



Washington State  
Department of Transportation

# SR 520 Bridge Replacement and HOV Program



## Phase II Environmental Site Assessment State Route (SR) 520 Eastbound Off-Ramp to Montlake Vicinity Seattle, Washington

Prepared for

Washington State Department of Transportation  
SR520 Bridge Replacement and HOV Program  
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December 8, 2016



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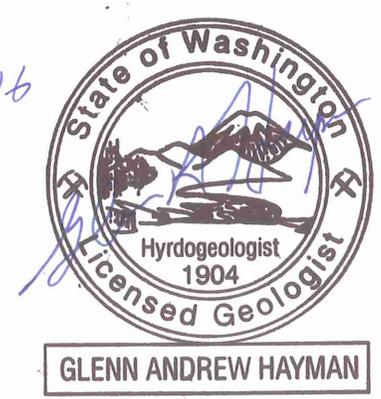
**Innovex Environmental Management, Inc.  
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We have performed a Phase II environmental site assessment of the property at (SR) 520 Eastbound Off-Ramp to Montlake Vicinity Seattle, Washington in conformance with the scope and limitations of ASTM Practice E 1903-11 and for the following objectives:

The Phase II ESA was conducted to determine if petroleum related contamination is present in the subsurface of the Washington State Department of Transportation (WSDOT) and City of Seattle right-of-way adjacent to the property at 2625 East Montlake Place East due to potential releases from the underground storage tanks (USTs) and/or the auto body/service station. The USTs and potential contaminant source was identified in the Limited Phase I ESA conducted by WSDOT (WSDOT, 2016).

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## Acronyms and Abbreviations

bgs	below ground surface
COPC	contaminant of potential concern
CSM	Conceptual Site Model
Dx	Diesel-range petroleum hydrocarbons
Ecology	Washington Department of Ecology
ESA	Environmental Site Assessment
Gx	Gasoline-range hydrocarbons
HCID	Hydrocarbon Identification
INNOVEX	Innovex Environmental Management, Inc.
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
MS	Matrix Spike
MSD	Matrix Spike Duplicate
MTCA	Model Toxics Control Act
NWTPH	Northwest Total Petroleum Hydrocarbons
OnSite	OnSite Environmental Inc.
PCB	Polychlorinated Biphenyl
PDI	Photoionization Detector
PP-13	Priority Pollutant Metals
Program	SR 520 Bridge Replacement and HOV Program
QA	Quality Assurance
QC	Quality Control
RPD	Relative Percent Difference
SR	State Route
SAP	Sampling and Analysis Plan
SDOT	Seattle Department of Transportation
SVOC	Semivolatile Organic Compound
TPHg	Total Petroleum Hydrocarbons as Gasoline
UST	Underground Storage Tank
VOC	Volatile Organic Compound
WAC	Washington Administrative Code
WSDOT	Washington State Department of Transportation



## 1.0 INTRODUCTION

### 1.1 Authorization

Innovex Environmental Management, Inc. (INNOVEX) has completed a Phase II Environmental Site Assessment (ESA) of City of Seattle and WSDOT right-of-way in the vicinity of 2625 East Montlake Place East, Seattle, Washington (Site) for the Washington State Department of Transportation (WSDOT) SR520 Bridge Replacement and HOV Program (PROGRAM). This work was conducted under Master Subconsultant Agreement Y-11848, Task Order AU00.

A Sampling and Analysis Plan (SAP) was prepared to describe the Phase II ESA activities for WSDOT and as a guide for field activities, including explorations, field testing, soil and groundwater sampling, and the handling of investigation-derived waste (IDW). Deviations from the SAP are summarized in Section 4.8 of this report.

### 1.2 Objective

The Phase II ESA was conducted to determine if petroleum related contamination was present in the subsurface of the right-of-ways adjacent to the property at 2625 East Montlake Place East due to the '76 gasoline and service station at this address. The USTs and potential contaminant source was identified in the Limited Phase I ESA conducted by WSDOT (WSDOT, 2016).

The scope of this investigation included the following tasks:

- Advancement of five soil borings to collect soil and, if encountered, groundwater samples for chemical analysis.
- Collection and analysis of soil and groundwater samples to determine if petroleum related contamination is present in the soil and groundwater beneath the Site.
- If petroleum related contamination is detected, additional samples were collected to determine the vertical and horizontal extent of the contamination plume when possible.
- Analysis of one soil sample from each boring for Site contaminants of concern. Additional samples were analyzed based on field observations and initial analytical results with the approval of the WSDOT Project Manager.
- The results of this Phase II investigation will be used to assist WSDOT in management of potential environmental risks associated with the reconstruction of the eastbound SR 520 off ramp to Montlake Boulevard East and East Montlake Place East.



## 2.0 SITE BACKGROUND

### 2.1 Site Description

The areas investigated for the Phase II ESA are the City of Seattle and WSDOT SR 520 right-of-ways in the vicinity of 2625 East Montlake Place East, Seattle, Washington. The investigation took place in the northeast quarter of Township 25 North, Range 4 East and Section 21.

### 2.2 Site History

East Montlake Place East was shown on the 1897 and the 1909 (reprinted 1936) topographic maps. The area of investigation was otherwise undeveloped. By 1949 the roads and streets in the area were largely in place, with the exception of SR 520. SR 520 was built in the early 1960's. The area adjacent to the Site is a residential neighborhood. The Montlake Playfield, a community center and sports field are a few blocks west of the Site.

#### 2.2.1 Limited Phase I Environmental Site Assessment (ESA) Results

WSDOT conducted a limited Phase I ESA of the property immediately south and west of the areas investigated. This property is occupied by a '76 gasoline and service station and the Montlake Boulevard Market. Historically the property has been operated as a gasoline station and grocery store since 1926. The findings of limited Phase I ESA relevant to this Phase II ESA are:

- There are three active single wall gasoline USTs on the property. The year of installation and the UST size are as follows: 1952, 5,000 gallons; 1962, 10,000 gallons; and 1975, 10,000 gallons. In addition, in 1975 a 300 gallon used waste oil UST was installed. This tank was closed in place at an unknown date.
- There was no known environmental documentation identified during the limited Phase I investigation indicating that an inadvertent spill or release was present at the Site.
- An oil and gasoline facility resided northwest of the Subject Property across West Montlake Place.

The Limited Phase I recommendations included that a Phase II ESA be conducted to determine if the USTs have failed and released petroleum hydrocarbons into the surrounding soils and/or groundwater.

### 2.3 Previous Environmental Sampling and Analysis

WSDOT provided INNOVEX with copies of analytical reports for soil samples from nearby geotechnical explorations. The explorations are H-609p-11, H-667p-15 and H-691p-16 (Figure 2). Hydrocarbons were not detected in the soil sample from each exploration that was analyzed. The analytical results for a groundwater sample collected from H-667p-15 were also provided. In the sample, total petroleum hydrocarbons as gasoline (TPHg) was detected at a concentration of 170 micrograms per liter (ug/L) and total xylenes were detected at a concentration of 1.6 ug/L. The concentrations of the detected compounds are below the cleanup levels of 1,000 ug/L and 1,000 ug/L respectively. Laboratory analytical reports are presented in Appendix B. The Limited Phase I ESA identified a former oil and gasoline facility near the location of boring H-667p-15 from a Sanborn Map. An enlarged portion of the Sanborn Map is contained in Appendix C



## **3.0 PHYSICAL SETTING AND GEOLOGY**

### **3.1 Physical Setting**

The Site lies within the Puget Sound Lowland, which consists of a broad, low-lying region situated between the Cascade Range to the east and the Olympic Mountains to the west. The Lowland depression is underlain by Tertiary volcanic and sedimentary bedrock and is filled to the present-day land surface with Quaternary glacial and non-glacial sediments.

The Puget Sound Lowland's present-day geomorphic features can be attributed to the last continental glacier, the Cordilleran ice sheet, which covered the region during the Fraser Glaciation. The ice sheet advanced from British Columbia 18,000 years ago to just south of Olympia and disappeared approximately 10,000 years ago (Lasmanis, 1991).

The area investigated was below paved surfaces within the right of way for East Montlake Place East and the SR 520 eastbound off ramp. The surface elevation was between approximately 50 and 60 feet above the NAVD 88 datum. The surface topography slopes to the west towards Portage Bay.

### **3.2 Site Geology and Hydrogeology**

The Site lies approximately 600 feet east of Portage Bay (Lake Union). Groundwater flow is inferred to follow surface topography and flow to the west. The topographic relief is less than 10 feet and primarily due to the slope of the SR 520 off ramp.

Subsurface conditions at the Site have been interpreted from the soil boring logs completed during this investigation, and from geotechnical boring logs provided by WSDOT. The borings were advanced as part of the geotechnical evaluation of the SR 520 Bridge Replacement and HOV Program. The borings are H-609p-15, H-667p-15 and H-691p-16 (Figure 2). Each boring had a transducer for water level measurements installed in it. The nominal depth to groundwater is generally 10 to 12 feet below ground surface (bgs). Boring logs, groundwater level measurements and a figure with boring locations are presented in Appendix B.

The near surface soils observed below the roadway subgrade were generally found to be silty sand with some fine gravel during the Phase II ESA fieldwork. Dense glacial till was encountered at a depth of between 15 and 24 feet bgs. The till was primarily fine silty sand and fine sandy silt with occasional fine to medium gravel and some clay. Boring logs are presented in Appendix A.



## **4.0 PHASE II ESA ACTIVITIES**

### **4.1 Scope of Assessment**

A sampling and analysis plan was developed (INNOVEX, 2016) to investigate recognized environmental conditions (RECs) identified in the Limited Phase I ESA. A total of five soil borings were conducted. Subsurface soil samples and groundwater samples were submitted to OnSite Environmental Inc. (OnSite) in Redmond, Washington for chemical analysis. Our rationale and the results for the exploration program are summarized below.

### **4.2 Conceptual Site Model and Sampling Plan**

In order to provide a framework for evaluating data gaps and subsequent analytical data, a conceptual site model (CSM) depicting potential sources of chemicals, release mechanisms, means of retention in or migration to exposure media, exposure routes, and receptors was developed for the Site. The CSM describes, in a generalized way, the interactions of potential contaminants, mechanisms of contaminant migration, and possible routes of human and ecological exposure under site-specific conditions.

Based on background information previously presented, the contaminants of potential concern (COPCs) identified for the Site included:

- Gasoline-range petroleum hydrocarbons
- Diesel-range petroleum hydrocarbons
- Oil-range petroleum hydrocarbons
- Polychlorinated Biphenyls (PCBs)
- Volatile Organic Compounds (VOCs)
- Semi-volatile Organics (SVOCs)
- Priority Pollutant Metals (PP-13)

The PP-13 metals are; antimony, arsenic, beryllium, cadmium, chromium, copper, lead, nickel, selenium, silver, thallium, and zinc.

### **4.3 Utility Location**

WSDOT coordinated the utility locating. There were no buried utilities at the boring locations. On East Montlake Place East there were overhead power lines for king County Metro electric trolleys. WSDOT arranged for the power to be turned off to allow soil borings to be advanced.

### **4.4 Permits and Traffic Control**

WSDOT obtained Street Use permits from WSDOT and SDOT for work on the SR 520 off ramp and East Montlake Place East respectively. In addition, they obtained a noise variance to allow the drilling to be conducted at night, and reducing the impacts to traffic and local businesses.

WSDOT prepared a traffic control plan. INNOVEX subcontracted GHD, Inc. to implement the plan.

## 4.5 Mud Rotary Drilling

Five soil borings were advanced and sampled (Figure 2) by a WSDOT drilling crew with a mud rotary drill rig. The drilling mud consisted of water with no bentonite or other additives. Drilling fieldwork was conducted on the nights of October 6, 7, and 8, 2016. Although there were no known releases identified by the Limited Phase I ESA, the Program determined it prudent to investigate subsurface conditions within City and State right-of-ways adjacent to the '76 gasoline and service station.

Fieldwork consisted of collecting subsurface soil samples with a 2-inch split spoon from a total of five soil boring locations. A petroleum odor and elevated photoionization detector (PID) readings were noted in soil samples from four of the borings. Soil samples from boring H-1-16 did not have a petroleum odor or elevated PID readings.

Downhole drilling equipment, hoses, and storage tank and pump on the rig were decontaminated at the WSDOT facility on South Corson Avenue in Seattle before the start of drilling and after each nights work. The drilling mud (water) was changed between each boring.

## 4.6 Soil and Groundwater Sampling

A total of 40 soil samples and one groundwater sample were collected from the soil borings. Eight soil samples were collected from each boring. The soil sample from each boring with the highest PID reading was selected for chemical analysis except as discussed in the next paragraph for samples from boring H-1-16. Additional samples were analyzed for selected analyses based on field screening and initial analytical results with the approval of the WSDOT Project Manager.

Soil samples from boring H-1-16 did not have elevated PID readings. The soil sample from above the contact with dense glacial till was selected for chemical analysis.

One groundwater sample was collected with the driller's bailer from boring H-3-16 located on the SR 520 off ramp. No other groundwater samples were collected.

## 4.7 Analytical Methods

The COPCs identified for the Site include petroleum hydrocarbon related constituents. Selected soil and groundwater samples were analyzed to determine the concentrations of these COPCs using the following methods:

- Hydrocarbon Identification (HCID) – Northwest Total Petroleum Hydrocarbon (NWTPH) HCID (Soil and water)
- Gasoline-range petroleum hydrocarbons –NWTPH-Gx (soil and water)
- Diesel-range petroleum hydrocarbons – NWTPH-Dx (soil and water)
- Oil-range petroleum hydrocarbons – NWTPH-Dx (soil and water)
- Polychlorinated Biphenyls (PCBs) – EPA Method 8082 (soil and water)
- Volatile Organic Compounds (VOCs) – EPA Method 8260 (soil and water)
- Semi-volatile Organics (SVOCs) – EPA Method 8270 (soil and water)
- Priority Pollutant Metals (PP-13) – EPA Method 6010 (soil and water)

Table 1 summarizes the samples submitted to OnSite Environmental (OnSite) for chemical analysis and the requested analyses.

#### 4.7.1 Soil Analytical Results

Detected analytes in soil are summarized in the following tables:

- Table 4 HCID Soil Analytical Results
- Table 5 Volatile Organic Compounds Detected in Soil Samples
- Table 6 Semivolatile Organic Compounds Detected in Soil Samples
- Table 7 Polychlorinated Biphenyls Detected in Soil Samples
- Table 8 Metals Detected in Soil Samples

The analytical reports are included in Appendix D.

HCID analysis did not detect total petroleum hydrocarbons as gasoline (TPHg) in samples from borings H-1-16, H-2-16, H-3-16, and H-5-16 (Table 4). TPHg was detected in two soil samples from boring H-4-16.

VOCs were not detected in the analyzed soil sample from boring H-1-16. VOCs were detected in soils samples from soil borings H-2-16, H-3-16, H-4-16 and H-5-16 (Table 5). VOCs above the applicable cleanup level were detected soil samples from borings H-3-16, H-4-16, and H-5-16. The compounds that had concentrations above the cleanup level are benzene, ethylbenzene, total xylene, naphthalene, and methylene chloride.

SVOCs were not detected in the analyzed soil sample from boring H-1-16. SVOCs were detected in soils samples from soil borings H-2-16, H-3-16, H-4-16 and H-5-16 (Table 6). One SVOC, naphthalene, was above the applicable cleanup level in a soil sample from borings H-4-16. Naphthalene is an analyte for both the VOC and SVOC analytical method.

PCBs were not detected in any of the soil samples analyzed (Table 7).

Chromium, copper, lead, nickel and zinc were present in soil samples from each of the five boring at concentrations above the analytical method reporting limit (Table 8). Copper and nickel were present at concentrations above the applicable cleanup level.

#### 4.7.2 Groundwater Analytical Results

One groundwater sample was collected as part of the investigation. It was collected from boring H-3-16 and was analyzed for VOCs, SVOCs and dissolved metals. The detected analytes are summarized in Table 2. Three VOCs, benzene, bromochloromethane, and chloroform; and two dissolved metals, antimony and arsenic had concentrations greater than the applicable cleanup level.

#### 4.8 Deviations from the Sampling and Analysis Plan

The approved SAP identified the drilling method as auger drilling. The drilling crew arrived on site with a mud rotary drill rig. The WSDOT and Innovex project managers talked and recognized that by using mud rotary drilling the data quality would be of reduced and that the data would not be legally defensible. Contaminant concentrations could be impacted because the recirculated drilling mud (water with no additives) could reduce high contaminant concentrations and also increase low

contaminant concentrations in soil or groundwater. However; the data would be of sufficient quality to determine if contaminants of concern were present in the soil and to meet the needs of WSDOT.

#### **4.9 Data Quality**

Data reports from OnSite were reviewed by INNOVEX. Laboratory provided data quality parameters were reviewed. Data qualifiers were applied as necessary. Data for VOCs, PAHs, PCBs, MTCA metals, and TPHs were determined by INNOVEX to be as qualified acceptable for all purposes following evaluation of the quality control specifications presented in the SAP; or equivalent requirements found in the contracted commercial laboratory analytical methods. Precision, accuracy, representativeness, comparability, and completeness parameters were evaluated for each method. In addition to laboratory control samples, the data were also reviewed for trip temperature and holding time.

The Method 5035A VOA vials provided for soil sample H-1-16-10 contained too much soil to perform the requested analysis. Therefore, the sample was extracted from an 8-ounce jar and analyzed. Some loss of volatiles may have occurred.

All four Internal Standards did not meet acceptance criteria for samples H-3-16-6 and H-3-16-8.5. The samples were re-analyzed with similar results. Leaks in the sealed VOA environment caused by grit between the VOA lip and VOA cap septum have been shown to cause low internal standard recovery. The samples were consequently extracted from their respective 8-ounce jars, analyzed, and reported. Some loss of volatiles may have occurred, and common laboratory solvents Acetone and Methylene Chloride may have been introduced during sample preparation. Acetone and Methylene Chloride were detected in sample H-3-16-8.5. Methylene Chloride was detected sample H-3-16-8.5. These results were “H” qualified indicating that the analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.

OnSite followed most recent version of the specified analytical methods. Precision was acceptable as demonstrated by the reported matrix spike/matrix spike duplicate (MS/MSD) laboratory control sample/laboratory control sample duplicate (LCS/LCSD) relative percent difference (RPD) values. Accuracy was also acceptable, as demonstrated by the reported surrogate, MS/MSD and LCS/LCD percent recovery values. Samples were collected and field activities were conducted in accordance with the SAP, with the exception of the deviations described in the above section.

#### **4.10 Disposal of Investigation Derived Waste**

Investigation Derived waste (IDW) including soil cuttings, decontamination water and drilling mud were containerized by the drilling crew and taken to the WSDOT facility on Corson Avenue, Seattle for temporary storage pending sample analytical results followed by proper disposal.

## 5.0 CONCLUSIONS AND RECOMMENDATIONS

Based on the above findings, deviations from the SAP and analytical results presented above we offer the following conclusions for this investigation.

- VOCs were the primary analytes detected in soil samples at concentrations above applicable MTCA cleanup levels.
- The only SVOC detected soil samples above the applicable MTCA cleanup level was naphthalene in one sample.
- PCBs were not detected in any of the soil or groundwater samples analyzed.
- Copper and Nickel are the only metals detected in soil samples at concentrations above the MTCA soil cleanup level for protection of groundwater.
- VOCs were present in the groundwater sample at concentrations above applicable MTCA cleanup levels.
- No SVOCs were detected in the groundwater sample at concentrations above applicable MTCA cleanup levels.
- PCBs were not detected in the groundwater sample.
- Antimony and arsenic were detected in the groundwater sample at concentrations above applicable MTCA cleanup levels.
- Soil and groundwater in the investigation area appear to be impacted by gasoline and related compounds at concentrations above applicable MTCA cleanup levels.
- Contamination was detected in soil samples collected from boring H-4-16 at depths of up to 25 feet bgs.

The source of the gasoline related analytes and contaminants detected in the soil and groundwater samples collected as part of this investigation is potentially the current or former USTs at the adjacent service station. It is recommended the additional investigation on the service station property be conducted to determine the extent of the contamination identified by this Phase II ESA.

To ensure the results of future environmental investigations are legally defensible INNOVEX recommends that mud rotary drilling not be used.



## **6.0 LIMITATIONS**

This report is based on the Site conditions, data, and other information available as of the date of the report, and the conclusions herein are applicable only to the time frame in which the report was prepared. Background information used to prepare this report including, but not limited to Site plans and other data have been furnished to INNOVEX by WSDOT and as available on Ecology's website. INNOVEX has relied on this information as furnished, and is neither responsible for nor has confirmed the accuracy of this information.



## 7.0 REFERENCES

WSDOT (Washington State Department of Transportation), 2016. Limited Phase I Environmental Site Assessment State Route (SR) 520 Montlake '76 Gasoline and Service Station, Seattle, Washington

Lasmanis, Raymond, 1991. The Geology of Washington Rocks and Minerals. Volume 66, No. 4, p. 262-277.



## TABLES



**Table 1 Summary of Sample Analyses, SR 520 Eastbound Off-Ramp to Montlake Vicinity Seattle, Washington**

Sample Name	Sample Depth (ft bgs)	HCID	TPHg	PCBs	VOCs	SVOCs	PP-13
<b>Soil Samples</b>							
H-1-16-10	10	X		X	X	X	X
H-2-16-13.5	13.5	X		X	X	X	X
H-3-16-3	3	X		X	X	X	X
H-3-16-6	6	X		X	X	X	X
H-3-16-8.5	8.5	X		X	X	X	X
H-4-16-3	3	X	X				
H-4-16-6	6	X	X				
H-4-16-8.5	8.5	X	X				
H-4-16-11	11	X	X				
H-4-16-16	16	X	X	X	X	X	X
H-4-16-18.5	18.5	X	X				
H-4-16-19.9	19.9	X	X				
H-4-16-25.4	25.4	X	X				
H-5-16-3	3	X					
H-5-16-6	6	X					
H-5-16-8.5	8.5	X					
H-5-16-11	11	X					
H-5-16-13.5	13.5	X		X	X	X	X
H-5-16-16	16	X					
H-5-16-18.5	18.5	X					
<b>Groundwater Sample</b>							
H-3-16	NA	X		X	X	X	X <sup>1</sup>

<sup>1</sup> Dissolved metals

**Table 2 Groundwater Cleanup Levels for Detected Analytes and Detected Analyte Concentrations in Sample H-3-16, SR 520 Eastbound Off-Ramp to Montlake Vicinity Seattle, Washington**

Chemical Name	Analyte Type	Sample H-3-16 (µg/L)	MTCA Cleanup Level (µg/L)		CAS #
acetone	VOC	6.5	7200	Method B Non cancer	67-64-1
benzene	VOC	<b>7.4</b>	0.80	Method B Cancer	71-43-2
bromodichloromethane	VOC	<b>1.0</b>	0.71	Method B Cancer	75-27-4
chloroform	VOC	<b>8.3</b>	1.41	Method B Cancer	67-66-3
isopropylbenzene (cumene)	VOC	0.025	800	Method B Non cancer	98-82-8
ethylbenzene	VOC	0.7	700	Method A Unrestricted Land Use	100-41-4
propylbenzene;n-	VOC	0.37	800	Method B Non cancer	103-65-1
trimethylbenzene;1,2,4-	VOC	0.44	--	No Cleanup Level Listed in CLARC Table	95-63-6
trimethylbenzene;1,3,5-	VOC	0.30	80	Method B Non cancer	108-67-8
xylenes, total	VOC	2.77	1000	Method A Unrestricted Land Use	1330-20-7
benzo[a]anthracene	SVOC	0.082	0.12	Method B Cancer	56-55-3
bis(2-ethylhexyl) phthalate	SVOC	5.7	6.25	Method B Cancer	117-81-7
chrysene	SVOC	0.012	12	Method B Cancer	218-01-9
di-n-butylphthalate (di-butyl phthalate)	SVOC	9.3	1600	Method B Non cancer	84-74-2
diethyl phthalate	SVOC	1.3	12800	Method B Non cancer	84-66-2
methyl naphthalene;1-	SVOC	0.13	560	Method B Non cancer	90-12-0
methyl naphthalene;2-	SVOC	0.23	32	Method B Non cancer	91-57-6
naphthalene	SVOC	0.25	160	Method A Unrestricted Land Use	91-20-3
phenanthrene	SVOC	0.11	--	No Cleanup Level Listed in CLARC Table	85-01-8
antimony	Metal	<b>18</b>	6.40	Method B Non cancer	7440-36-0
arsenic, inorganic	Metal	<b>3.3</b>	0.06	Method B Cancer	7440-38-2
lead	Metal	2.5	15	Method A Unrestricted Land Use	7439-92-1

Exceeds cleanup level

Table derived from Ecology's CLARC Master Table at <https://fortress.wa.gov/ecy/clarc/CLARCDATATables.aspx>

VOC - Volatile Organic Compound

SVOC - Semivolatile Organic Compound

Metal - Dissolved Metal

**Table 3 Soil Cleanup Levels for Detected Analytes, SR 520 Eastbound Off-Ramp to Montlake Vicinity Seattle, Washington**

Chemical Name	Sample Name	Highest Soil Concentration (mg/kg)	MTCA Cleanup Level (mg/kg)		CAS #
acetone	H-2-16-13.5	0.060	2	Protective of Groundwater	67-64-1
benzene	H-3-16-8.5	<b>0.35</b>	0.0017	Protective of Groundwater	71-43-2
chromium(III)	H-5-16-13.5	64	2000	Method A Unrestricted Land Use	16065-83-1
copper	H-5-16-13.5	<b>46</b>	14	Protective of Groundwater	7440-50-8
isopropylbenzene (cumene)	H-5-16-13.5	0.19	8000	Method B Non cancer	98-82-8
ethylbenzene	H-4-16-18.5	<b>1.4</b>	0.34	Protective of Groundwater	100-41-4
lead	H-2-16-13.5	11	150	Protective of Groundwater	7439-92-1
2-butanone (methyl ethyl ketone)	H-3-16-8.5	0.021	48000	Method B Non cancer	78-93-3
p-isopropyltoluene	H-5-16-13.5	0.32	--	No Cleanup Level Listed in CLARC Table	99-87-6
1-methyl naphthalene	H-4-16-16	0.37	34	Method B Cancer	90-12-0
2-methyl naphthalene	H-4-16-16	0.74	320	Method B Non cancer	91-57-6
methylene chloride	H-3-16-6	<b>0.053 H</b>	0.0015	Protective of Groundwater	75-09-2
naphthalene	H-4-16-16	<b>0.64</b>	0.24	Protective of Groundwater	91-20-3
n-butylbenzene	H-5-16-13.5	2.0	4000	Method B Non cancer	104-51-8
nickel soluble salts	H-5-16-13.5	<b>69</b>	6.53	Protective of Groundwater	7440-02-0
n-propylbenzene	H-5-16-13.5	1.2	8000	Method B Non cancer	103-65-1
sec-butylbenzene	H-5-16-13.5	0.51	8000	Method B Non cancer	135-98-8
toluene	H-4-16-18.5	0.090	0.27	Protective of Groundwater	108-88-3
TPHg with benzene present	H-4-16-19.9	<b>99</b>	30	Method A Unrestricted Land Use	unavailable25
1,2,4-trimethylbenzene	H-3-16-8.5	1.8	--	No Cleanup Level Listed in CLARC Table	95-63-6
1,3,5-trimethylbenzene	H-4-16-16	0.56	800	Method B Non cancer	108-67-8
xylene, total	H-4-16-18.5	<b>2.28</b>	1	Protective of Groundwater	1330-20-7
zinc	H-5-16-13.5	69	299	Protective of Groundwater	7440-66-6

Exceeds cleanup level

H - The result indicated is a common laboratory contaminant and may have been introduced during sample preparation.

