Arsenic	ND	12	EPA 6010D	2-27-18	2-27-18	
Cadmium	ND	0.62	EPA 6010D	2-27-18	2-27-18	
Chromium	15	0.62	EPA 6010D	2-27-18	2-27-18	
Lead	8.8	6.2	EPA 6010D	2-27-18	2-27-18	
Mercury	ND	0.31	EPA 7471B	2-27-18	2-27-18	
Client ID:	PP12-7					
	02-234-03					
Laboratory ID: Arsenic	16	15	EPA 6010D	2-27-18	2-27-18	
Cadmium	4.1	0.73	EPA 6010D	2-27-16 2-27-18	2-27-18	
Chromium	31	0.73	EPA 6010D	2-27-18	2-27-18	
Lead	500	7.3	EPA 6010D	2-27-18	2-27-18	
Mercury	0.76	0.37	EPA 7471B	2-27-18	2-27-18	
Worday	00	0.07	217(71718	2 27 10	2 27 10	
Client ID:	PP13-10					
Laboratory ID:	02-234-08					
Arsenic	ND	12	EPA 6010D	2-27-18	2-27-18	
Cadmium	3.5	0.62	EPA 6010D	2-27-18	2-27-18	
Chromium	29	0.62	EPA 6010D	2-27-18	2-27-18	
Lead	330	6.2	EPA 6010D	2-27-18	2-27-18	
Mercury	ND	0.31	EPA 7471B	2-27-18	2-27-18	
Client ID:	PP15-6					
Laboratory ID:	02-234-11					
Arsenic	ND	15	EPA 6010D	2-27-18	2-27-18	
Cadmium	0.79	0.76	EPA 6010D	2-27-18	2-27-18	
Chromium	22	0.76	EPA 6010D	2-27-18	2-27-18	
Lead	75	7.6	EPA 6010D	2-27-18	2-27-18	
Mercury	ND	0.38	EPA 7471B	2-27-18	2-27-18	

Laboratory Reference: 1802-234 Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B

Matrix: Soil

Offits. Hig/Kg (ppin)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP15-12			-		
Laboratory ID:	02-234-12					
Arsenic	24	12	EPA 6010D	2-27-18	2-27-18	
Cadmium	1.0	0.59	EPA 6010D	2-27-18	2-27-18	
Chromium	34	0.59	EPA 6010D	2-27-18	2-27-18	
Lead	270	5.9	EPA 6010D	2-27-18	2-27-18	
Mercury	ND	0.30	EPA 7471B	2-27-18	2-27-18	
Client ID:	PP19-7					
Laboratory ID:	02-234-15					
Arsenic	ND	13	EPA 6010D	2-27-18	2-27-18	
Cadmium	ND	0.63	EPA 6010D	2-27-18	2-27-18	
Chromium	17	0.63	EPA 6010D	2-27-18	2-27-18	
Lead	82	6.3	EPA 6010D	2-27-18	2-27-18	
Mercury	ND	0.32	EPA 7471B	2-27-18	2-27-18	
Client ID:	PP19-10					
Laboratory ID:	02-234-16					
Arsenic	ND	12	EPA 6010D	2-27-18	2-27-18	
Cadmium	0.95	0.59	EPA 6010D	2-27-18	2-27-18	
Chromium	30	0.59	EPA 6010D	2-27-18	2-27-18	
Lead	340	5.9	EPA 6010D	2-27-18	2-27-18	
Mercury	ND	0.29	EPA 7471B	2-27-18	2-27-18	
Client ID:	PP19-15					
Laboratory ID:	02-234-17					
Arsenic	ND	12	EPA 6010D	2-27-18	2-27-18	
Cadmium	3.2	0.60	EPA 6010D	2-27-18	2-27-18	
Chromium	17	0.60	EPA 6010D	2-27-18	2-27-18	
Lead	96	6.0	EPA 6010D	2-27-18	2-27-18	
Mercury	ND	0.30	EPA 7471B	2-27-18	2-27-18	
o.oury	110	0.00	217(71710	221 10	22110	

Laboratory Reference: 1802-234 Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B

Matrix: Soil

Offits. Hig/Kg (ppin)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP22-4					
Laboratory ID:	02-234-18					
Arsenic	ND	13	EPA 6010D	2-27-18	2-27-18	
Cadmium	ND	0.64	EPA 6010D	2-27-18	2-27-18	
Chromium	9.9	0.64	EPA 6010D	2-27-18	2-27-18	
Lead	ND	6.4	EPA 6010D	2-27-18	2-27-18	
Mercury	ND	0.32	EPA 7471B	2-27-18	2-27-18	
Client ID:	PP22-9					
Laboratory ID:	02-234-19					
Arsenic	ND	12	EPA 6010D	2-27-18	2-27-18	
Cadmium	ND	0.61	EPA 6010D	2-27-18	2-27-18	
Chromium	9.9	0.61	EPA 6010D	2-27-18	2-27-18	
Lead	ND	6.1	EPA 6010D	2-27-18	2-27-18	
Mercury	ND	0.31	EPA 7471B	2-27-18	2-27-18	
Client ID:	PP22-13					
Laboratory ID:	02-234-20					
Arsenic	ND	12	EPA 6010D	2-27-18	2-27-18	
Cadmium	2.4	0.58	EPA 6010D	2-27-18	2-27-18	
Chromium	36	0.58	EPA 6010D	2-27-18	2-27-18	
Lead	270	5.8	EPA 6010D	2-27-18	2-27-18	
Mercury	ND	0.29	EPA 7471B	2-27-18	2-27-18	
Client ID:	PP25-7					
Laboratory ID:	02-234-23	1.1	EDA 6010D	2 27 40	2 27 40	
Arsenic	ND	14	EPA 6010D	2-27-18	2-27-18	
Chromium	ND	0.68	EPA 6010D	2-27-18	2-27-18	
Chromium	13	0.68	EPA 6010D	2-27-18	2-27-18	
Lead	8.8 ND	6.8	EPA 6010D	2-27-18	2-27-18	
Mercury	ND	0.34	EPA 7471B	2-27-18	2-27-18	

Laboratory Reference: 1802-234 Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B

Matrix: Soil

Offits. Hig/Kg (ppin)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP25-13					
Laboratory ID:	02-234-24					
Arsenic	ND	17	EPA 6010D	2-27-18	2-27-18	
Cadmium	1.9	0.85	EPA 6010D	2-27-18	2-27-18	
Chromium	38	0.85	EPA 6010D	2-27-18	2-27-18	
Lead	140	8.5	EPA 6010D	2-27-18	2-27-18	
Mercury	ND	0.42	EPA 7471B	2-27-18	2-27-18	
Client ID:	PP26-11					
Laboratory ID:	02-234-28					
Arsenic	ND	12	EPA 6010D	2-27-18	2-27-18	
Cadmium	ND	0.58	EPA 6010D	2-27-18	2-27-18	
Chromium	21	0.58	EPA 6010D	2-27-18	2-27-18	
Lead	ND	5.8	EPA 6010D	2-27-18	2-27-18	
Mercury	ND	0.29	EPA 7471B	2-27-18	2-27-18	
Client ID:	PP26-17					
Laboratory ID:	02-234-29					
Arsenic	ND	15	EPA 6010D	2-27-18	2-27-18	
Cadmium	ND	0.77	EPA 6010D	2-27-18	2-27-18	
Chromium	15	0.77	EPA 6010D	2-27-18	2-27-18	
Lead	ND	7.7	EPA 6010D	2-27-18	2-27-18	
Mercury	ND	0.39	EPA 7471B	2-27-18	2-27-18	
Client ID:	DD24 0					
Client ID:	PP34-8					
Laboratory ID:	02-234-31	40	EDA 00405	0.07.10	0.07.10	
Arsenic	ND	12	EPA 6010D	2-27-18	2-27-18	
Cadmium	ND	0.61	EPA 6010D	2-27-18	2-27-18	
Chromium	12	0.61	EPA 6010D	2-27-18	2-27-18	
Lead	6.1	6.1	EPA 6010D	2-27-18	2-27-18	
Mercury	ND	0.30	EPA 7471B	2-27-18	2-27-18	

Laboratory Reference: 1802-234 Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B **QUALITY CONTROL**

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0227SM2					
Arsenic	ND	10	EPA 6010D	2-27-18	2-27-18	
Cadmium	ND	0.50	EPA 6010D	2-27-18	2-27-18	
Chromium	ND	0.50	EPA 6010D	2-27-18	2-27-18	
Lead	ND	5.0	EPA 6010D	2-27-18	2-27-18	
Laboratory ID:	MB0227S1					
Mercury	ND	0.25	EPA 7471B	2-27-18	2-27-18	

					Source	Pe	rcent	Recovery		RPD	
Analyte	Re	sult	Spike	Level	Result	Rec	overy	Limits	RPD	Limit	Flags
DUPLICATE											
Laboratory ID:	02-2	34-29									
	ORIG	DUP									
Arsenic	ND	ND	NA	NA		1	NA	NA	NA	20	
Cadmium	ND	ND	NA	NA		1	NΑ	NA	NA	20	
Chromium	9.55	9.05	NA	NA		1	NΑ	NA	5	20	
Lead	ND	ND	NA	NA		1	NΑ	NA	NA	20	
Laboratory ID:	02-2	34-29									
Mercury	ND	ND	NA	NA		1	NA	NA	NA	20	
MATRIX SPIKES											
Laboratory ID:	02-2	34-29									
Laboratory ID.	MS	MSD	MS	MSD		MS	MSD				
Arsenic	97.5	96.8	100	100	ND	98	97	75-125	1	20	
Cadmium	46.1	46.0	50.0	50.0	ND	92	92	75-125	0	20	
Chromium	104	105	100	100	9.55	95	95	75-125	0	20	
Lead	227	229	250	250	ND	91	92	75-125	1	20	
Laboratory ID:	02-2	34-29									
Mercury	0.575	0.571	0.500	0.500	0.00680	114	113	80-120	1	20	

Laboratory Reference: 1802-234

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B

Matrix: Soil

g/. tg (pp)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP12-10					
Laboratory ID:	02-234-04					
Arsenic	ND	13	EPA 6010D	3-2-18	3-2-18	
Cadmium	ND	0.64	EPA 6010D	3-2-18	3-2-18	
Chromium	25	0.64	EPA 6010D	3-2-18	3-2-18	
Lead	30	6.4	EPA 6010D	3-2-18	3-2-18	
Mercury	ND	0.32	EPA 7471B	3-7-18	3-7-18	
Client ID:	PP25-17					
Laboratory ID:	02-234-25					
Arsenic	ND	19	EPA 6010D	3-2-18	3-2-18	
Cadmium	ND	0.93	EPA 6010D	3-2-18	3-2-18	
Chromium	28	0.93	EPA 6010D	3-2-18	3-2-18	
Lead	ND	9.3	EPA 6010D	3-2-18	3-2-18	
Mercury	ND	0.47	EPA 7471B	3-7-18	3-7-18	

Laboratory Reference: 1802-234 Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B **QUALITY CONTROL**

Matrix: Soil

Result						
	PQL	Method	Prepared	Analyzed	Flags	
MB0302SM2						
ND	10	EPA 6010D	3-2-18	3-2-18		
ND	0.50	EPA 6010D	3-2-18	3-2-18		
ND	0.50	EPA 6010D	3-2-18	3-2-18		
ND	5.0	EPA 6010D	3-2-18	3-2-18		
MB0307S2						
ND	0.25	EPA 7471B	3-7-18	3-7-18	•	
	ND ND ND ND	ND 10 ND 0.50 ND 0.50 ND 5.0 MB0307S2	ND 10 EPA 6010D ND 0.50 EPA 6010D ND 0.50 EPA 6010D ND 5.0 EPA 6010D MB0307S2	ND 10 EPA 6010D 3-2-18 ND 0.50 EPA 6010D 3-2-18 ND 0.50 EPA 6010D 3-2-18 ND 5.0 EPA 6010D 3-2-18 MB0307S2	ND 10 EPA 6010D 3-2-18 3-2-18 ND 0.50 EPA 6010D 3-2-18 3-2-18 ND 0.50 EPA 6010D 3-2-18 3-2-18 ND 5.0 EPA 6010D 3-2-18 3-2-18 MB0307S2	

					Source	Pe	rcent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Rec	overy	Limits	RPD	Limit	Flags
DUPLICATE											
Laboratory ID:	02-20	65-04									
	ORIG	DUP									
Arsenic	ND	ND	NA	NA			NA	NA	NA	20	
Cadmium	ND	ND	NA	NA		I	NA	NA	NA	20	
Chromium	67.5	67.5	NA	NA		-	NA	NA	0	20	
Lead	8.25	6.70	NA	NA			NA	NA	21	20	С
Laboratory ID:	02-2	71-21									
Mercury	ND	ND	NA	NA			NA	NA	NA	20	
MATRIX SPIKES											
Laboratory ID:	02-20	65-04									
	MS	MSD	MS	MSD		MS	MSD				
Arsenic	90.0	89.4	100	100	ND	90	89	75-125	1	20	
Cadmium	44.5	43.9	50.0	50.0	ND	89	88	75-125	1	20	
Chromium	156	153	100	100	67.5	88	85	75-125	2	20	
Lead	219	219	250	250	8.25	84	84	75-125	0	20	
Laboratory ID:	02-2	71-21									
Mercury	0.605	0.590	0.500	0.500	0.0364	114	111	80-120	3	20	

TCLP LEAD EPA 1311/6010D

Matrix: TCLP Extract Units: mg/L (ppm)

····g/ = (pp····/				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP12-7					
Laboratory ID:	02-234-03					
Lead	0.23	0.20	EPA 6010D	3-16-18	3-16-18	
Client ID:	PP13-10					
Laboratory ID:	02-234-08					
Lead	ND	0.20	EPA 6010D	3-16-18	3-16-18	
Client ID:	PP15-12					
Laboratory ID:	02-234-12					
Lead	3.8	0.20	EPA 6010D	3-16-18	3-16-18	
Client ID:	PP19-10					
Laboratory ID:	02-234-16					
Lead	1.3	0.20	EPA 6010D	3-16-18	3-16-18	
Client ID:	PP22-13					
Laboratory ID:	02-234-20					
Lead	ND ND	0.20	EPA 6010D	3-16-18	3-16-18	
Client ID:	PP25-13					
Laboratory ID:	02-234-24					
Lead	ND	0.20	EPA 6010D	3-16-18	3-16-18	

Laboratory Reference: 1802-234 Project: 17-06520-000

TCLP LEAD EPA 1311/6010D **QUALITY CONTROL**

Matrix: TCLP Extract Units: mg/L (ppm)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0316TM1					
Lead	ND	0.20	EPA 6010D	3-16-18	3-16-18	

Analyte Resul		sult	Spike Le		Source Result		rcent	Recovery Limits	RPD	RPD Limit	Flags
DUPLICATE			-								
Laboratory ID:	02-23	34-03									
	ORIG	DUP									
Lead	0.228	ND	NA	NA		1	NA	NA	NA	20	
MATRIX SPIKES											
Laboratory ID:	02-234-03										
	MS	MSD	MS	MSD		MS	MSD				
Lead	9.06	9.12	10.0	10.0	0.228	88	89	75-125	1	20	•

% MOISTURE

Date Analyzed: 2-23&28&3-5-18

Client ID	Lab ID	% Moisture
PP12-2	02-234-02	19
PP12-7	02-234-03	32
PP12-10	02-234-04	22
PP13-10	02-234-08	20
PP15-6	02-234-11	34
PP15-12	02-234-12	16
PP19-7	02-234-15	21
PP19-10	02-234-16	15
PP19-15	02-234-17	16
PP22-4	02-234-18	22
PP22-9	02-234-19	19
PP22-13	02-234-20	14
PP25-7	02-234-23	27
PP25-13	02-234-24	41
PP25-17	02-234-25	46
PP26-11	02-234-28	14
PP26-17	02-234-29	35
PP34-8	02-234-31	18
PP34-15	02-234-33	28



Data Qualifiers and Abbreviations

- A Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
- B The analyte indicated was also found in the blank sample.
- C The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
- E The value reported exceeds the quantitation range and is an estimate.
- F Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
- H The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
- I Compound recovery is outside of the control limits.
- J The value reported was below the practical quantitation limit. The value is an estimate.
- K Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
- L The RPD is outside of the control limits.
- M Hydrocarbons in the gasoline range are impacting the diesel range result.
- M1 Hydrocarbons in the gasoline range (toluene-naphthalene) are present in the sample.
- N Hydrocarbons in the lube oil range are impacting the diesel range result.
- N1 Hydrocarbons in diesel range are impacting lube oil range results.
- O Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
- P The RPD of the detected concentrations between the two columns is greater than 40.
- Q Surrogate recovery is outside of the control limits.
- S Surrogate recovery data is not available due to the necessary dilution of the sample.
- T The sample chromatogram is not similar to a typical ______.
- U The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- U1 The practical quantitation limit is elevated due to interferences present in the sample.
- V Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
- W Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
- X Sample extract treated with a mercury cleanup procedure.
- X1- Sample extract treated with a sulfuric acid/silica gel cleanup procedure.
- Y The calibration verification for this analyte exceeded the 20% drift specified in method 8260C, and therefore the reported result should be considered an estimate. The overall performance of the calibration verification standard met the acceptance criteria of the method.

7 -

ND - Not Detected at PQL

PQL - Practical Quantitation Limit

RPD - Relative Percent Difference





Chain of Custody

Page _ | _ of _ 4

Analytical Laboratory Testing Services 14648 NE 95th Street • Redmond, WA 98052			Turnaround Request (in working days)			L	Laboratory Number: 02 - 234																
Project Name: Project Name: Project Manag Sampled by:	1-010520-000 Dic Pare	Same 2 Day		1 Day 3 Days ys)	Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX		NWTPH-Dx (Acid / SG Clean-up)	Volatiles 8260C Halogenated Volatiles 8260C	EDB EPA 8011 (Waters Only)	Semivolatiles 8270D/SIM (with low-level PAHs)	PAHS 8270D/SIM (low-level) CPAHS	PCBs 8082A	Organochlorine Pesticides 8081B	Organistical And Larbinidas 927 UL/SIIM	Total RCRA Metals	Total MTCA Metals	TCLP Metals LEAD ONLY	1 1			% Moisture
	P*10-10	2.21.18		soil	2	A	A		A				A	A				A					X
	712-2		9:50	1	2	0	A	1				(X					X)				0
3 P	712-7		10:00		2	0	A	1				(X					X	0				
4 PI	712-10		10:10		2	A	A		9									0	1				
5 P	712-16		10:20		2	A	A		A				A	A				A					X
6 PF	713-2		10:30		2	A	A		A				A	A				A					X
7 PP	13-6		10:40		2	A	A		A				A	A				Ø	+				X
8 PF	13-10		10150		2	0	A		A				X)	A				X	0				Q
9 7	713-15		11:00		2	A	A		A				A	A				A					X
10 PP	15-4	/	11:40	V	2	A	A		A				A	A				A					X
	Signature (mpany				Date			Time				ts/Spe				1	0				
Relinquished	Mayer his	0	terre	NOC			1	122		181		1	7	-	AI	ch) , V	en	for	15	0,	Don	25
Received Relinquished			OSE			_	2/22/18		8	1505		A-Archivefor 30 Days Added 2/27/18.DB(STA)											
Received									+			C) 4	dde	d	3/	3/1	9. [13	CST	A)	/
Relinquished					_				+											75	/		
Received									+			Data	Pac	kage:	Stan	dard		evel II		Level	IV 🗆		
Reviewed/Date			Reviewed/Date									Chromatograms with final report ☐ Electronic Data Deliverables (EDDs) ☐											



Chain of Custody

Page 2 of 4

	Analytical Laboratory Testing Services 14648 NE 95th Street • Redmond, WA 98052		rnaround Req in working da			L	abo	orato	ry l	Num	ber:			0	2	- 2	3	4							
Projec	Phone: (425) 883-3881 - www.onsite-env.com any: HOVEYA et Number: 17 - 0 10 5 20 - 000 et Name: Pauli Payk et Manager: BNCE Carpentor BWCE Carpentor BWCE Carpentor	(TPH	dard (7 Days) I analysis 5 Da (other)	1 Day 3 Days	Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX		NW I PH-Dx (Acid / SG Clean-up)	Halogenated Volatiles 8260C	EDB EPA 8011 (Waters Only)	Semivolatiles 8270D/SIM (with low-level PAHs)	S 8270D/SIM (low-level) CPAHS	PCBS 8082A	Organochlorine Pesticides 8081B	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	TCLP Metals LEAD ONLY	HEM (oil and grease) 1664A				% Moisture
Lab ID	And the second s	Date Sampled	Time Sampled	Matrix	Nun	NN	NM			Halo	EDB			L LCB	organ C	Orga	S	Total	1	TCLI	HEN			_	
11	PPIS-U	2-21.18		9011	1	X	A		0	-	+		X	2	-	-		_(X	/				-	8
12	PP15-12		12:05	-	2	(X)		-	A	+	+			A		-		((X)	0				4	\otimes
13	PP15-16		12:15		2	A	A		A	+	+		A	A		-			A					_	X
1+	PP19-3		12:20		2	A		-	A	-	-		A	A	4				A					1	X
15	PP19-7		12:30		2	X	A					1	X	A	4			(X)						\otimes
10	PP19-10		12:40		2	X	A		A				X)	A					\otimes	0					8
17	PP19-15		12:50		2	X	A		A				()	4				(D						X
18	PP22-4		1340		2	X	A		A				2	A					8						(>)
M	PP22-9		1350		2	0	A		4				D.	A				(X						0
20	PP 22-13	1	1355	V	2	X	A		4				DI.	4				(D	0					X
	Signature		ompany				Date			ime		Com	ments	s/Spe	cial li	nstruc	tion	S							
Relin	quished mayuu MF	**	Herre	ra				12																	
Rece			OSE				2/:	22/	8	15	05	+													
Relin	equished O																								
777.1	ived																								
	quished																								
-	ived											Data	Pack	age:	Stan	dard		Lev	el III		Leve	IV]		
Revi	ewed/Date		Reviewed/Dat	е								Chron	matog	grams	with	final	repo	ort 🗌	Elec	ctroni	c Data	a Deli	verables	s (EDDs	s) 🗆

OnSite Environmental Inc.

Chain of Custody

Page 3 of 4

Analytical Laboratory Testing Services 14648 NE 95th Street • Redmond, WA 98052	Turn (in	naround Req working da	uest ys)		L	abo	rato	ry N	luml	ber:		C	12	- 2	23	4								
Phone: (425) 883-3881 • www.onsite-env.com Company: HOVEVA Project Number: 17-00520-600 Project Name: Paulic Paul Project Manager: Sampled by: B WCL Caupertev Lab ID Sample Identification	Same 2 Day	,	1 Day	Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX		NWTPH-Dx (Acid / SG Clean-up)	Halogenated Volatiles 8260C	EDB EPA 8011 (Waters Only)	Semivolatiles 8270D/SIM (with low-level PAHs)	PAHS 8270D/SIM (low-level) CPAHS	PCBs 8082A	Organochlorine Pesticides 8081B	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	TCLP Metals LEAD ONLY	HEM (oil and grease) 1664A				% Moisture
21 PP2Z-17	2-21.18		Gril	2	A			A				A	A					A						X
22 PP 25-3	1	1420	(2	A			A				A	A					A						X
23 PP25-7		1425		2	8	A		A				X	A				((X))					(X)
24 PP25-13		1430		2	8	9						\otimes	0				($\widehat{\chi}$	0				(X
25 PP25-17		1445		2	A	10							9					•						
26 PP 210-3		1455		2	A	A		A				A	A					A						X
27 PP210-7		1505		2	A	A		A				A	A					A						X
28 PD 24-11		KID		2	X	A		A			((X)	A				(X					(\otimes
29 PP210-17		KB		2	X	A		A				X	A				(X					(XX
30 PP 34-4	Y	1540	V	2	A	A		A				A	A					A						X
Signature	\	mpany				Date	۸		ime		Cor	nmer	ts/Sp	ecial I	Instru	ıctior	ns							
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Received Relinquished	- 0	SE				21	22	1,8	150	25														
Received				-				-			-													
Relinquished					-			+			-													
Received								+	-		Data	a Pac	kage	Star	ndar	d 🗆	Lev	/el III		Level	IV 🗆			
Reviewed/Date		Reviewed/Da	te														_					rables (E	DDs) [



Chain of Custody

Page 4 of 4

	Analytical Laboratory Testing Services 14648 NE 95th Street • Redmond, WA 98052		naround Req n working da			L	abo	rato	ry l	lumb	er:		()2	-	23	3 4								
Projec	HUVEVA Number: 17-0USZU-000 Name: Paufic Payk Manager: BUML Carrenter	Same 2 Da Stand (TPH		1 Day 3 Days	Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX	NWTPH-Gx	White Spen	Halogenated Volatiles 8260C	EDB EPA 8011 (Waters Only)	Semivolatiles 8270D/SIM (with low-level PAHs)	PAHS 8270D/SIM (low-level) CPAHS.	PCBs 8082A	Organochlorine Pesticides 8081B	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	TCLP Metals	HEM (oil and grease) 1664A				% Moisture
31	PP34-8	2.21.5		soil	2	(2)	A		A			(3	A		-			1						(X)
32	PP34-13		1600		2	×	A		A				A	A					A						X
33	PP34-15		1005	/	2	A	A	4	4				•	A					A						•
									-														+		
Reline	guished Signature		HUVE OSP	ua				22/1	8	ime 150		Com	nmen	ts/Sp	ecial	Instru	iction	ıs					- Con	641	
Relin	quished		-66																						
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nevie	wed/ Date		neviewed/Da	le								Chro	mato	ogram	s wit	h fina	al rep	ort [Ele	ctronic	Data	Delive	erables	(EDDs)	



14648 NE 95th Street, Redmond, WA 98052 • (425) 883-3881

March 15, 2018

Bruce Carpenter Herrera Environmental Consultants, Inc. 2200 6th Avenue, Suite 1100 Seattle, WA 98121

Re: Analytical Data for Project 17-06520-000

Laboratory Reference No. 1802-278

Dear Bruce:

Enclosed are the analytical results and associated quality control data for samples submitted on February 28, 2018.

The standard policy of OnSite Environmental, Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

David Baumeister Project Manager

Enclosures

Project: 17-06520-000

Case Narrative

Samples were collected on February 26 and 27, 2018 and received by the laboratory on February 28, 2018. They were maintained at the laboratory at a temperature of 2°C to 6°C.

Please note that any and all soil sample results are reported on a dry-weight basis, unless otherwise noted below.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

Project: 17-06520-000

NWTPH-HCID

Matrix: Soil

Offits. Ing/Kg (ppm)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-11_2.5					
Laboratory ID:	02-278-01					
Gasoline Range Organics	ND	26	NWTPH-HCID	3-1-18	3-1-18	
Diesel Range Organics	ND	65	NWTPH-HCID	3-1-18	3-1-18	
Lube Oil	Detected	130	NWTPH-HCID	3-1-18	3-1-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	101	50-150				
Client ID:	B-11_12.5					
Laboratory ID:	02-278-04					
Gasoline Range Organics	ND	24	NWTPH-HCID	3-1-18	3-1-18	
Diesel Range Organics	ND	60	NWTPH-HCID	3-1-18	3-1-18	
Lube Oil Range Organics	ND	120	NWTPH-HCID	3-1-18	3-1-18	
Surrogate:	Percent Recovery	Control Limits	14W II II-IIOID	0-1-10	0-1-10	
o-Terphenyl	116	50-150				
o respicings	770	00 100				
Client ID:	B-10_7.5					
Laboratory ID:	02-278-06					
Gasoline Range Organics	ND	24	NWTPH-HCID	3-1-18	3-1-18	
Diesel Range Organics	ND	59	NWTPH-HCID	3-1-18	3-1-18	
Lube Oil	Detected	120	NWTPH-HCID	3-1-18	3-1-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	111	50-150				
Client ID:	B-08_5					
Laboratory ID:	02-278-09					
Gasoline Range Organics	ND	28	NWTPH-HCID	3-1-18	3-1-18	
Diesel Range Organics	ND	69	NWTPH-HCID	3-1-18	3-1-18	
Lube Oil Range Organics	ND	140	NWTPH-HCID	3-1-18	3-1-18	
Surrogate:	Percent Recovery	Control Limits		<u> </u>	<u> </u>	
o-Terphenyl	105	50-150				
σ . σ. μσ. η.		00 .00				
Client ID:	B-08_12.5					
Laboratory ID:	02-278-11					
Gasoline Range Organics	ND	27	NWTPH-HCID	3-1-18	3-1-18	
Diesel Range Organics	ND	67	NWTPH-HCID	3-1-18	3-1-18	
Lube Oil Range Organics	ND	130	NWTPH-HCID	3-1-18	3-1-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	107	50-150				

Project: 17-06520-000

NWTPH-HCID

Matrix: Soil

5 5 41 7				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-07_2.5					
Laboratory ID:	02-278-13					
Gasoline Range Organics	ND	23	NWTPH-HCID	3-1-18	3-1-18	
Diesel Range Organics	ND	180	NWTPH-HCID	3-1-18	3-1-18	U1
Lube Oil	Detected	120	NWTPH-HCID	3-1-18	3-1-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	118	50-150				
Client ID:	B-07_7.5					
Laboratory ID:	02-278-14					
Gasoline Range Organics	ND	25	NWTPH-HCID	3-1-18	3-1-18	
Diesel Range Organics	ND	290	NWTPH-HCID	3-1-18	3-1-18	U1
Lube Oil	Detected	120	NWTPH-HCID	3-1-18	3-1-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	114	50-150				

Project: 17-06520-000

NWTPH-HCID QUALITY CONTROL

Matrix: Soil

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
METHOD BLANK				-	_	
Laboratory ID:	MB0301S1					
Gasoline Range Organics	ND	20	NWTPH-HCID	3-1-18	3-1-18	
Diesel Range Organics	ND	50	NWTPH-HCID	3-1-18	3-1-18	
Lube Oil Range Organics	ND	100	NWTPH-HCID	3-1-18	3-1-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	116	50-150				

Project: 17-06520-000

NWTPH-Dx

Matrix: Soil

Analyto	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Analyte Client ID:	B-11 2.5	FWL	METHOR	Fiepaieu	Allalyzeu	ı ıayə
Laboratory ID:	02-278-01					
Diesel Range Organics	nd	33	NWTPH-Dx	3-9-18	3-9-18	U1
Lube Oil Range Organics	380	65	NWTPH-Dx	3-9-18	3-9-18	01
Surrogate:	Percent Recovery	Control Limits		00.0	0 0 10	
o-Terphenyl	95	50-150				
, ,						
Client ID:	B-10_7.5					
Laboratory ID:	02-278-06					
Diesel Range Organics	ND	30	NWTPH-Dx	3-9-18	3-12-18	
Lube Oil	88	59	NWTPH-Dx	3-9-18	3-12-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	51	50-150				
Client ID:	B-07 2.5					
Laboratory ID:	02-278-13					
Diesel Range Organics	ND	150	NWTPH-Dx	3-9-18	3-9-18	U1
Lube Oil	4400	290	NWTPH-Dx	3-9-18	3-9-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	81	50-150				
Client ID:	B-07_7.5					
Laboratory ID:	02-278-14					
Diesel Range Organics	ND	310	NWTPH-Dx	3-9-18	3-9-18	
Lube Oil	1800	620	NWTPH-Dx	3-9-18	3-9-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl		50-150				S
Client ID:	B-07_12.5					
Laboratory ID:	02-278-15					
Diesel Range Organics	ND	33	NWTPH-Dx	3-9-18	3-9-18	
Lube Oil	180	66	NWTPH-Dx	3-9-18	3-9-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	106	50-150				

Project: 17-06520-000

NWTPH-Dx QUALITY CONTROL

Matrix: Soil

Analyte	Result	PQL	Method	Date Prepared	Date Analvzed	Flags
METHOD BLANK	Result	1 &L	Hictioa	ricpared	Allalyzea	ı iays
Laboratory ID:	MB0309S1					
Diesel Range Organics	ND	25	NWTPH-Dx	3-9-18	3-9-18	
Lube Oil Range Organics	ND	50	NWTPH-Dx	3-9-18	3-9-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	91	50-150				

					Source	Percent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Recovery	Limits	RPD	Limit	Flags
DUPLICATE										
Laboratory ID:	02-27	78-15								
	ORIG	DUP								
Diesel Range	ND	ND	NA	NA		NA	NA	NA	NA	
Lube Oil	137	93.2	NA	NA		NA	NA	38	NA	
Surrogate:										
o-Terphenyl						106 75	50-150			

Project: 17-06520-000

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-11_2.5					
Laboratory ID:	02-278-01					
Benzo[a]anthracene	0.016	0.0086	EPA 8270D/SIM	3-5-18	3-5-18	
Chrysene	0.035	0.0086	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[b]fluoranthene	0.053	0.0086	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo(j,k)fluoranthene	0.014	0.0086	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[a]pyrene	0.019	0.0086	EPA 8270D/SIM	3-5-18	3-5-18	
Indeno(1,2,3-c,d)pyrene	0.017	0.0086	EPA 8270D/SIM	3-5-18	3-5-18	
Dibenz[a,h]anthracene	ND	0.0086	EPA 8270D/SIM	3-5-18	3-5-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	70	32 - 115				
Pyrene-d10	85	35 - 129				
Terphenyl-d14	88	33 - 114				

Project: 17-06520-000

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-11_12.5					
Laboratory ID:	02-278-04					
Benzo[a]anthracene	ND	0.0079	EPA 8270D/SIM	3-5-18	3-5-18	
Chrysene	ND	0.0079	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[b]fluoranthene	ND	0.0079	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo(j,k)fluoranthene	ND	0.0079	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[a]pyrene	ND	0.0079	EPA 8270D/SIM	3-5-18	3-5-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0079	EPA 8270D/SIM	3-5-18	3-5-18	
Dibenz[a,h]anthracene	ND	0.0079	EPA 8270D/SIM	3-5-18	3-5-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	62	32 - 115				
Pyrene-d10	77	35 - 129				
Terphenyl-d14	80	33 - 114				

Project: 17-06520-000

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-10_7.5					
Laboratory ID:	02-278-06					
Benzo[a]anthracene	ND	0.0079	EPA 8270D/SIM	3-5-18	3-5-18	
Chrysene	ND	0.0079	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[b]fluoranthene	ND	0.0079	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo(j,k)fluoranthene	ND	0.0079	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[a]pyrene	ND	0.0079	EPA 8270D/SIM	3-5-18	3-5-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0079	EPA 8270D/SIM	3-5-18	3-5-18	
Dibenz[a,h]anthracene	ND	0.0079	EPA 8270D/SIM	3-5-18	3-5-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	73	32 - 115				
Pyrene-d10	87	35 - 129				
Terphenyl-d14	92	33 - 114				

Project: 17-06520-000

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-08_5					
Laboratory ID:	02-278-09					
Benzo[a]anthracene	ND	0.0092	EPA 8270D/SIM	3-5-18	3-6-18	
Chrysene	ND	0.0092	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[b]fluoranthene	ND	0.0092	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo(j,k)fluoranthene	ND	0.0092	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[a]pyrene	ND	0.0092	EPA 8270D/SIM	3-5-18	3-6-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0092	EPA 8270D/SIM	3-5-18	3-6-18	
Dibenz[a,h]anthracene	ND	0.0092	EPA 8270D/SIM	3-5-18	3-6-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	77	32 - 115				
Pyrene-d10	87	35 - 129				
Terphenyl-d14	90	33 - 114				

Project: 17-06520-000

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-08_12.5					
Laboratory ID:	02-278-11					
Benzo[a]anthracene	ND	0.0089	EPA 8270D/SIM	3-5-18	3-5-18	
Chrysene	ND	0.0089	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[b]fluoranthene	ND	0.0089	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo(j,k)fluoranthene	ND	0.0089	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[a]pyrene	ND	0.0089	EPA 8270D/SIM	3-5-18	3-5-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0089	EPA 8270D/SIM	3-5-18	3-5-18	
Dibenz[a,h]anthracene	ND	0.0089	EPA 8270D/SIM	3-5-18	3-5-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	76	32 - 115				
Pyrene-d10	88	35 - 129				
Terphenyl-d14	98	33 - 114				

Project: 17-06520-000

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-07_2.5					
Laboratory ID:	02-278-13					
Benzo[a]anthracene	ND	0.039	EPA 8270D/SIM	3-5-18	3-5-18	
Chrysene	ND	0.039	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[b]fluoranthene	ND	0.039	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo(j,k)fluoranthene	ND	0.039	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[a]pyrene	ND	0.039	EPA 8270D/SIM	3-5-18	3-5-18	
Indeno(1,2,3-c,d)pyrene	ND	0.039	EPA 8270D/SIM	3-5-18	3-5-18	
Dibenz[a,h]anthracene	ND	0.039	EPA 8270D/SIM	3-5-18	3-5-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	78	32 - 115				
Pyrene-d10	89	35 - 129				
Terphenyl-d14	81	33 - 114				

Project: 17-06520-000

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-07_7.5					
Laboratory ID:	02-278-14					
Benzo[a]anthracene	ND	0.041	EPA 8270D/SIM	3-5-18	3-5-18	
Chrysene	ND	0.041	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[b]fluoranthene	ND	0.041	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo(j,k)fluoranthene	ND	0.041	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[a]pyrene	ND	0.041	EPA 8270D/SIM	3-5-18	3-5-18	
Indeno(1,2,3-c,d)pyrene	ND	0.041	EPA 8270D/SIM	3-5-18	3-5-18	
Dibenz[a,h]anthracene	ND	0.041	EPA 8270D/SIM	3-5-18	3-5-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	82	32 - 115				
Pyrene-d10	93	35 - 129				
Terphenyl-d14	92	33 - 114				
•						

Project: 17-06520-000

PAHS EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0305S1					
Benzo[a]anthracene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Chrysene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[a]pyrene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	84	32 - 115				
Pyrene-d10	94	35 - 129				
Terphenyl-d14	102	33 - 114				

Project: 17-06520-000

PAHS EPA 8270D/SIM SB/SBD QUALITY CONTROL

						Per	cent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	F	Reco	overy	Limits	RPD	Limit	Flags
SPIKE BLANKS											
Laboratory ID:	SB03	05S1									
	SB	SBD	SB	SBD	5	SB	SBD				
Benzo[a]anthracene	0.0878	0.0851	0.0833	0.0833	1	05	102	64 - 135	3	15	
Chrysene	0.0869	0.0844	0.0833	0.0833	1	04	101	70 - 119	3	15	
Benzo[b]fluoranthene	0.0826	0.0807	0.0833	0.0833	9	99	97	54 - 135	2	15	
Benzo(j,k)fluoranthene	0.0872	0.0839	0.0833	0.0833	1	05	101	66 - 122	4	15	
Benzo[a]pyrene	0.0816	0.0789	0.0833	0.0833	ç	8	95	62 - 125	3	15	
Indeno(1,2,3-c,d)pyrene	0.0752	0.0724	0.0833	0.0833	ç	90	87	55 - 129	4	15	
Dibenz[a,h]anthracene	0.0769	0.0756	0.0833	0.0833	ç	92	91	58 - 125	2	15	
Surrogate:											
2-Fluorobiphenyl					8	38	87	32 - 115			
Pyrene-d10					9	99	96	35 - 129			
Terphenyl-d14					1	05	102	33 - 114			

Project: 17-06520-000

PAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-07_12.5					
Laboratory ID:	02-278-15					
Benzo[a]anthracene	ND	0.0088	EPA 8270D/SIM	3-12-18	3-12-18	
Chrysene	ND	0.0088	EPA 8270D/SIM	3-12-18	3-12-18	
Benzo[b]fluoranthene	ND	0.0088	EPA 8270D/SIM	3-12-18	3-12-18	
Benzo(j,k)fluoranthene	ND	0.0088	EPA 8270D/SIM	3-12-18	3-12-18	
Benzo[a]pyrene	ND	0.0088	EPA 8270D/SIM	3-12-18	3-12-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0088	EPA 8270D/SIM	3-12-18	3-12-18	
Dibenz[a,h]anthracene	ND	0.0088	EPA 8270D/SIM	3-12-18	3-12-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	57	32 - 115				
Pyrene-d10	68	35 - 129				
Terphenyl-d14	73	33 - 114				

Project: 17-06520-000

PAHS EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0312S2					
Benzo[a]anthracene	ND	0.0067	EPA 8270D/SIM	3-12-18	3-12-18	_
Chrysene	ND	0.0067	EPA 8270D/SIM	3-12-18	3-12-18	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270D/SIM	3-12-18	3-12-18	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270D/SIM	3-12-18	3-12-18	
Benzo[a]pyrene	ND	0.0067	EPA 8270D/SIM	3-12-18	3-12-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0067	EPA 8270D/SIM	3-12-18	3-12-18	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270D/SIM	3-12-18	3-12-18	
Surrogate:	Percent Recovery	Control Limits				_
2-Fluorobiphenyl	64	32 - 115				
Pyrene-d10	71	35 - 129				
Terphenyl-d14	77	33 - 114				

Project: 17-06520-000

PAHS EPA 8270D/SIM MS/MSD QUALITY CONTROL

					Source	Per	cent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Rec	overy	Limits	RPD	Limit	Flags
MATRIX SPIKES											_
Laboratory ID:	03-0	78-01									
	MS	MSD	MS	MSD		MS	MSD				
Benzo[a]anthracene	0.0626	0.0696	0.0833	0.0833	ND	75	84	27 - 143	11	23	
Chrysene	0.0626	0.0678	0.0833	0.0833	0.00705	67	73	22 - 130	8	24	
Benzo[b]fluoranthene	0.0610	0.0687	0.0833	0.0833	0.00792	64	73	15 - 141	12	26	
Benzo(j,k)fluoranthene	0.0517	0.0547	0.0833	0.0833	ND	62	66	42 - 112	6	24	
Benzo[a]pyrene	0.0579	0.0645	0.0833	0.0833	ND	70	77	33 - 126	11	26	
Indeno(1,2,3-c,d)pyrene	0.0607	0.0657	0.0833	0.0833	ND	73	79	30 - 125	8	25	
Dibenz[a,h]anthracene	0.0543	0.0584	0.0833	0.0833	ND	65	70	31 - 124	7	22	
Surrogate:											
2-Fluorobiphenyl						58	61	32 - 115			
Pyrene-d10						73	76	35 - 129			
Terphenyl-d14						62	64	33 - 114			

Project: 17-06520-000

PCBs by EPA 8082A

Matrix: Soil

orins. Trig/reg (ppin)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-11_2.5					
Laboratory ID:	02-278-01					
Aroclor 1016	ND	0.26	EPA 8082A	3-9-18	3-9-18	
Aroclor 1221	ND	0.26	EPA 8082A	3-9-18	3-9-18	
Aroclor 1232	ND	0.26	EPA 8082A	3-9-18	3-9-18	
Aroclor 1242	ND	0.26	EPA 8082A	3-9-18	3-9-18	
Aroclor 1248	ND	0.26	EPA 8082A	3-9-18	3-9-18	
Aroclor 1254	ND	0.26	EPA 8082A	3-9-18	3-9-18	
Aroclor 1260	ND	0.26	EPA 8082A	3-9-18	3-9-18	
Surrogate:	Percent Recovery	Control Limits				
DCB	91	40-134				
Client ID:	B-10_7.5					
Laboratory ID:	02-278-06					
Aroclor 1016	ND	0.059	EPA 8082A	3-9-18	3-9-18	
Aroclor 1221	ND	0.059	EPA 8082A	3-9-18	3-9-18	
Aroclor 1232	ND	0.059	EPA 8082A	3-9-18	3-9-18	
Aroclor 1242	ND	0.059	EPA 8082A	3-9-18	3-9-18	
Aroclor 1248	ND	0.059	EPA 8082A	3-9-18	3-9-18	
Aroclor 1254	ND	0.059	EPA 8082A	3-9-18	3-9-18	
Aroclor 1260	ND	0.059	EPA 8082A	3-9-18	3-9-18	
Surrogate:	Percent Recovery	Control Limits				
DCB	87	40-134				
Client ID:	B-07_2.5					
Laboratory ID:	02-278-13					
Aroclor 1016	ND	0.058	EPA 8082A	3-9-18	3-9-18	
Aroclor 1221	ND	0.058	EPA 8082A	3-9-18	3-9-18	
Aroclor 1232	ND	0.058	EPA 8082A	3-9-18	3-9-18	
Aroclor 1242	ND	0.058	EPA 8082A	3-9-18	3-9-18	
Aroclor 1248	ND	0.058	EPA 8082A	3-9-18	3-9-18	
Aroclor 1254	ND	0.058	EPA 8082A	3-9-18	3-9-18	
Aroclor 1260	ND	0.058	EPA 8082A	3-9-18	3-9-18	
Surrogate:	Percent Recovery	Control Limits				
DCB	73	40-134				

Project: 17-06520-000

PCBs by EPA 8082A

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-07_7.5					
Laboratory ID:	02-278-14					
Aroclor 1016	ND	0.062	EPA 8082A	3-9-18	3-9-18	
Aroclor 1221	ND	0.062	EPA 8082A	3-9-18	3-9-18	
Aroclor 1232	ND	0.062	EPA 8082A	3-9-18	3-9-18	
Aroclor 1242	ND	0.062	EPA 8082A	3-9-18	3-9-18	
Aroclor 1248	ND	0.062	EPA 8082A	3-9-18	3-9-18	
Aroclor 1254	ND	0.062	EPA 8082A	3-9-18	3-9-18	
Aroclor 1260	ND	0.062	EPA 8082A	3-9-18	3-9-18	
Surrogate:	Percent Recovery	Control Limits				
DCB	74	40-134				
Client ID:	B-07_12.5					
Laboratory ID:	02-278-15					
Aroclor 1016	ND	0.066	EPA 8082A	3-9-18	3-9-18	
Aroclor 1221	ND	0.066	EPA 8082A	3-9-18	3-9-18	
Aroclor 1232	ND	0.066	EPA 8082A	3-9-18	3-9-18	
Aroclor 1242	ND	0.066	EPA 8082A	3-9-18	3-9-18	
Aroclor 1248	ND	0.066	EPA 8082A	3-9-18	3-9-18	
Aroclor 1254	ND	0.066	EPA 8082A	3-9-18	3-9-18	
Aroclor 1260	ND	0.066	EPA 8082A	3-9-18	3-9-18	
Surrogate:	Percent Recovery	Control Limits				
DCB	70	40-134				

Project: 17-06520-000

PCBs by EPA 8082A QUALITY CONTROL

Matrix: Soil

Units: mg/Kg (ppm)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0309S1					
Aroclor 1016	ND	0.050	EPA 8082A	3-9-18	3-9-18	
Aroclor 1221	ND	0.050	EPA 8082A	3-9-18	3-9-18	
Aroclor 1232	ND	0.050	EPA 8082A	3-9-18	3-9-18	
Aroclor 1242	ND	0.050	EPA 8082A	3-9-18	3-9-18	
Aroclor 1248	ND	0.050	EPA 8082A	3-9-18	3-9-18	
Aroclor 1254	ND	0.050	EPA 8082A	3-9-18	3-9-18	
Aroclor 1260	ND	0.050	EPA 8082A	3-9-18	3-9-18	
<u> </u>	D (D	0 , 11: "	•		•	_

Surrogate: Percent Recovery Control Limits DCB 80 40-134

Analyte	Re	sult	Spike	Level	Source Result		rcent	Recovery Limits	RPD	RPD Limit	Flags
MATRIX SPIKES							-				
Laboratory ID:	03-0	36-01									
-	MS	MSD	MS	MSD		MS	MSD				
Aroclor 1260	0.373	0.385	0.500	0.500	ND	75	77	34-126	3	16	
Surrogate:											
DCB						75	73	40-134			

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B

Matrix: Soil

Offits. Trig/Ng (ppin)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-11_2.5					
Laboratory ID:	02-278-01					
Arsenic	ND	13	EPA 6010D	3-7-18	3-7-18	
Cadmium	ND	0.65	EPA 6010D	3-7-18	3-7-18	
Chromium	16	0.65	EPA 6010D	3-7-18	3-7-18	
Lead	15	6.5	EPA 6010D	3-7-18	3-7-18	
Mercury	ND	0.32	EPA 7471B	3-2-18	3-2-18	
Client ID:	B-11_12.5					
	02-278-04					
Laboratory ID: Arsenic	02-278-04 ND	12	EPA 6010D	3-7-18	3-7-18	
Cadmium	ND	0.60	EPA 6010D	3-7-18 3-7-18	3-7-18 3-7-18	
Chromium	13	0.60				
Lead	ND	6.0	EPA 6010D EPA 6010D	3-7-18 3-7-18	3-7-18 3-7-18	
Mercury	ND ND	0.30	EPA 7471B	3-7-16 3-2-18	3-7-16 3-2-18	
Mercury	ND	0.50	LIATATIO	J-Z-10	J-Z-10	
Client ID:	B-10_7.5					
Laboratory ID:	02-278-06					
Arsenic	ND	12	EPA 6010D	3-7-18	3-7-18	
Cadmium	ND	0.59	EPA 6010D	3-7-18	3-7-18	
Chromium	8.0	0.59	EPA 6010D	3-7-18	3-7-18	
Lead	ND	5.9	EPA 6010D	3-7-18	3-7-18	
Mercury	ND	0.29	EPA 7471B	3-2-18	3-2-18	
Client ID:	B-08_5					
Laboratory ID:	02-278-09					
Arsenic	ND	14	EPA 6010D	3-7-18	3-7-18	
Cadmium	ND	0.69	EPA 6010D	3-7-18	3-7-18	
Chromium	16	0.69	EPA 6010D	3-7-18	3-7-18	
Lead	6.9	6.9	EPA 6010D	3-7-18	3-7-18	
Mercury	ND	0.35	EPA 7471B	3-2-18	3-2-18	

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B

Matrix: Soil

Onito. Ing/itg (ppin)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-08_12.5					
Laboratory ID:	02-278-11					
Arsenic	ND	13	EPA 6010D	3-7-18	3-7-18	
Cadmium	ND	0.67	EPA 6010D	3-7-18	3-7-18	
Chromium	18	0.67	EPA 6010D	3-7-18	3-7-18	
Lead	ND	6.7	EPA 6010D	3-7-18	3-7-18	
Mercury	ND	0.34	EPA 7471B	3-2-18	3-2-18	
Client ID:	D 07 2 5					
	B-07_2.5					
Laboratory ID:	02-278-13					
Arsenic	ND	12	EPA 6010D	3-7-18	3-7-18	
Cadmium	ND	0.58	EPA 6010D	3-7-18	3-7-18	
Chromium	15	0.58	EPA 6010D	3-7-18	3-7-18	
Lead	6.0	5.8	EPA 6010D	3-7-18	3-7-18	
Mercury	ND	0.29	EPA 7471B	3-2-18	3-2-18	
Client ID:	B-07_7.5					
Laboratory ID:	02-278-14					
Arsenic	ND	12	EPA 6010D	3-7-18	3-7-18	
Cadmium	ND	0.62	EPA 6010D	3-7-18	3-7-18	
Chromium	14	0.62	EPA 6010D	3-7-18	3-7-18	
Lead	ND	6.2	EPA 6010D	3-7-18	3-7-18	
Mercury	ND	0.31	EPA 7471B	3-2-18	3-2-18	

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B QUALITY CONTROL

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0307S2					
Arsenic	ND	10	EPA 6010D	3-7-18	3-7-18	
Cadmium	ND	0.50	EPA 6010D	3-7-18	3-7-18	
Chromium	ND	0.50	EPA 6010D	3-7-18	3-7-18	
Lead	ND	5.0	EPA 6010D	3-7-18	3-7-18	
Laboratory ID:	MB0302S1					
Mercury	ND	0.25	EPA 7471B	3-2-18	3-2-18	

					Source	Pe	rcent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Rec	overy	Limits	RPD	Limit	Flags
DUPLICATE											
Laboratory ID:	03-03	36-07									
	ORIG	DUP									
Arsenic	ND	ND	NA	NA			NA	NA	NA	20	
Cadmium	ND	ND	NA	NA		ı	NA	NA	NA	20	
Chromium	20.8	18.0	NA	NA		ı	NA	NA	14	20	
Lead	8.15	7.35	NA	NA			NA	NA	10	20	
Laboratory ID:	02-20	65-04									
Mercury	ND	ND	NA	NA		l	NA	NA	NA	20	
MATRIX SPIKES											
Laboratory ID:	03-0	36-07									
	MS	MSD	MS	MSD		MS	MSD				
Arsenic	92.5	96.0	100	100	ND	93	96	75-125	4	20	
Cadmium	44.8	46.3	50.0	50.0	ND	90	93	75-125	3	20	
Chromium	108	113	100	100	20.8	87	92	75-125	5	20	
Lead	226	233	250	250	8.15	87	90	75-125	3	20	
Laboratory ID:	02-20	65-04									
Mercury	0.548	0.546	0.500	0.500	0.0400	102 101		80-120	0	20	

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-07_12.5					
Laboratory ID:	02-278-15					
Arsenic	ND	13	EPA 6010D	3-9-18	3-9-18	
Cadmium	ND	0.66	EPA 6010D	3-9-18	3-9-18	
Chromium	10	0.66	EPA 6010D	3-9-18	3-9-18	
Lead	ND	6.6	EPA 6010D	3-9-18	3-9-18	
Mercury	ND	0.33	EPA 7471B	3-12-18	3-12-18	

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B QUALITY CONTROL

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0309SM1					
Arsenic	ND	5.0	EPA 6010D	3-9-18	3-9-18	
Cadmium	ND	0.50	EPA 6010D	3-9-18	3-9-18	
Chromium	ND	0.50	EPA 6010D	3-9-18	3-9-18	
Lead	ND	5.0	EPA 6010D	3-9-18	3-9-18	
Laboratory ID:	MB0312S1					
Mercury	ND	0.25	EPA 7471B	3-12-18	3-12-18	

					Source	Pe	rcent	Recovery		RPD			
Analyte	Res	sult	Spike	Level	Result	Rec	overy	Limits	RPD	Limit	Flags		
DUPLICATE													
Laboratory ID:	03-00	07-11											
	ORIG	DUP									_		
Arsenic	ND	ND	NA	NA		ı	NA	NA	NA	20			
Cadmium	ND	ND	NA	NA		ı	NA	NA	NA	20			
Chromium	16.5	15.5	NA	NA		ı	NA	NA	6	20			
Lead	ND	ND	NA	NA		l	NA	NA	NA	20			
Laboratory ID:	03-0	75-02											
Mercury	ND	ND	NA	NA		ı	NA	NA	NA	20			
MATRIX SPIKES													
Laboratory ID:	03-0	07-11											
	MS	MSD	MS	MSD		MS	MSD						
Arsenic	101	99.5	100	100	ND	101	100	75-125	1	20			
Cadmium	46.0	46.9	50.0	50.0	ND	92 94				75-125	2	20	
Chromium	109	115	100	100	16.5	93	98	75-125	5	20			
Lead	230	236	250	250	ND	92	94	75-125	2	20			
Laboratory ID:	03-0	75-02											
Mercury	0.545	0.547	0.500	0.500	0.0159	106 106		80-120	0	20			

Project: 17-06520-000

% MOISTURE

Date Analyzed: 3-1&9-18

Client ID	Lab ID	% Moisture
B-11_2.5	02-278-01	23
B-11_12.5	02-278-04	16
B-10_7.5	02-278-06	15
B-08_5	02-278-09	28
B-08_12.5	02-278-11	25
B-07_2.5	02-278-13	14
B-07_7.5	02-278-14	19
B-07_12.5	02-278-15	24



Data Qualifiers and Abbreviations

- A Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
- B The analyte indicated was also found in the blank sample.
- C The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
- E The value reported exceeds the quantitation range and is an estimate.
- F Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
- H The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
- I Compound recovery is outside of the control limits.
- J The value reported was below the practical quantitation limit. The value is an estimate.
- K Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
- L The RPD is outside of the control limits.
- M Hydrocarbons in the gasoline range are impacting the diesel range result.
- M1 Hydrocarbons in the gasoline range (toluene-naphthalene) are present in the sample.
- N Hydrocarbons in the lube oil range are impacting the diesel range result.
- N1 Hydrocarbons in diesel range are impacting lube oil range results.
- O Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
- P The RPD of the detected concentrations between the two columns is greater than 40.
- Q Surrogate recovery is outside of the control limits.
- S Surrogate recovery data is not available due to the necessary dilution of the sample.
- T The sample chromatogram is not similar to a typical
- U The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- U1 The practical quantitation limit is elevated due to interferences present in the sample.
- V Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
- W Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
- X Sample extract treated with a mercury cleanup procedure.
- X1- Sample extract treated with a sulfuric acid/silica gel cleanup procedure.
- Y The calibration verification for this analyte exceeded the 20% drift specified in method 8260C, and therefore the reported result should be considered an estimate. The overall performance of the calibration verification standard met the acceptance criteria of the method.