Table derived from Ecology's CLARC Master Table at <https://fortress.wa.gov/ecy/clarc/CLARCDATATables.aspx>

**Table 4 HCID Soil Analytical Results, SR 520 Eastbound Off-Ramp to Montlake Vicinity Seattle, Washington**

Sample ID	Sample Date	Sample depth (ft.)	TPHg (mg/kg)	TPHd (mg/kg)	TPHo (mg/kg)
H-1-16-10	10/6/2016	10	ND	ND	ND
H-2-16-13.5	10/7/2016	13.5	ND	ND	ND
H-3-16-3	10/7/2016	3	ND	ND	ND
H-3-16-6	10/7/2016	6	ND	ND	ND
H-3-16-8.5	10/7/2016	8.5	ND	ND	ND
H-4-16-3	10/8/2016	3	ND	ND	ND
H-4-16-6	10/8/2016	6	ND	ND	ND
H-4-16-8.5	10/8/2016	8.5	ND	ND	ND
H-4-16-11	10/8/2016	11	ND	ND	ND
H-4-16-16	10/8/2016	16	Detected	ND	ND
H-4-16-18.5	10/8/2016	18.5	Detected	ND	ND
H-4-16-19.9	10/8/2016	19.9	ND	ND	ND
H-4-16-25.4	10/8/2016	25.4	ND	ND	ND
H-5-16-3	10/8/2016	3	ND	ND	ND
H-5-16-6	10/8/2016	6	ND	ND	ND
H-5-16-8.5	10/8/2016	8.5	ND	ND	ND
H-5-16-11	10/8/2016	11	ND	ND	ND
H-5-16-13.5	10/8/2016	13.5	ND	ND	ND
H-5-16-16	10/8/2016	16	ND	ND	ND
H-5-16-18.5	10/8/2016	18.5	ND	ND	ND

ND = Not Detected

Table 5 Volatile Organic Compounds Detected in Soil Samples, SR 520 Eastbound Off-Ramp to Montlake Vicinity, Seattle, Washington

Sample ID	Sample Date	Sample depth (ft.)	TPHg (mg/kg)	Acetone (mg/kg)	Methylene Chloride (mg/kg)	2-Butanone (mg/kg)	Benzene (mg/kg)	Toluene1 (mg/kg)	Ethyl-benzene (mg/kg)	Total-xylenes (mg/kg)	Isopropyl-benzene (mg/kg)	n-Propyl-benzene (mg/kg)	1,3,5-Trimethyl-benzene (mg/kg)	1,2,4-Trimethyl-benzene (mg/kg)	sec-Butyl-benzene (mg/kg)	p-Isopropyl-toluene (mg/kg)	n-Butyl-benzene (mg/kg)	Napthalene (mg/kg)	
H-1-16-10	10/6/2016	10	--	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
H-2-16-13.5	10/7/2016	13.5	--	0.060	ND	0.021	0.0053	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
H-3-16-3	10/7/2016	3	--	ND	ND	ND	0.0055	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
H-3-16-6	10/7/2016	6	--	0.023 H	0.053 H	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
H-3-16-8.5	10/7/2016	8.5	--	ND	0.022 H	ND	0.038	ND	0.0050	0.016	0.0018	0.0032	0.0015	0.0018	0.0013	0.002	ND	ND	ND
H-4-16-3	10/8/2016	3	ND	--	--	--	ND	ND	ND	ND	--	--	--	--	--	--	--	--	--
H-4-16-6	10/8/2016	6	ND	--	--	--	0.024	ND	ND	ND	--	--	--	--	--	--	--	--	--
H-4-16-8.5	10/8/2016	8.5	ND	--	--	--	0.045	ND	ND	ND	--	--	--	--	--	--	--	--	--
H-4-16-11	10/8/2016	11	ND	--	--	--	0.026	ND	ND	ND	--	--	--	--	--	--	--	--	--
H-4-16-16	10/8/2016	16	69	ND	ND	ND	ND	ND	0.55	1.89	0.092	0.36	0.56	1.8	0.064	0.053	0.29	ND	ND
H-4-16-18.5	10/8/2016	18.5	30	--	--	--	0.13	0.074	0.76	2.28	--	--	--	--	--	--	--	--	--
H-4-16-19.9	10/8/2016	19.9	99	--	--	--	0.35	0.090	1.4	2.79	--	--	--	--	--	--	--	--	--
H-4-16-25.4	10/8/2016	25.4	ND	--	--	--	0.092	0.064	ND	0.088	--	--	--	--	--	--	--	--	--
H-5-16-13.5	10/8/2016	13.5	--	ND	ND	ND	ND	ND	0.089	ND	0.19	1.2	0.15	ND	0.51	0.32	2.0	0.64	0.64
MTCA Cleanup Level			30	2	0.0015	48000	0.0017	0.27	0.34	1	8000	8000	800	No CUL	8000	No CUL	4000	0.24	0.24

-- Not Analyzed

ND Not Detected

## Exceeds Cleanup Level

H H - The result indicated is a common laboratory contaminant and may have been introduced during sample preparation.

CUL Cleanup Level



**Table 6 Semi Volatile Organic Compounds Detected in Soil Samples, SR 520 Eastbound Off-Ramp to Montlake Vicinity, Seattle, Washington**

Sample ID	Sample Date	Sample depth (ft.)	Napthalene (mg/kg)	2-Methylnaphthalene (mg/kg)	1-Methylnaphthalene (mg/kg)1
H-1-16-10	10/6/2016	10	ND	ND	ND
H-2-16-13.5	10/7/2016	13.5	0.0096	ND	ND
H-3-16-8.5	10/7/2016	8.5	ND	0.018	0.013
H-4-16-16	10/8/2016	16	<b>0.59</b>	0.74	0.37
H-5-16-13.5	10/8/2016	13.5	0.039	0.35	0.037
MTCA Cleanup Level			0.24	320	34

ND = Not Detected

**Table 7 Polychlorinated Biphenyls Detected in Soil Samples, SR 520 Eastbound Off-Ramp to Montlake Vicinity, Seattle, Washington**

<b>Sample ID</b>	<b>Sample Date</b>	<b>Sample depth (ft.)</b>	<b>Aroclor 1016 (mg/kg)</b>	<b>Aroclor 1212 (mg/kg)</b>	<b>Aroclor 1232 (mg/kg)</b>	<b>Aroclor 1242 (mg/kg)</b>	<b>Aroclor 1248 (mg/kg)</b>	<b>Aroclor 1254 (mg/kg)</b>	<b>Aroclor 1260 (mg/kg)</b>
H-1-16-10	10/6/2016	10	ND						
H-2-16-13.5	10/7/2016	13.5	ND						
H-3-16-3	10/7/2016	3	ND						
H-3-16-6	10/7/2016	6	ND						
H-3-16-8.5	10/7/2016	8.5	ND						
H-4-16-16	10/8/2016	16	ND						
H-5-16-13.5	10/8/2016	13.5	ND						

ND = Not Detected

**Table 8 Metals Detected in Soil Samples, SR 520 Eastbound Off-Ramp to Montlake Vicinity, Seattle, Washington**

Sample ID	Sample Date	Sample depth (ft.)	Chromium (mg/kg)	Copper (mg/kg)	Lead (mg/kg)	Nickel (mg/kg)	Zinc (mg/kg)
H-1-16-10	10/6/2016	10	36	8.4	ND	27	20
H-2-16-13.5	10/7/2016	13.5	37	24	11	36	56
H-3-16-3	10/7/2016	3	28	13	ND	33	26
H-3-16-6	10/7/2016	6	27	11	ND	30	24
H-3-16-8.5	10/7/2016	8.5	29	8.6	ND	24	26
H-4-16-16	10/8/2016	16	28	11	ND	30	24
H-5-16-13.5	10/8/2016	13.5	64	46	ND	69	69
MTCA Cleanup Level			2,000	14	150	6.53	299

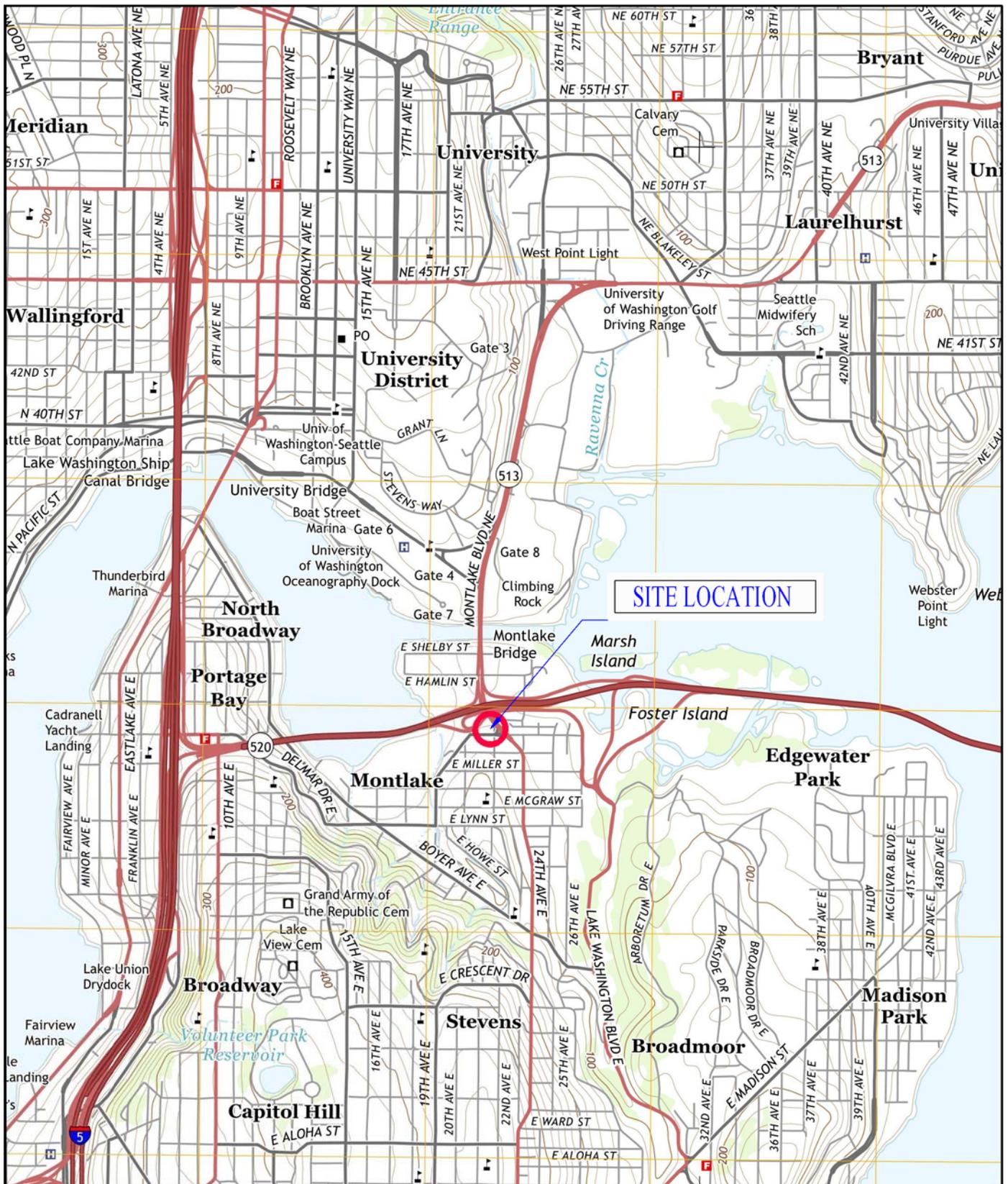
ND = Not Detected

## = Exceeds Cleanup Level



## FIGURES





**DESIGNED BY**  
 Innovex Environmental  
 Thuan Bui

**DRAWN BY**  
 ICD  
 October 5, 2016

0 1000 2000  
 SCALE IN FEET



16310 NE 80th St., Suite 300  
 Redmond, WA 98052  
 (800) 988-7880

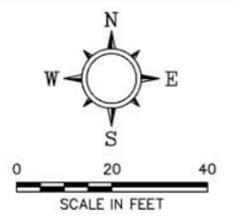
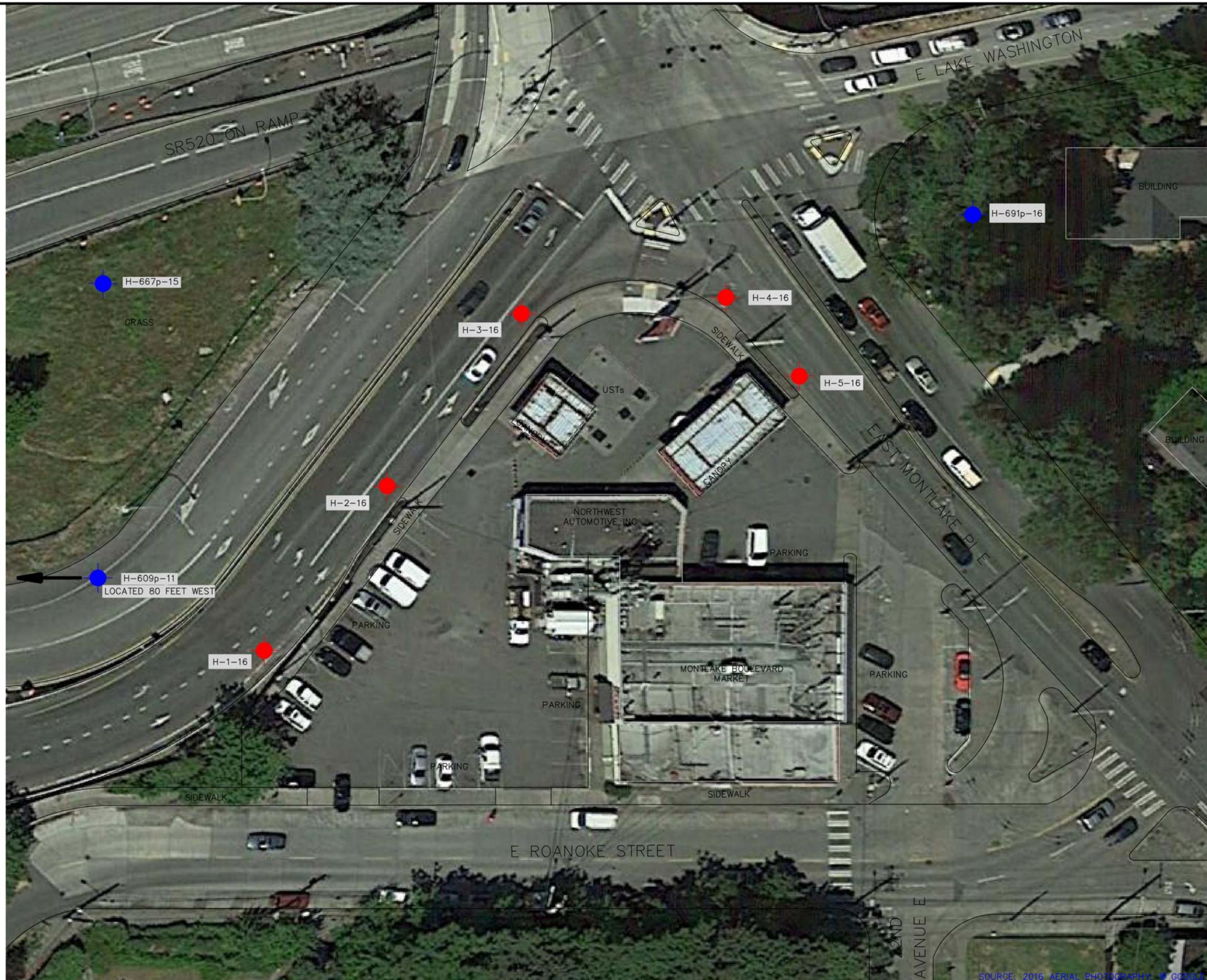
LATITUDE 47D 38M 38S NORTH  
 LONGITUDE 122D 18M 15S WEST

US GEOLOGICAL SURVEY - 2014  
 7.5 MINUTE QUADRANGLE MAP  
 SEATTLE NORTH, WASHINGTON

**FIGURE 1**  
**Site Location Map**

**STATE ROUTE (SR)520**  
**EASTBOUND OFF-RAMP**  
**TO MONTLAKE VICINITY**  
**2625 MONTLAKE PLACE EAST**  
**SEATTLE, WASHINGTON**





**LEGEND**

- H-1-16 SOIL BORING LOCATION
- H-691p-16 GEOTECHNICAL EXPLORATIONS

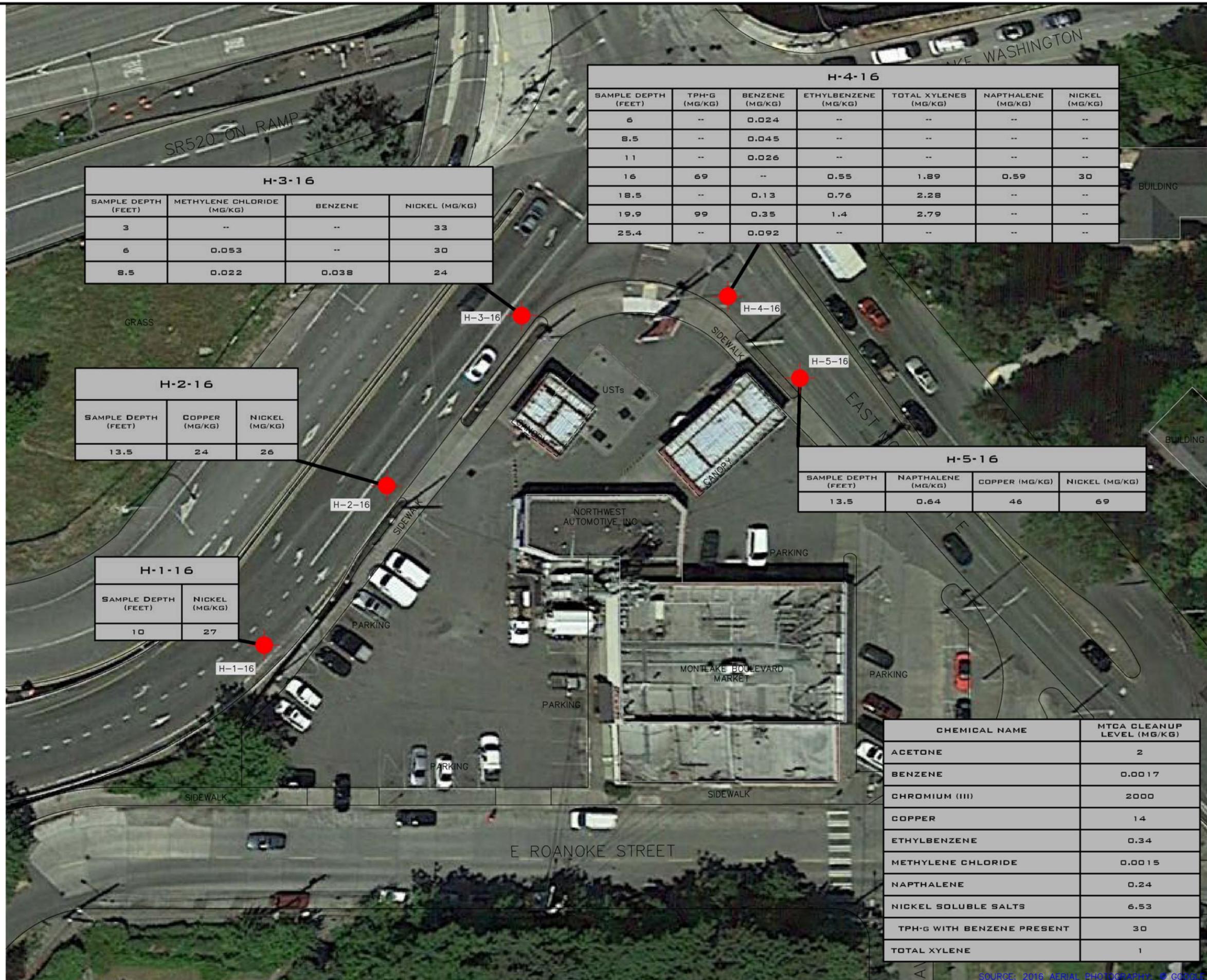
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 Innovex Environmental  
 Thuan Bui  
**DRAWN BY**  
 ICD  
 December 5, 2016

**FIGURE 2**  
**Boring Locations Map**  
 STATE ROUTE (SR) 520  
 EAST BOUND OFF-RAMP  
 TO MONTLAKE VICINITY  
 2625 MONTLAKE PLACE EAST  
 SEATTLE, WASHINGTON



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**H-3-16**

SAMPLE DEPTH (FEET)	METHYLENE CHLORIDE (MG/KG)	BENZENE	NICKEL (MG/KG)
3	--	--	33
6	0.053	--	30
8.5	0.022	0.038	24

**H-4-16**

SAMPLE DEPTH (FEET)	TPH-G (MG/KG)	BENZENE (MG/KG)	ETHYLBENZENE (MG/KG)	TOTAL XYLENES (MG/KG)	NAPHTHALENE (MG/KG)	NICKEL (MG/KG)
6	--	0.024	--	--	--	--
8.5	--	0.045	--	--	--	--
11	--	0.026	--	--	--	--
16	69	--	0.55	1.89	0.59	30
18.5	--	0.13	0.76	2.28	--	--
19.9	99	0.35	1.4	2.79	--	--
25.4	--	0.092	--	--	--	--

**H-2-16**

SAMPLE DEPTH (FEET)	COPPER (MG/KG)	NICKEL (MG/KG)
13.5	24	26

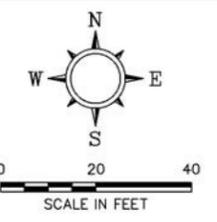
**H-5-16**

SAMPLE DEPTH (FEET)	NAPHTHALENE (MG/KG)	COPPER (MG/KG)	NICKEL (MG/KG)
13.5	0.64	46	69

**H-1-16**

SAMPLE DEPTH (FEET)	NICKEL (MG/KG)
10	27

CHEMICAL NAME	MTCA CLEANUP LEVEL (MG/KG)
ACETONE	2
BENZENE	0.0017
CHROMIUM (III)	2000
COPPER	14
ETHYLBENZENE	0.34
METHYLENE CHLORIDE	0.0015
NAPHTHALENE	0.24
NICKEL SOLUBLE SALTS	6.53
TPH-G WITH BENZENE PRESENT	30
TOTAL XYLENE	1



**LEGEND**

● H-1-16 SOIL BORING LOCATION

-- BELOW CLEANUP LEVEL

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Thuan Bui

**DRAWN BY**  
Thuan Bui  
November 1, 2016

**FIGURE 3**  
**Soil Sample Analytical Results Exceeding the MTCA Cleanup Levels**  
**STATE ROUTE (SR) 520 EASTBOUND OFF-RAMP TO MONTLAKE VICINITY 2625 MONTLAKE PLACE EAST SEATTLE, WASHINGTON**



16310 NE 80th St., Suite 300  
Redmond, WA 98052  
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**Appendix A**  
**Boring Logs**

---





Drilling Co. : WSDOT Job No. : 31008 Boring No. : 1-110  
 Drilling Rig Eq : \_\_\_\_\_ Job Name : SR 520 Exploration  
 Drilling Method : Mud Rotary Logged by : Jennifer Heise  
 Drill Diameter : \_\_\_\_\_ Location : Mentlake Exit Ramp  
 Weather Conditions : 55°F, rain, showers Start Date : 10/6/16 End Date : 10/6/16

Well Construction	Time	Blows/6 in.	Headspace PID/OVA	Sample ID (X = Lab Sample)	Depth (ft)	Sample Interval	USCS Code	FIELD CLASSIFICATION [Density/consistency, color, minor, MAJOR, then trace constituents; moisture; structure; other; (Geology USCS Classification)]
	2152							drilling began
		2			2			
		2						
		1						
		2						
			0.0		4			(4-6) medium brown silty sand w/ gravel, sample is <del>totally</del> graded, gravel is fine to 1/2" diameter, sub angular to subrounded, sample is damp, no odor, no debris
					6			
	2158				8			(9-11) top 1 foot is blue fine sand w/ little gravel, then lens of orange sand, then 1" of gray sand w/ silt and very little fine gravel; no odor; sample is damp
		1	0.0					
		11	0.0		10			
		11						
		13						
		20						
	2204				15			(15) very poor sample recovery (≈ 4") blue-gray fine sand w/ silt and gravel ≈ 1.5" diameter, dense, dry, no odor
		50/4	0.0					
	2223				20			(20) very poor sample recovery (≈ 6") similar to previous - blue-gray, partly graded sand w/ fine to 1" gravel, ≈ 3% gravel, subangular to subrounded; sample is dry; no odor observed
		50/3	0.0					

GROUNDWATER DATA			COMMENTS (i.e. materials used, visitors, problems etc.) :
Water Depth	Time	Date	

**SUMMARY OF TIME**

Boring/Sample : \_\_\_\_\_ hrs. Standby : \_\_\_\_\_ hrs.  
 Setup/Cleanup : \_\_\_\_\_ hrs. Decon : \_\_\_\_\_ hrs.  
 Boring No. : \_\_\_\_\_ Sheet \_\_\_\_\_ of \_\_\_\_\_

Drilling Co. : <u>WSDOT</u>	Job No. : _____	Boring No. : <u>H-1-16</u>
Drilling Rig Eq : _____	Job Name : _____	
Drilling Method : _____	Logged by : <u>Jennifer Haise</u>	
Drill Diameter : _____	Location : _____	
Weather Conditions : _____	Start Date : <u>161006</u>	End Date : <u>161006</u>

Well Construction	Time	Blows/6 in.	Headspace PID/OVA	Sample ID (X = Lab Sample)	Depth (ft)	Sample Interval	USCS Code	FIELD CLASSIFICATION [Density/consistency, color, minor, MAJOR, then trace constituents; moisture ; structure; other; (Geology USCS Classification)]
					21			
					22			
					23			
					24			24-25) (1 foot sample recovery) gray till-blue gray fine sand/silt layered w/ medium grained sand w/ gravel up to 3/4"; mostly fine gravel, up to 15% sample is dry, no odor
	2236	29 50/5			25			
					26			
					27			driller encountered gravel at 27'
					28			
					29			
	2248	50/5			30			(30) (poor sample recovery ≈ 4-5") till, blue gray fine sand w/ silt and gravel; poorly graded, gravel is fine to 3/4" diameter, subangular to subrounded, up to 35% gravel; sample is dense + dry
					31			
					32			
					33			
	2300	50/12			35			NO RECOVERY
					36			
					37			
					38			
					39			
	2313	50/12			40			NO RECOVERY

GROUNDWATER DATA			COMMENTS (i.e. materials used, visitors, problems etc.) :
Water Depth	Time	Date	

SUMMARY OF TIME	
Sample : _____ hrs.	Standby : _____ hrs.
_____ hrs.	Decon: _____ hrs.
Sheet _____ of _____	



# FIELD LOG OF BORING

Page 3 of 3

Drilling Co. : _____	Job No. : _____	Boring No. : <u>H-1-16</u>
Drilling Rig Eq : _____	Job Name : _____	
Drilling Method : _____	Logged by : _____	
Drill Diameter : _____	Location : _____	
Weather Conditions : _____	Start Date : <u>16/006</u>	End Date : <u>16/006</u>

Well Construction	Time	Blows/6 in.	Headspace PID/OVA	Sample ID (X = Lab Sample)	Depth (ft)	Sample Interval	USCS Code	FIELD CLASSIFICATION [Density/consistency, color, minor, MAJOR, then trace constituents; moisture ; structure; other; (Geology USCS Classification)]
					41			
	2330	50/B			45			(45) poor sample recovery (~2-3") fill-blue gray silty sand w/ little gravel, sample is dry and dense
		0.0						
	2350	50/B			50			(50) poor sample recovery (~2-3") fill-gray silty sand w/ some medium grained sand, very little fine gravel, no color, dense, dry
		0.0						
					55			boring terminated at 50' bgs

GROUNDWATER DATA			COMMENTS (i.e. materials used, visitors, problems etc.) :
Water Depth	Time	Date	
SUMMARY OF TIME			
Boring/Sample : _____ hrs.	Standby : _____ hrs.		
Setup/Cleanup : _____ hrs.	Decon: _____ hrs.		
Boring No. : _____	Sheet _____ of _____		



# FIELD LOG OF BORING

Page 1 of 2

H-2-16

Drilling Co. : <u>WSDOT</u>	Job No. : <u>31008</u>	Boring No. : _____
Drilling Rig Eq : _____	Job Name : <u>SR 520 @ Mountlake Place II</u>	
Drilling Method : <u>mid Rotary</u>	Logged by : <u>G. Hayward</u>	
Drill Diameter : _____	Location : <u>Seattle</u>	
Weather Conditions : <u>Raw showers</u>	Start Date : <u>10/7/16</u>	End Date : <u>10/7/16</u>

Well Construction	Time	Blows/6 in.	Headspace PID/OVA	Sample ID (X = Lab Sample)	Depth (ft)	Sample Interval	USCS Code	FIELD CLASSIFICATION [Density/consistency, color, minor, MAJOR, then trace constituents; moisture ; structure; other; (Geology USCS Classification)]
	21:20	2	0.0		1.5-3		SM	Asphalt Gray silty sand w/few gravel, moist, loose, no odor, PID=0.0
		2						
		2						
		2			4-6			no recovery
		1						
		2						
		1						
	21:35	2	0.7		7-8.5		SM	Gray silty sand, moist, v. loose, no odor, PID=0.7
		1						
	21:40	1	3.9		9-11			same, petroleum odor
		1						
		2						
		1						
	21:45	2	4.7		12-13.5		SM	same as above PID=4.7 wood debris @ 13 feet, petroleum odor
		1						
	21:50	2	0.0		14-16			Dark gray silty sand, stiff
		2						
		3						
		7						
	21:55	16	0.0		17-18.5			Gray silty sand, dense, moist, no odor
		20						
		31						
		50/4	10.3		19-19.3			Gray silty sand, very dense no odor
	22:00							

GROUNDWATER DATA			COMMENTS (i.e. materials used; visitors, problems etc.) :
Water Depth	Time	Date	

SUMMARY OF TIME	
Boring/Sample : _____ hrs.	Standby : _____ hrs.
Setup/Cleanup : _____ hrs.	Decon: _____ hrs.
Boring No. : _____	Sheet _____ of _____





# FIELD LOG OF BORING

14-3-16

Drilling Co. : <u>WSDOT</u>	Job No. : <u>31008</u>	Boring No. : _____
Drilling Rig Eq : _____	Job Name : <u>SR 520 @ E. Montlake Phase II</u>	
Drilling Method : <u>Mudrotary</u>	Logged by : <u>G. Hayman</u>	
Drill Diameter : _____	Location : <u>Seattle, WA</u>	
Weather Conditions : _____	Start Date : <u>10/7/16</u>	End Date : <u>10/8/16</u>

Well Construction	Time	Blows/6 in.	Headspace PID/OVA	Sample ID (X = Lab Sample)	Depth (ft)	Sample Interval	USCS Code	FIELD CLASSIFICATION [Density/consistency, color, minor, MAJOR, then trace constituents; moisture ; structure; other; (Geology USCS Classification)]
	23:25	7	10.5		1.5-3		SP	Asphalt Sand w/ fine gravel, moist, loose, petroleum odor PID=10.5
	23:30	3	21.5		4-6		SM	Gray silty sand, moist, very loose, petroleum odor, PID:21.5
	23:35	1	35.4		7-8.5			Brown silty fine sand, few fine gravel very loose, petroleum odor, PID=35.4
	23:45	5	-					NO recovery
	23:50	23	0		12-13.5		SM	Brown silty sand, wet, dense, no odor PID=0
	0:50	56.4	0		17-17.4			Bailed boring dry, allowed water to recover, 00:25 collected groundwater sample Silty sand w/ few fine gravel, no odor PID=0
								TD=17.4 feet

GROUNDWATER DATA			COMMENTS (i.e. materials used; visitors, problems etc.) :
Water Depth	Time	Date	

SUMMARY OF TIME	
Boring/Sample : _____ hrs.	Standby : _____ hrs.
Setup/Cleanup : _____ hrs.	Decon: _____ hrs.
Boring No. : _____	Sheet _____ of _____



# FIELD LOG OF BORING

Page 1 of 2

H-4-16

Drilling Co. : <u>WSDOT</u>	Job No. : <u>310080</u>	Boring No. : _____
Drilling Rig Eq : _____	Job Name : <u>SB 520 @ E. Montlake Phase II</u>	
Drilling Method : <u>Mud rotary</u>	Logged by : <u>G. Hayward</u>	
Drill Diameter : <u>3"</u>	Location : <u>Seattle, WA</u>	
Weather Conditions : <u>Rain showers</u>	Start Date : <u>10/8/16</u>	End Date : <u>10/9/16</u>

Well Construction	Time	Blows/6 in.	Headspace PID/OVA	Sample ID (X = Lab Sample)	Depth (ft)	Sample Interval	USCS Code	FIELD CLASSIFICATION [Density/consistency, color, minor, MAJOR, then trace constituents; moisture ; structure; other; (Geology USCS Classification)]
	21:45	4	20.2		1.5-3		SM	Asphalt gray silty sand, damp, loose, Petroleum odor, PID=20.2
	21:50	3	17		4-6			Gray silty sand, damp, very loose, petroleum odor, PID=17
	21:55	1	15.6		7-8.5			gray silty sand, very loose, damp, Petroleum odor, PID=15.6
	22:00		43		9-11		SM	Gray-brown silty sand, very stiff. strong Petroleum odor
		16						No recovery
	22:05	9	720				SM	gray silty sand, wet, medium dense, very strong petroleum odor PID=720
	22:10	16	380		17-18.5			same as above
	22:15	18	120		19-19.9			gray silty sand, very dense, petroleum odor, 2" till in tip of split spoon

GROUNDWATER DATA			COMMENTS (i.e. materials used; visitors, problems etc.) :
Water Depth	Time	Date	

SUMMARY OF TIME	
Boring/Sample : _____ hrs.	Standby : _____ hrs.
Setup/Cleanup : _____ hrs.	Decon: _____ hrs.
Boring No. : _____	Sheet _____ of _____



# FIELD LOG OF BORING

Page 2 of 2

H-4-16

Drilling Co. : <u>WSDOT</u>	Job No. : <u>31608</u>	Boring No. : _____
Drilling Rig Eq : _____	Job Name : <u>SR 520 @ E. Mount Lake Phase II</u>	
Drilling Method : <u>Mud Rotary</u>	Logged by : <u>G. Huggins</u>	
Drill Diameter : <u>3 inch</u>	Location : _____	
Weather Conditions : <u>rain showers</u>	Start Date : <u>10/8/16</u>	End Date : <u>10/8/16</u>

Well Construction	Time	Blows/6 in.	Headspace PID/OVA	Sample ID (X = Lab Sample)	Depth (ft)	Sample Interval	USCS Code	FIELD CLASSIFICATION [Density/consistency, color, minor, MAJOR, then trace constituents; moisture ; structure; other; (Geology USCS Classification)]
					21			
	22:35	26 45 50/5	3.4		24- 25.4		SM	Gray silty very fine sand, very dense no odor, damp PID=3.4
	22:45	50/2	0.3		29- 29.2		SM	Well graded gravel, sand and silt, very dense, PID=0.3  TD = 29.2 FT

GROUNDWATER DATA			COMMENTS (i.e. materials used; visitors, problems etc.) :
Water Depth	Time	Date	

SUMMARY OF TIME	
Boring/Sample : _____ hrs.	Standby : _____ hrs.
Setup/Cleanup : _____ hrs.	Decon: _____ hrs.
Boring No. : _____	Sheet _____ of _____



# FIELD LOG OF BORING

Page 1 of 2  
H-5-16

Drilling Co. : <u>WSDOT</u>	Job No. : <u>3100P</u>	Boring No. : _____
Drilling Rig Eq : _____	Job Name : <u>SR520 @ E. Montlake Phase II</u>	
Drilling Method : <u>Mud rotary</u>	Logged by : <u>G. Hayman</u>	
Drill Diameter : <u>3"</u>	Location : <u>Seattle, WA</u>	
Weather Conditions : <u>Rain shower</u>	Start Date : <u>10/8/16</u>	End Date : <u>10/9/16</u>

Well Construction	Time	Blows/6 in.	Headspace PID/OVA	Sample ID (X = Lab Sample)	Depth (ft)	Sample Interval	USCS Code	FIELD CLASSIFICATION [Density/consistency, color, minor, MAJOR, then trace constituents; moisture ; structure; other; (Geology USCS Classification)]
	23:55	6	3.5		1.5-3		SM	8" Asphalt & concrete gray silty sand w/ few gravel, moist loose, no odor PID=3.5
	23:59	4	16.5		4-6		SM	gray silty sand, few gravel, loose petroleum odor, PID=16.5
	0:05	1	1.6		7-8.5		SM	gray silty fine sand w/ few <sup>fine</sup> gravel loose, no odor PID=1.6
	0:10	3	16		9-11		SM	gray silty sand, dense, petroleum odor, PID=16
	0:15	2	380		12-13.5		SM	same as above, strong petroleum odor PID=380
	0:20	11	25		14-16		SM	light gray silty sand, medium dense & petroleum odor, PID=25
	0:25	8	2.5		17-18.5		SM	silty sand w/ few gravel, medium dense, no odor, PID=2.5
	0:30	8	1.1		19-21		SM	same as above

GROUNDWATER DATA			COMMENTS (i.e. materials used; visitors, problems etc.) :
Water Depth	Time	Date	

SUMMARY OF TIME	
Boring/Sample : _____ hrs.	Standby : _____ hrs.
Setup/Cleanup : _____ hrs.	Decon: _____ hrs.
Boring No. : _____	Sheet _____ of _____



# FIELD LOG OF BORING

Page 2 of 2

H-5-16

Drilling Co. : <u>WSDOT</u>	Job No. : <u>3100A</u>	Boring No. : _____
Drilling Rig Eq : _____	Job Name : <u>SR 520 @ E. Montlake Phase II</u>	
Drilling Method : <u>Mud Rotary</u>	Logged by : <u>G. Hayman</u>	
Drill Diameter : <u>3"</u>	Location : <u>Seattle, WA</u>	
Weather Conditions : <u>Rain shower</u>	Start Date : <u>10/8/16</u>	End Date : <u>10/9/16</u>

Well Construction	Time	Blows/6 in.	Headspace PID/OVA	Sample ID (X = Lab Sample)	Depth (ft)	Sample Interval	USCS Code	FIELD CLASSIFICATION [Density/consistency, color, minor, MAJOR, then trace constituents; moisture ; structure; other; (Geology USCS Classification)]
	0:35	10 11 12			22- 23.5		SM	gray silty fine sand, medium dense, no odor, PID=0.4
	0:40	5/0.0 14			24- 24.3			gray brown silty sand w/few gravel very dense, no odor, PID=0.0
	0:45	5/0.0 12			24- 24.2			same as above, no odor PID=0.0
								TD = 29.2

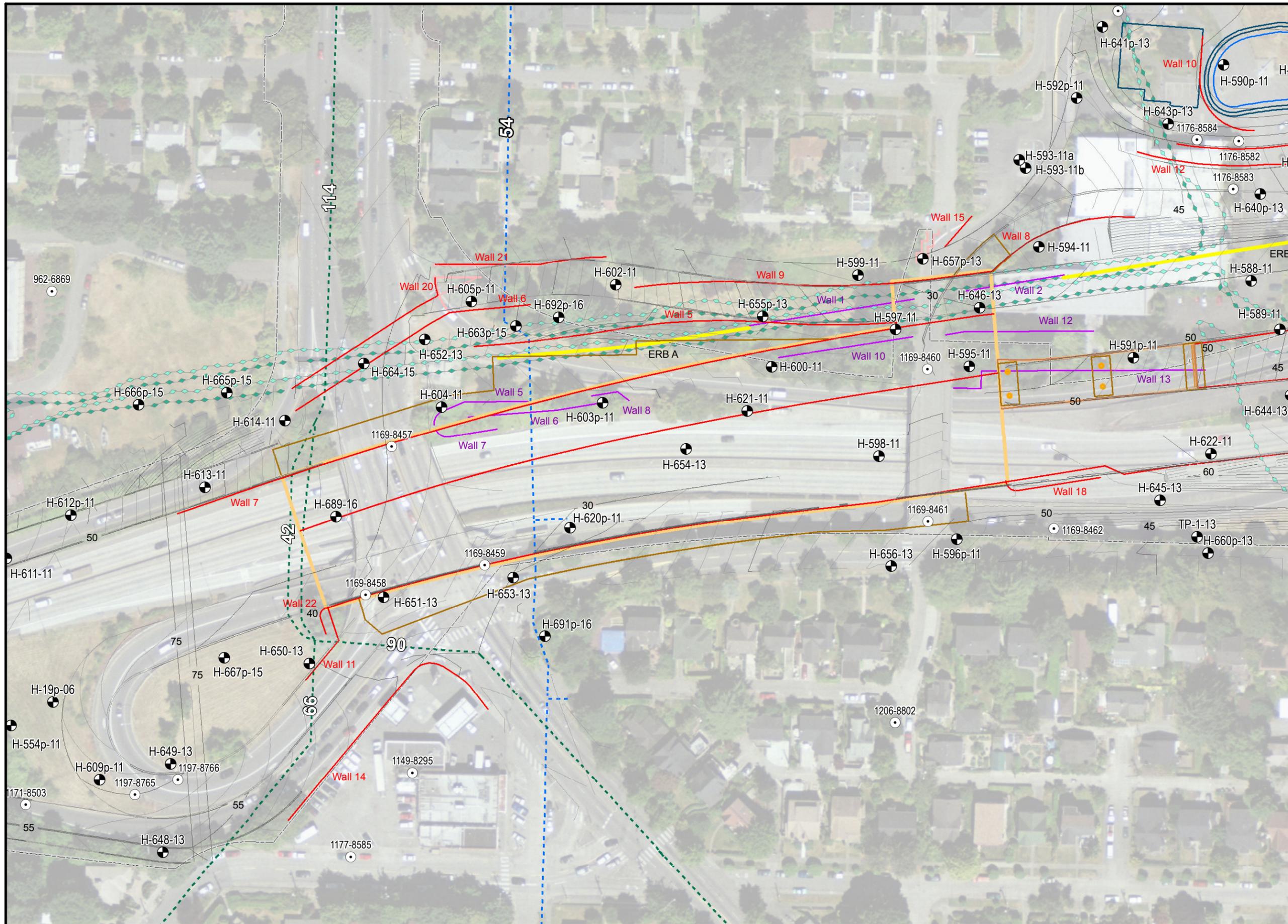
GROUNDWATER DATA			COMMENTS (i.e. materials used; visitors, problems etc.) :
Water Depth	Time	Date	

SUMMARY OF TIME	
Boring/Sample : _____ hrs.	Standby : _____ hrs.
Setup/Cleanup : _____ hrs.	Decon: _____ hrs.
Boring No. : _____	Sheet _____ of _____

**Appendix B**  
**WSDOT Boring Logs and Water Level Figures**

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- ### LEGEND
- RECENT EXPLORATIONS
  - HISTORIC EXPLORATIONS
  - Proposed Pier Location & Designation
  - Proposed Alignment
  - Proposed Structure
  - Proposed Retaining Wall
  - Proposed Bridge
  - Proposed Stormwater Facility
  - Geophysical Survey & Designation
  - WSDOT ROW
  - City of Seattle ROW
  - Historic Lake Washington Ship Canal & Shoreline
  - Historic Canal & Shoreline from WSDOT
  - Landslide Area

- ### NOTES
1. WSDOT = Washington State Department Of Transportation. ROW = Right Of Way
  2. Proposed Alignment Features from WSDOT, 2014-09-08.
  3. 'Historic Canal & Shoreline' digitized from WSDOT, n.d., Digitized Lake Union Plat map: AutoCAD drawing received from WSDOT May 16, 2013
  4. "Historic Lake Washington Ship Canal & Shoreline" from U. S. Army, Seattle District Engineer, 1952, Realestate - Lake Washington Ship Canal and Hiram M. Chittenden Locks: 1 sheet, October 22.



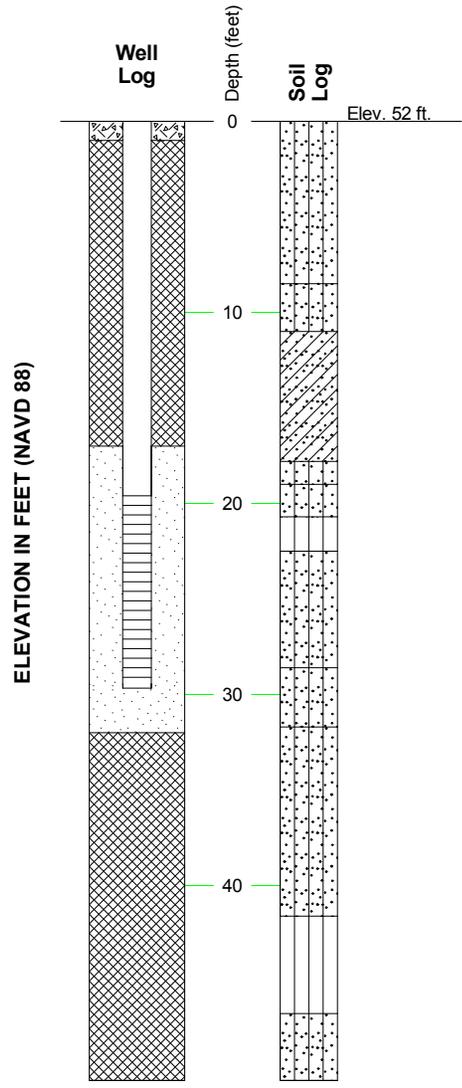
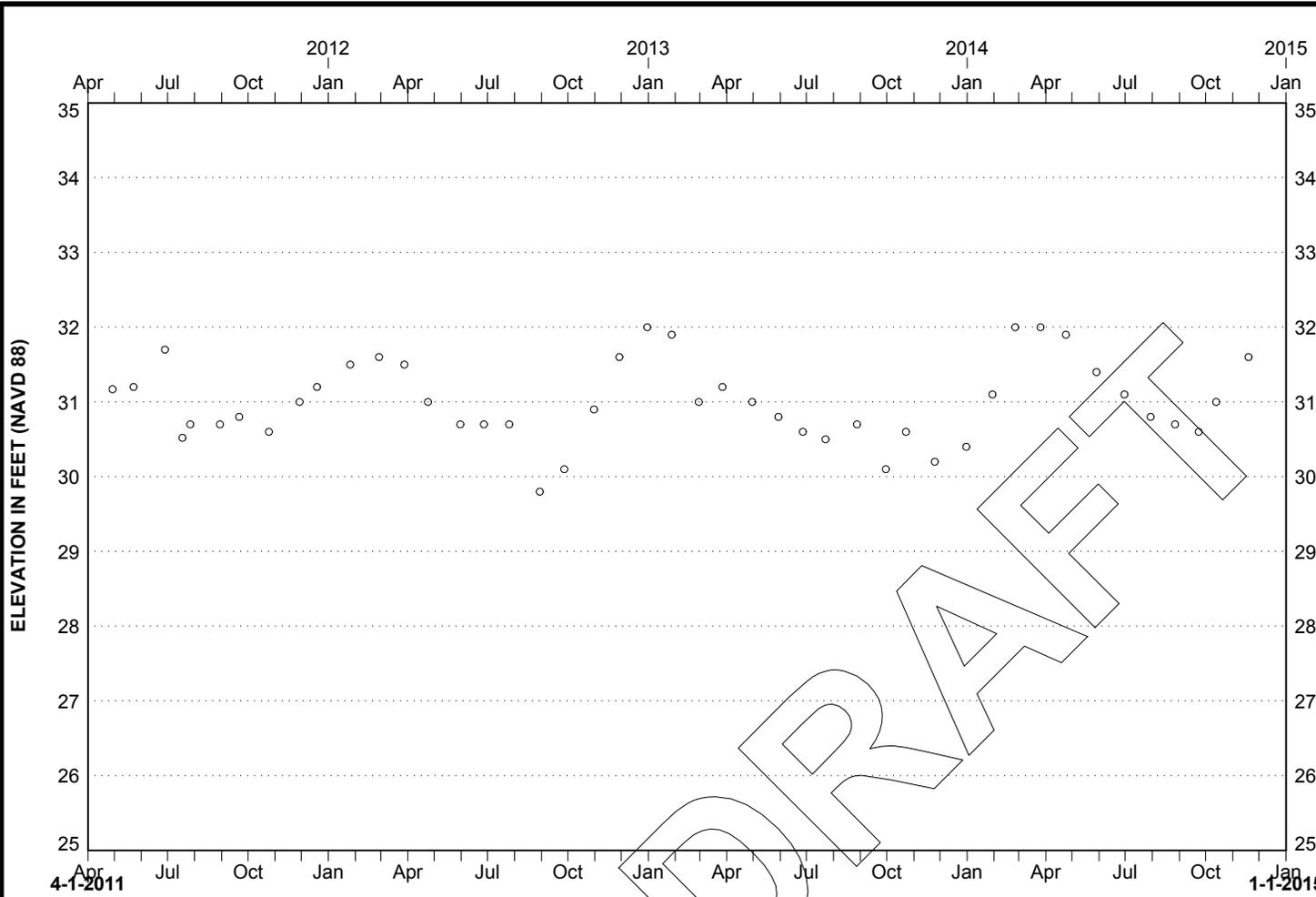
SR 520, I-5 to Medina  
Westside Final Conceptual Design  
Seattle, Washington

## SITE AND EXPLORATION PLAN

August 2016 21-1-20624-807

**SHANNON & WILSON, INC.**  
GEOTECHNICAL AND ENVIRONMENTAL CONSULTANTS





Well Identification	Geologic Unit(s) in VWP or OW Depth Range	Read Point (feet)		Water Level Elevation (feet)	
		Depth	Elevation	Low	High
○ H-609p-11 OW1	Qvat, Qvd	29.5	22.5	29.8	32.0

- NOTES:**
1. Refer to Interpreted Geologic Units and Descriptions exhibit for more information on geologic units.
  2. Read Point Depth or Elevation (as listed in table above) is equal to the depth or elevation of the tip of the vibrating wire pressure (VWP) transducer or the bottom of the screen of the observation well (OW). Measured groundwater elevations less than these values are considered as "Dry."
  3. WSDOT collected monthly groundwater measurements from April 2011 to December 2014.
  4. Well was destroyed January 2015.

SR 520 Bridge Replacement and HOV Program  
 Seattle Vulnerable Structures  
 Seattle, Washington

**GROUNDWATER LEVEL MEASUREMENTS**  
**BORING H-609p-11**  
 May 2015 21-1-20624-807

**SHANNON & WILSON, INC.**  
 Geotechnical and Environmental Consultants

**EXHIBIT 27**

EXHIBIT 27

SR520\_GW\_PL01\_ELEVANI\_21-20624\_JMW.EDIT.GPJ\_21-20624.GPJ\_5/17/15



Instrument	Screen Bottom Elev (ft)	Geologic Unit
OW1	29.5	Hf, Qpgt

<span style="color: green;">—</span>	OW1
<span style="color: red;">- - -</span>	OW1 (Dry)
<span style="color: green;">○</span>	OW1 Manual
<span style="color: red;">□</span>	OW1 Manual (Dry)

SR 520 Bridge Replacement and HOV Program  
I-5 to Lake Washington  
Seattle, Washington

**GROUNDWATER LEVEL  
MEASUREMENTS  
BORING H-667P-15**

July 2016

21-1-22242-005

**SHANNON & WILSON, INC.**  
Geotechnical and Environmental Consultants

**EXHIBIT 36**

**Notes:**

1. Groundwater data provided by WSDOT
2. Measurements are shown as dry for readings below instrument elevation of pressure transducer or bottom of screen interval
3. Refer to Shannon & Wilson boring logs for soil descriptions and well construction details

ver:0.7.7\_posi62\_dev0-q78fa448\_bv:MDG



Instrument	Screen Bottom Elev (ft)	Geologic Unit
OW1	34.0	QPGT, QPGL

<span style="color: green;">—</span> OW1
<span style="color: red;">- -</span> OW1 (Dry)

SR 520 Bridge Replacement and HOV Program  
I-5 to Lake Washington  
Seattle, Washington

**GROUNDWATER LEVEL MEASUREMENTS BORING H-691P-16**

July 2016 21-1-22242-005

**SHANNON & WILSON, INC.** Geotechnical and Environmental Consultants **EXHIBIT 47**

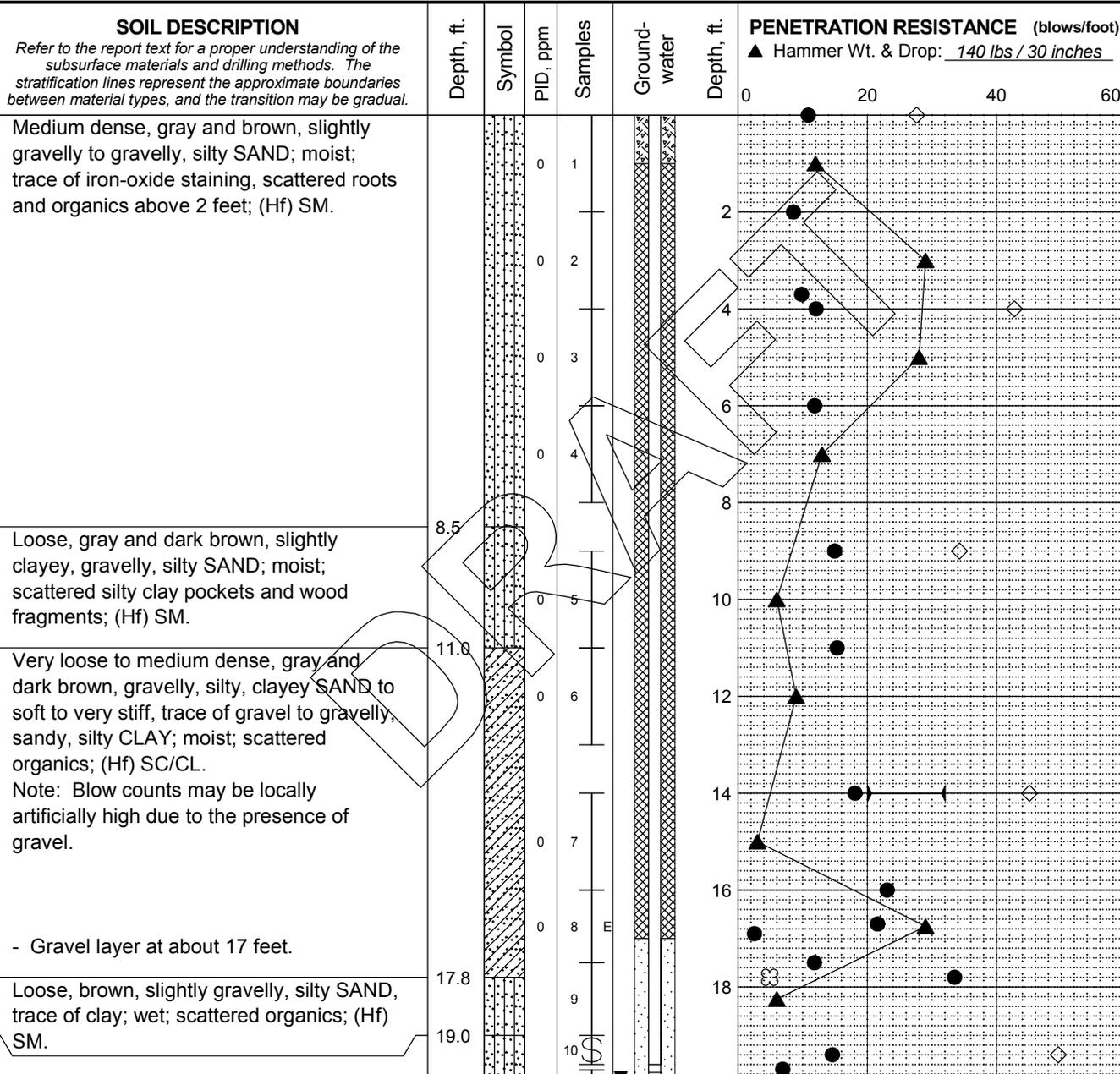
- Notes:
1. Groundwater data provided by WSDOT
  2. Measurements are shown as dry for readings below instrument elevation of pressure transducer or bottom of screen interval
  3. Refer to Shannon & Wilson boring logs for soil descriptions and well construction details

**EXHIBIT 47**

C:\Users\mndt\repos\qwd\plots\plots.py



Total Depth: <u>50.2 ft.</u>	Northing: <u>238,178 ft.</u>	Drilling Method: <u>Advanced Casing</u>	Hole Diam.: <u>5 in.</u>
Top Elevation: <u>52 ft.</u>	Easting: <u>1,277,493 ft.</u>	Drilling Company: <u>WSDOT</u>	Rod Diam.: <u>1 3/4-inch</u>
Vert. Datum: <u>NAVD 88</u>	Station: _____	Drill Rig Equipment: <u>CME 850 Track Rig 9C2-3</u>	Hammer Type: <u>Automatic</u>
Horiz. Datum: <u>NAD 83/91</u>	Offset: _____	Other Comments: _____	Hammer ER: <u>88 %</u>



CONTINUED NEXT SHEET

**LEGEND**

- \* Sample Not Recovered
- E Environmental Sample Obtained
- [Symbol] Standard Penetration Test
- [Symbol] 3" O.D. Thin-Walled Tube
- ▼ Groundwater Level in Well

- ◇ % Fines (<0.075mm)
- % Water Content
- Plastic Limit
- Liquid Limit
- Natural Water Content

**NOTES**

1. Refer to KEY for explanation of symbols, codes, abbreviations and definitions.
2. The stratification lines represent the approximate boundaries between soil types, and the transition may be gradual.
3. Discussions in the text of the geotechnical data reports are necessary for a proper understanding of the nature of the subsurface materials.
4. Groundwater level, if indicated above, is for the date specified and may vary. Groundwater level is the highest available measurement to date. Groundwater plots contain complete data sets.
5. USCS designation is based on visual-manual classification and selected lab testing.
6. Hammer ER = hammer energy ratio (efficiency) as a percentage.