7 -

ND - Not Detected at PQL

PQL - Practical Quantitation Limit

RPD - Relative Percent Difference





Chain of Custody

Page L of Z

Analytical Laboratory Testing Services 14648 NE 95th Street * Redmond, WA 98052	Turnaround Request (in working days)		L	abo	ratory	Numl	ber:				0	2.	-2	78			
Phone: (425) 883-3881 · www.onsite-env.com Company: Aspect Consulting Project Number: 1706520 - 000 Project Name: Pacific Right Bank Project Manager: Sampled by: 34F	(Check One) Same Day 1 Day 2 Days 3 Days Standard (7 Days) (TPH analysis 5 Days) (other) Date Time	Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX	NWTPH-Gx NWTPH-Dx (Acid / SG Clean-up)	Volatiles 8260C Halogenated Volatiles 8260C	EDB EPA 8011 (Waters Only)	Semivolatiles 8270D/SIM (with low-level PAHs) PAHS 8270D/SIM (low-level)	١	Organochlorine Pesticides 8081B	Organophosphorus Pesticides 8270D/SIM	Total RCRA Metals	Total MTCA Metals	TCLP Metals	HEM (oil and grease) 1664A		% Moisture
Lab ID Sample Identification	Sampled Sampled Matrix 2/26/18 8:56 50:\	三乙			N N	S E			1		ŏ 5	Tot			Ē		
2 B-11_2.5	2/26/18 9:04	T	A				-	-	X			+	A				
3 13-11_10	2/26/18 9:20	\forall	A				1										
4 13-11-12.5	2/26/18 9:25	\forall						•				1	№				0
5 B-10_2.5	2/26/10 10:34	П	A										1				
6 B-10_7.5	2/26/18 10:55				X				X				0				•
7 13-10-12.5	2/26/18 11:15		A										A				
8 13-10_117.5	2/26/18 11:30		A										N				
9 3-08-5	2/26/18 15:10		0					0					•				
10 B-08-75	2/24/18 15:15	4	•	- 40	6			-6	- na	>		-		- 40			-0
Relinquished Signature	Company			Date		Time	-/	Comme	nts/Sp	ecial I	nstruct	ions		No.			
Received National Relinquished	Aspect Consol active	T	2	2/2	28 18 'S 18	8:18	3:15 Added 2/28/18. DB (STA) & Added 3/8/18. DD (STA)										
Received															/		
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Received Reviewed/Date								Data Pa	ckage	: Star	ndard	□ Le	evel III		evel IV []	
	Reviewed/Date							Chroma	togran	ns with	final i	eport	Elec	ctronic	Data Deliv	verables (ED	DS)



Chain of Custody

Page 2 of 2

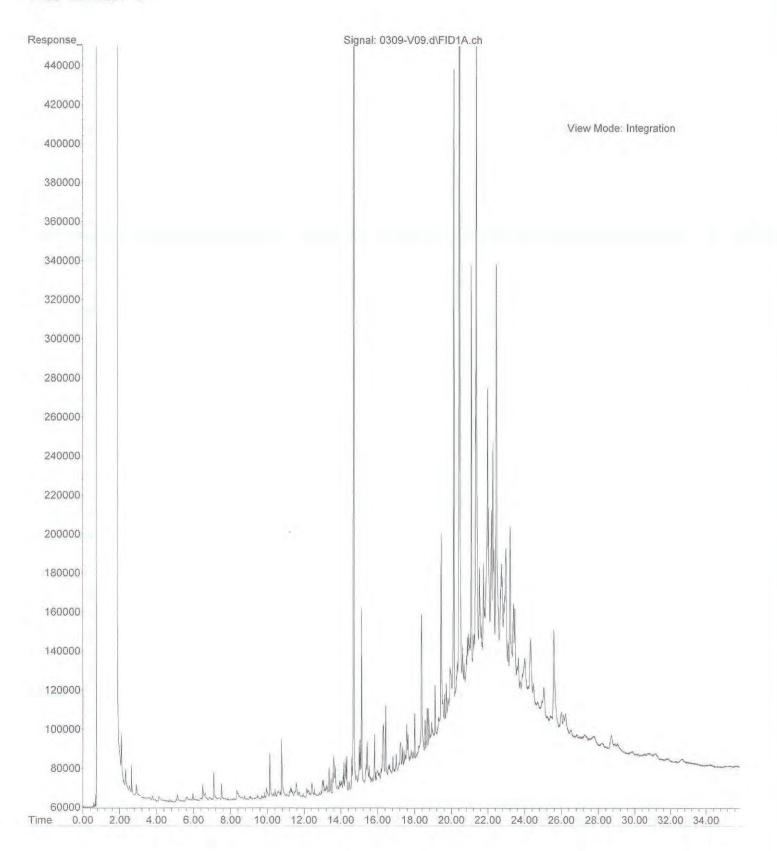
	Analytical Laboratory Testing Services 14648 NE 95th Street • Redmond, WA 98052 Phone: (425) 883-3881 • www.onsite-env.com		naround Re n working d			L	abo	orat	ory	Nur	nbe	er:						C	12	- 2	27	78				
	Bopact Consulting	Same Same Stand (TPH	1 2 2	1 Day 3 Days ays)	Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX	NWTPH-Gx	NWTPH-Dx (Acid / SG Clean-up)	Volatiles 8260C	Halogenated Volatiles 8260C	EDB EPA 8011 (Waters Only)	Semivolatiles 8270D/SIM (with low-level PAHs)	PAHS 8270D/SIM (low-level) CPAHS	PCBs 8082A	Organochlorine Pesticides 8081B	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	TCLP Metals	HEM (oil and grease) 1664A				% Mosture
11	13-08-12.5	2/26/18	12:50	50:1	2									•												
12	B-08_17.5	2/26/18	16:25		1	n														A						
13	B-07-2,5	2/22/18	8:50			•		(X					•	(X)											0
14	B-07-7.5	212718	8:55					((X)					1	X											•
15	B-07_125	2/24/8	9:05			A		((x)					(X)	\otimes					1	14	67				(X)
1le	B-07-17,5	2127/18	9:10		له	A														A						
	Signature		ompany		1		Date			Time			Con	nmen	ts/Sp	ecial	Instr	uctio	ns							
Relinquished Received	Tool of		Aspecd	Consul e ta	110	6	2/2	188	8	8	15															
Relinquishe	a a		16 261	E +20																						
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Reviewed/D	ate		Reviewed/D	ate									Chro	omato	ogran	ns wi	th fin	al rep	oort [Ele	ectron	ic Dat	a Deliv	/erable	s (EDD	s) 🗌

File :X:\DIESELS\VIGO\DATA\V180309\0309-V09.d Operator :JT

Operator : JT
Acquired : 9 Mar 2018 13:12 using AcqMethod V171020F.M
Instrument : Vigo

Sample Name: 02-278-01

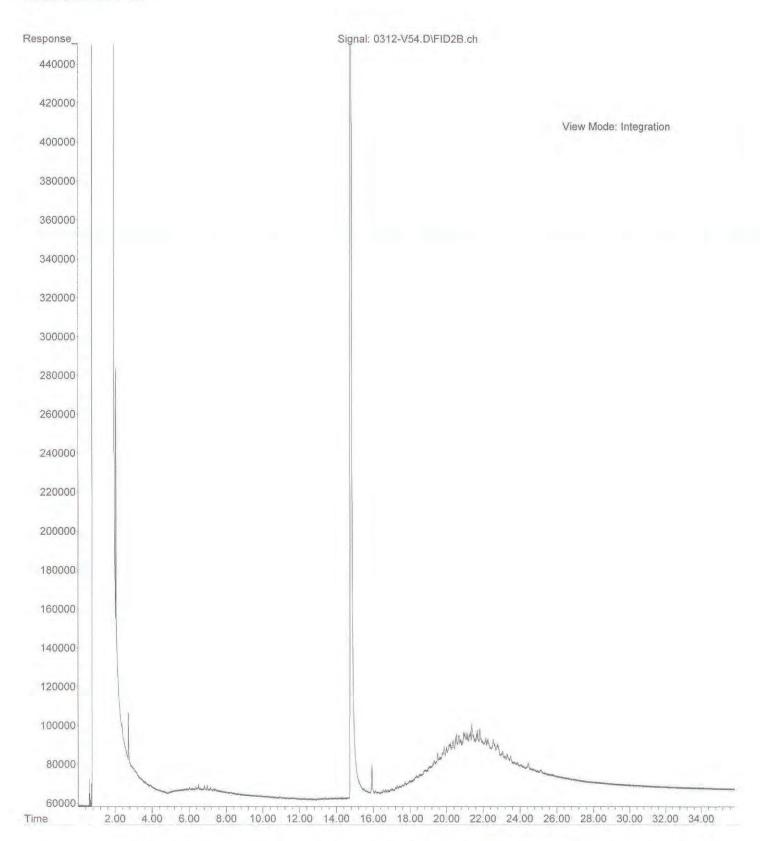
Misc Info : Vial Number: 9



File :X:\DIESELS\VIGO\DATA\V180312.SEC\0312-V54.D
Operator :JT

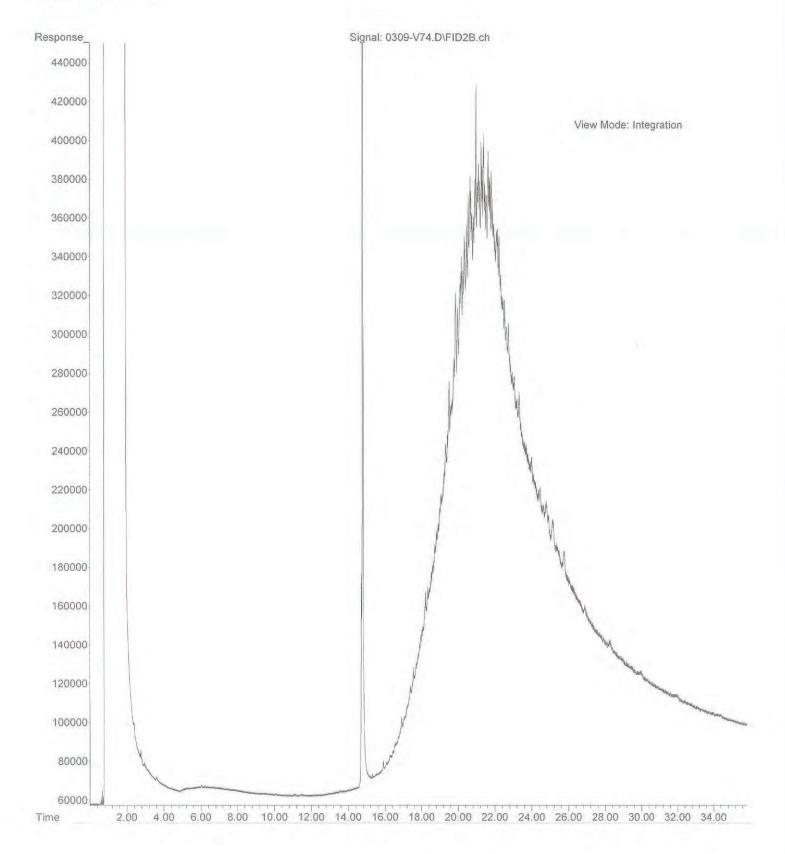
Operator : JT
Acquired : 12 Mar 2018 9:44
Instrument : Vigo using AcqMethod V171020F.M

Sample Name: 02-278-06 RR



File :X:\DIESELS\VIGO\DATA\V180309.SEC\0309-V74.D
Operator : JT
Acquired : 9 Mar 2018 23:29 using AcqMethod V171020F.M
Instrument : Vigo

Sample Name: 02-278-13 5X



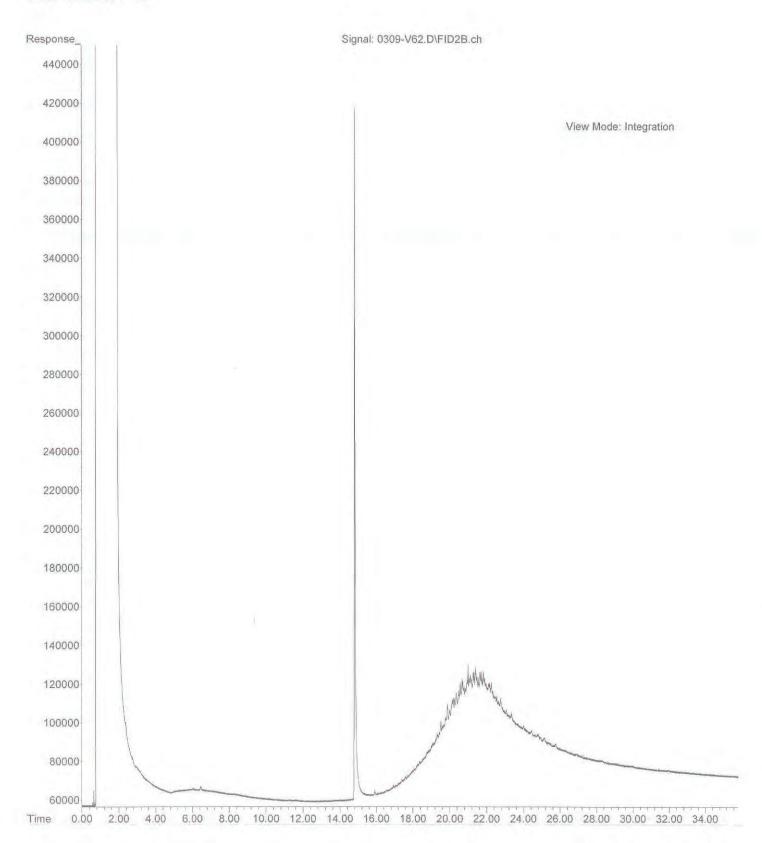
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Operator : JT

Acquired : 9 Mar 2018 15:32 using AcqMethod V171020F.M

Instrument : Vigo

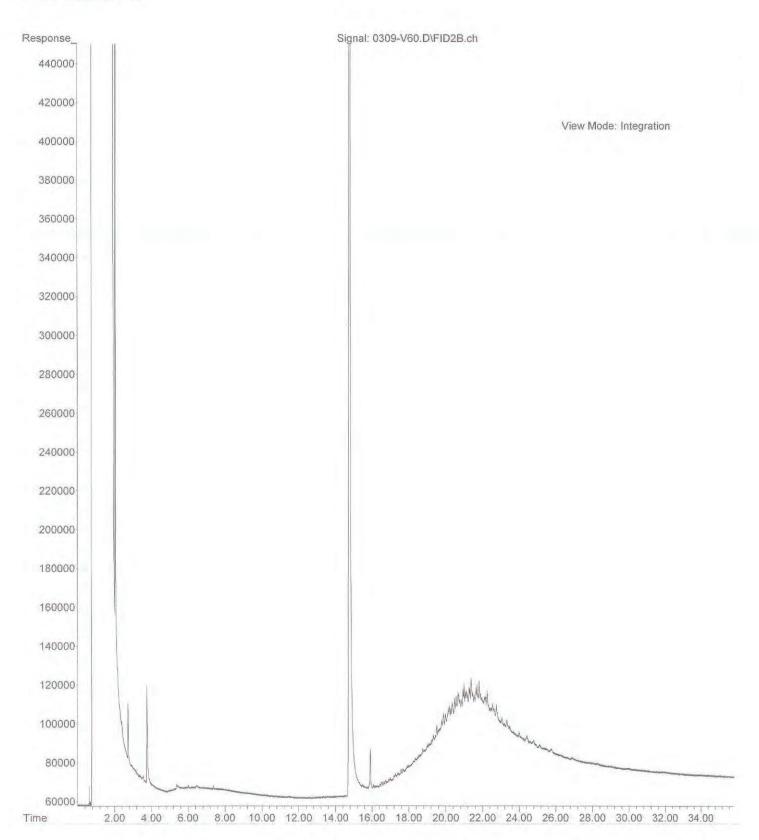
Sample Name: 02-278-14 10X



:X:\DIESELS\VIGO\DATA\V180309.SEC\0309-V60.D

File :X:\DIESELS\VIGO\DATA\V180309.SEC\0309-V60.D
Operator : JT
Acquired : 9 Mar 2018 13:52 using AcqMethod V171020F.M
Instrument : Vigo

Sample Name: 02-278-15 DUP





14648 NE 95th Street, Redmond, WA 98052 • (425) 883-3881

March 15, 2018

Bruce Carpenter Herrera Environmental Consultants, Inc. 2200 6th Avenue, Suite 1100 Seattle, WA 98121

Re: Analytical Data for Project 17-06520-000

Laboratory Reference No. 1803-007

Dear Bruce:

Enclosed are the analytical results and associated quality control data for samples submitted on March 1, 2018.

The standard policy of OnSite Environmental, Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

David Baumeister Project Manager

Enclosures

Project: 17-06520-000

Case Narrative

Samples were collected on February 28 and March 1, 2018 and received by the laboratory on March 1, 2018. They were maintained at the laboratory at a temperature of 2°C to 6°C.

Please note that any and all soil sample results are reported on a dry-weight basis, unless otherwise noted below.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

Project: 17-06520-000

NWTPH-HCID

Matrix: Soil

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP28-8	. ~=	mourou		7	ı iugo
Laboratory ID:	03-007-02					
Gasoline Range Organics	ND	25	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	63	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	130	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	108	50-150				
Client ID:	PP28-10					
Laboratory ID:	03-007-03					
Gasoline Range Organics	ND	25	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	62	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	120	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	111	50-150				
Client ID:	PP29-3					
Laboratory ID:	03-007-05					
Gasoline Range Organics	ND	450	NWTPH-HCID	3-2-18	3-5-18	
Diesel Range Organics	ND	1100	NWTPH-HCID	3-2-18	3-5-18	
Lube Oil	Detected	2300	NWTPH-HCID	3-2-18	3-5-18	
Surrogate:	Percent Recovery	Control Limits				•
o-Terphenyl		50-150				S
Client ID:	PP29-5					
Laboratory ID:	03-007-06					
Gasoline Range Organics	ND	26	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	65	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	130	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits	-	-		
o-Terphenyl	106	50-150				
Client ID:	PP29-10					
Laboratory ID:	03-007-07					
Laboratory ID.				0.0.40	0.0.40	
Gasoline Range Organics	ND	27	NWTPH-HCID	3-2-18	3-2-18	
	ND ND	27 66	NWTPH-HCID NWTPH-HCID	3-2-18 3-2-18	3-2-18 3-2-18	
Gasoline Range Organics						
Gasoline Range Organics Diesel Range Organics	ND	66	NWTPH-HCID	3-2-18	3-2-18	

Project: 17-06520-000

NWTPH-HCID

Matrix: Soil

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP32-4	. ~=	motriou	opa.oa	7a.y 20 a	. lugo
Laboratory ID:	03-007-09					
Gasoline Range Organics	ND	25	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	62	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	120	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits		0 2 .0	0 = .0	
o-Terphenyl	107	50-150				
· · · · · · · · · · · · · · · · · · ·						
Client ID:	PP32-7					
Laboratory ID:	03-007-10					
Gasoline Range Organics	ND	25	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	68	NWTPH-HCID	3-2-18	3-2-18	U1
Lube Oil	Detected	120	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	113	50-150				
Client ID:	PP31-3					
Laboratory ID:	03-007-13					
Gasoline Range Organics	ND	23	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND ND	57	NWTPH-HCID	3-2-18	3-2-18 3-2-18	
Lube Oil Range Organics	ND	110	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits	14W II II II IOID	0-2-10	0-2-10	
o-Terphenyl	112	50-150				
о тогрионут	7.72	00 100				
Client ID:	PP31-11					
Laboratory ID:	03-007-15					
Gasoline Range Organics	ND	25	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	62	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	130	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	112	50-150				
Client ID:	PP33-3					
Laboratory ID:	03-007-17					
Gasoline Range Organics	ND	450	NWTPH-HCID	3-2-18	3-5-18	
Diesel Range Organics	ND	1100	NWTPH-HCID	3-2-18	3-5-18	
Lube Oil	Detected	2300	NWTPH-HCID	3-2-18	3-5-18	
Surrogate:	Percent Recovery	Control Limits		0 2 10	0 0 10	
o-Terphenyl		50-150				S
o . orprioriyi		00 100				J

Project: 17-06520-000

NWTPH-HCID

Matrix: Soil

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP33-5	· QL	Motriou	Порагоа	Allalyzou	riugo
Laboratory ID:	03-007-18					
Gasoline Range Organics	ND	24	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	59	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	120	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	114	50-150				
Client ID:	PP33-10					
Laboratory ID:	03-007-19					
Gasoline Range Organics	ND	27	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	67	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	130	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	111	50-150				
011 - 4 15	DD00 0					
Client ID:	PP23-2					
Laboratory ID:	03-007-21					
Gasoline Range Organics	Detected	24	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	Detected	59	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil	Detected	120	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery 115	Control Limits 50-150				
o-Terphenyl	715	30-730				
Client ID:	PP23-5					
Laboratory ID:	03-007-22					
Gasoline Range Organics	ND	22	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	55	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	110	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	105	50-150				
Client ID:	PP23-10					
Laboratory ID:	03-007-23					
Gasoline Range Organics	ND	27	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	67	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	130	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	106	50-150				

Project: 17-06520-000

NWTPH-HCID

Matrix: Soil

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP23-15					
Laboratory ID:	03-007-24					
Gasoline Range Organics	ND	31	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	76	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	150	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	110	50-150				
Client ID:	PP21-2					
Laboratory ID:	03-007-25					
Gasoline Range Organics	ND	27	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	66	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil	Detected	130	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	106	50-150				
Client ID:	PP21-6					
Laboratory ID:	03-007-26	41	NWTPH-HCID	3-2-18	3-2-18	
Gasoline Range Organics	ND ND	4 I 100	NWTPH-HCID	3-2-18 3-2-18		U1
Diesel Range Organics					3-2-18 3-2-18	ΟI
Lube Oil Surrogate:	Detected Percent Recovery	210 Control Limits	NWTPH-HCID	3-2-18	3-2-10	
o-Terphenyl	105	50-150				
о-тегрпенуі	103	30-730				
Client ID:	PP21-10					
Laboratory ID:	03-007-27					
Gasoline Range Organics	ND	330	NWTPH-HCID	3-2-18	3-5-18	
Diesel Range Organics	ND	810	NWTPH-HCID	3-2-18	3-5-18	
Lube Oil	Detected	1600	NWTPH-HCID	3-2-18	3-5-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl		50-150				S
Client ID:	PP20-2					
Laboratory ID:	03-007-29					
Gasoline Range Organics	ND	23	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	58	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil	Detected	120	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	103	50-150				

Project: 17-06520-000

NWTPH-HCID

Matrix: Soil

Onits. Ing/Kg (ppin)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP20-5					
Laboratory ID:	03-007-30					
Gasoline Range Organics	ND	29	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	86	NWTPH-HCID	3-2-18	3-2-18	U1
Lube Oil	Detected	150	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	110	50-150				
Client ID:	PP20-10					
Laboratory ID:	03-007-31					
Gasoline Range Organics	ND	26	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	64	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil	Detected	130	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	111	50-150				
Client ID:	PP18-3					
Laboratory ID:	03-007-33					
Gasoline Range Organics	ND	25	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	61	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	120	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	104	50-150				
Client ID:	PP18-5					
Laboratory ID:	03-007-34					
Gasoline Range Organics	ND	40	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	99	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	200	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	107	50-150				
Client ID:	PP18-10					
Laboratory ID:	03-007-35					
Gasoline Range Organics	ND	27	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	67	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	130	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	105	50-150				
r - J						

Project: 17-06520-000

NWTPH-HCID

Matrix: Soil

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP17-1				<u>.</u>	
Laboratory ID:	03-007-37					
Gasoline Range Organics	ND	26	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	64	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	130	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	106	50-150				
Client ID:	PP17-5					
Laboratory ID:	03-007-38					
Gasoline Range Organics	ND	24	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	61	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	120	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	106	50-150				
Client ID:	PP17-10					
Laboratory ID:	03-007-39					
Gasoline Range Organics	ND	23	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	58	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	120	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	105	50-150				
Client ID:	PP16-1					
Laboratory ID:	03-007-41					
Gasoline Range Organics	ND	210	NWTPH-HCID	3-2-18	3-5-18	
Diesel Range Organics	ND	530	NWTPH-HCID	3-2-18	3-5-18	
Lube Oil	Detected	1100	NWTPH-HCID	3-2-18	3-5-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl		50-150				S
, ,						
Client ID:	PP16-11					
Laboratory ID:	03-007-42					
Gasoline Range Organics	ND	25	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	61	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	120	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	103	50-150				

Project: 17-06520-000

NWTPH-HCID

Matrix: Soil

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP14-5				<u>.</u>	90
Laboratory ID:	03-007-45					
Gasoline Range Organics	ND	25	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	62	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	130	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits	-			
o-Terphenyl	108	50-150				
Client ID:	PP14-12					
Laboratory ID:	03-007-46					
Gasoline Range Organics	ND	26	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	65	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	130	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	104	50-150				
Client ID:	PP11-5					
Laboratory ID:	03-007-49					
Gasoline Range Organics	ND	26	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	66	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	130	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	104	50-150				
Client ID:	PP11-10					
Laboratory ID:	03-007-50					
Gasoline Range Organics	ND	25	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND ND	63	NWTPH-HCID	3-2-18	3-2-18 3-2-18	
Lube Oil Range Organics	ND	130	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits	14W II II-IIOID	J-Z-10	J-Z-10	
o-Terphenyl	106	50-150				
о-тегрпенуі	700	30-130				
Client ID:	PP10-11					
Laboratory ID:	03-007-52					
Gasoline Range Organics	ND	33	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	82	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	170	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits		-	-	
o-Terphenyl	101	50-150				
, ,						

Project: 17-06520-000

NWTPH-HCID

Matrix: Soil

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP10-17	. ~=	motriou		7 11 a. 1 2 a a	ı iugo
Laboratory ID:	03-007-53					
Gasoline Range Organics	ND	25	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	62	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	120	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	99	50-150				
Client ID:	PP27-7					
Laboratory ID:	03-007-55					
Gasoline Range Organics	ND	29	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	73	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	150	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	106	50-150				
Client ID:	PP27-10					
Laboratory ID:	03-007-56					
Gasoline Range Organics	ND	24	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	61	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	120	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	102	50-150				
Client ID:	PP30-5					
Laboratory ID:	03-007-59					
Gasoline Range Organics	ND	26	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND Data at a d	64	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	Detected	130	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	98	50-150				
Client ID:	PP30-10					
Laboratory ID:	03-007-60					
Gasoline Range Organics	ND	23	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND ND	23 57	NWTPH-HCID	3-2-16 3-2-18	3-2-18 3-2-18	
Lube Oil Range Organics	ND ND	110	NWTPH-HCID	3-2-16 3-2-18	3-2-18 3-2-18	
Surrogate:	Percent Recovery	Control Limits	1444 II II-IICID	J-Z-10	J-Z-10	
o-Terphenyl	101	50-150				
o-i erprierryi	101	30-130				

Project: 17-06520-000

NWTPH-HCID

Matrix: Soil

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP24-7	1 42	Mictriou	Troparca	Analyzea	i iugo
Laboratory ID:	03-007-63					
Gasoline Range Organics	Detected	280	NWTPH-HCID	3-2-18	3-5-18	
Diesel Range Organics	ND	690	NWTPH-HCID	3-2-18	3-5-18	
Lube Oil	Detected	1400	NWTPH-HCID	3-2-18	3-5-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl		50-150				S
Client ID:	PP24-10					
Laboratory ID:	03-007-64					
Gasoline Range Organics	ND	23	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	58	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	120	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	104	50-150				

Project: 17-06520-000

NWTPH-HCID QUALITY CONTROL

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0302S2					
Gasoline Range Organics	ND	20	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	50	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	100	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	106	50-150				
Laboratory ID:	MB0302S3					
Gasoline Range Organics	ND	20	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	50	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	100	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	105	50-150				
Laboratory ID:	MB0302S4					
Gasoline Range Organics	ND	20	NWTPH-HCID	3-2-18	3-2-18	
Diesel Range Organics	ND	50	NWTPH-HCID	3-2-18	3-2-18	
Lube Oil Range Organics	ND	100	NWTPH-HCID	3-2-18	3-2-18	
Surrogate:	Percent Recovery	Control Limits	·	·		
o-Terphenyl	107	50-150				

Project: 17-06520-000

NWTPH-Gx/BTEX

Matrix: Soil

Analyte Result PQL Method Client ID: PP32-10 Laboratory ID: 03-007-11	l Prepared Analyzed Flaç
Laboratory ID: 03_007_11	
Laboratory ID. 03-007-11	
Benzene ND 0.026 EPA 8021	1B 3-9-18 3-9-18
Toluene ND 0.13 EPA 8021	1B 3-9-18 3-9-18
Ethyl Benzene ND 0.13 EPA 8021	1B 3-9-18 3-9-18
m,p-Xylene ND 0.26 EPA 8021	1B 3-9-18 3-9-18
o-Xylene ND 0.13 EPA 8021	1B 3-9-18 3-9-18
Gasoline ND 13 NWTPH-0	Gx 3-9-18 3-9-18
Surrogate: Percent Recovery Control Limits	
Fluorobenzene 87 66-130	
Client ID: PP33-3	
Laboratory ID: 03-007-17	
Benzene ND 0.024 EPA 8021	IB 3-9-18 3-9-18
Toluene ND 0.12 EPA 8021	1B 3-9-18 3-9-18
Ethyl Benzene ND 0.12 EPA 8021	1B 3-9-18 3-9-18
m,p-Xylene ND 0.24 EPA 8021	IB 3-9-18 3-9-18
o-Xylene ND 0.12 EPA 8021	1B 3-9-18 3-9-18
Gasoline ND 12 NWTPH-0	Gx 3-9-18 3-9-18
Surrogate: Percent Recovery Control Limits	
Fluorobenzene 85 66-130	
Client ID: PP23-2	
Laboratory ID: 03-007-21	
Benzene ND 0.025 EPA 8021	IB 3-9-18 3-9-18
Toluene ND 0.13 EPA 8021	1B 3-9-18 3-9-18
Ethyl Benzene ND 0.13 EPA 8021	IB 3-9-18 3-9-18
m,p-Xylene ND 0.25 EPA 8021	IB 3-9-18 3-9-18
o-Xylene ND 0.13 EPA 8021	IB 3-9-18 3-9-18
Gasoline ND 13 NWTPH-0	Gx 3-9-18 3-9-18
Surrogate: Percent Recovery Control Limits	
Fluorobenzene 104 66-130	

Project: 17-06520-000

NWTPH-Gx/BTEX

Matrix: Soil

0 0 (11)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP21-10					
Laboratory ID:	03-007-27					
Benzene	ND	0.024	EPA 8021B	3-9-18	3-9-18	
Toluene	ND	0.12	EPA 8021B	3-9-18	3-9-18	
Ethyl Benzene	ND	0.12	EPA 8021B	3-9-18	3-9-18	
m,p-Xylene	ND	0.24	EPA 8021B	3-9-18	3-9-18	
o-Xylene	ND	0.12	EPA 8021B	3-9-18	3-9-18	
Gasoline	39	12	NWTPH-Gx	3-9-18	3-9-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	98	66-130				
Client ID:	PP16-1					
Laboratory ID:	03-007-41					
Benzene	ND	0.020	EPA 8021B	3-9-18	3-9-18	
Toluene	ND	0.053	EPA 8021B	3-9-18	3-9-18	
Ethyl Benzene	ND	0.053	EPA 8021B	3-9-18	3-9-18	
m,p-Xylene	ND	0.11	EPA 8021B	3-9-18	3-9-18	
o-Xylene	ND	0.053	EPA 8021B	3-9-18	3-9-18	
Gasoline	ND	5.3	NWTPH-Gx	3-9-18	3-9-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	84	66-130				
Client ID:	PP24-7					
Laboratory ID:	03-007-63					
Benzene	ND	0.020	EPA 8021B	3-9-18	3-9-18	
Toluene	ND	0.075	EPA 8021B	3-9-18	3-9-18	
Ethyl Benzene	ND	0.075	EPA 8021B	3-9-18	3-9-18	
m,p-Xylene	ND	0.15	EPA 8021B	3-9-18	3-9-18	
o-Xylene	ND	0.075	EPA 8021B	3-9-18	3-9-18	
Gasoline	ND	7.5	NWTPH-Gx	3-9-18	3-9-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	103	66-130				

Project: 17-06520-000

NWTPH-Gx/BTEX QUALITY CONTROL

Matrix: Soil

Units: mg/kg (ppm)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0309S1					
Benzene	ND	0.020	EPA 8021B	3-9-18	3-9-18	
Toluene	ND	0.050	EPA 8021B	3-9-18	3-9-18	
Ethyl Benzene	ND	0.050	EPA 8021B	3-9-18	3-9-18	
m,p-Xylene	ND	0.10	EPA 8021B	3-9-18	3-9-18	
o-Xylene	ND	0.050	EPA 8021B	3-9-18	3-9-18	
Gasoline	ND	5.0	NWTPH-Gx	3-9-18	3-9-18	

Surrogate: Percent Recovery Control Limits Fluorobenzene 87 66-130

					Source	Pei	rcent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Rec	overy	Limits	RPD	Limit	Flags
DUPLICATE											
Laboratory ID:	03-07	73-01									
	ORIG	DUP									
Benzene	ND	ND	NA	NA		1	NΑ	NA	NA	30	
Toluene	ND	ND	NA	NA		1	NΑ	NA	NA	30	
Ethyl Benzene	ND	ND	NA	NA		1	NΑ	NA	NA	30	
m,p-Xylene	ND	ND	NA	NA		1	NΑ	NA	NA	30	
o-Xylene	ND	ND	NA	NA		1	NΑ	NA	NA	30	
Gasoline	ND	ND	NA	NA		1	NΑ	NA	NA	30	
Surrogate:											
Fluorobenzene						97	94	66-130			
SPIKE BLANKS											
Laboratory ID:	SB03	09S1									
	SB	SBD	SB	SBD		SB	SBD				
Benzene	0.961	0.980	1.00	1.00		96	98	70-120	2	11	
Toluene	0.980	0.991	1.00	1.00		98	99	73-121	1	14	
Ethyl Benzene	0.983	0.999	1.00	1.00		98	100	74-121	2	11	
m,p-Xylene	1.01	1.02	1.00	1.00		101	102	75-124	1	13	
o-Xylene	0.959	0.974	1.00	1.00		96	97	75-121	2	12	
Surrogate:	•	•	•			•	•		•	•	•
Fluorobenzene						93	94	66-130			

Project: 17-06520-000

NWTPH-Dx

Matrix: Soil

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	PP29-3			•	-	
Laboratory ID:	03-007-05					
Diesel Range Organics	ND	280	NWTPH-Dx	3-8-18	3-9-18	
Lube Oil	5900	560	NWTPH-Dx	3-8-18	3-9-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl		50-150				S
OII ID	DD00 7					
Client ID:	PP32-7					
Laboratory ID:	03-007-10	20	ANATOLI D	0.0.40	0.0.40	114
Diesel Range Organics	ND	60	NWTPH-Dx	3-8-18	3-8-18	U1
Lube Oil	650	61	NWTPH-Dx	3-8-18	3-8-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	95	50-150				
Client ID:	PP33-3					
Laboratory ID:	03-007-17					
Diesel Range Organics	ND	750	NWTPH-Dx	3-8-18	3-9-18	U1
Lube Oil	12000	1100	NWTPH-Dx	3-8-18	3-9-18	0.
Surrogate:	Percent Recovery	Control Limits	TWW THE DX	0 0 10	0 0 10	
o-Terphenyl		50-150				S
о тогриону.		00 /00				
Client ID:	PP23-2					
Laboratory ID:	03-007-21					
Diesel Range Organics	ND	29	NWTPH-Dx	3-8-18	3-8-18	
Lube Oil Range Organics	200	59	NWTPH-Dx	3-8-18	3-8-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	90	50-150				
011 / ID	2224					
Client ID:	PP21-2					
Laboratory ID:	03-007-25					
Diesel Range Organics	57	33	NWTPH-Dx	3-8-18	3-8-18	N
Lube Oil Range Organics	540	66	NWTPH-Dx	3-8-18	3-8-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	108	50-150				
Client ID:	PP21-6					
Laboratory ID:	03-007-26					
Diesel Range Organics	150	51	NWTPH-Dx	3-8-18	3-8-18	N
Lube Oil Range Organics	960	100	NWTPH-Dx	3-8-18	3-8-18	14
Surrogate:	Percent Recovery	Control Limits	TANA IL LI-DV	0-0-10	0-0-10	
o-Terphenyl	103	50-150				
υ- ι σιριισιιγι	103	JU-130				

Project: 17-06520-000

NWTPH-Dx

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP21-10					
Laboratory ID:	03-007-27					
Diesel Range Organics	1800	410	NWTPH-Dx	3-8-18	3-9-18	
Lube Oil Range Organics	10000	810	NWTPH-Dx	3-8-18	3-9-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl		50-150				S
Client ID:	PP21-15					
Laboratory ID:	03-007-28					
Diesel Range Organics	110	68	NWTPH-Dx	3-8-18	3-8-18	
Lube Oil Range Organics	1200	140	NWTPH-Dx	3-8-18	3-8-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	107	50-150				
Client ID:	PP20-2					
	03-007-29					
Laboratory ID:		00	NIM/TOLL Dec	0.0.40	0.0.40	
Diesel Range Organics	ND	29 58	NWTPH-Dx	3-8-18	3-8-18	
Lube Oil	380	Control Limits	NWTPH-Dx	3-8-18	3-8-18	
Surrogate:	Percent Recovery 92	50-150				
o-Terphenyl	92	50-150				
Client ID:	PP20-5					
Laboratory ID:	03-007-30					
Diesel Range Organics	ND	200	NWTPH-Dx	3-8-18	3-9-18	U1
Lube Oil	1300	370	NWTPH-Dx	3-8-18	3-9-18	.
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	103	50-150				
Client ID:	PP20-10					
Laboratory ID:	03-007-31					
Diesel Range Organics	ND	32	NWTPH-Dx	3-8-18	3-8-18	
Lube Oil	95	64	NWTPH-Dx	3-8-18	3-8-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	84	50-150				
Client ID:	PP20-15					
Laboratory ID:	03-007-32					
Diesel Range Organics	ND	58	NWTPH-Dx	3-8-18	3-13-18	U1
Lube Oil	320	70	NWTPH-Dx	3-8-18	3-13-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	97	50-150				

Project: 17-06520-000

NWTPH-Dx

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP16-1					
Laboratory ID:	03-007-41					
Diesel Range Organics	ND	320	NWTPH-Dx	3-8-18	3-9-18	U1
Lube Oil	3200	530	NWTPH-Dx	3-8-18	3-9-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl		50-150				S
Client ID:	PP30-5					
Laboratory ID:	03-007-59					
Diesel Range Organics	71	32	NWTPH-Dx	3-8-18	3-8-18	N
Lube Oil	630	64	NWTPH-Dx	3-8-18	3-8-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	91	50-150				
Client ID:	PP24-7					
Laboratory ID:	03-007-63					
Diesel Range Organics	ND	35	NWTPH-Dx	3-8-18	3-12-18	
Lube Oil	130	69	NWTPH-Dx	3-8-18	3-12-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	91	50-150				

Project: 17-06520-000

NWTPH-Dx QUALITY CONTROL

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0308S1					
Diesel Range Organics	ND	25	NWTPH-Dx	3-8-18	3-8-18	
Lube Oil Range Organics	ND	50	NWTPH-Dx	3-8-18	3-8-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	99	50-150				

					Source	Perce	ent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Recov	ery	Limits	RPD	Limit	Flags
DUPLICATE											
Laboratory ID:	03-00	7-25									
	ORIG	DUP									
Diesel Range Organics	43.0	29.9	NA	NA		NA		NA	36	NA	N
Lube Oil Range Organics	406	369	NA	NA		NA		NA	10	NA	
Surrogate:											
o-Terphenyl						108	86	50-150			
Laboratory ID:	03-00	7-32									
	ORIG	DUP									
Diesel Range	ND	ND	NA	NA		NA		NA	NA	NA	U1
Lube Oil	230	203	NA	NA		NA		NA	12	NA	
Surrogate: o-Terphenyl						97	79	50-150			

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP28-8					
Laboratory ID:	03-007-02					
Benzo[a]anthracene	ND	0.0083	EPA 8270D/SIM	3-5-18	3-5-18	
Chrysene	ND	0.0083	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[b]fluoranthene	ND	0.0083	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo(j,k)fluoranthene	ND	0.0083	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[a]pyrene	ND	0.0083	EPA 8270D/SIM	3-5-18	3-5-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0083	EPA 8270D/SIM	3-5-18	3-5-18	
Dibenz[a,h]anthracene	ND	0.0083	EPA 8270D/SIM	3-5-18	3-5-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	74	32 - 115				
Pyrene-d10	92	35 - 129				
Terphenyl-d14	94	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP28-10					
Laboratory ID:	03-007-03					
Benzo[a]anthracene	0.011	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
Chrysene	0.011	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[b]fluoranthene	0.0084	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo(j,k)fluoranthene	ND	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[a]pyrene	0.010	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
Dibenz[a,h]anthracene	ND	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	82	32 - 115				
Pyrene-d10	87	35 - 129				
Terphenyl-d14	90	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP29-3					
Laboratory ID:	03-007-05					
Benzo[a]anthracene	ND	0.038	EPA 8270D/SIM	3-5-18	3-5-18	
Chrysene	ND	0.038	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[b]fluoranthene	ND	0.038	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo(j,k)fluoranthene	ND	0.038	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[a]pyrene	ND	0.038	EPA 8270D/SIM	3-5-18	3-5-18	
Indeno(1,2,3-c,d)pyrene	ND	0.038	EPA 8270D/SIM	3-5-18	3-5-18	
Dibenz[a,h]anthracene	ND	0.038	EPA 8270D/SIM	3-5-18	3-5-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	69	32 - 115				
Pyrene-d10	77	35 - 129				
Terphenyl-d14	72	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP29-5					
Laboratory ID:	03-007-06					
Benzo[a]anthracene	ND	0.0087	EPA 8270D/SIM	3-5-18	3-5-18	
Chrysene	ND	0.0087	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[b]fluoranthene	ND	0.0087	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo(j,k)fluoranthene	ND	0.0087	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[a]pyrene	ND	0.0087	EPA 8270D/SIM	3-5-18	3-5-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0087	EPA 8270D/SIM	3-5-18	3-5-18	
Dibenz[a,h]anthracene	ND	0.0087	EPA 8270D/SIM	3-5-18	3-5-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	84	32 - 115				
Pyrene-d10	90	35 - 129				
Terphenyl-d14	98	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP29-10					
Laboratory ID:	03-007-07					
Benzo[a]anthracene	ND	0.0088	EPA 8270D/SIM	3-5-18	3-6-18	
Chrysene	ND	0.0088	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[b]fluoranthene	ND	0.0088	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo(j,k)fluoranthene	ND	0.0088	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[a]pyrene	ND	0.0088	EPA 8270D/SIM	3-5-18	3-6-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0088	EPA 8270D/SIM	3-5-18	3-6-18	
Dibenz[a,h]anthracene	ND	0.0088	EPA 8270D/SIM	3-5-18	3-6-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	81	32 - 115				
Pyrene-d10	88	35 - 129				
Terphenyl-d14	92	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP32-4					
Laboratory ID:	03-007-09					
Benzo[a]anthracene	0.010	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
Chrysene	0.015	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[b]fluoranthene	0.024	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo(j,k)fluoranthene	ND	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[a]pyrene	0.011	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
Indeno(1,2,3-c,d)pyrene	0.0095	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
Dibenz[a,h]anthracene	ND	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	74	32 - 115				
Pyrene-d10	81	35 - 129				
Terphenyl-d14	83	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP32-7					
Laboratory ID:	03-007-10					
Benzo[a]anthracene	0.038	0.0082	EPA 8270D/SIM	3-5-18	3-7-18	
Chrysene	0.055	0.0082	EPA 8270D/SIM	3-5-18	3-7-18	
Benzo[b]fluoranthene	0.051	0.0082	EPA 8270D/SIM	3-5-18	3-7-18	
Benzo(j,k)fluoranthene	0.013	0.0082	EPA 8270D/SIM	3-5-18	3-7-18	
Benzo[a]pyrene	0.036	0.0082	EPA 8270D/SIM	3-5-18	3-7-18	
Indeno(1,2,3-c,d)pyrene	0.025	0.0082	EPA 8270D/SIM	3-5-18	3-7-18	
Dibenz[a,h]anthracene	0.0091	0.0082	EPA 8270D/SIM	3-5-18	3-7-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	79	32 - 115				
Pyrene-d10	91	35 - 129				
Terphenyl-d14	86	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP31-3					
Laboratory ID:	03-007-13					
Benzo[a]anthracene	ND	0.0075	EPA 8270D/SIM	3-5-18	3-6-18	
Chrysene	ND	0.0075	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[b]fluoranthene	0.0093	0.0075	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo(j,k)fluoranthene	ND	0.0075	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[a]pyrene	ND	0.0075	EPA 8270D/SIM	3-5-18	3-6-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0075	EPA 8270D/SIM	3-5-18	3-6-18	
Dibenz[a,h]anthracene	ND	0.0075	EPA 8270D/SIM	3-5-18	3-6-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	83	32 - 115				
Pyrene-d10	90	35 - 129				
Terphenyl-d14	92	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP31-11					
Laboratory ID:	03-007-15					
Benzo[a]anthracene	ND	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
Chrysene	ND	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[b]fluoranthene	ND	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo(j,k)fluoranthene	ND	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[a]pyrene	ND	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
Dibenz[a,h]anthracene	ND	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	85	32 - 115				
Pyrene-d10	95	35 - 129				
Terphenyl-d14	100	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP33-3					
Laboratory ID:	03-007-17					
Benzo[a]anthracene	ND	0.075	EPA 8270D/SIM	3-5-18	3-7-18	
Chrysene	0.19	0.075	EPA 8270D/SIM	3-5-18	3-7-18	
Benzo[b]fluoranthene	0.14	0.075	EPA 8270D/SIM	3-5-18	3-7-18	
Benzo(j,k)fluoranthene	ND	0.075	EPA 8270D/SIM	3-5-18	3-7-18	
Benzo[a]pyrene	ND	0.075	EPA 8270D/SIM	3-5-18	3-7-18	
Indeno(1,2,3-c,d)pyrene	ND	0.075	EPA 8270D/SIM	3-5-18	3-7-18	
Dibenz[a,h]anthracene	ND	0.075	EPA 8270D/SIM	3-5-18	3-7-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	71	32 - 115				
Pyrene-d10	94	35 - 129				
Terphenyl-d14	82	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP33-5					
Laboratory ID:	03-007-18					
Benzo[a]anthracene	ND	0.0079	EPA 8270D/SIM	3-5-18	3-6-18	
Chrysene	ND	0.0079	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[b]fluoranthene	ND	0.0079	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo(j,k)fluoranthene	ND	0.0079	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[a]pyrene	ND	0.0079	EPA 8270D/SIM	3-5-18	3-6-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0079	EPA 8270D/SIM	3-5-18	3-6-18	
Dibenz[a,h]anthracene	ND	0.0079	EPA 8270D/SIM	3-5-18	3-6-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	<i>7</i> 5	32 - 115				
Pyrene-d10	84	35 - 129				
Terphenyl-d14	81	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP33-10					
Laboratory ID:	03-007-19					
Benzo[a]anthracene	ND	0.0090	EPA 8270D/SIM	3-5-18	3-6-18	
Chrysene	ND	0.0090	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[b]fluoranthene	ND	0.0090	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo(j,k)fluoranthene	ND	0.0090	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[a]pyrene	ND	0.0090	EPA 8270D/SIM	3-5-18	3-6-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0090	EPA 8270D/SIM	3-5-18	3-6-18	
Dibenz[a,h]anthracene	ND	0.0090	EPA 8270D/SIM	3-5-18	3-6-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	74	32 - 115				
Pyrene-d10	90	35 - 129				
Terphenyl-d14	93	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP23-2					
Laboratory ID:	03-007-21					
Benzo[a]anthracene	0.012	0.0078	EPA 8270D/SIM	3-5-18	3-6-18	
Chrysene	0.017	0.0078	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[b]fluoranthene	0.017	0.0078	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo(j,k)fluoranthene	ND	0.0078	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[a]pyrene	0.011	0.0078	EPA 8270D/SIM	3-5-18	3-6-18	
Indeno(1,2,3-c,d)pyrene	0.0082	0.0078	EPA 8270D/SIM	3-5-18	3-6-18	
Dibenz[a,h]anthracene	ND	0.0078	EPA 8270D/SIM	3-5-18	3-6-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	<i>7</i> 5	32 - 115				
Pyrene-d10	85	35 - 129				
Terphenyl-d14	85	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP23-5					
Laboratory ID:	03-007-22					
Benzo[a]anthracene	0.18	0.015	EPA 8270D/SIM	3-5-18	3-7-18	
Chrysene	0.044	0.015	EPA 8270D/SIM	3-5-18	3-7-18	
Benzo[b]fluoranthene	0.10	0.015	EPA 8270D/SIM	3-5-18	3-7-18	
Benzo(j,k)fluoranthene	ND	0.015	EPA 8270D/SIM	3-5-18	3-7-18	
Benzo[a]pyrene	0.044	0.015	EPA 8270D/SIM	3-5-18	3-7-18	
Indeno(1,2,3-c,d)pyrene	0.026	0.015	EPA 8270D/SIM	3-5-18	3-7-18	
Dibenz[a,h]anthracene	0.019	0.015	EPA 8270D/SIM	3-5-18	3-7-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	84	32 - 115				
Pyrene-d10	98	35 - 129				
Terphenyl-d14	79	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP23-10					
Laboratory ID:	03-007-23					
Benzo[a]anthracene	ND	0.0090	EPA 8270D/SIM	3-5-18	3-6-18	
Chrysene	ND	0.0090	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[b]fluoranthene	ND	0.0090	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo(j,k)fluoranthene	ND	0.0090	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[a]pyrene	ND	0.0090	EPA 8270D/SIM	3-5-18	3-6-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0090	EPA 8270D/SIM	3-5-18	3-6-18	
Dibenz[a,h]anthracene	ND	0.0090	EPA 8270D/SIM	3-5-18	3-6-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	78	32 - 115				
Pyrene-d10	83	35 - 129				
Terphenyl-d14	87	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP23-15					
Laboratory ID:	03-007-24					
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	3-5-18	3-6-18	
Chrysene	ND	0.010	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	3-5-18	3-6-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	3-5-18	3-6-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	3-5-18	3-6-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	85	32 - 115				
Pyrene-d10	92	35 - 129				
Terphenyl-d14	94	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP21-2					
Laboratory ID:	03-007-25					
Benzo[a]anthracene	0.076	0.0089	EPA 8270D/SIM	3-5-18	3-7-18	
Chrysene	0.097	0.0089	EPA 8270D/SIM	3-5-18	3-7-18	
Benzo[b]fluoranthene	0.11	0.0089	EPA 8270D/SIM	3-5-18	3-7-18	
Benzo(j,k)fluoranthene	0.033	0.0089	EPA 8270D/SIM	3-5-18	3-7-18	
Benzo[a]pyrene	0.062	0.0089	EPA 8270D/SIM	3-5-18	3-7-18	
Indeno(1,2,3-c,d)pyrene	0.040	0.0089	EPA 8270D/SIM	3-5-18	3-7-18	
Dibenz[a,h]anthracene	0.0097	0.0089	EPA 8270D/SIM	3-5-18	3-7-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	79	32 - 115				
Pyrene-d10	95	35 - 129				
Terphenyl-d14	87	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP21-6					
Laboratory ID:	03-007-26					
Benzo[a]anthracene	0.14	0.014	EPA 8270D/SIM	3-5-18	3-7-18	
Chrysene	0.18	0.014	EPA 8270D/SIM	3-5-18	3-7-18	
Benzo[b]fluoranthene	0.22	0.014	EPA 8270D/SIM	3-5-18	3-7-18	
Benzo(j,k)fluoranthene	0.077	0.014	EPA 8270D/SIM	3-5-18	3-7-18	
Benzo[a]pyrene	0.13	0.014	EPA 8270D/SIM	3-5-18	3-7-18	
Indeno(1,2,3-c,d)pyrene	0.11	0.014	EPA 8270D/SIM	3-5-18	3-7-18	
Dibenz[a,h]anthracene	0.024	0.014	EPA 8270D/SIM	3-5-18	3-7-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	72	32 - 115				
Pyrene-d10	88	35 - 129				
Terphenyl-d14	82	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP21-10					
Laboratory ID:	03-007-27					
Benzo[a]anthracene	0.90	0.011	EPA 8270D/SIM	3-5-18	3-7-18	
Chrysene	1.0	0.011	EPA 8270D/SIM	3-5-18	3-7-18	
Benzo[b]fluoranthene	1.3	0.011	EPA 8270D/SIM	3-5-18	3-7-18	
Benzo(j,k)fluoranthene	0.30	0.011	EPA 8270D/SIM	3-5-18	3-7-18	
Benzo[a]pyrene	0.85	0.011	EPA 8270D/SIM	3-5-18	3-7-18	
Indeno(1,2,3-c,d)pyrene	0.71	0.011	EPA 8270D/SIM	3-5-18	3-7-18	
Dibenz[a,h]anthracene	0.19	0.011	EPA 8270D/SIM	3-5-18	3-7-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	76	32 - 115				
Pyrene-d10	98	35 - 129				
Terphenyl-d14	77	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP20-2					
Laboratory ID:	03-007-29					
Benzo[a]anthracene	0.0080	0.0077	EPA 8270D/SIM	3-5-18	3-6-18	
Chrysene	0.019	0.0077	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[b]fluoranthene	0.023	0.0077	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo(j,k)fluoranthene	ND	0.0077	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[a]pyrene	0.013	0.0077	EPA 8270D/SIM	3-5-18	3-6-18	
Indeno(1,2,3-c,d)pyrene	0.013	0.0077	EPA 8270D/SIM	3-5-18	3-6-18	
Dibenz[a,h]anthracene	ND	0.0077	EPA 8270D/SIM	3-5-18	3-6-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	78	32 - 115				
Pyrene-d10	90	35 - 129				
Terphenyl-d14	89	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP20-5					
Laboratory ID:	03-007-30					
Benzo[a]anthracene	0.23	0.098	EPA 8270D/SIM	3-5-18	3-7-18	
Chrysene	0.37	0.098	EPA 8270D/SIM	3-5-18	3-7-18	
Benzo[b]fluoranthene	0.20	0.098	EPA 8270D/SIM	3-5-18	3-7-18	
Benzo(j,k)fluoranthene	ND	0.098	EPA 8270D/SIM	3-5-18	3-7-18	
Benzo[a]pyrene	0.22	0.098	EPA 8270D/SIM	3-5-18	3-7-18	
Indeno(1,2,3-c,d)pyrene	0.12	0.098	EPA 8270D/SIM	3-5-18	3-7-18	
Dibenz[a,h]anthracene	ND	0.098	EPA 8270D/SIM	3-5-18	3-7-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	74	32 - 115				
Pyrene-d10	98	35 - 129				
Terphenyl-d14	85	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP20-10					
Laboratory ID:	03-007-31					
Benzo[a]anthracene	ND	0.043	EPA 8270D/SIM	3-5-18	3-5-18	
Chrysene	ND	0.043	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[b]fluoranthene	ND	0.043	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo(j,k)fluoranthene	ND	0.043	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[a]pyrene	ND	0.043	EPA 8270D/SIM	3-5-18	3-5-18	
Indeno(1,2,3-c,d)pyrene	ND	0.043	EPA 8270D/SIM	3-5-18	3-5-18	
Dibenz[a,h]anthracene	ND	0.043	EPA 8270D/SIM	3-5-18	3-5-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	80	32 - 115				
Pyrene-d10	89	35 - 129				
Terphenyl-d14	89	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP18-3					
Laboratory ID:	03-007-33					
Benzo[a]anthracene	ND	0.0082	EPA 8270D/SIM	3-5-18	3-6-18	
Chrysene	ND	0.0082	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[b]fluoranthene	ND	0.0082	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo(j,k)fluoranthene	ND	0.0082	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[a]pyrene	ND	0.0082	EPA 8270D/SIM	3-5-18	3-6-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0082	EPA 8270D/SIM	3-5-18	3-6-18	
Dibenz[a,h]anthracene	ND	0.0082	EPA 8270D/SIM	3-5-18	3-6-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	73	32 - 115				
Pyrene-d10	79	35 - 129				
Terphenyl-d14	83	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP18-5					
Laboratory ID:	03-007-34					
Benzo[a]anthracene	ND	0.013	EPA 8270D/SIM	3-5-18	3-6-18	
Chrysene	ND	0.013	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[b]fluoranthene	ND	0.013	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo(j,k)fluoranthene	ND	0.013	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[a]pyrene	ND	0.013	EPA 8270D/SIM	3-5-18	3-6-18	
Indeno(1,2,3-c,d)pyrene	ND	0.013	EPA 8270D/SIM	3-5-18	3-6-18	
Dibenz[a,h]anthracene	ND	0.013	EPA 8270D/SIM	3-5-18	3-6-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	81	32 - 115				
Pyrene-d10	88	35 - 129				
Terphenyl-d14	90	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP18-10					
Laboratory ID:	03-007-35					
Benzo[a]anthracene	ND	0.0089	EPA 8270D/SIM	3-5-18	3-6-18	
Chrysene	ND	0.0089	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[b]fluoranthene	ND	0.0089	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo(j,k)fluoranthene	ND	0.0089	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[a]pyrene	ND	0.0089	EPA 8270D/SIM	3-5-18	3-6-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0089	EPA 8270D/SIM	3-5-18	3-6-18	
Dibenz[a,h]anthracene	ND	0.0089	EPA 8270D/SIM	3-5-18	3-6-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	82	32 - 115				
Pyrene-d10	93	35 - 129				
Terphenyl-d14	93	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP17-1					
Laboratory ID:	03-007-37					
Benzo[a]anthracene	ND	0.0086	EPA 8270D/SIM	3-5-18	3-6-18	
Chrysene	ND	0.0086	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[b]fluoranthene	ND	0.0086	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo(j,k)fluoranthene	ND	0.0086	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[a]pyrene	ND	0.0086	EPA 8270D/SIM	3-5-18	3-6-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0086	EPA 8270D/SIM	3-5-18	3-6-18	
Dibenz[a,h]anthracene	ND	0.0086	EPA 8270D/SIM	3-5-18	3-6-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	82	32 - 115				
Pyrene-d10	90	35 - 129				
Terphenyl-d14	91	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

			Date	Date	
Result	PQL	Method	Prepared	Analyzed	Flags
PP17-5					
03-007-38					
ND	0.0081	EPA 8270D/SIM	3-5-18	3-6-18	
ND	0.0081	EPA 8270D/SIM	3-5-18	3-6-18	
ND	0.0081	EPA 8270D/SIM	3-5-18	3-6-18	
ND	0.0081	EPA 8270D/SIM	3-5-18	3-6-18	
ND	0.0081	EPA 8270D/SIM	3-5-18	3-6-18	
ND	0.0081	EPA 8270D/SIM	3-5-18	3-6-18	
ND	0.0081	EPA 8270D/SIM	3-5-18	3-6-18	
Percent Recovery	Control Limits				
79	32 - 115				
92	35 - 129				
94	33 - 114				
	PP17-5 03-007-38 ND ND ND ND ND ND ND Percent Recovery 79 92	PP17-5 03-007-38 ND 0.0081 Percent Recovery Control Limits 79 32 - 115 92 35 - 129	PP17-5 03-007-38 ND 0.0081 EPA 8270D/SIM Percent Recovery Control Limits 79 32 - 115 92 35 - 129	Result PQL Method Prepared PP17-5 03-007-38 8 3-5-18 3-5-18 ND 0.0081 EPA 8270D/SIM 3-5-18 Percent Recovery Control Limits 79 32 - 115 92 35 - 129 35 - 129	Result PQL Method Prepared Analyzed PP17-5 03-007-38 8 3-6-18

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP17-10					
Laboratory ID:	03-007-39					
Benzo[a]anthracene	ND	0.0077	EPA 8270D/SIM	3-5-18	3-6-18	
Chrysene	ND	0.0077	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[b]fluoranthene	ND	0.0077	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo(j,k)fluoranthene	ND	0.0077	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[a]pyrene	ND	0.0077	EPA 8270D/SIM	3-5-18	3-6-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0077	EPA 8270D/SIM	3-5-18	3-6-18	
Dibenz[a,h]anthracene	ND	0.0077	EPA 8270D/SIM	3-5-18	3-6-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	86	32 - 115				
Pyrene-d10	96	35 - 129				
Terphenyl-d14	98	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

			Date	Date	
Result	PQL	Method	Prepared	Analyzed	Flags
PP16-1					
03-007-41					
0.20	0.036	EPA 8270D/SIM	3-5-18	3-7-18	
0.060	0.036	EPA 8270D/SIM	3-5-18	3-7-18	
0.086	0.036	EPA 8270D/SIM	3-5-18	3-7-18	
ND	0.036	EPA 8270D/SIM	3-5-18	3-7-18	
0.058	0.036	EPA 8270D/SIM	3-5-18	3-7-18	
0.036	0.036	EPA 8270D/SIM	3-5-18	3-7-18	
ND	0.036	EPA 8270D/SIM	3-5-18	3-7-18	
Percent Recovery	Control Limits				
<i>7</i> 5	32 - 115				
95	35 - 129				
76	33 - 114				
	PP16-1 03-007-41 0.20 0.060 0.086 ND 0.058 0.036 ND Percent Recovery 75 95	PP16-1 03-007-41 0.036 0.060 0.036 0.086 0.036 ND 0.036 0.058 0.036 0.036 0.036 ND 0.036 Percent Recovery Control Limits 75 32 - 115 95 35 - 129	PP16-1 03-007-41 0.20 0.036 EPA 8270D/SIM 0.060 0.036 EPA 8270D/SIM 0.086 0.036 EPA 8270D/SIM ND 0.036 EPA 8270D/SIM 0.036 EPA 8270D/SIM ND 0.036 EPA 8270D/SIM ND 0.036 EPA 8270D/SIM Percent Recovery Control Limits 75 32 - 115 95 35 - 129	Result PQL Method Prepared PP16-1 03-007-41 03-007-41 3-5-18 0.20 0.036 EPA 8270D/SIM 3-5-18 0.060 0.036 EPA 8270D/SIM 3-5-18 0.086 0.036 EPA 8270D/SIM 3-5-18 ND 0.036 EPA 8270D/SIM 3-5-18 0.036 EPA 8270D/SIM 3-5-18 ND 0.036 EPA 8270D/SIM 3-5-18 ND 0.036 EPA 8270D/SIM 3-5-18 Percent Recovery Control Limits 75 32 - 115 35 - 129	Result PQL Method Prepared Analyzed PP16-1 03-007-41 3-007-41 3-5-18 3-7-18 0.20 0.036 EPA 8270D/SIM 3-5-18 3-7-18 0.060 0.036 EPA 8270D/SIM 3-5-18 3-7-18 0.086 0.036 EPA 8270D/SIM 3-5-18 3-7-18 ND 0.036 EPA 8270D/SIM 3-5-18 3-7-18 0.036 EPA 8270D/SIM 3-5-18 3-7-18 ND 0.036 EPA 8270D/SIM 3-5-18 3-7-18 Percent Recovery Control Limits 75 32 - 115 35 - 129 35 - 129

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP16-11					
Laboratory ID:	03-007-42					
Benzo[a]anthracene	0.12	0.0082	EPA 8270D/SIM	3-5-18	3-6-18	
Chrysene	0.11	0.0082	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[b]fluoranthene	0.15	0.0082	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo(j,k)fluoranthene	0.049	0.0082	EPA 8270D/SIM	3-5-18	3-6-18	
Benzo[a]pyrene	0.10	0.0082	EPA 8270D/SIM	3-5-18	3-6-18	
Indeno(1,2,3-c,d)pyrene	0.069	0.0082	EPA 8270D/SIM	3-5-18	3-6-18	
Dibenz[a,h]anthracene	0.015	0.0082	EPA 8270D/SIM	3-5-18	3-6-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	83	32 - 115				
Pyrene-d10	92	35 - 129				
Terphenyl-d14	91	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

			Date	Date	
Result	PQL	Method	Prepared	Analyzed	Flags
PP14-5					
03-007-45					
ND	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
ND	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
ND	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
ND	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
ND	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
ND	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
ND	0.0083	EPA 8270D/SIM	3-5-18	3-6-18	
Percent Recovery	Control Limits				
80	32 - 115				
90	35 - 129				
93	33 - 114				
	PP14-5 03-007-45 ND ND ND ND ND ND ND ND Percent Recovery 80 90	PP14-5 03-007-45 ND 0.0083 Percent Recovery Control Limits 80 32 - 115 90 35 - 129	PP14-5 03-007-45 ND 0.0083 EPA 8270D/SIM Percent Recovery Control Limits 80 32 - 115 90 35 - 129	Result PQL Method Prepared PP14-5 03-007-45 3-5-18 3-5-18 ND 0.0083 EPA 8270D/SIM 3-5-18 Percent Recovery Control Limits 80 32 - 115 32 - 115 90 35 - 129 35 - 129	Result PQL Method Prepared Analyzed PP14-5 03-007-45 8270D/SIM 3-5-18 3-6-18 ND 0.0083 EPA 8270D/SIM 3-5-18 3-6-18 Percent Recovery Control Limits 80 32 - 115 30 35 - 129

Project: 17-06520-000

cPAHs EPA 8270D/SIM

			Date	Date	
Result	PQL	Method	Prepared	Analyzed	Flags
PP14-12					
03-007-46					
ND	0.0086	EPA 8270D/SIM	3-5-18	3-6-18	
ND	0.0086	EPA 8270D/SIM	3-5-18	3-6-18	
ND	0.0086	EPA 8270D/SIM	3-5-18	3-6-18	
ND	0.0086	EPA 8270D/SIM	3-5-18	3-6-18	
ND	0.0086	EPA 8270D/SIM	3-5-18	3-6-18	
0.0089	0.0086	EPA 8270D/SIM	3-5-18	3-6-18	
ND	0.0086	EPA 8270D/SIM	3-5-18	3-6-18	
Percent Recovery	Control Limits				
88	32 - 115				
96	35 - 129				
98	33 - 114				
	PP14-12 03-007-46 ND ND ND ND ND ND Percent Recovery 88 96	PP14-12 03-007-46 0.0086 ND 0.0086 Percent Recovery Control Limits 88 32 - 115 96 35 - 129	PP14-12 03-007-46 ND 0.0086 EPA 8270D/SIM Percent Recovery Control Limits 88 32 - 115 96 35 - 129	Result PQL Method Prepared PP14-12 03-007-46 FPA 8270D/SIM 3-5-18 ND 0.0086 EPA 8270D/SIM 3-5-18 Percent Recovery Control Limits 88 32 - 115 96 35 - 129	Result PQL Method Prepared Analyzed PP14-12 03-007-46 3-007-46 3-5-18 3-6-18 ND 0.0086 EPA 8270D/SIM 3-5-18 3-6-18 0.0089 0.0086 EPA 8270D/SIM 3-5-18 3-6-18 ND 0.0086 EPA 8270D/SIM 3-5-18 3-6-18 Percent Recovery Control Limits 88 32 - 115 3-6-18 96 35 - 129 35 - 129

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP11-5					
Laboratory ID:	03-007-49					
Benzo[a]anthracene	ND	0.0088	EPA 8270D/SIM	3-7-18	3-9-18	
Chrysene	ND	0.0088	EPA 8270D/SIM	3-7-18	3-9-18	
Benzo[b]fluoranthene	ND	0.0088	EPA 8270D/SIM	3-7-18	3-9-18	
Benzo(j,k)fluoranthene	ND	0.0088	EPA 8270D/SIM	3-7-18	3-9-18	
Benzo[a]pyrene	ND	0.0088	EPA 8270D/SIM	3-7-18	3-9-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0088	EPA 8270D/SIM	3-7-18	3-9-18	
Dibenz[a,h]anthracene	ND	0.0088	EPA 8270D/SIM	3-7-18	3-9-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	90	32 - 115				
Pyrene-d10	90	35 - 129				
Terphenyl-d14	96	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP11-10					
Laboratory ID:	03-007-50					
Benzo[a]anthracene	ND	0.0084	EPA 8270D/SIM	3-7-18	3-7-18	
Chrysene	ND	0.0084	EPA 8270D/SIM	3-7-18	3-7-18	
Benzo[b]fluoranthene	ND	0.0084	EPA 8270D/SIM	3-7-18	3-7-18	
Benzo(j,k)fluoranthene	ND	0.0084	EPA 8270D/SIM	3-7-18	3-7-18	
Benzo[a]pyrene	ND	0.0084	EPA 8270D/SIM	3-7-18	3-7-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0084	EPA 8270D/SIM	3-7-18	3-7-18	
Dibenz[a,h]anthracene	ND	0.0084	EPA 8270D/SIM	3-7-18	3-7-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	77	32 - 115				
Pyrene-d10	87	35 - 129				
Terphenyl-d14	94	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

			Date	Date	
Result	PQL	Method	Prepared	Analyzed	Flags
PP10-11					
03-007-52					
ND	0.011	EPA 8270D/SIM	3-7-18	3-7-18	
ND	0.011	EPA 8270D/SIM	3-7-18	3-7-18	
ND	0.011	EPA 8270D/SIM	3-7-18	3-7-18	
ND	0.011	EPA 8270D/SIM	3-7-18	3-7-18	
ND	0.011	EPA 8270D/SIM	3-7-18	3-7-18	
ND	0.011	EPA 8270D/SIM	3-7-18	3-7-18	
ND	0.011	EPA 8270D/SIM	3-7-18	3-7-18	
Percent Recovery	Control Limits				
70	32 - 115				
79	35 - 129				
85	33 - 114				
	PP10-11 03-007-52 ND ND ND ND ND ND ND ND Percent Recovery 70 79	PP10-11 03-007-52 ND 0.011 ND 0.011 ND 0.011 ND 0.011 ND 0.011 ND 0.011 Percent Recovery Control Limits 70 32 - 115 79 35 - 129	PP10-11 03-007-52 ND 0.011 EPA 8270D/SIM Percent Recovery Control Limits 70 32 - 115 79 35 - 129	Result PQL Method Prepared PP10-11 03-007-52	Result PQL Method Prepared Analyzed PP10-11 3-007-52 3-7-18 3-7-18 3-7-18 ND 0.011 EPA 8270D/SIM 3-7-18 3-7-18 Percent Recovery Control Limits 70 32 - 115 35 - 129