SR 520 Bridge Replacement and HOV Program  
West Approach Bridge  
Seattle, Washington

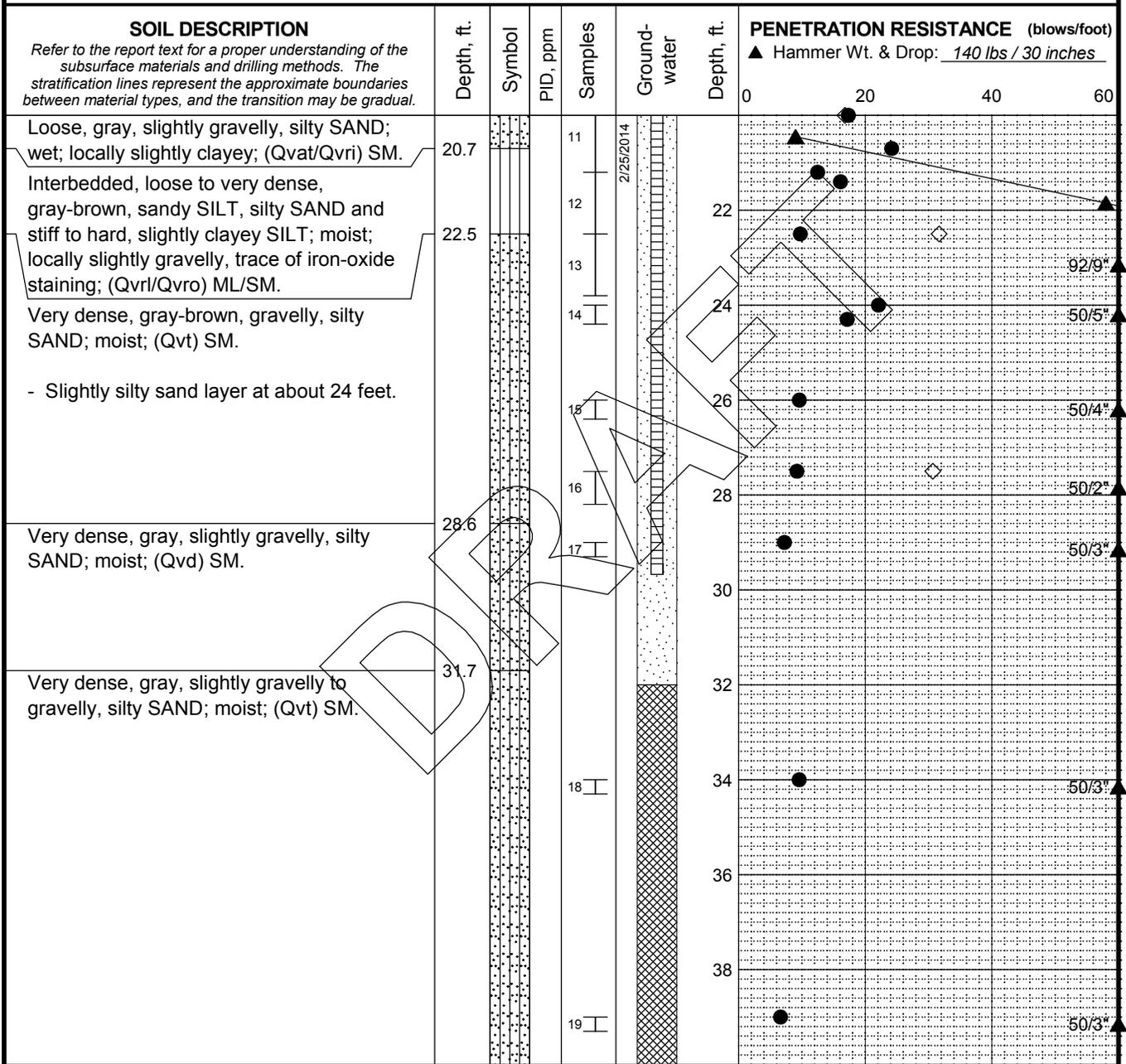
**LOG OF BORING H-609p-11**

June 2015 21-1-20624-807

<b>SHANNON &amp; WILSON, INC.</b> Geotechnical and Environmental Consultants	<b>EXHIBIT 132</b> Sheet 1 of 3
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SR520\_MASTER\_LOG\_E\_SMP\_21-20624-GPJ\_SHAN\_WIL\_GDT\_9/14/16 Log: BRC Rev: JKP Typ: CLP

Total Depth: 50.2 ft. Northing: 238,178 ft. Drilling Method: Advanced Casing Hole Diam.: 5 in.  
 Top Elevation: 52 ft. Easting: 1,277,493 ft. Drilling Company: WSDOT Rod Diam.: 1 3/4-inch  
 Vert. Datum: NAVD 88 Station: \_\_\_\_\_ Drill Rig Equipment: CME 850 Track Rig 9C2-3 Hammer Type: Automatic  
 Horiz. Datum: NAD 83/91 Offset: \_\_\_\_\_ Other Comments: \_\_\_\_\_ Hammer ER: 88 %



CONTINUED NEXT SHEET

**LEGEND**

- \* Sample Not Recovered
- E Environmental Sample Obtained
- Standard Penetration Test
- 3" O.D. Thin-Walled Tube
- Groundwater Level in Well

- % Organics
- % Fines (<0.075mm)
- % Water Content
- Plastic Limit
- Liquid Limit
- Natural Water Content

**NOTES**

1. Refer to KEY for explanation of symbols, codes, abbreviations and definitions.
2. The stratification lines represent the approximate boundaries between soil types, and the transition may be gradual.
3. Discussions in the text of the geotechnical data reports are necessary for a proper understanding of the nature of the subsurface materials.
4. Groundwater level, if indicated above, is for the date specified and may vary. Groundwater level is the highest available measurement to date. Groundwater plots contain complete data sets.
5. USCS designation is based on visual-manual classification and selected lab testing.
6. Hammer ER = hammer energy ratio (efficiency) as a percentage.

SR 520 Bridge Replacement and HOV Program  
 West Approach Bridge  
 Seattle, Washington

**LOG OF BORING H-609p-11**

June 2015

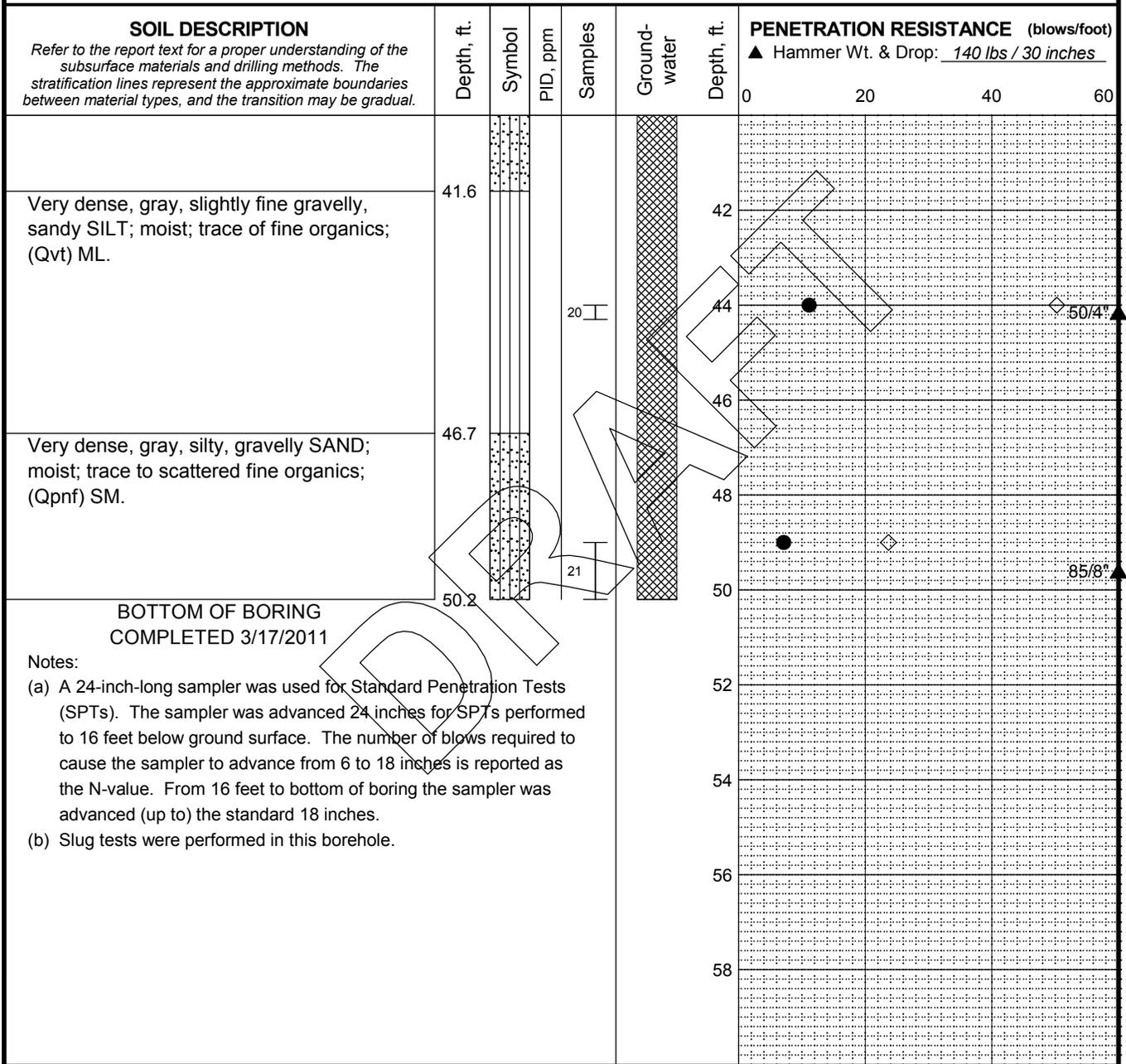
21-1-20624-807

**SHANNON & WILSON, INC.**  
 Geotechnical and Environmental Consultants

**EXHIBIT 132**  
 Sheet 2 of 3

SR520\_MASTER\_LOG\_E\_SMP\_21-20624-GPJ\_SHAN\_WIL\_GDT\_9/14/16 Log: BRC Rev: JKP Typ: CLP

Total Depth: 50.2 ft. Northing: 238,178 ft. Drilling Method: Advanced Casing Hole Diam.: 5 in.  
 Top Elevation: 52 ft. Easting: 1,277,493 ft. Drilling Company: WSDOT Rod Diam.: 1 3/4-inch  
 Vert. Datum: NAVD 88 Station: \_\_\_\_\_ Drill Rig Equipment: CME 850 Track Rig 9C2-3 Hammer Type: Automatic  
 Horiz. Datum: NAD 83/91 Offset: \_\_\_\_\_ Other Comments: \_\_\_\_\_ Hammer ER: 88 %



**Notes:**

- (a) A 24-inch-long sampler was used for Standard Penetration Tests (SPTs). The sampler was advanced 24 inches for SPTs performed to 16 feet below ground surface. The number of blows required to cause the sampler to advance from 6 to 18 inches is reported as the N-value. From 16 feet to bottom of boring the sampler was advanced (up to) the standard 18 inches.
- (b) Slug tests were performed in this borehole.

**LEGEND**

- \* Sample Not Recovered
- E Environmental Sample Obtained
- ┌ Standard Penetration Test
- └ 3" O.D. Thin-Walled Tube
- ▼ Groundwater Level in Well

**NOTES**

1. Refer to KEY for explanation of symbols, codes, abbreviations and definitions.
2. The stratification lines represent the approximate boundaries between soil types, and the transition may be gradual.
3. Discussions in the text of the geotechnical data reports are necessary for a proper understanding of the nature of the subsurface materials.
4. Groundwater level, if indicated above, is for the date specified and may vary. Groundwater level is the highest available measurement to date. Groundwater plots contain complete data sets.
5. USCS designation is based on visual-manual classification and selected lab testing.
6. Hammer ER = hammer energy ratio (efficiency) as a percentage.

SR 520 Bridge Replacement and HOV Program  
 West Approach Bridge  
 Seattle, Washington

**LOG OF BORING H-609p-11**

June 2015

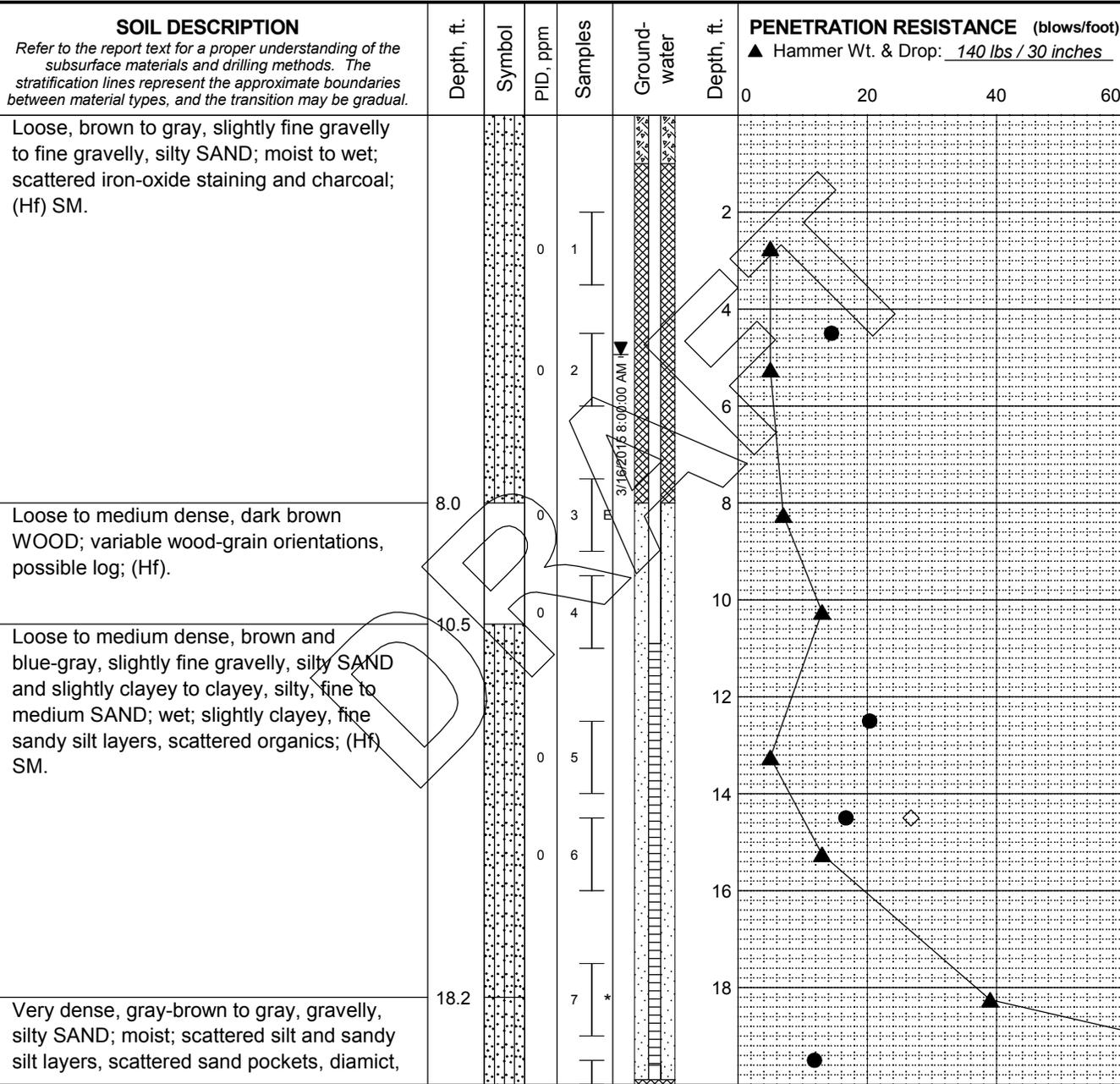
21-1-20624-807

**SHANNON & WILSON, INC.**  
 Geotechnical and Environmental Consultants

**EXHIBIT 132**  
 Sheet 3 of 3

SR520\_MASTER\_LOG\_E\_SMP\_21-20624-GPJ\_SHAN\_WIL\_GDT\_9/14/16 Log: BRC Rev: JKP Typ: CLP

Total Depth: 50.2 ft. Northing: 238,310 ft. Drilling Method: Advanced Casing Hole Diam.: 5 in.  
 Top Elevation: 49 ft. Easting: 1,277,628 ft. Drilling Company: WSDOT Rod Diam.: 1 3/4-inch  
 Vert. Datum: NAVD 88 Station: \_\_\_\_\_ Drill Rig Equipment: CME LC55 Track Rig 9C7-1 Hammer Type: Automatic  
 Horiz. Datum: NAD 83/91 Offset: \_\_\_\_\_ Other Comments: \_\_\_\_\_ Hammer ER: 81 %



CONTINUED NEXT SHEET

**LEGEND**

- \* Sample Not Recovered
- E Environmental Sample Obtained
- C Corrosion and/or Cation Exchange Sample Obtained
- [Symbol] Standard Penetration Test
- [Symbol] 2.5" O.D. Split Spoon Sample
- ▼ Groundwater Level in Well
- ◇ % Fines (<0.075mm)
- % Water Content
- Liquid Limit
- Natural Water Content

**NOTES**

1. Refer to KEY for explanation of symbols, codes, abbreviations and definitions.
2. The stratification lines represent the approximate boundaries between soil types, and the transition may be gradual.
3. Discussions in the text of the geotechnical data reports are necessary for a proper understanding of the nature of the subsurface materials.
4. Groundwater level, if indicated above, is for the date specified and may vary. Groundwater level is the highest available measurement to date. Groundwater plots contain complete data sets.
5. USCS designation is based on visual-manual classification and selected lab testing.
6. Hammer ER = hammer energy ratio (efficiency) as a percentage.

SR 520 Bridge Replacement and HOV Program  
West Approach to Montlake Lid  
Seattle, Washington

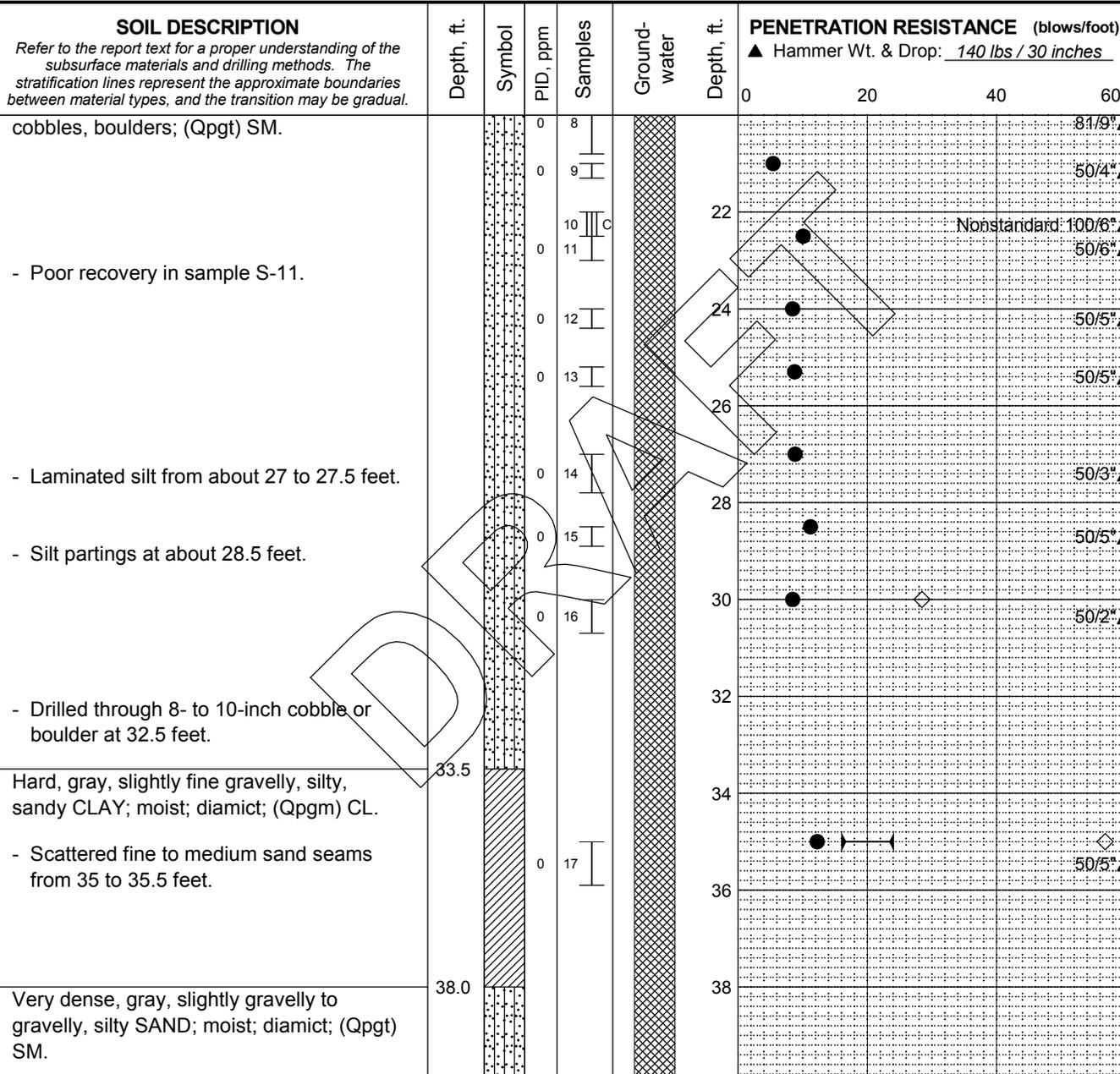
**LOG OF BORING H-667p-15**

September 2016 21-1-22242-002

<b>SHANNON &amp; WILSON, INC.</b> Geotechnical and Environmental Consultants	<b>EXHIBIT A-73</b> Sheet 1 of 3
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SR520\_MASTER\_LOG\_E\_SMP\_21-20624.GPJ\_SHAN\_WIL\_GDT\_9/14/16 Log: JMW Rev: DPO Typ: CLP

Total Depth: 50.2 ft. Northing: 238,310 ft. Drilling Method: Advanced Casing Hole Diam.: 5 in.  
 Top Elevation: 49 ft. Easting: 1,277,628 ft. Drilling Company: WSDOT Rod Diam.: 1 3/4-inch  
 Vert. Datum: NAVD 88 Station: \_\_\_\_\_ Drill Rig Equipment: CME LC55 Track Rig 9C7-1 Hammer Type: Automatic  
 Horiz. Datum: NAD 83/91 Offset: \_\_\_\_\_ Other Comments: \_\_\_\_\_ Hammer ER: 81 %



CONTINUED NEXT SHEET

**LEGEND**

- \* Sample Not Recovered
- E Environmental Sample Obtained
- C Corrosion and/or Cation Exchange Sample Obtained
- ⊥ Standard Penetration Test
- ⊥ 2.5" O.D. Split Spoon Sample
- ▼ Groundwater Level in Well
- ◇ % Fines (<0.075mm)
- % Water Content
- Plastic Limit
- Liquid Limit
- Natural Water Content

**NOTES**

1. Refer to KEY for explanation of symbols, codes, abbreviations and definitions.
2. The stratification lines represent the approximate boundaries between soil types, and the transition may be gradual.
3. Discussions in the text of the geotechnical data reports are necessary for a proper understanding of the nature of the subsurface materials.
4. Groundwater level, if indicated above, is for the date specified and may vary. Groundwater level is the highest available measurement to date. Groundwater plots contain complete data sets.
5. USCS designation is based on visual-manual classification and selected lab testing.
6. Hammer ER = hammer energy ratio (efficiency) as a percentage.

SR 520 Bridge Replacement and HOV Program  
West Approach to Montlake Lid  
Seattle, Washington

**LOG OF BORING H-667p-15**

September 2016 21-1-22242-002

**SHANNON & WILSON, INC.**  
Geotechnical and Environmental Consultants

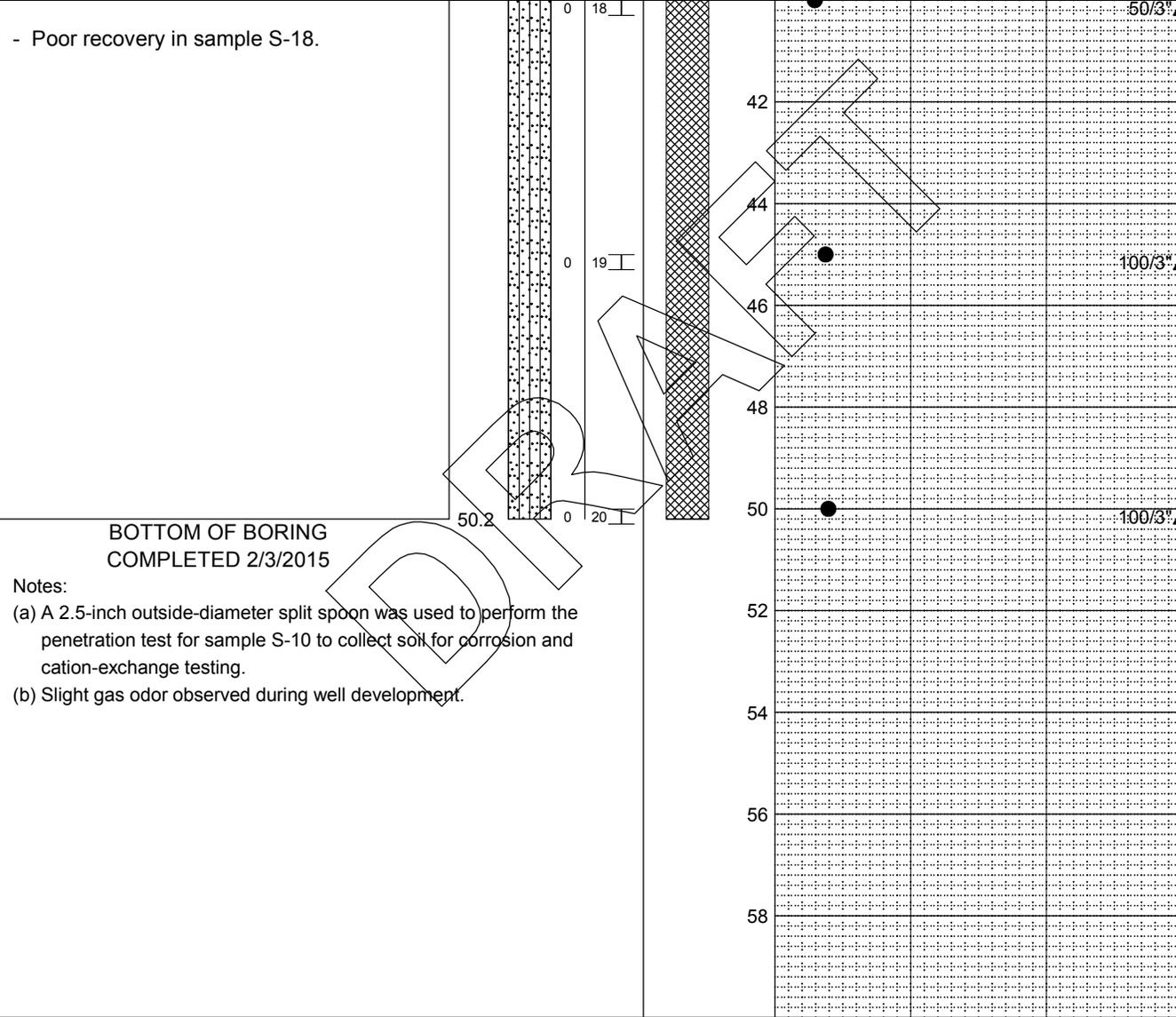
**EXHIBIT A-73**  
Sheet 2 of 3

SR520\_MASTER\_LOG\_E\_SMP\_21-20624.GPJ\_SHAN\_WIL\_GDT\_9/14/16 Log: JMW Rev: DPO Typ: CLP

Total Depth: 50.2 ft. Northing: 238,310 ft. Drilling Method: Advanced Casing Hole Diam.: 5 in.  
 Top Elevation: 49 ft. Easting: 1,277,628 ft. Drilling Company: WSDOT Rod Diam.: 1 3/4-inch  
 Vert. Datum: NAVD 88 Station: \_\_\_\_\_ Drill Rig Equipment: CME LC55 Track Rig 9C7-1 Hammer Type: Automatic  
 Horiz. Datum: NAD 83/91 Offset: \_\_\_\_\_ Other Comments: \_\_\_\_\_ Hammer ER: 81 %

**SOIL DESCRIPTION**  
 Refer to the report text for a proper understanding of the subsurface materials and drilling methods. The stratification lines represent the approximate boundaries between material types, and the transition may be gradual.

**PENETRATION RESISTANCE (blows/foot)**  
 ▲ Hammer Wt. & Drop: 140 lbs / 30 inches



**LEGEND**

\* Sample Not Recovered      ▼ Groundwater Level in Well      ◇ % Fines (<0.075mm)  
 E Environmental Sample Obtained      ● % Water Content  
 C Corrosion and/or Cation Exchange Sample Obtained      Plastic Limit —●— Liquid Limit  
 | Standard Penetration Test      Natural Water Content  
 ||| 2.5" O.D. Split Spoon Sample

**NOTES**

1. Refer to KEY for explanation of symbols, codes, abbreviations and definitions.
2. The stratification lines represent the approximate boundaries between soil types, and the transition may be gradual.
3. Discussions in the text of the geotechnical data reports are necessary for a proper understanding of the nature of the subsurface materials.
4. Groundwater level, if indicated above, is for the date specified and may vary. Groundwater level is the highest available measurement to date. Groundwater plots contain complete data sets.
5. USCS designation is based on visual-manual classification and selected lab testing.
6. Hammer ER = hammer energy ratio (efficiency) as a percentage.

SR 520 Bridge Replacement and HOV Program  
 West Approach to Montlake Lid  
 Seattle, Washington

**LOG OF BORING H-667p-15**

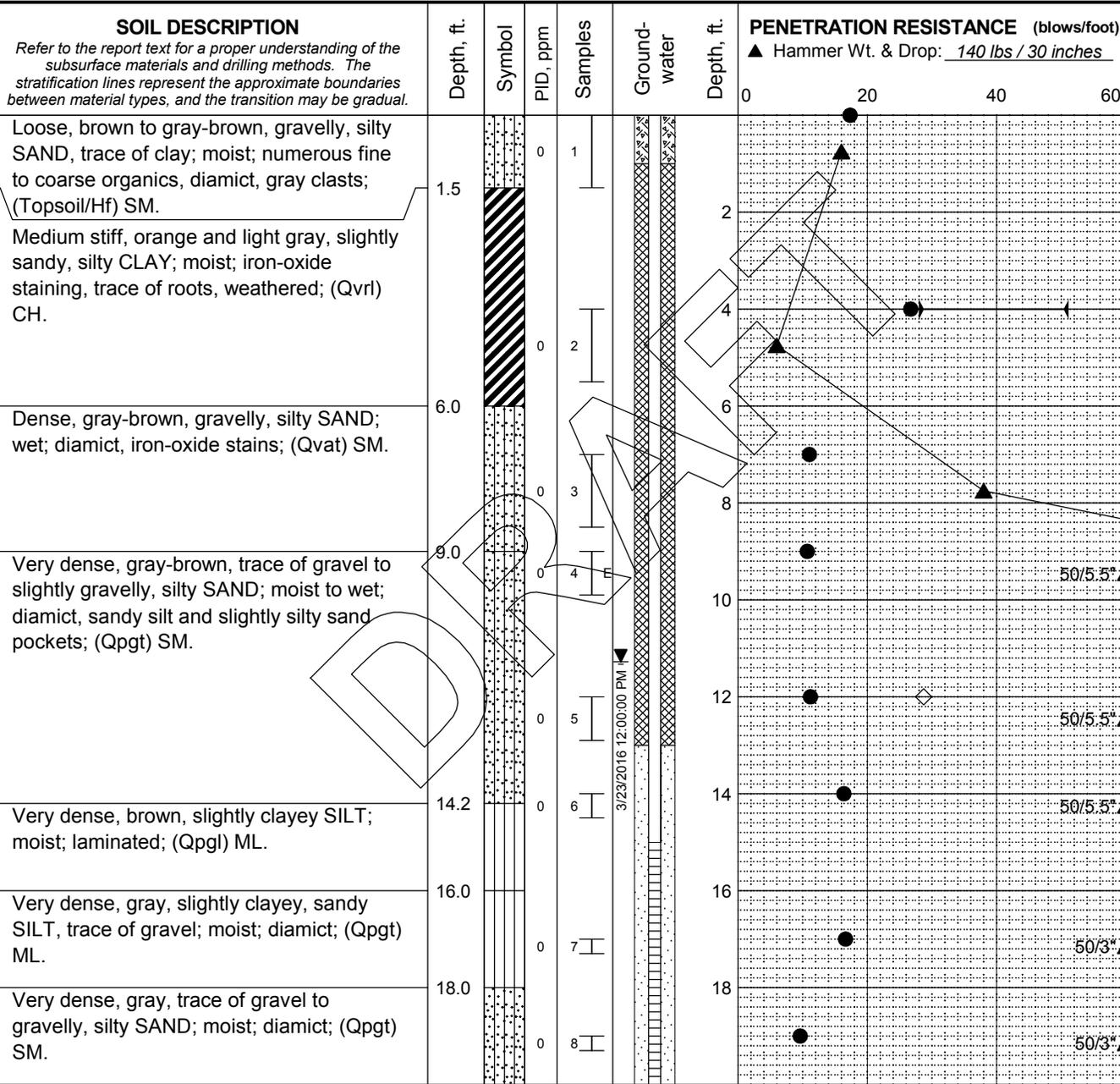
September 2016      21-1-22242-002

**SHANNON & WILSON, INC.**  
 Geotechnical and Environmental Consultants

**EXHIBIT A-73**  
 Sheet 3 of 3

SR520\_MASTER\_LOG\_E\_SMP\_21-20624.GPJ\_SHAN\_WIL\_GDT\_9/14/16 Log: JMW Rev: DPO Typ: CLP

Total Depth: 75.5 ft. Northing: 238,334 ft. Drilling Method: Advanced Casing Hole Diam.: 4 in.  
 Top Elevation: 59 ft. Easting: 1,277,977 ft. Drilling Company: WSDOT Rod Diam.: AWJ 1-3/4"  
 Vert. Datum: NAVD 88 Station: \_\_\_\_\_ Drill Rig Equipment: CME LC55 Track Rig 9C7-1 Hammer Type: Automatic  
 Horiz. Datum: NAD 83/91 Offset: \_\_\_\_\_ Other Comments: \_\_\_\_\_ Hammer ER: \_\_\_\_\_



CONTINUED NEXT SHEET

**LEGEND**

- \* Sample Not Recovered
- E Environmental Sample Obtained
- ┌ Standard Penetration Test
- ▼ Groundwater Level in Well
- ◇ % Fines (<0.075mm)
- % Water Content
- Plastic Limit —●— Liquid Limit
- Natural Water Content

**NOTES**

1. Refer to KEY for explanation of symbols, codes, abbreviations and definitions.
2. The stratification lines represent the approximate boundaries between soil types, and the transition may be gradual.
3. Discussions in the text of the geotechnical data reports are necessary for a proper understanding of the nature of the subsurface materials.
4. Groundwater level, if indicated above, is for the date specified and may vary. Groundwater level is the highest available measurement to date. Groundwater plots contain complete data sets.
5. USCS designation is based on visual-manual classification and selected lab testing.
6. Hammer ER = hammer energy ratio (efficiency) as a percentage.

SR 520 Bridge Replacement and HOV Program  
West Approach to Montlake Lid  
Seattle, Washington

**LOG OF BORING H-691p-16**

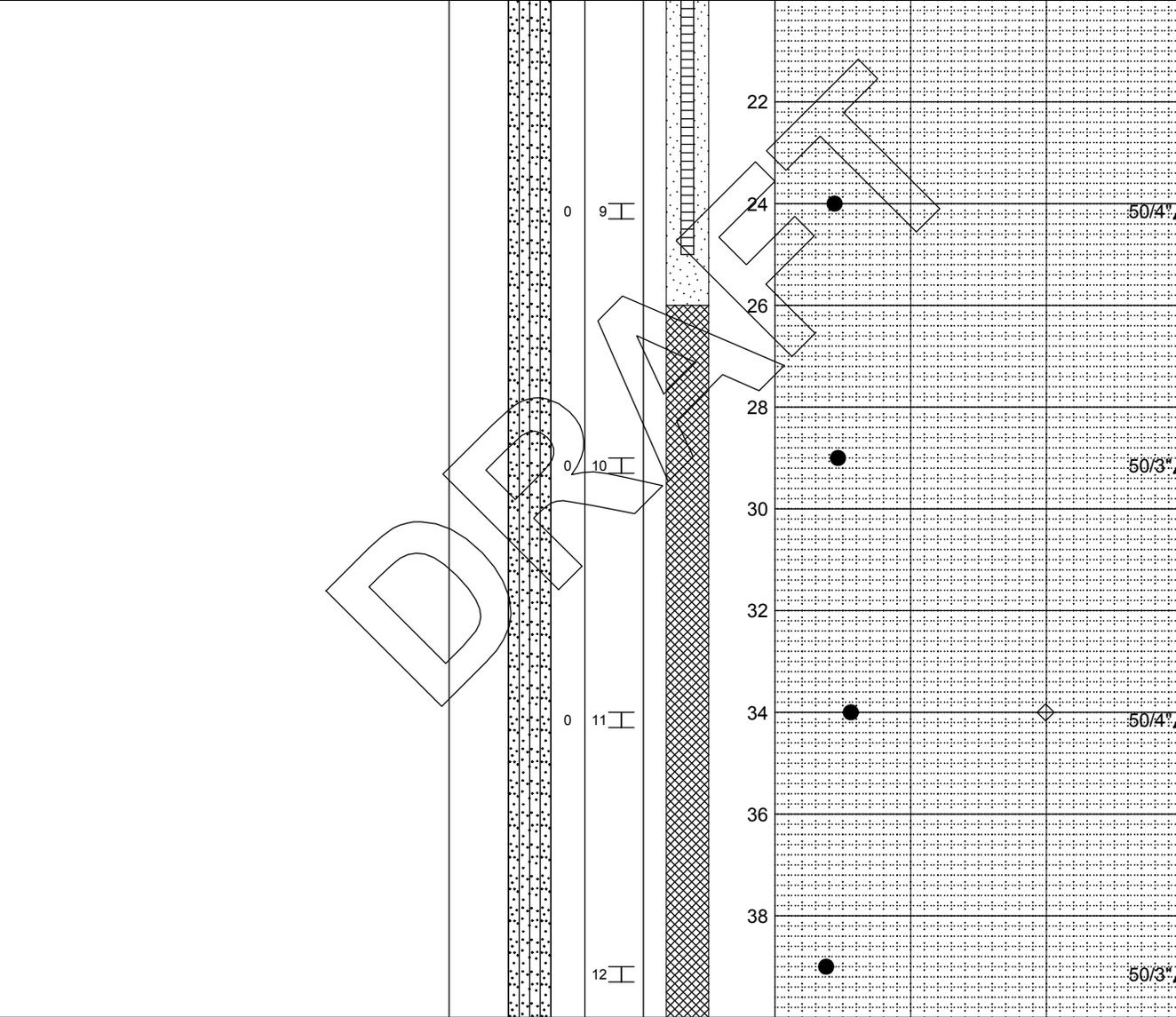
September 2016 21-1-22242-002

<b>SHANNON &amp; WILSON, INC.</b> Geotechnical and Environmental Consultants	<b>EXHIBIT</b> Sheet 1 of 4
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SR520\_MASTER\_LOG\_E\_SMP\_21-20624-GPJ\_SHAN\_WIL\_GDT\_9/14/16 Log: JMW Rev: JKP Typ: JKP

Total Depth: <u>75.5 ft.</u>	Northing: <u>238,334 ft.</u>	Drilling Method: <u>Advanced Casing</u>	Hole Diam.: <u>4 in.</u>
Top Elevation: <u>59 ft.</u>	Easting: <u>1,277,977 ft.</u>	Drilling Company: <u>WSDOT</u>	Rod Diam.: <u>AWJ 1-3/4"</u>
Vert. Datum: <u>NAVD 88</u>	Station: _____	Drill Rig Equipment: <u>CME LC55 Track Rig 9C7-1</u>	Hammer Type: <u>Automatic</u>
Horiz. Datum: <u>NAD 83/91</u>	Offset: _____	Other Comments: _____	Hammer ER: _____

<b>SOIL DESCRIPTION</b> <i>Refer to the report text for a proper understanding of the subsurface materials and drilling methods. The stratification lines represent the approximate boundaries between material types, and the transition may be gradual.</i>	Depth, ft.	Symbol	PID, ppm	Samples	Ground-water	Depth, ft.	<b>PENETRATION RESISTANCE (blows/foot)</b>
							▲ Hammer Wt. & Drop: <u>140 lbs / 30 inches</u>



CONTINUED NEXT SHEET

**LEGEND**

- \* Sample Not Recovered
- E Environmental Sample Obtained
- ┌ Standard Penetration Test
- ▼ Groundwater Level in Well
- ◇ % Fines (<0.075mm)
- % Water Content
- Plastic Limit —●— Liquid Limit
- Natural Water Content

**NOTES**

1. Refer to KEY for explanation of symbols, codes, abbreviations and definitions.
2. The stratification lines represent the approximate boundaries between soil types, and the transition may be gradual.
3. Discussions in the text of the geotechnical data reports are necessary for a proper understanding of the nature of the subsurface materials.
4. Groundwater level, if indicated above, is for the date specified and may vary. Groundwater level is the highest available measurement to date. Groundwater plots contain complete data sets.
5. USCS designation is based on visual-manual classification and selected lab testing.
6. Hammer ER = hammer energy ratio (efficiency) as a percentage.

SR 520 Bridge Replacement and HOV Program  
West Approach to Montlake Lid  
Seattle, Washington

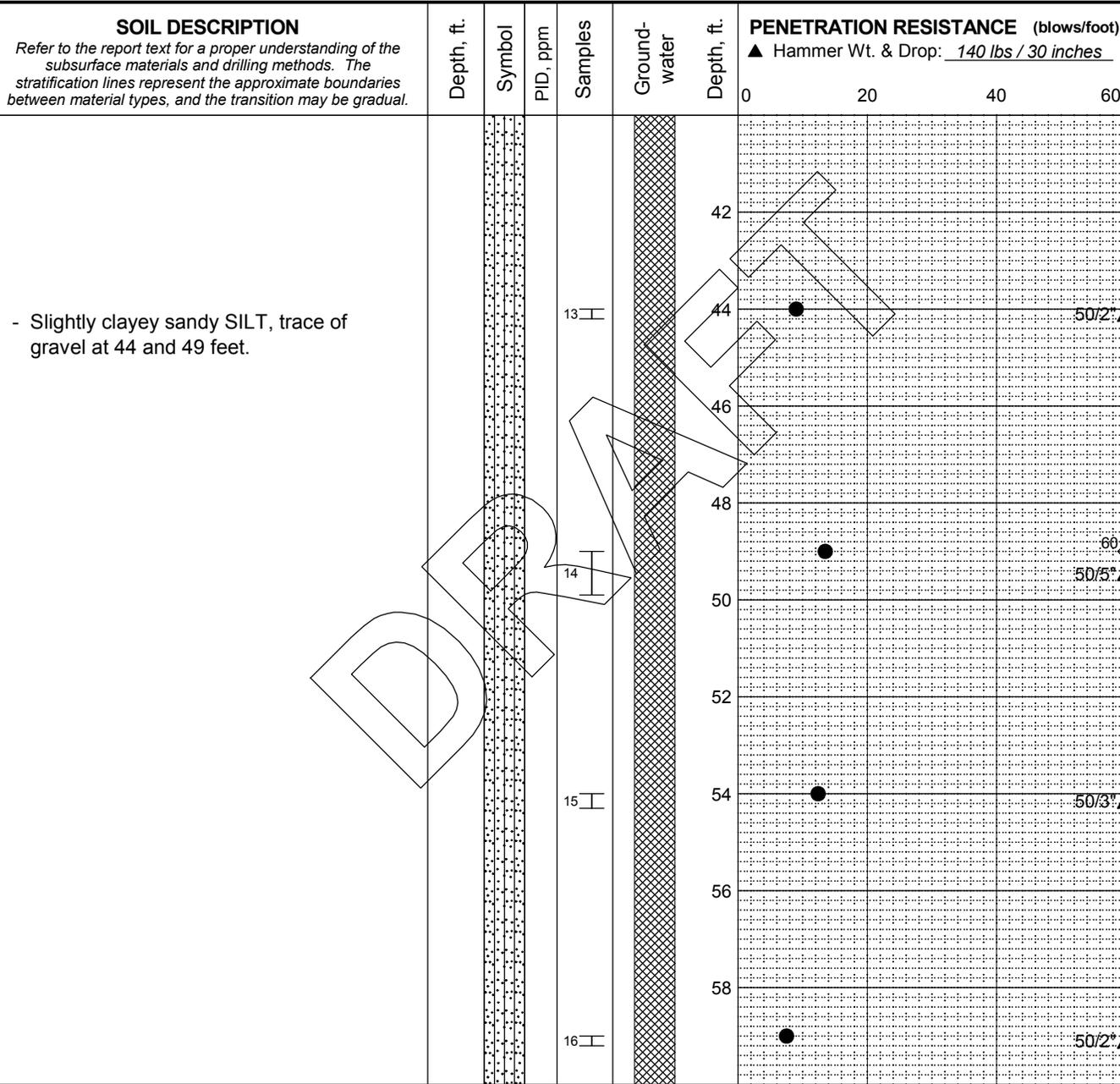
**LOG OF BORING H-691p-16**

September 2016 21-1-22242-002

<b>SHANNON &amp; WILSON, INC.</b> Geotechnical and Environmental Consultants	<b>EXHIBIT</b> Sheet 2 of 4
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SR520\_MASTER\_LOG\_E\_SMP\_21-20624-GPJ\_SHAN\_WIL\_GDT\_9/14/16 Log: JMW Rev: JKP Typ: JKP

Total Depth: 75.5 ft. Northing: 238,334 ft. Drilling Method: Advanced Casing Hole Diam.: 4 in.  
 Top Elevation: 59 ft. Easting: 1,277,977 ft. Drilling Company: WSDOT Rod Diam.: AWJ 1-3/4"  
 Vert. Datum: NAVD 88 Station: \_\_\_\_\_ Drill Rig Equipment: CME LC55 Track Rig 9C7-1 Hammer Type: Automatic  
 Horiz. Datum: NAD 83/91 Offset: \_\_\_\_\_ Other Comments: \_\_\_\_\_ Hammer ER: \_\_\_\_\_



CONTINUED NEXT SHEET

**LEGEND**

- \* Sample Not Recovered
- E Environmental Sample Obtained
- ┌ Standard Penetration Test
- ▼ Groundwater Level in Well
- ◇ % Fines (<0.075mm)
- % Water Content
- Plastic Limit —●— Liquid Limit
- Natural Water Content

**NOTES**

1. Refer to KEY for explanation of symbols, codes, abbreviations and definitions.
2. The stratification lines represent the approximate boundaries between soil types, and the transition may be gradual.
3. Discussions in the text of the geotechnical data reports are necessary for a proper understanding of the nature of the subsurface materials.
4. Groundwater level, if indicated above, is for the date specified and may vary. Groundwater level is the highest available measurement to date. Groundwater plots contain complete data sets.
5. USCS designation is based on visual-manual classification and selected lab testing.
6. Hammer ER = hammer energy ratio (efficiency) as a percentage.

SR 520 Bridge Replacement and HOV Program  
 West Approach to Montlake Lid  
 Seattle, Washington

**LOG OF BORING H-691p-16**

September 2016 21-1-22242-002

**SHANNON & WILSON, INC.**  
 Geotechnical and Environmental Consultants

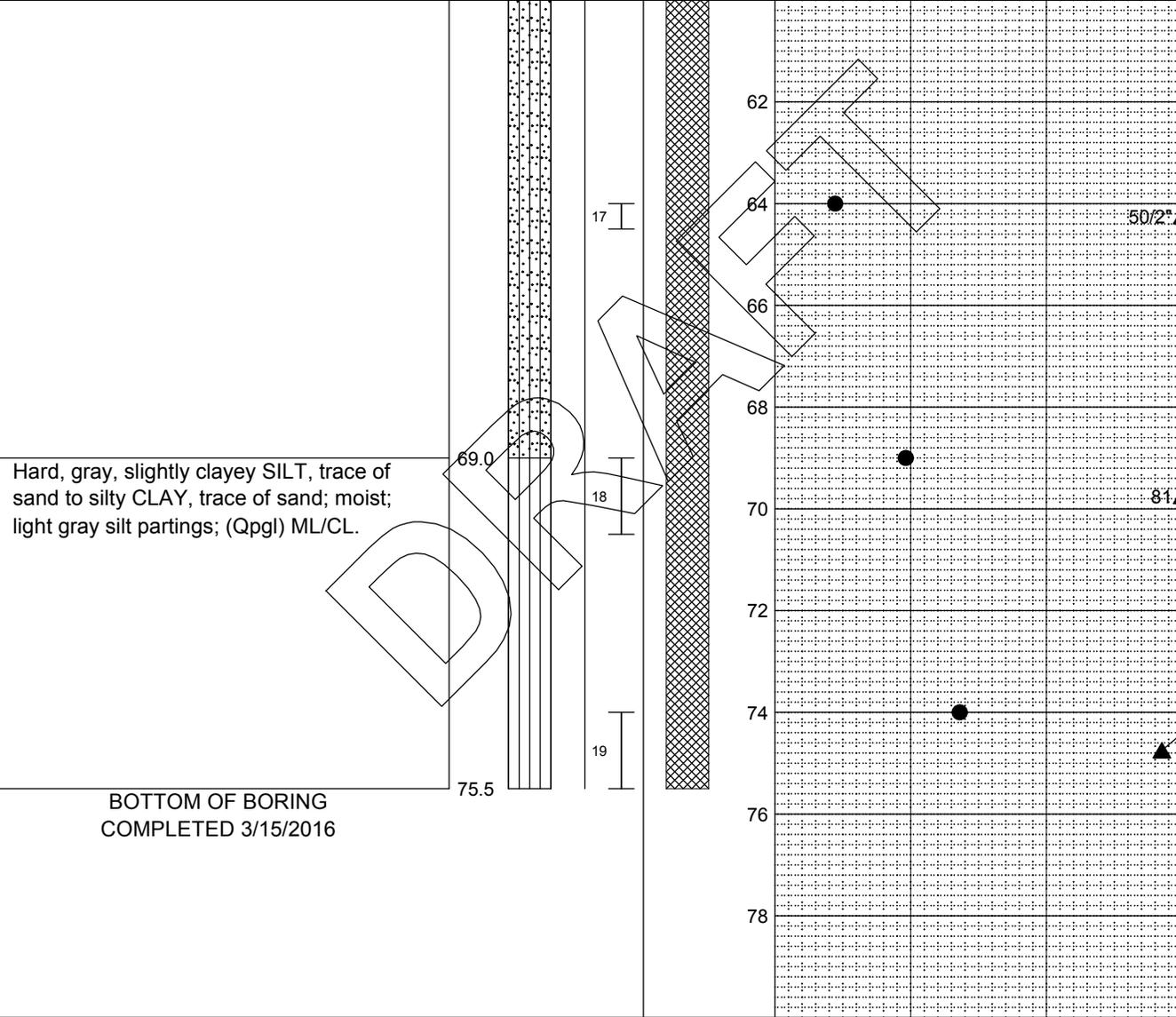
**EXHIBIT**  
 Sheet 3 of 4

SR520\_MASTER\_LOG\_E\_SMP\_21-20624-GPJ\_SHAN\_WIL\_GDT\_9/14/16 Log: JMW Rev: JKP Typ: JKP

Total Depth: 75.5 ft. Northing: 238,334 ft. Drilling Method: Advanced Casing Hole Diam.: 4 in.  
 Top Elevation: 59 ft. Easting: 1,277,977 ft. Drilling Company: WSDOT Rod Diam.: AWJ 1-3/4"  
 Vert. Datum: NAVD 88 Station: \_\_\_\_\_ Drill Rig Equipment: CME LC55 Track Rig 9C7-1 Hammer Type: Automatic  
 Horiz. Datum: NAD 83/91 Offset: \_\_\_\_\_ Other Comments: \_\_\_\_\_ Hammer ER: \_\_\_\_\_

**SOIL DESCRIPTION**  
 Refer to the report text for a proper understanding of the subsurface materials and drilling methods. The stratification lines represent the approximate boundaries between material types, and the transition may be gradual.

**PENETRATION RESISTANCE (blows/foot)**  
 ▲ Hammer Wt. & Drop: 140 lbs / 30 inches



**LEGEND**

\* Sample Not Recovered      ▼ Groundwater Level in Well  
 E Environmental Sample Obtained      ◇ % Fines (<0.075mm)  
 I Standard Penetration Test      ● % Water Content  
 Plastic Limit —●— Liquid Limit  
 Natural Water Content

**NOTES**

1. Refer to KEY for explanation of symbols, codes, abbreviations and definitions.
2. The stratification lines represent the approximate boundaries between soil types, and the transition may be gradual.
3. Discussions in the text of the geotechnical data reports are necessary for a proper understanding of the nature of the subsurface materials.
4. Groundwater level, if indicated above, is for the date specified and may vary. Groundwater level is the highest available measurement to date. Groundwater plots contain complete data sets.
5. USCS designation is based on visual-manual classification and selected lab testing.
6. Hammer ER = hammer energy ratio (efficiency) as a percentage.

SR 520 Bridge Replacement and HOV Program  
 West Approach to Montlake Lid  
 Seattle, Washington

**LOG OF BORING H-691p-16**

September 2016      21-1-22242-002

**SHANNON & WILSON, INC.**  
 Geotechnical and Environmental Consultants

**EXHIBIT**  
 Sheet 4 of 4

SR520\_MASTER\_LOG\_E\_SMP\_21-20624-GPJ\_SHAN\_WIL\_GDT\_9/14/16 Log: JMW Rev: JKP Typ: JKP



14648 NE 95<sup>th</sup> Street, Redmond, WA 98052 • (425) 883-3881

March 30, 2011

Cody Johnson  
Shannon & Wilson, Inc.  
400 N 34th Street, Suite 100  
Seattle, WA 98103

Re: Analytical Data for Project 21-1-20624-612  
Laboratory Reference No. 1103-188

Dear Cody:

Enclosed are the analytical results and associated quality control data for samples submitted on March 21, 2011.

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,

A handwritten signature in black ink, appearing to read "DB", with a long horizontal flourish extending to the right.

David Baumeister  
Project Manager

Enclosures

Date of Report: March 30, 2011  
Samples Submitted: March 21, 2011  
Laboratory Reference: 1103-188  
Project: 21-1-20624-612

### **Case Narrative**

Samples were collected on March 15, 17 and 19, 2011 and received by the laboratory on March 21, 2011. They were maintained at the laboratory at a temperature of 2°C to 6°C.

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.