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP10-17					
Laboratory ID:	03-007-53					
Benzo[a]anthracene	ND	0.0083	EPA 8270D/SIM	3-7-18	3-7-18	
Chrysene	ND	0.0083	EPA 8270D/SIM	3-7-18	3-7-18	
Benzo[b]fluoranthene	ND	0.0083	EPA 8270D/SIM	3-7-18	3-7-18	
Benzo(j,k)fluoranthene	ND	0.0083	EPA 8270D/SIM	3-7-18	3-7-18	
Benzo[a]pyrene	ND	0.0083	EPA 8270D/SIM	3-7-18	3-7-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0083	EPA 8270D/SIM	3-7-18	3-7-18	
Dibenz[a,h]anthracene	ND	0.0083	EPA 8270D/SIM	3-7-18	3-7-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	72	32 - 115				
Pyrene-d10	86	35 - 129				
Terphenyl-d14	96	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP27-7					
Laboratory ID:	03-007-55					
Benzo[a]anthracene	0.036	0.0097	EPA 8270D/SIM	3-7-18	3-8-18	
Chrysene	0.051	0.0097	EPA 8270D/SIM	3-7-18	3-8-18	
Benzo[b]fluoranthene	0.049	0.0097	EPA 8270D/SIM	3-7-18	3-8-18	
Benzo(j,k)fluoranthene	0.014	0.0097	EPA 8270D/SIM	3-7-18	3-8-18	
Benzo[a]pyrene	0.038	0.0097	EPA 8270D/SIM	3-7-18	3-8-18	
Indeno(1,2,3-c,d)pyrene	0.025	0.0097	EPA 8270D/SIM	3-7-18	3-8-18	
Dibenz[a,h]anthracene	ND	0.0097	EPA 8270D/SIM	3-7-18	3-8-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	61	32 - 115				
Pyrene-d10	71	35 - 129				
Terphenyl-d14	75	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP27-10					
Laboratory ID:	03-007-56					
Benzo[a]anthracene	ND	0.0081	EPA 8270D/SIM	3-7-18	3-8-18	
Chrysene	ND	0.0081	EPA 8270D/SIM	3-7-18	3-8-18	
Benzo[b]fluoranthene	ND	0.0081	EPA 8270D/SIM	3-7-18	3-8-18	
Benzo(j,k)fluoranthene	ND	0.0081	EPA 8270D/SIM	3-7-18	3-8-18	
Benzo[a]pyrene	ND	0.0081	EPA 8270D/SIM	3-7-18	3-8-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0081	EPA 8270D/SIM	3-7-18	3-8-18	
Dibenz[a,h]anthracene	ND	0.0081	EPA 8270D/SIM	3-7-18	3-8-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	88	32 - 115				
Pyrene-d10	87	35 - 129				
Terphenyl-d14	89	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP30-5					
Laboratory ID:	03-007-59					
Benzo[a]anthracene	ND	0.085	EPA 8270D/SIM	3-7-18	3-9-18	
Chrysene	ND	0.085	EPA 8270D/SIM	3-7-18	3-9-18	
Benzo[b]fluoranthene	ND	0.085	EPA 8270D/SIM	3-7-18	3-9-18	
Benzo(j,k)fluoranthene	ND	0.085	EPA 8270D/SIM	3-7-18	3-9-18	
Benzo[a]pyrene	ND	0.085	EPA 8270D/SIM	3-7-18	3-9-18	
Indeno(1,2,3-c,d)pyrene	ND	0.085	EPA 8270D/SIM	3-7-18	3-9-18	
Dibenz[a,h]anthracene	ND	0.085	EPA 8270D/SIM	3-7-18	3-9-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	76	32 - 115				
Pyrene-d10	85	35 - 129				
Terphenyl-d14	83	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

			Date	Date	
Result	PQL	Method	Prepared	Analyzed	Flags
PP30-10					
03-007-60					
ND	0.0076	EPA 8270D/SIM	3-7-18	3-8-18	
ND	0.0076	EPA 8270D/SIM	3-7-18	3-8-18	
ND	0.0076	EPA 8270D/SIM	3-7-18	3-8-18	
ND	0.0076	EPA 8270D/SIM	3-7-18	3-8-18	
ND	0.0076	EPA 8270D/SIM	3-7-18	3-8-18	
ND	0.0076	EPA 8270D/SIM	3-7-18	3-8-18	
ND	0.0076	EPA 8270D/SIM	3-7-18	3-8-18	
Percent Recovery	Control Limits				
88	32 - 115				
90	35 - 129				
95	33 - 114				
	PP30-10 03-007-60 ND ND ND ND ND ND ND ND Percent Recovery 88 90	PP30-10 03-007-60 ND 0.0076 Percent Recovery Control Limits 88 32 - 115 90 35 - 129	PP30-10 03-007-60 ND 0.0076 EPA 8270D/SIM Percent Recovery Control Limits 88 32 - 115 90 35 - 129	Result PQL Method Prepared PP30-10 03-007-60 3-007-60 3-7-18 3-7-18 ND 0.0076 EPA 8270D/SIM 3-7-18 <td>Result PQL Method Prepared Analyzed PP30-10 03-007-60 BPA 8270D/SIM 3-7-18 3-8-18 ND 0.0076 EPA 8270D/SIM 3-7-18 3-8-18 Percent Recovery Control Limits 88 32 - 115 90 35 - 129 35 - 129</td>	Result PQL Method Prepared Analyzed PP30-10 03-007-60 BPA 8270D/SIM 3-7-18 3-8-18 ND 0.0076 EPA 8270D/SIM 3-7-18 3-8-18 Percent Recovery Control Limits 88 32 - 115 90 35 - 129 35 - 129

Project: 17-06520-000

cPAHs EPA 8270D/SIM

			Date	Date	
Result	PQL	Method	Prepared	Analyzed	Flags
PP24-7					
03-007-63					
ND	0.018	EPA 8270D/SIM	3-7-18	3-9-18	
0.035	0.018	EPA 8270D/SIM	3-7-18	3-9-18	
0.031	0.018	EPA 8270D/SIM	3-7-18	3-9-18	
ND	0.018	EPA 8270D/SIM	3-7-18	3-9-18	
ND	0.018	EPA 8270D/SIM	3-7-18	3-9-18	
ND	0.018	EPA 8270D/SIM	3-7-18	3-9-18	
ND	0.018	EPA 8270D/SIM	3-7-18	3-9-18	
Percent Recovery	Control Limits				
71	32 - 115				
86	35 - 129				
84	33 - 114				
	PP24-7 03-007-63 ND 0.035 0.031 ND ND ND ND Percent Recovery 71 86	PP24-7 03-007-63 ND 0.018 0.035 0.018 0.031 0.018 ND 0.018 ND 0.018 ND 0.018 ND 0.018 ND 0.018 Percent Recovery Control Limits 71 32 - 115 86 35 - 129	PP24-7 03-007-63 ND 0.018 EPA 8270D/SIM 0.035 0.018 EPA 8270D/SIM 0.031 0.018 EPA 8270D/SIM ND 0.018 EPA 8270D/SIM ND 0.018 EPA 8270D/SIM ND 0.018 EPA 8270D/SIM ND 0.018 EPA 8270D/SIM Percent Recovery Control Limits 71 32 - 115 86 35 - 129	PP24-7 03-007-63 ND 0.018 EPA 8270D/SIM 3-7-18 0.035 0.018 EPA 8270D/SIM 3-7-18 0.031 0.018 EPA 8270D/SIM 3-7-18 ND 0.018 EPA 8270D/SIM 3-7-18 Percent Recovery Control Limits 71 32 - 115 35 - 129	Result PQL Method Prepared Analyzed PP24-7 03-007-63 </td

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP24-10					
Laboratory ID:	03-007-64					
Benzo[a]anthracene	ND	0.0077	EPA 8270D/SIM	3-7-18	3-8-18	
Chrysene	ND	0.0077	EPA 8270D/SIM	3-7-18	3-8-18	
Benzo[b]fluoranthene	ND	0.0077	EPA 8270D/SIM	3-7-18	3-8-18	
Benzo(j,k)fluoranthene	ND	0.0077	EPA 8270D/SIM	3-7-18	3-8-18	
Benzo[a]pyrene	ND	0.0077	EPA 8270D/SIM	3-7-18	3-8-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0077	EPA 8270D/SIM	3-7-18	3-8-18	
Dibenz[a,h]anthracene	ND	0.0077	EPA 8270D/SIM	3-7-18	3-8-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	85	32 - 115				
Pyrene-d10	88	35 - 129				
Terphenyl-d14	94	33 - 114				

Project: 17-06520-000

CPAHS EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0305S1					
Benzo[a]anthracene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Chrysene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[a]pyrene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	84	32 - 115				
Pyrene-d10	94	35 - 129				
Terphenyl-d14	102	33 - 114				

Project: 17-06520-000

CPAHS EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0305S2					
Benzo[a]anthracene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Chrysene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Benzo[a]pyrene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270D/SIM	3-5-18	3-5-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	91	32 - 115				
Pyrene-d10	103	35 - 129				
Terphenyl-d14	108	33 - 114				

Project: 17-06520-000

CPAHS EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
						_
Laboratory ID:	MB0307S1					
Benzo[a]anthracene	ND	0.0067	EPA 8270D/SIM	3-7-18	3-7-18	
Chrysene	ND	0.0067	EPA 8270D/SIM	3-7-18	3-7-18	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270D/SIM	3-7-18	3-7-18	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270D/SIM	3-7-18	3-7-18	
Benzo[a]pyrene	ND	0.0067	EPA 8270D/SIM	3-7-18	3-7-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0067	EPA 8270D/SIM	3-7-18	3-7-18	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270D/SIM	3-7-18	3-7-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	80	32 - 115				
Pyrene-d10	92	35 - 129				
Terphenyl-d14	90	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM SB/SBD QUALITY CONTROL

					Per	cent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Reco	overy	Limits	RPD	Limit	Flags
SPIKE BLANKS										
Laboratory ID:	SB03	805S1								
	SB	SBD	SB	SBD	SB	SBD				
Benzo[a]anthracene	0.0878	0.0851	0.0833	0.0833	105	102	64 - 135	3	15	
Chrysene	0.0869	0.0844	0.0833	0.0833	104	101	70 - 119	3	15	
Benzo[b]fluoranthene	0.0826	0.0807	0.0833	0.0833	99	97	54 - 135	2	15	
Benzo(j,k)fluoranthene	0.0872	0.0839	0.0833	0.0833	105	101	66 - 122	4	15	
Benzo[a]pyrene	0.0816	0.0789	0.0833	0.0833	98	95	62 - 125	3	15	
Indeno(1,2,3-c,d)pyrene	0.0752	0.0724	0.0833	0.0833	90	87	55 - 129	4	15	
Dibenz[a,h]anthracene	0.0769	0.0756	0.0833	0.0833	92	91	58 - 125	2	15	
Surrogate:										
2-Fluorobiphenyl					88	87	32 - 115			
Pyrene-d10					99	96	35 - 129			
Terphenyl-d14					105	102	33 - 114			

Project: 17-06520-000

cPAHs EPA 8270D/SIM MS/MSD QUALITY CONTROL

					Source	Per	cent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Rec	overy	Limits	RPD	Limit	Flags
MATRIX SPIKES											
Laboratory ID:	03-0	07-31									
	MS	MSD	MS	MSD		MS	MSD				
Benzo[a]anthracene	0.0785	0.0813	0.0833	0.0833	ND	94	98	27 - 143	4	23	
Chrysene	0.0770	0.0834	0.0833	0.0833	ND	92	100	22 - 130	8	24	
Benzo[b]fluoranthene	0.0753	0.0798	0.0833	0.0833	ND	90	96	15 - 141	6	26	
Benzo(j,k)fluoranthene	0.0700	0.0742	0.0833	0.0833	ND	84	89	42 - 112	6	24	
Benzo[a]pyrene	0.0718	0.0759	0.0833	0.0833	ND	86	91	33 - 126	6	26	
Indeno(1,2,3-c,d)pyrene	0.0683	0.0752	0.0833	0.0833	ND	82	90	30 - 125	10	25	
Dibenz[a,h]anthracene	0.0684	0.0734	0.0833	0.0833	ND	82	88	31 - 124	7	22	
Surrogate:											
2-Fluorobiphenyl						77	81	32 - 115			
Pyrene-d10						85	90	35 - 129			
Terphenyl-d14						86	91	33 - 114			

Project: 17-06520-000

cPAHs EPA 8270D/SIM MS/MSD QUALITY CONTROL

					Source	Per	cent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Rec	overy	Limits	RPD	Limit	Flags
MATRIX SPIKES											
Laboratory ID:	03-03	36-01									
	MS	MSD	MS	MSD		MS	MSD				
Benzo[a]anthracene	0.0816	0.0895	0.0833	0.0833	0.00697	90	99	27 - 143	9	23	
Chrysene	0.0782	0.0842	0.0833	0.0833	0.00958	82	90	22 - 130	7	24	
Benzo[b]fluoranthene	0.0811	0.0890	0.0833	0.0833	0.0122	83	92	15 - 141	9	26	
Benzo(j,k)fluoranthene	0.0729	0.0791	0.0833	0.0833	ND	88	95	42 - 112	8	24	
Benzo[a]pyrene	0.0746	0.0820	0.0833	0.0833	ND	90	98	33 - 126	9	26	
Indeno(1,2,3-c,d)pyrene	0.0768	0.0876	0.0833	0.0833	ND	92	105	30 - 125	13	25	
Dibenz[a,h]anthracene	0.0664	0.0774	0.0833	0.0833	ND	80	93	31 - 124	15	22	
Surrogate:											
2-Fluorobiphenyl						70	77	32 - 115			
Pyrene-d10						81	90	35 - 129			
Terphenyl-d14						84	93	33 - 114			

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP32-10					
Laboratory ID:	03-007-11					
Benzo[a]anthracene	ND	0.0081	EPA 8270D/SIM	3-8-18	3-9-18	
Chrysene	ND	0.0081	EPA 8270D/SIM	3-8-18	3-9-18	
Benzo[b]fluoranthene	ND	0.0081	EPA 8270D/SIM	3-8-18	3-9-18	
Benzo(j,k)fluoranthene	ND	0.0081	EPA 8270D/SIM	3-8-18	3-9-18	
Benzo[a]pyrene	ND	0.0081	EPA 8270D/SIM	3-8-18	3-9-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0081	EPA 8270D/SIM	3-8-18	3-9-18	
Dibenz[a,h]anthracene	ND	0.0081	EPA 8270D/SIM	3-8-18	3-9-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	68	32 - 115				
Pyrene-d10	70	35 - 129				
Terphenyl-d14	<i>7</i> 5	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP21-15					
Laboratory ID:	03-007-28					
Benzo[a]anthracene	ND	0.018	EPA 8270D/SIM	3-12-18	3-15-18	
Chrysene	ND	0.018	EPA 8270D/SIM	3-12-18	3-15-18	
Benzo[b]fluoranthene	ND	0.018	EPA 8270D/SIM	3-12-18	3-15-18	
Benzo(j,k)fluoranthene	ND	0.018	EPA 8270D/SIM	3-12-18	3-15-18	
Benzo[a]pyrene	ND	0.018	EPA 8270D/SIM	3-12-18	3-15-18	
Indeno(1,2,3-c,d)pyrene	ND	0.018	EPA 8270D/SIM	3-12-18	3-15-18	
Dibenz[a,h]anthracene	ND	0.018	EPA 8270D/SIM	3-12-18	3-15-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	74	32 - 115				
Pyrene-d10	89	35 - 129				
Terphenyl-d14	89	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP20-15					
Laboratory ID:	03-007-32					
Benzo[a]anthracene	ND	0.046	EPA 8270D/SIM	3-8-18	3-9-18	
Chrysene	ND	0.046	EPA 8270D/SIM	3-8-18	3-9-18	
Benzo[b]fluoranthene	ND	0.046	EPA 8270D/SIM	3-8-18	3-9-18	
Benzo(j,k)fluoranthene	ND	0.046	EPA 8270D/SIM	3-8-18	3-9-18	
Benzo[a]pyrene	ND	0.046	EPA 8270D/SIM	3-8-18	3-9-18	
Indeno(1,2,3-c,d)pyrene	ND	0.046	EPA 8270D/SIM	3-8-18	3-9-18	
Dibenz[a,h]anthracene	ND	0.046	EPA 8270D/SIM	3-8-18	3-9-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	77	32 - 115				
Pyrene-d10	72	35 - 129				
Terphenyl-d14	76	33 - 114				

Project: 17-06520-000

CPAHS EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0308S1					
Benzo[a]anthracene	ND	0.0067	EPA 8270D/SIM	3-8-18	3-9-18	
Chrysene	ND	0.0067	EPA 8270D/SIM	3-8-18	3-9-18	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270D/SIM	3-8-18	3-9-18	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270D/SIM	3-8-18	3-9-18	
Benzo[a]pyrene	ND	0.0067	EPA 8270D/SIM	3-8-18	3-9-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0067	EPA 8270D/SIM	3-8-18	3-9-18	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270D/SIM	3-8-18	3-9-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	70	32 - 115				
Pyrene-d10	83	35 - 129				
Terphenyl-d14	83	33 - 114				

Project: 17-06520-000

CPAHS EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
						_
Laboratory ID:	MB0312S1					
Benzo[a]anthracene	ND	0.0040	EPA 8270D/SIM	3-12-18	3-12-18	_
Chrysene	ND	0.0040	EPA 8270D/SIM	3-12-18	3-12-18	
Benzo[b]fluoranthene	ND	0.0040	EPA 8270D/SIM	3-12-18	3-12-18	
Benzo(j,k)fluoranthene	ND	0.0040	EPA 8270D/SIM	3-12-18	3-12-18	
Benzo[a]pyrene	ND	0.0040	EPA 8270D/SIM	3-12-18	3-12-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0040	EPA 8270D/SIM	3-12-18	3-12-18	
Dibenz[a,h]anthracene	ND	0.0040	EPA 8270D/SIM	3-12-18	3-12-18	
Surrogate:	Percent Recovery	Control Limits				_
2-Fluorobiphenyl	63	32 - 115				
Pyrene-d10	<i>7</i> 5	35 - 129				
Terphenyl-d14	81	33 - 114				

Project: 17-06520-000

cPAHs EPA 8270D/SIM SB/SBD QUALITY CONTROL

					Per	cent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Rec	overy	Limits	RPD	Limit	Flags
SPIKE BLANKS										
Laboratory ID:	SB03	808S1								
	SB	SBD	SB	SBD	SB	SBD				
Benzo[a]anthracene	0.0694	0.0713	0.0833	0.0833	83	86	64 - 135	3	15	
Chrysene	0.0645	0.0662	0.0833	0.0833	77	79	70 - 119	3	15	
Benzo[b]fluoranthene	0.0653	0.0677	0.0833	0.0833	78	81	54 - 135	4	15	
Benzo(j,k)fluoranthene	0.0658	0.0688	0.0833	0.0833	79	83	66 - 122	4	15	
Benzo[a]pyrene	0.0641	0.0667	0.0833	0.0833	77	80	62 - 125	4	15	
Indeno(1,2,3-c,d)pyrene	0.0637	0.0675	0.0833	0.0833	76	81	55 - 129	6	15	
Dibenz[a,h]anthracene	0.0623	0.0640	0.0833	0.0833	75	77	58 - 125	3	15	
Surrogate:										
2-Fluorobiphenyl					73	69	32 - 115			
Pyrene-d10					76	<i>7</i> 5	35 - 129			
Terphenyl-d14					83	82	33 - 114			

Project: 17-06520-000

cPAHs EPA 8270D/SIM SB/SBD QUALITY CONTROL

					Pei	cent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Rec	overy	Limits	RPD	Limit	Flags
SPIKE BLANKS										
Laboratory ID:	SB03	12S1								
	SB	SBD	SB	SBD	SB	SBD				
Benzo[a]anthracene	0.0397	0.0412	0.0500	0.0500	79	82	64 - 135	4	15	
Chrysene	0.0372	0.0391	0.0500	0.0500	74	78	70 - 119	5	15	
Benzo[b]fluoranthene	0.0375	0.0399	0.0500	0.0500	75	80	54 - 135	6	15	
Benzo(j,k)fluoranthene	0.0381	0.0389	0.0500	0.0500	76	78	66 - 122	2	15	
Benzo[a]pyrene	0.0334	0.0344	0.0500	0.0500	67	69	62 - 125	3	15	
Indeno(1,2,3-c,d)pyrene	0.0382	0.0396	0.0500	0.0500	76	79	55 - 129	4	15	
Dibenz[a,h]anthracene	0.0371	0.0384	0.0500	0.0500	74	77	58 - 125	3	15	
Surrogate:										
2-Fluorobiphenyl					63	64	32 - 115			
Pyrene-d10					74	<i>7</i> 5	35 - 129			
Terphenyl-d14					79	80	33 - 114			

Project: 17-06520-000

PCBs EPA 8082A

Matrix: Soil

critis. Trig/rkg (ppini)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP29-3					
Laboratory ID:	03-007-05					
Aroclor 1016	ND	0.056	EPA 8082A	3-12-18	3-13-18	
Aroclor 1221	ND	0.056	EPA 8082A	3-12-18	3-13-18	
Aroclor 1232	ND	0.056	EPA 8082A	3-12-18	3-13-18	
Aroclor 1242	ND	0.056	EPA 8082A	3-12-18	3-13-18	
Aroclor 1248	ND	0.056	EPA 8082A	3-12-18	3-13-18	
Aroclor 1254	ND	0.056	EPA 8082A	3-12-18	3-13-18	
Aroclor 1260	ND	0.056	EPA 8082A	3-12-18	3-13-18	
Surrogate:	Percent Recovery	Control Limits				
DCB	81	40-134				
Client ID:	PP32-7					
Laboratory ID:	03-007-10					
Aroclor 1016	ND	0.061	EPA 8082A	3-12-18	3-13-18	
Aroclor 1221	ND	0.061	EPA 8082A	3-12-18	3-13-18	
Aroclor 1232	ND	0.061	EPA 8082A	3-12-18	3-13-18	
Aroclor 1242	ND	0.061	EPA 8082A	3-12-18	3-13-18	
Aroclor 1248	ND	0.061	EPA 8082A	3-12-18	3-13-18	
Aroclor 1254	ND	0.061	EPA 8082A	3-12-18	3-13-18	
Aroclor 1260	ND	0.061	EPA 8082A	3-12-18	3-13-18	
Surrogate:	Percent Recovery	Control Limits				
DCB	85	40-134				
Client ID:	PP32-10					
Laboratory ID:	03-007-11					
Aroclor 1016	ND	0.061	EPA 8082A	3-12-18	3-13-18	
Aroclor 1221	ND	0.061	EPA 8082A	3-12-18	3-13-18	
Aroclor 1232	ND	0.061	EPA 8082A	3-12-18	3-13-18	
Aroclor 1242	ND	0.061	EPA 8082A	3-12-18	3-13-18	
Aroclor 1248	ND	0.061	EPA 8082A	3-12-18	3-13-18	
Aroclor 1254	ND	0.061	EPA 8082A	3-12-18	3-13-18	
Aroclor 1260	ND	0.061	EPA 8082A	3-12-18	3-13-18	
Surrogate:	Percent Recovery	Control Limits			· ·	
DCB	70	40-134				

Project: 17-06520-000

PCBs EPA 8082A

Matrix: Soil

Units: mg/Kg (ppm)

ome. mg/rtg (ppm)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP33-3					
Laboratory ID:	03-007-17					
Aroclor 1016	ND	0.056	EPA 8082A	3-12-18	3-13-18	
Aroclor 1221	ND	0.056	EPA 8082A	3-12-18	3-13-18	
Aroclor 1232	ND	0.056	EPA 8082A	3-12-18	3-13-18	
Aroclor 1242	ND	0.056	EPA 8082A	3-12-18	3-13-18	
Aroclor 1248	ND	0.056	EPA 8082A	3-12-18	3-13-18	
Aroclor 1254	ND	0.056	EPA 8082A	3-12-18	3-13-18	
Aroclor 1260	ND	0.056	EPA 8082A	3-12-18	3-13-18	
Surrogate:	Percent Recovery	Control Limits				
DCB	72	40-134				
Client ID:	PP23-2					
Laboratory ID:	03-007-21					
Aroclor 1016	ND	0.059	EPA 8082A	3-12-18	3-13-18	
Aroclor 1221	ND	0.059	EPA 8082A	3-12-18	3-13-18	
Aroclor 1232	ND	0.059	EPA 8082A	3-12-18	3-13-18	
Aroclor 1242	ND	0.059	EPA 8082A	3-12-18	3-13-18	
Aroclor 1248	ND	0.059	EPA 8082A	3-12-18	3-13-18	
Aroclor 1254	ND	0.059	EPA 8082A	3-12-18	3-13-18	
Aroclor 1260	ND	0.059	EPA 8082A	3-12-18	3-13-18	
Surrogate:	Percent Recovery	Control Limits				
DCB	70	40-134				
Client ID:	PP21-2					
Laboratory ID:	03-007-25					
Aroclor 1016	ND	0.066	EPA 8082A	3-12-18	3-13-18	
Aroclor 1221	ND	0.066	EPA 8082A	3-12-18	3-13-18	
Aroclor 1232	ND	0.066	EPA 8082A	3-12-18	3-13-18	
Aroclor 1242	ND	0.066	EPA 8082A	3-12-18	3-13-18	
Aroclor 1248	ND	0.066	EPA 8082A	3-12-18	3-13-18	
Aroclor 1254	0.42	0.066	EPA 8082A	3-12-18	3-13-18	
Aroclor 1260	0.10	0.066	EPA 8082A	3-12-18	3-13-18	
Surrogate:	Percent Recovery	Control Limits				

DCB 74 40-134

Project: 17-06520-000

PCBs EPA 8082A

Matrix: Soil

Units: mg/Kg (ppm)

Analysta	Popult	PQL	Mathad	Date	Date	Eloss
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP21-6					
Laboratory ID:	03-007-26					
Aroclor 1016	ND	0.10	EPA 8082A	3-12-18	3-14-18	X
Aroclor 1221	ND	0.10	EPA 8082A	3-12-18	3-14-18	X
Aroclor 1232	ND	0.10	EPA 8082A	3-12-18	3-14-18	X
Aroclor 1242	ND	0.10	EPA 8082A	3-12-18	3-14-18	X
Aroclor 1248	ND	0.10	EPA 8082A	3-12-18	3-14-18	X
Aroclor 1254	0.27	0.10	EPA 8082A	3-12-18	3-14-18	X
Aroclor 1260	ND	0.10	EPA 8082A	3-12-18	3-14-18	X
Surrogate:	Percent Recovery	Control Limits				
DCB	76	40-134				
Client ID:	PP21-10					
Laboratory ID:	03-007-27					
Aroclor 1016	ND	0.081	EPA 8082A	3-12-18	3-14-18	X
Aroclor 1221	ND	0.081	EPA 8082A	3-12-18	3-14-18	Χ
Aroclor 1232	ND	0.081	EPA 8082A	3-12-18	3-14-18	Χ
Aroclor 1242	0.11	0.081	EPA 8082A	3-12-18	3-14-18	Χ
Aroclor 1248	ND	0.081	EPA 8082A	3-12-18	3-14-18	X
Aroclor 1254	0.99	0.081	EPA 8082A	3-12-18	3-14-18	X
Aroclor 1260	0.23	0.081	EPA 8082A	3-12-18	3-14-18	X
Surrogate:	Percent Recovery	Control Limits				
DCB	87	40-134				
Client ID:	PP21-15					
Laboratory ID:	03-007-28					
Aroclor 1016	ND	0.14	EPA 8082A	3-12-18	3-13-18	
Aroclor 1221	ND	0.14	EPA 8082A	3-12-18	3-13-18	
Aroclor 1232	ND	0.14	EPA 8082A	3-12-18	3-13-18	
Aroclor 1242	ND	0.14	EPA 8082A	3-12-18	3-13-18	
Aroclor 1248	ND	0.14	EPA 8082A	3-12-18	3-13-18	
Aroclor 1254	ND	0.14	EPA 8082A	3-12-18	3-13-18	
Aroclor 1260	ND	0.14	EPA 8082A	3-12-18	3-13-18	
Surrogate:	Percent Recovery	Control Limits				
DOD	00	10 101				

DCB 68 40-134

Project: 17-06520-000

PCBs EPA 8082A

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP20-2					
Laboratory ID:	03-007-29					
Aroclor 1016	ND	0.058	EPA 8082A	3-12-18	3-13-18	
Aroclor 1221	ND	0.058	EPA 8082A	3-12-18	3-13-18	
Aroclor 1232	ND	0.058	EPA 8082A	3-12-18	3-13-18	
Aroclor 1242	ND	0.058	EPA 8082A	3-12-18	3-13-18	
Aroclor 1248	ND	0.058	EPA 8082A	3-12-18	3-13-18	
Aroclor 1254	ND	0.058	EPA 8082A	3-12-18	3-13-18	
Aroclor 1260	ND	0.058	EPA 8082A	3-12-18	3-13-18	
Surrogate:	Percent Recovery	Control Limits				
DCB	83	40-134				
Client ID:	PP20-5					
Laboratory ID:	03-007-30					
Aroclor 1016	ND	0.073	EPA 8082A	3-12-18	3-14-18	Х
Aroclor 1221	ND	0.073	EPA 8082A	3-12-18	3-14-18	Х
Aroclor 1232	ND	0.073	EPA 8082A	3-12-18	3-14-18	Χ
Aroclor 1242	0.16	0.073	EPA 8082A	3-12-18	3-14-18	Χ
Aroclor 1248	ND	0.073	EPA 8082A	3-12-18	3-14-18	X
Aroclor 1254	0.48	0.073	EPA 8082A	3-12-18	3-14-18	Χ
Aroclor 1260	0.10	0.073	EPA 8082A	3-12-18	3-14-18	Χ
Surrogate:	Percent Recovery	Control Limits				
DCB	123	40-134				
Client ID:	PP20-10					
Laboratory ID:	03-007-31					
Aroclor 1016	ND	0.064	EPA 8082A	3-12-18	3-13-18	
Aroclor 1221	ND	0.064	EPA 8082A	3-12-18	3-13-18	
Aroclor 1232	ND	0.064	EPA 8082A	3-12-18	3-13-18	
Aroclor 1242	ND	0.064	EPA 8082A	3-12-18	3-13-18	
Aroclor 1248	ND	0.064	EPA 8082A	3-12-18	3-13-18	
Aroclor 1254	ND	0.064	EPA 8082A	3-12-18	3-13-18	
Aroclor 1260	ND	0.064	EPA 8082A	3-12-18	3-13-18	
Surrogate:	Percent Recovery	Control Limits				
DCB	93	40-134				

Project: 17-06520-000

PCBs EPA 8082A

Matrix: Soil

oring. Trig/reg (ppin)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP20-15					
Laboratory ID:	03-007-32					
Aroclor 1016	ND	0.070	EPA 8082A	3-12-18	3-13-18	
Aroclor 1221	ND	0.070	EPA 8082A	3-12-18	3-13-18	
Aroclor 1232	ND	0.070	EPA 8082A	3-12-18	3-13-18	
Aroclor 1242	ND	0.070	EPA 8082A	3-12-18	3-13-18	
Aroclor 1248	ND	0.070	EPA 8082A	3-12-18	3-13-18	
Aroclor 1254	ND	0.070	EPA 8082A	3-12-18	3-13-18	
Aroclor 1260	ND	0.070	EPA 8082A	3-12-18	3-13-18	
Surrogate:	Percent Recovery	Control Limits				
DCB	67	40-134				
Client ID:	PP16-1					
Laboratory ID:	03-007-41					
Aroclor 1016	ND	0.053	EPA 8082A	3-12-18	3-13-18	
Aroclor 1221	ND	0.053	EPA 8082A	3-12-18	3-13-18	
Aroclor 1232	ND	0.053	EPA 8082A	3-12-18	3-13-18	
Aroclor 1242	ND	0.053	EPA 8082A	3-12-18	3-13-18	
Aroclor 1248	ND	0.053	EPA 8082A	3-12-18	3-13-18	
Aroclor 1254	ND	0.053	EPA 8082A	3-12-18	3-13-18	
Aroclor 1260	ND	0.053	EPA 8082A	3-12-18	3-13-18	
Surrogate:	Percent Recovery	Control Limits				
DCB	59	40-134				
Client ID:	PP30-5					
Laboratory ID:	03-007-59					
Aroclor 1016	ND	0.064	EPA 8082A	3-12-18	3-13-18	
Aroclor 1221	ND	0.064	EPA 8082A	3-12-18	3-13-18	
Aroclor 1232	ND	0.064	EPA 8082A	3-12-18	3-13-18	
Aroclor 1242	ND	0.064	EPA 8082A	3-12-18	3-13-18	
Aroclor 1248	ND	0.064	EPA 8082A	3-12-18	3-13-18	
Aroclor 1254	ND	0.064	EPA 8082A	3-12-18	3-13-18	
Aroclor 1260	ND	0.064	EPA 8082A	3-12-18	3-13-18	
Surrogate:	Percent Recovery	Control Limits		-		
DCB	65	40-134				

Project: 17-06520-000

PCBs EPA 8082A

Matrix: Soil

Units: mg/Kg (ppm)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP24-7					
Laboratory ID:	03-007-63					
Aroclor 1016	ND	0.069	EPA 8082A	3-12-18	3-13-18	
Aroclor 1221	ND	0.069	EPA 8082A	3-12-18	3-13-18	
Aroclor 1232	ND	0.069	EPA 8082A	3-12-18	3-13-18	
Aroclor 1242	ND	0.069	EPA 8082A	3-12-18	3-13-18	
Aroclor 1248	ND	0.069	EPA 8082A	3-12-18	3-13-18	
Aroclor 1254	ND	0.069	EPA 8082A	3-12-18	3-13-18	
Aroclor 1260	ND	0.069	EPA 8082A	3-12-18	3-13-18	
		•			•	•

Surrogate: Percent Recovery Control Limits DCB 75 40-134

Project: 17-06520-000

PCBs EPA 8082A QUALITY CONTROL

Matrix: Soil

5 5 (1)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0312S1					
Aroclor 1016	ND	0.050	EPA 8082A	3-12-18	3-13-18	
Aroclor 1221	ND	0.050	EPA 8082A	3-12-18	3-13-18	
Aroclor 1232	ND	0.050	EPA 8082A	3-12-18	3-13-18	
Aroclor 1242	ND	0.050	EPA 8082A	3-12-18	3-13-18	
Aroclor 1248	ND	0.050	EPA 8082A	3-12-18	3-13-18	
Aroclor 1254	ND	0.050	EPA 8082A	3-12-18	3-13-18	
Aroclor 1260	ND	0.050	EPA 8082A	3-12-18	3-13-18	
Surrogate:	Percent Recovery	Control Limits				
DCB	81	40-134				
Laboratory ID:	MB0312S1					
Aroclor 1016	ND	0.050	EPA 8082A	3-12-18	3-13-18	Χ
Aroclor 1221	ND	0.050	EPA 8082A	3-12-18	3-13-18	Χ
Aroclor 1232	ND	0.050	EPA 8082A	3-12-18	3-13-18	Χ
Aroclor 1242	ND	0.050	EPA 8082A	3-12-18	3-13-18	Χ
Aroclor 1248	ND	0.050	EPA 8082A	3-12-18	3-13-18	Χ
Aroclor 1254	ND	0.050	EPA 8082A	3-12-18	3-13-18	Χ
Aroclor 1260	ND	0.050	EPA 8082A	3-12-18	3-13-18	Х
Surrogate:	Percent Recovery	Control Limits		·		
DCB	78	40-134				

Analyte	Re	sult	Spike	Level	Source Result		rcent	Recovery Limits	RPD	RPD Limit	Flags
MATRIX SPIKES											
Laboratory ID:	03-0	75-05									
	MS	MSD	MS	MSD		MS	MSD				
Aroclor 1260	0.422	0.372	0.500	0.500	ND	84	74	34-126	13	16	
Surrogate:											
DCB						86	76	40-134			

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B

Matrix: Soil

Office. Hig/Kg (ppin)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP28-8					
Laboratory ID:	03-007-02					
Arsenic	ND	13	EPA 6010D	3-8-18	3-8-18	
Cadmium	ND	0.63	EPA 6010D	3-8-18	3-8-18	
Chromium	25	0.63	EPA 6010D	3-8-18	3-8-18	
Lead	49	6.3	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.31	EPA 7471B	3-5-18	3-5-18	
Client ID:	PP28-10					
Laboratory ID:	03-007-03					
Arsenic	ND	12	EPA 6010D	3-8-18	3-8-18	
Cadmium	ND	0.62	EPA 6010D	3-8-18	3-8-18	
Chromium	23	0.62	EPA 6010D	3-8-18	3-8-18	
Lead	6.8	6.2	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.31	EPA 7471B	3-5-18	3-5-18	
-						
Client ID:	PP29-3					
Laboratory ID:	03-007-05					
Arsenic	ND	11	EPA 6010D	3-8-18	3-8-18	
Cadmium	ND	0.56	EPA 6010D	3-8-18	3-8-18	
Chromium	25	0.56	EPA 6010D	3-8-18	3-8-18	
Lead	8.8	5.6	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.28	EPA 7471B	3-5-18	3-5-18	
Client ID:	PP29-5					
Laboratory ID:	03-007-06					
Arsenic	ND	13	EPA 6010D	3-8-18	3-8-18	
Cadmium	ND	0.65	EPA 6010D	3-8-18	3-8-18	
Chromium	15	0.65	EPA 6010D	3-8-18	3-8-18	
Lead	ND	6.5	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.32	EPA 7471B	3-5-18	3-5-18	

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B

Matrix: Soil

onits. Hig/Ng (ppin)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP29-10					
Laboratory ID:	03-007-07					
Arsenic	ND	13	EPA 6010D	3-8-18	3-8-18	
Cadmium	ND	0.66	EPA 6010D	3-8-18	3-8-18	
Chromium	16	0.66	EPA 6010D	3-8-18	3-8-18	
Lead	ND	6.6	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.33	EPA 7471B	3-5-18	3-5-18	
Client ID:	PP32-4					
Laboratory ID:	03-007-09					
Arsenic	ND	12	EPA 6010D	3-8-18	3-8-18	
Cadmium	ND	0.62	EPA 6010D	3-8-18	3-8-18	
Chromium	16	0.62	EPA 6010D	3-8-18	3-8-18	
Lead	12	6.2	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.31	EPA 7471B	3-5-18	3-5-18	
Client ID:	PP32-7					
	03-007-10					
Laboratory ID: Arsenic	ND	12	EPA 6010D	3-8-18	3-8-18	
Cadmium	ND ND	0.61	EPA 6010D	3-0-10 3-8-18	3-0-10 3-8-18	
Chromium	15	0.61	EPA 6010D	3-6-16 3-8-18	3-8-18	
Lead	15	6.1	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.1	EPA 7471B	3-5-18	3-5-18	
ivier cur y	ND	0.51	LIATATID	3-3-10	3-3-10	
Client ID:	PP31-3					
Laboratory ID:	03-007-13					
Arsenic	ND	11	EPA 6010D	3-8-18	3-8-18	
Cadmium	ND	0.56	EPA 6010D	3-8-18	3-8-18	
Chromium	21	0.56	EPA 6010D	3-8-18	3-8-18	
Lead	12	5.6	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.28	EPA 7471B	3-5-18	3-5-18	

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B

Matrix: Soil

onits. Hig/Ng (ppin)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP31-11					
Laboratory ID:	03-007-15					
Arsenic	ND	12	EPA 6010D	3-8-18	3-8-18	
Cadmium	ND	0.62	EPA 6010D	3-8-18	3-8-18	
Chromium	12	0.62	EPA 6010D	3-8-18	3-8-18	
Lead	ND	6.2	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.31	EPA 7471B	3-5-18	3-5-18	
Client ID:	PP33-3					
Laboratory ID:	03-007-17					
Arsenic	ND	11	EPA 6010D	3-8-18	3-8-18	
Cadmium	ND	0.56	EPA 6010D	3-8-18	3-8-18	
Chromium	17	0.56	EPA 6010D	3-8-18	3-8-18	
Lead	8.1	5.6	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.28	EPA 7471B	3-5-18	3-5-18	
Client ID:	PP33-5					
Laboratory ID:	03-007-18					
Arsenic	ND	12	EPA 6010D	3-8-18	3-8-18	
Cadmium	ND	0.59	EPA 6010D	3-8-18	3-8-18	
Chromium	19	0.59	EPA 6010D	3-8-18	3-8-18	
Lead	7.3	5.9	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.30	EPA 7471B	3-5-18	3-5-18	
Client ID:	PP33-10					
Laboratory ID:	03-007-19					
Arsenic	ND	13	EPA 6010D	3-8-18	3-8-18	
Cadmium	ND	0.67	EPA 6010D	3-8-18	3-8-18	
Chromium	13	0.67	EPA 6010D	3-8-18	3-8-18	
Lead	ND	6.7	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.34	EPA 7471B	3-5-18	3-5-18	

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B

Matrix: Soil

Units: mg/kg (ppm)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP23-2			•	-	
Laboratory ID:	03-007-21					
Arsenic	ND	12	EPA 6010D	3-8-18	3-8-18	
Cadmium	ND	0.59	EPA 6010D	3-8-18	3-8-18	
Chromium	19	0.59	EPA 6010D	3-8-18	3-8-18	
Lead	460	5.9	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.29	EPA 7471B	3-5-18	3-5-18	
Client ID:	PP23-5					
Laboratory ID:	03-007-22					
Arsenic	ND	11	EPA 6010D	3-8-18	3-8-18	
Cadmium	ND	0.55	EPA 6010D	3-8-18	3-8-18	
Chromium	15	0.55	EPA 6010D	3-8-18	3-8-18	
Lead	97	5.5	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.27	EPA 7471B	3-5-18	3-5-18	
Client ID:	PP23-10					
Laboratory ID:	03-007-23					
Arsenic	ND	13	EPA 6010D	3-8-18	3-8-18	
Cadmium	ND	0.67	EPA 6010D	3-8-18	3-8-18	
Chromium	13	0.67	EPA 6010D	3-8-18	3-8-18	
Lead	ND	6.7	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.34	EPA 7471B	3-5-18	3-5-18	
Client ID:	PP23-15					
Laboratory ID:	03-007-24					
Arsenic	ND	15	EPA 6010D	3-8-18	3-8-18	
Cadmium	ND	0.76	EPA 6010D	3-8-18	3-8-18	
Chromium	18	0.76	EPA 6010D	3-8-18	3-8-18	
Lead	220	7.6	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.38	EPA 7471B	3-5-18	3-5-18	

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B

Matrix: Soil

Units: mg/kg (ppm)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP21-2					
Laboratory ID:	03-007-25					
Arsenic	ND	13	EPA 6010D	3-8-18	3-9-18	
Cadmium	0.90	0.66	EPA 6010D	3-8-18	3-9-18	
Chromium	30	0.66	EPA 6010D	3-8-18	3-9-18	
Lead	740	6.6	EPA 6010D	3-8-18	3-9-18	
Mercury	0.87	0.33	EPA 7471B	3-5-18	3-5-18	
Client ID:	PP21-6					
Laboratory ID:	03-007-26					
Arsenic	17	10	EPA 6010D	3-8-18	3-8-18	
Cadmium	27	1.0	EPA 6010D	3-8-18	3-8-18	
Chromium	74	1.0	EPA 6010D	3-8-18	3-8-18	
Lead	2800	10	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.51	EPA 7471B	3-5-18	3-5-18	
Client ID:	PP21-10					
Laboratory ID:	03-007-27					
Arsenic	ND	16	EPA 6010D	3-8-18	3-8-18	
Cadmium	6.2	0.81	EPA 6010D	3-8-18	3-8-18	
Chromium	23	0.81	EPA 6010D	3-8-18	3-8-18	
Lead	180	8.1	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.41	EPA 7471B	3-5-18	3-5-18	
OII ID	DD00.0					
Client ID:	PP20-2					
Laboratory ID:	03-007-29	10	EDA 00465	0.0.40	0.0.10	
Arsenic	ND	12	EPA 6010D	3-8-18	3-8-18	
Cadmium	ND	0.58	EPA 6010D	3-8-18	3-8-18	
Chromium	15	0.58	EPA 6010D	3-8-18	3-8-18	
Lead 	28	5.8	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.29	EPA 7471B	3-5-18	3-5-18	

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B

Matrix: Soil

Offits. Hig/Kg (ppiff)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP20-5					
Laboratory ID:	03-007-30					
Arsenic	ND	15	EPA 6010D	3-8-18	3-8-18	
Cadmium	3.7	0.73	EPA 6010D	3-8-18	3-8-18	
Chromium	53	0.73	EPA 6010D	3-8-18	3-8-18	
Lead	630	7.3	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.37	EPA 7471B	3-7-18	3-7-18	
Client ID:	PP20-10					
Laboratory ID:	03-007-31					
Arsenic	ND	13	EPA 6010D	3-8-18	3-8-18	
Cadmium	ND	0.64	EPA 6010D	3-8-18	3-8-18	
Chromium	22	0.64	EPA 6010D	3-8-18	3-8-18	
Lead	270	6.4	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.32	EPA 7471B	3-7-18	3-7-18	
Client ID:	PP18-3					
Laboratory ID:	03-007-33					
Arsenic	ND	12	EPA 6010D	3-8-18	3-8-18	
Cadmium	2.4	0.61	EPA 6010D	3-8-18	3-8-18	
Chromium	63	0.61	EPA 6010D	3-8-18	3-8-18	
Lead	130	6.1	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.31	EPA 7471B	3-7-18	3-7-18	
Client ID:	PP18-5					
Laboratory ID:	03-007-34					
Arsenic	ND	20	EPA 6010D	3-8-18	3-8-18	
Cadmium	3.1	0.99	EPA 6010D	3-8-18	3-8-18	
Chromium	3.1 39	0.99	EPA 6010D	3-0-10 3-8-18	3-8-18	
Lead	230	0.99 9.9	EPA 6010D EPA 6010D	3-8-18 3-8-18		
					3-8-18	
Mercury	ND	0.49	EPA 7471B	3-7-18	3-7-18	

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B

Matrix: Soil

Offits. Hig/Kg (ppHI)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP18-10					
Laboratory ID:	03-007-35					
Arsenic	ND	13	EPA 6010D	3-8-18	3-8-18	
Cadmium	1.2	0.67	EPA 6010D	3-8-18	3-8-18	
Chromium	40	0.67	EPA 6010D	3-8-18	3-8-18	
Lead	97	6.7	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.33	EPA 7471B	3-7-18	3-7-18	
Client ID:	PP17-1					
Laboratory ID:	03-007-37					
Arsenic	19	13	EPA 6010D	3-8-18	3-8-18	
Cadmium	4.0	0.64	EPA 6010D	3-8-18	3-8-18	
Chromium	57	0.64	EPA 6010D	3-8-18	3-8-18	
Lead	7300	32	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.32	EPA 7471B	3-7-18	3-7-18	
Client ID:	PP17-5					
Laboratory ID:	03-007-38					
Arsenic	ND	12	EPA 6010D	3-8-18	3-8-18	
Cadmium	4.5	0.61	EPA 6010D	3-8-18	3-8-18	
Chromium	59	0.61	EPA 6010D	3-8-18	3-8-18	
Lead	380	6.1	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.30	EPA 7471B	3-7-18	3-7-18	
Client ID:	PP17-10					
Laboratory ID:	03-007-39					
Arsenic	ND	12	EPA 6010D	3-8-18	3-8-18	
Cadmium	1.6	0.58	EPA 6010D	3-8-18	3-8-18	
Chromium	24	0.58	EPA 6010D	3-8-18	3-8-18	
Lead	29	5.8	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.29	EPA 7471B	3-7-18	3-7-18	

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B

Matrix: Soil

Units: mg/Kg (ppm)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP16-1					
Laboratory ID:	03-007-41					
Arsenic	ND	11	EPA 6010D	3-8-18	3-8-18	
Cadmium	ND	0.53	EPA 6010D	3-8-18	3-8-18	
Chromium	20	0.53	EPA 6010D	3-8-18	3-8-18	
Lead	17	5.3	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.27	EPA 7471B	3-7-18	3-7-18	
Client ID:	PP16-11					
Laboratory ID:	03-007-42					
Arsenic	ND	12	EPA 6010D	3-8-18	3-8-18	
Cadmium	1.2	0.61	EPA 6010D	3-8-18	3-8-18	
Chromium	24	0.61	EPA 6010D	3-8-18	3-8-18	
Lead	200	6.1	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.31	EPA 7471B	3-7-18	3-7-18	
Client ID:	PP14-5					
Laboratory ID:	03-007-45					
Arsenic	ND	12	EPA 6010D	3-8-18	3-8-18	
Cadmium	ND	0.62	EPA 6010D	3-8-18	3-8-18	
Chromium	12	0.62	EPA 6010D	3-8-18	3-8-18	
Lead	ND	6.2	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.31	EPA 7471B	3-7-18	3-7-18	
	5 54446					
Client ID:	PP14-12					
_aboratory ID:	03-007-46					
Arsenic	ND	13	EPA 6010D	3-8-18	3-8-18	
Cadmium	ND	0.65	EPA 6010D	3-8-18	3-8-18	
Chromium	11	0.65	EPA 6010D	3-8-18	3-8-18	
Lead	ND	6.5	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.32	EPA 7471B	3-9-18	3-9-18	

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B

Matrix: Soil

Units: mg/Kg (ppm)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP11-5					
Laboratory ID:	03-007-49					
Arsenic	19	13	EPA 6010D	3-8-18	3-8-18	
Cadmium	1.7	0.66	EPA 6010D	3-8-18	3-8-18	
Chromium	39	0.66	EPA 6010D	3-8-18	3-8-18	
Lead	840	6.6	EPA 6010D	3-8-18	3-8-18	
Mercury	0.41	0.33	EPA 7471B	3-9-18	3-9-18	
Client ID:	PP11-10					
Laboratory ID:	03-007-50					
Arsenic	ND	13	EPA 6010D	3-8-18	3-8-18	
Cadmium	0.72	0.63	EPA 6010D	3-8-18	3-8-18	
Chromium	36	0.63	EPA 6010D	3-8-18	3-8-18	
Lead	480	6.3	EPA 6010D	3-8-18	3-8-18	
Mercury	1.2	0.31	EPA 7471B	3-9-18	3-9-18	
Client ID:	PP10-11					
Laboratory ID:	03-007-52					
Arsenic	ND	16	EPA 6010D	3-8-18	3-8-18	
Cadmium	ND	0.82	EPA 6010D	3-8-18	3-8-18	
Chromium	22	0.82	EPA 6010D	3-8-18	3-8-18	
Lead	15	8.2	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.41	EPA 7471B	3-9-18	3-9-18	
Client ID:	PP10-17					
Laboratory ID:	03-007-53					
Arsenic	ND	12	EPA 6010D	3-8-18	3-8-18	
Cadmium	ND	0.62	EPA 6010D	3-8-18	3-8-18	
Chromium	15	0.62	EPA 6010D	3-8-18	3-8-18	
Lead	ND	6.2	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.31	EPA 7471B	3-9-18	3-9-18	

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B

Matrix: Soil

Offits. Hig/Ng (ppiff)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP27-7					
Laboratory ID:	03-007-55					
Arsenic	ND	15	EPA 6010D	3-8-18	3-8-18	
Cadmium	0.88	0.73	EPA 6010D	3-8-18	3-8-18	
Chromium	28	0.73	EPA 6010D	3-8-18	3-8-18	
Lead	180	7.3	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.36	EPA 7471B	3-9-18	3-9-18	
Client ID:	PP27-10					
Laboratory ID:	03-007-56					
Arsenic	ND	12	EPA 6010D	3-8-18	3-8-18	
Cadmium	ND	0.61	EPA 6010D	3-8-18	3-8-18	
Chromium	12	0.61	EPA 6010D	3-8-18	3-8-18	
Lead	ND	6.1	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.30	EPA 7471B	3-9-18	3-9-18	
Client ID:	PP30-5					
Laboratory ID:	03-007-59					
Arsenic	ND	13	EPA 6010D	3-8-18	3-8-18	
Cadmium	0.83	0.64	EPA 6010D	3-8-18	3-8-18	
Chromium	16	0.64	EPA 6010D	3-8-18	3-8-18	
Lead	31	6.4	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.32	EPA 7471B	3-9-18	3-9-18	
Client ID:	PP30-10					
Laboratory ID:	03-007-60					
Arsenic	ND	11	EPA 6010D	3-8-18	3-8-18	
Cadmium	ND	0.57	EPA 6010D	3-8-18	3-8-18	
Chromium	14	0.57	EPA 6010D	3-8-18	3-8-18	
Lead	ND	5.7	EPA 6010D	3-8-18	3-8-18	
Mercury	ND	0.28	EPA 7471B	3-9-18	3-9-18	

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B

Matrix: Soil

5 5 ,				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP24-7					
Laboratory ID:	03-007-63					
Arsenic	ND	14	EPA 6010D	3-9-18	3-9-18	
Cadmium	2.3	0.69	EPA 6010D	3-9-18	3-9-18	
Chromium	18	0.69	EPA 6010D	3-9-18	3-9-18	
Lead	480	6.9	EPA 6010D	3-9-18	3-9-18	
Mercury	ND	0.35	EPA 7471B	3-9-18	3-9-18	
Client ID:	PP24-10					
Laboratory ID:	03-007-64					
Arsenic	ND	12	EPA 6010D	3-9-18	3-9-18	
Cadmium	ND	0.58	EPA 6010D	3-9-18	3-9-18	
Chromium	10	0.58	EPA 6010D	3-9-18	3-9-18	
Lead	ND	5.8	EPA 6010D	3-9-18	3-9-18	
Mercury	ND	0.29	EPA 7471B	3-9-18	3-9-18	

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B METHOD BLANK QUALITY CONTROL

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0309SM1					
Arsenic	ND	5.0	EPA 6010D	3-9-18	3-9-18	
Cadmium	ND	0.50	EPA 6010D	3-9-18	3-9-18	
Chromium	ND	0.50	EPA 6010D	3-9-18	3-9-18	
Lead	ND	5.0	EPA 6010D	3-9-18	3-9-18	
Laboratory ID:	MB0309S1					
Mercury	ND	0.25	EPA 7471B	3-9-18	3-9-18	
Laboratory ID:	MB0308SM2					
Arsenic	ND	10	EPA 6010D	3-8-18	3-8-18	
Cadmium	ND	0.50	EPA 6010D	3-8-18	3-8-18	
Chromium	ND	0.50	EPA 6010D	3-8-18	3-8-18	
Lead	ND	5.0	EPA 6010D	3-8-18	3-8-18	
Laboratory ID:	MB0307S2					
Mercury	ND	0.25	EPA 7471B	3-7-18	3-7-18	
Laboratory ID:	MB0308SM1					
Arsenic	ND	5.0	EPA 6010D	3-8-18	3-8-18	
Cadmium	ND	0.50	EPA 6010D	3-8-18	3-8-18	
Chromium	ND	0.50	EPA 6010D	3-8-18	3-8-18	
Lead	ND	5.0	EPA 6010D	3-8-18	3-8-18	
Laboratory ID:	MB0305S1					
Mercury	ND	0.25	EPA 7471B	3-5-18	3-5-18	

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B DUPLICATE QUALITY CONTROL

Matrix: Soil

	,				Source	Percent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Recovery	Limits	RPD	Limit	Flags
DUPLICATE										
Laboratory ID:	03-00)7-11								
	ORIG	DUP								
Arsenic	ND	ND	NA	NA		NA	NA	NA	20	
Cadmium	ND	ND	NA	NA		NA	NA	NA	20	
Chromium	16.5	15.5	NA	NA		NA	NA	6	20	
Lead	ND	ND	NA	NA		NA	NA	NA	20	
Laboratory ID:	03-03	36-07								
Mercury	ND	ND	NA	NA		NA	NA	NA	20	
Laboratory ID:	03-00)7-41								
	ORIG	DUP								
Arsenic	ND	ND	NA	NA		NA	NA	NA	20	
Cadmium	ND	ND	NA	NA		NA	NA	NA	20	
Chromium	18.8	19.3	NA	NA		NA	NA	2	20	
Lead	15.8	12.9	NA	NA		NA	NA	20	20	
Laboratory ID:	02-27	' 1-21								
Mercury	ND	ND	NA	NA		NA	NA	NA	20	
Laboratory ID:	03-00	7-09								
	ORIG	DUP								
Arsenic	ND	ND	NA	NA		NA	NA	NA	20	
Cadmium	ND	ND	NA	NA		NA	NA	NA	20	
Chromium	12.6	13.0	NA	NA		NA	NA	4	20	
Lead	9.65	10.9	NA	NA		NA	NA	12	20	
Laboratory ID:	03-00	7-09								
Mercury	ND	ND	NA	NA		NA	NA	NA	20	

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B MS/MSD QUALITY CONTROL

Matrix: Soil

			Source	Per	cent	Recovery		RPD			
Analyte	Res	sult	Spike	Level	Result	Rec	overy	Limits	RPD	Limit	Flags
MATRIX SPIKES											
Laboratory ID:	03-00	07-11									
	MS	MSD	MS	MSD		MS	MSD				
Arsenic	101	99.5	100	100	ND	101	100	75-125	1	20	
Cadmium	46.0	46.9	50.0	50.0	ND	92	94	75-125	2	20	
Chromium	109	115	100	100	16.5	93	98	75-125	5	20	
Lead	230	236	250	250	ND	92	94	75-125	2	20	
Laboratory ID:	03-03	36-07									
Mercury	0.564	0.529	0.500	0.500	0.0553	102	95	80-120	6	20	
Laboratory ID:	03-00	17 -41									
Laboratory ID.	MS	MSD	MS	MSD		MS	MSD				
Arsenic	99.5	101	100	100	ND	100	101	75-125	1	20	
Cadmium	48.4	48.2	50.0	50.0	ND	97	96	75-125	1	20	
Chromium	115	117	100	100	18.8	96	98	75-125	1	20	
Lead	250	249	250	250	15.8	94	93	75-125	0	20	
Laboratory ID:	02-2	72-21									
Mercury	0.605	0.590	0.500	0.500	0.0364	114	111	80-120	3	20	
Laboratory ID:	03-00	07-09									
	MS	MSD	MS	MSD		MS	MSD				
Arsenic	98.0	99.0	100	100	ND	98	99	75-125	1	20	
Cadmium	45.8	45.7	50.0	50.0	ND	92	91	75-125	0	20	
Chromium	107	107	100	100	12.6	95	94	75-125	0	20	
Lead	234	236	250	250	9.65	90	90	75-125	1	20	
Laboratory ID:	03-00	07-09									
Mercury	0.568	0.575	0.500	0.500	0.0317	107	109	80-120	1	20	
							107 109 00-				

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B

Matrix: Soil

omie. mg/rtg (ppm/				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP32-10					
Laboratory ID:	03-007-11					
Arsenic	ND	12	EPA 6010D	3-9-18	3-9-18	
Cadmium	ND	0.61	EPA 6010D	3-9-18	3-9-18	
Chromium	20	0.61	EPA 6010D	3-9-18	3-9-18	
Lead	ND	6.1	EPA 6010D	3-9-18	3-9-18	
Mercury	ND	0.31	EPA 7471B	3-12-18	3-12-18	
Client ID:	PP21-15					
Laboratory ID:	03-007-28					
Arsenic	ND	14	EPA 6010D	3-9-18	3-9-18	
Cadmium	ND	1.4	EPA 6010D	3-9-18	3-9-18	
Chromium	20	1.4	EPA 6010D	3-9-18	3-9-18	
Lead	ND	14	EPA 6010D	3-9-18	3-9-18	
Mercury	ND	0.68	EPA 7471B	3-12-18	3-12-18	
Client ID:	PP20-15					
Laboratory ID:	03-007-32					
Arsenic	ND	14	EPA 6010D	3-9-18	3-9-18	
Cadmium	ND	0.70	EPA 6010D	3-9-18	3-9-18	
Chromium	15	0.70	EPA 6010D	3-9-18	3-9-18	
Lead	330	7.0	EPA 6010D	3-9-18	3-9-18	
Mercury	ND	0.35	EPA 7471B	3-12-18	3-12-18	

Project: 17-06520-000

TOTAL METALS EPA 6010D/7471B QUALITY CONTROL

Matrix: Soil

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0309SM1					
Arsenic	ND	5.0	EPA 6010D	3-9-18	3-9-18	
Cadmium	ND	0.50	EPA 6010D	3-9-18	3-9-18	
Chromium	ND	0.50	EPA 6010D	3-9-18	3-9-18	
Lead	ND	5.0	EPA 6010D	3-9-18	3-9-18	
Laboratory ID:	MB0312S1					
Mercury	ND	0.25	EPA 7471B	3-12-18	3-12-18	

					Source	Pe	rcent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Rec	overy	Limits	RPD	Limit	Flags
DUPLICATE											,
Laboratory ID:	03-00	07-11									
	ORIG	DUP									,
Arsenic	ND	ND	NA	NA		ı	NA	NA	NA	20	
Cadmium	ND	ND	NA	NA		ı	NA	NA	NA	20	
Chromium	16.5	15.5	NA	NA		ı	NA	NA	6	20	
Lead	ND	ND	NA	NA		I	NA	NA	NA	20	
Laboratory ID:	03-0	75-02									
Mercury	ND	ND	NA	NA			NA	NA	NA	20	
MATRIX SPIKES											
Laboratory ID:	03-00	07-11									
	MS	MSD	MS	MSD		MS	MSD				
Arsenic	101	99.5	100	100	ND	101	100	75-125	1	20	
Cadmium	46.0	46.9	50.0	50.0	ND	92	94	75-125	2	20	
Chromium	109	115	100	100	16.5	93	98	98 75-125		20	
Lead	230	236	250	250	ND	92	94	75-125	2	20	
Laboratory ID:	03-075-02										
Mercury	0.545	0.547	0.500	0.500	0.0159	106	106	80-120	0	20	

Project: 17-06520-000

% MOISTURE

Date Analyzed: 3-2&7-18

Client ID	Lab ID	% Moisture
PP28-8	03-007-02	20
PP28-10	03-007-03	19
PP29-3	03-007-05	11
PP29-5	03-007-06	23
PP29-10	03-007-07	25
PP32-4	03-007-09	20
PP32-7	03-007-10	19
PP32-10	03-007-11	18
PP31-3	03-007-13	11
PP31-11	03-007-15	20
PP33-3	03-007-17	11
PP33-5	03-007-18	16
PP33-10	03-007-19	26
PP23-2	03-007-21	15
PP23-5	03-007-22	9
PP23-10	03-007-23	26
PP23-15	03-007-24	34
PP21-2	03-007-25	25
PP21-6	03-007-26	51
PP21-10	03-007-27	39
PP21-15	03-007-28	63
PP20-2	03-007-29	13
PP20-5	03-007-30	32
PP20-10	03-007-31	22
PP20-15	03-007-32	28
PP18-3	03-007-33	19
PP18-5	03-007-34	49

Project: 17-06520-000

% MOISTURE

Date Analyzed: 3-2&7-18

Client ID	Lab ID	% Moisture
PP18-10	03-007-35	25
PP17-1	03-007-37	22
PP17-5	03-007-38	18
PP17-10	03-007-39	14
PP16-1	03-007-41	6
PP16-11	03-007-42	19
PP14-5	03-007-45	20
PP14-12	03-007-46	23
PP11-5	03-007-49	24
PP11-10	03-007-50	20
PP10-11	03-007-52	39
PP10-17	03-007-53	20
PP27-7	03-007-55	31
PP27-10	03-007-56	17
PP30-5	03-007-59	22
PP30-10	03-007-60	12
PP24-7	03-007-63	28
PP24-10	03-007-64	14



Data Qualifiers and Abbreviations

- A Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
- B The analyte indicated was also found in the blank sample.
- C The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
- E The value reported exceeds the quantitation range and is an estimate.
- F Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
- H The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
- I Compound recovery is outside of the control limits.
- J The value reported was below the practical quantitation limit. The value is an estimate.
- K Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
- L The RPD is outside of the control limits.
- M Hydrocarbons in the gasoline range are impacting the diesel range result.
- M1 Hydrocarbons in the gasoline range (toluene-naphthalene) are present in the sample.
- N Hydrocarbons in the lube oil range are impacting the diesel range result.
- N1 Hydrocarbons in diesel range are impacting lube oil range results.
- O Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
- P The RPD of the detected concentrations between the two columns is greater than 40.
- Q Surrogate recovery is outside of the control limits.
- S Surrogate recovery data is not available due to the necessary dilution of the sample.
- T The sample chromatogram is not similar to a typical
- U The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- U1 The practical quantitation limit is elevated due to interferences present in the sample.
- V Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
- W Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
- X Sample extract treated with a mercury cleanup procedure.
- X1- Sample extract treated with a sulfuric acid/silica gel cleanup procedure.
- Y The calibration verification for this analyte exceeded the 20% drift specified in method 8260C, and therefore the reported result should be considered an estimate. The overall performance of the calibration verification standard met the acceptance criteria of the method.

7 -

ND - Not Detected at PQL

PQL - Practical Quantitation Limit

RPD - Relative Percent Difference





Chain of Custody

	1	XV	7
Page _	1	of_	

	14648 NE !	aboratory Testing Services 95th Street • Redmond, WA 98052		Turnaround Req (in working da			L	abo	orat	ory	Nur	nbe	er:	0	3	-() (7									
Project Project Sample	Number: 17-06526 Name: Pull C Po	0-000	□ 2	Days [andard (7 Days) PH analysis 5 Da (other)	1 Day	Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX	NWTPH-Gx	NWTPH-Dx (☐ Acid / SG Clean-up)	Volatiles 8260C	Halogenated Volatiles 8260C	EDB EPA 8011 (Waters Only)	Semivolatiles 8270D/SIM (with low-level PAHs)	PAHS 8270D/SIM (low-level)	CBs 8082A	Organochlorine Pesticides 8081B	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	TCLP Metals	HEM (oil and grease) 1664A				% Moisture
1	PP28-		3.11		Soil	2	_	_	_	_				0, 0	40			0	0		-		_			+	%
2	PP7-8 -	8	1	1035	1	2	X								X						X						X
3	PP28-	10		1040		Z	X								X						X						X
4	PPZ8-	B		1045		2																					
5	PP29-			1140		2	X			(X)					X	8					X						X
0	PP29-:	2-10-1		1145		2	X								X						X						X
7	PP29-1	0		1150		2	X								X						X						X
8	PPZ9-	15		1155		Z																					
9	PP32-L			9120		2	X							1	X						X						X
[0]	PP32-7			9130	1	2	X			8				,	X	8					X						X
		Signature		Company		SI-II-		Date			Time	1.3			A CONTRACTOR OF THE PARTY OF TH	s/Spe			-	ALC: UNKNOWN							miles - 1
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Chain of Custody

Page 2 of 7

	Analytical Laboratory Testing Services 14648 NE 95th Street • Redmond, WA 98052		Turnaround Req (in working da			L	abo	orat	tory	Nu	umb	er:	0	3	- (0 (7								
Project Project Sample	Number: 17-06520-000 Name: Paulic Pork Manager: Bruce Carpenter			1 Day	r of Containers	HCID	NWTPH-Gx/BTEX	-Gx	NWTPH-Dx (Acid / SG Clean-up)	Volatiles 8260C	Halogenated Volatiles 8260C	EDB EPA 8011 (Waters Only)	Semivolatiles 8270D/SIM (with low-level PAHs)	CAHS 8270D/SIM (low-level) CPAHS		Organochlorine Pesticides 8081B	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	etals	and grease) 1664A			ure
Lab ID	Sample Identification	Date Sample		Matrix	Number	NWTPH-HCID	NWTP	NWTPH-Gx	NWTP	Volatile	Haloger	EDB EF	Semivo (with lo	を表	PCBs 8	Organo	Organo	Chlorina	Total RC	Total M	TCLP Metals	HEM (oil			% Moisture
11	PP32-10	3.1.1	8 9:35	Soil	2		(8						(\otimes	X					X) .				X
12	PP3Z-15		940		2																Z				
13	PP31-3		1000		2	X								X						X	X				X
14	PP31-5		1005		2																/1				
15	PP31-11		1010		2	X								X						X					X
16	PP31-15		1015		2																, 2				
17	PP33-3		1105		2	X	X)	8					X	X					X	1				X
18	PP33-5		1110		2	X								X						x	X				X
19	PP33 - 10		115		2	X								X						X	X				X
20	PP33-15	1	1120	V	2																1				
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Revie	wed/Date		Reviewed/Dat	te										_	_	_		_	_					EDDs)	

OnSite Environmental Inc.

Chain of Custody

Page 3 of 7

Analytical Laboratory Testing Services 14648 NE 95th Street • Redmond, WA 98052	Turnaround Request (in working days)				L	Laboratory Number: 03-007												Ì					
Phone: (425) 883-3881 • www.onsite-env.com Company: HUVEVA Project Number: Project Name: Project Manager: Sampled by: Sampled by:	Same 2 Day Stand (TPH	,	1 Day 3 Days	er of Containers	NWTPH-HCID	NWTPH-Gx/BTEX		NWTPH-Dx (☐ Acid / SG Clean-up)	Volatiles 8260C Halogenated Volatiles 8260C	EDB EPA 8011 (Waters Only)	Semivolatiles 8270D/SIM	w-level PAHs) 2270D/SIM (low-level)	PCBs 8082A	Organochlorine Pesticides 8081B	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	TCLP Metals	il and grease) 1664A			ture
ab ID Sample Identification	Date Sampled	Time Sampled	Matrix	Number	NWTP	NWTP	NWTPH-Gx	_	Volatil	EDB E	Semiv	_	-	Organi	Organi	Chlorir	Total R	Total N	TCLP	HEM (oil		70	% Moisture
U PP23-2	2.28.18	1400	Soil	2	X	(X)		\otimes				X	X					X					X
12 PP23-S	1	1405		2	X							X						X)	<
23 PP23-10		1415		2	X							X						X)	0
24 PP23-15		1420		2	X							X						X				k	1
25 PP21-2		1330		Z	χ			$\widehat{\mathbb{X}}$				X	X					X				X	
210 PPZI-10		1335		2	X			\otimes				X	0					X				R	
27 pp21-10		1345		Z	X	8		8				X	(X)					X				1	X
28 PP21-K		1350		2				X				0	(3)				((X)					X
19 PP20-2		1250		2	X			D				X	(3)					X)	1
30 PPZO-S	V	1255	V	2	X			8)			X	(V)					X				7	
Relinquished Received Received Received Received		Herve	va SE			Date 3			1 400 1	10	Co	omme	ints/Sp	ecial	Instr	uction	18				1		
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Reviewed/Date		Reviewed/Date									Data Package: Standard ☐ Level III ☐ Level IV ☐ Chromatograms with final report ☐ Electronic Data Deliverables (EDDs) ☐											-	



Chain of Custody

Page 4 of 3

	Analytical Laboratory Testing Services 14648 NE 95th Street • Redmond, WA 98052		Turnaround Request (in working days)					Laboratory Number: 03-007																	
Project Project Samp	tonera IT-010520-000 St Name: PACADIC POWK St Manager: BINCLE CAUPENTU BINCLE CAUPENTU BINCLE SUPERIOR BINCLE	Date	ays adard (7 Days) I analysis 5 Da (other)	1 Day 3 Days	Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX	NWTPH-Gx	NWTPH-Dx (☐ Acid / SG Clean-up)	Volatiles 8260C	Halogenated Volatiles 8260C	EDB EPA 8011 (Waters Only)	Semivolatiles 8270D/SIM (with low-level PAHs) PAHS 8270D/SIM (low-level)	2011	Organochlorine Pesticides 8081B	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	TCLP Metals	HEM (oil and grease) 1664A			% Moisture	Moisture
Ab ID		Sampled 2:28 to	Sampled 1305	SO!	7	X	Z	Z	(X)		I	ш	4	X	0	0	0	P	X	Ť.	I	+		3	1
32	22.001	1	1310	1	2				X				Q	Q					X					()	2
	PP13-3		1150		2	x							X						X					X	1
	PP18-S		1200		2	X							X						X					×	1
	PP18-10		1205		2	X							X						X)	X
36	PP18-15		1210		Z																				
37	PP17-1		1100		2	X							X						X					7	4
38	PPI7-S		1110		2	X							X						X						X
39			1120		2	X							X						X					1	X
40	PP17-18		1128	V	2																				
Relin	signature squished Shipman Market		ompany Hevra	va	,		Date		78	Time	e 34	0	Comme	nts/Sp	ecial	Instr	uctio	ns							
Rece			(0)	275			7	1.1	18		30														
Relir	nquished						21		5		2		1												
Rece	eived																								
Relin	nquished																								
Rece	sived											Data Package: Standard ☐ Level III ☐ Level IV ☐													
Revi	ewed/Date		Reviewed/Date										Chromatograms with final report ☐ Electronic Data Deliverables (EDDs) ☐												



Page S of 7

	Analytical Laboratory Testing Services 14648 NE 95th Street • Redmond, WA 98052	Turi (ir	naround Req working da	uest ys)		L	abo	rato	ry N	luml	oer:		03	} -	0	07	7								
Project Project Sample	Number: 17-00520-000 Name: POULLIC POWN Manager: BYLLE CALPOLUTEV	Same 2 Day Stand (TPH		1 Day	Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX	NWTPH-Gx	Notatiles 8260C	Halogenated Volatiles 8260C	EDB EPA 8011 (Waters Only)		PAHS 8270D/SIM (low-level)		Organochlorine Pesticides 8081B	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	TCLP Metals	HEM (oil and grease) 1664A				% Moisture
Lab ID	PPIG-1	2.28.18			2		(X)		(X	1	Ш	0, 2	X	0		0	0	-	X	-	_				X
42	PP16-11		1045		Z	X							X						X						X
43	PP 16-15		1050		Z	-					T														F
44	PP14-1		950		Z																				
45	PP14-S		955		2	X							X						X						X
	PP14-12		1000		2	X							X						X						Y
	PP14-14		1005		2														,,						^
48	PP11-2		915		2																				
49	PPII - 5		920		2	X							X						X						X
50	PP11-10	/	925	V	2	X							χ						X						X
	Signature		mpany	0114			Date			īme	1.	Con	mmei	nts/Sp	ecial	Instr	uction	ns							
Relino	quished Maywalla	K	HUVA	NA)		3	111		131	10	-													
	quished	> (980				31	110	1	134	0	-													
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Revie	wed/Date		Reviewed/Da	te					-			Chr	omat	ogran	ns wi	th fina	al rep	ort [Ele	ctroni	c Data	a Delive	erables	(EDDs)	



Page 0 of 7

Analytical Laboratory Testing Services 14648 NE 95th Street • Redmond, WA 98052		naround Req working da			La	abo	rato	ry N	Num	ber:		0	3 -	0	07	7								
Phone: (425) 883-3881 * www.onsite-env.com Company: Project Number: Project Name: Project Manager: BYCE CAMPUTU Sampled by BYWAN BYWAN Sampled by	Same 2 Day		1 Day 3 Days	Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX		NW I PH-DX (Acid / SG Clean-up)	voiatiles 6200C Halogenated Volatiles 8260C	EDB EPA 8011 (Waters Only)	Semivolatiles 8270D/SIM	ow-level PAHs) 8270D/SIM (low-level)	PCBs 8082A	Organochlorine Pesticides 8081B	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	TCLP Metals	HEM (oil and grease) 1664A				sture
Lab ID Sample Identification	Date Sampled	Time Sampled	Matrix	Numl	NWT	NWT	HWN	T NN	Halog	EDB	Semi	with PAHS	38	Organ	Organ	Chlor	Total	Total	TCLP	HEM				% Moisture
51 PP11-1S	2.28.18		Soil	2				4		-	+	+	+	-	_	_								
52 PP10-11		840		2	X			4			-	X	5					X						X
53 PP10-17		8402		2	X							X						X						X
54 PPZA -4		1545		2																				
SS PP 27-7		1550		2	X							X						X						X
Sie PPZ9-10		1600		2	X							X						X						X
S7 PP27-15		1610		Z																				
58 PP20-3		1505		2																				
59 PP30-S		1510		2	X		(X				X	Q	5				X						X
60 PP30-10		1520	/	2	X							>						X						X
Signature		mpany			11	Date		1	Time		C	omm	ents/S	pecia	Inst	ructio	ns							
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Received											Da	ata P	ackag	e: St	tanda	rd 🗆	Le	vel III		Leve	I IV 🗆			
Reviewed/Date		Reviewed/Da	te								Ch	roma	atogra	ms w	ith fir	nal rep	oort [] Ele	ctron	ic Data	a Delive	rables (E	DDs)]

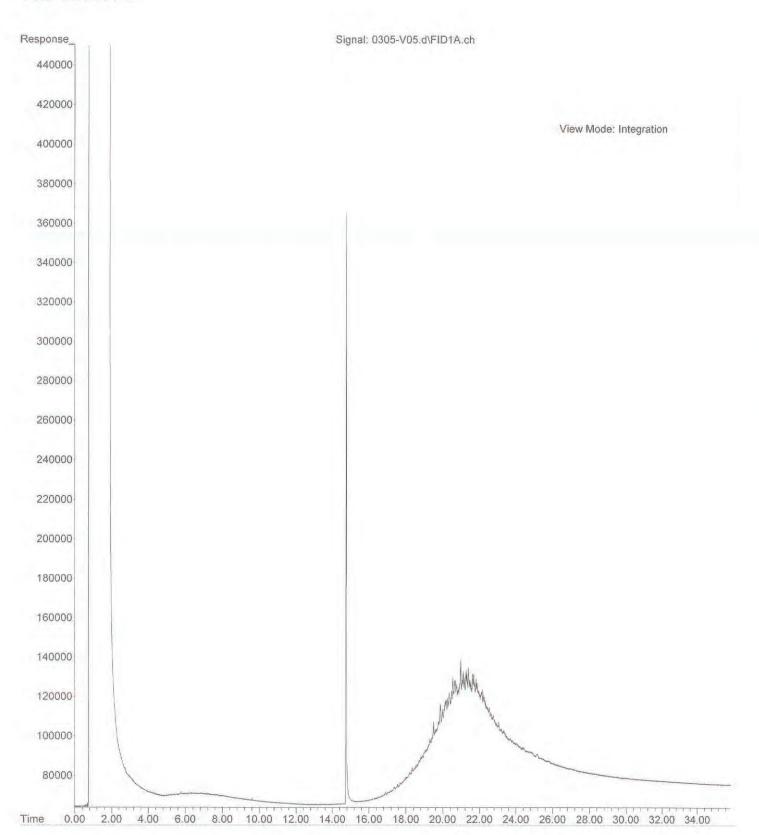


Page 7 of 7

	Analytical Laboratory Testing Services 14648 NE 95th Street • Redmond, WA 98052	Tur (i	rnaround Rec n working da	juest lys)		La	abo	rato	ry	Num	ber:	(13	-	nı	17									
Projec	Phone: (425) 883-3881 · www.onsite-env.com Herrywa t Number: 17-04520-060 t Name: Paulic Paul t Manager: Bruce Carpentel ed by: Brunna Bland	Sam 2 Da Stan (TPI-	,	1 Day	Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX		NVV I PH-DX (Acid / SG Ciean-up)	Volatiles 8260C Halogenated Volatiles 8260C	EDB EPA 8011 (Waters Only)	Semivolatiles 8270D/SIM (with low-level PAHs)			Organochlorine Pesticides 8081B	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	TCLP Metals	HEM (oil and grease) 1664A				sture
Lab ID	Sample Identification	Date Sampled	Time Sampled	Matrix	Num	NWT	NWT	NWT	I AAN	Volati	EDB 6	Semiv (with 1	野	PCBs	Organ	Organ	Chlori	Total F	Total	TCLP	HEM				% Moisture
[Ol	PP30-15	2.28.18	1525	301	2																				
102	PP24-2		1436		2																				
63	PP24-7		1435		2	X	(8)	(3				X	X)				X						X
104	PPZ4-10	V	1445	V	2	X							X						X						X
					Н						-												+	_	
	Signature	C	ompany				Date			Time		Cor	mmer	ıts/Sp	ecial	Instr	uctio	ns							-2-911
Rece			Herrie	NE SE			31	1/4	3	13	40														
	quished								4																
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	quished								1												_				
Rece			B 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1					-				-	_	_	_						Leve				
Hevie	ewed/Date		Reviewed/Da	ite								Chr	omate	ogran	ns wi	th fin	al rep	ort [Ele	ectron	ic Data	a Deliv	erables	(EDDs)	

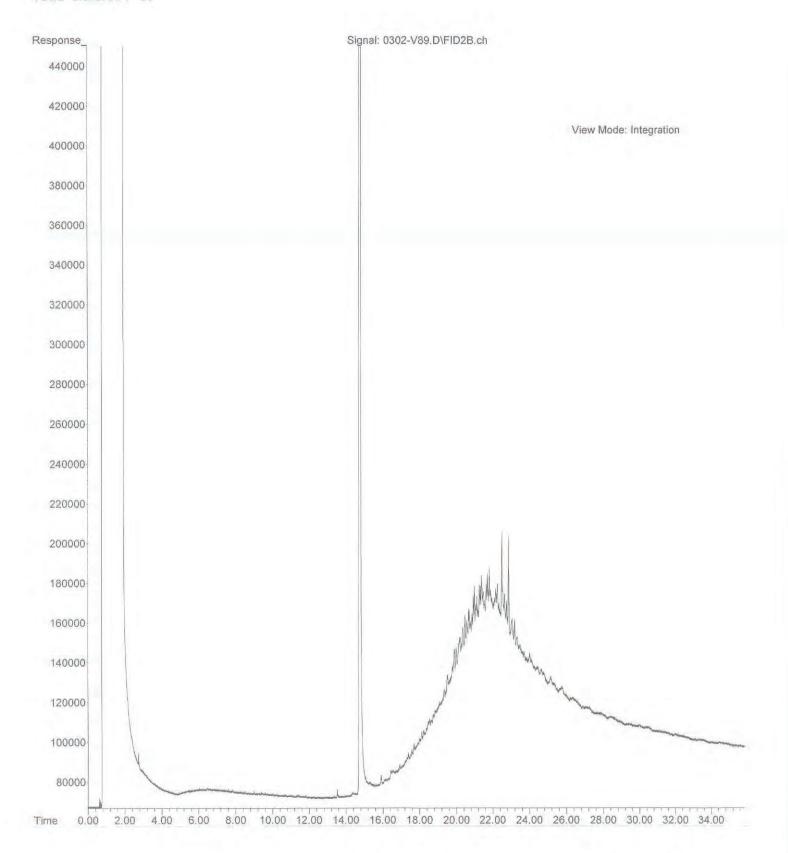
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Instrument : Vigo

Sample Name: 03-007-05 20X

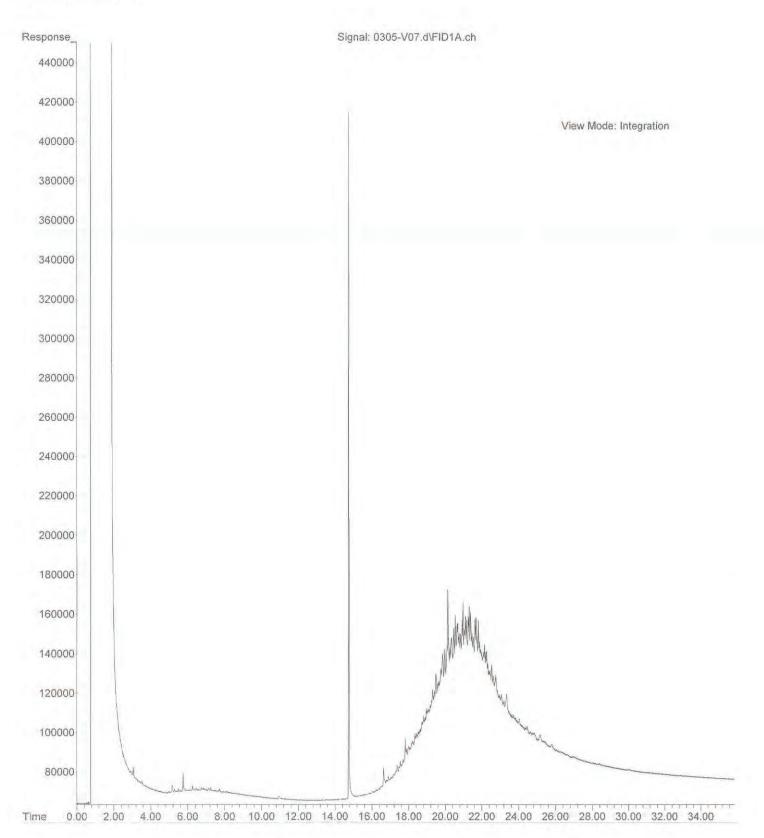


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Operator : JT
Acquired : 3 Mar 2018 9:11 using AcqMethod V171020F.M
Instrument : Vigo

Sample Name: 03-007-10



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Acquired : 5 Mar 2018 13:32 using AcqMethod
Instrument : Vigo
Sample Name: 03-007-17 20X using AcqMethod V171020F.M

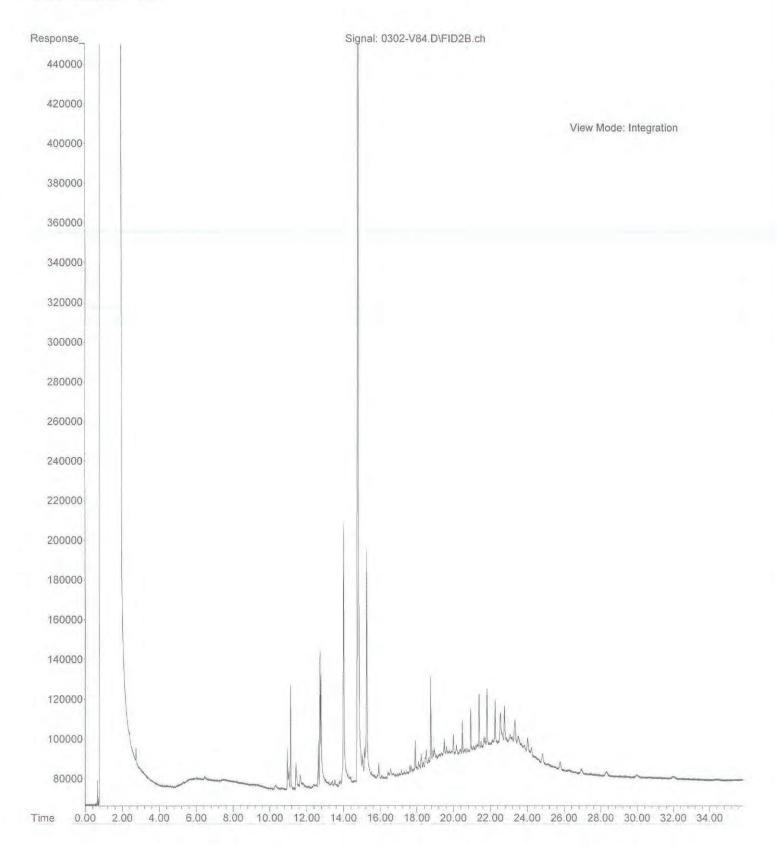


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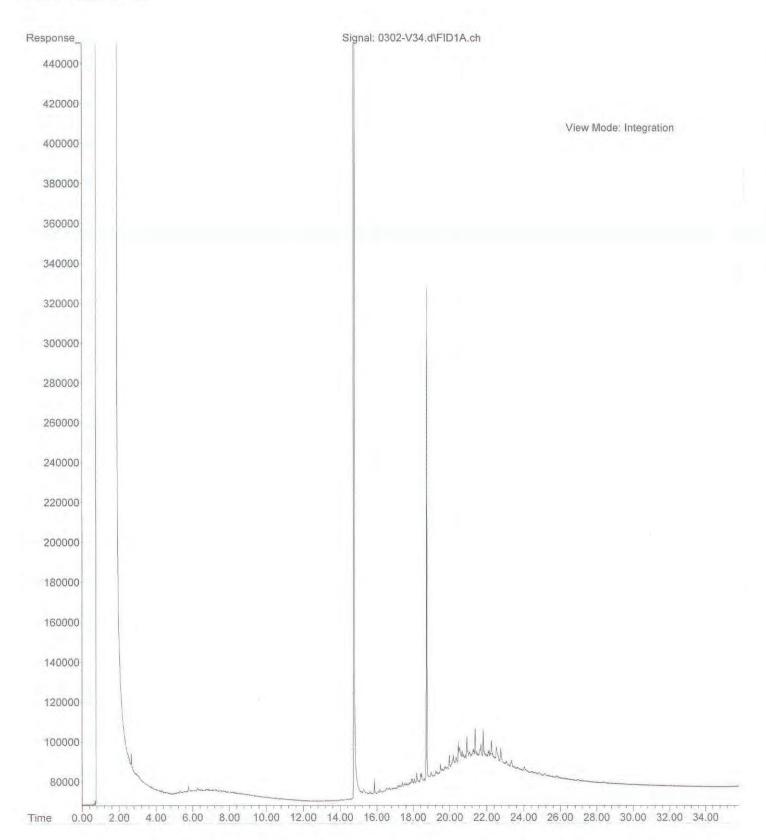
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Instrument: Vigo Sample Name: 03-007-21

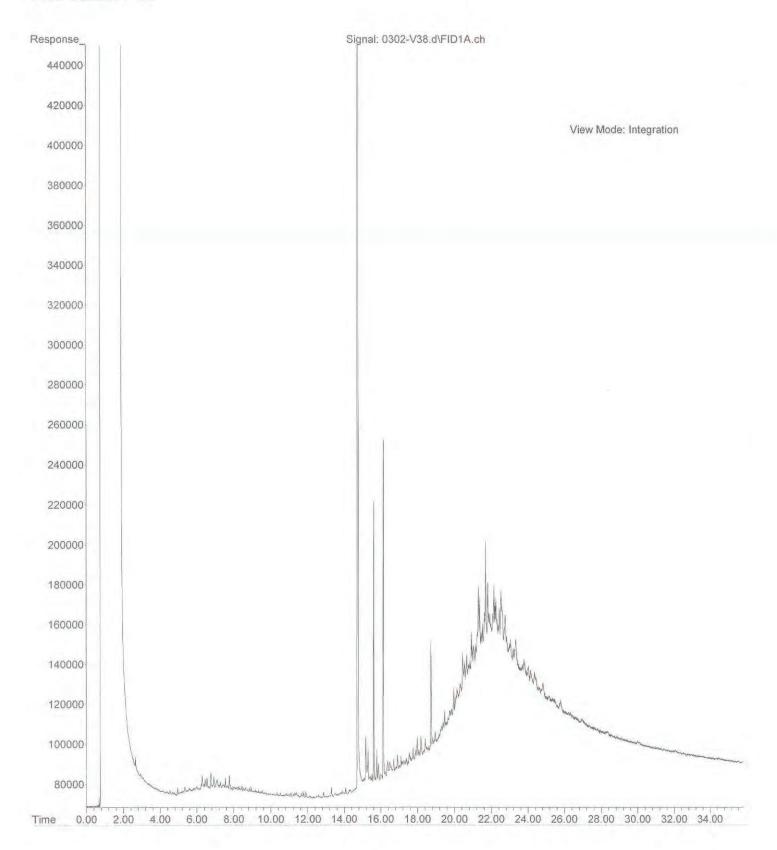


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Instrument : Vigo
Sample Name: 03-007-25



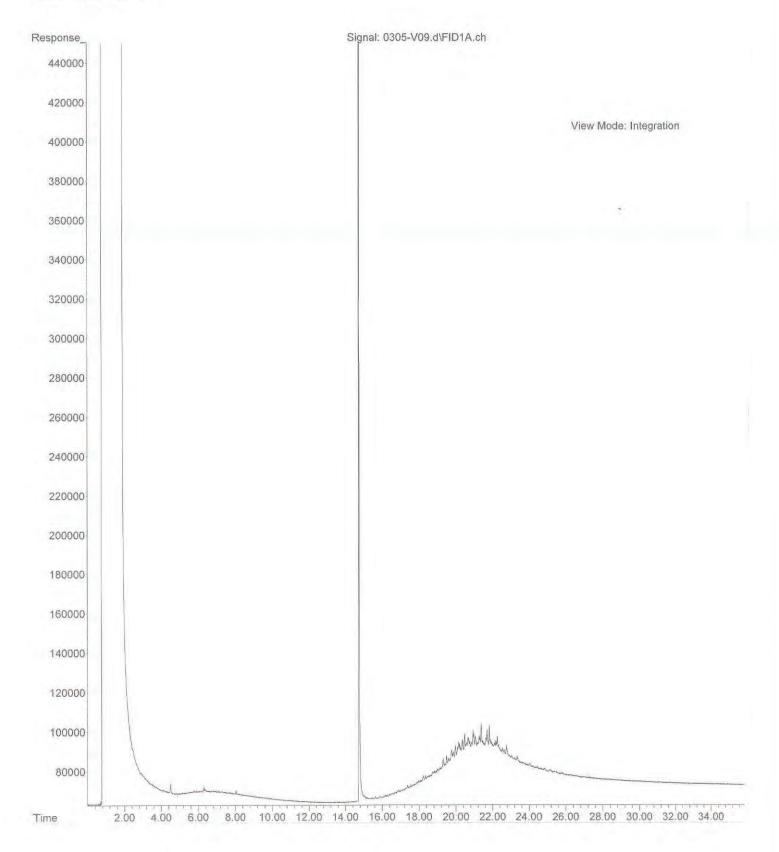
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Instrument : Vigo

Sample Name: 03-007-26

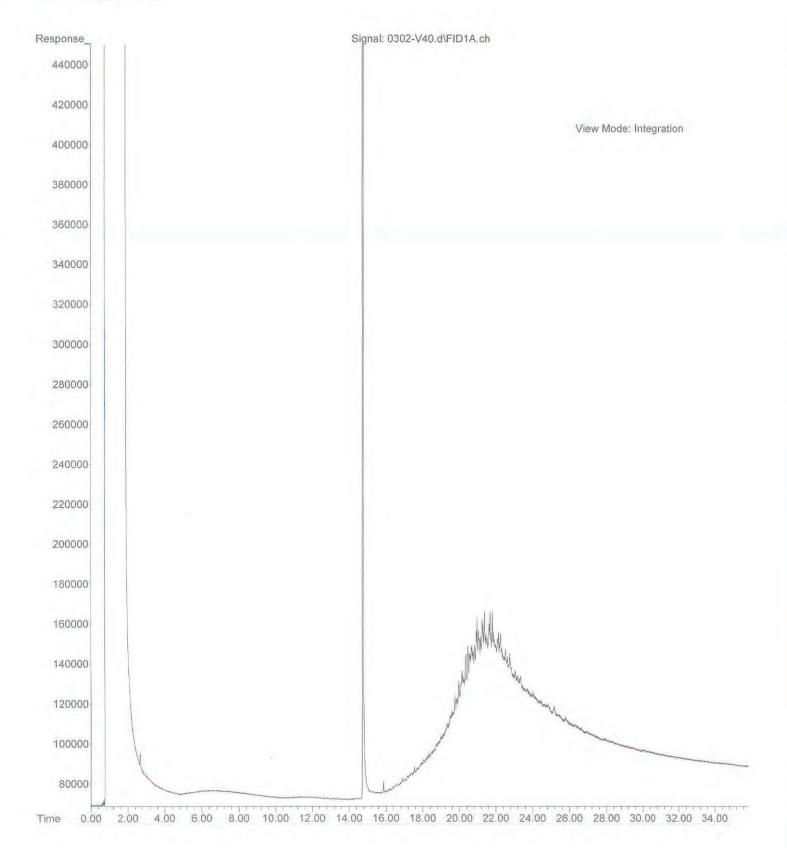


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Acquired : 5 Mar 2018 14:52 using AcqMethod
Instrument : Vigo
Sample Name: 03-007-27 10X

using AcqMethod V171020F.M

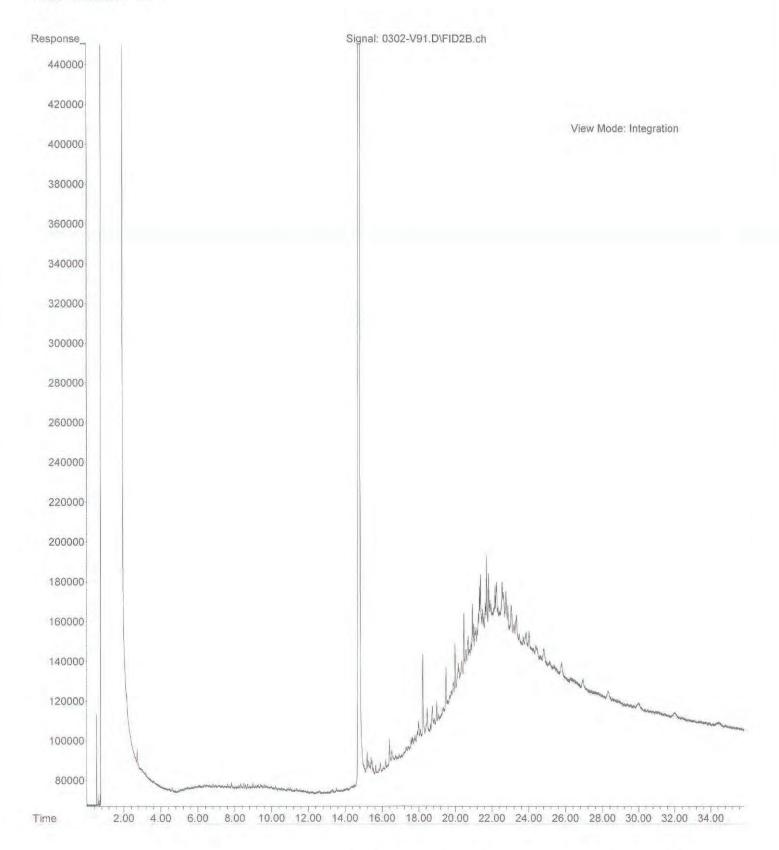


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Sample Name: 03-007-29 using AcqMethod V171020F.M



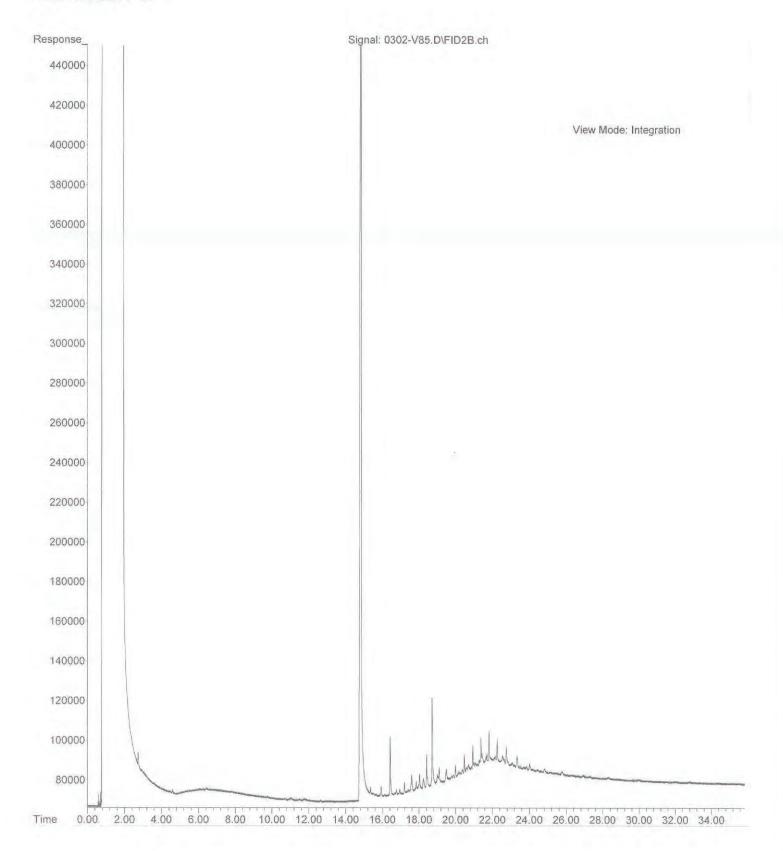
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Operator : JT
Acquired : 3 Mar 2018 10:30 using AcqMethod V171020F.M
Instrument : Vigo

Sample Name: 03-007-30



File :X:\DIESELS\VIGO\DATA\V180302.SEC\0302-V85.D
Operator : JT
Acquired : 3 Mar 2018 6:33 using AcqMethod V171020F.M
Instrument : Vigo

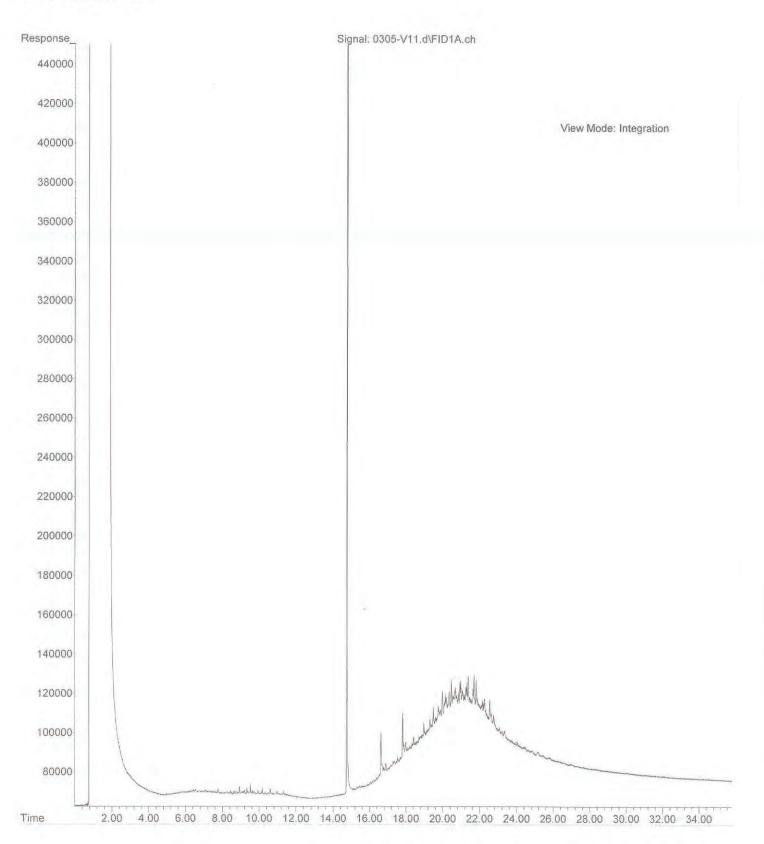
Sample Name: 03-007-31



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Acquired : 5 Mar 2018 16:11 using AcqMethod V171020F.M Instrument : Vigo

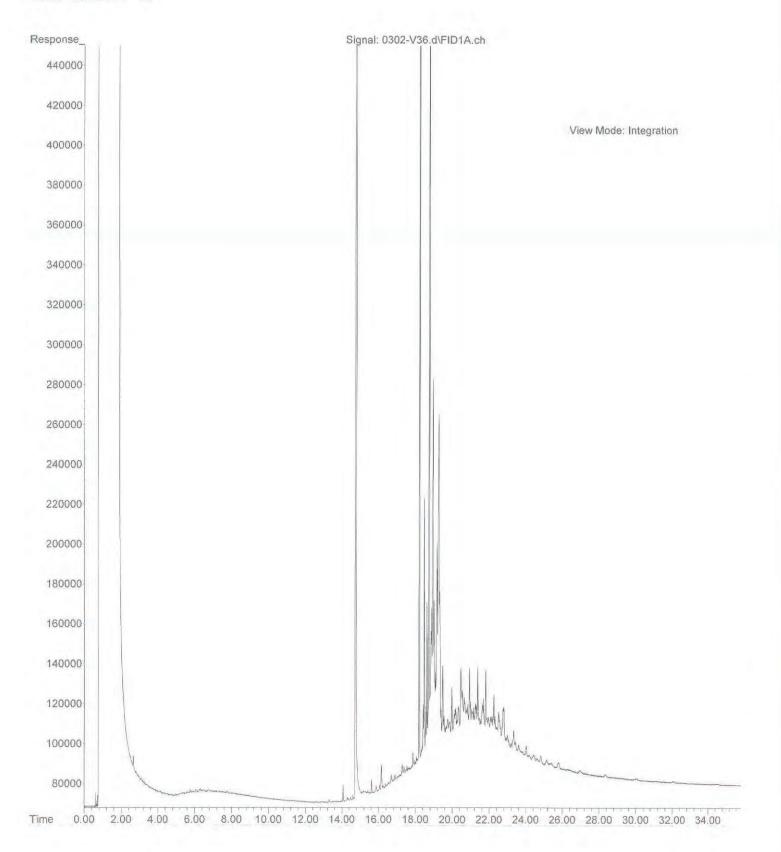
Sample Name: 03-007-41 10X



:X:\DIESELS\VIGO\DATA\V180302\0302-V36.d

7:12 using AcqMethod V171020F.M

Operator : JT
Acquired : 3 Mar 2018
Instrument : Vigo
Sample Name: 03-007-59

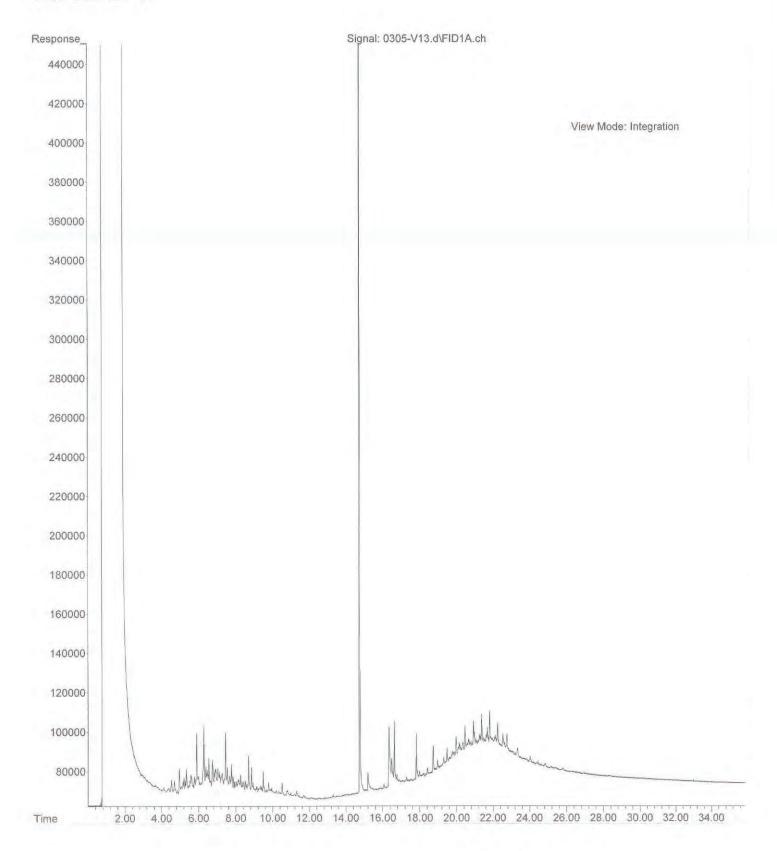


File :X:\DIESELS\VIGO\DATA\V180305\0305-V13.d Operator :JT

Acquired : 5 Mar 2018 17:31 using AcqMethod V171020F.M

Instrument : Vigo

Sample Name: 03-007-63 10X





14648 NE 95th Street, Redmond, WA 98052 • (425) 883-3881

March 26, 2018

Bruce Carpenter Herrera Environmental Consultants, Inc. 2200 6th Avenue, Suite 1100 Seattle, WA 98121

Re: Analytical Data for Project 17-06520-000

Laboratory Reference No. 1803-007B

Dear Bruce:

Enclosed are the analytical results and associated quality control data for samples submitted on March 1, 2018.

The standard policy of OnSite Environmental, Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

David Baumeister Project Manager

Enclosures



Project: 17-06520-000

Case Narrative

Samples were collected on February 28 and March 1, 2018 and received by the laboratory on March 1, 2018. They were maintained at the laboratory at a temperature of 2°C to 6°C.

Please note that any and all soil sample results are reported on a dry-weight basis, unless otherwise noted below.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

TCLP Lead EPA 1311/6010D Analysis

Due to a limited amount of sample, less than the required 100g was tumbled for TCLP analysis for samples PP31-6, PP21-10 and PP17-1 (03-007-26, 03-007-27, and 03-007-37) The amount of sample used was: 50g.

Any other QA/QC issues associated with this extraction and analysis will be indicated with a footnote reference and discussed in detail on the Data Qualifier page.

Project: 17-06520-000

TCLP LEAD EPA 1311/6010D

Offits. Hig/L (ppiff)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP23-2					
Laboratory ID:	03-007-21					
Lead	0.27	0.20	EPA 6010D	3-23-18	3-23-18	
Client ID:	PP23-15					
Laboratory ID:	03-007-24					
Lead	7.4	0.20	EPA 6010D	3-23-18	3-23-18	
Client ID:	PP21-2					
Laboratory ID:	03-007-25					
Lead	1.1	0.20	EPA 6010D	3-23-18	3-23-18	
Client ID:	PP21-6					
Laboratory ID:	03-007-26					
Lead	0.67	0.20	EPA 6010D	3-23-18	3-23-18	
Client ID:	PP21-10					
Laboratory ID:	03-007-27					
Lead	0.72	0.20	EPA 6010D	3-23-18	3-23-18	
Client ID:	PP20-5					
Laboratory ID:	03-007-30					
Lead	8.4	0.20	EPA 6010D	3-23-18	3-23-18	
Client ID:	PP20-10					
Laboratory ID:	03-007-31					
Lead	0.35	0.20	EPA 6010D	3-23-18	3-23-18	
Client ID:	DD00.45					
Client ID: Laboratory ID:	PP20-15 03-007-32					
	0.44	0.20	EPA 6010D	3-23-18	3-23-18	
Lead	V. 44	0.20	EFA 00 10D	3-23-10	3-23-10	

Project: 17-06520-000

TCLP LEAD EPA 1311/6010D

- J. ((1)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP18-3					
Laboratory ID:	03-007-33					
Lead	ND	0.20	EPA 6010D	3-23-18	3-23-18	
Client ID:	PP18-5					
Laboratory ID:	03-007-34					
	ND	0.20	EPA 6010D	3-23-18	3-23-18	
Lead	ND	0.20	EPA 60 10D	3-23-10	3-23-10	
Client ID:	PP17-1					
Laboratory ID:	03-007-37					
Lead	0.79	0.20	EPA 6010D	3-23-18	3-23-18	
Client ID:	PP17-5					
Laboratory ID:	03-007-38					
Lead	3.5	0.20	EPA 6010D	3-23-18	3-23-18	
Client ID:	PP16-11					
Laboratory ID:	03-007-42					
Lead	0.31	0.20	EPA 6010D	3-23-18	3-23-18	
Client ID:	PP11-5					
Laboratory ID:	03-007-49	0.00	EDA 0040D	0.00.40	0.00.40	
Lead	3.0	0.20	EPA 6010D	3-23-18	3-23-18	
Client ID:	PP11-10					
Laboratory ID:	03-007-50					
Lead	2.9	0.20	EPA 6010D	3-23-18	3-23-18	
Client ID:	PP27-7					
Laboratory ID:	03-007-55					
Lead	0.92	0.20	EPA 6010D	3-23-18	3-23-18	

Project: 17-06520-000

TCLP LEAD EPA 1311/6010D

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	PP24-7					_
Laboratory ID:	03-007-63					
Lead	16	0.20	EPA 6010D	3-23-18	3-23-18	

Project: 17-06520-000

TCLP LEAD EPA 1311/6010D QUALITY CONTROL

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0323TM1					
Lead	ND	0.20	EPA 6010D	3-23-18	3-23-18	
Laboratory ID:	MB0323TM2					
Lead	ND	0.20	EPA 6010D	3-23-18	3-23-18	

					Source	Pe	rcent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Red	covery	Limits	RPD	Limit	Flags
DUPLICATE											
Laboratory ID:	03-00	07-24									
	ORIG	DUP									
Lead	7.41	7.60	NA	NA			NA	NA	2	20	
Laboratory ID:	03-00	07-38									
	ORIG	DUP									
Lead	3.54	3.60	NA	NA			NA	NA	2	20	
MATRIX SPIKES											
Laboratory ID:	03-00	07-24									
	MS	MSD	MS	MSD		MS	MSD				
Lead	16.4	16.7	10.0	10.0	7.41	90	93	75-125	2	20	
Laboratory ID:	03-00	07-38									
	MS	MSD	MS	MSD		MS	MSD				
Lead	13.0	13.0	10.0	10.0	3.54	94	95	75-125	1	20	



Data Qualifiers and Abbreviations

- A Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
- B The analyte indicated was also found in the blank sample.
- C The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
- E The value reported exceeds the quantitation range and is an estimate.
- F Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
- H The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
- I Compound recovery is outside of the control limits.
- J The value reported was below the practical quantitation limit. The value is an estimate.
- K Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
- L The RPD is outside of the control limits.
- M Hydrocarbons in the gasoline range are impacting the diesel range result.
- M1 Hydrocarbons in the gasoline range (toluene-naphthalene) are present in the sample.
- N Hydrocarbons in the lube oil range are impacting the diesel range result.
- N1 Hydrocarbons in diesel range are impacting lube oil range results.
- O Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
- P The RPD of the detected concentrations between the two columns is greater than 40.
- Q Surrogate recovery is outside of the control limits.
- S Surrogate recovery data is not available due to the necessary dilution of the sample.
- T The sample chromatogram is not similar to a typical
- U The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- U1 The practical quantitation limit is elevated due to interferences present in the sample.
- V Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
- W Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
- X Sample extract treated with a mercury cleanup procedure.
- X1- Sample extract treated with a sulfuric acid/silica gel cleanup procedure.
- Y The calibration verification for this analyte exceeded the 20% drift specified in method 8260C, and therefore the reported result should be considered an estimate. The overall performance of the calibration verification standard met the acceptance criteria of the method.

7 -

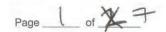
ND - Not Detected at PQL

PQL - Practical Quantitation Limit

RPD - Relative Percent Difference







	Analytical Laboratory Testing Services 14648 NE 95th Street • Redmond, WA 98052	Tur (i	naround Req n working day	uest /s)		L	abo	rate	ory	Nun	nbe	r:	0	3	-() (7									
Project N	HOVERA Jumber: 2-06520-000 Jame: Paulic Poule Janager: SWCE Campenter	Same 2 Date Sampled	_	1 Day 3 Days ys)	Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX		NWTPH-Dx (☐ Acid / SG Clean-up)	Volatiles 8260C	Halogenated Volatiles 8260C	EDB EPA 8011 (Waters Only)	Semivolatiles 8270D/SIM (with low-level PAHs)	PHTS 8270D/SIM (low-level)	CBS 8082A	Organochlorine Pesticides 8081B	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	TCLP Metals	HEM (oil and grease) 1664A				% Moisture
	PP28-3	3.1.18		Soil	2																					01
2	PP28-8	1	1035	1	2	X							,	X						X						X
3	PP28-10		1040		Z	X							1	X						X						X
	PPZ8-15		1045		2																					
-	PP29-3		1140		2	X			8				,	X	8					X						X
0	PP29-5		1145		2	X								X						X						X
-	PP29-10		1150		2	X								X						X						X
8 1	PP29-15		1155		2																					
9 7	PP32-4		9120		2	X							,	X						X						X
10 9	PP32-7		9130	V	2	X			\otimes				1	X	8					X						X
	Signature		ompany				Date			Time					s/Spe											
Relinque Receive Relinque Receive	ed wished ed) +	terre	86 86			3/	1/4	8		40		(8	A	lde rd	d dec	3/	3/10	18	. 5	多				
Receive	ed											1	Data	Pack	age:	Sta	ndar	d \square	Lev	/el III		Leve	IV [
Review	ed/Date		Reviewed/Dat	e									Chror	nato	grams	s wit	h fina	ıl rep	ort [Ele	ctroni	c Data	a Deliv	erables ((EDDs) [

OnSite Environmental Inc.

Chain of Custody

Page 2 of 7

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riojec	t Number: 17 - 0 VS t Name: Paulli t Manager: BWL	883-3881 · www.onsite-env.com The second of	Sar 2 D Sta		1 Day	Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX	NWTPH-Gx	NWTPH-Dx (☐ Acid / SG Clean-up)	Volatiles 8260C	Halogenated Volatiles 8260C	EDB EPA 8011 (Waters Only)		CHAIS 8270D/SIM (low-level) CPAHS	3s 8082A	Organochlorine Pesticides 8081B	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	TCLP Metals	M (oil and grease) 1664A				% Moisture
Lab ID		imple Identification	Sample	1 Sampled	Matrix		WN	1	N N	NN NN	Vols	Hal	ED	1		1		Org	- CH	Tota	1	TCI.	HEM		+	+	1
11	PP32-11		3.1.1		Soil	2		X	_	-				-	XX	X)	_		(X	2			-		X
12	PP32-15	>		940		Z	L									_	\dashv					1				-	
13	PP31-3			1000		2	X								*						X	A				_	X
14	PP31-5			1005		2																					
15	PP31-11			1010		2	X								X						X						X
16	PP31-15		1	1015		2																2					
17	-			IIOS		2	X	0		(8)					X	X					X	1					X
18				1110		2	X								X						X	X					X
19	PP33 -			115		2	X								X						X	X					X
20	PP33-1	15	1	1120	1	2																1					
	11 30	Signature		Company				Date			Time	е		Com	ment	s/Spe	ecial	Instr	uction	18							
Relin	quished	mianuhus	8	Herre	ia			13	31	815	1	34	0														
Rece	eived			~	M	>		3/	1/1	8	1	34	w														
Relin	nquished				0																						
Rec	eived																										
Reli	nquished							1																			
Rec	eived													Data	Pack	kage	Sta	ndar	rd 🗆	Le	vel III		Leve	el IV 🗆]		
Revi	ewed/Date			Reviewed/Da	ite									Chro	mato	gram	s wit	h fin	al rep	ort [Ele	ectror	ic Dat	a Deliv	rerables	s (EDDs)	



Page 3 of 3

Analytical Laboratory Testing Services 14648 NE 95th Street • Redmond, WA 98052		naround Req working day			La	abo	rato	ry N	umk	er:	(3	-(0 0	7									
Phone: (425) 883-3881 • www.onsite-env.com Company: Here Project Number: Project Name: Project Manager: Project Manager: Sampled by: Sampled by: Sampled by: Sampled by: Sampled by: Sampled Manager: Sampled Bland Sampled Bland	Date		1 Day	Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX	NWTPH-Gx	Volatiles 8260C	Halogenated Volatiles 8260C	EDB EPA 8011 (Waters Only)	Semivolatiles 8270D/SIM (with low-level PAHs)	8270D/SIM (low-level)	CBs 8082A	Organochlorine Pesticides 8081B	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	TCLP Metals Load	HEM (oil and grease) 1664A				% Moisture
Lab ID Sample Identification W PP23-2	2.28.18		Soil	2	X(N		2	1	Ш	0) 3		<u>X</u>		O	O		X	0	_	1			X
22 PP23-S	1	1405	1	2	X							X	_					X						X
23 PP23-10		1415		Z	X							X						X						X
24 PP23-15		1420		2	X							X						X	0				1	X
25 PP21-Z		1330		Z	χ		((X)					(\$)					X	0					X
20 PPZI-L		1335		2	X		(8				X	0					X	0					X
27 pp21-10		1345		2	X	8		8				X	X					X	0					X
28 PP21-K		1350		2				(X			((X)	(3)				(\otimes					(X
29 PP20-2		1250		2	X			X)				X	(3)					X						X
30 PP20-5	1	1255	V	2	X		(3)				X	$\langle y \rangle$					X	0					X
Relinquished Received Relinquished		Herve	va DE			Date 3			134 34	-	Cor	nmen	ts/Spi	ecial	Instru	uction	ns							
Received																								
Relinquished																								
Received											Data	a Pac	kage	Sta	ndar	d 🗆	Lev	vel III		Leve	IV 🗆			
Reviewed/Date		Reviewed/Dat	te								Chro	omato	gram	s wit	h fina	al rep	ort [Elec	ctroni	c Data	Delive	erables (E	EDDs) [



Page 4 of 7

Analytical Laboratory Testing Services 14648 NE 95th Street • Redmond, WA 98052		naround Req n working da			La	abo	orat	ory	Nu	mb	er:	03	-	nı	77							
Phone: (425) 883-3881 • www.onsite-env.com Company: Project Number: The project Name: Project Manager: Project Manager: Brough Campental Sampled by: Brough Brown Brow	Same 2 Day Stance (TPH		1 Day	Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX	NWTPH-Gx	NWTPH-Dx (☐ Acid / SG Clean-up)	Volatiles 8260C	Halogenated Volatiles 8260C	EDB EPA 8011 (Waters Only)	Semivolatiles 8270D/SIM (with low-level PAHs)	2/1/2	Organochlorine Pesticides 8081B	MIS/QI	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	TCLP-Wetels Lead HEM (oil and grease) 1664A			% Moisture
31 PP20-10	Sampled	Sampled	Sol	N	X	Ž	2	X	9	H	日	S & &	X	ō	ō	ਠ	2	X	0			% X
31 PP20-10 32 PP20-15	2,28,18	1305	001	2	^			X				Q	1					X	0			X
33 PP18-3		1150		2	x							X	-					X	0			X
34 PP18-S		1200		Z	X							X						X	0			X
35 PP18 - 10		1205		Z	X							X	1					X				X
36 PP18-15		1210		Z																		
37 PP17-1		1100		2	X							X						X	0			X
38 PP17-5		1110		2	X							X						X	0			X
39 PP17-10		1120		2	X							X						X				X
40 PP17-18		1128	V	2																		
Relinquished Received Relinquished		Herra COS	eva NE)		Date 3	1.1	78 ls	A	34	-	Comme	nts/Sp	ecial	Instr	uction	ns					
Received																						
Relinquished Received												200	4.0						П .			
Reviewed/Date		Reviewed/Da	te							_										/el IV □		D-1/17
(1021)110(1011)		Control of the Control	-									Griroma	logran	iis Wi	in in	ai rep	Ort L	FIE	ctronic Da	ata Deliver	ables (EDI	JS) [_]

OnSite Environmental Inc.

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Page 5 of 7

Analytical Laboratory Testing Services 14648 NE 95th Street • Redmond,	NA 98052 (ir	naround Requ working day			Li	abo	rato	ry N	Num	ber	:	0	3 -	0	07	7								
Phone: (425) 883-3881 • www.onsiti Company: HOVEYON Project Number: Project Name: POUL & Powe Project Manager: BYCL Carperter Sampled by:	Same 2 Day Stand (TPH		1 Day 3 Days ys)	Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX		NVIPH-DX (Acid / Sci Clean-up)	Volatilies ocoolo Halogenated Volatiles 8260C	EDB EPA 8011 (Waters Only)	Semivolatiles 8270D/SIM		PCBs 8082A	Organochlorine Pesticides 8081B	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	TCLP Metale—Lead	HEM (oil and grease) 1664A				% Moisture
Lab ID Sample Identification 41 PPI LO-1	2:28:12		Goil	2	X	X		(X	1		03.	>	(X	1				X	-	_				X
42 PP16-11	T	1045		Z	X							>	<					X	0					X
43 PP 10-15		1050		Z														13						
44 PP14-1		950		Z								1												
45 PP14-S		955		2	X							X						X						X
44 PP14-12		1000		2	X							X						X						χ
47 PP14-14		1005		2																				
48 PP11-2		915		2																				
49 PP11 - 5		920		2	X							X						X	0					X
50 PP11-10		925	V	2	X							X						X	0					X
Relinquished Signature	11	HW4	2111	,		Date			Time	.1.	C	omm	ents/S	pecial	Instr	ructio	ns		1000					
Relinquished May	wyman	101 202	N'AC	_		2/	110		134	40	-													
Relinquished		you				21	110	1) (/0	9													
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Page 4 of 7

Analytical Laboratory Testing Services 14648 NE 95th Street • Redmond, WA 98052		naround Req working da			L	abo	rato	ry l	Num	ber:		03	} -	0 (07	1								
Phone: (425) 883-3881 • www.onsite-env.com Company: Heyrera Project Number: 17-04520-000 Project Name: Project Manager: Byce Carpentul Sampled by Byanna Bland	Same 2 Day Stand (TPH	diard (7 Days) analysis 5 Da (other)	- I	Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX		NWTPH-Dx (Acid / SG Clean-up)	Volatilies ozouc. Halogenated Volatiles 8260C	EDB EPA 8011 (Waters Only)	Semivolatiles 8270D/SIM	PAHs 8270D/SIM (low-level)	CBs 8082A	Organochlorine Pesticides 8081B	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	TCLP Metals Lead	HEM (oil and grease) 1664A				% Moisture
Si PPII-IS	Sampled 2.28.12	93S	Matrix	2	Z	Z	Z	2	> I	Ш	00 3	2 OE (٦٩	0	0	0	12	F	E	I			\forall	%
52 PP10-11	1000	840	1	2	X				1	T	T	X						X					\Box	X
53 PP10-17		8400		2	X							X						X						X
54 PP27 -4		1545		2																				
SS PP 27-7		1550		2	X							X						X	0					X
Sie PPZ9-10		1600		2	X							X						X						X
S7 PP27-15		1610		2																				
58 PP30-3		1505		2																				
59 PP30-S		1510		2	X			X)				X	X					X						X
60 PP30-10	V	1520	V	2	X							X						X						X
Received Relinquished Relinquished		HWV4	SE)		3/3/		8	13 13	40	,	omme	ents/Sp	pecial	Instr	uctio	ns							
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Relinquished Received				_				-			Do	ta Da	ckac	o. Ct	anda	rd 🗆	Los	الا امر	n	Leve	I IV 🗆			
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Page T of 7

	14648 NE 9	poratory Testing Services 15th Street • Redmond, WA 98052	Turi (ir	naround Req n working da	uest ys)		L	abo	rato	ry	Nur	nbe	er:	C	13	_	n	17								
Project	Number: 7-0 Name: Paul	Si 883-3881 · www.onsite-env.com EVA EVSZO-OCO LIC Park Ce Carpertel INA Bland	(ТРН	/s [dard (7 Days) analysis 5 Da	1 Day 3 Days	Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX	NWTPH-Gx	NWTPH-Dx (☐ Acid / SG Clean-up)	Volatiles 8260C	Halogenated Volatiles 8260C	EDB EPA 8011 (Waters Only)	Semivolatiles 8270D/SIM (with low-level PAHs)	PATS 8270D/SIM (low-level)	s 8082A	Organochlorine Pesticides 8081B	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	TCLP Metals Lead	HEM (oil and grease) 1664A			% Moisture
Lab ID		imple Identification	Date Sampled	Time Sampled	Matrix		WN	NN NN	WN	NZ	Vola	Halo	EDB	Sem (with	村	PCB	Orga	Orga	Chlo	Tota	Tota	TCL	HEN		-	× ×
Col	PP30-K		2.28.18		801	2																				
62	PP24-2			1436		2		4								_										
63	PP 24-2	+		1435		2	X	8) (X					X	X					X	0				X
104	PPZ4-1	O	V	1445	V	2	X								X						X					X
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14648 NE 95th Street, Redmond, WA 98052 • (425) 883-3881

April 3, 2018

George Iftner Herrera Environmental Consultants, Inc. 2200 6th Avenue, Suite 1100 Seattle, WA 98121

Re: Analytical Data for Project 17-06520-000

Laboratory Reference No. 1803-248

Dear George:

Enclosed are the analytical results and associated quality control data for samples submitted on March 23, 2018.