Date of Report: March 30, 2011  
 Samples Submitted: March 21, 2011  
 Laboratory Reference: 1103-188  
 Project: 21-1-20624-612

### NWTPH-HCID

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-552-11:4</b>					
Laboratory ID:	03-188-03					
Gasoline Range Organics	<b>ND</b>	23	NWTPH-HCID	3-21-11	3-21-11	
Diesel Range Organics	<b>ND</b>	57	NWTPH-HCID	3-21-11	3-21-11	
Lube Oil Range Organics	<b>ND</b>	110	NWTPH-HCID	3-21-11	3-21-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>136</i>	<i>50-150</i>				

<b>Client ID:</b>	<b>H-609p-11:16.7</b>					
Laboratory ID:	03-188-04					
Gasoline Range Organics	<b>ND</b>	24	NWTPH-HCID	3-21-11	3-21-11	
Diesel Range Organics	<b>ND</b>	61	NWTPH-HCID	3-21-11	3-21-11	
Lube Oil Range Organics	<b>ND</b>	120	NWTPH-HCID	3-21-11	3-21-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>128</i>	<i>50-150</i>				

Date of Report: March 30, 2011  
 Samples Submitted: March 21, 2011  
 Laboratory Reference: 1103-188  
 Project: 21-1-20624-612

**NWTPH-HCID  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg (ppm)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>METHOD BLANK</b>						
Laboratory ID:	MB0321S1					
Gasoline Range Organics	<b>ND</b>	20	NWTPH-HCID	3-21-11	3-21-11	
Diesel Range Organics	<b>ND</b>	50	NWTPH-HCID	3-21-11	3-21-11	
Lube Oil Range Organics	<b>ND</b>	100	NWTPH-HCID	3-21-11	3-21-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>136</i>	<i>50-150</i>				

Date of Report: March 30, 2011  
 Samples Submitted: March 21, 2011  
 Laboratory Reference: 1103-188  
 Project: 21-1-20624-612

**NWTPH-Dx**  
 (with acid/silica gel clean-up)

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-536-11:8</b>					
Laboratory ID:	03-188-01					
Diesel Range Organics	<b>ND</b>	140	NWTPH-Dx	3-25-11	3-25-11	
Lube Oil Range Organics	<b>ND</b>	290	NWTPH-Dx	3-25-11	3-25-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>97</i>	<i>50-150</i>				
<b>Client ID:</b>	<b>H-561-11:4</b>					
Laboratory ID:	03-188-02					
Diesel Range Organics	<b>ND</b>	54	NWTPH-Dx	3-25-11	3-25-11	
Lube Oil	<b>140</b>	110	NWTPH-Dx	3-25-11	3-25-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>93</i>	<i>50-150</i>				

Date of Report: March 30, 2011  
 Samples Submitted: March 21, 2011  
 Laboratory Reference: 1103-188  
 Project: 21-1-20624-612

**NWTPH-Dx  
 QUALITY CONTROL  
 (with acid/silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB0325S1					
Diesel Range Organics	<b>ND</b>	25	NWTPH-Dx	3-25-11	3-25-11	
Lube Oil Range Organics	<b>ND</b>	50	NWTPH-Dx	3-25-11	3-25-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	105	50-150				

Analyte	Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>						
Laboratory ID:	03-224-01					
	ORIG	DUP				
Diesel Range Organics	<b>ND</b>	<b>ND</b>		NA	NA	
Lube Oil	<b>54.4</b>	<b>53.9</b>		1	NA	
<i>Surrogate:</i>						
<i>o-Terphenyl</i>			103 86	50-150		

Date of Report: March 30, 2011  
 Samples Submitted: March 21, 2011  
 Laboratory Reference: 1103-188  
 Project: 21-1-20624-612

**SEMIVOLATILES by EPA 8270D/SIM**  
 page 1 of 2

Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-536-11:8</b>					
Laboratory ID:	03-188-01					
n-Nitrosodimethylamine	ND	0.19	EPA 8270	3-28-11	3-28-11	
Pyridine	ND	1.9	EPA 8270	3-28-11	3-28-11	
Phenol	ND	0.19	EPA 8270	3-28-11	3-28-11	
Aniline	ND	0.19	EPA 8270	3-28-11	3-28-11	
bis(2-Chloroethyl)ether	ND	0.19	EPA 8270	3-28-11	3-28-11	
2-Chlorophenol	ND	0.19	EPA 8270	3-28-11	3-28-11	
1,3-Dichlorobenzene	ND	0.19	EPA 8270	3-28-11	3-28-11	
1,4-Dichlorobenzene	ND	0.19	EPA 8270	3-28-11	3-28-11	
Benzyl alcohol	ND	0.19	EPA 8270	3-28-11	3-28-11	
1,2-Dichlorobenzene	ND	0.19	EPA 8270	3-28-11	3-28-11	
2-Methylphenol (o-Cresol)	ND	0.19	EPA 8270	3-28-11	3-28-11	
bis(2-Chloroisopropyl)ether	ND	0.19	EPA 8270	3-28-11	3-28-11	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.19	EPA 8270	3-28-11	3-28-11	
n-Nitroso-di-n-propylamine	ND	0.19	EPA 8270	3-28-11	3-28-11	
Hexachloroethane	ND	0.19	EPA 8270	3-28-11	3-28-11	
Nitrobenzene	ND	0.19	EPA 8270	3-28-11	3-28-11	
Isophorone	ND	0.19	EPA 8270	3-28-11	3-28-11	
2-Nitrophenol	ND	0.19	EPA 8270	3-28-11	3-28-11	
2,4-Dimethylphenol	ND	4.7	EPA 8270	3-28-11	3-28-11	
bis(2-Chloroethoxy)methane	ND	0.19	EPA 8270	3-28-11	3-28-11	
2,4-Dichlorophenol	ND	0.19	EPA 8270	3-28-11	3-28-11	
1,2,4-Trichlorobenzene	ND	0.19	EPA 8270	3-28-11	3-28-11	
Naphthalene	ND	0.038	EPA 8270/SIM	3-28-11	3-29-11	
4-Chloroaniline	ND	0.19	EPA 8270	3-28-11	3-28-11	
Hexachlorobutadiene	ND	0.19	EPA 8270	3-28-11	3-28-11	
4-Chloro-3-methylphenol	ND	0.19	EPA 8270	3-28-11	3-28-11	
2-Methylnaphthalene	ND	0.038	EPA 8270/SIM	3-28-11	3-29-11	
1-Methylnaphthalene	ND	0.038	EPA 8270/SIM	3-28-11	3-29-11	
Hexachlorocyclopentadiene	ND	0.19	EPA 8270	3-28-11	3-28-11	
2,4,6-Trichlorophenol	ND	0.19	EPA 8270	3-28-11	3-28-11	
2,3-Dichloroaniline	ND	0.19	EPA 8270	3-28-11	3-28-11	
2,4,5-Trichlorophenol	ND	0.19	EPA 8270	3-28-11	3-28-11	
2-Chloronaphthalene	ND	0.19	EPA 8270	3-28-11	3-28-11	
2-Nitroaniline	ND	0.19	EPA 8270	3-28-11	3-28-11	
1,4-Dinitrobenzene	ND	0.19	EPA 8270	3-28-11	3-28-11	
Dimethylphthalate	ND	0.19	EPA 8270	3-28-11	3-28-11	
1,3-Dinitrobenzene	ND	0.95	EPA 8270	3-28-11	3-28-11	
2,6-Dinitrotoluene	ND	0.19	EPA 8270	3-28-11	3-28-11	
1,2-Dinitrobenzene	ND	0.19	EPA 8270	3-28-11	3-28-11	
Acenaphthylene	ND	0.038	EPA 8270/SIM	3-28-11	3-29-11	
3-Nitroaniline	ND	0.19	EPA 8270	3-28-11	3-28-11	

Date of Report: March 30, 2011  
 Samples Submitted: March 21, 2011  
 Laboratory Reference: 1103-188  
 Project: 21-1-20624-612

**SEMIVOLATILES by EPA 8270D/SIM**  
 page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-536-11:8</b>					
Laboratory ID:	03-188-01					
2,4-Dinitrophenol	ND	0.95	EPA 8270	3-28-11	3-28-11	
Acenaphthene	ND	0.038	EPA 8270/SIM	3-28-11	3-29-11	
4-Nitrophenol	ND	0.19	EPA 8270	3-28-11	3-28-11	
2,4-Dinitrotoluene	ND	0.19	EPA 8270	3-28-11	3-28-11	
Dibenzofuran	ND	0.19	EPA 8270	3-28-11	3-28-11	
2,3,5,6-Tetrachlorophenol	ND	0.19	EPA 8270	3-28-11	3-28-11	
2,3,4,6-Tetrachlorophenol	ND	0.19	EPA 8270	3-28-11	3-28-11	
Diethylphthalate	ND	0.95	EPA 8270	3-28-11	3-28-11	
4-Chlorophenyl-phenylether	ND	0.19	EPA 8270	3-28-11	3-28-11	
4-Nitroaniline	ND	0.19	EPA 8270	3-28-11	3-28-11	
Fluorene	ND	0.038	EPA 8270/SIM	3-28-11	3-29-11	
4,6-Dinitro-2-methylphenol	ND	0.95	EPA 8270	3-28-11	3-28-11	
n-Nitrosodiphenylamine	ND	0.19	EPA 8270	3-28-11	3-28-11	
1,2-Diphenylhydrazine	ND	0.19	EPA 8270	3-28-11	3-28-11	
4-Bromophenyl-phenylether	ND	0.19	EPA 8270	3-28-11	3-28-11	
Hexachlorobenzene	ND	0.19	EPA 8270	3-28-11	3-28-11	
Pentachlorophenol	ND	0.95	EPA 8270	3-28-11	3-28-11	
Phenanthrene	0.079	0.038	EPA 8270/SIM	3-28-11	3-29-11	
Anthracene	ND	0.038	EPA 8270/SIM	3-28-11	3-29-11	
Carbazole	ND	0.19	EPA 8270	3-28-11	3-28-11	
Di-n-butylphthalate	ND	0.19	EPA 8270	3-28-11	3-28-11	
Fluoranthene	0.10	0.038	EPA 8270/SIM	3-28-11	3-29-11	
Benzidine	ND	1.9	EPA 8270	3-28-11	3-28-11	
Pyrene	0.10	0.038	EPA 8270/SIM	3-28-11	3-29-11	
Butylbenzylphthalate	ND	0.19	EPA 8270	3-28-11	3-28-11	
bis(2-Ethylhexyl)adipate	ND	0.19	EPA 8270	3-28-11	3-28-11	
3,3'-Dichlorobenzidine	ND	1.9	EPA 8270	3-28-11	3-28-11	
Benz[a]anthracene	0.050	0.038	EPA 8270/SIM	3-28-11	3-29-11	
Chrysene	0.057	0.038	EPA 8270/SIM	3-28-11	3-29-11	
bis(2-Ethylhexyl)phthalate	0.44	0.19	EPA 8270	3-28-11	3-28-11	
Di-n-octylphthalate	ND	0.19	EPA 8270	3-28-11	3-28-11	
Benzo(b)fluoranthene	0.039	0.038	EPA 8270/SIM	3-28-11	3-29-11	
Benzo(k)fluoranthene	ND	0.038	EPA 8270/SIM	3-28-11	3-29-11	
Benzo(a)pyrene	0.050	0.038	EPA 8270/SIM	3-28-11	3-29-11	
Indeno[1,2,3-cd]pyrene	0.038	0.038	EPA 8270/SIM	3-28-11	3-29-11	
Dibenzo(a,h)anthracene	ND	0.038	EPA 8270/SIM	3-28-11	3-29-11	
Benzo(ghi)perylene	0.043	0.038	EPA 8270/SIM	3-28-11	3-29-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorophenol</i>	<i>37</i>	<i>30 - 97</i>				
<i>Phenol-d6</i>	<i>54</i>	<i>40 - 104</i>				
<i>D5-Nitrobenzene</i>	<i>45</i>	<i>35 - 102</i>				
<i>2-Fluorobiphenyl</i>	<i>57</i>	<i>44 - 97</i>				
<i>2,4,6-Tribromophenol</i>	<i>69</i>	<i>41 - 110</i>				
<i>D14-Terphenyl</i>	<i>67</i>	<i>53 - 107</i>				

Date of Report: March 30, 2011  
 Samples Submitted: March 21, 2011  
 Laboratory Reference: 1103-188  
 Project: 21-1-20624-612

**SEMIVOLATILES by EPA 8270D/SIM**  
 page 1 of 2

Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-561-11:4</b>					
Laboratory ID:	03-188-02					
n-Nitrosodimethylamine	ND	0.072	EPA 8270	3-28-11	3-28-11	
Pyridine	ND	0.72	EPA 8270	3-28-11	3-28-11	
Phenol	ND	0.072	EPA 8270	3-28-11	3-28-11	
Aniline	ND	0.072	EPA 8270	3-28-11	3-28-11	
bis(2-Chloroethyl)ether	ND	0.072	EPA 8270	3-28-11	3-28-11	
2-Chlorophenol	ND	0.072	EPA 8270	3-28-11	3-28-11	
1,3-Dichlorobenzene	ND	0.072	EPA 8270	3-28-11	3-28-11	
1,4-Dichlorobenzene	ND	0.072	EPA 8270	3-28-11	3-28-11	
Benzyl alcohol	ND	0.072	EPA 8270	3-28-11	3-28-11	
1,2-Dichlorobenzene	ND	0.072	EPA 8270	3-28-11	3-28-11	
2-Methylphenol (o-Cresol)	ND	0.072	EPA 8270	3-28-11	3-28-11	
bis(2-Chloroisopropyl)ether	ND	0.072	EPA 8270	3-28-11	3-28-11	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.072	EPA 8270	3-28-11	3-28-11	
n-Nitroso-di-n-propylamine	ND	0.072	EPA 8270	3-28-11	3-28-11	
Hexachloroethane	ND	0.072	EPA 8270	3-28-11	3-28-11	
Nitrobenzene	ND	0.072	EPA 8270	3-28-11	3-28-11	
Isophorone	ND	0.072	EPA 8270	3-28-11	3-28-11	
2-Nitrophenol	ND	0.072	EPA 8270	3-28-11	3-28-11	
2,4-Dimethylphenol	ND	1.8	EPA 8270	3-28-11	3-28-11	
bis(2-Chloroethoxy)methane	ND	0.072	EPA 8270	3-28-11	3-28-11	
2,4-Dichlorophenol	ND	0.072	EPA 8270	3-28-11	3-28-11	
1,2,4-Trichlorobenzene	ND	0.072	EPA 8270	3-28-11	3-28-11	
Naphthalene	ND	0.014	EPA 8270/SIM	3-28-11	3-29-11	
4-Chloroaniline	ND	0.072	EPA 8270	3-28-11	3-28-11	
Hexachlorobutadiene	ND	0.072	EPA 8270	3-28-11	3-28-11	
4-Chloro-3-methylphenol	ND	0.072	EPA 8270	3-28-11	3-28-11	
2-Methylnaphthalene	ND	0.014	EPA 8270/SIM	3-28-11	3-29-11	
1-Methylnaphthalene	ND	0.014	EPA 8270/SIM	3-28-11	3-29-11	
Hexachlorocyclopentadiene	ND	0.072	EPA 8270	3-28-11	3-28-11	
2,4,6-Trichlorophenol	ND	0.072	EPA 8270	3-28-11	3-28-11	
2,3-Dichloroaniline	ND	0.072	EPA 8270	3-28-11	3-28-11	
2,4,5-Trichlorophenol	ND	0.072	EPA 8270	3-28-11	3-28-11	
2-Chloronaphthalene	ND	0.072	EPA 8270	3-28-11	3-28-11	
2-Nitroaniline	ND	0.072	EPA 8270	3-28-11	3-28-11	
1,4-Dinitrobenzene	ND	0.072	EPA 8270	3-28-11	3-28-11	
Dimethylphthalate	ND	0.072	EPA 8270	3-28-11	3-28-11	
1,3-Dinitrobenzene	ND	0.36	EPA 8270	3-28-11	3-28-11	
2,6-Dinitrotoluene	ND	0.072	EPA 8270	3-28-11	3-28-11	
1,2-Dinitrobenzene	ND	0.072	EPA 8270	3-28-11	3-28-11	
Acenaphthylene	ND	0.014	EPA 8270/SIM	3-28-11	3-29-11	
3-Nitroaniline	ND	0.072	EPA 8270	3-28-11	3-28-11	

Date of Report: March 30, 2011  
 Samples Submitted: March 21, 2011  
 Laboratory Reference: 1103-188  
 Project: 21-1-20624-612

**SEMIVOLATILES by EPA 8270D/SIM**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-561-11:4</b>					
Laboratory ID:	03-188-02					
2,4-Dinitrophenol	ND	0.36	EPA 8270	3-28-11	3-28-11	
Acenaphthene	ND	0.014	EPA 8270/SIM	3-28-11	3-29-11	
4-Nitrophenol	ND	0.072	EPA 8270	3-28-11	3-28-11	
2,4-Dinitrotoluene	ND	0.072	EPA 8270	3-28-11	3-28-11	
Dibenzofuran	ND	0.072	EPA 8270	3-28-11	3-28-11	
2,3,5,6-Tetrachlorophenol	ND	0.072	EPA 8270	3-28-11	3-28-11	
2,3,4,6-Tetrachlorophenol	ND	0.072	EPA 8270	3-28-11	3-28-11	
Diethylphthalate	ND	0.36	EPA 8270	3-28-11	3-28-11	
4-Chlorophenyl-phenylether	ND	0.072	EPA 8270	3-28-11	3-28-11	
4-Nitroaniline	ND	0.072	EPA 8270	3-28-11	3-28-11	
Fluorene	ND	0.014	EPA 8270/SIM	3-28-11	3-29-11	
4,6-Dinitro-2-methylphenol	ND	0.36	EPA 8270	3-28-11	3-28-11	
n-Nitrosodiphenylamine	ND	0.072	EPA 8270	3-28-11	3-28-11	
1,2-Diphenylhydrazine	ND	0.072	EPA 8270	3-28-11	3-28-11	
4-Bromophenyl-phenylether	ND	0.072	EPA 8270	3-28-11	3-28-11	
Hexachlorobenzene	ND	0.072	EPA 8270	3-28-11	3-28-11	
Pentachlorophenol	ND	0.36	EPA 8270	3-28-11	3-28-11	
Phenanthrene	0.016	0.014	EPA 8270/SIM	3-28-11	3-29-11	
Anthracene	ND	0.014	EPA 8270/SIM	3-28-11	3-29-11	
Carbazole	ND	0.072	EPA 8270	3-28-11	3-28-11	
Di-n-butylphthalate	ND	0.072	EPA 8270	3-28-11	3-28-11	
Fluoranthene	0.023	0.014	EPA 8270/SIM	3-28-11	3-29-11	
Benzidine	ND	0.72	EPA 8270	3-28-11	3-28-11	
Pyrene	0.023	0.014	EPA 8270/SIM	3-28-11	3-29-11	
Butylbenzylphthalate	ND	0.072	EPA 8270	3-28-11	3-28-11	
bis-2-Ethylhexyladipate	ND	0.072	EPA 8270	3-28-11	3-28-11	
3,3'-Dichlorobenzidine	ND	0.72	EPA 8270	3-28-11	3-28-11	
Benz[a]anthracene	ND	0.014	EPA 8270/SIM	3-28-11	3-29-11	
Chrysene	0.020	0.014	EPA 8270/SIM	3-28-11	3-29-11	
bis(2-Ethylhexyl)phthalate	0.15	0.072	EPA 8270	3-28-11	3-28-11	
Di-n-octylphthalate	ND	0.072	EPA 8270	3-28-11	3-28-11	
Benzo(b)fluoranthene	ND	0.014	EPA 8270/SIM	3-28-11	3-29-11	
Benzo(k)fluoranthene	ND	0.014	EPA 8270/SIM	3-28-11	3-29-11	
Benzo(a)pyrene	ND	0.014	EPA 8270/SIM	3-28-11	3-29-11	
Indeno[1,2,3-cd]pyrene	ND	0.014	EPA 8270/SIM	3-28-11	3-29-11	
Dibenzo(a,h)anthracene	ND	0.014	EPA 8270/SIM	3-28-11	3-29-11	
Benzo(ghi)perylene	0.016	0.014	EPA 8270/SIM	3-28-11	3-29-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorophenol</i>	<i>40</i>	<i>30 - 97</i>				
<i>Phenol-d6</i>	<i>58</i>	<i>40 - 104</i>				
<i>D5-Nitrobenzene</i>	<i>54</i>	<i>35 - 102</i>				
<i>2-Fluorobiphenyl</i>	<i>64</i>	<i>44 - 97</i>				
<i>2,4,6-Tribromophenol</i>	<i>78</i>	<i>41 - 110</i>				
<i>D14-Terphenyl</i>	<i>72</i>	<i>53 - 107</i>				

Date of Report: March 30, 2011  
 Samples Submitted: March 21, 2011  
 Laboratory Reference: 1103-188  
 Project: 21-1-20624-612

**SEMIVOLATILES by EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
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Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0328S1					
n-Nitrosodimethylamine	ND	0.033	EPA 8270	3-28-11	3-28-11	
Pyridine	ND	0.33	EPA 8270	3-28-11	3-28-11	
Phenol	ND	0.033	EPA 8270	3-28-11	3-28-11	
Aniline	ND	0.033	EPA 8270	3-28-11	3-28-11	
bis(2-Chloroethyl)ether	ND	0.033	EPA 8270	3-28-11	3-28-11	
2-Chlorophenol	ND	0.033	EPA 8270	3-28-11	3-28-11	
1,3-Dichlorobenzene	ND	0.033	EPA 8270	3-28-11	3-28-11	
1,4-Dichlorobenzene	ND	0.033	EPA 8270	3-28-11	3-28-11	
Benzyl alcohol	ND	0.033	EPA 8270	3-28-11	3-28-11	
1,2-Dichlorobenzene	ND	0.033	EPA 8270	3-28-11	3-28-11	
2-Methylphenol (o-Cresol)	ND	0.033	EPA 8270	3-28-11	3-28-11	
bis(2-Chloroisopropyl)ether	ND	0.033	EPA 8270	3-28-11	3-28-11	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.033	EPA 8270	3-28-11	3-28-11	
n-Nitroso-di-n-propylamine	ND	0.033	EPA 8270	3-28-11	3-28-11	
Hexachloroethane	ND	0.033	EPA 8270	3-28-11	3-28-11	
Nitrobenzene	ND	0.033	EPA 8270	3-28-11	3-28-11	
Isophorone	ND	0.033	EPA 8270	3-28-11	3-28-11	
2-Nitrophenol	ND	0.033	EPA 8270	3-28-11	3-28-11	
2,4-Dimethylphenol	ND	0.83	EPA 8270	3-28-11	3-28-11	
bis(2-Chloroethoxy)methane	ND	0.033	EPA 8270	3-28-11	3-28-11	
2,4-Dichlorophenol	ND	0.033	EPA 8270	3-28-11	3-28-11	
1,2,4-Trichlorobenzene	ND	0.033	EPA 8270	3-28-11	3-28-11	
Naphthalene	ND	0.0067	EPA 8270/SIM	3-28-11	3-29-11	
4-Chloroaniline	ND	0.033	EPA 8270	3-28-11	3-28-11	
Hexachlorobutadiene	ND	0.033	EPA 8270	3-28-11	3-28-11	
4-Chloro-3-methylphenol	ND	0.033	EPA 8270	3-28-11	3-28-11	
2-Methylnaphthalene	ND	0.0067	EPA 8270/SIM	3-28-11	3-29-11	
1-Methylnaphthalene	ND	0.0067	EPA 8270/SIM	3-28-11	3-29-11	
Hexachlorocyclopentadiene	ND	0.033	EPA 8270	3-28-11	3-28-11	
2,4,6-Trichlorophenol	ND	0.033	EPA 8270	3-28-11	3-28-11	
2,3-Dichloroaniline	ND	0.033	EPA 8270	3-28-11	3-28-11	
2,4,5-Trichlorophenol	ND	0.033	EPA 8270	3-28-11	3-28-11	
2-Chloronaphthalene	ND	0.033	EPA 8270	3-28-11	3-28-11	
2-Nitroaniline	ND	0.033	EPA 8270	3-28-11	3-28-11	
1,4-Dinitrobenzene	ND	0.033	EPA 8270	3-28-11	3-28-11	
Dimethylphthalate	ND	0.033	EPA 8270	3-28-11	3-28-11	
1,3-Dinitrobenzene	ND	0.17	EPA 8270	3-28-11	3-28-11	
2,6-Dinitrotoluene	ND	0.033	EPA 8270	3-28-11	3-28-11	
1,2-Dinitrobenzene	ND	0.033	EPA 8270	3-28-11	3-28-11	
Acenaphthylene	ND	0.0067	EPA 8270/SIM	3-28-11	3-29-11	
3-Nitroaniline	ND	0.033	EPA 8270	3-28-11	3-28-11	

Date of Report: March 30, 2011  
 Samples Submitted: March 21, 2011  
 Laboratory Reference: 1103-188  
 Project: 21-1-20624-612

**SEMIVOLATILES by EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0328S1					
2,4-Dinitrophenol	ND	0.17	EPA 8270	3-28-11	3-28-11	
Acenaphthene	ND	0.0067	EPA 8270/SIM	3-28-11	3-29-11	
4-Nitrophenol	ND	0.033	EPA 8270	3-28-11	3-28-11	
2,4-Dinitrotoluene	ND	0.033	EPA 8270	3-28-11	3-28-11	
Dibenzofuran	ND	0.033	EPA 8270	3-28-11	3-28-11	
2,3,5,6-Tetrachlorophenol	ND	0.033	EPA 8270	3-28-11	3-28-11	
2,3,4,6-Tetrachlorophenol	ND	0.033	EPA 8270	3-28-11	3-28-11	
Diethylphthalate	ND	0.17	EPA 8270	3-28-11	3-28-11	
4-Chlorophenyl-phenylether	ND	0.033	EPA 8270	3-28-11	3-28-11	
4-Nitroaniline	ND	0.033	EPA 8270	3-28-11	3-28-11	
Fluorene	ND	0.0067	EPA 8270/SIM	3-28-11	3-29-11	
4,6-Dinitro-2-methylphenol	ND	0.17	EPA 8270	3-28-11	3-28-11	
n-Nitrosodiphenylamine	ND	0.033	EPA 8270	3-28-11	3-28-11	
1,2-Diphenylhydrazine	ND	0.033	EPA 8270	3-28-11	3-28-11	
4-Bromophenyl-phenylether	ND	0.033	EPA 8270	3-28-11	3-28-11	
Hexachlorobenzene	ND	0.033	EPA 8270	3-28-11	3-28-11	
Pentachlorophenol	ND	0.17	EPA 8270	3-28-11	3-28-11	
Phenanthrene	ND	0.0067	EPA 8270/SIM	3-28-11	3-29-11	
Anthracene	ND	0.0067	EPA 8270/SIM	3-28-11	3-29-11	
Carbazole	ND	0.033	EPA 8270	3-28-11	3-28-11	
Di-n-butylphthalate	ND	0.033	EPA 8270	3-28-11	3-28-11	
Fluoranthene	ND	0.0067	EPA 8270/SIM	3-28-11	3-29-11	
Benzidine	ND	0.33	EPA 8270	3-28-11	3-28-11	
Pyrene	ND	0.0067	EPA 8270/SIM	3-28-11	3-29-11	
Butylbenzylphthalate	ND	0.033	EPA 8270	3-28-11	3-28-11	
bis-2-Ethylhexyladipate	ND	0.033	EPA 8270	3-28-11	3-28-11	
3,3'-Dichlorobenzidine	ND	0.33	EPA 8270	3-28-11	3-28-11	
Benz[a]anthracene	ND	0.0067	EPA 8270/SIM	3-28-11	3-29-11	
Chrysene	ND	0.0067	EPA 8270/SIM	3-28-11	3-29-11	
bis(2-Ethylhexyl)phthalate	ND	0.033	EPA 8270	3-28-11	3-28-11	
Di-n-octylphthalate	ND	0.033	EPA 8270	3-28-11	3-28-11	
Benzo(b)fluoranthene	ND	0.0067	EPA 8270/SIM	3-28-11	3-29-11	
Benzo(k)fluoranthene	ND	0.0067	EPA 8270/SIM	3-28-11	3-29-11	
Benzo(a)pyrene	ND	0.0067	EPA 8270/SIM	3-28-11	3-29-11	
Indeno[1,2,3-cd]pyrene	ND	0.0067	EPA 8270/SIM	3-28-11	3-29-11	
Dibenzo(a,h)anthracene	ND	0.0067	EPA 8270/SIM	3-28-11	3-29-11	
Benzo(ghi)perylene	ND	0.0067	EPA 8270/SIM	3-28-11	3-29-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorophenol</i>	<i>49</i>	<i>30 - 97</i>				
<i>Phenol-d6</i>	<i>53</i>	<i>40 - 104</i>				
<i>D5-Nitrobenzene</i>	<i>53</i>	<i>35 - 102</i>				
<i>2-Fluorobiphenyl</i>	<i>58</i>	<i>44 - 97</i>				
<i>2,4,6-Tribromophenol</i>	<i>80</i>	<i>41 - 110</i>				
<i>D14-Terphenyl</i>	<i>79</i>	<i>53 - 107</i>				

Date of Report: March 30, 2011  
 Samples Submitted: March 21, 2011  
 Laboratory Reference: 1103-188  
 Project: 21-1-20624-612

**SEMIVOLATILES by EPA 8270D/SIM  
 MS/MSD QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result		Spike Level		Source	Percent		Recovery	RPD	RPD	Flags
	MS	MSD	MS	MSD	Result	Recovery	Limits	RPD	Limit		
<b>MATRIX SPIKES</b>											
Laboratory ID:	03-191-06										
	MS	MSD	MS	MSD		MS	MSD				
Phenol	<b>0.985</b>	<b>1.07</b>	1.33	1.33	ND	74	80	41 - 106	8	29	
2-Chlorophenol	<b>0.987</b>	<b>1.08</b>	1.33	1.33	ND	74	81	43 - 104	9	36	
1,4-Dichlorobenzene	<b>0.410</b>	<b>0.445</b>	0.667	0.667	ND	61	67	25 - 94	8	40	
n-Nitroso-di-n-propylamine	<b>0.469</b>	<b>0.506</b>	0.667	0.667	ND	70	76	40 - 100	8	34	
1,2,4-Trichlorobenzene	<b>0.431</b>	<b>0.479</b>	0.667	0.667	ND	65	72	39 - 86	11	34	
4-Chloro-3-methylphenol	<b>1.17</b>	<b>1.19</b>	1.33	1.33	ND	88	89	60 - 102	2	25	
Acenaphthene	<b>0.522</b>	<b>0.535</b>	0.667	0.667	ND	78	80	54 - 94	2	23	
4-Nitrophenol	<b>1.18</b>	<b>1.19</b>	1.33	1.33	ND	89	89	30 - 133	1	25	
2,4-Dinitrotoluene	<b>0.630</b>	<b>0.645</b>	0.667	0.667	ND	94	97	46 - 107	2	26	
Pentachlorophenol	<b>1.24</b>	<b>1.28</b>	1.33	1.33	ND	93	96	54 - 111	3	29	
Pyrene	<b>0.704</b>	<b>0.719</b>	0.667	0.667	0.193	77	79	54 - 108	2	21	
<i>Surrogate:</i>											
2-Fluorophenol						60	67	30 - 97			
Phenol-d6						69	74	40 - 104			
D5-Nitrobenzene						67	76	35 - 102			
2-Fluorobiphenyl						71	75	44 - 97			
2,4,6-Tribromophenol						85	90	41 - 110			
D14-Terphenyl						80	83	53 - 107			

Date of Report: March 30, 2011  
 Samples Submitted: March 21, 2011  
 Laboratory Reference: 1103-188  
 Project: 21-1-20624-612

### PCBs by EPA 8082

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-536-11:8</b>					
Laboratory ID:	03-188-01					
PCB-aroclor 1016	<b>ND</b>	0.28	EPA 8082	3-23-11	3-23-11	
PCB-aroclor 1221	<b>ND</b>	0.28	EPA 8082	3-23-11	3-23-11	
PCB-aroclor 1232	<b>ND</b>	0.28	EPA 8082	3-23-11	3-23-11	
PCB-aroclor 1242	<b>ND</b>	0.28	EPA 8082	3-23-11	3-23-11	
PCB-aroclor 1248	<b>ND</b>	0.28	EPA 8082	3-23-11	3-23-11	
PCB-aroclor 1254	<b>ND</b>	0.28	EPA 8082	3-23-11	3-23-11	
PCB-aroclor 1260	<b>ND</b>	0.28	EPA 8082	3-23-11	3-23-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>DCB</i>	<i>76</i>	<i>42-123</i>				
<b>Client ID:</b>	<b>H-561-11:4</b>					
Laboratory ID:	03-188-02					
PCB-aroclor 1016	<b>ND</b>	0.11	EPA 8082	3-23-11	3-23-11	
PCB-aroclor 1221	<b>ND</b>	0.11	EPA 8082	3-23-11	3-23-11	
PCB-aroclor 1232	<b>ND</b>	0.11	EPA 8082	3-23-11	3-23-11	
PCB-aroclor 1242	<b>ND</b>	0.11	EPA 8082	3-23-11	3-23-11	
PCB-aroclor 1248	<b>ND</b>	0.11	EPA 8082	3-23-11	3-23-11	
PCB-aroclor 1254	<b>ND</b>	0.11	EPA 8082	3-23-11	3-23-11	
PCB-aroclor 1260	<b>ND</b>	0.11	EPA 8082	3-23-11	3-23-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>DCB</i>	<i>75</i>	<i>42-123</i>				