Please note the data for the additionally requested analyses will follow in the final report

The standard policy of OnSite Environmental, Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely

David Baumeister Project Manager

Enclosures

Project: 17-06520-000

Case Narrative

Samples were collected on March 23, 2018 and received by the laboratory on March 23, 2018. They were maintained at the laboratory at a temperature of 2°C to 6°C.

Please note that any and all soil sample results are reported on a dry-weight basis, unless otherwise noted below.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

NWTPH-HCID

Matrix: Water Units: mg/L (ppm)

Onits. Hig/L (ppin)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW1					
Laboratory ID:	03-248-01					
Gasoline Range Organics	ND	0.10	NWTPH-HCID	3-27-18	3-27-18	
Diesel Range Organics	ND	0.26	NWTPH-HCID	3-27-18	3-27-18	
Lube Oil Range Organics	ND	0.41	NWTPH-HCID	3-27-18	3-27-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	81	50-150				
Client ID:	MW2					
Laboratory ID:	03-248-02					
Gasoline Range Organics	ND	0.11	NWTPH-HCID	3-27-18	3-27-18	
Diesel Range Organics	ND	0.11	NWTPH-HCID	3-27-18	3-27-18	
Lube Oil Range Organics	ND	0.42	NWTPH-HCID	3-27-18	3-27-18	
Surrogate:	Percent Recovery	Control Limits	IWWIIIIIOID	0 27 10	0 27 10	
o-Terphenyl	84	50-150				
0-Telphenyi	04	30-130				
Client ID:	MW3					
Laboratory ID:	03-248-03					
Gasoline Range Organics	ND	0.10	NWTPH-HCID	3-27-18	3-27-18	
Diesel Range Organics	ND	0.26	NWTPH-HCID	3-27-18	3-27-18	
Lube Oil Range Organics	ND	0.41	NWTPH-HCID	3-27-18	3-27-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	85	50-150				
Client ID:	MW4					
Laboratory ID:	03-248-04					
Gasoline Range Organics	ND	0.11	NWTPH-HCID	3-27-18	3-27-18	
Diesel Range Organics	ND	0.27	NWTPH-HCID	3-27-18	3-27-18	
Lube Oil Range Organics	ND	0.44	NWTPH-HCID	3-27-18	3-27-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	89	50-150				
Client ID:	MW5					
Laboratory ID:	03-248-05					
Gasoline Range Organics	ND	0.10	NWTPH-HCID	3-27-18	3-27-18	
Diesel Range Organics	ND	0.10	NWTPH-HCID	3-27-18	3-27-18	
Lube Oil Range Organics	ND ND	0.28	NWTPH-HCID	3-27-16 3-27-18	3-27-18 3-27-18	
Surrogate:	Percent Recovery	Control Limits	NVVIFILLIOID	3-21-10	J-21-10	
o-Terphenyl	84	50-150				
о-тегрпенуі	04	30-130				

NWTPH-HCID

Matrix: Water Units: mg/L (ppm)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW6					
Laboratory ID:	03-248-06					
Gasoline Range Organics	ND	0.11	NWTPH-HCID	3-27-18	3-27-18	
Diesel Range Organics	ND	0.28	NWTPH-HCID	3-27-18	3-27-18	
Lube Oil Range Organics	ND	0.45	NWTPH-HCID	3-27-18	3-27-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	84	50-150				
Client ID:	B-06					
Laboratory ID:	03-248-07					
Gasoline Range Organics	ND	0.10	NWTPH-HCID	3-27-18	3-27-18	
Diesel Range Organics	ND	0.10	NWTPH-HCID	3-27-18	3-27-18	
Lube Oil Range Organics	ND	0.41	NWTPH-HCID	3-27-18	3-27-18	
Surrogate:	Percent Recovery	Control Limits	NW II II II OID	0 27 10	0 27 10	
o-Terphenyl	123	50-150				
o respirent	720	00 700				
Client ID:	B-09					
Laboratory ID:	03-248-08					
Gasoline Range Organics	ND	0.10	NWTPH-HCID	3-27-18	3-27-18	
Diesel Range Organics	ND	0.26	NWTPH-HCID	3-27-18	3-27-18	
Lube Oil Range Organics	ND	0.42	NWTPH-HCID	3-27-18	3-27-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	84	50-150				
Client ID.	D 44					
Client ID:	B-11					
Laboratory ID:	03-248-09	0.40	NA/TRULLIC:S	0.07.10	0.07.10	
Gasoline Range Organics	ND	0.10	NWTPH-HCID	3-27-18	3-27-18	
Diesel Range Organics	ND	0.26	NWTPH-HCID	3-27-18	3-27-18	
Lube Oil Range Organics	ND	0.42	NWTPH-HCID	3-27-18	3-27-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	92	50-150				

NWTPH-HCID QUALITY CONTROL

Matrix: Water Units: mg/L (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0327W1					
Gasoline Range Organics	ND	0.040	NWTPH-HCID	3-27-18	3-27-18	
Diesel Range Organics	ND	0.10	NWTPH-HCID	3-27-18	3-27-18	
Lube Oil Range Organics	ND	0.16	NWTPH-HCID	3-27-18	3-27-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	80	50-150				

TOTAL METALS EPA 200.8/7470A

Matrix: Water Units: ug/L (ppb)

3 (11)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW1					
Laboratory ID:	03-248-01					
Arsenic	ND	3.3	EPA 200.8	3-27-18	3-27-18	
Cadmium	ND	4.4	EPA 200.8	3-27-18	3-27-18	
Chromium	ND	11	EPA 200.8	3-27-18	3-27-18	
Lead	ND	1.1	EPA 200.8	3-27-18	3-27-18	
Mercury	ND	0.50	EPA 7470A	3-30-18	3-30-18	
Client ID:	MW2					
Laboratory ID:	03-248-02					
Arsenic	3.4	3.3	EPA 200.8	3-27-18	3-27-18	
Cadmium	ND	4.4	EPA 200.8	3-27-18	3-27-18	
Chromium	ND	11	EPA 200.8	3-27-18	3-27-18	
Lead	ND	1.1	EPA 200.8	3-27-18	3-27-18	
Mercury	ND	0.50	EPA 7470A	3-30-18	3-30-18	
Client ID:	MW3					
Laboratory ID:	03-248-03					
Arsenic	ND	3.3	EPA 200.8	3-27-18	3-27-18	
Cadmium	ND	4.4	EPA 200.8	3-27-18	3-27-18	
Chromium	ND	11	EPA 200.8	3-27-18	3-27-18	
Lead	ND	1.1	EPA 200.8	3-27-18	3-27-18	
Mercury	ND	0.50	EPA 7470A	3-30-18	3-30-18	
Client ID:	MW4					
Laboratory ID:	03-248-04					
Arsenic	6.0	3.3	EPA 200.8	3-27-18	3-27-18	
Cadmium	ND	4.4	EPA 200.8	3-27-18	3-27-18	
Chromium	ND	11	EPA 200.8	3-27-18	3-27-18	
Lead	ND	1.1	EPA 200.8	3-27-18	3-27-18	
Mercury	ND	0.50	EPA 7470A	3-30-18	3-30-18	

TOTAL METALS EPA 200.8/7470A

Matrix: Water Units: ug/L (ppb)

J (11)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW5					
Laboratory ID:	03-248-05					
Arsenic	ND	3.3	EPA 200.8	3-27-18	3-27-18	
Cadmium	ND	4.4	EPA 200.8	3-27-18	3-27-18	
Chromium	ND	11	EPA 200.8	3-27-18	3-27-18	
Lead	3.2	1.1	EPA 200.8	3-27-18	3-27-18	
Mercury	ND	0.50	EPA 7470A	3-30-18	3-30-18	
Client ID:	MW6					
Laboratory ID:	03-248-06					
Arsenic	ND	3.3	EPA 200.8	3-27-18	3-27-18	
Cadmium	ND ND	3.3 4.4	EPA 200.8 EPA 200.8	3-27-16 3-27-18	3-27-16 3-27-18	
Chromium	ND ND	4.4 11	EPA 200.8	3-27-16 3-27-18	3-27-18 3-27-18	
Lead	ND	1.1	EPA 200.8	3-27-18	3-27-18	
Mercury	ND	0.50	EPA 7470A	3-30-18	3-30-18	
Moroary		0.00	LITTITOT	0 00 10	0 00 10	
011 10	D 00					
Client ID:	B-06					
Laboratory ID:	03-248-07					
Arsenic	ND	3.3	EPA 200.8	3-27-18	3-27-18	
Cadmium	ND	4.4	EPA 200.8	3-27-18	3-27-18	
Chromium	ND	11	EPA 200.8	3-27-18	3-27-18	
Lead	1.9	1.1	EPA 200.8	3-27-18	3-27-18	
Mercury	ND	0.50	EPA 7470A	3-30-18	3-30-18	
Client ID:	B-09					
Laboratory ID:	03-248-08					
Arsenic	3.7	3.3	EPA 200.8	3-27-18	3-27-18	
Cadmium	ND	4.4	EPA 200.8	3-27-18	3-27-18	
Chromium	ND	11	EPA 200.8	3-27-18	3-27-18	
Lead	ND	1.1	EPA 200.8	3-27-18	3-27-18	
Mercury	ND	0.50	EPA 7470A	3-30-18	3-30-18	

TOTAL METALS EPA 200.8/7470A

Matrix: Water Units: ug/L (ppb)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-11					
Laboratory ID:	03-248-09					
Arsenic	ND	3.3	EPA 200.8	3-27-18	3-27-18	
Cadmium	ND	4.4	EPA 200.8	3-27-18	3-27-18	
Chromium	ND	11	EPA 200.8	3-27-18	3-27-18	
Lead	ND	1.1	EPA 200.8	3-27-18	3-27-18	
Mercury	ND	0.50	EPA 7470A	3-30-18	3-30-18	

Project: 17-06520-000

TOTAL METALS EPA 200.8/7470A QUALITY CONTROL

Matrix: Water
Units: ug/L (ppb)

			Date	Date	
Result	PQL	Method	Prepared	Analyzed	Flags
MB0327WM1					
ND	8.3	EPA 200.8	3-27-18	3-27-18	
ND	11	EPA 200.8	3-27-18	3-27-18	
ND	28	EPA 200.8	3-27-18	3-27-18	
ND	2.8	EPA 200.8	3-27-18	3-27-18	
MB0330W1					
ND	0.50	EPA 7470A	3-30-18	3-30-18	
	MB0327WM1 ND ND ND ND ND MD ND	MB0327WM1 ND 8.3 ND 11 ND 28 ND 2.8 MB0330W1	MB0327WM1 ND 8.3 EPA 200.8 ND 11 EPA 200.8 ND 28 EPA 200.8 ND 2.8 EPA 200.8 MB0330W1	Result PQL Method Prepared MB0327WM1 ND 8.3 EPA 200.8 3-27-18 ND 11 EPA 200.8 3-27-18 ND 28 EPA 200.8 3-27-18 ND 2.8 EPA 200.8 3-27-18 MD 2.8 EPA 200.8 3-27-18	Result PQL Method Prepared Analyzed MB0327WM1 ND 8.3 EPA 200.8 3-27-18 3-27-18 ND 11 EPA 200.8 3-27-18 3-27-18 ND 28 EPA 200.8 3-27-18 3-27-18 ND 2.8 EPA 200.8 3-27-18 3-27-18 MB0330W1 MB0330W1 MB0330W1 MB0330W1

					Source	Percent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Recovery	Limits	RPD	Limit	Flags
DUPLICATE										
Laboratory ID:	03-06	88-09								
	ORIG	DUP								
Arsenic	ND	ND	NA	NA		NA	NA	NA	20	
Cadmium	ND	ND	NA	NA		NA	NA	NA	20	
Chromium	ND	ND	NA	NA		NA	NA	NA	20	
Lead	ND	ND	NA	NA		NA	NA	NA	20	
Laboratory ID:	03-24	18-01								
Mercury	ND	ND	NA	NA		NA	NA	NA	20	

MATRIX SPIKES

Laboratory ID:	03-0	68-09									
	MS	MSD	MS	MSD		MS	MSD				
Arsenic	114	110	111	111	ND	102	99	75-125	3	20	
Cadmium	123	120	111	111	ND	111	108	75-125	3	20	
Chromium	109	108	111	111	ND	99	97	75-125	1	20	
Lead	125	123	111	111	ND	113	111	75-125	11	20	
Laboratory ID:	03-2	48-01									
Mercury	12.2	11.4	12.5	12.5	ND	97	91	75-125	7	20	

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW1					
Laboratory ID:	03-248-01					
Benzo[a]anthracene	ND	0.0094	EPA 8270D/SIM	3-26-18	3-26-18	
Chrysene	ND	0.0094	EPA 8270D/SIM	3-26-18	3-26-18	
Benzo[b]fluoranthene	ND	0.0094	EPA 8270D/SIM	3-26-18	3-26-18	
Benzo(j,k)fluoranthene	ND	0.0094	EPA 8270D/SIM	3-26-18	3-26-18	
Benzo[a]pyrene	ND	0.0094	EPA 8270D/SIM	3-26-18	3-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0094	EPA 8270D/SIM	3-26-18	3-26-18	
Dibenz[a,h]anthracene	ND	0.0094	EPA 8270D/SIM	3-26-18	3-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	34	25 - 107				
Pyrene-d10	53	28 - 103				
Terphenyl-d14	52	36 - 129				

cPAHs EPA 8270D/SIM

			Date	Date	
Result	PQL	Method	Prepared	Analyzed	Flags
MW2					
03-248-02					
ND	0.011	EPA 8270D/SIM	3-26-18	3-26-18	
ND	0.011	EPA 8270D/SIM	3-26-18	3-26-18	
ND	0.011	EPA 8270D/SIM	3-26-18	3-26-18	
ND	0.011	EPA 8270D/SIM	3-26-18	3-26-18	
ND	0.011	EPA 8270D/SIM	3-26-18	3-26-18	
ND	0.011	EPA 8270D/SIM	3-26-18	3-26-18	
ND	0.011	EPA 8270D/SIM	3-26-18	3-26-18	
Percent Recovery	Control Limits				
53	25 - 107				
68	28 - 103				
68	36 - 129				
	MW2 03-248-02 ND ND ND ND ND ND ND ND ND SD Percent Recovery 53 68	MW2 03-248-02 ND 0.011 ND 0.011 ND 0.011 ND 0.011 ND 0.011 ND 0.011 Percent Recovery Control Limits 53 25 - 107 68 28 - 103	MW2 03-248-02 0.011 EPA 8270D/SIM ND 0.011 EPA 8270D/SIM Percent Recovery Control Limits 53 25 - 107 68 28 - 103	MW2 03-248-02 ND 0.011 EPA 8270D/SIM 3-26-18 Percent Recovery Control Limits 53 25 - 107 68 28 - 103	Result PQL Method Prepared Analyzed MW2 03-248-02 ND 0.011 EPA 8270D/SIM 3-26-18 3-26-18 Percent Recovery Control Limits 53 25 - 107 68 28 - 103 40 40 40

cPAHs EPA 8270D/SIM

J				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW3					
Laboratory ID:	03-248-03					
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Chrysene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	62	25 - 107				
Pyrene-d10	84	28 - 103				
Terphenyl-d14	81	36 - 129				

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW4					
Laboratory ID:	03-248-04					
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Chrysene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	48	25 - 107				
Pyrene-d10	71	28 - 103				
Terphenyl-d14	68	36 - 129				

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW5					
Laboratory ID:	03-248-05					
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Chrysene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	43	25 - 107				
Pyrene-d10	68	28 - 103				
Terphenyl-d14	64	36 - 129				

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW6					
Laboratory ID:	03-248-06					
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Chrysene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	58	25 - 107				
Pyrene-d10	77	28 - 103				
Terphenyl-d14	74	36 - 129				

cPAHs EPA 8270D/SIM

· ·				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-06					
Laboratory ID:	03-248-07					
Benzo[a]anthracene	ND	0.0094	EPA 8270D/SIM	3-26-18	3-26-18	
Chrysene	ND	0.0094	EPA 8270D/SIM	3-26-18	3-26-18	
Benzo[b]fluoranthene	ND	0.0094	EPA 8270D/SIM	3-26-18	3-26-18	
Benzo(j,k)fluoranthene	ND	0.0094	EPA 8270D/SIM	3-26-18	3-26-18	
Benzo[a]pyrene	ND	0.0094	EPA 8270D/SIM	3-26-18	3-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0094	EPA 8270D/SIM	3-26-18	3-26-18	
Dibenz[a,h]anthracene	ND	0.0094	EPA 8270D/SIM	3-26-18	3-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	51	25 - 107				
Pyrene-d10	70	28 - 103				
Terphenyl-d14	65	36 - 129				

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-09					
Laboratory ID:	03-248-08					
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Chrysene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	55	25 - 107				
Pyrene-d10	74	28 - 103				
Terphenyl-d14	70	36 - 129				

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-11					
Laboratory ID:	03-248-09					
Benzo[a]anthracene	ND	0.011	EPA 8270D/SIM	3-26-18	3-26-18	
Chrysene	ND	0.011	EPA 8270D/SIM	3-26-18	3-26-18	
Benzo[b]fluoranthene	ND	0.011	EPA 8270D/SIM	3-26-18	3-26-18	
Benzo(j,k)fluoranthene	ND	0.011	EPA 8270D/SIM	3-26-18	3-26-18	
Benzo[a]pyrene	ND	0.011	EPA 8270D/SIM	3-26-18	3-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.011	EPA 8270D/SIM	3-26-18	3-26-18	
Dibenz[a,h]anthracene	ND	0.011	EPA 8270D/SIM	3-26-18	3-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	44	25 - 107				
Pyrene-d10	62	28 - 103				
Terphenyl-d14	59	36 - 129				

cPAHs EPA 8270D/SIM **METHOD BLANK QUALITY CONTROL**

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0326W1					
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Chrysene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	3-26-18	3-26-18	
Surrogate:	Percent Recovery	Control Limits				_
2-Fluorobiphenyl	56	25 - 107				
Pyrene-d10	82	28 - 103				
Terphenyl-d14	79	36 - 129				

cPAHs EPA 8270D/SIM SB/SBD QUALITY CONTROL

					Pei	rcent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Rec	overy	Limits	RPD	Limit	Flags
SPIKE BLANKS										
Laboratory ID:	SB03	26W1								
	SB	SBD	SB	SBD	SB	SBD				
Benzo[a]anthracene	0.391	0.424	0.500	0.500	78	85	47 - 131	8	27	
Chrysene	0.415	0.450	0.500	0.500	83	90	48 - 120	8	29	
Benzo[b]fluoranthene	0.383	0.449	0.500	0.500	77	90	42 - 128	16	29	
Benzo(j,k)fluoranthene	0.420	0.429	0.500	0.500	84	86	46 - 121	2	27	
Benzo[a]pyrene	0.368	0.402	0.500	0.500	74	80	34 - 121	9	29	
Indeno(1,2,3-c,d)pyrene	0.407	0.445	0.500	0.500	81	89	39 - 128	9	28	
Dibenz[a,h]anthracene	0.410	0.448	0.500	0.500	82	90	39 - 125	9	30	
Surrogate:										
2-Fluorobiphenyl					56	61	25 - 107			
Pyrene-d10					80	85	28 - 103			
Terphenyl-d14					78	82	36 - 129			

Project: 17-06520-000

VOLATILES by EPA 8260C Page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW1					
Laboratory ID:	03-248-01					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chloromethane	ND	1.3	EPA 8260C	4-2-18	4-2-18	
Vinyl Chloride	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromomethane	ND	0.25	EPA 8260C	4-2-18	4-2-18	
Chloroethane	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Acetone	ND	5.0	EPA 8260C	4-2-18	4-2-18	
Iodomethane	ND	1.5	EPA 8260C	4-2-18	4-2-18	
Carbon Disulfide	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methylene Chloride	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Vinyl Acetate	ND	1.0	EPA 8260C	4-2-18	4-2-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Butanone	ND	5.0	EPA 8260C	4-2-18	4-2-18	
Bromochloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chloroform	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Benzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Trichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Dibromomethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromodichloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	4-2-18	4-2-18	
Toluene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	

Project: 17-06520-000

VOLATILES by EPA 8260C Page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW1					
Laboratory ID:	03-248-01					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Tetrachloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Hexanone	ND	2.0	EPA 8260C	4-2-18	4-2-18	
Dibromochloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Ethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
m,p-Xylene	ND	0.40	EPA 8260C	4-2-18	4-2-18	
o-Xylene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Styrene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromoform	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Isopropylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
n-Propylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
n-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	4-2-18	4-2-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Naphthalene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Surrogate:	Percent Recovery					

Surrogate: Percent Recovery Control Limit Dibromofluoromethane 96 75-127
Toluene-d8 96 80-127
4-Bromofluorobenzene 95 78-125

Project: 17-06520-000

VOLATILES by EPA 8260C Page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW2					
Laboratory ID:	03-248-02					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chloromethane	ND	1.3	EPA 8260C	4-2-18	4-2-18	
Vinyl Chloride	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromomethane	ND	0.25	EPA 8260C	4-2-18	4-2-18	
Chloroethane	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Acetone	ND	5.0	EPA 8260C	4-2-18	4-2-18	
Iodomethane	ND	1.5	EPA 8260C	4-2-18	4-2-18	
Carbon Disulfide	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methylene Chloride	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Vinyl Acetate	ND	1.0	EPA 8260C	4-2-18	4-2-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Butanone	ND	5.0	EPA 8260C	4-2-18	4-2-18	
Bromochloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chloroform	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Benzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Trichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Dibromomethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromodichloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	4-2-18	4-2-18	
Toluene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	

Project: 17-06520-000

VOLATILES by EPA 8260C Page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW2	FQL	Wethou	Frepareu	Allalyzeu	гіауз
Laboratory ID:	03-248-02					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Tetrachloroethene	ND	0.20	EPA 8260C	4-2-18 4-2-18	4-2-16 4-2-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	4-2-18 4-2-18	4-2-18 4-2-18	
2-Hexanone	ND	2.0	EPA 8260C EPA 8260C	4-2-16 4-2-18	4-2-16 4-2-18	
Dibromochloromethane	ND	0.20	EPA 8260C EPA 8260C	4-2-16 4-2-18	4-2-16 4-2-18	
	ND	0.20				
1,2-Dibromoethane			EPA 8260C	4-2-18	4-2-18	
Chlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Ethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
m,p-Xylene	ND	0.40	EPA 8260C	4-2-18	4-2-18	
o-Xylene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Styrene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromoform	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Isopropylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
n-Propylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
n-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	4-2-18	4-2-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Naphthalene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Surrogate:	Percent Recovery	Control Limits		-	-	

Surrogate: Percent Recovery Control Limits
Dibromofluoromethane 96 75-127
Toluene-d8 95 80-127
4-Bromofluorobenzene 94 78-125

Project: 17-06520-000

VOLATILES by EPA 8260C Page 1 of 2

omis. ug/L				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW3					
Laboratory ID:	03-248-03					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chloromethane	ND	1.3	EPA 8260C	4-2-18	4-2-18	
Vinyl Chloride	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromomethane	ND	0.25	EPA 8260C	4-2-18	4-2-18	
Chloroethane	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Acetone	ND	5.0	EPA 8260C	4-2-18	4-2-18	
Iodomethane	ND	1.5	EPA 8260C	4-2-18	4-2-18	
Carbon Disulfide	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methylene Chloride	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Vinyl Acetate	ND	1.0	EPA 8260C	4-2-18	4-2-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
(cis) 1,2-Dichloroethene	0.22	0.20	EPA 8260C	4-2-18	4-2-18	
2-Butanone	ND	5.0	EPA 8260C	4-2-18	4-2-18	
Bromochloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chloroform	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Benzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Trichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Dibromomethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromodichloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	4-2-18	4-2-18	
Toluene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	

Project: 17-06520-000

VOLATILES by EPA 8260C Page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW3					
Laboratory ID:	03-248-03					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Tetrachloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Hexanone	ND	2.0	EPA 8260C	4-2-18	4-2-18	
Dibromochloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Ethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
m,p-Xylene	ND	0.40	EPA 8260C	4-2-18	4-2-18	
o-Xylene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Styrene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromoform	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Isopropylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
n-Propylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
n-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	4-2-18	4-2-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Naphthalene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Surrogate:	Percent Recovery	Control Limits				

Surrogate: Percent Recovery Control Limits
Dibromofluoromethane 98 75-127
Toluene-d8 95 80-127
4-Bromofluorobenzene 96 78-125

Project: 17-06520-000

VOLATILES by EPA 8260C Page 1 of 2

omis. ug/L				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW4					
Laboratory ID:	03-248-04					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chloromethane	ND	1.3	EPA 8260C	4-2-18	4-2-18	
Vinyl Chloride	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromomethane	ND	0.25	EPA 8260C	4-2-18	4-2-18	
Chloroethane	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Acetone	ND	5.0	EPA 8260C	4-2-18	4-2-18	
Iodomethane	ND	1.5	EPA 8260C	4-2-18	4-2-18	
Carbon Disulfide	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methylene Chloride	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Vinyl Acetate	ND	1.0	EPA 8260C	4-2-18	4-2-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Butanone	ND	5.0	EPA 8260C	4-2-18	4-2-18	
Bromochloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chloroform	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Benzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Trichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Dibromomethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromodichloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	4-2-18	4-2-18	
Toluene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	

Project: 17-06520-000

VOLATILES by EPA 8260C Page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW4					
Laboratory ID:	03-248-04					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Tetrachloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Hexanone	ND	2.0	EPA 8260C	4-2-18	4-2-18	
Dibromochloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chlorobenzene	1.1	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Ethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
m,p-Xylene	ND	0.40	EPA 8260C	4-2-18	4-2-18	
o-Xylene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Styrene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromoform	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Isopropylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
n-Propylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
n-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	4-2-18	4-2-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Naphthalene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Surrogate:	Percent Recovery	Control Limits				

Surrogate: Percent Recovery Control Limit
Dibromofluoromethane 97 75-127
Toluene-d8 97 80-127
4-Bromofluorobenzene 95 78-125

Project: 17-06520-000

VOLATILES by EPA 8260C Page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW5					
Laboratory ID:	03-248-05					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chloromethane	ND	1.3	EPA 8260C	4-2-18	4-2-18	
Vinyl Chloride	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromomethane	ND	0.25	EPA 8260C	4-2-18	4-2-18	
Chloroethane	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Acetone	ND	5.0	EPA 8260C	4-2-18	4-2-18	
Iodomethane	ND	1.5	EPA 8260C	4-2-18	4-2-18	
Carbon Disulfide	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methylene Chloride	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Vinyl Acetate	ND	1.0	EPA 8260C	4-2-18	4-2-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Butanone	ND	5.0	EPA 8260C	4-2-18	4-2-18	
Bromochloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chloroform	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Benzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Trichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Dibromomethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromodichloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	4-2-18	4-2-18	
Toluene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	

Project: 17-06520-000

VOLATILES by EPA 8260C Page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW5	FQL	Wethou	Frepareu	Allalyzeu	гіауз
Laboratory ID:	03-248-05					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Tetrachloroethene	ND	0.20	EPA 8260C EPA 8260C	4-2-16 4-2-18	4-2-16 4-2-18	
	ND ND	0.20				
1,3-Dichloropropane		2.0	EPA 8260C	4-2-18	4-2-18	
2-Hexanone	ND ND	0.20	EPA 8260C	4-2-18	4-2-18	
Dibromochloromethane			EPA 8260C	4-2-18	4-2-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Ethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
m,p-Xylene	ND	0.40	EPA 8260C	4-2-18	4-2-18	
o-Xylene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Styrene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromoform	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Isopropylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
n-Propylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
n-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	4-2-18	4-2-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Naphthalene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Surrogate:	Percent Recovery	Control Limits				

Surrogate: Percent Recovery Control Limits
Dibromofluoromethane 99 75-127
Toluene-d8 97 80-127
4-Bromofluorobenzene 95 78-125

VOLATILES by EPA 8260C Page 1 of 2

Analyte	Result	PQL	Method	Date	Date	
				Prepared	Analyzed	Flags
Client ID:	MW6					
Laboratory ID:	03-248-06					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chloromethane	ND	1.3	EPA 8260C	4-2-18	4-2-18	
Vinyl Chloride	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromomethane	ND	0.25	EPA 8260C	4-2-18	4-2-18	
Chloroethane	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Acetone	ND	5.0	EPA 8260C	4-2-18	4-2-18	
lodomethane	ND	1.5	EPA 8260C	4-2-18	4-2-18	
Carbon Disulfide	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methylene Chloride	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Vinyl Acetate	ND	1.0	EPA 8260C	4-2-18	4-2-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Butanone	ND	5.0	EPA 8260C	4-2-18	4-2-18	
Bromochloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chloroform	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Benzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Trichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Dibromomethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromodichloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	4-2-18	4-2-18	
Toluene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	

VOLATILES by EPA 8260C Page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW6					
Laboratory ID:	03-248-06					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Tetrachloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Hexanone	ND	2.0	EPA 8260C	4-2-18	4-2-18	
Dibromochloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Ethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
m,p-Xylene	ND	0.40	EPA 8260C	4-2-18	4-2-18	
o-Xylene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Styrene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromoform	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Isopropylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
n-Propylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
n-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	4-2-18	4-2-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Naphthalene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Surrogate:	Percent Recovery	Control Limits				

Dibromofluoromethane 101 75-127 Toluene-d8 99 80-127 4-Bromofluorobenzene 96 78-125

Project: 17-06520-000

VOLATILES by EPA 8260C Page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-06					
Laboratory ID:	03-248-07					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chloromethane	ND	1.3	EPA 8260C	4-2-18	4-2-18	
Vinyl Chloride	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromomethane	ND	0.25	EPA 8260C	4-2-18	4-2-18	
Chloroethane	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Acetone	ND	5.0	EPA 8260C	4-2-18	4-2-18	
Iodomethane	ND	1.5	EPA 8260C	4-2-18	4-2-18	
Carbon Disulfide	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methylene Chloride	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Vinyl Acetate	ND	1.0	EPA 8260C	4-2-18	4-2-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Butanone	ND	5.0	EPA 8260C	4-2-18	4-2-18	
Bromochloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chloroform	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Benzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Trichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Dibromomethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromodichloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	4-2-18	4-2-18	
Toluene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	

Project: 17-06520-000

VOLATILES by EPA 8260C Page 2 of 2

Analyte	Result	PQL	Method	Date	Date	
				Prepared	Analyzed	Flags
Client ID:	B-06					
Laboratory ID:	03-248-07					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Tetrachloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Hexanone	ND	2.0	EPA 8260C	4-2-18	4-2-18	
Dibromochloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Ethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
m,p-Xylene	ND	0.40	EPA 8260C	4-2-18	4-2-18	
o-Xylene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Styrene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromoform	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Isopropylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
n-Propylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
n-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	4-2-18	4-2-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Naphthalene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Surrogate:	Percent Recovery	Control Limits				

Surrogate: Percent Recovery Control Limi
Dibromofluoromethane 101 75-127
Toluene-d8 101 80-127
4-Bromofluorobenzene 97 78-125

Project: 17-06520-000

VOLATILES by EPA 8260C Page 1 of 2

omis. ug/L				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-09					
Laboratory ID:	03-248-08					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chloromethane	ND	1.3	EPA 8260C	4-2-18	4-2-18	
Vinyl Chloride	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromomethane	ND	0.25	EPA 8260C	4-2-18	4-2-18	
Chloroethane	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Acetone	ND	5.0	EPA 8260C	4-2-18	4-2-18	
Iodomethane	ND	1.5	EPA 8260C	4-2-18	4-2-18	
Carbon Disulfide	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methylene Chloride	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Vinyl Acetate	ND	1.0	EPA 8260C	4-2-18	4-2-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Butanone	ND	5.0	EPA 8260C	4-2-18	4-2-18	
Bromochloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chloroform	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Benzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Trichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Dibromomethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromodichloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	4-2-18	4-2-18	
Toluene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	

Project: 17-06520-000

VOLATILES by EPA 8260C Page 2 of 2

Analyte		PQL		Date	Date	
	Result		Method	Prepared	Analyzed	Flags
Client ID:	B-09					
Laboratory ID:	03-248-08					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Tetrachloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Hexanone	ND	2.0	EPA 8260C	4-2-18	4-2-18	
Dibromochloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Ethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
m,p-Xylene	ND	0.40	EPA 8260C	4-2-18	4-2-18	
o-Xylene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Styrene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromoform	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Isopropylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
n-Propylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
n-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dibromo-3-chloropropane		1.0	EPA 8260C	4-2-18	4-2-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Naphthalene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Surrogate:	Percent Recovery	Control Limits		-	-	

Surrogate: Percent Recovery Control Limits
Dibromofluoromethane 101 75-127
Toluene-d8 97 80-127
4-Bromofluorobenzene 96 78-125

Project: 17-06520-000

VOLATILES by EPA 8260C Page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-11					
Laboratory ID:	03-248-09					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chloromethane	ND	1.3	EPA 8260C	4-2-18	4-2-18	
Vinyl Chloride	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromomethane	ND	0.25	EPA 8260C	4-2-18	4-2-18	
Chloroethane	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Acetone	ND	5.0	EPA 8260C	4-2-18	4-2-18	
Iodomethane	ND	1.5	EPA 8260C	4-2-18	4-2-18	
Carbon Disulfide	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methylene Chloride	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Vinyl Acetate	ND	1.0	EPA 8260C	4-2-18	4-2-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Butanone	ND	5.0	EPA 8260C	4-2-18	4-2-18	
Bromochloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chloroform	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Benzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Trichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Dibromomethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromodichloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	4-2-18	4-2-18	
Toluene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	

VOLATILES by EPA 8260C Page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	B-11					
Laboratory ID:	03-248-09					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Tetrachloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Hexanone	ND	2.0	EPA 8260C	4-2-18	4-2-18	
Dibromochloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chlorobenzene	0.36	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Ethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
m,p-Xylene	ND	0.40	EPA 8260C	4-2-18	4-2-18	
o-Xylene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Styrene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromoform	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Isopropylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
n-Propylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
n-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dibromo-3-chloropropane	e ND	1.0	EPA 8260C	4-2-18	4-2-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Naphthalene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Surrogate:	Percent Recovery	Control Limits				

Dibromofluoromethane 100 75-127 80-127 Toluene-d8 98 4-Bromofluorobenzene 96 78-125

VOLATILES by EPA 8260C METHOD BLANK QUALITY CONTROL

Page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0402W1					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chloromethane	ND	1.3	EPA 8260C	4-2-18	4-2-18	
Vinyl Chloride	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromomethane	ND	0.25	EPA 8260C	4-2-18	4-2-18	
Chloroethane	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Acetone	ND	5.0	EPA 8260C	4-2-18	4-2-18	
Iodomethane	ND	1.5	EPA 8260C	4-2-18	4-2-18	
Carbon Disulfide	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methylene Chloride	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Vinyl Acetate	ND	1.0	EPA 8260C	4-2-18	4-2-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Butanone	ND	5.0	EPA 8260C	4-2-18	4-2-18	
Bromochloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chloroform	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Benzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Trichloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Dibromomethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromodichloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	4-2-18	4-2-18	
Toluene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	4-2-18	4-2-18	

Project: 17-06520-000

VOLATILES by EPA 8260C METHOD BLANK QUALITY CONTROL

Page 2 of 2

Analyta	Pocult	PQL	Mothod	Date	Date Analyzed	Flage
Analyte	Result	PQL	Method	Prepared	Anaryzeu	Flags
Laboratory ID:	MB0402W1					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Tetrachloroethene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Hexanone	ND	2.0	EPA 8260C	4-2-18	4-2-18	
Dibromochloromethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Chlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Ethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
m,p-Xylene	ND	0.40	EPA 8260C	4-2-18	4-2-18	
o-Xylene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Styrene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromoform	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Isopropylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Bromobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	4-2-18	4-2-18	
n-Propylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
n-Butylbenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	4-2-18	4-2-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
Naphthalene	ND	1.0	EPA 8260C	4-2-18	4-2-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	4-2-18	4-2-18	
Surrogate:		· · · · · · · · · · · · · · · · · · ·				

Surrogate: Percent Recovery Control Limit
Dibromofluoromethane 96 75-127
Toluene-d8 97 80-127
4-Bromofluorobenzene 96 78-125



VOLATILES by EPA 8260C SB/SBD QUALITY CONTROL

	Result				Per	Percent		RPD		
Analyte			Spike Level		Recovery		Limits	RPD	Limit	Flags
SPIKE BLANKS										
Laboratory ID:	SB0402W1									
	SB	SBD	SB	SBD	SB	SBD				
1,1-Dichloroethene	9.74	9.66	10.0	10.0	97	97	63-126	1	21	
Benzene	9.79	9.70	10.0	10.0	98	97	78-122	1	19	
Trichloroethene	9.69	9.60	10.0	10.0	97	96	63-120	1	20	
Toluene	10.2	10.0	10.0	10.0	102	100	79-124	2	19	
Chlorobenzene	9.36	9.28	10.0	10.0	94	93	78-120	1	19	
Surrogate:										
Dibromofluoromethane					97	100	75-127			
Toluene-d8					99	102	80-127			
4-Bromofluorobenzene					98	101	78-125			



Data Qualifiers and Abbreviations

- A Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
- B The analyte indicated was also found in the blank sample.
- C The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
- E The value reported exceeds the quantitation range and is an estimate.
- F Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
- H The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
- I Compound recovery is outside of the control limits.
- J The value reported was below the practical quantitation limit. The value is an estimate.
- K Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
- L The RPD is outside of the control limits.
- M Hydrocarbons in the gasoline range are impacting the diesel range result.
- M1 Hydrocarbons in the gasoline range (toluene-naphthalene) are present in the sample.
- N Hydrocarbons in the lube oil range are impacting the diesel range result.
- N1 Hydrocarbons in diesel range are impacting lube oil range results.
- O Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
- P The RPD of the detected concentrations between the two columns is greater than 40.
- Q Surrogate recovery is outside of the control limits.
- S Surrogate recovery data is not available due to the necessary dilution of the sample.
- T The sample chromatogram is not similar to a typical ______.
- U The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- U1 The practical quantitation limit is elevated due to interferences present in the sample.
- V Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
- W Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
- X Sample extract treated with a mercury cleanup procedure.
- X1- Sample extract treated with a sulfuric acid/silica gel cleanup procedure.
- Y The calibration verification for this analyte exceeded the 20% drift specified in method 8260C, and therefore the reported result should be considered an estimate. The overall performance of the calibration verification standard met the acceptance criteria of the method.

7 -

ND - Not Detected at PQL

PQL - Practical Quantitation Limit

RPD - Relative Percent Difference





Chain of Custody

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	14648 NE	aboratory Testing Services 95th Street • Redmond, WA 98052		irnaround Req (in working da			La	abo	rato	ory	Nur	nbe	er:	()3	-	24	18									
Project Project	Number: OFIC P Manager: DVAL d by:	ark	2 D	ays [ndard (7 Days) H analysis 5 Da (other)	1 Day 3 Days	Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX	NWTPH-Gx	NWTPH-Dx (Acid / SG Clean-up)	Volatiles 8260C	Halogenated Volatiles 8260C	EDB EPA 8011 (Waters Only)	Semivolatiles 8270D/SIM (with low-level PAHs)	PAHs 8270D/SIM (low-level)	PCBs 8082A	Organochlorine Pesticides 8081B	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	TCLP Metals	HEM (oil and grease) 1664A	CPAHS by 82700/SIM			% Moisture
1	MWI		3/23/19	16 11	water	8	1	_		-	(X)										/			/	T	T	6,
2	MW2		1	12:16	1	8	V				$\langle \hat{\mathbf{x}} \rangle$										V			~	T		
3	MW3			15:26		8	V				(3)										V			V			
4	mw 4			14:10		8	1				8										V			1/			
5	MWS			16:05		8	/				(3)										1			V			
6	MW6			15:15		8	/				3										/			V			
7	B-06			11:32		8	V				3										/			V			
8	B-09			14:18		8	V			(3										V			V			
9	B-11		V	13:10	-	8	V			(3										/						
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-	uished		>	-0	286	_	-	51	23	ILX	1	15	S	-	-	en	2 f	CE	35,	Pt	we	lin	19	HC	WIP		
Recei	ved		1																				10				
Relino	uished													0	add	ed	3/	301	18	D-1	36	STA)				
Recei	ved																			Lev		. /		IV 🗆			
Revie	wed/Date			Reviewed/Da	te									Chro	matc	gram	s wit	h fina	l rep	ort 🗌	Elec	tronic	Data	Delive	rables	(EDDs) 🗆



14648 NE 95th Street, Redmond, WA 98052 • (425) 883-3881

July 2, 2018

Mark Ewbank Herrera Environmental Consultants, Inc. 2200 6th Avenue, Suite 1100 Seattle, WA 98121

Re: Analytical Data for Project 17-06520-000

Laboratory Reference No. 1806-222

Dear Mark:

Enclosed are the analytical results and associated quality control data for samples submitted on June 21, 2018.

The standard policy of OnSite Environmental, Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

David Baumeister Project Manager

Enclosures

Project: 17-06520-000

Case Narrative

Samples were collected on June 21, 2018 and received by the laboratory on June 21, 2018. They were maintained at the laboratory at a temperature of 2°C to 6°C.

Please note that any and all soil sample results are reported on a dry-weight basis, unless otherwise noted below.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

Project: 17-06520-000

NWTPH-Gx

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW1					
Laboratory ID:	06-222-01					
Gasoline	ND	100	NWTPH-Gx	6-25-18	6-25-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	71	66-117				
Client ID:	MW2					
Laboratory ID:	06-222-02					
Gasoline	ND	100	NWTPH-Gx	6-25-18	6-25-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	79	66-117				
Client ID:	MW3					
Laboratory ID:	06-222-03					
Gasoline	ND	100	NWTPH-Gx	6-25-18	6-25-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	75	66-117				
Client ID:	MW4					
Laboratory ID:	06-222-04					
Gasoline	ND	100	NWTPH-Gx	6-25-18	6-25-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	72	66-117				
Client ID:	MW5					
Laboratory ID:	06-222-05					
Gasoline	ND	100	NWTPH-Gx	6-25-18	6-25-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	75	66-117				
Client ID:	MW6					
Laboratory ID:	06-222-06					
Gasoline	ND	100	NWTPH-Gx	6-25-18	6-25-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	75	66-117				

Project: 17-06520-000

NWTPH-Gx

Omic. 49/2 (pps)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW7					
Laboratory ID:	06-222-07					
Gasoline	ND	100	NWTPH-Gx	6-25-18	6-25-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	75	66-117				
Client ID:	MW8					
Laboratory ID:	06-222-08					
Gasoline	ND	100	NWTPH-Gx	6-25-18	6-25-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	75	66-117				
Client ID:	MW9					
Laboratory ID:	06-222-09					
Gasoline	ND	100	NWTPH-Gx	6-25-18	6-25-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	77	66-117				
Client ID:	Trip Blank					
Laboratory ID:	06-222-10					
Gasoline	ND	100	NWTPH-Gx	6-25-18	6-25-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	75	66-117				

Project: 17-06520-000

NWTPH-Gx QUALITY CONTROL

Matrix: Water Units: ug/L (ppb)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0625W2					
Gasoline	ND	100	NWTPH-Gx	6-25-18	6-25-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	76	66-117				

					Source	Percent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Recovery	Limits	RPD	Limit	Flags
DUPLICATE										
Laboratory ID:	06-22	22-01								
	ORIG	DUP								
Gasoline	ND	ND	NA	NA		NA	NA	NA	30	
Surrogate:		•			•				•	•
Elugrobenzene						71 73	66-117			

73 66-117 Fluorobenzene

Project: 17-06520-000

NWTPH-Dx

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW1					
Laboratory ID:	06-222-01					
Diesel Range Organics	ND	0.26	NWTPH-Dx	6-22-18	6-22-18	
Lube Oil Range Organics	ND	0.41	NWTPH-Dx	6-22-18	6-22-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	86	50-150				
						
Client ID:	MW2					
Laboratory ID:	06-222-02					
Diesel Range Organics	ND	0.27	NWTPH-Dx	6-22-18	6-22-18	
Lube Oil Range Organics	ND -	0.43	NWTPH-Dx	6-22-18	6-22-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	85	50-150				
Client ID:	MW3					
Laboratory ID:	06-222-03					
Diesel Range Organics	ND	0.26	NWTPH-Dx	6-22-18	6-22-18	
Lube Oil Range Organics	ND ND	0.20	NWTPH-Dx	6-22-18	6-22-18	
Surrogate:	Percent Recovery	Control Limits	INVVIFII-DX	0-22-10	0-22-10	
o-Terphenyl	90	50-150				
0-Terprierryi	90	30-130				
Client ID:	MW4					
Laboratory ID:	06-222-04					
Diesel Range Organics	ND	0.27	NWTPH-Dx	6-22-18	6-22-18	
Lube Oil Range Organics	ND	0.43	NWTPH-Dx	6-22-18	6-22-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	88	50-150				
. ,						
Client ID:	MW5					
Laboratory ID:	06-222-05					
Diesel Range Organics	ND	0.26	NWTPH-Dx	6-22-18	6-22-18	
Lube Oil Range Organics	ND	0.41	NWTPH-Dx	6-22-18	6-22-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	94	50-150				
Oliana ID.	BANAZO					
Client ID:	MW6					
Laboratory ID:	06-222-06	0.00	ANATELLE	0.00.10	0.00.10	
Diesel Range Organics	ND	0.26	NWTPH-Dx	6-22-18	6-22-18	
Lube Oil Range Organics	ND 15	0.41	NWTPH-Dx	6-22-18	6-22-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	96	50-150				

Project: 17-06520-000

NWTPH-Dx

			Date	Date	
Result	PQL	Method	Prepared	Analyzed	Flags
MW7					
06-222-07					
ND	0.26	NWTPH-Dx	6-22-18	6-22-18	
ND	0.41	NWTPH-Dx	6-22-18	6-22-18	
Percent Recovery	Control Limits				
87	50-150				
06-222-08					
ND	0.26	NWTPH-Dx	6-22-18	6-22-18	
ND	0.41	NWTPH-Dx	6-22-18	6-22-18	
Percent Recovery	Control Limits				
86	50-150				
MW9					
06-222-09					
ND	0.26	NWTPH-Dx	6-22-18	6-22-18	
ND	0.41	NWTPH-Dx	6-22-18	6-22-18	
Percent Recovery	Control Limits				
92	50-150				
	MW7 06-222-07 ND ND Percent Recovery 87 MW8 06-222-08 ND ND Percent Recovery 86 MW9 06-222-09 ND ND ND Percent Recovery	MW7 06-222-07 ND 0.26 ND 0.41 Percent Recovery 87 Control Limits 50-150 MW8 06-222-08 ND 0.26 ND 0.41 Percent Recovery 86 Control Limits 50-150 MW9 06-222-09 ND 0.26 ND 0.26 ND 0.26 ND 0.41 Percent Recovery Control Limits Control Limits Control Limits	MW7 06-222-07 06-222-07 ND 0.26 NWTPH-Dx Percent Recovery 87 Control Limits 50-150 NWTPH-Dx MW8 06-222-08 NWTPH-Dx ND 0.26 NWTPH-Dx Percent Recovery 86 Control Limits 50-150 NWTPH-Dx ND 0.26 NWTPH-Dx ND 0.26 NWTPH-Dx ND 0.41 NWTPH-Dx Percent Recovery Control Limits NWTPH-Dx Percent Recovery Control Limits Control Limits	Result PQL Method Prepared MW7 06-222-07 ND 0.26 NWTPH-Dx 6-22-18 ND 0.41 NWTPH-Dx 6-22-18 Percent Recovery 87 50-150 ND 0.26 NWTPH-Dx 6-22-18 ND 0.41 NWTPH-Dx 6-22-18 Percent Recovery 86 50-150 ND 0.26 NWTPH-Dx 6-22-18 ND 0.26 NWTPH-Dx 6-22-18 ND 0.26 NWTPH-Dx 6-22-18 ND 0.41 NWTPH-Dx 6-22-18 Percent Recovery Control Limits 6-22-18	Result PQL Method Prepared Analyzed MW7 06-222-07 06-222-07 6-22-18 6-22-18 ND 0.26 NWTPH-Dx 6-22-18 6-22-18 ND 0.41 NWTPH-Dx 6-22-18 6-22-18 Percent Recovery 06-222-08 50-150 NWTPH-Dx 6-22-18 6-22-18 ND 0.41 NWTPH-Dx 6-22-18 6-22-18 Percent Recovery 06-222-09 Control Limits 50-150 6-22-18 6-22-18 6-22-18 ND 0.26 NWTPH-Dx 6-22-18 6-22-18 ND 0.26 NWTPH-Dx 6-22-18 6-22-18 ND 0.41 NWTPH-Dx 6-22-18 6-22-18 Percent Recovery Control Limits 6-22-18 6-22-18

Project: 17-06520-000

NWTPH-Dx QUALITY CONTROL

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0622W1					
Diesel Range Organics	ND	0.25	NWTPH-Dx	6-22-18	6-22-18	
Lube Oil Range Organics	ND	0.40	NWTPH-Dx	6-22-18	6-22-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	87	50-150				

Analyte	Res	sult	Spike	Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
DUPLICATE										_
Laboratory ID:	06-22	22-02								
	ORIG	DUP								
Diesel Range	ND	ND	NA	NA		NA	NA	NA	NA	
Lube Oil Range	ND	ND	NA	NA		NA	NA	NA	NA	
Surrogate: o-Terphenyl						85 93	50-150			

Project: 17-06520-000

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW1					
Laboratory ID:	06-222-01					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chloromethane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Vinyl Chloride	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromomethane	ND	0.31	EPA 8260C	6-25-18	6-25-18	
Chloroethane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Acetone	ND	5.0	EPA 8260C	6-25-18	6-25-18	
Iodomethane	ND	1.9	EPA 8260C	6-25-18	6-25-18	
Carbon Disulfide	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methylene Chloride	ND	1.0	EPA 8260C	6-25-18	6-25-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Vinyl Acetate	ND	1.0	EPA 8260C	6-25-18	6-25-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Butanone	ND	5.0	EPA 8260C	6-25-18	6-25-18	
Bromochloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chloroform	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Benzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Trichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Dibromomethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromodichloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	6-25-18	6-25-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	6-25-18	6-25-18	
Toluene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	

Project: 17-06520-000

VOLATILES EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW1					
Laboratory ID:	06-222-01					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Tetrachloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Hexanone	ND	2.0	EPA 8260C	6-25-18	6-25-18	
Dibromochloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Ethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
m,p-Xylene	ND	0.40	EPA 8260C	6-25-18	6-25-18	
o-Xylene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Styrene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromoform	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Isopropylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2,3-Trichloropropane	ND	0.27	EPA 8260C	6-25-18	6-25-18	
n-Propylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
n-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Naphthalene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	98	75-127				

 Dibromofluoromethane
 98
 75-127

 Toluene-d8
 98
 80-127

 4-Bromofluorobenzene
 98
 78-125



Project: 17-06520-000

VOLATILES EPA 8260C page 1 of 2

Client ID: MW2 Laboratory ID: 06-222-02 Dichlorodiffuoromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloromethane ND 1.0 EPA 8260C 6-25-18 6-25-18 Chloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromomethane ND 0.31 EPA 8260C 6-25-18 6-25-18 Bromomethane ND 1.0 EPA 8260C 6-25-18 6-25-18 Chloroethane ND 1.0 EPA 8260C 6-25-18 6-25-18 Chloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Trichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Iodomethane ND 1.9 EPA 8260C 6-25-18 6-25-18 Carbon Disulfide ND 0.20 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 (trans) 1					Date	Date	
Laboratory ID:	Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Dichlorodiffluoromethane	Client ID:	MW2					
Chloromethane ND 1.0 EPA 8260C 6-25-18 6-25-18 Vinyl Chloride ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromomethane ND 0.31 EPA 8260C 6-25-18 6-25-18 Chloroethane ND 1.0 EPA 8260C 6-25-18 6-25-18 Trichlorofluoromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Trichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Acetone ND 5.0 EPA 8260C 6-25-18 6-25-18 Iodomethane ND 1.9 EPA 8260C 6-25-18 6-25-18 Carbon Disulfide ND 1.9 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 0.20 EPA 8260C 6-25-18 6-25-18	Laboratory ID:	06-222-02					
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Bromomethane	Chloromethane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Chloroethane ND 1.0 EPA 8260C 6-25-18 6-25-18 Trichlorofluoromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Acetone ND 5.0 EPA 8260C 6-25-18 6-25-18 Iodomethane ND 1.9 EPA 8260C 6-25-18 6-25-18 Carbon Disulfide ND 1.0 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 0.20 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyle Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18	Vinyl Chloride	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Trichlorofluoromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Acetone ND 5.0 EPA 8260C 6-25-18 6-25-18 Iodomethane ND 1.9 EPA 8260C 6-25-18 6-25-18 Carbon Disulfide ND 0.20 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Minyl Acetate ND 1.0 EPA 8260C 6-25-18 6-25-18 Q-2-Dichloroethane ND 0.20 EPA 8260C 6-25-18	Bromomethane	ND	0.31	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Acetone ND 5.0 EPA 8260C 6-25-18 6-25-18 Iodomethane ND 1.9 EPA 8260C 6-25-18 6-25-18 Carbon Disulfide ND 0.20 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 0.20 EPA 8260C 6-25-18 6-25-18 (trans) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Vinyl Acetate ND 0.20 EPA 8260C 6-25-18 6-25-18 Vinyl Acetate ND 0.20 EPA 8260C 6-25-18 <t< td=""><td>Chloroethane</td><td>ND</td><td>1.0</td><td>EPA 8260C</td><td>6-25-18</td><td>6-25-18</td><td></td></t<>	Chloroethane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Acetone ND 5.0 EPA 8260C 6-25-18 6-25-18 Iodomethane ND 1.9 EPA 8260C 6-25-18 6-25-18 Carbon Disulfide ND 0.20 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroptopane ND 0.20 EPA 8260C 6-25-18 6-25-18 2,2-Dichloroptopenee ND 0.20 EPA 8260C 6-25-18	Trichlorofluoromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Indomethane ND	1,1-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Carbon Disulfide ND 0.20 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 (trans) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Vinyl Acetate ND 0.20 EPA 8260C 6-25-18 6-25-18 Q-2-Dichloroptane ND 0.20 EPA 8260C 6-25-18 6-25-18 Q-2-Dichloroptane ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromochloromethane ND 0.20 EPA 8260C	Acetone	ND	5.0	EPA 8260C	6-25-18	6-25-18	
Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 (trans) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Vinyl Acetate ND 1.0 EPA 8260C 6-25-18 6-25-18 2,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 (cis) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromochloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloropthane ND 0.20 EPA 8260C 6-25-18 <td< td=""><td>Iodomethane</td><td>ND</td><td>1.9</td><td>EPA 8260C</td><td>6-25-18</td><td>6-25-18</td><td></td></td<>	Iodomethane	ND	1.9	EPA 8260C	6-25-18	6-25-18	
(trans) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Vinyl Acetate ND 1.0 EPA 8260C 6-25-18 6-25-18 2,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 (cis) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 5.0 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromochloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 <	Carbon Disulfide	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Vinyl Acetate ND 1.0 EPA 8260C 6-25-18 6-25-18 2,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 (cis) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 5.0 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 <t< td=""><td>Methylene Chloride</td><td>ND</td><td>1.0</td><td>EPA 8260C</td><td>6-25-18</td><td>6-25-18</td><td></td></t<>	Methylene Chloride	ND	1.0	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Vinyl Acetate ND 1.0 EPA 8260C 6-25-18 6-25-18 2,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 4,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 5.0 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromochloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 <td>(trans) 1,2-Dichloroethene</td> <td>ND</td> <td>0.20</td> <td>EPA 8260C</td> <td>6-25-18</td> <td>6-25-18</td> <td></td>	(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
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2,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 (cis) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 5.0 EPA 8260C 6-25-18 6-25-18 Bromochloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1,1-Trichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA 8260C <td>1,1-Dichloroethane</td> <td>ND</td> <td>0.20</td> <td>EPA 8260C</td> <td>6-25-18</td> <td>6-25-18</td> <td></td>	1,1-Dichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
(cis) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 5.0 EPA 8260C 6-25-18 6-25-18 Bromochloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1,1-Trichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Benzene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18	Vinyl Acetate	ND	1.0	EPA 8260C	6-25-18	6-25-18	
2-Butanone ND 5.0 EPA 8260C 6-25-18 6-25-18 Bromochloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1,1-Trichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Benzene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA 8260C 6-25-18	2,2-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromochloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1,1-Trichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Benzene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Trichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromodichloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 <td>(cis) 1,2-Dichloroethene</td> <td>ND</td> <td>0.20</td> <td>EPA 8260C</td> <td>6-25-18</td> <td>6-25-18</td> <td></td>	(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1,1-Trichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Benzene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Trichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 </td <td>2-Butanone</td> <td>ND</td> <td>5.0</td> <td>EPA 8260C</td> <td>6-25-18</td> <td>6-25-18</td> <td></td>	2-Butanone	ND	5.0	EPA 8260C	6-25-18	6-25-18	
1,1,1-Trichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Benzene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Trichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromodichloromethane ND 1.0 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 1.0	Bromochloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Benzene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Trichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromodichloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18	Chloroform	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Benzene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Trichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromodichloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	1,1,1-Trichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Benzene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Trichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromodichloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	Carbon Tetrachloride	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Trichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromodichloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	1,1-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Trichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromodichloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	Benzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromodichloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	1,2-Dichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Dibromomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromodichloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	Trichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromodichloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	1,2-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	Dibromomethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
(cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	Bromodichloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
(cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	6-25-18	6-25-18	
(trans) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18	Toluene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
	(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	

Project: 17-06520-000

VOLATILES EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW2					
Laboratory ID:	06-222-02					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Tetrachloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Hexanone	ND	2.0	EPA 8260C	6-25-18	6-25-18	
Dibromochloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Ethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
m,p-Xylene	ND	0.40	EPA 8260C	6-25-18	6-25-18	
o-Xylene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Styrene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromoform	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Isopropylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2,3-Trichloropropane	ND	0.27	EPA 8260C	6-25-18	6-25-18	
n-Propylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
n-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Naphthalene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	96	75-127				

4-Bromofluorobenzene

Toluene-d8

80-127

78-125

97

95

Project: 17-06520-000

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW3					
Laboratory ID:	06-222-03					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chloromethane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Vinyl Chloride	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromomethane	ND	0.31	EPA 8260C	6-25-18	6-25-18	
Chloroethane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Acetone	ND	5.0	EPA 8260C	6-25-18	6-25-18	
lodomethane	ND	1.9	EPA 8260C	6-25-18	6-25-18	
Carbon Disulfide	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methylene Chloride	ND	1.0	EPA 8260C	6-25-18	6-25-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Vinyl Acetate	ND	1.0	EPA 8260C	6-25-18	6-25-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
(cis) 1,2-Dichloroethene	0.33	0.20	EPA 8260C	6-25-18	6-25-18	
2-Butanone	ND	5.0	EPA 8260C	6-25-18	6-25-18	
Bromochloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chloroform	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Benzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Trichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Dibromomethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromodichloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	6-25-18	6-25-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	6-25-18	6-25-18	
Toluene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	

Project: 17-06520-000

VOLATILES EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW3					
Laboratory ID:	06-222-03					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Tetrachloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Hexanone	ND	2.0	EPA 8260C	6-25-18	6-25-18	
Dibromochloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Ethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
m,p-Xylene	ND	0.40	EPA 8260C	6-25-18	6-25-18	
o-Xylene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Styrene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromoform	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Isopropylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2,3-Trichloropropane	ND	0.27	EPA 8260C	6-25-18	6-25-18	
n-Propylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
n-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Naphthalene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	99	75-127				
Talvana dO	00	00.407				



Project: 17-06520-000

VOLATILES EPA 8260C page 1 of 2

Analyte Result PQL Method Prepared Analyzed Flags Client ID: MW4 Laboratory ID: 06-222-04 □ <					Date	Date	
Laboratory ID: 06-222-04 Dichlorodifluoromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloromethane ND 1.0 EPA 8260C 6-25-18 6-25-18 Vinyl Chloride ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromomethane ND 0.31 EPA 8260C 6-25-18 6-25-18 Chloroethane ND 1.0 EPA 8260C 6-25-18 6-25-18 Chloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Trichlorofluoromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Acetone ND 5.0 EPA 8260C 6-25-18 6-25-18 Iodomethane ND 1.9 EPA 8260C 6-25-18 6-25-18 Iodomethane ND 1.9 EPA 8260C 6-25-18 6-25-18 Iodomethane ND 1.0 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 1.0 EPA 8260C	Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Dichlorodifluoromethane	Client ID:	MW4					
Chloromethane ND 1.0 EPA 8260C 6-25-18 6-25-18 Vinyl Chloride ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromomethane ND 0.31 EPA 8260C 6-25-18 6-25-18 Chloroethane ND 1.0 EPA 8260C 6-25-18 6-25-18 Trichlorofluoromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Acetone ND 5.0 EPA 8260C 6-25-18 6-25-18 Idodomethane ND 1.9 EPA 8260C 6-25-18 6-25-18 Carbon Disulfide ND 1.9 EPA 8260C 6-25-18 6-25-18 Carbon Disulfide ND 1.0 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 0.20 EPA 8260C 6-25-18 6-25-18	Laboratory ID:	06-222-04					
Vinyl Chloride ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromomethane ND 0.31 EPA 8260C 6-25-18 6-25-18 Chloroethane ND 1.0 EPA 8260C 6-25-18 6-25-18 Trichlorofluoromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Acetone ND 5.0 EPA 8260C 6-25-18 6-25-18 Iodomethane ND 1.9 EPA 8260C 6-25-18 6-25-18 Iodomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Iodomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl L-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Vinyl Acetate ND 0.20 EPA 8260C 6-25-18 6-25-18 <t< td=""><td>Dichlorodifluoromethane</td><td>ND</td><td>0.20</td><td>EPA 8260C</td><td>6-25-18</td><td>6-25-18</td><td></td></t<>	Dichlorodifluoromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromomethane ND 0.31 EPA 8260C 6-25-18 6-25-18 Chloroethane ND 1.0 EPA 8260C 6-25-18 6-25-18 Trichlorofluoromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Acetone ND 5.0 EPA 8260C 6-25-18 6-25-18 Iodomethane ND 1.9 EPA 8260C 6-25-18 6-25-18 Carbon Disulfide ND 0.20 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 (trans) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 (trans) 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 (trans) 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 (cis) 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-1	Chloromethane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Chloroethane ND 1.0 EPA 8260C 6-25-18 6-25-18 Trichlorofluoromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Acetone ND 5.0 EPA 8260C 6-25-18 6-25-18 Iodomethane ND 1.9 EPA 8260C 6-25-18 6-25-18 Carbon Disulfide ND 1.0 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 0.20 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyle Ebury Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethane ND 0.20 EPA 8260C 6-25-18	Vinyl Chloride	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Trichlorofluoromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Acetone ND 5.0 EPA 8260C 6-25-18 6-25-18 Iodomethane ND 1.9 EPA 8260C 6-25-18 6-25-18 Carbon Disulfide ND 0.20 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethane ND 0.20 EPA 8260C 6-25-18	Bromomethane	ND	0.31	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Acetone ND 5.0 EPA 8260C 6-25-18 6-25-18 Iodomethane ND 1.9 EPA 8260C 6-25-18 6-25-18 Carbon Disulfide ND 0.20 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 (trans) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Vinyl Acetate ND 1.0 EPA 8260C 6-25-18 6-25-18 Vinyl Acetate ND 0.20 EPA 8260C 6-25-18 6-25-18 (cis) 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 (cis) 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18	Chloroethane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Acetone ND 5.0 EPA 8260C 6-25-18 6-25-18 Iodomethane ND 1.9 EPA 8260C 6-25-18 6-25-18 Carbon Disulfide ND 0.20 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 (trans) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 2,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 2,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18	Trichlorofluoromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Iodomethane ND 1.9 EPA 8260C 6-25-18 6-25-18 Carbon Disulfide ND 0.20 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 Methyl Erbury Lether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl L-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Vinyl Acetate ND 1.0 EPA 8260C 6-25-18 6-25-18 2,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 2,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 2,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Businer ND 0.20 EPA 8260C 6-25-18	1,1-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Carbon Disulfide ND 0.20 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 (trans) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 2,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 2,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 2,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 0.20 EPA 8260C 6-25-18 6-25-18 3-Bromochloromethane ND 0.20 EPA 8260C 6-	Acetone	ND	5.0	EPA 8260C	6-25-18	6-25-18	
Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 (trans) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Vinyl Acetate ND 1.0 EPA 8260C 6-25-18 6-25-18 2,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 (cis) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 0.20 EPA 8260C 6-25-18 6-25-18 3-Bromochloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 4-LITATION Contraction ND 0.20 EPA 8260C 6-25-18	Iodomethane	ND	1.9	EPA 8260C	6-25-18	6-25-18	
(trans) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Vinyl Acetate ND 1.0 EPA 8260C 6-25-18 6-25-18 2,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 (cis) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 5.0 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1,1-Trichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18	Carbon Disulfide	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Vinyl Acetate ND 1.0 EPA 8260C 6-25-18 6-25-18 2,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 (cis) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 5.0 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromochloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18	Methylene Chloride	ND	1.0	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Vinyl Acetate ND 1.0 EPA 8260C 6-25-18 6-25-18 2,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 (cis) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 5.0 EPA 8260C 6-25-18 6-25-18 Bromochloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Benzene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 </td <td>(trans) 1,2-Dichloroethene</td> <td>ND</td> <td>0.20</td> <td>EPA 8260C</td> <td>6-25-18</td> <td>6-25-18</td> <td></td>	(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
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2,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 (cis) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 5.0 EPA 8260C 6-25-18 6-25-18 Bromochloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1,1-Trichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA	1,1-Dichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
(cis) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 5.0 EPA 8260C 6-25-18 6-25-18 Bromochloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1,1-Trichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromodichloromethane ND 0.20 EPA 8260C 6-25	Vinyl Acetate	ND	1.0	EPA 8260C	6-25-18	6-25-18	
2-Butanone ND 5.0 EPA 8260C 6-25-18 6-25-18 Bromochloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1,1-Trichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 <td>2,2-Dichloropropane</td> <td>ND</td> <td>0.20</td> <td>EPA 8260C</td> <td>6-25-18</td> <td>6-25-18</td> <td></td>	2,2-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromochloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1,1-Trichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C	(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
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1,1,1-Trichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Benzene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Trichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromodichloromethane ND 1.0 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260	Bromochloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
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1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Trichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromodichloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	1,1-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Trichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromodichloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	Benzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromodichloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	1,2-Dichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Dibromomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromodichloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	Trichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromodichloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	1,2-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	Dibromomethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
(cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	Bromodichloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
	Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	6-25-18	6-25-18	
(trans) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18	Toluene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
	(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	

Project: 17-06520-000

VOLATILES EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW4					
Laboratory ID:	06-222-04					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Tetrachloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Hexanone	ND	2.0	EPA 8260C	6-25-18	6-25-18	
Dibromochloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chlorobenzene	2.5	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Ethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
m,p-Xylene	ND	0.40	EPA 8260C	6-25-18	6-25-18	
o-Xylene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Styrene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromoform	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Isopropylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2,3-Trichloropropane	ND	0.27	EPA 8260C	6-25-18	6-25-18	
n-Propylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
n-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Naphthalene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	100	75-127				

 Dibromofluoromethane
 100
 75-127

 Toluene-d8
 100
 80-127

 4-Bromofluorobenzene
 97
 78-125



Project: 17-06520-000

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW5					
Laboratory ID:	06-222-05					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chloromethane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Vinyl Chloride	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromomethane	ND	0.31	EPA 8260C	6-25-18	6-25-18	
Chloroethane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Acetone	ND	5.0	EPA 8260C	6-25-18	6-25-18	
lodomethane	ND	1.9	EPA 8260C	6-25-18	6-25-18	
Carbon Disulfide	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methylene Chloride	ND	1.0	EPA 8260C	6-25-18	6-25-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Vinyl Acetate	ND	1.0	EPA 8260C	6-25-18	6-25-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Butanone	ND	5.0	EPA 8260C	6-25-18	6-25-18	
Bromochloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chloroform	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Benzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Trichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Dibromomethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromodichloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	6-25-18	6-25-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	6-25-18	6-25-18	
Toluene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	

Project: 17-06520-000

VOLATILES EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW5					
Laboratory ID:	06-222-05					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Tetrachloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Hexanone	ND	2.0	EPA 8260C	6-25-18	6-25-18	
Dibromochloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Ethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
m,p-Xylene	ND	0.40	EPA 8260C	6-25-18	6-25-18	
o-Xylene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Styrene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromoform	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Isopropylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2,3-Trichloropropane	ND	0.27	EPA 8260C	6-25-18	6-25-18	
n-Propylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
n-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Naphthalene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	99	75-127				

 Dibromofluoromethane
 99
 75-127

 Toluene-d8
 100
 80-127

 4-Bromofluorobenzene
 97
 78-125

Project: 17-06520-000

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW6					
Laboratory ID:	06-222-06					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chloromethane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Vinyl Chloride	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromomethane	ND	0.31	EPA 8260C	6-25-18	6-25-18	
Chloroethane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Acetone	ND	5.0	EPA 8260C	6-25-18	6-25-18	
lodomethane	ND	1.9	EPA 8260C	6-25-18	6-25-18	
Carbon Disulfide	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methylene Chloride	ND	1.0	EPA 8260C	6-25-18	6-25-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Vinyl Acetate	ND	1.0	EPA 8260C	6-25-18	6-25-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Butanone	ND	5.0	EPA 8260C	6-25-18	6-25-18	
Bromochloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chloroform	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Benzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Trichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Dibromomethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromodichloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	6-25-18	6-25-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	6-25-18	6-25-18	
Toluene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	

Project: 17-06520-000

VOLATILES EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW6					
Laboratory ID:	06-222-06					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Tetrachloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Hexanone	ND	2.0	EPA 8260C	6-25-18	6-25-18	
Dibromochloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Ethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
m,p-Xylene	ND	0.40	EPA 8260C	6-25-18	6-25-18	
o-Xylene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Styrene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromoform	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Isopropylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2,3-Trichloropropane	ND	0.27	EPA 8260C	6-25-18	6-25-18	
n-Propylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
n-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Naphthalene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	100	75-127				
-						

4-Bromofluorobenzene

Toluene-d8

80-127

78-125

99

97

Project: 17-06520-000

VOLATILES EPA 8260C page 1 of 2

Analyte				Date	Date	
	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW7					
Laboratory ID:	06-222-07					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chloromethane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Vinyl Chloride	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromomethane	ND	0.31	EPA 8260C	6-25-18	6-25-18	
Chloroethane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Acetone	ND	5.0	EPA 8260C	6-25-18	6-25-18	
lodomethane	ND	1.9	EPA 8260C	6-25-18	6-25-18	
Carbon Disulfide	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methylene Chloride	ND	1.0	EPA 8260C	6-25-18	6-25-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Vinyl Acetate	ND	1.0	EPA 8260C	6-25-18	6-25-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Butanone	ND	5.0	EPA 8260C	6-25-18	6-25-18	
Bromochloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chloroform	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Benzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Trichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Dibromomethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromodichloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	6-25-18	6-25-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	6-25-18	6-25-18	
Toluene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	

Project: 17-06520-000

VOLATILES EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW7					
Laboratory ID:	06-222-07					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Tetrachloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Hexanone	ND	2.0	EPA 8260C	6-25-18	6-25-18	
Dibromochloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Ethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
m,p-Xylene	ND	0.40	EPA 8260C	6-25-18	6-25-18	
o-Xylene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Styrene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromoform	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Isopropylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2,3-Trichloropropane	ND	0.27	EPA 8260C	6-25-18	6-25-18	
n-Propylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
n-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Naphthalene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	99	75-127				
-						

4-Bromofluorobenzene

Toluene-d8

80-127

78-125

100

99

Project: 17-06520-000

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW8					
Laboratory ID:	06-222-08					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chloromethane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Vinyl Chloride	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromomethane	ND	0.31	EPA 8260C	6-25-18	6-25-18	
Chloroethane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Acetone	ND	5.0	EPA 8260C	6-25-18	6-25-18	
lodomethane	ND	1.9	EPA 8260C	6-25-18	6-25-18	
Carbon Disulfide	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methylene Chloride	ND	1.0	EPA 8260C	6-25-18	6-25-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Vinyl Acetate	ND	1.0	EPA 8260C	6-25-18	6-25-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Butanone	ND	5.0	EPA 8260C	6-25-18	6-25-18	
Bromochloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chloroform	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Benzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Trichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Dibromomethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromodichloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	6-25-18	6-25-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	6-25-18	6-25-18	
Toluene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	

Project: 17-06520-000

VOLATILES EPA 8260C page 2 of 2

Date Date Analyte Result **PQL** Method **Prepared Analyzed** Flags Client ID: 8WM Laboratory ID: 06-222-08 1,1,2-Trichloroethane 0.20 **EPA 8260C** ND 6-25-18 6-25-18 Tetrachloroethene ND 0.20 **EPA 8260C** 6-25-18 6-25-18 1,3-Dichloropropane ND 0.20 **EPA 8260C** 6-25-18 6-25-18 2-Hexanone ND 2.0 **EPA 8260C** 6-25-18 6-25-18 Dibromochloromethane ND 0.20 **EPA 8260C** 6-25-18 6-25-18 1.2-Dibromoethane ND 0.20 6-25-18 **EPA 8260C** 6-25-18 ND 0.20 **EPA 8260C** 6-25-18 Chlorobenzene 6-25-18 1,1,1,2-Tetrachloroethane ND 0.20 **EPA 8260C** 6-25-18 6-25-18 Ethylbenzene ND 0.20 **EPA 8260C** 6-25-18 6-25-18 m,p-Xylene ND 0.40 **EPA 8260C** 6-25-18 6-25-18 ND 0.20 6-25-18 6-25-18 o-Xylene **EPA 8260C** Styrene ND 0.20 **EPA 8260C** 6-25-18 6-25-18 Bromoform ND 1.0 **EPA 8260C** 6-25-18 6-25-18 Isopropylbenzene ND 0.20 **EPA 8260C** 6-25-18 6-25-18 Bromobenzene ND 0.20 **EPA 8260C** 6-25-18 6-25-18 1,1,2,2-Tetrachloroethane ND 0.20 **EPA 8260C** 6-25-18 6-25-18 1,2,3-Trichloropropane ND 0.27 6-25-18 **EPA 8260C** 6-25-18 ND 0.20 **EPA 8260C** 6-25-18 6-25-18 n-Propylbenzene ND 0.20 **EPA 8260C** 6-25-18 2-Chlorotoluene 6-25-18 4-Chlorotoluene ND 0.20 **EPA 8260C** 6-25-18 6-25-18 1.3.5-Trimethylbenzene ND 0.20 **EPA 8260C** 6-25-18 6-25-18 tert-Butylbenzene ND 0.20 **EPA 8260C** 6-25-18 6-25-18 1,2,4-Trimethylbenzene ND 0.20 **EPA 8260C** 6-25-18 6-25-18 ND 0.20 **EPA 8260C** sec-Butylbenzene 6-25-18 6-25-18 ND 0.20 **EPA 8260C** 6-25-18 6-25-18 1,3-Dichlorobenzene p-Isopropyltoluene ND 0.20 **EPA 8260C** 6-25-18 6-25-18 1,4-Dichlorobenzene ND 0.20 **EPA 8260C** 6-25-18 6-25-18 1,2-Dichlorobenzene ND 0.20 **EPA 8260C** 6-25-18 6-25-18 ND 0.20 **EPA 8260C** 6-25-18 n-Butylbenzene 6-25-18 1,2-Dibromo-3-chloropropane ND 1.0 **EPA 8260C** 6-25-18 6-25-18 1.2.4-Trichlorobenzene ND 0.20 **EPA 8260C** 6-25-18 6-25-18 Hexachlorobutadiene ND 1.0 **EPA 8260C** 6-25-18 6-25-18 Naphthalene ND 1.0 **EPA 8260C** 6-25-18 6-25-18 1,2,3-Trichlorobenzene ND 0.20 **EPA 8260C** 6-25-18 6-25-18 Surrogate: Percent Recovery Control Limits Dibromofluoromethane 101 75-127

4-Bromofluorobenzene

Toluene-d8

80-127

78-125

100

98

Project: 17-06520-000

VOLATILES EPA 8260C

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW9					
Laboratory ID:	06-222-09					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chloromethane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Vinyl Chloride	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromomethane	ND	0.31	EPA 8260C	6-25-18	6-25-18	
Chloroethane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Acetone	ND	5.0	EPA 8260C	6-25-18	6-25-18	
lodomethane	ND	1.9	EPA 8260C	6-25-18	6-25-18	
Carbon Disulfide	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methylene Chloride	ND	1.0	EPA 8260C	6-25-18	6-25-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Vinyl Acetate	ND	1.0	EPA 8260C	6-25-18	6-25-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Butanone	ND	5.0	EPA 8260C	6-25-18	6-25-18	
Bromochloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chloroform	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Benzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Trichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Dibromomethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromodichloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	6-25-18	6-25-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	6-25-18	6-25-18	
Toluene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	

Project: 17-06520-000

VOLATILES EPA 8260C

page 2 of 2

Analyte Client ID: Laboratory ID: 1,1,2-Trichloroethane Tetrachloroethene	Result MW9 06-222-09 ND ND ND ND ND ND ND ND ND	0.20 0.20 0.20	Method EPA 8260C EPA 8260C	Prepared 6-25-18	Analyzed 6-25-18	Flags
Laboratory ID: 1,1,2-Trichloroethane	06-222-09 ND ND ND ND	0.20 0.20			6-25-18	
1,1,2-Trichloroethane	ND ND ND ND	0.20 0.20			6-25-18	
, ,	ND ND ND	0.20 0.20			6-25-18	
Tetrachloroethene	ND ND	0.20	EPA 8260C			
	ND			6-25-18	6-25-18	
1,3-Dichloropropane			EPA 8260C	6-25-18	6-25-18	
2-Hexanone	ND	2.0	EPA 8260C	6-25-18	6-25-18	
Dibromochloromethane		0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chlorobenzene	0.36	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Ethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
m,p-Xylene	ND	0.40	EPA 8260C	6-25-18	6-25-18	
o-Xylene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Styrene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromoform	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Isopropylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2,3-Trichloropropane	ND	0.27	EPA 8260C	6-25-18	6-25-18	
n-Propylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
n-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Naphthalene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	99	75-127				

 Dibromofluoromethane
 99
 75-127

 Toluene-d8
 100
 80-127

 4-Bromofluorobenzene
 98
 78-125



Project: 17-06520-000

VOLATILES EPA 8260C page 1 of 2

Client ID: Trip Blank Laboratory ID: 06-222-10 Dichlorodifluoromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloromethane ND 1.0 EPA 8260C 6-25-18 6-25-18 Chloromethane ND 1.0 EPA 8260C 6-25-18 6-25-18 Bromomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromomethane ND 1.0 EPA 8260C 6-25-18 6-25-18 Chloroethane ND 1.0 EPA 8260C 6-25-18 6-25-18 Chloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Trichlorofluoromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Iclodomethane ND 5.0 EPA 8260C 6-25-18 6-25-18 Iclodomethane ND 1.9 EPA 8260C 6-25-18 6-25-18 Iclodomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Ic					Date	Date	
Laboratory ID: 06-222-10 Dichlorodifluoromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloromethane ND 1.0 EPA 8260C 6-25-18 6-25-18 Vinyl Chloride ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromomethane ND 0.31 EPA 8260C 6-25-18 6-25-18 Chloroethane ND 1.0 EPA 8260C 6-25-18 6-25-18 Chloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Acetone ND 5.0 EPA 8260C 6-25-18 6-25-18 Indomethane ND 1.9 EPA 8260C 6-25-18 6-25-18 Carbon Disulfide ND 0.20 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 Kirans) 1,2-Dichloroethene ND 0.20	Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Dichlorodifluoromethane	Client ID:	Trip Blank					
Chloromethane ND 1.0 EPA 8260C 6-25-18 6-25-18 Vinyl Chloride ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromomethane ND 0.31 EPA 8260C 6-25-18 6-25-18 Chloroethane ND 1.0 EPA 8260C 6-25-18 6-25-18 Trichlorofluoromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Acetone ND 5.0 EPA 8260C 6-25-18 6-25-18 Iodomethane ND 1.9 EPA 8260C 6-25-18 6-25-18 Carbon Disulfide ND 0.20 EPA 8260C 6-25-18 6-25-18	Laboratory ID:	06-222-10					
Vinyl Chloride ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromomethane ND 0.31 EPA 8260C 6-25-18 6-25-18 Chloroethane ND 1.0 EPA 8260C 6-25-18 6-25-18 Trichlorofluoromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Acetone ND 5.0 EPA 8260C 6-25-18 6-25-18 Acetone ND 1.9 EPA 8260C 6-25-18 6-25-18 Iodomethane ND 1.9 EPA 8260C 6-25-18 6-25-18 Iodomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyliched ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyliched ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl L-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 <td< td=""><td>Dichlorodifluoromethane</td><td>ND</td><td>0.20</td><td>EPA 8260C</td><td>6-25-18</td><td>6-25-18</td><td></td></td<>	Dichlorodifluoromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromomethane	Chloromethane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Chloroethane ND 1.0 EPA 8260C 6-25-18 6-25-18 Trichlorofluoromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Acetone ND 5.0 EPA 8260C 6-25-18 6-25-18 Iodomethane ND 1.9 EPA 8260C 6-25-18 6-25-18 Carbon Disulfide ND 1.0 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 0.20 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 0.20 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyle Ether ND 0.20 EPA 8260C 6-25-18 6-25-18	Vinyl Chloride	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Trichlorofluoromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Acetone ND 5.0 EPA 8260C 6-25-18 6-25-18 Iodomethane ND 1.9 EPA 8260C 6-25-18 6-25-18 Carbon Disulfide ND 0.20 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 Methyl EButyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Minyl Acetate ND 1.0 EPA 8260C 6-25-18 6-25-18 Q-2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 <	Bromomethane	ND	0.31	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Acetone ND 5.0 EPA 8260C 6-25-18 6-25-18 Iodomethane ND 1.9 EPA 8260C 6-25-18 6-25-18 Carbon Disulfide ND 0.20 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 0.20 EPA 8260C 6-25-18 6-25-18 (trans) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Vinyl Acetate ND 1.0 EPA 8260C 6-25-18 6-25-18 Vinyl Acetate ND 0.20 EPA 8260C 6-25-18 <td< td=""><td>Chloroethane</td><td>ND</td><td>1.0</td><td>EPA 8260C</td><td>6-25-18</td><td>6-25-18</td><td></td></td<>	Chloroethane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Acetone ND 5.0 EPA 8260C 6-25-18 6-25-18 Iodomethane ND 1.9 EPA 8260C 6-25-18 6-25-18 Carbon Disulfide ND 0.20 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Cis) 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 <td>Trichlorofluoromethane</td> <td>ND</td> <td>0.20</td> <td>EPA 8260C</td> <td>6-25-18</td> <td>6-25-18</td> <td></td>	Trichlorofluoromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Indomethane ND	1,1-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Carbon Disulfide ND 0.20 EPA 8260C 6-25-18 6-25-18 Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 Vinyl Acetate ND 0.20 EPA 8260C 6-25-18 6-25-18 2,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Gebata ND 0.20 EPA 8260C 6-25-18 </td <td>Acetone</td> <td>ND</td> <td>5.0</td> <td>EPA 8260C</td> <td>6-25-18</td> <td>6-25-18</td> <td></td>	Acetone	ND	5.0	EPA 8260C	6-25-18	6-25-18	
Methylene Chloride ND 1.0 EPA 8260C 6-25-18 6-25-18 (trans) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Vinyl Acetate ND 1.0 EPA 8260C 6-25-18 6-25-18 2,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 (cis) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromochloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6	Iodomethane	ND	1.9	EPA 8260C	6-25-18	6-25-18	
(trans) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Vinyl Acetate ND 1.0 EPA 8260C 6-25-18 6-25-18 2,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 (cis) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 5.0 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromochloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 </td <td>Carbon Disulfide</td> <td>ND</td> <td>0.20</td> <td>EPA 8260C</td> <td>6-25-18</td> <td>6-25-18</td> <td></td>	Carbon Disulfide	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methyl t-Butyl Ether ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Vinyl Acetate ND 1.0 EPA 8260C 6-25-18 6-25-18 2,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 (cis) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 5.0 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromochloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chlorofrm ND 0.20 EPA 8260C 6-25-18 6-25-18 Chlorofrom ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18	Methylene Chloride	ND	1.0	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Vinyl Acetate ND 1.0 EPA 8260C 6-25-18 6-25-18 2,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 (cis) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 5.0 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromochloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Trichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 <td>(trans) 1,2-Dichloroethene</td> <td>ND</td> <td>0.20</td> <td>EPA 8260C</td> <td>6-25-18</td> <td>6-25-18</td> <td></td>	(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Vinyl Acetate ND 1.0 EPA 8260C 6-25-18 6-25-18 2,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 (cis) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 5.0 EPA 8260C 6-25-18 6-25-18 Bromochloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Benzene ND 0.20 EPA 8260C 6-25-18 6-25-18	Methyl t-Butyl Ether	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 (cis) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 5.0 EPA 8260C 6-25-18 6-25-18 Bromochloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1,1-Trichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Benzene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Trichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA 8260C 6-25	1,1-Dichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
(cis) 1,2-Dichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Butanone ND 5.0 EPA 8260C 6-25-18 6-25-18 Bromochloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1,1-Trichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C	Vinyl Acetate	ND	1.0	EPA 8260C	6-25-18	6-25-18	
2-Butanone ND 5.0 EPA 8260C 6-25-18 6-25-18 Bromochloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1,1-Trichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Benzene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Tichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA 8260C 6-25-18 6-25	2,2-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromochloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1,1-Trichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Benzene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Trichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromodichloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18	(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chloroform ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1,1-Trichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Benzene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Trichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 </td <td>2-Butanone</td> <td>ND</td> <td>5.0</td> <td>EPA 8260C</td> <td>6-25-18</td> <td>6-25-18</td> <td></td>	2-Butanone	ND	5.0	EPA 8260C	6-25-18	6-25-18	
1,1,1-Trichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Benzene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Trichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromodichloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 1.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA	Bromochloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Carbon Tetrachloride ND 0.20 EPA 8260C 6-25-18 6-25-18 1,1-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Benzene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Trichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromodichloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18	Chloroform	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Benzene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Trichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromodichloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	1,1,1-Trichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Benzene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Trichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromodichloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	Carbon Tetrachloride	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichloroethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Trichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromodichloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	1,1-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Trichloroethene ND 0.20 EPA 8260C 6-25-18 6-25-18 1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromodichloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	Benzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichloropropane ND 0.20 EPA 8260C 6-25-18 6-25-18 Dibromomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromodichloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	1,2-Dichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Dibromomethane ND 0.20 EPA 8260C 6-25-18 6-25-18 Bromodichloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	Trichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromodichloromethane ND 0.20 EPA 8260C 6-25-18 6-25-18 2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	1,2-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Chloroethyl Vinyl Ether ND 1.0 EPA 8260C 6-25-18 6-25-18 (cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	Dibromomethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
(cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	Bromodichloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
(cis) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18 Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Methyl Isobutyl Ketone ND 2.0 EPA 8260C 6-25-18 6-25-18 Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Toluene ND 1.0 EPA 8260C 6-25-18 6-25-18	Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	6-25-18	6-25-18	
(trans) 1,3-Dichloropropene ND 0.20 EPA 8260C 6-25-18 6-25-18	Toluene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
	(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	

Project: 17-06520-000

VOLATILES EPA 8260C page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	Trip Blank					
Laboratory ID:	06-222-10					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Tetrachloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Hexanone	ND	2.0	EPA 8260C	6-25-18	6-25-18	
Dibromochloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Ethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
m,p-Xylene	ND	0.40	EPA 8260C	6-25-18	6-25-18	
o-Xylene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Styrene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromoform	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Isopropylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2,3-Trichloropropane	ND	0.27	EPA 8260C	6-25-18	6-25-18	
n-Propylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
n-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Naphthalene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	100	75-127				
Toluene-d8	102	80-127				

4-Bromofluorobenzene

78-125

98

Project: 17-06520-000

VOLATILES by EPA 8260C METHOD BLANK QUALITY CONTROL

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0625W1					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chloromethane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Vinyl Chloride	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromomethane	ND	0.31	EPA 8260C	6-25-18	6-25-18	
Chloroethane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Acetone	ND	5.0	EPA 8260C	6-25-18	6-25-18	
lodomethane	ND	1.9	EPA 8260C	6-25-18	6-25-18	
Carbon Disulfide	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methylene Chloride	ND	1.0	EPA 8260C	6-25-18	6-25-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Vinyl Acetate	ND	1.0	EPA 8260C	6-25-18	6-25-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Butanone	ND	5.0	EPA 8260C	6-25-18	6-25-18	
Bromochloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chloroform	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Benzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Trichloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Dibromomethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromodichloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	6-25-18	6-25-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	6-25-18	6-25-18	
Toluene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	6-25-18	6-25-18	

Project: 17-06520-000

VOLATILES by EPA 8260C METHOD BLANK QUALITY CONTROL

page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Analyte	Result	1 &L	Metrica	Trepared	Anaryzea	i iugo
Laboratory ID:	MB0625W1					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Tetrachloroethene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Hexanone	ND	2.0	EPA 8260C	6-25-18	6-25-18	
Dibromochloromethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Chlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Ethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
m,p-Xylene	ND	0.40	EPA 8260C	6-25-18	6-25-18	
o-Xylene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Styrene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromoform	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Isopropylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Bromobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2,3-Trichloropropane	ND	0.27	EPA 8260C	6-25-18	6-25-18	
n-Propylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
n-Butylbenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	6-25-18	6-25-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
Naphthalene	ND	1.0	EPA 8260C	6-25-18	6-25-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	6-25-18	6-25-18	
Surrogate:	Percent Recovery	Control Limits				

Surrogate: Percent Recovery Control Limit
Dibromofluoromethane 100 75-127
Toluene-d8 102 80-127
4-Bromofluorobenzene 102 78-125



Project: 17-06520-000

VOLATILES by EPA 8260C SB/SBD QUALITY CONTROL

					Per	cent	Recovery		RPD	
Analyte	Result		Spike	Spike Level		overy	Limits	RPD	Limit	Flags
SPIKE BLANKS										
Laboratory ID:	SB06	25W1								
	SB	SBD	SB	SBD	SB	SBD				
1,1-Dichloroethene	9.02	8.85	10.0	10.0	90	89	62-129	2	15	
Benzene	9.49	9.38	10.0	10.0	95	94	77-127	1	15	
Trichloroethene	9.80	9.61	10.0	10.0	98	96	70-120	2	15	
Toluene	10.2	10.1	10.0	10.0	102	101	82-123	1	15	
Chlorobenzene	9.74	9.58	10.0	10.0	97	96	79-120	2	15	
Surrogate:										
Dibromofluoromethane					98	99	75-127			
Toluene-d8					104	104	80-127			
4-Bromofluorobenzene					103	102	78-125			

Project: 17-06520-000

TOTAL METALS EPA 200.8/7470A

0 (11)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW1					
Laboratory ID:	06-222-01					
Arsenic	ND	3.3	EPA 200.8	6-27-18	6-27-18	
Cadmium	ND	4.4	EPA 200.8	6-27-18	6-27-18	
Chromium	ND	11	EPA 200.8	6-27-18	6-27-18	
Lead	ND	1.1	EPA 200.8	6-27-18	6-27-18	
Mercury	ND	0.50	EPA 7470A	6-29-18	6-29-18	
Client ID:	MW2					
Laboratory ID:	06-222-02					
Arsenic	ND	3.3	EPA 200.8	6-27-18	6-27-18	
Cadmium	ND	4.4	EPA 200.8	6-27-18	6-27-18	
Chromium	ND	11	EPA 200.8	6-27-18	6-27-18	
Lead	ND	1.1	EPA 200.8	6-27-18	6-27-18	
Mercury	ND	0.50	EPA 7470A	6-29-18	6-29-18	
	.,_	0.00		0 20 .0	0 20 .0	
Client ID:	MW3					
Laboratory ID:	06-222-03					
Arsenic	ND	3.3	EPA 200.8	6-27-18	6-27-18	
Cadmium	ND	4.4	EPA 200.8	6-27-18	6-27-18	
Chromium	ND	11	EPA 200.8	6-27-18	6-27-18	
Lead	ND	1.1	EPA 200.8	6-27-18	6-27-18	
Mercury	ND	0.50	EPA 7470A	6-29-18	6-29-18	
Client ID:	MW4					
Laboratory ID:	06-222-04					
Arsenic	11	3.3	EPA 200.8	6-27-18	6-27-18	
Cadmium	ND	4.4	EPA 200.8	6-27-18	6-27-18	
Chromium	ND	11	EPA 200.8	6-27-18	6-27-18	
Lead	ND	1.1	EPA 200.8	6-27-18	6-27-18	
Mercury	ND	0.50	EPA 7470A	6-29-18	6-29-18	

Project: 17-06520-000

TOTAL METALS EPA 200.8/7470A

Offits. ug/L (ppb)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW5			•	-	_
Laboratory ID:	06-222-05					
Arsenic	ND	3.3	EPA 200.8	6-27-18	6-27-18	
Cadmium	ND	4.4	EPA 200.8	6-27-18	6-27-18	
Chromium	ND	11	EPA 200.8	6-27-18	6-27-18	
Lead	5.6	1.1	EPA 200.8	6-27-18	6-27-18	
Mercury	ND	0.50	EPA 7470A	6-29-18	6-29-18	
Client ID:	MW6					
Laboratory ID:	06-222-06					
Arsenic	ND	3.3	EPA 200.8	6-27-18	6-27-18	
Cadmium	ND	4.4	EPA 200.8	6-27-18	6-27-18	
Chromium	ND	11	EPA 200.8	6-27-18	6-27-18	
Lead	ND	1.1	EPA 200.8	6-27-18	6-27-18	
Mercury	ND	0.50	EPA 7470A	6-29-18	6-29-18	
Client ID:	MW7					
Laboratory ID:	06-222-07					
Arsenic	4.6	3.3	EPA 200.8	6-27-18	6-27-18	
Cadmium	ND	4.4	EPA 200.8	6-27-18	6-27-18	
Chromium	ND	11	EPA 200.8	6-27-18	6-27-18	
Lead	2.0	1.1	EPA 200.8	6-27-18	6-27-18	
Mercury	ND	0.50	EPA 7470A	6-29-18	6-29-18	
011	1414/O					
Client ID:	MW8					
Laboratory ID:	06-222-08		ED4 222 2	0.07.10	0.0= 10	
Arsenic	3.9	3.3	EPA 200.8	6-27-18	6-27-18	
Cadmium	ND	4.4	EPA 200.8	6-27-18	6-27-18	
Chromium	ND	11	EPA 200.8	6-27-18	6-27-18	
Lead	ND	1.1	EPA 200.8	6-27-18	6-27-18	
Mercury	ND	0.50	EPA 7470A	6-29-18	6-29-18	

Project: 17-06520-000

TOTAL METALS EPA 200.8/7470A

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW9					
Laboratory ID:	06-222-09					
Arsenic	ND	3.3	EPA 200.8	6-27-18	6-27-18	
Cadmium	ND	4.4	EPA 200.8	6-27-18	6-27-18	
Chromium	ND	11	EPA 200.8	6-27-18	6-27-18	
Lead	ND	1.1	EPA 200.8	6-27-18	6-27-18	
Mercury	ND	0.50	EPA 7470A	6-29-18	6-29-18	

Project: 17-06520-000

TOTAL METALS EPA 200.8/7470A QUALITY CONTROL

Matrix: Water
Units: ug/L (ppb)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0627WM1					
Arsenic	ND	3.3	EPA 200.8	6-27-18	6-27-18	
Cadmium	ND	4.4	EPA 200.8	6-27-18	6-27-18	
Chromium	ND	11	EPA 200.8	6-27-18	6-27-18	
Lead	ND	1.1	EPA 200.8	6-27-18	6-27-18	
Laboratory ID:	MB0629W1					
Mercury	ND	0.50	EPA 7470A	6-29-18	6-29-18	

					Source	Pe	rcent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Rec	overy	Limits	RPD	Limit	Flags
DUPLICATE											,
Laboratory ID:	06-10	00-09									
	ORIG	DUP									
Arsenic	ND	ND	NA	NA		ı	NA	NA	NA	20	
Cadmium	ND	ND	NA	NA		ı	NA	NA	NA	20	
Chromium	ND	ND	NA	NA		ı	NA	NA	NA	20	
Lead	ND	ND	NA	NA		l	NA	NA	NA	20	
Laboratory ID:	06-29	95-01									
Mercury	ND	ND	NA	NA			NA	NA	NA	20	
MATRIX SPIKES											
Laboratory ID:	06-10	00-09									
,	MS	MSD	MS	MSD		MS	MSD				
Arsenic	219	220	222	222	ND	99	99	75-125	0	20	
Cadmium	222	221	222	222	ND	100	99	75-125	1	20	
Chromium	202	208	222	222	ND	91	94	75-125	3	20	
Lead	231	237	222	222	ND	104	107	75-125	3	20	
Laboratory ID:	06-29	95-01									
Mercury	12.4	12.4	12.5	12.5	ND	99	99	75-125	0	20	

Project: 17-06520-000

cPAHs EPA 8270D/SIM

			Date	Date	
Result	PQL	Method	Prepared	Analyzed	Flags
MW1					
06-222-01					
ND	0.0095	EPA 8270D/SIM	6-25-18	6-26-18	
ND	0.0095	EPA 8270D/SIM	6-25-18	6-26-18	
ND	0.0095	EPA 8270D/SIM	6-25-18	6-26-18	
ND	0.0095	EPA 8270D/SIM	6-25-18	6-26-18	
ND	0.0095	EPA 8270D/SIM	6-25-18	6-26-18	
ND	0.0095	EPA 8270D/SIM	6-25-18	6-26-18	
ND	0.0095	EPA 8270D/SIM	6-25-18	6-26-18	
Percent Recovery	Control Limits				
58	21 - 110				
72	19 - 111				
74	32 - 137				
	MW1 06-222-01 ND ND ND ND ND ND ND ND Percent Recovery 58 72	MW1 06-222-01 ND 0.0095 Percent Recovery Control Limits 58 21 - 110 72 19 - 111	MW1 06-222-01 0.0095 EPA 8270D/SIM ND 0.0095 EPA 8270D/SIM Percent Recovery Control Limits 58 21 - 110 72 19 - 111	Result PQL Method Prepared MW1 06-222-01 6-25-18 ND 0.0095 EPA 8270D/SIM 6-25-18 Percent Recovery Control Limits 58 21 - 110 72 19 - 111	Result PQL Method Prepared Analyzed MW1 06-222-01 6-25-18 6-26-18 ND 0.0095 EPA 8270D/SIM 6-25-18 6-26-18 Percent Recovery Control Limits 58 21 - 110 21 - 110 21 - 110 72 19 - 111 111 111

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW2					
Laboratory ID:	06-222-02					
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Chrysene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	57	21 - 110				
Pyrene-d10	73	19 - 111				
Terphenyl-d14	73	32 - 137				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

·				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW3					
Laboratory ID:	06-222-03					
Benzo[a]anthracene	ND	0.0097	EPA 8270D/SIM	6-25-18	6-26-18	
Chrysene	ND	0.0097	EPA 8270D/SIM	6-25-18	6-26-18	
Benzo[b]fluoranthene	ND	0.0097	EPA 8270D/SIM	6-25-18	6-26-18	
Benzo(j,k)fluoranthene	ND	0.0097	EPA 8270D/SIM	6-25-18	6-26-18	
Benzo[a]pyrene	ND	0.0097	EPA 8270D/SIM	6-25-18	6-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0097	EPA 8270D/SIM	6-25-18	6-26-18	
Dibenz[a,h]anthracene	ND	0.0097	EPA 8270D/SIM	6-25-18	6-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	60	21 - 110				
Pyrene-d10	82	19 - 111				
Terphenyl-d14	80	32 - 137				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW4					
Laboratory ID:	06-222-04					
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Chrysene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	51	21 - 110				
Pyrene-d10	66	19 - 111				
Terphenyl-d14	66	32 - 137				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

			Date	Date	
Result	PQL	Method	Prepared	Analyzed	Flags
MW5					
06-222-05					
ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Percent Recovery	Control Limits				
71	21 - 110				
69	19 - 111				
95	32 - 137				
	MW5 06-222-05 ND ND ND ND ND ND ND ND Percent Recovery 71 69	MW5 06-222-05 ND 0.010 ND 0.010 ND 0.010 ND 0.010 ND 0.010 ND 0.010 Percent Recovery Control Limits 71 21 - 110 69 19 - 111	MW5 06-222-05 0.010 EPA 8270D/SIM ND 0.010 EPA 8270D/SIM Percent Recovery Control Limits 71 21 - 110 69 19 - 111	Result PQL Method Prepared MW5 06-222-05	Result PQL Method Prepared Analyzed MW5 06-222-05 6-25-18 6-26-18 ND 0.010 EPA 8270D/SIM 6-25-18 6-26-18 Percent Recovery Control Limits 71 21 - 110 69 19 - 111 11

Project: 17-06520-000

cPAHs EPA 8270D/SIM

			Date	Date	
Result	PQL	Method	Prepared	Analyzed	Flags
MW6					
06-222-06					
0.014	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
0.014	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
0.012	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
0.012	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
0.011	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
0.012	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
0.011	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Percent Recovery	Control Limits				
60	21 - 110				
72	19 - 111				
71	32 - 137				
	MW6 06-222-06 0.014 0.014 0.012 0.012 0.011 0.012 0.011 Percent Recovery 60 72	MW6 06-222-06 0.014 0.010 0.012 0.010 0.012 0.010 0.011 0.010 0.012 0.010 0.011 0.010 Percent Recovery Control Limits 60 21 - 110 72 19 - 111	MW6 06-222-06 0.014 0.010 EPA 8270D/SIM 0.014 0.010 EPA 8270D/SIM 0.012 0.010 EPA 8270D/SIM 0.012 0.010 EPA 8270D/SIM 0.011 0.010 EPA 8270D/SIM 0.012 0.010 EPA 8270D/SIM 0.011 0.010 EPA 8270D/SIM Percent Recovery Control Limits 60 21 - 110 72 19 - 111	Result PQL Method Prepared MW6 06-222-06	Result PQL Method Prepared Analyzed MW6 06-222-06 6-25-18 6-26-18 0.014 0.010 EPA 8270D/SIM 6-25-18 6-26-18 0.014 0.010 EPA 8270D/SIM 6-25-18 6-26-18 0.012 0.010 EPA 8270D/SIM 6-25-18 6-26-18 0.011 0.010 EPA 8270D/SIM 6-25-18 6-26-18 0.012 0.010 EPA 8270D/SIM 6-25-18 6-26-18 0.012 0.010 EPA 8270D/SIM 6-25-18 6-26-18 0.011 0.010 EPA 8270D/SIM 6-25-18 6-26-18 Percent Recovery Control Limits 6-25-18 6-26-18 60 21 - 110 21 - 110 21 - 110 72 19 - 111 111 111

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW7					
Laboratory ID:	06-222-07					
Benzo[a]anthracene	ND	0.0096	EPA 8270D/SIM	6-25-18	6-26-18	
Chrysene	ND	0.0096	EPA 8270D/SIM	6-25-18	6-26-18	
Benzo[b]fluoranthene	ND	0.0096	EPA 8270D/SIM	6-25-18	6-26-18	
Benzo(j,k)fluoranthene	ND	0.0096	EPA 8270D/SIM	6-25-18	6-26-18	
Benzo[a]pyrene	ND	0.0096	EPA 8270D/SIM	6-25-18	6-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0096	EPA 8270D/SIM	6-25-18	6-26-18	
Dibenz[a,h]anthracene	ND	0.0096	EPA 8270D/SIM	6-25-18	6-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	57	21 - 110				
Pyrene-d10	68	19 - 111				
Terphenyl-d14	68	32 - 137				
l erphenyl-d14	68	32 - 137				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW8					
Laboratory ID:	06-222-08					
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Chrysene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	55	21 - 110				
Pyrene-d10	71	19 - 111				
Terphenyl-d14	70	32 - 137				
i erpnenyi-a14	70	32 - 137				

Project: 17-06520-000

cPAHs EPA 8270D/SIM

			Date	Date	
Result	PQL	Method	Prepared	Analyzed	Flags
MW9					
06-222-09					
ND	0.0098	EPA 8270D/SIM	6-25-18	6-26-18	
ND	0.0098	EPA 8270D/SIM	6-25-18	6-26-18	
ND	0.0098	EPA 8270D/SIM	6-25-18	6-26-18	
ND	0.0098	EPA 8270D/SIM	6-25-18	6-26-18	
ND	0.0098	EPA 8270D/SIM	6-25-18	6-26-18	
ND	0.0098	EPA 8270D/SIM	6-25-18	6-26-18	
ND	0.0098	EPA 8270D/SIM	6-25-18	6-26-18	
Percent Recovery	Control Limits				
50	21 - 110				
69	19 - 111				
68	32 - 137				
	MW9 06-222-09 ND ND ND ND ND ND ND ND Percent Recovery 50 69	MW9 06-222-09 ND 0.0098 Percent Recovery Control Limits 50 21 - 110 69 19 - 111	MW9 06-222-09 ND 0.0098 EPA 8270D/SIM Percent Recovery Control Limits 50 21 - 110 69 19 - 111	Result PQL Method Prepared MW9 06-222-09 6-25-18 ND 0.0098 EPA 8270D/SIM 6-25-18 Percent Recovery Control Limits 50 21 - 110 69 19 - 111	Result PQL Method Prepared Analyzed MW9 06-222-09 6-25-18 6-26-18 ND 0.0098 EPA 8270D/SIM 6-25-18 6-26-18 Percent Recovery Control Limits 50 21 - 110 69 19 - 111

Project: 17-06520-000

CPAHS EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0625W1					
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Chrysene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	6-25-18	6-26-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	64	21 - 110				
Pyrene-d10	84	19 - 111				
Terphenyl-d14	83	32 - 137				

Project: 17-06520-000

cPAHs EPA 8270D/SIM SB/SBD QUALITY CONTROL

					Per	cent	Recovery		RPD	
Analyte	Re	sult	Spike	Level	Rec	overy	Limits	RPD	Limit	Flags
SPIKE BLANKS										
Laboratory ID:	SB06	25W1								
	SB	SBD	SB	SBD	SB	SBD				
Benzo[a]anthracene	0.395	0.425	0.500	0.500	79	85	57 - 127	7	15	
Chrysene	0.354	0.371	0.500	0.500	71	74	51 - 120	5	15	
Benzo[b]fluoranthene	0.398	0.422	0.500	0.500	80	84	54 - 124	6	17	
Benzo(j,k)fluoranthene	0.369	0.392	0.500	0.500	74	78	50 - 127	6	18	
Benzo[a]pyrene	0.398	0.436	0.500	0.500	80	87	50 - 120	9	16	
Indeno(1,2,3-c,d)pyrene	0.417	0.443	0.500	0.500	83	89	46 - 132	6	20	
Dibenz[a,h]anthracene	0.404	0.421	0.500	0.500	81	84	49 - 129	4	18	
Surrogate:										
2-Fluorobiphenyl					57	53	21 - 110			
Pyrene-d10					73	77	19 - 111			
Terphenyl-d14					70	76	32 - 137			



Data Qualifiers and Abbreviations

- A Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
- B The analyte indicated was also found in the blank sample.
- C The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
- E The value reported exceeds the quantitation range and is an estimate.
- F Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
- H The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
- I Compound recovery is outside of the control limits.
- J The value reported was below the practical quantitation limit. The value is an estimate.
- K Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
- L The RPD is outside of the control limits.
- M Hydrocarbons in the gasoline range are impacting the diesel range result.
- M1 Hydrocarbons in the gasoline range (toluene-naphthalene) are present in the sample.
- N Hydrocarbons in the lube oil range are impacting the diesel range result.
- N1 Hydrocarbons in diesel range are impacting lube oil range results.
- O Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
- P The RPD of the detected concentrations between the two columns is greater than 40.
- Q Surrogate recovery is outside of the control limits.
- S Surrogate recovery data is not available due to the necessary dilution of the sample.
- T The sample chromatogram is not similar to a typical
- U The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- U1 The practical quantitation limit is elevated due to interferences present in the sample.
- V Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
- W Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
- X Sample extract treated with a mercury cleanup procedure.
- X1- Sample extract treated with a sulfuric acid/silica gel cleanup procedure.
- Y The calibration verification for this analyte exceeded the 20% drift specified in method 8260C, and therefore the reported result should be considered an estimate. The overall performance of the calibration verification standard met the acceptance criteria of the method.

7 -

ND - Not Detected at PQL

PQL - Practical Quantitation Limit

RPD - Relative Percent Difference





Chain of Custody

Page _____ of ____

Analytical Laboratory Testing Services 14648 NE 95th Street • Redmond, WA 98052		naround Req n working da			L	abo	rate	ory	Nun	nbe	r:	0	<u>6</u> ·	- 2	2	2									
Phone: (425) 883-3881 • www.onsite-env.com		(Check One)														Σ									
Project Number:	Same	Day [1 Day					(0								8270D/SIM	4					M			
17-06520-600	2 Day	/s [3 Days					Clean-up)							081B		8151A					15			
Project Name: Paufic Park Project Manager:		dard (7 Days) analysis 5 Da	ave)	50				SG	0000	Duozo	s Only	MIS	-level)		sides 8	esticid	oicides				grease) 1664A	by 8270 D			
Project Manager: May K & W bank	(IPA	analysis 5 Da	195)	tainer	ķ.	×		Acid /	lite!	Mater	Water	270D/S	M (low		Pestic	orus P	d Hert	tals	tals		ease)	82			
Sampled by:		(other)		of Con	HCID	Gx/BTI	×	□) ×(8260C	med vo	8011	tiles 8; -level l	1S/Q0/	82A	hlorine	hosph	ed Aci	RA Me	CA Me	etals	and gi	by	-		Ire
Sampled by: George Ifther	Date	Time		mber	/TPH-I	NWTPH-Gx/BTEX	NWTPH-Gx	NWTPH-Dx	Volatiles 8260C	Halogenated Volatiles ozooc	EDB EPA 8011 (Waters Only,	Semivolatiles 8270D/SIM (with low-level PAHs)	Hs 82	PCBs 8082A	Organochlorine Pesticides 8081B	Organophosphorus Pesticides	Chlorinated Acid Herbicides	Total RCRA Metals	Total MTCA Metals	TCLP Metals	HEM (oil and	CPAHS			% Moisture
Lab ID Sample Identification	Sampled	Sampled	Matrix	10	Ž	ž	Ž	Ž	9	E G		§ §	PA	PC	ŏ	ŏ	ਠ	oT	P /	DT	포		+	+	%
· MWI	6/21/18	1251	water	世			/	V	V										V			V			
2 MW2	1	13:25	1	111			V	V	V										/			V			
3 MW3		1152					V	V	V										V			/			
4 mn4		12:25	1				V	1	V										V			1			
S MW5		1355					V	V	V										1			1			
v MW6		1445					V	1	V										V			V			
7 MW7.		1030					V	1	V										V			V			
3 MW8		1535					V	1	0										1			V			
9 mw9		1430		V			V	/	/										V	fl.o.		V			
10 Trip Blank	A		V	3			V		1									200	E	AC.	-	1			
Signature /	C	ompany				Date	е		Time			Con	nmen	ts/Sp	ecial	Instr	uction	ns					**********		1
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July 10, 2018

Mark Ewbank Herrera Environmental Consultants, Inc. 2200 6th Avenue, Suite 1100 Seattle, WA 98121

Re: Analytical Data for Project 17-06520-000

Laboratory Reference No. 1806-310

Dear Mark:

Enclosed are the analytical results and associated quality control data for samples submitted on June 29, 2018.

The standard policy of OnSite Environmental, Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

David Baumeister Project Manager

Enclosures



Project: 17-06520-000

Case Narrative

Samples were collected on June 29, 2018 and received by the laboratory on June 29, 2018. They were maintained at the laboratory at a temperature of 2°C to 6°C.

Please note that any and all soil sample results are reported on a dry-weight basis, unless otherwise noted below.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

NWTPH-Gx

Matrix: Water Units: ug/L (ppb)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	SW1					
Laboratory ID:	06-310-01					
Gasoline	ND	100	NWTPH-Gx	7-3-18	7-3-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	75	66-117				
Client ID:	SW2					
Laboratory ID:	06-310-02					
Gasoline	ND	100	NWTPH-Gx	7-3-18	7-3-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	76	66-117				
Client ID:	SW3					
Laboratory ID:	06-310-03					
Gasoline	ND	100	NWTPH-Gx	7-3-18	7-3-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	76	66-117				

Project: 17-06520-000

NWTPH-Gx QUALITY CONTROL

Matrix: Water Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
METHOD BLANK				•	•	
Laboratory ID:	MB0703W3					
Gasoline	ND	100	NWTPH-Gx	7-3-18	7-3-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	74	66-117				

Analyte	Res	sult	Spike	Level	Source Result	Pero Reco		Recovery Limits	RPD	RPD Limit	Flags
DUPLICATE											
Laboratory ID:	07-00	05-01									
	ORIG	DUP									
Gasoline	ND	ND	NA	NA		N	A	NA	NA	30	
Surrogate:											
Fluorobenzene						76	79	66-117			

NWTPH-Dx

Matrix: Water Units: mg/L (ppm)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	SW1					
Laboratory ID:	06-310-01					
Diesel Range Organics	ND	0.25	NWTPH-Dx	7-2-18	7-2-18	
Lube Oil Range Organics	ND	0.41	NWTPH-Dx	7-2-18	7-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	103	50-150				
Client ID:	SW2					
Laboratory ID:	06-310-02					
Diesel Range Organics	ND	0.26	NWTPH-Dx	7-2-18	7-2-18	
Lube Oil Range Organics	ND	0.42	NWTPH-Dx	7-2-18	7-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	95	50-150				
Client ID:	SW3					
Laboratory ID:	06-310-03					
Diesel Range Organics	ND	0.26	NWTPH-Dx	7-2-18	7-2-18	
Lube Oil Range Organics	ND	0.42	NWTPH-Dx	7-2-18	7-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	94	50-150				

NWTPH-Dx QUALITY CONTROL

Matrix: Water Units: mg/L (ppm)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0702W1					
Diesel Range Organics	ND	0.25	NWTPH-Dx	7-2-18	7-2-18	_
Lube Oil Range Organics	ND	0.40	NWTPH-Dx	7-2-18	7-2-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	87	50-150				

					Source	Per	cent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Reco	overy	Limits	RPD	Limit	Flags
DUPLICATE											
Laboratory ID:	06-30	03-01									
	ORIG	DUP									
Diesel Range	ND	ND	NA	NA		Ν	lΑ	NA	NA	NA	
Lube Oil Range	ND	ND	NA	NA		N	IA	NA	NA	NA	
Surrogate:											
o-Terphenyl						99	103	50-150			

VOLATILES EPA 8260C page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	SW1					
Laboratory ID:	06-310-01					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Chloromethane	ND	1.0	EPA 8260C	7-3-18	7-3-18	
Vinyl Chloride	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Bromomethane	ND	2.0	EPA 8260C	7-3-18	7-3-18	
Chloroethane	ND	1.0	EPA 8260C	7-3-18	7-3-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Acetone	ND	5.0	EPA 8260C	7-3-18	7-3-18	
Iodomethane	ND	5.0	EPA 8260C	7-3-18	7-3-18	
Carbon Disulfide	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Methylene Chloride	ND	2.0	EPA 8260C	7-3-18	7-3-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Vinyl Acetate	ND	1.0	EPA 8260C	7-3-18	7-3-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
2-Butanone	ND	5.0	EPA 8260C	7-3-18	7-3-18	
Bromochloromethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Chloroform	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Benzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Trichloroethene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Dibromomethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Bromodichloromethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	7-3-18	7-3-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	7-3-18	7-3-18	
Toluene	ND	1.0	EPA 8260C	7-3-18	7-3-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	7-3-18	7-3-18	

Project: 17-06520-000

VOLATILES EPA 8260C

page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	SW1	1 42	Wethou	rrepared	Analyzeu	i iags
Laboratory ID:	06-310-01					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Tetrachloroethene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
2-Hexanone	ND	2.0	EPA 8260C	7-3-18	7-3-18	
Dibromochloromethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Chlorobenzene	ND ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Ethylbenzene	ND ND	0.20	EPA 8260C	7-3-18 7-3-18	7-3-18 7-3-18	
m,p-Xylene	ND ND	0.40	EPA 8260C	7-3-18 7-3-18	7-3-18 7-3-18	
o-Xylene	ND ND	0.40	EPA 8260C	7-3-16 7-3-18	7-3-18 7-3-18	
Styrene	ND ND	0.20	EPA 8260C	7-3-16 7-3-18	7-3-18 7-3-18	
Bromoform	ND ND	1.0	EPA 8260C	7-3-16 7-3-18	7-3-16 7-3-18	
Isopropylbenzene	ND ND	0.20	EPA 8260C	7-3-16 7-3-18	7-3-16 7-3-18	
	ND ND	0.20		7-3-16 7-3-18	7-3-16 7-3-18	
Bromobenzene			EPA 8260C			
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
n-Propylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
n-Butylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2-Dibromo-3-chloropropane		1.0	EPA 8260C	7-3-18	7-3-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	7-3-18	7-3-18	
Naphthalene	ND	1.0	EPA 8260C	7-3-18	7-3-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Surrogate:	Percent Recovery	Control Limits				

Surrogate: Percent Recovery Control Limit Dibromofluoromethane 99 75-127
Toluene-d8 94 80-127
4-Bromofluorobenzene 95 78-125



Project: 17-06520-000

VOLATILES EPA 8260C page 1 of 2

omis. ug/L				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	SW2					
Laboratory ID:	06-310-02					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Chloromethane	ND	1.0	EPA 8260C	7-3-18	7-3-18	
Vinyl Chloride	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Bromomethane	ND	2.0	EPA 8260C	7-3-18	7-3-18	
Chloroethane	ND	1.0	EPA 8260C	7-3-18	7-3-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Acetone	ND	5.0	EPA 8260C	7-3-18	7-3-18	
Iodomethane	ND	5.0	EPA 8260C	7-3-18	7-3-18	
Carbon Disulfide	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Methylene Chloride	ND	2.0	EPA 8260C	7-3-18	7-3-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Vinyl Acetate	ND	1.0	EPA 8260C	7-3-18	7-3-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
2-Butanone	ND	5.0	EPA 8260C	7-3-18	7-3-18	
Bromochloromethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Chloroform	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Benzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Trichloroethene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Dibromomethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Bromodichloromethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	7-3-18	7-3-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	7-3-18	7-3-18	
Toluene	ND	1.0	EPA 8260C	7-3-18	7-3-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	7-3-18	7-3-18	

VOLATILES EPA 8260C

page 2 of 2

Analyta	Result	PQL	Method	Date Prepared	Date Analyzed	Flogo
Analyte Client ID:	SW2	PQL	Wethou	Prepared	Anaryzeu	Flags
Laboratory ID:	06-310-02					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Tetrachloroethene	ND ND	0.20	EPA 8260C	7-3-16 7-3-18	7-3-18 7-3-18	
	ND	0.20	EPA 8260C	7-3-18 7-3-18	7-3-18 7-3-18	
1,3-Dichloropropane 2-Hexanone	ND ND	2.0	EPA 8260C EPA 8260C	7-3-16 7-3-18	7-3-16 7-3-18	
Z-nexanone Dibromochloromethane	ND ND	0.20		7-3-16 7-3-18	7-3-16 7-3-18	
	ND ND	0.20	EPA 8260C EPA 8260C	7-3-16 7-3-18	7-3-16 7-3-18	
1,2-Dibromoethane						
Chlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Ethylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
m,p-Xylene	ND	0.40	EPA 8260C	7-3-18	7-3-18	
o-Xylene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Styrene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Bromoform	ND	1.0	EPA 8260C	7-3-18	7-3-18	
Isopropylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Bromobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
n-Propylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
n-Butylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	7-3-18	7-3-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	7-3-18	7-3-18	
Naphthalene	ND	1.0	EPA 8260C	7-3-18	7-3-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Surrogate:	Percent Recovery					

Surrogate:	Percent Recovery	Control Limits
Dibromofluoromethane	101	75-127
Toluene-d8	95	80-127
4-Bromofluorobenzene	95	78-125



Project: 17-06520-000

VOLATILES EPA 8260C

page 1 of 2

omis. ug/L				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	SW3					
Laboratory ID:	06-310-03					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Chloromethane	ND	1.0	EPA 8260C	7-3-18	7-3-18	
Vinyl Chloride	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Bromomethane	ND	2.0	EPA 8260C	7-3-18	7-3-18	
Chloroethane	ND	1.0	EPA 8260C	7-3-18	7-3-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Acetone	ND	5.0	EPA 8260C	7-3-18	7-3-18	
Iodomethane	ND	5.0	EPA 8260C	7-3-18	7-3-18	
Carbon Disulfide	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Methylene Chloride	ND	2.0	EPA 8260C	7-3-18	7-3-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Vinyl Acetate	ND	1.0	EPA 8260C	7-3-18	7-3-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
2-Butanone	ND	5.0	EPA 8260C	7-3-18	7-3-18	
Bromochloromethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Chloroform	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Benzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Trichloroethene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Dibromomethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Bromodichloromethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	7-3-18	7-3-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	7-3-18	7-3-18	
Toluene	ND	1.0	EPA 8260C	7-3-18	7-3-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	7-3-18	7-3-18	

Project: 17-06520-000

VOLATILES EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	SW3					
Laboratory ID:	06-310-03					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Tetrachloroethene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
2-Hexanone	ND	2.0	EPA 8260C	7-3-18	7-3-18	
Dibromochloromethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Chlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Ethylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
m,p-Xylene	ND	0.40	EPA 8260C	7-3-18	7-3-18	
o-Xylene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Styrene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Bromoform	ND	1.0	EPA 8260C	7-3-18	7-3-18	
Isopropylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Bromobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
n-Propylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
n-Butylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	7-3-18	7-3-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	7-3-18	7-3-18	
Naphthalene	ND	1.0	EPA 8260C	7-3-18	7-3-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Surrogate:	Percent Recovery	Control Limits				

Surrogate: Percent Recovery Control Limits
Dibromofluoromethane 98 75-127
Toluene-d8 94 80-127
4-Bromofluorobenzene 94 78-125



Project: 17-06520-000

VOLATILES EPA 8260C page 1 of 2

			Date	Date	
Result	PQL	Method	Prepared	Analyzed	Flags
Trip Blank					
06-310-04					
ND	0.20	EPA 8260C	7-3-18	7-3-18	
ND	1.0	EPA 8260C	7-3-18	7-3-18	
ND	0.20	EPA 8260C	7-3-18	7-3-18	
ND	2.0	EPA 8260C	7-3-18	7-3-18	
ND	1.0	EPA 8260C	7-3-18	7-3-18	
ND	0.20	EPA 8260C	7-3-18	7-3-18	
ND	0.20	EPA 8260C	7-3-18	7-3-18	
ND	5.0	EPA 8260C	7-3-18	7-3-18	
ND	5.0	EPA 8260C	7-3-18	7-3-18	
ND	0.20	EPA 8260C	7-3-18	7-3-18	
ND	2.0	EPA 8260C	7-3-18	7-3-18	
ND	0.20	EPA 8260C	7-3-18	7-3-18	
ND	0.20	EPA 8260C	7-3-18	7-3-18	
ND	0.20	EPA 8260C	7-3-18	7-3-18	
ND	1.0	EPA 8260C	7-3-18	7-3-18	
ND	0.20	EPA 8260C	7-3-18	7-3-18	
ND	0.20	EPA 8260C	7-3-18	7-3-18	
ND	5.0	EPA 8260C	7-3-18	7-3-18	
ND	0.20	EPA 8260C	7-3-18	7-3-18	
ND	0.20	EPA 8260C	7-3-18	7-3-18	
ND	0.20	EPA 8260C	7-3-18	7-3-18	
ND	0.20	EPA 8260C	7-3-18	7-3-18	
ND	0.20	EPA 8260C	7-3-18	7-3-18	
ND	0.20	EPA 8260C	7-3-18	7-3-18	
ND	0.20	EPA 8260C	7-3-18	7-3-18	
ND	0.20	EPA 8260C	7-3-18	7-3-18	
ND	0.20	EPA 8260C	7-3-18	7-3-18	
ND	0.20	EPA 8260C	7-3-18	7-3-18	
ND	0.20	EPA 8260C	7-3-18	7-3-18	
ND	1.0	EPA 8260C	7-3-18	7-3-18	
ND	0.20	EPA 8260C	7-3-18	7-3-18	
ND	2.0	EPA 8260C	7-3-18	7-3-18	
ND	1.0	EPA 8260C	7-3-18	7-3-18	
ND	0.20	EPA 8260C	7-3-18	7-3-18	
	Trip Blank 06-310-04 ND	Trip Blank 06-310-04 ND 0.20 ND 1.0 ND 0.20 ND 2.0 ND 1.0 ND 0.20 ND 0.20 ND 5.0 ND 5.0 ND 0.20 ND	Trip Blank 06-310-04 ND 0.20 EPA 8260C ND 1.0 EPA 8260C ND 0.20 EPA 8260C ND 0.20 EPA 8260C ND 1.0 EPA 8260C ND 0.20 EPA 8260C ND 0.20 EPA 8260C ND 5.0 EPA 8260C ND 5.0 EPA 8260C ND 5.0 EPA 8260C ND 0.20 EPA 8260C ND <t< td=""><td>Result PQL Method Prepared Trip Blank 06-310-04 06-310-04 FPA 8260C 7-3-18 ND 0.20 EPA 8260C 7-3-18 ND 0.20 EPA 8260C 7-3-18 ND 0.20 EPA 8260C 7-3-18 ND 1.0 EPA 8260C 7-3-18 ND 0.20 EPA 8260C 7-3-18 ND 0.20 EPA 8260C 7-3-18 ND 0.20 EPA 8260C 7-3-18 ND 5.0 EPA 8260C 7-3-18 ND 5.0 EPA 8260C 7-3-18 ND 5.0 EPA 8260C 7-3-18 ND 0.20 EPA 8260C 7-3-18 <tr< td=""><td>Result PQL Method Prepared Analyzed Trip Blank 06-310-04 06-310-04 FPA 8260C 7-3-18 7-3-18 ND 0.20 EPA 8260C 7-3-18 7-3-18 ND 1.0 EPA 8260C 7-3-18 7-3-18 ND 2.0 EPA 8260C 7-3-18 7-3-18 ND 1.0 EPA 8260C 7-3-18 7-3-18 ND 1.0 EPA 8260C 7-3-18 7-3-18 ND 0.20 EPA 8260C 7-3-18 7-3-18 ND 0.20 EPA 8260C 7-3-18 7-3-18 ND 0.20 EPA 8260C 7-3-18 7-3-18 ND 5.0 EPA 8260C 7-3-18 7-3-18 ND 5.0 EPA 8260C 7-3-18 7-3-18 ND 0.20 EPA 8260C 7-3-18 7-3-18 ND 0.20 EPA 8260C 7-3-18 7-3-18 ND 0.20 EPA 8260C 7-3-18 7-3-1</td></tr<></td></t<>	Result PQL Method Prepared Trip Blank 06-310-04 06-310-04 FPA 8260C 7-3-18 ND 0.20 EPA 8260C 7-3-18 ND 0.20 EPA 8260C 7-3-18 ND 0.20 EPA 8260C 7-3-18 ND 1.0 EPA 8260C 7-3-18 ND 0.20 EPA 8260C 7-3-18 ND 0.20 EPA 8260C 7-3-18 ND 0.20 EPA 8260C 7-3-18 ND 5.0 EPA 8260C 7-3-18 ND 5.0 EPA 8260C 7-3-18 ND 5.0 EPA 8260C 7-3-18 ND 0.20 EPA 8260C 7-3-18 <tr< td=""><td>Result PQL Method Prepared Analyzed Trip Blank 06-310-04 06-310-04 FPA 8260C 7-3-18 7-3-18 ND 0.20 EPA 8260C 7-3-18 7-3-18 ND 1.0 EPA 8260C 7-3-18 7-3-18 ND 2.0 EPA 8260C 7-3-18 7-3-18 ND 1.0 EPA 8260C 7-3-18 7-3-18 ND 1.0 EPA 8260C 7-3-18 7-3-18 ND 0.20 EPA 8260C 7-3-18 7-3-18 ND 0.20 EPA 8260C 7-3-18 7-3-18 ND 0.20 EPA 8260C 7-3-18 7-3-18 ND 5.0 EPA 8260C 7-3-18 7-3-18 ND 5.0 EPA 8260C 7-3-18 7-3-18 ND 0.20 EPA 8260C 7-3-18 7-3-18 ND 0.20 EPA 8260C 7-3-18 7-3-18 ND 0.20 EPA 8260C 7-3-18 7-3-1</td></tr<>	Result PQL Method Prepared Analyzed Trip Blank 06-310-04 06-310-04 FPA 8260C 7-3-18 7-3-18 ND 0.20 EPA 8260C 7-3-18 7-3-18 ND 1.0 EPA 8260C 7-3-18 7-3-18 ND 2.0 EPA 8260C 7-3-18 7-3-18 ND 1.0 EPA 8260C 7-3-18 7-3-18 ND 1.0 EPA 8260C 7-3-18 7-3-18 ND 0.20 EPA 8260C 7-3-18 7-3-18 ND 0.20 EPA 8260C 7-3-18 7-3-18 ND 0.20 EPA 8260C 7-3-18 7-3-18 ND 5.0 EPA 8260C 7-3-18 7-3-18 ND 5.0 EPA 8260C 7-3-18 7-3-18 ND 0.20 EPA 8260C 7-3-18 7-3-18 ND 0.20 EPA 8260C 7-3-18 7-3-18 ND 0.20 EPA 8260C 7-3-18 7-3-1