Date of Report: March 30, 2011  
 Samples Submitted: March 21, 2011  
 Laboratory Reference: 1103-188  
 Project: 21-1-20624-612

**PCBs by EPA 8082  
 QUALITY CONTROL**

Matrix: Solid  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB0323S1					
PCB-aroclor 1016	ND	0.050	EPA 8082	3-23-11	3-23-11	
PCB-aroclor 1221	ND	0.050	EPA 8082	3-23-11	3-23-11	
PCB-aroclor 1232	ND	0.050	EPA 8082	3-23-11	3-23-11	
PCB-aroclor 1242	ND	0.050	EPA 8082	3-23-11	3-23-11	
PCB-aroclor 1248	ND	0.050	EPA 8082	3-23-11	3-23-11	
PCB-aroclor 1254	ND	0.050	EPA 8082	3-23-11	3-23-11	
PCB-aroclor 1260	ND	0.050	EPA 8082	3-23-11	3-23-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
DCB	88		42-123			

Analyte	Result		Spike Level		Source Result	Percent Recovery		Recovery Limits	RPD	RPD Limit	Flags
<b>MATRIX SPIKES</b>											
Laboratory ID:	03-191-06										
	MS	MSD	MS	MSD		MS	MSD				
PCB-aroclor 1260	0.442	0.435	0.500	0.500	ND	88	87	44-125	2	15	
<i>Surrogate:</i>											
DCB						87	85	42-123			

Date of Report: March 30, 2011  
 Samples Submitted: March 21, 2011  
 Laboratory Reference: 1103-188  
 Project: 21-1-20624-612

**TOTAL METALS**  
**EPA 6010B/6020/7471A**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	EPA Method	Date	Date	Flags
				Prepared	Analyzed	
Lab ID:	03-188-01					
<b>Client ID:</b>	<b>H-536-11:8</b>					
Arsenic	<b>ND</b>	14	6020	3-22-11	3-28-11	
Barium	<b>130</b>	14	6010B	3-22-11	3-22-11	
Cadmium	<b>ND</b>	1.4	6020	3-22-11	3-28-11	
Chromium	<b>66</b>	2.8	6010B	3-22-11	3-22-11	
Lead	<b>37</b>	28	6010B	3-22-11	3-22-11	
Mercury	<b>ND</b>	1.4	7471A	3-23-11	3-23-11	
Selenium	<b>ND</b>	14	6020	3-22-11	3-28-11	
Silver	<b>ND</b>	2.8	6010B	3-22-11	3-22-11	

Lab ID:	03-188-02					
<b>Client ID:</b>	<b>H-561-11:4</b>					
Arsenic	<b>ND</b>	11	6010B	3-22-11	3-22-11	
Barium	<b>48</b>	5.4	6010B	3-22-11	3-22-11	
Cadmium	<b>ND</b>	1.1	6010B	3-22-11	3-22-11	
Chromium	<b>32</b>	1.1	6010B	3-22-11	3-22-11	
Lead	<b>ND</b>	11	6010B	3-22-11	3-22-11	
Mercury	<b>ND</b>	0.54	7471A	3-23-11	3-23-11	
Selenium	<b>ND</b>	22	6010B	3-22-11	3-22-11	
Silver	<b>ND</b>	1.1	6010B	3-22-11	3-22-11	

Date of Report: March 30, 2011  
 Samples Submitted: March 21, 2011  
 Laboratory Reference: 1103-188  
 Project: 21-1-20624-612

**TOTAL METALS  
 EPA 6010B/7471A**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	EPA Method	Date	Date	Flags
				Prepared	Analyzed	
Lab ID:	03-188-03					
Client ID:	H-552-11:4					
Arsenic	ND	5.7	6010B	3-22-11	3-22-11	
Barium	47	2.8	6010B	3-22-11	3-22-11	
Cadmium	ND	0.57	6010B	3-22-11	3-22-11	
Chromium	47	0.57	6010B	3-22-11	3-22-11	
Lead	ND	5.7	6010B	3-22-11	3-22-11	
Mercury	ND	0.28	7471A	3-23-11	3-23-11	
Selenium	ND	11	6010B	3-22-11	3-22-11	
Silver	ND	0.57	6010B	3-22-11	3-22-11	

Lab ID: 03-188-04  
 Client ID: H-609p-11:16.7

Arsenic	ND	6.1	6010B	3-22-11	3-22-11	
Barium	86	3.0	6010B	3-22-11	3-22-11	
Cadmium	ND	0.61	6010B	3-22-11	3-22-11	
Chromium	48	0.61	6010B	3-22-11	3-22-11	
Lead	28	6.1	6010B	3-22-11	3-22-11	
Mercury	ND	0.30	7471A	3-23-11	3-23-11	
Selenium	ND	12	6010B	3-22-11	3-22-11	
Silver	ND	0.61	6010B	3-22-11	3-22-11	

Date of Report: March 30, 2011  
 Samples Submitted: March 21, 2011  
 Laboratory Reference: 1103-188  
 Project: 21-1-20624-612

**TOTAL METALS  
 EPA 6010B/7471A  
 METHOD BLANK QUALITY CONTROL**

Date Extracted: 3-22&23-11  
 Date Analyzed: 3-22&23-11  
  
 Matrix: Soil  
 Units: mg/kg (ppm)  
  
 Lab ID: MB0322S2,MB0322S3&MB0323S1

Analyte	Method	Result	PQL
Arsenic	6010B	<b>ND</b>	5.0
Barium	6010B	<b>ND</b>	2.5
Cadmium	6010B	<b>ND</b>	0.50
Chromium	6010B	<b>ND</b>	0.50
Lead	6010B	<b>ND</b>	5.0
Mercury	7471A	<b>ND</b>	0.25
Selenium	6010B	<b>ND</b>	10
Silver	6010B	<b>ND</b>	0.50

Date of Report: March 30, 2011  
Samples Submitted: March 21, 2011  
Laboratory Reference: 1103-188  
Project: 21-1-20624-612

**TOTAL METALS  
EPA 6010B  
METHOD BLANK QUALITY CONTROL**

Date Extracted: 3-22-11  
Date Analyzed: 3-28-11  
  
Matrix: Soil  
Units: mg/kg (ppm)  
  
Lab ID: MB0322S2

Analyte	Method	Result	PQL
Arsenic	6010B	<b>ND</b>	2.5
Cadmium	6010B	<b>ND</b>	0.25
Selenium	6010B	<b>ND</b>	2.5

Date of Report: March 30, 2011  
 Samples Submitted: March 21, 2011  
 Laboratory Reference: 1103-188  
 Project: 21-1-20624-612

**TOTAL METALS  
 EPA 6010B/7471A  
 DUPLICATE QUALITY CONTROL**

Date Extracted: 3-22&23-11

Date Analyzed: 3-22&23-11

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-162-05

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Arsenic	<b>ND</b>	<b>ND</b>	NA	5.0	
Barium	<b>19.0</b>	<b>15.5</b>	20	2.5	
Cadmium	<b>ND</b>	<b>ND</b>	NA	0.50	
Chromium	<b>8.55</b>	<b>7.55</b>	12	0.50	
Lead	<b>ND</b>	<b>ND</b>	NA	5.0	
Mercury	<b>ND</b>	<b>ND</b>	NA	0.25	
Selenium	<b>ND</b>	<b>ND</b>	NA	10	
Silver	<b>ND</b>	<b>ND</b>	NA	0.50	

Date of Report: March 30, 2011  
Samples Submitted: March 21, 2011  
Laboratory Reference: 1103-188  
Project: 21-1-20624-612

**TOTAL METALS  
EPA 6010B  
DUPLICATE QUALITY CONTROL**

Date Extracted: 3-22-11  
Date Analyzed: 3-28-11  
  
Matrix: Soil  
Units: mg/kg (ppm)  
  
Lab ID: 03-162-05

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Arsenic	<b>2.50</b>	<b>ND</b>	NA	2.5	
Cadmium	<b>ND</b>	<b>ND</b>	NA	0.25	
Selenium	<b>ND</b>	<b>ND</b>	NA	2.5	

Date of Report: March 30, 2011  
 Samples Submitted: March 21, 2011  
 Laboratory Reference: 1103-188  
 Project: 21-1-20624-612

**TOTAL METALS  
 EPA 6010B/7471A  
 MS/MSD QUALITY CONTROL**

Date Extracted: 3-22&23-11

Date Analyzed: 3-22&23-11

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-162-05

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Arsenic	100	<b>93.5</b>	93	<b>96.1</b>	96	3	
Barium	100	<b>121</b>	102	<b>119</b>	100	2	
Cadmium	50	<b>46.4</b>	93	<b>47.4</b>	95	2	
Chromium	100	<b>101</b>	93	<b>104</b>	95	2	
Lead	250	<b>231</b>	92	<b>237</b>	95	2	
Mercury	0.50	<b>0.528</b>	106	<b>0.513</b>	103	3	
Selenium	100	<b>93.5</b>	93	<b>96.9</b>	97	4	
Silver	25	<b>23.4</b>	94	<b>23.6</b>	94	1	

Date of Report: March 30, 2011  
 Samples Submitted: March 21, 2011  
 Laboratory Reference: 1103-188  
 Project: 21-1-20624-612

**TOTAL METALS  
 EPA 6010B  
 MS/MSD QUALITY CONTROL**

Date Extracted: 3-22-11

Date Analyzed: 3-28-11

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-162-05

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Arsenic	100	<b>100</b>	98	<b>101</b>	99	1	
Cadmium	50	<b>48.3</b>	97	<b>51.0</b>	102	6	
Selenium	100	<b>103</b>	103	<b>103</b>	103	0	

Date of Report: March 30, 2011  
Samples Submitted: March 21, 2011  
Laboratory Reference: 1103-188  
Project: 21-1-20624-612

**% MOISTURE**

Date Analyzed: 3-21&23-11

Client ID	Lab ID	% Moisture
H-536-11:8	03-188-01	82
H-561-11:4	03-188-02	54
H-552-11:4	03-188-03	12
H-609p-11:16.7	03-188-04	17



### 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.
- Y - Sample extract treated with an acid/silica gel cleanup procedure.
- Z -
- ND - Not Detected at PQL  
 PQL - Practical Quantitation Limit  
 RPD - Relative Percent Difference





14648 NE 95<sup>th</sup> Street, Redmond, WA 98052 • (425) 883-3881

February 6, 2015

Edwin Ptak  
Shannon & Wilson, Inc.  
400 N 34th Street, Suite 100  
Seattle, WA 98103

Re: Analytical Data for Project 21-1-20624  
Laboratory Reference No. 1501-213

Dear Edwin:

Enclosed are the analytical results and associated quality control data for samples submitted on January 29, 2015.

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,

A handwritten signature in black ink, appearing to read "DB", with a long horizontal stroke extending to the right.

David Baumeister  
Project Manager

Enclosures

Date of Report: February 6, 2015  
Samples Submitted: January 29, 2015  
Laboratory Reference: 1501-213  
Project: 21-1-20624

### **Case Narrative**

Samples were collected on January 28, 2015 and received by the laboratory on January 29, 2015. They were maintained at the laboratory at a temperature of 2°C to 6°C.

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.

Date of Report: February 6, 2015  
 Samples Submitted: January 29, 2015  
 Laboratory Reference: 1501-213  
 Project: 21-1-20624

### NWTPH-HCID

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-662p-15:7.5</b>					
Laboratory ID:	01-213-01					
Gasoline Range Organics	<b>ND</b>	27	NWTPH-HCID	1-29-15	1-29-15	
Diesel Range Organics	<b>ND</b>	67	NWTPH-HCID	1-29-15	1-29-15	
Lube Oil Range Organics	<b>ND</b>	140	NWTPH-HCID	1-29-15	1-29-15	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	94	50-150				

<b>Client ID:</b>	<b>H-667p-15:7.5</b>					
Laboratory ID:	01-213-02					
Gasoline Range Organics	<b>ND</b>	23	NWTPH-HCID	1-29-15	1-29-15	
Diesel Range Organics	<b>ND</b>	57	NWTPH-HCID	1-29-15	1-29-15	
Lube Oil Range Organics	<b>ND</b>	110	NWTPH-HCID	1-29-15	1-29-15	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	99	50-150				

Date of Report: February 6, 2015  
 Samples Submitted: January 29, 2015  
 Laboratory Reference: 1501-213  
 Project: 21-1-20624

**NWTPH-HCID  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg (ppm)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>METHOD BLANK</b>						
Laboratory ID:	MB0129S1					
Gasoline Range Organics	<b>ND</b>	20	NWTPH-HCID	1-29-15	1-29-15	
Diesel Range Organics	<b>ND</b>	50	NWTPH-HCID	1-29-15	1-29-15	
Lube Oil Range Organics	<b>ND</b>	100	NWTPH-HCID	1-29-15	1-29-15	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>104</i>	<i>50-150</i>				

Date of Report: February 6, 2015  
 Samples Submitted: January 29, 2015  
 Laboratory Reference: 1501-213  
 Project: 21-1-20624

**TOTAL METALS  
 EPA 6010C/7471B**

Matrix: Soil  
 Units: mg/kg (ppm)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>EPA Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
Lab ID:	01-213-01					
<b>Client ID:</b>	<b>H-662p-15:7.5</b>					
Arsenic	ND	13	6010C	2-3-15	2-3-15	
Barium	160	3.4	6010C	2-3-15	2-3-15	
Cadmium	ND	0.67	6010C	2-3-15	2-3-15	
Chromium	79	0.67	6010C	2-3-15	2-3-15	
Lead	ND	6.7	6010C	2-3-15	2-3-15	
Mercury	ND	0.34	7471B	2-4-15	2-4-15	
Selenium	ND	13	6010C	2-3-15	2-3-15	
Silver	ND	1.3	6010C	2-3-15	2-3-15	

Lab ID:	01-213-02					
<b>Client ID:</b>	<b>H-667p-15:7.5</b>					
Arsenic	ND	11	6010C	2-3-15	2-3-15	
Barium	49	2.9	6010C	2-3-15	2-3-15	
Cadmium	ND	0.57	6010C	2-3-15	2-3-15	
Chromium	44	0.57	6010C	2-3-15	2-3-15	
Lead	20	5.7	6010C	2-3-15	2-3-15	
Mercury	ND	0.29	7471B	2-4-15	2-4-15	
Selenium	ND	11	6010C	2-3-15	2-3-15	
Silver	ND	1.1	6010C	2-3-15	2-3-15	

Date of Report: February 6, 2015  
 Samples Submitted: January 29, 2015  
 Laboratory Reference: 1501-213  
 Project: 21-1-20624

**TOTAL METALS  
 EPA 6010C/7471B  
 METHOD BLANK QUALITY CONTROL**

Date Extracted: 2-3&4-15  
 Date Analyzed: 2-3&4-15  
  
 Matrix: Soil  
 Units: mg/kg (ppm)  
  
 Lab ID: MB0203SM1&MB0204S1

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
Mercury	7471B	ND	0.25
Selenium	6010C	ND	10
Silver	6010C	ND	1.0

Date of Report: February 6, 2015  
 Samples Submitted: January 29, 2015  
 Laboratory Reference: 1501-213  
 Project: 21-1-20624

**TOTAL METALS  
 EPA 6010C/7471B  
 DUPLICATE QUALITY CONTROL**

Date Extracted: 2-3&4-15

Date Analyzed: 2-3&4-15

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 01-203-16

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Arsenic	ND	ND	NA	10	
Barium	33.2	29.0	14	2.5	
Cadmium	ND	ND	NA	0.50	
Chromium	23.8	18.5	25	0.50	
Lead	ND	ND	NA	5.0	
Mercury	ND	ND	NA	0.25	
Selenium	ND	ND	NA	10	
Silver	ND	ND	NA	1.0	

Date of Report: February 6, 2015  
 Samples Submitted: January 29, 2015  
 Laboratory Reference: 1501-213  
 Project: 21-1-20624

**TOTAL METALS  
 EPA 6010C/7471B  
 MS/MSD QUALITY CONTROL**

Date Extracted: 2-3&4-15

Date Analyzed: 2-3&4-15

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 01-203-16

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Arsenic	100	<b>93.1</b>	93	<b>94.4</b>	94	1	
Barium	100	<b>126</b>	93	<b>122</b>	89	4	
Cadmium	50.0	<b>46.0</b>	92	<b>45.8</b>	92	0	
Chromium	100	<b>112</b>	88	<b>109</b>	86	2	
Lead	250	<b>236</b>	94	<b>235</b>	94	0	
Mercury	0.500	<b>0.539</b>	108	<b>0.524</b>	105	3	
Selenium	100	<b>94.9</b>	95	<b>94.3</b>	94	1	
Silver	25.0	<b>21.9</b>	88	<b>21.9</b>	88	0	

Date of Report: February 6, 2015  
Samples Submitted: January 29, 2015  
Laboratory Reference: 1501-213  
Project: 21-1-20624

### % MOISTURE

Date Analyzed: 1-29-15

Client ID	Lab ID	% Moisture
H-662p-15:7.5	01-213-01	26
H-667p-15:7.5	01-213-02	12



### 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.
  - Z -
- ND - Not Detected at PQL  
 PQL - Practical Quantitation Limit  
 RPD - Relative Percent Difference





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February 24, 2015

Edwin Ptak  
Shannon & Wilson, Inc.  
400 N 34th Street, Suite 100  
Seattle, WA 98103

Re: Analytical Data for Project 21-1-20624-802  
Laboratory Reference No. 1502-107

Dear Edwin:

Enclosed are the analytical results and associated quality control data for samples submitted on February 11, 2015.

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,

A handwritten signature in black ink, appearing to read "D. Baumeister", with a long horizontal flourish extending to the right.

David Baumeister  
Project Manager

Enclosures

Date of Report: February 24, 2015  
Samples Submitted: February 11, 2015  
Laboratory Reference: 1502-107  
Project: 21-1-20624-802

### **Case Narrative**

Samples were collected on February 10, 2015 and received by the laboratory on February 11, 2015. They were maintained at the laboratory at a temperature of 2°C to 6°C.

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.

Date of Report: February 24, 2015  
 Samples Submitted: February 11, 2015  
 Laboratory Reference: 1502-107  
 Project: 21-1-20624-802

**NWTPH-HCID**

Matrix: Water  
 Units: mg/L (ppm)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>H-667p-15:GW</b>					
Laboratory ID:	02-107-01					
Gasoline Range Organics	<b>Detected</b>	0.10	NWTPH-HCID	2-13-15	2-13-15	
Diesel Range Organics	<b>ND</b>	0.26	NWTPH-HCID	2-13-15	2-13-15	
Lube Oil Range Organics	<b>ND</b>	0.41	NWTPH-HCID	2-13-15	2-13-15	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	79	50-150				

Date of Report: February 24, 2015  
 Samples Submitted: February 11, 2015  
 Laboratory Reference: 1502-107  
 Project: 21-1-20624-802

**NWTPH-HCID  
 QUALITY CONTROL**

Matrix: Water  
 Units: mg/L (ppm)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>METHOD BLANK</b>						
Laboratory ID:	MB0213W1					
Gasoline Range Organics	<b>ND</b>	0.10	NWTPH-HCID	2-13-15	2-13-15	
Diesel Range Organics	<b>ND</b>	0.25	NWTPH-HCID	2-13-15	2-13-15	
Lube Oil Range Organics	<b>ND</b>	0.40	NWTPH-HCID	2-13-15	2-13-15	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>87</i>	<i>50-150</i>				

Date of Report: February 24, 2015  
 Samples Submitted: February 11, 2015  
 Laboratory Reference: 1502-107  
 Project: 21-1-20624-802

**TOTAL PP METALS  
 EPA 200.8/7470A**

Matrix: Water  
 Units: ug/L (ppb)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>EPA Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
Lab ID:	02-107-01					
<b>Client ID:</b>	<b>H-667p-15:GW</b>					
Antimony	ND	5.6	200.8	2-17-15	2-17-15	
Arsenic	5.8	3.3	200.8	2-17-15	2-17-15	
Beryllium	ND	11	200.8	2-17-15	2-17-15	
Cadmium	ND	4.4	200.8	2-17-15	2-17-15	
Chromium	ND	11	200.8	2-17-15	2-17-15	
Copper	ND	11	200.8	2-17-15	2-17-15	
Lead	ND	1.1	200.8	2-17-15	2-17-15	
Mercury	ND	0.50	7470A	2-12-15	2-12-15	
Nickel	ND	22	200.8	2-17-15	2-17-15	
Selenium	ND	5.6	200.8	2-17-15	2-17-15	
Silver	ND	11	200.8	2-17-15	2-17-15	
Thallium	ND	5.6	200.8	2-17-15	2-17-15	
Zinc	ND	28	200.8	2-17-15	2-17-15	

Date of Report: February 24, 2015  
 Samples Submitted: February 11, 2015  
 Laboratory Reference: 1502-107  
 Project: 21-1-20624-802

**TOTAL PP METALS  
 EPA 200.8  
 METHOD BLANK QUALITY CONTROL**

Date Extracted: 2-17-15  
 Date Analyzed: 2-17-15  
 Matrix: Water  
 Units: ug/L (ppb)  
 Lab ID: MB0217WM1

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

Date of Report: February 24, 2015  
Samples Submitted: February 11, 2015  
Laboratory Reference: 1502-107  
Project: 21-1-20624-802

**TOTAL MERCURY  
EPA 7470A  
METHOD BLANK QUALITY CONTROL**

Date Extracted: 2-12-15  
Date Analyzed: 2-12-15  
  
Matrix: Water  
Units: ug/L (ppb)  
  
Lab ID: MB0212W1

Analyte	Method	Result	PQL
Mercury	7470A	<b>ND</b>	0.50

Date of Report: February 24, 2015  
 Samples Submitted: February 11, 2015  
 Laboratory Reference: 1502-107  
 Project: 21-1-20624-802

**TOTAL PP METALS  
 EPA 200.8  
 DUPLICATE QUALITY CONTROL**

Date Extracted: 2-17-15

Date Analyzed: 2-17-15

Matrix: Water

Units: ug/L (ppb)

Lab ID: 02-055-05

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	ND	ND	NA	28	

Date of Report: February 24, 2015  
Samples Submitted: February 11, 2015  
Laboratory Reference: 1502-107  
Project: 21-1-20624-802

**TOTAL MERCURY  
EPA 7470A  
DUPLICATE QUALITY CONTROL**

Date Extracted: 2-12-15

Date Analyzed: 2-12-15

Matrix: Water

Units: ug/L (ppb)

Lab ID: 02-107-01

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Mercury	<b>ND</b>	<b>ND</b>	NA	0.50	

Date of Report: February 24, 2015  
 Samples Submitted: February 11, 2015  
 Laboratory Reference: 1502-107  
 Project: 21-1-20624-802

**TOTAL PP METALS  
 EPA 200.8  
 MS/MSD QUALITY CONTROL**

Date Extracted: 2-17-15

Date Analyzed: 2-17-15

Matrix: Water

Units: ug/L (ppb)

Lab ID: 02-055-05

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Antimony	111	<b>115</b>	104	<b>115</b>	104	0	
Arsenic	111	<b>115</b>	103	<b>117</b>	105	2	
Beryllium	111	<b>109</b>	98	<b>109</b>	98	0	
Cadmium	111	<b>112</b>	101	<b>113</b>	102	0	
Chromium	111	<b>103</b>	93	<b>101</b>	91	2	
Copper	111	<b>100</b>	90	<b>98.8</b>	89	2	
Lead	111	<b>107</b>	97	<b>108</b>	98	1	
Nickel	111	<b>101</b>	91	<b>104</b>	93	2	
Selenium	111	<b>118</b>	107	<b>121</b>	109	2	
Silver	111	<b>107</b>	96	<b>106</b>	96	0	
Thallium	111	<b>110</b>	99	<b>112</b>	100	1	
Zinc	111	<b>120</b>	108	<b>118</b>	106	1	

Date of Report: February 24, 2015  
Samples Submitted: February 11, 2015  
Laboratory Reference: 1502-107  
Project: 21-1-20624-802

**TOTAL MERCURY  
EPA 7470A  
MS/MSD QUALITY CONTROL**

Date Extracted: 2-12-15

Date Analyzed: 2-12-15

Matrix: Water

Units: ug/L (ppb)

Lab ID: 02-107-01

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Mercury	12.5	<b>11.4</b>	91	<b>11.6</b>	93	2	

Date of Report: February 24, 2015  
 Samples Submitted: February 11, 2015  
 Laboratory Reference: 1502-107  
 Project: 21-1-20624-802

**DISSOLVED PP METALS**  
**EPA 200.8/7470A**

Matrix: Water  
 Units: ug/L (ppb)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>EPA Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
Lab ID:	02-107-01					
<b>Client ID:</b>	<b>H-667p-15:GW</b>					
Antimony	ND	5.0	200.8	2-11-15	2-19-15	
Arsenic	3.4	3.0	200.8	2-11-15	2-19-15	
Beryllium	ND	10	200.8	2-11-15	2-19-15	
Cadmium	ND	4.0	200.8	2-11-15	2-19-15	
Chromium	ND	10	200.8	2-11-15	2-19-15	
Copper	ND	10	200.8	2-11-15	2-19-15	
Lead	ND	1.0	200.8	2-11-15	2-19-15	
Mercury	ND	0.50	7470A	2-11-15	2-12-15	
Nickel	ND	20	200.8	2-11-15	2-19-15	
Selenium	ND	5.0	200.8	2-11-15	2-19-15	
Silver	ND	10	200.8	2-11-15	2-19-15	
Thallium	ND	5.0	200.8	2-11-15	2-19-15	
Zinc	ND	25	200.8	2-11-15	2-19-15	

Date of Report: February 24, 2015  
 Samples Submitted: February 11, 2015  
 Laboratory Reference: 1502-107  
 Project: 21-1-20624-802

**DISSOLVED PP METALS  
 EPA 200.8/7470A  
 METHOD BLANK QUALITY CONTROL**

Date Filtered: 2-11-15  
 Date Analyzed: 2-12&19-15  
 Matrix: Water  
 Units: ug/L (ppb)  
 Lab ID: MB0211F1

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

Date of Report: February 24, 2015  
 Samples Submitted: February 11, 2015  
 Laboratory Reference: 1502-107  
 Project: 21-1-20624-802

**DISSOLVED PP METALS  
 EPA 200.8/7470A  
 DUPLICATE QUALITY CONTROL**

Date Filtered: 2-11-15  
 Date Analyzed: 2-12&19-15

Matrix: Water  
 Units: ug/L (ppb)

Lab ID: 02-107-01

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Antimony	ND	ND	NA	5.0	
Arsenic	3.36	3.32	1	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	ND	ND	NA	25	

Date of Report: February 24, 2015  
 Samples Submitted: February 11, 2015  
 Laboratory Reference: 1502-107  
 Project: 21-1-20624-802

**DISSOLVED PP METALS  
 EPA 200.8/7470A  
 MS/MSD QUALITY CONTROL**

Date Filtered: 2-11-15  
 Date Analyzed: 2-12&19-15

Matrix: Water  
 Units: ug/L (ppb)

Lab ID: 02-107-01

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Antimony	200	<b>206</b>	103	<b>211</b>	105	2	
Arsenic	200	<b>203</b>	100	<b>205</b>	101	1	
Beryllium	200	<b>199</b>	99	<b>197</b>	98	1	
Cadmium	200	<b>199</b>	100	<b>200</b>	100	0	
Chromium	200	<b>192</b>	96	<b>191</b>	95	1	
Copper	200	<b>185</b>	92	<b>186</b>	93	0	
Lead	200	<b>196</b>	98	<b>196</b>	98	0	
Mercury	12.5	<b>11.2</b>	89	<b>11.6</b>	93	4	
Nickel	200	<b>192</b>	96	<b>191</b>	95	1	
Selenium	200	<b>223</b>	112	<b>223</b>	111	0	
Silver	200	<b>174</b>	87	<b>177</b>	89	2	
Thallium	200	<b>201</b>	101	<b>202</b>	101	0	
Zinc	200	<b>205</b>	103	<b>211</b>	105	3	

Date of Report: February 24, 2015  
 Samples Submitted: February 11, 2015  
 Laboratory Reference: 1502-107  
 Project: 21-1-20624-802

**NWTPH-Gx/BTEX**

Matrix: Water  
 Units: ug/L (ppb)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>H-667p-15:GW</b>					
Laboratory ID:	02-107-01					
Benzene	<b>ND</b>	1.0	EPA 8021B	2-20-15	2-20-15	
Toluene	<b>ND</b>	1.0	EPA 8021B	2-20-15	2-20-15	
Ethyl Benzene	<b>ND</b>	1.0	EPA 8021B	2-20-15	2-20-15	
m,p-Xylene	<b>1.6</b>	1.0	EPA 8021B	2-20-15	2-20-15	
o-Xylene	<b>ND</b>	1.0	EPA 8021B	2-20-15	2-20-15	
Gasoline	<b>170</b>	100	NWTPH-Gx	2-20-15	2-20-15	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	<i>104</i>	<i>71-113</i>				