Project: 17-06520-000

VOLATILES EPA 8260C

page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	Trip Blank	PQL	Wethou	Frepareu	Allalyzeu	riays
Laboratory ID:	06-310-04					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Tetrachloroethene	ND	0.20	EPA 8260C	7-3-18 7-3-18	7-3-18 7-3-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	7-3-18 7-3-18	7-3-18 7-3-18	
2-Hexanone	ND	2.0	EPA 8260C	7-3-18 7-3-18	7-3-18 7-3-18	
Dibromochloromethane	ND	0.20	EPA 8260C	7-3-18 7-3-18	7-3-18 7-3-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	7-3-18 7-3-18	7-3-18 7-3-18	
Chlorobenzene	ND	0.20	EPA 8260C	7-3-18 7-3-18	7-3-18 7-3-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	7-3-18 7-3-18	7-3-18 7-3-18	
	ND ND	0.20	EPA 8260C EPA 8260C	7-3-16 7-3-18	7-3-16 7-3-18	
Ethylbenzene	ND ND	0.40	EPA 8260C EPA 8260C	7-3-16 7-3-18	7-3-16 7-3-18	
m,p-Xylene o-Xylene	ND ND	0.40	EPA 8260C EPA 8260C	7-3-16 7-3-18	7-3-16 7-3-18	
· ·	ND ND	0.20	EPA 8260C EPA 8260C	7-3-16 7-3-18	7-3-16 7-3-18	
Styrene	ND ND	1.0				
Bromoform	ND ND	0.20	EPA 8260C	7-3-18 7-3-18	7-3-18 7-3-18	
Isopropylbenzene			EPA 8260C			
Bromobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
n-Propylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
n-Butylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	7-3-18	7-3-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	7-3-18	7-3-18	
Naphthalene	ND	1.0	EPA 8260C	7-3-18	7-3-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Surrogate:	Percent Recovery	Control Limits				

Surrogate: Percent Recovery Control Limits
Dibromofluoromethane 99 75-127
Toluene-d8 95 80-127
4-Bromofluorobenzene 93 78-125



VOLATILES by EPA 8260C METHOD BLANK QUALITY CONTROL

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0703W1					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Chloromethane	ND	1.0	EPA 8260C	7-3-18	7-3-18	
Vinyl Chloride	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Bromomethane	ND	2.0	EPA 8260C	7-3-18	7-3-18	
Chloroethane	ND	1.0	EPA 8260C	7-3-18	7-3-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Acetone	ND	5.0	EPA 8260C	7-3-18	7-3-18	
Iodomethane	ND	5.0	EPA 8260C	7-3-18	7-3-18	
Carbon Disulfide	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Methylene Chloride	ND	2.0	EPA 8260C	7-3-18	7-3-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Vinyl Acetate	ND	1.0	EPA 8260C	7-3-18	7-3-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
2-Butanone	ND	5.0	EPA 8260C	7-3-18	7-3-18	
Bromochloromethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Chloroform	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Benzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Trichloroethene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Dibromomethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Bromodichloromethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	7-3-18	7-3-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	7-3-18	7-3-18	
Toluene	ND	1.0	EPA 8260C	7-3-18	7-3-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	7-3-18	7-3-18	

Project: 17-06520-000

VOLATILES by EPA 8260C METHOD BLANK QUALITY CONTROL

page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Allalyte	Nesuit	r Q L	Wetriou	гтератец	Allalyzeu	i iags
Laboratory ID:	MB0703W1					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Tetrachloroethene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
2-Hexanone	ND	2.0	EPA 8260C	7-3-18	7-3-18	
Dibromochloromethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Chlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Ethylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
m,p-Xylene	ND	0.40	EPA 8260C	7-3-18	7-3-18	
o-Xylene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Styrene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Bromoform	ND	1.0	EPA 8260C	7-3-18	7-3-18	
Isopropylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Bromobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	7-3-18	7-3-18	
n-Propylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
n-Butylbenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	7-3-18	7-3-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	7-3-18	7-3-18	
Naphthalene	ND	1.0	EPA 8260C	7-3-18	7-3-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	7-3-18	7-3-18	
Surrogate:	Percent Recovery	Control Limits				

Surrogate: Percent Recovery Control Limit
Dibromofluoromethane 100 75-127
Toluene-d8 98 80-127
4-Bromofluorobenzene 98 78-125



VOLATILES by EPA 8260C MS/MSD QUALITY CONTROL

					Source	Per	cent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Rec	overy	Limits	RPD	Limit	Flags
MATRIX SPIKES											
Laboratory ID:	06-3	11-06									
	MS	MSD	MS	MSD		MS	MSD				
1,1-Dichloroethene	10.4	10.6	10.0	10.0	ND	104	106	60-124	2	17	
Benzene	10.1	10.4	10.0	10.0	ND	101	104	67-130	3	22	
Trichloroethene	9.72	9.60	10.0	10.0	ND	97	96	71-120	1	15	
Toluene	10.2	10.0	10.0	10.0	ND	102	100	79-118	2	24	
Chlorobenzene	9.93	9.82	10.0	10.0	ND	99	98	74-120	1	17	
Surrogate:											
Dibromofluoromethane						100	99	75-127			
Toluene-d8						97	96	80-127			
4-Bromofluorobenzene						96	95	78-125			

Project: 17-06520-000

TOTAL METALS EPA 200.8/7470A

Matrix: Water
Units: ug/L (ppb)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	SW1					
Laboratory ID:	06-310-01					
Arsenic	ND	3.3	EPA 200.8	7-5-18	7-5-18	
Cadmium	ND	4.4	EPA 200.8	7-5-18	7-5-18	
Chromium	ND	11	EPA 200.8	7-5-18	7-5-18	
Lead	ND	1.1	EPA 200.8	7-5-18	7-5-18	
Mercury	ND	0.50	EPA 7470A	7-5-18	7-5-18	
Client ID:	SW2					
Laboratory ID:	06-310-02					
Arsenic	ND	3.3	EPA 200.8	7-5-18	7-5-18	
Cadmium	ND ND	3.3 4.4	EPA 200.8 EPA 200.8	7-5-18 7-5-18	7-5-18 7-5-18	
Chromium	ND ND	11	EPA 200.8 EPA 200.8	7-5-18 7-5-18	7-5-18 7-5-18	
Lead	ND ND	1.1	EPA 200.8	7-5-18 7-5-18	7-5-18 7-5-18	
Mercury	ND	0.50	EPA 7470A	7-5-18	7-5-18	
Client ID:	SW3					
Laboratory ID:	06-310-03					
Arsenic	ND	3.3	EPA 200.8	7-5-18	7-5-18	
Cadmium	ND	4.4	EPA 200.8	7-5-18	7-5-18	
Chromium	ND	11	EPA 200.8	7-5-18	7-5-18	
Lead	ND	1.1	EPA 200.8	7-5-18	7-5-18	
Mercury	ND	0.50	EPA 7470A	7-5-18	7-5-18	

Project: 17-06520-000

TOTAL METALS EPA 200.8/7470A QUALITY CONTROL

Matrix: Water
Units: ug/L (ppb)

								Date	Date		
Analyte		Result		PQL	Me	ethod		Prepared	Analyzed		Flags
METHOD BLANK											
Laboratory ID:	N	/IB0705WM1									
Arsenic		ND		3.3	EPA	A 200.	8	7-5-18	7-5-18	3	
Cadmium		ND		4.4	EPA	A 200.	8	7-5-18	7-5-18	3	
Chromium		ND		11	EP/	\ 200.	8	7-5-18	7-5-18	3	
Lead		ND		1.1	EPA	200.	8	7-5-18	7-5-18	3	
Laboratory ID:		MB0705W1									
Mercury		ND		0.50	EPA	7470	Α	7-5-18	7-5-18	3	
					Source	Pe	rcent	Recovery		RPD	
Analyte	Res	sult	Spike	e Level	Result	Rec	overy	Limits	RPD	Limit	Flags
DUPLICATE											
Laboratory ID:	07-00	06-02									
	ORIG	DUP									
Arsenic	5.67	6.06	NA	NA		I	NA	NA	7	20	
Cadmium	ND	ND	NA	NA		I	NA	NA	NA	20	
Chromium	ND	ND	NA	NA			NA	NA	NA	20	
Lead	ND	ND	NA	NA			NA	NA	NA	20	
Laboratory ID:	07-00	05-01									
Mercury	ND	ND	NA	NA			NA	NA	NA	20	
MATRIX SPIKES											
Laboratory ID:	07-00	06-02									
	MS	MSD	MS	MSD		MS	MSD)			
Arsenic	233	243	222	222	5.67	102	107	75-125	4	20	
Cadmium	228	235	222	222	ND	103	106	75-125	3	20	
Chromium	211	216	222	222	ND	95	97	75-125	2	20	
Lead	219	221	222	222	ND	99	100	75-125	1	20	
Laboratory ID:	07-00	05-01									

85

75-125

20

89

Mercury

11.1

10.6

12.5

12.5

ND

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	SW1					_
Laboratory ID:	06-310-01					
Benzo[a]anthracene	ND	0.011	EPA 8270D/SIM	7-2-18	7-5-18	
Chrysene	ND	0.011	EPA 8270D/SIM	7-2-18	7-5-18	
Benzo[b]fluoranthene	ND	0.011	EPA 8270D/SIM	7-2-18	7-5-18	
Benzo(j,k)fluoranthene	ND	0.011	EPA 8270D/SIM	7-2-18	7-5-18	
Benzo[a]pyrene	ND	0.011	EPA 8270D/SIM	7-2-18	7-5-18	
Indeno(1,2,3-c,d)pyrene	ND	0.011	EPA 8270D/SIM	7-2-18	7-5-18	
Dibenz[a,h]anthracene	ND	0.011	EPA 8270D/SIM	7-2-18	7-5-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	66	21 - 110				
Pyrene-d10	85	19 - 111				
Terphenyl-d14	84	32 - 137				

cPAHs EPA 8270D/SIM

			Date	Date	
Result	PQL	Method	Prepared	Analyzed	Flags
SW2					
06-310-02					
ND	0.010	EPA 8270D/SIM	7-2-18	7-5-18	
ND	0.010	EPA 8270D/SIM	7-2-18	7-5-18	
ND	0.010	EPA 8270D/SIM	7-2-18	7-5-18	
ND	0.010	EPA 8270D/SIM	7-2-18	7-5-18	
ND	0.010	EPA 8270D/SIM	7-2-18	7-5-18	
ND	0.010	EPA 8270D/SIM	7-2-18	7-5-18	
ND	0.010	EPA 8270D/SIM	7-2-18	7-5-18	
Percent Recovery	Control Limits				
65	21 - 110				
85	19 - 111				
82	32 - 137				
	\$W2 06-310-02 ND ND ND ND ND ND ND ND Percent Recovery 65 85	SW2 06-310-02 0.010 ND 0.010 ND 0.010 ND 0.010 ND 0.010 ND 0.010 ND 0.010 Percent Recovery Control Limits 65 21 - 110 85 19 - 111	SW2 06-310-02 ND 0.010 EPA 8270D/SIM Percent Recovery Control Limits 65 21 - 110 85 19 - 111	Result PQL Method Prepared SW2 06-310-02 ND 0.010 EPA 8270D/SIM 7-2-18 Percent Recovery Control Limits 65 21 - 110 85 19 - 111	Result PQL Method Prepared Analyzed SW2 06-310-02

cPAHs EPA 8270D/SIM

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	06-310-03					
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	7-2-18	7-5-18	
Chrysene	ND	0.010	EPA 8270D/SIM	7-2-18	7-5-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	7-2-18	7-5-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	7-2-18	7-5-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	7-2-18	7-5-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	7-2-18	7-5-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	7-2-18	7-5-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	66	21 - 110				
Pyrene-d10	81	19 - 111				
Terphenyl-d14	80	32 - 137				

Date of Report: July 10, 2018 Samples Submitted: June 29, 2018 Laboratory Reference: 1806-310

Project: 17-06520-000

CPAHS EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0702W1					
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	7-2-18	7-5-18	
Chrysene	ND	0.010	EPA 8270D/SIM	7-2-18	7-5-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	7-2-18	7-5-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	7-2-18	7-5-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	7-2-18	7-5-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	7-2-18	7-5-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	7-2-18	7-5-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	71	21 - 110				
Pyrene-d10	86	19 - 111				
Terphenyl-d14	90	32 - 137				

Date of Report: July 10, 2018 Samples Submitted: June 29, 2018 Laboratory Reference: 1806-310 Project: 17-06520-000

cPAHs EPA 8270D/SIM SB/SBD QUALITY CONTROL

3					Р	ercent	Recovery		RPD	
Analyte	Re	sult	Spike	Level	Re	covery	Limits	RPD	Limit	Flags
SPIKE BLANKS										
Laboratory ID:	SB07	02W1								
	SB	SBD	SB	SBD	SE	SBD				
Benzo[a]anthracene	0.472	0.511	0.500	0.500	94	102	57 - 127	8	15	
Chrysene	0.466	0.499	0.500	0.500	93	100	51 - 120	7	15	
Benzo[b]fluoranthene	0.467	0.509	0.500	0.500	93	102	54 - 124	9	17	
Benzo(j,k)fluoranthene	0.483	0.502	0.500	0.500	97	100	50 - 127	4	18	
Benzo[a]pyrene	0.450	0.483	0.500	0.500	90	97	50 - 120	7	16	
Indeno(1,2,3-c,d)pyrene	0.483	0.500	0.500	0.500	97	100	46 - 132	3	20	
Dibenz[a,h]anthracene	0.480	0.516	0.500	0.500	96	103	49 - 129	7	18	
Surrogate:										
2-Fluorobiphenyl					64	78	21 - 110			
Pyrene-d10					92	95	19 - 111			
Terphenyl-d14					95	99	32 - 137			

Date of Report: July 10, 2018 Samples Submitted: June 29, 2018 Laboratory Reference: 1806-310

Project: 17-06520-000

HARDNESS EPA 6010D/SM 2340B

Matrix: Water

Units: mg eqt. CaCO3/L (ppm)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	SW1					
Laboratory ID:	06-310-01					
Hardness	66	1.0	6010D/SM 2340B	7-3-18	7-5-18	
Client ID:	SW2					
Laboratory ID:	06-310-02					
Hardness	71	1.0	6010D/SM 2340B	7-3-18	7-5-18	
Client ID:	SW3					
Laboratory ID:	06-310-03					
Hardness	74	1.0	6010D/SM 2340B	7-3-18	7-5-18	

Date of Report: July 10, 2018 Samples Submitted: June 29, 2018 Laboratory Reference: 1806-310 Project: 17-06520-000

HARDNESS EPA 6010D/SM 2340B **QUALITY CONTROL**

Matrix: Water

Units: mg eqt. CaCO3/L (ppm)

Analista	Decult	DOL	Mathad	Date	Date	Flores
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0703WH1					
Hardness	ND	1.0	6010D/SM 2340B	7-3-18	7-5-18	

					Source	Per	cent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Rec	overy	Limits	RPD	Limit	Flags
DUPLICATE											
Laboratory ID:	06-3	10-01									
	ORIG	DUP									
Hardness	65.6	64.6	N	IA	NA	N	NΑ	NA	2	20	
MATRIX SPIKES											
Laboratory ID:	06-3	10-01									
	MS	MSD	MS	MSD		MS	MSD				
Hardness	200	201	132	132	65.6	102	103	75-125	0	20	
SPIKE BLANK											
Laboratory ID:	SB070	3WH1									
	S	В	S	BB	•	5	SB		•	•	
Hardness	14	40	1	32	NA	1	06	80-120	NA	NA	



Data Qualifiers and Abbreviations

- A Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
- B The analyte indicated was also found in the blank sample.
- C The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
- E The value reported exceeds the quantitation range and is an estimate.
- F Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
- H The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
- I Compound recovery is outside of the control limits.
- J The value reported was below the practical quantitation limit. The value is an estimate.
- K Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
- L The RPD is outside of the control limits.
- M Hydrocarbons in the gasoline range are impacting the diesel range result.
- M1 Hydrocarbons in the gasoline range (toluene-naphthalene) are present in the sample.
- N Hydrocarbons in the lube oil range are impacting the diesel range result.
- N1 Hydrocarbons in diesel range are impacting lube oil range results.
- O Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
- P The RPD of the detected concentrations between the two columns is greater than 40.
- Q Surrogate recovery is outside of the control limits.
- S Surrogate recovery data is not available due to the necessary dilution of the sample.
- T The sample chromatogram is not similar to a typical ______.
- U The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- U1 The practical quantitation limit is elevated due to interferences present in the sample.
- V Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
- W Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
- X Sample extract treated with a mercury cleanup procedure.
- X1- Sample extract treated with a sulfuric acid/silica gel cleanup procedure.
- Y The calibration verification for this analyte exceeded the 20% drift specified in method 8260C, and therefore the reported result should be considered an estimate. The overall performance of the calibration verification standard met the acceptance criteria of the method.

7 -

ND - Not Detected at PQL

PQL - Practical Quantitation Limit

RPD - Relative Percent Difference



MA	OnSite	Inc
	Environmental	inc.

Chain of Custody

	1		1	
Page	1	of	1	

Analytical Laboratory Testing Services 14648 NE 95th Street • Redmond, WA 98052	Turnaround Request (in working days)		La	abo	rato	ry N	lumb	er:			06	-3	3 1	0					197			
Phone: (425) 883-3881 • www.onsite-env.com Company: Herrera Env, Cons. Project Number - 065 20 - 000 Project Name: Pacific Powk Project Manager: BAKK EWBANK Sampled by George IAHR Lab ID Sample Identification	(Check One) Same Day 1 Day 2 Days 3 Days Standard (7 Days) (TPH analysis 5 Days) (other) Date Time Sampled Sampled Matrix	Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX		NWTPH-Dx (Acid / SG Clean-up) Volatiles 8260C	Halogenated Volatiles 8260C	EDB EPA 8011 (Waters Only)	Semivolatiles 8270D/SIM (with low-level PAHs)	PAHS 8270D/SIM (low-level)	PUBS 8082A Organochlorine Pesticides 8081B	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	TCLP Metals	HEM (oil and grease) 1664A	CPAHS 8270D/SIM	HWINOSS		% Moisture	
SW/	6/29/18 9120 water				1	1	/								1			V	/			
1 SW7	18:45	11			V	V	V								1			V	V			
3 SW3	Y, 8:15 V	n			V	/	/								1			V	V			
4 Trip Blank	6/29/18 10:00 water	-3				1																
		Ħ																		1		
Signature	Company			Date	/		ime			ment	s/Spec	ial Ins	tructio	ons								
Received Received Received Relinquished Received Relinquished	Heyreva Env ALPHA ALPHA OSE	. G	15,	6/2	29/1	81	12:14	8/11	H	DA	e f	es	p nl·	CE	35,	pe	en e	dia	ng			
Received									Data	Pack	kage:	Stand	lard [Le	evel III		Leve	I IV [1			
Reviewed/Date	Reviewed/Date								Chro	mato	grams	with f	inal re	port [☐ Ele	ectroni	c Data	a Deliv	erables	s (EDD	s) 🗌	



14040 NE 33 Otteet, Neumona, WA 30032 (423) 003-3001

October 5, 2018

Mark Ewbank Herrera Environmental Consultants, Inc. 2200 6th Avenue, Suite 1100 Seattle, WA 98121

Re: Analytical Data for Project 17-06520-000

Laboratory Reference No. 1809-277

Dear Mark:

Enclosed are the analytical results and associated quality control data for samples submitted on September 27, 2018.

The standard policy of OnSite Environmental, Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

David Baumeister Project Manager

Enclosures



Date of Report: October 5, 2018

Samples Submitted: September 27, 2018 Laboratory Reference: 1809-277

Project: 17-06520-000

Case Narrative

Samples were collected on September 26, 2018 and received by the laboratory on September 27, 2018. They were maintained at the laboratory at a temperature of 2°C to 6°C.

Please note that any and all soil sample results are reported on a dry-weight basis, unless otherwise noted below.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

cPAHs EPA 8270D/SIM Analysis

Samples MW5 and method blank MB0927W1 each had one surrogate recovery out of control limits. This is within allowance of our standard operating procedure as long as the recovery is above 10%.

Any other QA/QC issues associated with this extraction and analysis will be indicated with a footnote reference and discussed in detail on the Data Qualifier page.

Laboratory Reference: 1809-277 Project: 17-06520-000

GASOLINE RANGE ORGANICS NWTPH-Gx

Matrix: Water Units: ug/L (ppb)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW1					
Laboratory ID:	09-277-01					
Gasoline	ND	100	NWTPH-Gx	9-28-18	9-28-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	88	66-117				
Client ID:	MW2					
Laboratory ID:	09-277-02					
Gasoline	ND	100	NWTPH-Gx	9-28-18	9-28-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	90	66-117				
Client ID:	MW3					
Laboratory ID:	09-277-03					
Gasoline	ND	100	NWTPH-Gx	9-28-18	9-28-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	90	66-117				
Client ID:	MW4					
Laboratory ID:	09-277-04					
Gasoline	ND	100	NWTPH-Gx	9-28-18	9-28-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	89	66-117				
Client ID:	MW5					
Laboratory ID:	09-277-05					
Gasoline	ND	100	NWTPH-Gx	9-28-18	9-28-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	88	66-117				
Client ID:	MW6					
Laboratory ID:	09-277-06					
Gasoline	ND	100	NWTPH-Gx	9-28-18	9-28-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	92	66-117				

Date of Report: October 5, 2018

Samples Submitted: September 27, 2018

Laboratory Reference: 1809-277 Project: 17-06520-000

GASOLINE RANGE ORGANICS NWTPH-Gx

Matrix: Water Units: ug/L (ppb)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW7					
Laboratory ID:	09-277-07					
Gasoline	ND	100	NWTPH-Gx	9-28-18	9-28-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	89	66-117				
Client ID:	MW8					
Laboratory ID:	09-277-08					
Gasoline	ND	100	NWTPH-Gx	9-28-18	9-28-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	90	66-117				
Client ID:	MW9					
Laboratory ID:	09-277-09					
Gasoline	ND	100	NWTPH-Gx	9-28-18	9-28-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	88	66-117				

Date of Report: October 5, 2018

Samples Submitted: September 27, 2018

Laboratory Reference: 1809-277 Project: 17-06520-000

GASOLINE RANGE ORGANICS NWTPH-Gx QUALITY CONTROL

Matrix: Water Units: ug/L (ppb)

Analyto	Result	PQL	Method	Date Prepared	Date Analvzed	Flage
Analyte METHOD BLANK	Result	FQL	METHOR	гтератец	Allalyzeu	Flags
	MDOOONAA					
Laboratory ID:	MB0928W1					
Gasoline	ND	100	NWTPH-Gx	9-28-18	9-28-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	89	66-117				

Analyte	Re	sult	Spike	Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
DUPLICATE			•							
Laboratory ID:	09-2	77-01								
	ORIG	DUP								
Gasoline	ND	ND	NA	NA		NA	NA	NA	30	
Surrogate:			•				_	•		•
Fluorobenzene						88 89	66-117			

Laboratory Reference: 1809-277

Project: 17-06520-000

DIESEL AND HEAVY OIL RANGE ORGANICS NWTPH-Dx

Matrix: Water Units: mg/L (ppm)

Offics. Hig/L (ppiff)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW1					
Laboratory ID:	09-277-01					
Diesel Range Organics	ND	0.27	NWTPH-Dx	9-28-18	9-28-18	
Lube Oil Range Organics	ND	0.43	NWTPH-Dx	9-28-18	9-28-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	87	50-150				
Client ID:	MW2					
Laboratory ID:	09-277-02					
Diesel Range Organics	ND	0.26	NWTPH-Dx	9-28-18	9-28-18	
Lube Oil Range Organics	ND	0.41	NWTPH-Dx	9-28-18	9-28-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	84	50-150				
Client ID:	MW3					
Laboratory ID:	09-277-03					
Diesel Range Organics	ND	0.26	NWTPH-Dx	9-28-18	9-28-18	
Lube Oil Range Organics	ND	0.41	NWTPH-Dx	9-28-18	9-28-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	87	50-150				
Client ID:	MW4					
Laboratory ID:	09-277-04					
Diesel Range Organics	ND	0.25	NWTPH-Dx	9-28-18	9-28-18	
Lube Oil Range Organics	ND	0.41	NWTPH-Dx	9-28-18	9-28-18	
Surrogate:	Percent Recovery	Control Limits		0 20 .0	0 20 .0	
o-Terphenyl	90	50-150				
o respirent	00	00 700				
Client ID:	MW5					
Laboratory ID:	09-277-05					
Diesel Range Organics	ND	0.26	NWTPH-Dx	9-28-18	9-28-18	
Lube Oil Range Organics	ND	0.20	NWTPH-Dx	9-28-18	9-28-18	
Surrogate:	Percent Recovery	Control Limits		0 20 10	0 20 10	
o-Terphenyl	86	50-150				
o . orphonyi	30	00 100				
Client ID:	MW6					
Laboratory ID:	09-277-06					
Diesel Range Organics	ND	0.26	NWTPH-Dx	9-28-18	9-28-18	
Lube Oil Range Organics	ND	0.41	NWTPH-Dx	9-28-18	9-28-18	
Surrogate:	Percent Recovery	Control Limits	TANA IL IL-DV	J 20-10	J 20-10	
o-Terphenyl	89	50-150				
o respiretly)	Uð.	00 100				

Laboratory Reference: 1809-277 Project: 17-06520-000

DIESEL AND HEAVY OIL RANGE ORGANICS NWTPH-Dx

Matrix: Water Units: mg/L (ppm)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW7					
Laboratory ID:	09-277-07					
Diesel Range Organics	ND	0.26	NWTPH-Dx	9-28-18	9-28-18	
Lube Oil Range Organics	ND	0.41	NWTPH-Dx	9-28-18	9-28-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	103	50-150				
Client ID:	MW8					
••						
Laboratory ID:	09-277-08					
Diesel Range Organics	ND	0.26	NWTPH-Dx	9-28-18	9-28-18	
Lube Oil Range Organics	ND	0.41	NWTPH-Dx	9-28-18	9-28-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	98	50-150				
Client ID:	MW9					
Laboratory ID:	09-277-09					
	ND	0.25	NWTPH-Dx	9-28-18	10-1-18	
Diesel Range Organics		0.25				
Lube Oil Range Organics	ND	0.41	NWTPH-Dx	9-28-18	10-1-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	82	50-150				

Laboratory Reference: 1809-277 Project: 17-06520-000

DIESEL AND HEAVY OIL RANGE ORGANICS NWTPH-Dx QUALITY CONTROL

Matrix: Water Units: mg/L (ppm)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0928W1					
Diesel Range Organics	ND	0.25	NWTPH-Dx	9-28-18	10-1-18	
Lube Oil Range Organics	ND	0.40	NWTPH-Dx	9-28-18	10-1-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	83	50-150				

					Source	Perd	ent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Reco	very	Limits	RPD	Limit	Flags
DUPLICATE											
Laboratory ID:	09-27	77-01									
	ORIG	DUP									
Diesel Range	ND	ND	NA	NA		N	A	NA	NA	NA	
Lube Oil Range	ND	ND	NA	NA		N	A	NA	NA	NA	
Surrogate:											
o-Terphenyl						87	92	50-150			

Laboratory Reference: 1809-277 Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW1					
Laboratory ID:	09-277-01					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloromethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Vinyl Chloride	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromomethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloroethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Acetone	ND	5.0	EPA 8260C	9-29-18	9-29-18	
Iodomethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Carbon Disulfide	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methylene Chloride	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Vinyl Acetate	ND	1.0	EPA 8260C	9-29-18	9-29-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Butanone	ND	5.0	EPA 8260C	9-29-18	9-29-18	
Bromochloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloroform	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Benzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Trichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Dibromomethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromodichloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	9-29-18	9-29-18	
Toluene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	

Laboratory Reference: 1809-277 Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW1					
Laboratory ID:	09-277-01					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Tetrachloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Hexanone	ND	2.0	EPA 8260C	9-29-18	9-29-18	
Dibromochloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Ethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
m,p-Xylene	ND	0.40	EPA 8260C	9-29-18	9-29-18	
o-Xylene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Styrene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromoform	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Isopropylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
n-Propylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
n-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Naphthalene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	100	75-127				

Date of Report: October 5, 2018 Samples Submitted: September 27, 2018 Laboratory Reference: 1809-277

Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 1 of 2

omis. ug/L				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW2					
Laboratory ID:	09-277-02					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloromethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Vinyl Chloride	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromomethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloroethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Acetone	ND	5.0	EPA 8260C	9-29-18	9-29-18	
Iodomethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Carbon Disulfide	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methylene Chloride	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Vinyl Acetate	ND	1.0	EPA 8260C	9-29-18	9-29-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Butanone	ND	5.0	EPA 8260C	9-29-18	9-29-18	
Bromochloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloroform	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Benzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Trichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Dibromomethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromodichloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	9-29-18	9-29-18	
Toluene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	

Laboratory Reference: 1809-277

Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW2					
Laboratory ID:	09-277-02					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Tetrachloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Hexanone	ND	2.0	EPA 8260C	9-29-18	9-29-18	
Dibromochloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Ethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
m,p-Xylene	ND	0.40	EPA 8260C	9-29-18	9-29-18	
o-Xylene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Styrene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromoform	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Isopropylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
n-Propylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
n-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Naphthalene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	102	75-127				

Dibromofluoromethane 102 75-127 101 80-127 Toluene-d8 4-Bromofluorobenzene 99 78-125

Laboratory Reference: 1809-277 Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW3					
Laboratory ID:	09-277-03					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloromethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Vinyl Chloride	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromomethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloroethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Acetone	ND	5.0	EPA 8260C	9-29-18	9-29-18	
lodomethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Carbon Disulfide	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methylene Chloride	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Vinyl Acetate	ND	1.0	EPA 8260C	9-29-18	9-29-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
(cis) 1,2-Dichloroethene	0.35	0.20	EPA 8260C	9-29-18	9-29-18	
2-Butanone	ND	5.0	EPA 8260C	9-29-18	9-29-18	
Bromochloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloroform	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Benzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Trichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Dibromomethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromodichloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	9-29-18	9-29-18	
Toluene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	

Date of Report: October 5, 2018 Samples Submitted: September 27, 2018 Laboratory Reference: 1809-277

Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW3					
Laboratory ID:	09-277-03					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Tetrachloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Hexanone	ND	2.0	EPA 8260C	9-29-18	9-29-18	
Dibromochloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Ethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
m,p-Xylene	ND	0.40	EPA 8260C	9-29-18	9-29-18	
o-Xylene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Styrene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromoform	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Isopropylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
n-Propylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
n-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dibromo-3-chloropropane		1.0	EPA 8260C	9-29-18	9-29-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Naphthalene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Surrogate:	Percent Recovery		2.7102000	0 20 10	0 20 10	
Surrogate.	1 Groom Necovery	COILLOI LIIIIIG				

Dibromofluoromethane 99 75-127 Toluene-d8 99 80-127 4-Bromofluorobenzene 95 78-125

Laboratory Reference: 1809-277 Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW4					
Laboratory ID:	09-277-04					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloromethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Vinyl Chloride	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromomethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloroethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Acetone	ND	5.0	EPA 8260C	9-29-18	9-29-18	
lodomethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Carbon Disulfide	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methylene Chloride	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Vinyl Acetate	ND	1.0	EPA 8260C	9-29-18	9-29-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Butanone	ND	5.0	EPA 8260C	9-29-18	9-29-18	
Bromochloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloroform	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Benzene	0.22	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Trichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Dibromomethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromodichloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	9-29-18	9-29-18	
Toluene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	

Date of Report: October 5, 2018 Samples Submitted: September 27, 2018 Laboratory Reference: 1809-277

Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW4	PQL	Wethou	Prepared	Anaryzeu	riays
Laboratory ID:	09-277-04					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Tetrachloroethene	ND ND	0.20		9-29-16 9-29-18	9-29-16	
	ND ND	0.20	EPA 8260C			
1,3-Dichloropropane			EPA 8260C	9-29-18	9-29-18	
2-Hexanone	ND	2.0	EPA 8260C	9-29-18	9-29-18	
Dibromochloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chlorobenzene	4.6	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Ethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
m,p-Xylene	ND	0.40	EPA 8260C	9-29-18	9-29-18	
o-Xylene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Styrene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromoform	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Isopropylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
n-Propylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
n-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dibromo-3-chloropropane		1.0	EPA 8260C	9-29-18	9-29-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Naphthalene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Surrogate:	Percent Recovery	Control Limits				

Surrogate: Percent Recovery Control Limits
Dibromofluoromethane 98 75-127
Toluene-d8 100 80-127
4-Bromofluorobenzene 95 78-125

Laboratory Reference: 1809-277 Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW5					
Laboratory ID:	09-277-05					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloromethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Vinyl Chloride	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromomethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloroethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Acetone	ND	5.0	EPA 8260C	9-29-18	9-29-18	
Iodomethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Carbon Disulfide	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methylene Chloride	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Vinyl Acetate	ND	1.0	EPA 8260C	9-29-18	9-29-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Butanone	ND	5.0	EPA 8260C	9-29-18	9-29-18	
Bromochloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloroform	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Benzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Trichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Dibromomethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromodichloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	9-29-18	9-29-18	
Toluene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	

Date of Report: October 5, 2018 Samples Submitted: September 27, 2018 Laboratory Reference: 1809-277

Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW5					
Laboratory ID:	09-277-05					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Tetrachloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Hexanone	ND	2.0	EPA 8260C	9-29-18	9-29-18	
Dibromochloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Ethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
m,p-Xylene	ND	0.40	EPA 8260C	9-29-18	9-29-18	
o-Xylene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Styrene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromoform	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Isopropylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
n-Propylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
n-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dibromo-3-chloropropane		1.0	EPA 8260C	9-29-18	9-29-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Naphthalene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	98	75-127				

Dibromofluoromethane 75-127 98 80-127 Toluene-d8 99 4-Bromofluorobenzene 95 78-125

Laboratory Reference: 1809-277 Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW6					
Laboratory ID:	09-277-06					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloromethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Vinyl Chloride	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromomethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloroethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Acetone	ND	5.0	EPA 8260C	9-29-18	9-29-18	
Iodomethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Carbon Disulfide	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methylene Chloride	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Vinyl Acetate	ND	1.0	EPA 8260C	9-29-18	9-29-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Butanone	ND	5.0	EPA 8260C	9-29-18	9-29-18	
Bromochloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloroform	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Benzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Trichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Dibromomethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromodichloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	9-29-18	9-29-18	
Toluene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	

Laboratory Reference: 1809-277 Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW6					
Laboratory ID:	09-277-06					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Tetrachloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Hexanone	ND	2.0	EPA 8260C	9-29-18	9-29-18	
Dibromochloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Ethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
m,p-Xylene	ND	0.40	EPA 8260C	9-29-18	9-29-18	
o-Xylene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Styrene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromoform	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Isopropylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
n-Propylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,4-Dichlorobenzene	0.20	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
n-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dibromo-3-chloropropane	e ND	1.0	EPA 8260C	9-29-18	9-29-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Naphthalene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Surrogate:	Percent Recovery	Control Limits				

Dibromofluoromethane 99 75-127 Toluene-d8 101 80-127 4-Bromofluorobenzene 97 78-125 Date of Report: October 5, 2018 Samples Submitted: September 27, 2018 Laboratory Reference: 1809-277 Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW7					
Laboratory ID:	09-277-07					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloromethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Vinyl Chloride	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromomethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloroethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Acetone	ND	5.0	EPA 8260C	9-29-18	9-29-18	
Iodomethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Carbon Disulfide	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methylene Chloride	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Vinyl Acetate	ND	1.0	EPA 8260C	9-29-18	9-29-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Butanone	ND	5.0	EPA 8260C	9-29-18	9-29-18	
Bromochloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloroform	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Benzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Trichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Dibromomethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromodichloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	9-29-18	9-29-18	
Toluene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	

Laboratory Reference: 1809-277

Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW7					
Laboratory ID:	09-277-07					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Tetrachloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Hexanone	ND	2.0	EPA 8260C	9-29-18	9-29-18	
Dibromochloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Ethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
m,p-Xylene	ND	0.40	EPA 8260C	9-29-18	9-29-18	
o-Xylene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Styrene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromoform	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Isopropylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
n-Propylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
n-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Naphthalene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Surrogate:	Percent Recovery	Control Limits				
Dibromoflyoromothono	06	75 107				

Surrogate: Percent Recovery Control Limit Dibromofluoromethane 96 75-127 Toluene-d8 101 80-127 4-Bromofluorobenzene 95 78-125



Laboratory Reference: 1809-277 Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW8					
Laboratory ID:	09-277-08					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloromethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Vinyl Chloride	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromomethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloroethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Acetone	ND	5.0	EPA 8260C	9-29-18	9-29-18	
Iodomethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Carbon Disulfide	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methylene Chloride	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Vinyl Acetate	ND	1.0	EPA 8260C	9-29-18	9-29-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Butanone	ND	5.0	EPA 8260C	9-29-18	9-29-18	
Bromochloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloroform	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Benzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Trichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Dibromomethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromodichloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	9-29-18	9-29-18	
Toluene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	

Laboratory Reference: 1809-277

Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW8					
Laboratory ID:	09-277-08					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Tetrachloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Hexanone	ND	2.0	EPA 8260C	9-29-18	9-29-18	
Dibromochloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Ethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
m,p-Xylene	ND	0.40	EPA 8260C	9-29-18	9-29-18	
o-Xylene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Styrene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromoform	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Isopropylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
n-Propylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
n-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Naphthalene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Surrogate:	Percent Recovery	Control Limits				<u> </u>

Surrogate: Percent Recovery Control Limit Dibromofluoromethane 100 75-127
Toluene-d8 102 80-127
4-Bromofluorobenzene 98 78-125



Laboratory Reference: 1809-277 Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW9					
Laboratory ID:	09-277-09					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloromethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Vinyl Chloride	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromomethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloroethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Acetone	ND	5.0	EPA 8260C	9-29-18	9-29-18	
lodomethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Carbon Disulfide	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methylene Chloride	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Vinyl Acetate	ND	1.0	EPA 8260C	9-29-18	9-29-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Butanone	ND	5.0	EPA 8260C	9-29-18	9-29-18	
Bromochloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloroform	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Benzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Trichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Dibromomethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromodichloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	9-29-18	9-29-18	
Toluene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	

Laboratory Reference: 1809-277

Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW9					
Laboratory ID:	09-277-09					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Tetrachloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Hexanone	ND	2.0	EPA 8260C	9-29-18	9-29-18	
Dibromochloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chlorobenzene	0.38	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Ethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
m,p-Xylene	ND	0.40	EPA 8260C	9-29-18	9-29-18	
o-Xylene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Styrene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromoform	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Isopropylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
n-Propylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
n-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Naphthalene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Surrogate:	Percent Recovery	Control Limits				
Dibramafluaramathana	00	75 107				

Dibromofluoromethane 99 75-127 Toluene-d8 101 80-127 4-Bromofluorobenzene 97 78-125



Laboratory Reference: 1809-277 Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	Trip Blank					
Laboratory ID:	09-277-10					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	9-29-18	9-30-18	
Chloromethane	ND	1.0	EPA 8260C	9-29-18	9-30-18	
Vinyl Chloride	ND	0.20	EPA 8260C	9-29-18	9-30-18	
Bromomethane	ND	0.20	EPA 8260C	9-29-18	9-30-18	
Chloroethane	ND	1.0	EPA 8260C	9-29-18	9-30-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	9-29-18	9-30-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
Acetone	ND	5.0	EPA 8260C	9-29-18	9-30-18	
Iodomethane	ND	1.0	EPA 8260C	9-29-18	9-30-18	
Carbon Disulfide	ND	0.20	EPA 8260C	9-29-18	9-30-18	
Methylene Chloride	ND	1.0	EPA 8260C	9-29-18	9-30-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	9-29-18	9-30-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	9-29-18	9-30-18	
Vinyl Acetate	ND	1.0	EPA 8260C	9-29-18	9-30-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-30-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
2-Butanone	ND	5.0	EPA 8260C	9-29-18	9-30-18	
Bromochloromethane	ND	0.20	EPA 8260C	9-29-18	9-30-18	
Chloroform	ND	0.20	EPA 8260C	9-29-18	9-30-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	9-29-18	9-30-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	9-29-18	9-30-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
Benzene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	9-29-18	9-30-18	
Trichloroethene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-30-18	
Dibromomethane	ND	0.20	EPA 8260C	9-29-18	9-30-18	
Bromodichloromethane	ND	0.20	EPA 8260C	9-29-18	9-30-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	9-29-18	9-30-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	9-29-18	9-30-18	
Toluene	ND	1.0	EPA 8260C	9-29-18	9-30-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-30-18	

Date of Report: October 5, 2018 Samples Submitted: September 27, 2018 Laboratory Reference: 1809-277

Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	Trip Blank					
Laboratory ID:	09-277-10					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	9-29-18	9-30-18	
Tetrachloroethene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-30-18	
2-Hexanone	ND	2.0	EPA 8260C	9-29-18	9-30-18	
Dibromochloromethane	ND	0.20	EPA 8260C	9-29-18	9-30-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	9-29-18	9-30-18	
Chlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-29-18	9-30-18	
Ethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
m,p-Xylene	ND	0.40	EPA 8260C	9-29-18	9-30-18	
o-Xylene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
Styrene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
Bromoform	ND	1.0	EPA 8260C	9-29-18	9-30-18	
Isopropylbenzene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
Bromobenzene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-29-18	9-30-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	9-29-18	9-30-18	
n-Propylbenzene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
n-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	9-29-18	9-30-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	9-29-18	9-30-18	
Naphthalene	ND	1.0	EPA 8260C	9-29-18	9-30-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-30-18	
Surrogate:	Percent Recovery	Control Limits				
Dibramafluaramathana	105	75 107				

Dibromofluoromethane 105 75-127 Toluene-d8 102 80-127 4-Bromofluorobenzene 97 78-125

Laboratory Reference: 1809-277 Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C METHOD BLANK QUALITY CONTROL

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0929W1					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloromethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Vinyl Chloride	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromomethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloroethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Acetone	ND	5.0	EPA 8260C	9-29-18	9-29-18	
Iodomethane	ND	1.0	EPA 8260C	9-29-18	9-29-18	
Carbon Disulfide	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methylene Chloride	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Vinyl Acetate	ND	1.0	EPA 8260C	9-29-18	9-29-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Butanone	ND	5.0	EPA 8260C	9-29-18	9-29-18	
Bromochloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chloroform	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Benzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Trichloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Dibromomethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromodichloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	9-29-18	9-29-18	
Toluene	ND	1.0	EPA 8260C	9-29-18	9-29-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-29-18	9-29-18	

Date of Report: October 5, 2018 Samples Submitted: September 27, 2018 Laboratory Reference: 1809-277

Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C METHOD BLANK QUALITY CONTROL

page 2 of 2

Analyte		PQL		Date	Date	
	Result		Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0929W1					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Tetrachloroethene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3-Dichloropropane	ND ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Hexanone	ND	2.0	EPA 8260C	9-29-18	9-29-18	
Dibromochloromethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Chlorobenzene	ND ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Ethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
m,p-Xylene	ND ND	0.40	EPA 8260C	9-29-18	9-29-18	
o-Xylene	ND ND	0.40	EPA 8260C	9-29-18	9-29-18	
Styrene	ND ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromoform	ND ND	1.0	EPA 8260C	9-29-18	9-29-18	
Isopropylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
Bromobenzene	ND ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,1,2,2-Tetrachloroethane	ND ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	9-29-18	9-29-18	
n-Propylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
2-Chlorotoluene	ND ND	0.20	EPA 8260C	9-29-18	9-29-18	
4-Chlorotoluene	ND ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2,4-Trimethylbenzene	ND ND	0.20	EPA 8260C	9-29-18	9-29-18	
sec-Butylbenzene	ND ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	9-29-18	9-29-18	
p-isopropyitoluerie 1,4-Dichlorobenzene	ND ND	0.20	EPA 8260C EPA 8260C	9-29-18	9-29-16	
1,4-Dichlorobenzene	ND ND	0.20	EPA 8260C	9-29-16	9-29-18	
n-Butylbenzene	ND ND	0.20	EPA 8260C	9-29-18	9-29-18	
1,2-Dibromo-3-chloropropane		1.0	EPA 8260C	9-29-18	9-29-18	
1,2,4-Trichlorobenzene	ND ND	0.20	EPA 8260C EPA 8260C	9-29-18	9-29-16	
Hexachlorobutadiene	ND ND	1.0	EPA 8260C EPA 8260C	9-29-16	9-29-16	
Naphthalene	ND ND	1.0	EPA 8260C	9-29-18	9-29-18	
1,2,3-Trichlorobenzene	ND ND	0.20	EPA 8260C EPA 8260C	9-29-18	9-29-18	
_	Percent Recovery	Control Limits	EFA 02000	3-23-10	3-23-10	
Surrogate:	111	CONTROL LIMINS				

Dibromofluoromethane 111 75-127 Toluene-d8 105 80-127 4-Bromofluorobenzene 102 78-125

Laboratory Reference: 1809-277 Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C SB/SBD QUALITY CONTROL

					Per	cent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Rec	overy	Limits	RPD	Limit	Flags
SPIKE BLANKS										
Laboratory ID:	SB09:	29W1								
	SB	SBD	SB	SBD	SB	SBD				
1,1-Dichloroethene	8.87	9.06	10.0	10.0	89	91	62-129	2	15	
Benzene	9.32	9.49	10.0	10.0	93	95	77-127	2	15	
Trichloroethene	9.55	9.81	10.0	10.0	96	98	70-120	3	15	
Toluene	9.80	10.1	10.0	10.0	98	101	82-123	3	15	
Chlorobenzene	9.04	9.21	10.0	10.0	90	92	79-120	2	15	
Surrogate:										
Dibromofluoromethane					105	104	75-127			
Toluene-d8					103	103	80-127			
4-Bromofluorobenzene					100	101	78-125			

Laboratory Reference: 1809-277 Project: 17-06520-000

TOTAL METALS EPA 200.8/7470A

σ.mo. αg/2 (ppε/)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW1					
Laboratory ID:	09-277-01					
Arsenic	ND	3.3	EPA 200.8	9-28-18	9-28-18	
Cadmium	ND	4.4	EPA 200.8	9-28-18	9-28-18	
Chromium	ND	11	EPA 200.8	9-28-18	9-28-18	
Lead	2.4	1.1	EPA 200.8	9-28-18	9-28-18	
Mercury	ND	0.50	EPA 7470A	9-28-18	9-28-18	
Client ID:	MW2					
Laboratory ID:	09-277-02					
Arsenic	4.9	3.3	EPA 200.8	9-28-18	9-28-18	
Cadmium	ND	4.4	EPA 200.8	9-28-18	9-28-18	
Chromium	ND	11	EPA 200.8	9-28-18	9-28-18	
Lead	ND	1.1	EPA 200.8	9-28-18	9-28-18	
Mercury	ND	0.50	EPA 7470A	9-28-18	9-28-18	
			-			
Client ID:	MW3					
Laboratory ID:	09-277-03					
Arsenic	ND	3.3	EPA 200.8	9-28-18	9-28-18	
Cadmium	ND	4.4	EPA 200.8	9-28-18	9-28-18	
Chromium	ND	11	EPA 200.8	9-28-18	9-28-18	
Lead	ND	1.1	EPA 200.8	9-28-18	9-28-18	
Mercury	ND	0.50	EPA 7470A	9-28-18	9-28-18	
Client ID:	MW4					
Laboratory ID:	09-277-04					
Arsenic	14	3.3	EPA 200.8	9-28-18	9-28-18	
Cadmium	ND	4.4	EPA 200.8	9-28-18	9-28-18	
Chromium	ND	11	EPA 200.8	9-28-18	9-28-18	
Lead	ND	1.1	EPA 200.8	9-28-18	9-28-18	
Mercury	ND	0.50	EPA 7470A	9-28-18	9-28-18	

Laboratory Reference: 1809-277 Project: 17-06520-000

TOTAL METALS EPA 200.8/7470A

σ.mo. αg/2 (ppε/)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW5					
Laboratory ID:	09-277-05					
Arsenic	ND	3.3	EPA 200.8	9-28-18	9-28-18	
Cadmium	ND	4.4	EPA 200.8	9-28-18	9-28-18	
Chromium	ND	11	EPA 200.8	9-28-18	9-28-18	
Lead	1.9	1.1	EPA 200.8	9-28-18	9-28-18	
Mercury	ND	0.50	EPA 7470A	9-28-18	9-28-18	
Client ID:	MW6					
Laboratory ID:	09-277-06					
Arsenic	4.5	3.3	EPA 200.8	9-28-18	9-28-18	
Cadmium	ND	4.4	EPA 200.8	9-28-18	9-28-18	
Chromium	ND	11	EPA 200.8	9-28-18	9-28-18	
Lead	ND	1.1	EPA 200.8	9-28-18	9-28-18	
Mercury	ND	0.50	EPA 7470A	9-28-18	9-28-18	
•						
Client ID:	MW7					
Laboratory ID:	09-277-07					
Arsenic	5.5	3.3	EPA 200.8	9-28-18	9-28-18	
Cadmium	ND	4.4	EPA 200.8	9-28-18	9-28-18	
Chromium	ND	11	EPA 200.8	9-28-18	9-28-18	
Lead	ND	1.1	EPA 200.8	9-28-18	9-28-18	
Mercury	ND	0.50	EPA 7470A	9-28-18	9-28-18	
Client ID:	MW8					
Laboratory ID:	09-277-08					
Arsenic	ND	3.3	EPA 200.8	9-28-18	9-28-18	
Cadmium	ND	4.4	EPA 200.8	9-28-18	9-28-18	
Chromium	ND	11	EPA 200.8	9-28-18	9-28-18	
Lead	ND	1.1	EPA 200.8	9-28-18	9-28-18	
Mercury	ND	0.50	EPA 7470A	9-28-18	9-28-18	

Date of Report: October 5, 2018

Samples Submitted: September 27, 2018

Laboratory Reference: 1809-277 Project: 17-06520-000

TOTAL METALS EPA 200.8/7470A

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW9					
Laboratory ID:	09-277-09					
Arsenic	3.6	3.3	EPA 200.8	9-28-18	9-28-18	
Cadmium	ND	4.4	EPA 200.8	9-28-18	9-28-18	
Chromium	ND	11	EPA 200.8	9-28-18	9-28-18	
Lead	ND	1.1	EPA 200.8	9-28-18	9-28-18	
Mercury	ND	0.50	EPA 7470A	9-28-18	9-28-18	

Laboratory Reference: 1809-277 Project: 17-06520-000

TOTAL METALS EPA 200.8/7470A **QUALITY CONTROL**

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0928WM1					
Arsenic	ND	3.3	EPA 200.8	9-28-18	9-28-18	
Cadmium	ND	4.4	EPA 200.8	9-28-18	9-28-18	
Chromium	ND	11	EPA 200.8	9-28-18	9-28-18	
Lead	ND	1.1	EPA 200.8	9-28-18	9-28-18	
Laboratory ID:	MB0928W1					
Mercury	ND	0.50	EPA 7470A	9-28-18	9-28-18	

					Source	Pe	rcent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Rec	overy	Limits	RPD	Limit	Flags
DUPLICATE											
Laboratory ID:	09-24	14-01									
	ORIG	DUP									
Arsenic	ND	ND	NA	NA		I	NA	NA	NA	20	
Cadmium	ND	ND	NA	NA		ı	NA	NA	NA	20	
Chromium	ND	ND	NA	NA		ı	NA	NA	NA	20	
Lead	ND	ND	NA	NA		l	NA	NA	NA	20	
Laboratory ID:	09-27	77-01									
Mercury	ND	ND	NA	NA		l	NA	NA	NA	20	
MATRIX SPIKES											
Laboratory ID:	09-24	14-01									
Laboratory ID.	MS	MSD	MS	MSD		MS	MSD				
Arsenic	222	219	222	222	ND	100	99	75-125	1	20	
Cadmium	211	204	222	222	ND	95	92	75-125	3	20	
Chromium	202	199	222	222	ND	91	90	75-125	2	20	
Lead	212	210	222	222	ND	96	95	75-125	1	20	
Laboratory ID:	09-27	77-01									
Mercury	12.1	12.0	12.5	12.5	ND	96	96	75-125	0	20	

Laboratory Reference: 1809-277

Project: 17-06520-000

cPAHs EPA 8270D/SIM

-				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW1					
Laboratory ID:	09-277-01					
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Chrysene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	73	21 - 110				
Pyrene-d10	89	19 - 111				
Terphenyl-d14	111	32 - 137				

Laboratory Reference: 1809-277 Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW2					
Laboratory ID:	09-277-02					
Benzo[a]anthracene	ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
Chrysene	ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
Benzo[b]fluoranthene	ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
Benzo(j,k)fluoranthene	ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
Benzo[a]pyrene	ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
Dibenz[a,h]anthracene	ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	59	21 - 110				
Pyrene-d10	78	19 - 111				
Terphenyl-d14	97	32 - 137				

Laboratory Reference: 1809-277

Project: 17-06520-000

cPAHs EPA 8270D/SIM

· ·				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW3					·
Laboratory ID:	09-277-03					
Benzo[a]anthracene	ND	0.0095	EPA 8270D/SIM	9-28-18	9-28-18	
Chrysene	ND	0.0095	EPA 8270D/SIM	9-28-18	9-28-18	
Benzo[b]fluoranthene	ND	0.0095	EPA 8270D/SIM	9-28-18	9-28-18	
Benzo(j,k)fluoranthene	ND	0.0095	EPA 8270D/SIM	9-28-18	9-28-18	
Benzo[a]pyrene	ND	0.0095	EPA 8270D/SIM	9-28-18	9-28-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0095	EPA 8270D/SIM	9-28-18	9-28-18	
Dibenz[a,h]anthracene	ND	0.0095	EPA 8270D/SIM	9-28-18	9-28-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	69	21 - 110				
Pyrene-d10	83	19 - 111				
Terphenyl-d14	100	32 - 137				

Laboratory Reference: 1809-277

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW4					
Laboratory ID:	09-277-04					
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Chrysene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	81	21 - 110				
Pyrene-d10	85	19 - 111				
Terphenyl-d14	134	32 - 137				

Laboratory Reference: 1809-277 Project: 17-06520-000

cPAHs EPA 8270D/SIM

J				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW5					
Laboratory ID:	09-277-05					
Benzo[a]anthracene	ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
Chrysene	ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
Benzo[b]fluoranthene	ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
Benzo(j,k)fluoranthene	ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
Benzo[a]pyrene	ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
Dibenz[a,h]anthracene	ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	91	21 - 110				
Pyrene-d10	86	19 - 111				
Terphenyl-d14	176	32 - 137				Q

Laboratory Reference: 1809-277 Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW6					
Laboratory ID:	09-277-06					
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Chrysene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	71	21 - 110				
Pyrene-d10	89	19 - 111				
Terphenyl-d14	122	32 - 137				
respirently a 14	122	0 <u>2</u> 101				

Laboratory Reference: 1809-277

Project: 17-06520-000

cPAHs EPA 8270D/SIM

			Date	Date	
Result	PQL	Method	Prepared	Analyzed	Flags
MW7					
09-277-07					
ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
Percent Recovery	Control Limits				
62	21 - 110				
86	19 - 111				
105	32 - 137				
	MW7 09-277-07 ND ND ND ND ND ND ND ND Percent Recovery 62 86	MW7 09-277-07 ND 0.0094 Percent Recovery Control Limits 62 21 - 110 86 19 - 111	MW7 09-277-07 ND 0.0094 EPA 8270D/SIM Percent Recovery Control Limits 62 21 - 110 86 19 - 111	Result PQL Method Prepared MW7	Result PQL Method Prepared Analyzed MW7 9-277-07 9-277-07 9-27-18 9-27-18 ND 0.0094 EPA 8270D/SIM 9-27-18 9-27-18 Percent Recovery Control Limits 62 21 - 110 86 19 - 111 111 111

Laboratory Reference: 1809-277

Project: 17-06520-000

cPAHs EPA 8270D/SIM

· ·				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW8					_
Laboratory ID:	09-277-08					
Benzo[a]anthracene	ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
Chrysene	ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
Benzo[b]fluoranthene	ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
Benzo(j,k)fluoranthene	ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
Benzo[a]pyrene	ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
Dibenz[a,h]anthracene	ND	0.0094	EPA 8270D/SIM	9-27-18	9-27-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	57	21 - 110				
Pyrene-d10	71	19 - 111				
Terphenyl-d14	86	32 - 137				

Laboratory Reference: 1809-277

Project: 17-06520-000

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW9					
Laboratory ID:	09-277-09					
Benzo[a]anthracene	ND	0.0098	EPA 8270D/SIM	9-27-18	9-27-18	
Chrysene	ND	0.0098	EPA 8270D/SIM	9-27-18	9-27-18	
Benzo[b]fluoranthene	ND	0.0098	EPA 8270D/SIM	9-27-18	9-27-18	
Benzo(j,k)fluoranthene	ND	0.0098	EPA 8270D/SIM	9-27-18	9-27-18	
Benzo[a]pyrene	ND	0.0098	EPA 8270D/SIM	9-27-18	9-27-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0098	EPA 8270D/SIM	9-27-18	9-27-18	
Dibenz[a,h]anthracene	ND	0.0098	EPA 8270D/SIM	9-27-18	9-27-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	77	21 - 110				
Pyrene-d10	82	19 - 111				
Terphenyl-d14	112	32 - 137				

Date of Report: October 5, 2018

Samples Submitted: September 27, 2018 Laboratory Reference: 1809-277

Project: 17-06520-000

CPAHS EPA 8270D/SIM
METHOD BLANK QUALITY CONTROL

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0927W1					
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Chrysene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	9-27-18	9-27-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	105	21 - 110				
Pyrene-d10	86	19 - 111				
Terphenyl-d14	143	32 - 137				Q

Laboratory Reference: 1809-277

Project: 17-06520-000

cPAHs EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB0928W1					
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	9-28-18	9-28-18	
Chrysene	ND	0.010	EPA 8270D/SIM	9-28-18	9-28-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	9-28-18	9-28-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	9-28-18	9-28-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	9-28-18	9-28-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	9-28-18	9-28-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	9-28-18	9-28-18	
Surrogate:	Percent Recovery	Control Limits				_
2-Fluorobiphenyl	68	21 - 110				
Pyrene-d10	83	19 - 111				
Terphenyl-d14	100	32 - 137				

Laboratory Reference: 1809-277 Project: 17-06520-000

cPAHs EPA 8270D/SIM **SB/SBD QUALITY CONTROL**

					Pe	rcent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Rec	overy	Limits	RPD	Limit	Flags
SPIKE BLANKS										
Laboratory ID:	SB09	27W1								
	SB	SBD	SB	SBD	SB	SBD				
Naphthalene	0.430	0.341	0.500	0.500	86	68	28 - 109	23	38	
Acenaphthylene	0.478	0.397	0.500	0.500	96	79	37 - 111	19	26	
Acenaphthene	0.496	0.434	0.500	0.500	99	87	41 - 113	13	33	
Fluorene	0.489	0.467	0.500	0.500	98	93	47 - 114	5	23	
Phenanthrene	0.465	0.461	0.500	0.500	93	92	50 - 113	1	18	
Anthracene	0.468	0.465	0.500	0.500	94	93	50 - 117	1	18	
Fluoranthene	0.490	0.489	0.500	0.500	98	98	52 - 120	0	15	
Pyrene	0.589	0.488	0.500	0.500	118	98	51 - 128	19	31	
Benzo[a]anthracene	0.541	0.539	0.500	0.500	108	108	57 - 127	0	15	
Chrysene	0.507	0.514	0.500	0.500	101	103	51 - 120	1	15	
Benzo[b]fluoranthene	0.502	0.506	0.500	0.500	100	101	54 - 124	1	17	
Benzo(j,k)fluoranthene	0.527	0.539	0.500	0.500	105	108	50 - 127	2	18	
Benzo[a]pyrene	0.503	0.509	0.500	0.500	101	102	50 - 120	1	16	
Indeno(1,2,3-c,d)pyrene	0.491	0.492	0.500	0.500	98	98	46 - 132	0	20	
Dibenz[a,h]anthracene	0.489	0.515	0.500	0.500	98	103	49 - 129	5	18	
Benzo[g,h,i]perylene	0.468	0.486	0.500	0.500	94	97	45 - 130	4	19	
Surrogate:										
2-Fluorobiphenyl					78	62	21 - 110			
Pyrene-d10					88	90	19 - 111			
Terphenyl-d14					118	108	32 - 137			

Laboratory Reference: 1809-277 Project: 17-06520-000

cPAHs EPA 8270D/SIM **SB/SBD QUALITY CONTROL**

					Per	cent	Recovery		RPD	
Analyte	Re	sult	Spike	Level	Rec	overy	Limits	RPD	Limit	Flags
SPIKE BLANKS										
Laboratory ID:	SB09	28W1								
	SB	SBD	SB	SBD	SB	SBD				
Naphthalene	0.373	0.392	0.500	0.500	75	78	28 - 109	5	38	
Acenaphthylene	0.430	0.437	0.500	0.500	86	87	37 - 111	2	26	
Acenaphthene	0.463	0.448	0.500	0.500	93	90	41 - 113	3	33	
Fluorene	0.429	0.439	0.500	0.500	86	88	47 - 114	2	23	
Phenanthrene	0.407	0.431	0.500	0.500	81	86	50 - 113	6	18	
Anthracene	0.409	0.436	0.500	0.500	82	87	50 - 117	6	18	
Fluoranthene	0.431	0.460	0.500	0.500	86	92	52 - 120	7	15	
Pyrene	0.451	0.453	0.500	0.500	90	91	51 - 128	0	31	
Benzo[a]anthracene	0.480	0.504	0.500	0.500	96	101	57 - 127	5	15	
Chrysene	0.453	0.488	0.500	0.500	91	98	51 - 120	7	15	
Benzo[b]fluoranthene	0.478	0.499	0.500	0.500	96	100	54 - 124	4	17	
Benzo(j,k)fluoranthene	0.451	0.491	0.500	0.500	90	98	50 - 127	8	18	
Benzo[a]pyrene	0.452	0.477	0.500	0.500	90	95	50 - 120	5	16	
Indeno(1,2,3-c,d)pyrene	0.454	0.484	0.500	0.500	91	97	46 - 132	6	20	
Dibenz[a,h]anthracene	0.446	0.495	0.500	0.500	89	99	49 - 129	10	18	
Benzo[g,h,i]perylene	0.458	0.481	0.500	0.500	92	96	45 - 130	5	19	
Surrogate:										
2-Fluorobiphenyl					73	87	21 - 110			
Pyrene-d10					81	86	19 - 111			
Terphenyl-d14					104	104	32 - 137			



Data Qualifiers and Abbreviations

- A Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
- B The analyte indicated was also found in the blank sample.
- C The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
- E The value reported exceeds the quantitation range and is an estimate.
- F Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
- H The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
- I Compound recovery is outside of the control limits.
- J The value reported was below the practical quantitation limit. The value is an estimate.
- K Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
- L The RPD is outside of the control limits.
- M Hydrocarbons in the gasoline range are impacting the diesel range result.
- M1 Hydrocarbons in the gasoline range (toluene-naphthalene) are present in the sample.
- N Hydrocarbons in the lube oil range are impacting the diesel range result.
- N1 Hydrocarbons in diesel range are impacting lube oil range results.
- O Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
- P The RPD of the detected concentrations between the two columns is greater than 40.
- Q Surrogate recovery is outside of the control limits.
- S Surrogate recovery data is not available due to the necessary dilution of the sample.
- T The sample chromatogram is not similar to a typical ______.
- U The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- U1 The practical quantitation limit is elevated due to interferences present in the sample.
- V Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
- W Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
- X Sample extract treated with a mercury cleanup procedure.
- X1- Sample extract treated with a sulfuric acid/silica gel cleanup procedure.
- Y The calibration verification for this analyte exceeded the 20% drift specified in method 8260C, and therefore the reported result should be considered an estimate. The overall performance of the calibration verification standard met the acceptance criteria of the method.

7 -

ND - Not Detected at PQL

PQL - Practical Quantitation Limit

RPD - Relative Percent Difference



MIR	OnSite		
	Environmental	inc.	

Chain of Custody

	1	1
Page	of_	1

Analytical Laboratory Testing Services	Turnaround Request (in working days)		Lak	oor	ator	y N	umb	er:		l	19	- 7	21	1							
14648 NE 95th Street • Redmond, WA 98052 Phone: (425) 883-3881 • www.onsite-env.com	(Check One)			T									_					Z	2		
Company: Helyera Env. Cons.	Same Day 1 Day				(ar							m	Organophosphorus Pesticides 8270D/SIM	1A	.,			101	-		
Project Number: 17 - 0 6520 - 6 00	2 Days 3 Days				ean-r		0	0		_		3081E	des 8%	s 815			4	1			
Project Name: Pacific Pavk	Standard (7 Days) (TPH analysis 5 Days)	SIL			/ SG Clean-up)		Halogenated Volatiles 8260C	ers Only)	Semivolatiles 8270D/SIM (with low-level PAHs)	w-level)	- 1	Pesticides 8081B	Pesticio	Chlorinated Acid Herbicides 8151A			TCLP Metals HEM (oil and grease) 1664A	0			
Project Manager:	(II I analysis s zaya)	itaine	1	X	Acid /		olatile	(Wate	270D PAHs	IM (lo	女	e Pes	orus	sid He	etals	etals	grease	-	2		
Sampled by:	(other)	of Containers	GIO	NWTPH-Gx/BTEX	NWTPH-GX	Volatiles 8260C	Ited V	EDB EPA 8011 (Waters	tiles 8-level	S/00/	82A	Organochlorine	hospi	ted A	Total RCRA Metals	Total MTCA Metals	etals	5	2		ture
B, Bland, G. Ifthey	(Garai)	Number	NWTPH-HCID	TPH-	NWTPH-GX	atiles	ogena	B EPA	mivola th low	Hs 82.	PCBs 8082A	ganoc	ganop	lorina	tal RC	tal MT	TCLP Metals HEM (oil and	1	LA MS		% Moisture
Lab ID Sample Identification	Outriples .		NN N	Š Z	A 2	Non	Hal	8	Sel	PA	6	ŏ	ō	ਠ	2	5	= =		4		%
1 MWI	9/26/8 11:00 water	W,			X	X ?	X									X		X			
2 MWZ	1 12:35	1			X	t	X	-	-	-						X	-	X		-	
3 MW3	14:40				X	K)	<		-							X	-	\\ \\ \\ \\ \\ \\ \\ \\ \\ \\ \\ \\ \\		-	
4 MW4	13:45				XX	7	X									X	-		V		\vdash
5 MWS	13:45				X	X	x									X		2			\mathbf{H}
6 MW6	12:35				X	X	X									X		-	0	-	\vdash
7 MW7	10:35				X	X	X									X		-		-	
8 MW8	11:35				X,	X	X									t		-	7	-	H
9 MW9	V 15:00 V	V			X	T	t									X	nes.		x well		
10 Trip Blank	9/26/18 - Water	3					X				1					100	MAR	a	Marie	(
Signature	Company			Date	-		Time		C	omme	nts/S	pecia	I Inst	ructio	ons		٨				
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14648 NE 95th Street, Redmond, WA 98052 • (425) 883-3881

October 17, 2018

Mark Ewbank Herrera Environmental Consultants, Inc. 2200 6th Avenue, Suite 1100 Seattle, WA 98121

Re: Analytical Data for Project 17-06520-000

Laboratory Reference No. 1810-129

Dear Mark:

Enclosed are the analytical results and associated quality control data for samples submitted on October 9, 2018.

The standard policy of OnSite Environmental, Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

David Baumeister Project Manager

Enclosures

Project: 17-06520-000

Case Narrative

Samples were collected on October 9, 2018 and received by the laboratory on October 9, 2018. They were maintained at the laboratory at a temperature of 2°C to 6°C.

Please note that any and all soil sample results are reported on a dry-weight basis, unless otherwise noted below.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

Project: 17-06520-000

GASOLINE RANGE ORGANICS NWTPH-Gx

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	SW1					
Laboratory ID:	10-129-01					
Gasoline	ND	100	NWTPH-Gx	10-9-18	10-9-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	92	66-117				
Client ID:	SW2					
Laboratory ID:	10-129-02					
Gasoline	ND	100	NWTPH-Gx	10-9-18	10-9-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	91	66-117				
Client ID:	SW3					
Laboratory ID:	10-129-03					
Gasoline	ND	100	NWTPH-Gx	10-9-18	10-9-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	93	66-117				
Client ID:	SW4					
Laboratory ID:	10-129-04					
Gasoline	ND	100	NWTPH-Gx	10-9-18	10-9-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	92	66-117				

Date of Report: October 17, 2018 Samples Submitted: October 9, 2018 Laboratory Reference: 1810-129 Project: 17-06520-000

GASOLINE RANGE ORGANICS NWTPH-Gx QUALITY CONTROL

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB1009W1					
Gasoline	ND	100	NWTPH-Gx	10-9-18	10-9-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	91	66-117				

Analyte	Res	sult	Spike	Level	Source Result	Perce Recov		Recovery Limits	RPD	RPD Limit	Flags
DUPLICATE			•								
Laboratory ID:	10-12	29-01									
	ORIG	DUP									
Gasoline	ND	ND	NA	NA		NA		NA	NA	30	
Surrogate:											
Fluorobenzene						92	92	66-117			

Date of Report: October 17, 2018 Samples Submitted: October 9, 2018 Laboratory Reference: 1810-129 Project: 17-06520-000

DIESEL AND HEAVY OIL RANGE ORGANICS NWTPH-Dx

5 (II)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	SW1					
Laboratory ID:	10-129-01					
Diesel Range Organics	ND	0.25	NWTPH-Dx	10-11-18	10-11-18	
Lube Oil Range Organics	ND	0.41	NWTPH-Dx	10-11-18	10-11-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	80	50-150				
Client ID:	SW2					
Laboratory ID:	10-129-02					
Diesel Range Organics	ND	0.26	NWTPH-Dx	10-11-18	10-11-18	
Lube Oil Range Organics	ND	0.42	NWTPH-Dx	10-11-18	10-11-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	78	50-150				
Client ID:	SW3					
Laboratory ID:	10-129-03					
Diesel Range Organics	ND	0.26	NWTPH-Dx	10-11-18	10-11-18	
Lube Oil Range Organics	ND	0.41	NWTPH-Dx	10-11-18	10-11-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	86	50-150				
Client ID:	SW4					
Laboratory ID:	10-129-04					
Diesel Range Organics	ND	0.25	NWTPH-Dx	10-11-18	10-11-18	
Lube Oil Range Organics	ND	0.40	NWTPH-Dx	10-11-18	10-11-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	83	50-150				
· ·						

Project: 17-06520-000

DIESEL AND HEAVY OIL RANGE ORGANICS NWTPH-Dx QUALITY CONTROL

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						_
Laboratory ID:	MB1011W1					
Diesel Range Organics	ND	0.25	NWTPH-Dx	10-11-18	10-11-18	
Lube Oil Range Organics	ND	0.40	NWTPH-Dx	10-11-18	10-11-18	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	84	50-150				

					Source	Perce	ent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Recov	ery	Limits	RPD	Limit	Flags
DUPLICATE											
Laboratory ID:	10-12	29-01									
	ORIG	DUP									
Diesel Range	ND	ND	NA	NA		NA		NA	NA	NA	
Lube Oil Range	ND	ND	NA	NA		NA		NA	NA	NA	
Surrogate:											
o-Terphenyl						80	80	50-150			

Date of Report: October 17, 2018 Samples Submitted: October 9, 2018 Laboratory Reference: 1810-129 Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	SW1					
Laboratory ID:	10-129-01					
Dichlorodifluoromethane	ND	0.30	EPA 8260C	10-11-18	10-11-18	
Chloromethane	ND	1.0	EPA 8260C	10-11-18	10-11-18	
Vinyl Chloride	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Bromomethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Chloroethane	ND	1.0	EPA 8260C	10-11-18	10-11-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Acetone	ND	5.0	EPA 8260C	10-11-18	10-11-18	
Iodomethane	ND	1.4	EPA 8260C	10-11-18	10-11-18	
Carbon Disulfide	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Methylene Chloride	ND	2.0	EPA 8260C	10-11-18	10-11-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Vinyl Acetate	ND	1.0	EPA 8260C	10-11-18	10-11-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
2-Butanone	ND	5.0	EPA 8260C	10-11-18	10-11-18	
Bromochloromethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Chloroform	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Benzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Trichloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Dibromomethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Bromodichloromethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	10-11-18	10-11-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	10-11-18	10-11-18	
Toluene	ND	1.0	EPA 8260C	10-11-18	10-11-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	10-11-18	10-11-18	

Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 2 of 2

Amalada	Decel	DOL	Mada al	Date	Date	F1
Analyte Client ID:	Result SW1	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	10-129-01	0.00	EDA 02000	10 11 10	40.44.40	
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Tetrachloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
2-Hexanone	ND	2.0	EPA 8260C	10-11-18	10-11-18	
Dibromochloromethane	ND	0.30	EPA 8260C	10-11-18	10-11-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Chlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1,1,2-Tetrachloroethane	ND	0.28	EPA 8260C	10-11-18	10-11-18	
Ethylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
m,p-Xylene	ND	0.40	EPA 8260C	10-11-18	10-11-18	
o-Xylene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Styrene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Bromoform	ND	2.0	EPA 8260C	10-11-18	10-11-18	
Isopropylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Bromobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
n-Propylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
n-Butylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2-Dibromo-3-chloropropane		1.9	EPA 8260C	10-11-18	10-11-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	10-11-18	10-11-18	
Naphthalene	ND	1.0	EPA 8260C	10-11-18	10-11-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Surrogate:	Percent Recovery	Control Limits	2 32000			

 Dibromofluoromethane
 97
 75-127

 Toluene-d8
 98
 80-127

 4-Bromofluorobenzene
 95
 78-125



Date of Report: October 17, 2018 Samples Submitted: October 9, 2018 Laboratory Reference: 1810-129 Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	SW2					
Laboratory ID:	10-129-02					
Dichlorodifluoromethane	ND	0.30	EPA 8260C	10-11-18	10-11-18	
Chloromethane	ND	1.0	EPA 8260C	10-11-18	10-11-18	
Vinyl Chloride	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Bromomethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Chloroethane	ND	1.0	EPA 8260C	10-11-18	10-11-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Acetone	ND	5.0	EPA 8260C	10-11-18	10-11-18	
Iodomethane	ND	1.4	EPA 8260C	10-11-18	10-11-18	
Carbon Disulfide	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Methylene Chloride	ND	2.0	EPA 8260C	10-11-18	10-11-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Vinyl Acetate	ND	1.0	EPA 8260C	10-11-18	10-11-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
2-Butanone	ND	5.0	EPA 8260C	10-11-18	10-11-18	
Bromochloromethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Chloroform	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Benzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Trichloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Dibromomethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Bromodichloromethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	10-11-18	10-11-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	10-11-18	10-11-18	
Toluene	ND	1.0	EPA 8260C	10-11-18	10-11-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	10-11-18	10-11-18	

Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	SW2					
Laboratory ID:	10-129-02					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Tetrachloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
2-Hexanone	ND	2.0	EPA 8260C	10-11-18	10-11-18	
Dibromochloromethane	ND	0.30	EPA 8260C	10-11-18	10-11-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Chlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1,1,2-Tetrachloroethane	ND	0.28	EPA 8260C	10-11-18	10-11-18	
Ethylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
m,p-Xylene	ND	0.40	EPA 8260C	10-11-18	10-11-18	
o-Xylene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Styrene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Bromoform	ND	2.0	EPA 8260C	10-11-18	10-11-18	
Isopropylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Bromobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
n-Propylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
n-Butylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2-Dibromo-3-chloropropane	ND	1.9	EPA 8260C	10-11-18	10-11-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	10-11-18	10-11-18	
Naphthalene	ND	1.0	EPA 8260C	10-11-18	10-11-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Surrogate:	Percent Recovery	Control Limits				

Surrogate: Percent Recovery Control Limits
Dibromofluoromethane 99 75-127
Toluene-d8 100 80-127
4-Bromofluorobenzene 98 78-125



Date of Report: October 17, 2018 Samples Submitted: October 9, 2018 Laboratory Reference: 1810-129 Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	SW3					
Laboratory ID:	10-129-03					
Dichlorodifluoromethane	ND	0.30	EPA 8260C	10-11-18	10-11-18	
Chloromethane	ND	1.0	EPA 8260C	10-11-18	10-11-18	
Vinyl Chloride	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Bromomethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Chloroethane	ND	1.0	EPA 8260C	10-11-18	10-11-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Acetone	ND	5.0	EPA 8260C	10-11-18	10-11-18	
Iodomethane	ND	1.4	EPA 8260C	10-11-18	10-11-18	
Carbon Disulfide	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Methylene Chloride	ND	2.0	EPA 8260C	10-11-18	10-11-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Vinyl Acetate	ND	1.0	EPA 8260C	10-11-18	10-11-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
2-Butanone	ND	5.0	EPA 8260C	10-11-18	10-11-18	
Bromochloromethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Chloroform	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Benzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Trichloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Dibromomethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Bromodichloromethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	10-11-18	10-11-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	10-11-18	10-11-18	
Toluene	1.1	1.0	EPA 8260C	10-11-18	10-11-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	10-11-18	10-11-18	

Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	SW3					
Laboratory ID:	10-129-03					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Tetrachloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
2-Hexanone	ND	2.0	EPA 8260C	10-11-18	10-11-18	
Dibromochloromethane	ND	0.30	EPA 8260C	10-11-18	10-11-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Chlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1,1,2-Tetrachloroethane	ND	0.28	EPA 8260C	10-11-18	10-11-18	
Ethylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
m,p-Xylene	ND	0.40	EPA 8260C	10-11-18	10-11-18	
o-Xylene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Styrene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Bromoform	ND	2.0	EPA 8260C	10-11-18	10-11-18	
Isopropylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Bromobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
n-Propylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
n-Butylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2-Dibromo-3-chloropropane	ND	1.9	EPA 8260C	10-11-18	10-11-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	10-11-18	10-11-18	
Naphthalene	ND	1.0	EPA 8260C	10-11-18	10-11-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Surrogate:	Percent Recovery	Control Limits				

Surrogate: Percent Recovery Control Limits
Dibromofluoromethane 98 75-127
Toluene-d8 99 80-127
4-Bromofluorobenzene 98 78-125



Date of Report: October 17, 2018 Samples Submitted: October 9, 2018 Laboratory Reference: 1810-129 Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 1 of 2

Ğ				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	SW4					
Laboratory ID:	10-129-04					
Dichlorodifluoromethane	ND	0.30	EPA 8260C	10-11-18	10-11-18	
Chloromethane	ND	1.0	EPA 8260C	10-11-18	10-11-18	
Vinyl Chloride	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Bromomethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Chloroethane	ND	1.0	EPA 8260C	10-11-18	10-11-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Acetone	ND	5.0	EPA 8260C	10-11-18	10-11-18	
Iodomethane	ND	1.4	EPA 8260C	10-11-18	10-11-18	
Carbon Disulfide	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Methylene Chloride	ND	2.0	EPA 8260C	10-11-18	10-11-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Vinyl Acetate	ND	1.0	EPA 8260C	10-11-18	10-11-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
2-Butanone	ND	5.0	EPA 8260C	10-11-18	10-11-18	
Bromochloromethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Chloroform	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Benzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Trichloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Dibromomethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Bromodichloromethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	10-11-18	10-11-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	10-11-18	10-11-18	
Toluene	ND	1.0	EPA 8260C	10-11-18	10-11-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	10-11-18	10-11-18	

Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 2 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	SW4					
Laboratory ID:	10-129-04					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Tetrachloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
2-Hexanone	ND	2.0	EPA 8260C	10-11-18	10-11-18	
Dibromochloromethane	ND	0.30	EPA 8260C	10-11-18	10-11-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Chlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1,1,2-Tetrachloroethane	ND	0.28	EPA 8260C	10-11-18	10-11-18	
Ethylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
m,p-Xylene	ND	0.40	EPA 8260C	10-11-18	10-11-18	
o-Xylene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Styrene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Bromoform	ND	2.0	EPA 8260C	10-11-18	10-11-18	
Isopropylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Bromobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
n-Propylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
n-Butylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2-Dibromo-3-chloropropane	ND	1.9	EPA 8260C	10-11-18	10-11-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	10-11-18	10-11-18	
Naphthalene	ND	1.0	EPA 8260C	10-11-18	10-11-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Surrogate:	Percent Recovery	Control Limits				

Surrogate: Percent Recovery Control Limits
Dibromofluoromethane 101 75-127
Toluene-d8 100 80-127
4-Bromofluorobenzene 96 78-125



Date of Report: October 17, 2018 Samples Submitted: October 9, 2018 Laboratory Reference: 1810-129 Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 1 of 2

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	Trip Blank					
Laboratory ID:	10-129-05					
Dichlorodifluoromethane	ND	0.30	EPA 8260C	10-11-18	10-11-18	
Chloromethane	ND	1.0	EPA 8260C	10-11-18	10-11-18	
Vinyl Chloride	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Bromomethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Chloroethane	ND	1.0	EPA 8260C	10-11-18	10-11-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Acetone	ND	5.0	EPA 8260C	10-11-18	10-11-18	
Iodomethane	ND	1.4	EPA 8260C	10-11-18	10-11-18	
Carbon Disulfide	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Methylene Chloride	ND	2.0	EPA 8260C	10-11-18	10-11-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Vinyl Acetate	ND	1.0	EPA 8260C	10-11-18	10-11-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
2-Butanone	ND	5.0	EPA 8260C	10-11-18	10-11-18	
Bromochloromethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Chloroform	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Benzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Trichloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Dibromomethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Bromodichloromethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	10-11-18	10-11-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	10-11-18	10-11-18	
Toluene	ND	1.0	EPA 8260C	10-11-18	10-11-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	10-11-18	10-11-18	

Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C

page 2 of 2

Analyte				Date	Date	
	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	Trip Blank					
Laboratory ID:	10-129-05					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Tetrachloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
2-Hexanone	ND	2.0	EPA 8260C	10-11-18	10-11-18	
Dibromochloromethane	ND	0.30	EPA 8260C	10-11-18	10-11-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Chlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1,1,2-Tetrachloroethane	ND	0.28	EPA 8260C	10-11-18	10-11-18	
Ethylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
m,p-Xylene	ND	0.40	EPA 8260C	10-11-18	10-11-18	
o-Xylene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Styrene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Bromoform	ND	2.0	EPA 8260C	10-11-18	10-11-18	
Isopropylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Bromobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
n-Propylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
n-Butylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2-Dibromo-3-chloropropane	ND	1.9	EPA 8260C	10-11-18	10-11-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	10-11-18	10-11-18	
Naphthalene	ND	1.0	EPA 8260C	10-11-18	10-11-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Surrogate:	Percent Recovery					

Surrogate: Percent Recovery Control Limits
Dibromofluoromethane 100 75-127
Toluene-d8 102 80-127
4-Bromofluorobenzene 97 78-125

Date of Report: October 17, 2018 Samples Submitted: October 9, 2018 Laboratory Reference: 1810-129

Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C METHOD BLANK QUALITY CONTROL

page 1 of 2

·				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB1011W1					
Dichlorodifluoromethane	ND	0.30	EPA 8260C	10-11-18	10-11-18	
Chloromethane	ND	1.0	EPA 8260C	10-11-18	10-11-18	
Vinyl Chloride	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Bromomethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Chloroethane	ND	1.0	EPA 8260C	10-11-18	10-11-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Acetone	ND	5.0	EPA 8260C	10-11-18	10-11-18	
Iodomethane	ND	1.4	EPA 8260C	10-11-18	10-11-18	
Carbon Disulfide	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Methylene Chloride	ND	2.0	EPA 8260C	10-11-18	10-11-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Vinyl Acetate	ND	1.0	EPA 8260C	10-11-18	10-11-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
2-Butanone	ND	5.0	EPA 8260C	10-11-18	10-11-18	
Bromochloromethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Chloroform	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Benzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Trichloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Dibromomethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Bromodichloromethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	10-11-18	10-11-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	10-11-18	10-11-18	
Toluene	ND	1.0	EPA 8260C	10-11-18	10-11-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	10-11-18	10-11-18	

Date of Report: October 17, 2018 Samples Submitted: October 9, 2018 Laboratory Reference: 1810-129

Project: 17-06520-000

VOLATILE ORGANICS EPA 8260C METHOD BLANK QUALITY CONTROL

page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Allalyte	Result	FQL	Wethou	гтератец	Allalyzeu	i iags
Laboratory ID:	MB1011W1					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Tetrachloroethene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
2-Hexanone	ND	2.0	EPA 8260C	10-11-18	10-11-18	
Dibromochloromethane	ND	0.30	EPA 8260C	10-11-18	10-11-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Chlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1,1,2-Tetrachloroethane	ND	0.28	EPA 8260C	10-11-18	10-11-18	
Ethylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
m,p-Xylene	ND	0.40	EPA 8260C	10-11-18	10-11-18	
o-Xylene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Styrene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Bromoform	ND	2.0	EPA 8260C	10-11-18	10-11-18	
Isopropylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Bromobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	10-11-18	10-11-18	
n-Propylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
n-Butylbenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
1,2-Dibromo-3-chloropropane	ND	1.9	EPA 8260C	10-11-18	10-11-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Hexachlorobutadiene	ND	1.0	EPA 8260C	10-11-18	10-11-18	
Naphthalene	ND	1.0	EPA 8260C	10-11-18	10-11-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	10-11-18	10-11-18	
Surrogate:	Percent Perceyany					

Surrogate: Percent Recovery Control Limits
Dibromofluoromethane 98 75-127
Toluene-d8 98 80-127
4-Bromofluorobenzene 95 78-125

VOLATILE ORGANICS EPA 8260C SB/SBD QUALITY CONTROL

					Per	cent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Reco	overy	Limits	RPD	Limit	Flags
SPIKE BLANKS										
Laboratory ID:	SB10	11W1								
	SB	SBD	SB	SBD	SB	SBD				
1,1-Dichloroethene	11.7	11.9	10.0	10.0	117	119	62-129	2	15	
Benzene	11.0	11.4	10.0	10.0	110	114	77-127	4	15	
Trichloroethene	10.1	10.2	10.0	10.0	101	102	70-120	1	15	
Toluene	11.0	11.3	10.0	10.0	110	113	82-123	3	15	
Chlorobenzene	9.97	10.1	10.0	10.0	100	101	79-120	1	15	
Surrogate:										
Dibromofluoromethane					95	101	75-127			
Toluene-d8					97	102	80-127			
4-Bromofluorobenzene					94	98	78-125			

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	SW1					
Laboratory ID:	10-129-01					
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	10-12-18	10-12-18	
Chrysene	ND	0.010	EPA 8270D/SIM	10-12-18	10-12-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	10-12-18	10-12-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	10-12-18	10-12-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	10-12-18	10-12-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	10-12-18	10-12-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	10-12-18	10-12-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	57	21 - 110				
Pyrene-d10	67	19 - 111				
Terphenyl-d14	83	32 - 137				

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	SW2					
Laboratory ID:	10-129-02					
Benzo[a]anthracene	ND	0.0099	EPA 8270D/SIM	10-12-18	10-12-18	
Chrysene	ND	0.0099	EPA 8270D/SIM	10-12-18	10-12-18	
Benzo[b]fluoranthene	ND	0.0099	EPA 8270D/SIM	10-12-18	10-12-18	
Benzo(j,k)fluoranthene	ND	0.0099	EPA 8270D/SIM	10-12-18	10-12-18	
Benzo[a]pyrene	ND	0.0099	EPA 8270D/SIM	10-12-18	10-12-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0099	EPA 8270D/SIM	10-12-18	10-12-18	
Dibenz[a,h]anthracene	ND	0.0099	EPA 8270D/SIM	10-12-18	10-12-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	65	21 - 110				
Pyrene-d10	78	19 - 111				
Terphenyl-d14	92	32 - 137				

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	SW3					_
Laboratory ID:	10-129-03					
Benzo[a]anthracene	ND	0.0098	EPA 8270D/SIM	10-12-18	10-12-18	
Chrysene	ND	0.0098	EPA 8270D/SIM	10-12-18	10-12-18	
Benzo[b]fluoranthene	ND	0.0098	EPA 8270D/SIM	10-12-18	10-12-18	
Benzo(j,k)fluoranthene	ND	0.0098	EPA 8270D/SIM	10-12-18	10-12-18	
Benzo[a]pyrene	ND	0.0098	EPA 8270D/SIM	10-12-18	10-12-18	
Indeno(1,2,3-c,d)pyrene	ND	0.0098	EPA 8270D/SIM	10-12-18	10-12-18	
Dibenz[a,h]anthracene	ND	0.0098	EPA 8270D/SIM	10-12-18	10-12-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	57	21 - 110				
Pyrene-d10	68	19 - 111				
Terphenyl-d14	82	32 - 137				

cPAHs EPA 8270D/SIM

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	SW4					
Laboratory ID:	10-129-04					
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	10-12-18	10-12-18	
Chrysene	ND	0.010	EPA 8270D/SIM	10-12-18	10-12-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	10-12-18	10-12-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	10-12-18	10-12-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	10-12-18	10-12-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	10-12-18	10-12-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	10-12-18	10-12-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	60	21 - 110				
Pyrene-d10	72	19 - 111				
Terphenyl-d14	90	32 - 137				
Chrysene Benzo[b]fluoranthene Benzo(j,k)fluoranthene Benzo[a]pyrene Indeno(1,2,3-c,d)pyrene Dibenz[a,h]anthracene Surrogate: 2-Fluorobiphenyl Pyrene-d10	ND ND ND ND ND Percent Recovery 60 72	0.010 0.010 0.010 0.010 0.010 0.010 Control Limits 21 - 110 19 - 111	EPA 8270D/SIM EPA 8270D/SIM EPA 8270D/SIM EPA 8270D/SIM EPA 8270D/SIM	10-12-18 10-12-18 10-12-18 10-12-18 10-12-18	10-12-18 10-12-18 10-12-18 10-12-18 10-12-18	

cPAHs EPA 8270D/SIM **METHOD BLANK QUALITY CONTROL**

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Laboratory ID:	MB1012W1					
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	10-12-18	10-12-18	
Chrysene	ND	0.010	EPA 8270D/SIM	10-12-18	10-12-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	10-12-18	10-12-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	10-12-18	10-12-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	10-12-18	10-12-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	10-12-18	10-12-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	10-12-18	10-12-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	53	21 - 110				
Pyrene-d10	67	19 - 111				
Terphenyl-d14	81	32 - 137				

cPAHs EPA 8270D/SIM SB/SBD QUALITY CONTROL

3					Per	cent	Recovery		RPD	
Analyte	Re	sult	Spike	Level	Rece	overy	Limits	RPD	Limit	Flags
SPIKE BLANKS										
Laboratory ID:	SB10	12W1								
	SB	SBD	SB	SBD	SB	SBD				
Benzo[a]anthracene	0.490	0.495	0.500	0.500	98	99	57 - 127	1	15	
Chrysene	0.446	0.429	0.500	0.500	89	86	51 - 120	4	15	
Benzo[b]fluoranthene	0.457	0.496	0.500	0.500	91	99	54 - 124	8	17	
Benzo(j,k)fluoranthene	0.477	0.459	0.500	0.500	95	92	50 - 127	4	18	
Benzo[a]pyrene	0.460	0.452	0.500	0.500	92	90	50 - 120	2	16	
Indeno(1,2,3-c,d)pyrene	0.442	0.441	0.500	0.500	88	88	46 - 132	0	20	
Dibenz[a,h]anthracene	0.465	0.460	0.500	0.500	93	92	49 - 129	1	18	
Surrogate:										
2-Fluorobiphenyl					56	60	21 - 110			
Pyrene-d10					77	76	19 - 111			
Terphenyl-d14					90	91	32 - 137			

Date of Report: October 17, 2018 Samples Submitted: October 9, 2018 Laboratory Reference: 1810-129

Project: 17-06520-000

TOTAL METALS EPA 200.8/7470A

Matrix: Water Units: ug/L (ppb)

omio. ug/2 (ppb)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	SW1					
Laboratory ID:	10-129-01					
Arsenic	ND	3.3	EPA 200.8	10-16-18	10-16-18	
Cadmium	ND	4.4	EPA 200.8	10-16-18	10-16-18	
Chromium	ND	11	EPA 200.8	10-16-18	10-16-18	
Lead	ND	1.1	EPA 200.8	10-16-18	10-16-18	
Mercury	ND	0.50	EPA 7470A	10-11-18	10-11-18	
Client ID:	SW2					
Laboratory ID:	10-129-02					
Arsenic	ND	3.3	EPA 200.8	10-16-18	10-16-18	
Cadmium	ND	4.4	EPA 200.8	10-16-18	10-16-18	
Chromium	ND	11	EPA 200.8	10-16-18	10-16-18	
Lead	ND	1.1	EPA 200.8	10-16-18	10-16-18	
Mercury	ND	0.50	EPA 7470A	10-11-18	10-11-18	
,						
Client ID:	SW3					
Laboratory ID:	10-129-03					
Arsenic	ND	3.3	EPA 200.8	10-16-18	10-16-18	
Cadmium	ND	4.4	EPA 200.8	10-16-18	10-16-18	
Chromium	ND	11	EPA 200.8	10-16-18	10-16-18	
Lead	ND	1.1	EPA 200.8	10-16-18	10-16-18	
Mercury	ND	0.50	EPA 7470A	10-11-18	10-11-18	
Client ID:	SW4					
Laboratory ID:	10-129-04					
Arsenic	ND	3.3	EPA 200.8	10-16-18	10-16-18	
Cadmium	ND	4.4	EPA 200.8	10-16-18	10-16-18	
Chromium	ND	11	EPA 200.8	10-16-18	10-16-18	
Lead	ND	1.1	EPA 200.8	10-16-18	10-16-18	
Mercury	ND	0.50	EPA 7470A	10-11-18	10-11-18	

TOTAL METALS EPA 200.8/7470A **QUALITY CONTROL**

Matrix: Water Units: ug/L (ppb)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB1016WM1					
Arsenic	ND	3.3	EPA 200.8	10-16-18	10-16-18	
Cadmium	ND	4.4	EPA 200.8	10-16-18	10-16-18	
Chromium	ND	11	EPA 200.8	10-16-18	10-16-18	
Lead	ND	1.1	EPA 200.8	10-16-18	10-16-18	
Laboratory ID:	MB1011W1					
Mercury	ND	0.50	EPA 7470A	10-11-18	10-11-18	

					Source	Pe	rcent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Rec	overy	Limits	RPD	Limit	Flags
DUPLICATE											
Laboratory ID:	10-1	14-01									
	ORIG	DUP									
Arsenic	ND	ND	NA	NA		1	NA	NA	NA	20	
Cadmium	ND	ND	NA	NA		1	NA	NA	NA	20	
Chromium	ND	ND	NA	NA		ı	NA	NA	NA	20	
Lead	ND	ND	NA	NA		<u> </u>	NA	NA	NA	20	
Laboratory ID:	10-12	29-01									
Mercury	ND	ND	NA	NA		1	NA	NA	NA	20	
MATRIX SPIKES											
Laboratory ID:	10-1	14-01									
	MS	MSD	MS	MSD		MS	MSD				
Arsenic	226	229	222	222	ND	102	103	75-125	1	20	
Cadmium	219	226	222	222	ND	99	102	75-125	3	20	
Chromium	206	208	222	222	ND	93	94	75-125	1	20	
Lead	229	227	222	222	ND	103	102	75-125	1	20	
Laboratory ID:	10-12	29-01									
Mercury	12.4	12.3	12.5	12.5	ND	99	99	75-125	0	20	

HARDNESS EPA 200.7/SM 2340B

Matrix: Water

Units: mg eqt. CaCO3/L (ppm)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	SW1					
Laboratory ID:	10-129-01					
Hardness	62	1.0	200.7/SM 2340B	10-10-18	10-10-18	
Client ID:	SW2					
Laboratory ID:	10-129-02					
Hardness	51	1.0	200.7/SM 2340B	10-10-18	10-10-18	
Client ID:	SW3					
Laboratory ID:	10-129-03					
Hardness	51	1.0	200.7/SM 2340B	10-10-18	10-10-18	
Client ID:	SW4					
Laboratory ID:	10-129-04					
Hardness	45	1.0	200.7/SM 2340B	10-10-18	10-10-18	

HARDNESS EPA 200.7/SM 2340B **QUALITY CONTROL**

Matrix: Water

Units: mg eqt. CaCO3/L (ppm)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB1010WH2					
Hardness	ND	1.0	200.7/SM 2340B	10-10-18	10-10-18	

					C	D-		December		RPD	
Analyte	Res	sult	Spike	Level	Source Result		rcent covery	Recovery Limits	RPD	Limit	Flags
DUPLICATE											
Laboratory ID:	10-0	70-01									
	ORIG	DUP									
Hardness	90.4	88.0	١	1A	NA		NA	NA	3	20	
MATRIX SPIKES											
Laboratory ID:	10-0	70-01									
	MS	MSD	MS	MSD		MS	MSD				
Hardness	215	225	132	132	90.4	94	102	75-125	5	20	
SPIKE BLANK											
Laboratory ID:	SB101	I0WH2									
	S	B	5	SB	•		SB		•		•
Hardness	12	27	1	32	NA		96	80-120	NA	NA	



Data Qualifiers and Abbreviations

- A Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
- B The analyte indicated was also found in the blank sample.
- C The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
- E The value reported exceeds the quantitation range and is an estimate.
- F Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
- H The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
- I Compound recovery is outside of the control limits.
- J The value reported was below the practical quantitation limit. The value is an estimate.
- K Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
- L The RPD is outside of the control limits.
- M Hydrocarbons in the gasoline range are impacting the diesel range result.
- M1 Hydrocarbons in the gasoline range (toluene-naphthalene) are present in the sample.
- N Hydrocarbons in the lube oil range are impacting the diesel range result.
- N1 Hydrocarbons in diesel range are impacting lube oil range results.
- O Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
- P The RPD of the detected concentrations between the two columns is greater than 40.
- Q Surrogate recovery is outside of the control limits.
- S Surrogate recovery data is not available due to the necessary dilution of the sample.
- T The sample chromatogram is not similar to a typical ______.
- U The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- U1 The practical quantitation limit is elevated due to interferences present in the sample.
- V Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
- W Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
- X Sample extract treated with a mercury cleanup procedure.
- X1- Sample extract treated with a sulfuric acid/silica gel cleanup procedure.
- Y The calibration verification for this analyte exceeded the 20% drift specified in method 8260C, and therefore the reported result should be considered an estimate. The overall performance of the calibration verification standard met the acceptance criteria of the method.

7 -

ND - Not Detected at PQL

PQL - Practical Quantitation Limit

RPD - Relative Percent Difference





Chain of Custody

	Analytical Laboratory Testing Services 14648 NE 95th Street - Redmond, WA 98052	Turnaround Request (in working days)			Turnaround Request (in working days)							1	0 -	12	9									
Sampled	Hervera 1-06520-000 ane: Dacific Park anager: Mark Flahounk	Sam 2 Date		1 Day 3 Days	Number of Containers	NWTPH-HCID	/BTEX		Acid / SG Clean-up)	Volatiles 8260C Halogenated Volatiles 8260C	EDB EPA 8011 (Waters Only)	olatiles 8270D/SIM ow-level PAHs)	2270D/SIM (low-level)	Organochlorine Pesticides 8081B	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	TCLP Metals	HEM (oil and grease) 1664A	Hardwag5			% Moisture
Lab ID	Sample Identification	Sampled	Sampled	Matrix		NN N	NZ	Ž	Š :	F	EDI	Ser	PAIRs PAIRs PAIRs	Org	Org	S	Tota	Tota	107	H .	7	+		% %
7		10/9/18		Wolfy				V	V	V	+		V	+	-			V			V			
2	5WZ 5W3	-	10:00	9	12			V	V		-	1	V	-				V			/	-		
3		-	10:30	-	12			V	V	V	-		V	-	-			V		1	/			
4	SW4	-	11:00		12			V	V	V	+		V	-				V		1	/			
5	Trip Blank	Y	_	×	3		_		-	V	-			-						_		_		
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APPENDIX H

Data Quality Assurance Review Memorandum



Herrera Environmental Consultants, Inc.

Internal Memorandum

Date: September 12, 2018

To: Project File 17-06520-000

Copy To:

From: Gina Catarra

Subject: Data Quality Assurance Review of the Pacific Park/Dumpsite Data

This memorandum presents a review of data quality for 92 soil samples collected from the Pacific Park/Dumpsite property between February 20 and March 1, 2018. OnSite Environmental, Inc. (OnSite), of Redmond, Washington analyzed the samples for

- Gasoline-range petroleum hydrocarbons (TPH-G) by Ecology's NWPTH-Gx method
- Diesel- and lube oil-range petroleum hydrocarbons by Ecology's NWTPH-Dx method
- Benzene, toluene, ethylbenzene, and xylenes (BTEX) by EPA method 8021B
- Polycyclic aromatic hydrocarbons (PAHs) by EPA method 8270D/SIM
- Polychlorinated biphenyls (PCBs) by EPA method 8082A
- Total Model Toxics Control Act (MTCA) metals (arsenic, cadmium, chromium, lead, and mercury) by EPA 6010D/7471B
- Total chromium and lead by toxicity characteristic leaching procedure (TCLP) by EPA methods 1311/6010D.

Results for the following samples were validated.

Sample ID	Lab Ref. No.	Date Collected	Analyses
B-04_12.5	1802-233	2/21/18	HCID, cPAHs, metals
B-05_7.5	1802-233	2/20/18	HCID, Dx, cPAHs, PCBs, metals, TCLP Cr



Sample ID	Lab Ref. No.	Date Collected	Analyses
B-05_12.5	1802-233	2/20/18	HCID, cPAHs, metals
B-14_5	1802-233	2/21/18	HCID, cPAHs, metals
B-14_10	1802-233	2/21/18	HCID, cPAHs, metals
B-15_5	1802-233	2/21/18	HCID, cPAHs, metals
B-15_7.5	1802-233	2/21/18	HCID, cPAHs, metals
B-15_15	1802-233	2/21/18	HCID, cPAHs, metals
B-17_10	1802-233	2/20/18	HCID, cPAHs, metals
B-17_15	1802-233	2/20/18	HCID, cPAHs, metals
PP12-2	1802-234	2/21/18	HCID, Dx, cPAHs, PCBs, metals
PP12-7	1802-234	2/21/18	HCID, Dx, cPAHs, PCBs, metals, TCLP Pb
PP12-10	1802-234	2/21/18	Dx, cPAHs, PCBs, metals
PP13-10	1802-234	2/21/18	HCID, cPAHs, metals, TCLP Pb
PP15-6	1802-234	2/21/18	HCID, Dx, cPAHs, PCBs, metals
PP15-12	1802-234	2/21/18	HCID, cPAHs, metals, TCLP Pb
PP19-7	1802-234	2/21/18	HCID, Dx, cPAHs, metals
PP19-10	1802-234	2/21/18	HCID, cPAHs, metals, TCLP Pb
PP19-15	1802-234	2/21/18	HCID, cPAHs, metals
PP22-4	1802-234	2/21/18	HCID, cPAHs, metals
PP22-9	1802-234	2/21/18	HCID, cPAHs, metals
PP22-13	1802-234	2/21/18	HCID, cPAHs, metals, TCLP Pb
PP25-7	1802-234	2/21/18	HCID, cPAHs, metals
PP25-13	1802-234	2/21/18	HCID, Gx/BTEX, Dx, cPAHs, PCBs, metals, TCLP Pb
PP25-17	1802-234	2/21/18	GX/BTEX, Dx, cPAHs, PCBs, metals
PP26-11	1802-234	2/21/18	HCID, cPAHs, metals
PP26-17	1802-234	2/21/18	HCID, cPAHs, metals
PP34-8	1802-234	2/21/18	HCID, cPAHs, metals
PP34-15	1802-234	2/21/18	cPAHs
B-06_2.5	1802-249	2/22/18	HCID, Dx, cPAHs, PCBs, metals
B-06_10	1802-249	2/22/18	Dx, cPAHs, metals
B-06_12.5	1802-249	2/22/18	HCID, Dx, cPAHs, PCBs, metals
B-09_5	1802-249	2/22/18	HCID, cPAHs, metals
B-09_15	1802-249	2/22/18	HCID, cPAHs, metals
B-13_7.5	1802-249	2/23/18	HCID, cPAHs, metals
B-13_15	1802-249	2/23/18	HCID, cPAHs, metals
B16_2.5	1802-249	2/23/18	HCID, cPAHs, metals
B-16_7.5	1802-249	2/23/18	HCID, cPAHs, metals
B-16_17.5	1802-249	2/23/18	HCID, cPAHs, metals



Sample ID	Lab Ref. No.	Date Collected	Analyses
B-11_2.5	1802-278	2/26/18	HCID, Dx, cPAHs, PCBs, metals
B-11_12.5	1802-278	2/26/18	HCID, cPAHs, metals
B-10_7.5	1802-278	2/26/18	HCID, Dx, cPAHs, PCBs, metals
B-05_5	1802-278	2/26/18	HCID, cPAHs, metals
B-08_12.5	1802-278	2/26/18	HCID, cPAHs, metals
B-07_2.5	1802-278	2/27/18	HCID, Dx, cPAHs, PCBs, metals
B-07_7.5	1802-278	2/27/18	HCID, Dx, cPAHs, PCBs, metals
B-07_12.5	1802-278	2/27/18	Dx, cPAHs, PCBs, metals
PP28-8	1803-007	3/1/18	HCID, cPAHs, metals
PP28-10	1803-007	3/1/18	HCID, cPAHs, metals
PP29-3	1803-007	3/1/18	HCID, Dx, cPAHs, PCBs, metals
PP29-5	1803-007	3/1/18	HCID, cPAHs, metals
PP29-10	1803-007	3/1/18	HCID, cPAHs, metals
PP32-4	1803-007	3/1/18	HCID, cPAHs, metals
PP32-7	1803-007	3/1/18	HCID, Dx, cPAHs, PCBs, metals
PP32-10	1803-007	3/1/18	Dx, cPAHs, PCBs, metals
PP31-3	1803-007	3/1/18	HCID, cPAHs, metals
PP31-11	1803-007	3/1/18	HCID, cPAHs, metals
PP33-3	1803-007	3/1/18	HCID, Gx/BTEX, Dx, cPAHs, PCBs, metals
PP33-5	1803-007	3/1/18	HCID, cPAHs, metals
PP33-10	1803-007	3/1/18	HCID, cPAHs, metals
PP23-2	1803-007	2/28/18	HCID, Gx/BTEX, Dx, cPAHs, PCBs, metals, TCLP Pb
PP23-5	1803-007	2/28/18	HCID, cPAHs, metals
PP23-10	1803-007	2/28/18	HCID, cPAHs, metals
PP23-15	1803-007	2/28/18	HCID, cPAHs, metals, TCLP Pb
PP21-2	1803-007	2/28/18	HCID, Dx, cPAHs, PCBs, metals, TCLP Pb
PP21-6	1803-007	2/28/18	HCID, Dx, cPAHs, PCBs, metals, TCLP Pb
PP21-10	1803-007	2/28/18	HCID, Gx/BTEX, Dx, cPAHs, PCBs, metals, TCLP Pb
PP21-15	1803-007	2/28/18	Dx, cPAHs, PCBs, metals
PP20-2	1803-007	2/28/18	HCID, Dx, cPAHs, PCBs, metals
PP20-5	1803-007	2/28/18	HCID, Dx, cPAHs, PCBs, metals, TCLP Pb
PP20-10	1803-007	2/28/18	HCID, Dx, cPAHs, PCBs, metals, TCLP Pb
PP20-15	1803-007	2/28/18	Dx, cPAHs, PCBs, metals, TCLP Pb
PP18-3	1803-007	2/28/18	HCID, cPAHs, metals, TCLP Pb
PP18-5	1803-007	2/28/18	HCID, cPAHs, metals, TCLP Pb
PP18-10	1803-007	2/28/18	HCID, cPAHs, metals
PP17-1	1803-007	2/28/18	HCID, cPAHs, metals, TCLP Pb



Sample ID	Lab Ref. No.	Date Collected	Analyses
PP17-5	1803-007	2/28/18	HCID, cPAHs, metals, TCLP Pb
PP17-10	1803-007	2/28/18	HCID, cPAHs, metals
PP16-1	1803-007	2/28/18	HCID, Gx/BTEX, Dx, cPAHs, PCBs, metals
PP16-11	1803-007	2/28/18	HCID, cPAHs, metals, TCLP Pb
PP14-5	1803-007	2/28/18	HCID, cPAHs, metals
PP14-12	1803-007	2/28/18	HCID, cPAHs, metals
PP11-5	1803-007	2/28/18	HCID, cPAHs, metals, TCLP Pb
PP11-10	1803-007	2/28/18	HCID, cPAHs, metals, TCLP Pb
PP10-11	1803-007	2/28/18	HCID, cPAHs, metals
PP10-17	1803-007	2/28/18	HCID, cPAHs, metals
PP27-7	1803-007	2/28/18	HCID, cPAHs, metals, TCLP Pb
PP27-10	1803-007	2/28/18	HCID, cPAHs, metals
PP30-5	1803-007	2/28/18	HCID, Dx, cPAHs, PCBs, metals
PP30-10	1803-007	2/28/18	HCID, cPAHs, metals
PP24-7	1803-007	2/28/18	HCID, Gx/BTEX, Dx, cPAHs, PCBs, metals, TCLP Pb
PP24-10	1803-007	2/28/18	HCID, cPAHs, metals

The laboratory's performance was reviewed in accordance with quality control (QC) criteria established in the *Pacific Park/Dumpsite Environmental and Economic Assessment Sampling and Analysis Plan* (SAP) (Herrera 2017), by the laboratory, and in the specified methods.

Quality control data summaries submitted by the laboratory were reviewed; raw data were not submitted by the laboratory. Data qualifiers (flags) were added to the sample results in the laboratory reports. Data validation results are summarized below, followed by definitions of data qualifiers.

Custody, Preservation, Holding Times, and Completeness—Acceptable

The samples were properly preserved and sample custody was maintained from sample collection to receipt at the laboratory. Samples were analyzed within the required method holding times. The laboratory reports were complete and contained results for all samples and tests requested on the chain-of-custody (COC) forms.



Laboratory Reporting Limits—Acceptable

The laboratory reporting limits were reasonable for the specified methods and were below relevant comparison criteria. No data were qualified based on laboratory reporting limits.

Method Blank Analysis—Acceptable

Method blanks were analyzed at the required frequency. Method blanks did not contain levels of target analytes above the laboratory reporting limits.

Laboratory Control Sample Analysis—Acceptable

Blank spike (BS) or blank spike/blank spike duplicate (BS/BSD) samples were analyzed for gasoline/BTEX and cPAHs. The percent recovery values met the criteria established by the laboratory.

Surrogate Spike Analysis—Acceptable with Discussion

Surrogate compounds were added to all samples and laboratory QC samples for all NWTPH-HCID, NWTPH-Gx/BTEX, NWTPH-Dx, cPAHs, and PCBs analyses. With the exceptions noted below, all surrogate percent recoveries met the control limits specified by the laboratory or method.

Surrogate o-terphenyl was not recovered for the analysis of samples B-07_7.5. PP29-3, PP33-3, PP21-10, or PP16-1 by method NWTPH-Dx, or for the analysis of sample PP33-3 by method NWTPH-HCID due to the necessary dilutions. No data were qualified.

Matrix Spike Analysis—Acceptable with Discussion

Matrix spike samples were analyzed for metals; matrix spike/matrix spike duplicate (MS/MSD) samples were analyzed for cPAHs, PCBs and metals. With the exception noted below, the percent recovery values met the control limits established by the methods.

A MS/MSD performed on a batch sample resulted in percent recovery values (0 and 218 percent) for chromium that exceeded the 75 to 125 percent criteria. No data were qualified because the sample was not a project sample and all other criteria were met.

Laboratory Duplicate Analysis—Acceptable with Discussion

Laboratory duplicate samples were analyzed for metals; MS/MSD samples were analyzed for cPAHs and PCBs. The relative percent difference (RPD) was calculated for each analyte where both duplicate values were greater than five times the reporting limit (RL). The difference



between duplicate values was calculated if the detected compound concentration was less than five times the RL in either the sample or the duplicate. The relative percent difference (RPD) values or difference values met the control limits established by the laboratory or specified method, except as noted below.

A MS/MSD performed on a batch sample resulted in a RPD values (88 percent) for chromium that exceeded the less than 20 percent criterion. No data were qualified because the sample was not a project sample and all other criteria were met.

DEFINITION OF DATA QUALIFIERS

The following are data qualifier definitions applied for this project.

Data Qualifier	Definition
J	Value is an estimate based on analytical results
R	Value is rejected based on analytical results
U	Value is below the reporting limit
UJ	Value is below the reporting limit and is an estimate based on analytical results

REFERENCES

Herrera. 2017. Pacific Park/Dumpsite Environmental and Economic Assessment, Pacific, Washington, Sampling and Analysis Plan. Prepared by Herrera Environmental Consultants for River and Floodplain Management Section, King County Water and Land Resources Division. April 2017.

APPENDIX I

Soil Vapor Monitoring Data



	MW9 (B-11)	
Gas Probe ID:	home (d)	
Sample ID: NA	1 /	

Canister ID:

NA

Initial Canister Pressure: Final Canister Pressure:

NA NA

Date & Time: 3/23/18011-45Total Casing Volume (cc): 6/8 cc/ft, x/6 = 3,708cc = 1 well vol,

Field Personnel: _G. Iftner_

Casing Volume Purged	Volume Purged (cc)	Purge Rate (ml/min)	Purge Time	CH₄ (% volume)	CO ₂ (% volume)	O ₂ (% volume)	H ₂ S (ppmv)
0	0	300	O Static	0,60	0,00	22.2	GLOD
1/4	0921	300	185 purpl	0.00	0.00	22.4	0-00
1/2	ø 1954	300	370 sec	0,00	0.00	22.5	0,00
3/4	02781	300	556 sec	0.00	0,00	22.5	0.00
1	03708	300	741 sec	0,00	0,00	22,6	0.00
1 1/4	04635	300	927 sec	0.00	0,00	22,5	0,00
1 1/2	05562	300	1112 sec	0,00	0.00	22.6.	0,00
1 3/4	06489	300	1298 sec	0.00	0,00	22.6	0,00
2	07416	300	1483 sec	0,00	0.00	22,6	0,00
2 1/4	0	300	sec				
2 1/2	0	300	sec		y		
2 3/4	0	300	sec				
3	0	300	sec				*.

3,708 ce/300 mil purge = Static WL = 5,85 rounded up to 6' Comments: 12,36 Min = 79

Equipment Used: Gem 2000+, Water Level Meter

hw6 Gas Probe ID:

Sample ID: NA

Sample ID: NA Date & Time: $\frac{3/23/18}{2} = \frac{9}{12.05}$ Total Casing Volume (cc): 618 α /ft × 6.29 = 3,893 cc = 1 Well vol.

Canister ID:

NA

Initial Canister Pressure: Final Canister Pressure:

NA NA

Field Personnel: _G. Iftner_

Casing Volume Purged	Volume Purged (cc)	Purge Rate (ml/min)	Purge Time	CH ₄ (% volume)	CO ₂ (% volume)	O ₂ (% volume)	H ₂ S (ppmv)
- 0	0	300	Static O sec	0100	0.00	22.2	0,00
1/4	0973	300	194 pural	6.00	0,00	22,1	0,00
1/2	01946	300	389 sec	0.00	0.00	22,0	0,00
3/4	02919	300	584 sec	0.00	0.60	22.0	0,00
1	ø 3893	300	778 sec	0,00	6.00	22,1	0,00
1 1/4	04866	300	972 sec	0.00	0,00	22,1	0,00
1 1/2	ø 5839	300	1,167 sec	0.00	0.00	22,1	0,00
1 3/4	\$6812	300	1,362 sec	0.00	0,00	22,1	0,00
2	e7185	300	1,556 sec	0.00	0,00	22,1	0.00
2 1/4	0	300	sec				
2 1/2	0	300	sec		:		
2 3/4	0	300	sec				
3	0	300	sec				

Static WL = 6.29' - 3,893/300 m/min = 129 min = 778 seconds Comments:

Equipment Used: Gem 2000+, Water Level Meter

Gae	Probe	ın.	M	W	6
uas.	LIUDE	112.			

Sample ID: NA 12/18 @ 11:00 AM

Canister ID:

NA

Initial Canister Pressure: Final Canister Pressure:

NA NA

Total Casing Volume (cc): 618 cc/ft x 6.24 = 3,856 cc = 1 Well Volume

Field Personnel: _G. Iftner_

Casing Volume Purged	Volume Purged (cc)	Purge Rate (ml/mln)	Purge Time	CH₄ (% volume)	CO ₂ (% volume)	O ₂ (% volume)	H ₂ S (ppmv)
0	0	3000	O Static	0.00	0,1	19,6	0,00
1/4	\$ 964	3000	19 Purge	0,00	6,3	19,1	0.00
1/2	8928	3000	38 sec	0.00	1.5	16.4	0,00
3/4	0 2892	3000	57 sec	0,08	3.1	14.8	0.00
. 1	x 3856	3000	76 sec	0,00	3.6	14.6	0,00
1 1/4	84820	3000	95 sec	0,00	3.7	14.4	0,00
1 1/2	85784	3000	114 sec	0.00	3,8	14.3	0.00
1 3/4	06748	3000	133 sec	0.00	3.8	14.2	0,00
2	67712	3000	154 sec	0,00	3,9	14.1	0,00
2 1/4	0	3000	sec				
2 1/2	0	3000	sec				
2 3/4	0	3000	sec				
3	0	3000	sec				

Static WL = 6.24' 3,856 (3000 milmin = 1,28 min = 76.8 seconds. Comments:

Equipment Used: SKC Pump, Gem 2000+, Water Level Meter

Gas Probe ID: MW 9 Sample ID: NA Date & Time: 6/21/18 © U:20

Canister ID: NA
Initial Canister Pressure: NA
Final Canister Pressure: NA

Final Canister Pressure:
Field Personnel: _G. Iftner_

Total Casing Volume (cc): 6/8 α /ff x 6.62 = 3,708 cc = 1 well Vol.

Casing Volume Purged	Volume Purged (cc)	Purge Rate (ml/min)	Purge Time	CH ₄ (% volume)	CO ₂ (% volume)	O ₂ (% volume)	H₂S (ppmv)
0	0	3000	O sec	6,00	0,1	14,6	
1/4	2927	3000	18 PWHE sec	0,00	0.1	19,7	
1/2	ø 1854	3000	36 sec	6.00	0.1	19,6	
3/4	g 2781	3000	SY sec	6.04	0.1	19,6	
1	ø318	3000	72 sec	0,00	01	19,6	
1 1/4	p 4635	3000	90 sec	0,00	oil	19,6	
1 1/2	ø5562	3000	log sec	0,00	0,1	19.6	
1 3/4	06489	3000	126 sec	0,00	0, (19.6	
2	07416	3000	144 sec	0.00	0,1	19.6	
2 1/4	0	3000	sec				
2 1/2	0	3000	sec				
2 3/4	0	3000	sec				
3	0	3000	sec				

Comments: 3,708 cs/3,000 milmin = 1,2 min = 72 seconds (1 will volume)

Equipment Used: SKC Pump, Gem 2000+, Water Level Meter

Gas Probe ID:	MW9		
Sample ID: NA Date & Time: _	9/26/18	0	14:Z0

Canister ID:

NA

Initial Canister Pressure: Final Canister Pressure:

NA NA

Total Casing Volume (cc): 6/8cc/ff x 6.98= 4,3/4 = 1 well Vol.

Field Personnel: _G. Iftner____

Casing Volume Purged	Volume Purged (cc)	Purge Rate (ml/min)	Purge Time	CH₄ (% volume)	CO ₂ (% volume)	O₂ (% volume)	H ₂ S (ppmv
0	0	3000	Static sec	0.0	0.4	19,4	0,0
1/4	ø1018.5	3000	22 sec	5.0	1.6	17.9	0.0
1/2	02151	3000	44 sec	0,0	2.6	16.6	0,0
3/4	031355	3000	66 sec	0.0	3,7	15.3	0, 0
1	94314	3000	86 sec	0.0	3,8	15.	0,0
1 1/4	\$392.5	3000	[08 sec	0,0	3-9	15.0	0.0
1 1/2	06471	3000	131 sec	S-0	3,9	15.0	0.0
1 3/4	ø7549.\$	3000	153 sec	0.0	3.9	14-9	0.0
2	68628	3000	172 sec	0.0	3.9	14.9	0.0
2 1/4	0	3000	sec				
2 1/2	0	3000	sec				
2 3/4	0	3000	sec				
3	0	3000	sec				

comments: , = 86 Static Water Level = 6,981: 4,314/3000 4/min(SKC proge vate) = 1.44 min

Burd Metri C Pressure 36.11 49. Equipment Used: SKC Pump, Gem 2000+, Water Level Meter

MW6 Gas Probe ID:

1 1/2

1 3/4

2 1/4

2 1/2

2 3/4

2

Canister ID:

0.0

0.0

0.0

7,6

7,5

NA

Sample ID: NA 9 /26/18 11:40

Initial Canister Pressure: Final Canister Pressure:

NA NA

Total Casing Volume (cc): 618 cc/ff x 8,53 = 5,272 (c = 1 well Vol.

\$ 7908

09226

00544

0

0

0

3000

3000

3000

3000

3000

3000

Field Personnel: _G. Iftner_

Volume Purged **Purge Rate Purge Time** CH₄ CO, O2 H₂S **Casing Volume Purged** (ml/min) (cc) (% volume) (% volume) (% volume) (ppmv) 0.0 Static 00 0 0 3000 sec er 1318 27 0.0 6,3 000 1/4 3000 sec Ø2636 0,0 6.0 1/2 3000 sec 0.0 8 0.0 ø 3954 3/4 3000 sec 11.0 Ø 5277 0.0 000 3000 sec 133 ø6590 7.5 0.0 10.9 0.0 1 1/4 3000 sec 7.6

60

187

212

sec

sec

sec

sec

sec

sec

3 0 3000 sec comments: Static Water Level = 8.53', 5,272 (c/3000 m/min (SKC purge rate) = 1.76 min Comments:

Equipment Used: SKC Pump, Gem 2000+, Water Level Meter

10,9

10,9

10,8

0,0

0.0

0.0

MWI Gas Probe ID:

Sample ID: NA

Date & Time: 9/26/18 10:00 Am

Total Casing Volume (cc): 618 cc/ff x 5.80= 3584 cc = 1 well vu |.

Canister ID:

Initial Canister Pressure:

NA NA

Final Canister Pressure:

NA

Field Personnel: __G. Iftner_

Casing Volume Purged	Volume Purged (cc)	Purge Rate (ml/min)	Purge Ti	me	CH₄ (% volume)	CO ₂ (% volume)	O ₂ (% volume)	H ₂ S (ppmv)
0	0	3000	Static	sec	6.0	0.0	20.6	0,0
1/4	8996	3000	18	sec	0.0	1.7	20.1	0.0
1/2	×1792	3000	36	sec	0.0	1.9	19.0	0.0
3/4	ø2698	3000	54	sec	0,0	2.0	18.9	0.0
1	03584	3000	72	sec	0.0	2,0	18.9	0.0
1 1/4	04480	3000	90	sec	0,0	2.1	18.8	0.0
1 1/2	ø5316	3000	108	sec	0:0	211	[8.8]	0,0
1 3/4	06272	3000	126	sec	0.0	2.1	18.7	0-0
2	27168	3000	144	sec	0,0	211	18.7	0.0
2 1/4	0	3000		sec				
2 1/2	0	3000		sec				
2 3/4	0	3000		sec				
3	0	3000		sec				

Statil W1 = 5.80' 3584 cc/3000 m1/min (SKC Purge rate) = 1.20 min = 72 Seconds Comments:

Equipment Used: SKC Pump, Gem 2000+, Water Level Meter

Page 1