Date of Report: February 24, 2015  
 Samples Submitted: February 11, 2015  
 Laboratory Reference: 1502-107  
 Project: 21-1-20624-802

**NWTPH-Gx/BTEX  
 QUALITY CONTROL**

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB0220W1					
Benzene	ND	1.0	EPA 8021B	2-20-15	2-20-15	
Toluene	ND	1.0	EPA 8021B	2-20-15	2-20-15	
Ethyl Benzene	ND	1.0	EPA 8021B	2-20-15	2-20-15	
m,p-Xylene	ND	1.0	EPA 8021B	2-20-15	2-20-15	
o-Xylene	ND	1.0	EPA 8021B	2-20-15	2-20-15	
Gasoline	ND	100	NWTPH-Gx	2-20-15	2-20-15	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	100	71-113				

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>								
Laboratory ID:	02-173-05							
	ORIG	DUP						
Benzene	ND	ND	NA	NA	NA	NA	NA	30
Toluene	ND	ND	NA	NA	NA	NA	NA	30
Ethyl Benzene	ND	ND	NA	NA	NA	NA	NA	30
m,p-Xylene	ND	ND	NA	NA	NA	NA	NA	30
o-Xylene	ND	ND	NA	NA	NA	NA	NA	30
Gasoline	ND	ND	NA	NA	NA	NA	NA	30
<i>Surrogate:</i>								
<i>Fluorobenzene</i>				103	103	71-113		

**SPIKE BLANKS**

Laboratory ID:	SB0220W1								
	SB	SBD	SB	SBD	SB	SBD			
Benzene	55.0	52.7	50.0	50.0	110	105	80-118	4	11
Toluene	56.8	54.4	50.0	50.0	114	109	81-119	4	11
Ethyl Benzene	55.7	53.2	50.0	50.0	111	106	80-121	5	12
m,p-Xylene	56.0	53.7	50.0	50.0	112	107	81-121	4	12
o-Xylene	54.8	52.7	50.0	50.0	110	105	81-119	4	12
<i>Surrogate:</i>									
<i>Fluorobenzene</i>					105	105	71-113		



### 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.
  - Z -
- ND - Not Detected at PQL  
 PQL - Practical Quantitation Limit  
 RPD - Relative Percent Difference



Analytical Laboratory Testing Services  
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# Chain of Custody

Laboratory Number: **02-107**

Company: Shannon & Wilson  
 Project Number: 21-20624-802  
 Project Name: SR 520  
 Project Manager: Ed Ptak  
 Sampled by: Jeff Weirs

Turnaround Request  
(in working days)

- (Check One)
- Same Day  1 Day
  - 2 Days  3 Days
  - Standard (7 Days)  
(TPH analysis 5 Days)
  - \_\_\_\_\_ (other)

Number of Containers

Date Sampled: 2/10/15 Time Sampled: 1600 Matrix: W

9  X

Lab ID: 1 H-667p-15:5W

Parameter	Result
% Moisture	
HEM (oil and grease) 1664A	
TCLP Metals	
Total MTCA Metals	
Total RCRA Metals	
Chlorinated Acid Herbicides 8151A	
Organophosphorus Pesticides 8270D/SIM	
Organochlorine Pesticides 8081B	
PCBs 8082A	
PAHs 8270D/SIM (low-level)	
Semivolatiles 8270D/SIM (with low-level PAHs)	
Halogenated Volatiles 8260C	
Volatiles 8260C	
NWTPH-Dx	
NWTPH-Gx	
NWTPH-Gx/BTEX	
NWTPH-HCID	
Total Priority Metals	<u>X</u>
Dissolved Metals	<u>X</u>

Signature	Company	Date	Time	Comments/Special Instructions
<u>[Signature]</u>	<u>SKW</u>	<u>2/10/15</u>	<u>700</u>	<u>Added 2/20/15. P13 (STA)</u>
<u>[Signature]</u>	<u>SPZOP</u>	<u>2/11/15</u>	<u>1016</u>	
<u>[Signature]</u>	<u>SPASH</u>	<u>2/11/15</u>	<u>1205</u>	
<u>[Signature]</u>	<u>QBE</u>	<u>2/11/15</u>	<u>1205</u>	
<u>[Signature]</u>				
Relinquished				
Received				
Relinquished				
Received				
Relinquished				
Received				
Reviewed/Date				Chromatograms with final report <input type="checkbox"/>



14648 NE 95<sup>th</sup> Street, Redmond, WA 98052 • (425) 883-3881

March 24, 2016

Edwin Ptak  
Shannon & Wilson, Inc.  
400 N 34th Street, Suite 100  
Seattle, WA 98103

Re: Analytical Data for Project 21-1-20624-822  
Laboratory Reference No. 1603-150

Dear Edwin:

Enclosed are the analytical results and associated quality control data for samples submitted on March 16, 2016.

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,

A handwritten signature in black ink, appearing to read "DB", with a long horizontal stroke extending to the right.

David Baumeister  
Project Manager

Enclosures

Date of Report: March 24, 2016  
Samples Submitted: March 16, 2016  
Laboratory Reference: 1603-150  
Project: 21-1-20624-822

### **Case Narrative**

Samples were collected on March 15, 2016 and received by the laboratory on March 16, 2016. 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 6010C/7471B Analysis

The duplicate RPD for Chromium is outside control limits due to sample inhomogeneity. The sample was 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.

Date of Report: March 24, 2016  
 Samples Submitted: March 16, 2016  
 Laboratory Reference: 1603-150  
 Project: 21-1-20624-822

### NWTPH-HCID

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-691p-16:9</b>					
Laboratory ID:	03-150-01					
Gasoline Range Organics	<b>ND</b>	22	NWTPH-HCID	3-17-16	3-17-16	
Diesel Range Organics	<b>ND</b>	56	NWTPH-HCID	3-17-16	3-17-16	
Lube Oil Range Organics	<b>ND</b>	110	NWTPH-HCID	3-17-16	3-17-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>119</i>	<i>50-150</i>				

Date of Report: March 24, 2016  
 Samples Submitted: March 16, 2016  
 Laboratory Reference: 1603-150  
 Project: 21-1-20624-822

**NWTPH-HCID  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg (ppm)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>METHOD BLANK</b>						
Laboratory ID:	MB0317S2					
Gasoline Range Organics	<b>ND</b>	20	NWTPH-HCID	3-17-16	3-17-16	
Diesel Range Organics	<b>ND</b>	50	NWTPH-HCID	3-17-16	3-17-16	
Lube Oil Range Organics	<b>ND</b>	100	NWTPH-HCID	3-17-16	3-17-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>125</i>	<i>50-150</i>				

Date of Report: March 24, 2016  
 Samples Submitted: March 16, 2016  
 Laboratory Reference: 1603-150  
 Project: 21-1-20624-822

**TOTAL METALS  
 EPA 6010C/7471B**

Matrix: Soil  
 Units: mg/kg (ppm)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>EPA Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
Lab ID:	03-150-01					
<b>Client ID:</b>	<b>H-691p-16:9</b>					
Arsenic	<b>ND</b>	11	6010C	3-23-16	3-23-16	
Barium	<b>37</b>	2.8	6010C	3-23-16	3-23-16	
Cadmium	<b>ND</b>	0.56	6010C	3-23-16	3-23-16	
Chromium	<b>32</b>	0.56	6010C	3-23-16	3-23-16	
Lead	<b>ND</b>	5.6	6010C	3-23-16	3-23-16	
Mercury	<b>ND</b>	0.28	7471B	3-21-16	3-21-16	
Selenium	<b>ND</b>	11	6010C	3-23-16	3-23-16	
Silver	<b>ND</b>	1.1	6010C	3-23-16	3-23-16	

Date of Report: March 24, 2016  
 Samples Submitted: March 16, 2016  
 Laboratory Reference: 1603-150  
 Project: 21-1-20624-822

**TOTAL METALS  
 EPA 6010C/7471B  
 METHOD BLANK QUALITY CONTROL**

Date Extracted: 3-21&23-16  
 Date Analyzed: 3-21&23-16  
 Matrix: Soil  
 Units: mg/kg (ppm)  
 Lab ID: MB0323SM1&MB0321S1

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
Mercury	7471B	ND	0.25
Selenium	6010C	ND	10
Silver	6010C	ND	1.0

Date of Report: March 24, 2016  
 Samples Submitted: March 16, 2016  
 Laboratory Reference: 1603-150  
 Project: 21-1-20624-822

**TOTAL METALS  
 EPA 6010C/7471B  
 DUPLICATE QUALITY CONTROL**

Date Extracted: 3-21&23-16

Date Analyzed: 3-21&23-16

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-150-01

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Arsenic	ND	ND	NA	10	
Barium	32.9	30.9	6	2.5	
Cadmium	ND	ND	NA	0.50	
Chromium	28.3	36.9	27	0.50	K
Lead	ND	ND	NA	5.0	
Mercury	ND	ND	NA	0.25	
Selenium	ND	ND	NA	10	
Silver	ND	ND	NA	1.0	

Date of Report: March 24, 2016  
 Samples Submitted: March 16, 2016  
 Laboratory Reference: 1603-150  
 Project: 21-1-20624-822

**TOTAL METALS  
 EPA 6010C/7471B  
 MS/MSD QUALITY CONTROL**

Date Extracted: 3-21&23-16

Date Analyzed: 3-21&23-16

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-150-01

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Arsenic	100	<b>100</b>	100	<b>99.0</b>	99	1	
Barium	100	<b>123</b>	91	<b>123</b>	90	0	
Cadmium	50.0	<b>48.6</b>	97	<b>48.2</b>	96	1	
Chromium	100	<b>112</b>	84	<b>112</b>	84	0	
Lead	250	<b>217</b>	87	<b>216</b>	86	0	
Mercury	0.500	<b>0.458</b>	92	<b>0.461</b>	92	1	
Selenium	100	<b>93.4</b>	93	<b>90.6</b>	91	3	
Silver	25.0	<b>22.8</b>	91	<b>22.4</b>	90	2	

Date of Report: March 24, 2016  
Samples Submitted: March 16, 2016  
Laboratory Reference: 1603-150  
Project: 21-1-20624-822

### % MOISTURE

Date Analyzed: 3-18-16

Client ID	Lab ID	% Moisture
H-691p-16:9	03-150-01	10



### 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.
  - Z -
- ND - Not Detected at PQL  
 PQL - Practical Quantitation Limit  
 RPD - Relative Percent Difference

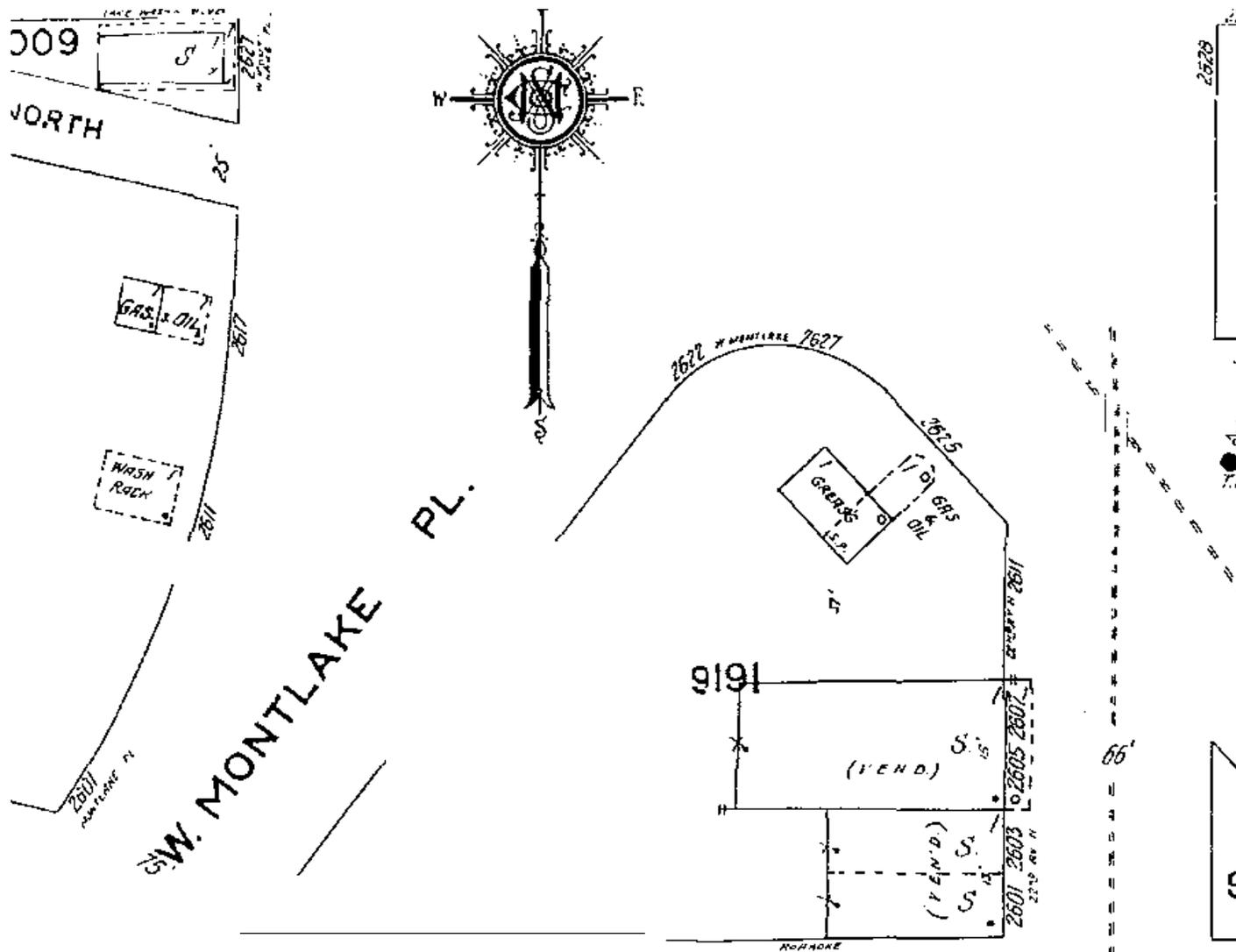




**Appendix C**  
**1930 Sanborn Map**

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Portion of 1930 Sanborn Map.



**Appendix D**  
**Analytical Reports**

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14648 NE 95<sup>th</sup> Street, Redmond, WA 98052 • (425) 883-3881

October 17, 2016

Glenn Hayman  
INNOVEX Environmental Mgt., Inc.  
16310 NE 80th St., Suite 300  
Redmond, WA 98052

Re: Analytical Data for Project 31008  
Laboratory Reference No. 1610-071

Dear Glenn:

Enclosed are the analytical results and associated quality control data for samples submitted on October 7, 2016.

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,

A handwritten signature in black ink, appearing to read "DB", with a long horizontal flourish extending to the right.

David Baumeister  
Project Manager

Enclosures



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OnSite Environmental, Inc. 14648 NE 95<sup>th</sup> Street, Redmond, WA 98052 (425) 883-3881

This report pertains to the samples analyzed in accordance with the chain of custody, and is intended only for the use of the individual or company to whom it is addressed.

Date of Report: October 17, 2016  
Samples Submitted: October 7, 2016  
Laboratory Reference: 1610-071  
Project: 31008

### Case Narrative

Samples were collected on October 6, 2016 and received by the laboratory on October 7, 2016. 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.

#### Volatiles EPA 8260C Analysis

The Method 5035A VOA vials provided for sample H-1-16-10 contained too much soil to perform the requested analysis. Therefore, the sample was extracted from an 8-ounce jar and analyzed. Some loss of volatiles may have occurred.

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.



Date of Report: October 17, 2016  
 Samples Submitted: October 7, 2016  
 Laboratory Reference: 1610-071  
 Project: 31008

### NWTPH-HCID

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-1-16-10</b>					
Laboratory ID:	10-071-02					
Gasoline Range Organics	<b>ND</b>	23	NWTPH-HCID	10-10-16	10-10-16	
Diesel Range Organics	<b>ND</b>	58	NWTPH-HCID	10-10-16	10-10-16	
Lube Oil Range Organics	<b>ND</b>	120	NWTPH-HCID	10-10-16	10-10-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>101</i>	<i>50-150</i>				



Date of Report: October 17, 2016  
 Samples Submitted: October 7, 2016  
 Laboratory Reference: 1610-071  
 Project: 31008

**NWTPH-HCID  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg (ppm)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>METHOD BLANK</b>						
Laboratory ID:	MB1010S2					
Gasoline Range Organics	<b>ND</b>	20	NWTPH-HCID	10-10-16	10-10-16	
Diesel Range Organics	<b>ND</b>	50	NWTPH-HCID	10-10-16	10-10-16	
Lube Oil Range Organics	<b>ND</b>	100	NWTPH-HCID	10-10-16	10-10-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>84</i>	<i>50-150</i>				



Date of Report: October 17, 2016  
 Samples Submitted: October 7, 2016  
 Laboratory Reference: 1610-071  
 Project: 31008

**VOLATILES EPA 8260C**  
 page 1 of 2

Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>H-1-16-10</b>					
<b>Laboratory ID:</b>	<b>10-071-02</b>					
Dichlorodifluoromethane	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
Chloromethane	ND	0.0095	EPA 8260C	10-7-16	10-7-16	
Vinyl Chloride	ND	0.0019	EPA 8260C	10-7-16	10-7-16	
Bromomethane	ND	0.0015	EPA 8260C	10-7-16	10-7-16	
Chloroethane	ND	0.0097	EPA 8260C	10-7-16	10-7-16	
Trichlorofluoromethane	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
1,1-Dichloroethene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
Acetone	ND	0.0083	EPA 8260C	10-7-16	10-7-16	
Iodomethane	ND	0.0058	EPA 8260C	10-7-16	10-7-16	
Carbon Disulfide	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
Methylene Chloride	ND	0.0058	EPA 8260C	10-7-16	10-7-16	
(trans) 1,2-Dichloroethene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
Methyl t-Butyl Ether	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
1,1-Dichloroethane	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
Vinyl Acetate	ND	0.0058	EPA 8260C	10-7-16	10-7-16	
2,2-Dichloropropane	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
(cis) 1,2-Dichloroethene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
2-Butanone	ND	0.0058	EPA 8260C	10-7-16	10-7-16	
Bromochloromethane	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
Chloroform	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
1,1,1-Trichloroethane	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
Carbon Tetrachloride	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
1,1-Dichloropropene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
Benzene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
1,2-Dichloroethane	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
Trichloroethene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
1,2-Dichloropropane	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
Dibromomethane	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
Bromodichloromethane	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
2-Chloroethyl Vinyl Ether	ND	0.0058	EPA 8260C	10-7-16	10-7-16	
(cis) 1,3-Dichloropropene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
Methyl Isobutyl Ketone	ND	0.0058	EPA 8260C	10-7-16	10-7-16	
Toluene	ND	0.0058	EPA 8260C	10-7-16	10-7-16	
(trans) 1,3-Dichloropropene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	



Date of Report: October 17, 2016  
 Samples Submitted: October 7, 2016  
 Laboratory Reference: 1610-071  
 Project: 31008

**VOLATILES EPA 8260C**  
 page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-1-16-10</b>					
Laboratory ID:	10-071-02					
1,1,2-Trichloroethane	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
Tetrachloroethene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
1,3-Dichloropropane	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
2-Hexanone	ND	0.0058	EPA 8260C	10-7-16	10-7-16	
Dibromochloromethane	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
1,2-Dibromoethane	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
Chlorobenzene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
1,1,1,2-Tetrachloroethane	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
Ethylbenzene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
m,p-Xylene	ND	0.0023	EPA 8260C	10-7-16	10-7-16	
o-Xylene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
Styrene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
Bromoform	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
Isopropylbenzene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
Bromobenzene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
1,1,2,2-Tetrachloroethane	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
1,2,3-Trichloropropane	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
n-Propylbenzene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
2-Chlorotoluene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
4-Chlorotoluene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
1,3,5-Trimethylbenzene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
tert-Butylbenzene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
1,2,4-Trimethylbenzene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
sec-Butylbenzene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
1,3-Dichlorobenzene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
p-Isopropyltoluene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
1,4-Dichlorobenzene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
1,2-Dichlorobenzene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
n-Butylbenzene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
1,2-Dibromo-3-chloropropane	ND	0.0058	EPA 8260C	10-7-16	10-7-16	
1,2,4-Trichlorobenzene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
Hexachlorobutadiene	ND	0.0058	EPA 8260C	10-7-16	10-7-16	
Naphthalene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
1,2,3-Trichlorobenzene	ND	0.0012	EPA 8260C	10-7-16	10-7-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>99</i>	<i>76-131</i>				
<i>Toluene-d8</i>	<i>104</i>	<i>80-126</i>				
<i>4-Bromofluorobenzene</i>	<i>101</i>	<i>60-146</i>				



Date of Report: October 17, 2016  
 Samples Submitted: October 7, 2016  
 Laboratory Reference: 1610-071  
 Project: 31008

**VOLATILES by EPA 8260C**  
**METHOD BLANK QUALITY CONTROL**  
 page 1 of 2

Matrix: Soil  
 Units: mg/kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1007S1					
Dichlorodifluoromethane	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
Chloromethane	ND	0.0082	EPA 8260C	10-7-16	10-7-16	
Vinyl Chloride	ND	0.0016	EPA 8260C	10-7-16	10-7-16	
Bromomethane	ND	0.0013	EPA 8260C	10-7-16	10-7-16	
Chloroethane	ND	0.0083	EPA 8260C	10-7-16	10-7-16	
Trichlorofluoromethane	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
1,1-Dichloroethene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
Acetone	ND	0.0071	EPA 8260C	10-7-16	10-7-16	
Iodomethane	ND	0.0050	EPA 8260C	10-7-16	10-7-16	
Carbon Disulfide	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
Methylene Chloride	ND	0.0050	EPA 8260C	10-7-16	10-7-16	
(trans) 1,2-Dichloroethene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
Methyl t-Butyl Ether	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
1,1-Dichloroethane	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
Vinyl Acetate	ND	0.0050	EPA 8260C	10-7-16	10-7-16	
2,2-Dichloropropane	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
(cis) 1,2-Dichloroethene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
2-Butanone	ND	0.0050	EPA 8260C	10-7-16	10-7-16	
Bromochloromethane	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
Chloroform	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
1,1,1-Trichloroethane	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
Carbon Tetrachloride	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
1,1-Dichloropropene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
Benzene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
1,2-Dichloroethane	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
Trichloroethene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
1,2-Dichloropropane	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
Dibromomethane	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
Bromodichloromethane	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
2-Chloroethyl Vinyl Ether	ND	0.0050	EPA 8260C	10-7-16	10-7-16	
(cis) 1,3-Dichloropropene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
Methyl Isobutyl Ketone	ND	0.0050	EPA 8260C	10-7-16	10-7-16	
Toluene	ND	0.0050	EPA 8260C	10-7-16	10-7-16	
(trans) 1,3-Dichloropropene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	



Date of Report: October 17, 2016  
 Samples Submitted: October 7, 2016  
 Laboratory Reference: 1610-071  
 Project: 31008

**VOLATILES by EPA 8260C**  
**METHOD BLANK QUALITY CONTROL**  
 page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:		MB1007S1				
1,1,2-Trichloroethane	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
Tetrachloroethene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
1,3-Dichloropropane	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
2-Hexanone	ND	0.0050	EPA 8260C	10-7-16	10-7-16	
Dibromochloromethane	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
1,2-Dibromoethane	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
Chlorobenzene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
1,1,1,2-Tetrachloroethane	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
Ethylbenzene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
m,p-Xylene	ND	0.0020	EPA 8260C	10-7-16	10-7-16	
o-Xylene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
Styrene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
Bromoform	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
Isopropylbenzene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
Bromobenzene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
1,1,1,2-Tetrachloroethane	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
1,2,3-Trichloropropane	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
n-Propylbenzene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
2-Chlorotoluene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
4-Chlorotoluene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
1,3,5-Trimethylbenzene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
tert-Butylbenzene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
1,2,4-Trimethylbenzene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
sec-Butylbenzene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
1,3-Dichlorobenzene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
p-Isopropyltoluene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
1,4-Dichlorobenzene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
1,2-Dichlorobenzene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
n-Butylbenzene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
1,2-Dibromo-3-chloropropane	ND	0.0050	EPA 8260C	10-7-16	10-7-16	
1,2,4-Trichlorobenzene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
Hexachlorobutadiene	ND	0.0050	EPA 8260C	10-7-16	10-7-16	
Naphthalene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
1,2,3-Trichlorobenzene	ND	0.0010	EPA 8260C	10-7-16	10-7-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>108</i>	<i>76-131</i>				
<i>Toluene-d8</i>	<i>111</i>	<i>80-126</i>				
<i>4-Bromofluorobenzene</i>	<i>108</i>	<i>60-146</i>				



Date of Report: October 17, 2016  
 Samples Submitted: October 7, 2016  
 Laboratory Reference: 1610-071  
 Project: 31008

**VOLATILES by EPA 8260C  
 SB/SBD QUALITY CONTROL**

Matrix: Soil  
 Units: mg/kg

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD		Flags
					SB	SBD	Limits	RPD	Limit	
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB1007S1									
	SB	SBD	SB	SBD	SB	SBD				
1,1-Dichloroethene	<b>0.0402</b>	<b>0.0409</b>	0.0500	0.0500	80	82	68-126	2	15	
Benzene	<b>0.0488</b>	<b>0.0523</b>	0.0500	0.0500	98	105	70-121	7	15	
Trichloroethene	<b>0.0470</b>	<b>0.0494</b>	0.0500	0.0500	94	99	75-120	5	15	
Toluene	<b>0.0500</b>	<b>0.0525</b>	0.0500	0.0500	100	105	80-120	5	15	
Chlorobenzene	<b>0.0475</b>	<b>0.0502</b>	0.0500	0.0500	95	100	76-120	6	15	
<i>Surrogate:</i>										
Dibromofluoromethane					100	102	76-131			
Toluene-d8					101	101	80-126			
4-Bromofluorobenzene					102	99	60-146			



Date of Report: October 17, 2016  
 Samples Submitted: October 7, 2016  
 Laboratory Reference: 1610-071  
 Project: 31008

### SEMIVOLATILES EPA 8270D/SIM

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Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-1-16-10</b>					
<b>Laboratory ID:</b>	<b>10-071-02</b>					
n-Nitrosodimethylamine	ND	0.039	EPA 8270D	10-10-16	10-11-16	
Pyridine	ND	0.39	EPA 8270D	10-10-16	10-11-16	
Phenol	ND	0.039	EPA 8270D	10-10-16	10-11-16	
Aniline	ND	0.19	EPA 8270D	10-10-16	10-11-16	
bis(2-Chloroethyl)ether	ND	0.039	EPA 8270D	10-10-16	10-11-16	
2-Chlorophenol	ND	0.039	EPA 8270D	10-10-16	10-11-16	
1,3-Dichlorobenzene	ND	0.039	EPA 8270D	10-10-16	10-11-16	
1,4-Dichlorobenzene	ND	0.039	EPA 8270D	10-10-16	10-11-16	
Benzyl alcohol	ND	0.19	EPA 8270D	10-10-16	10-11-16	
1,2-Dichlorobenzene	ND	0.039	EPA 8270D	10-10-16	10-11-16	
2-Methylphenol (o-Cresol)	ND	0.039	EPA 8270D	10-10-16	10-11-16	
bis(2-Chloroisopropyl)ether	ND	0.039	EPA 8270D	10-10-16	10-11-16	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.039	EPA 8270D	10-10-16	10-11-16	
n-Nitroso-di-n-propylamine	ND	0.039	EPA 8270D	10-10-16	10-11-16	
Hexachloroethane	ND	0.039	EPA 8270D	10-10-16	10-11-16	
Nitrobenzene	ND	0.039	EPA 8270D	10-10-16	10-11-16	
Isophorone	ND	0.039	EPA 8270D	10-10-16	10-11-16	
2-Nitrophenol	ND	0.039	EPA 8270D	10-10-16	10-11-16	
2,4-Dimethylphenol	ND	0.039	EPA 8270D	10-10-16	10-11-16	
bis(2-Chloroethoxy)methane	ND	0.039	EPA 8270D	10-10-16	10-11-16	
2,4-Dichlorophenol	ND	0.039	EPA 8270D	10-10-16	10-11-16	
1,2,4-Trichlorobenzene	ND	0.039	EPA 8270D	10-10-16	10-11-16	
Naphthalene	ND	0.0077	EPA 8270D/SIM	10-10-16	10-10-16	
4-Chloroaniline	ND	0.19	EPA 8270D	10-10-16	10-11-16	
Hexachlorobutadiene	ND	0.039	EPA 8270D	10-10-16	10-11-16	
4-Chloro-3-methylphenol	ND	0.039	EPA 8270D	10-10-16	10-11-16	
2-Methylnaphthalene	ND	0.0077	EPA 8270D/SIM	10-10-16	10-10-16	
1-Methylnaphthalene	ND	0.0077	EPA 8270D/SIM	10-10-16	10-10-16	
Hexachlorocyclopentadiene	ND	0.039	EPA 8270D	10-10-16	10-11-16	
2,4,6-Trichlorophenol	ND	0.039	EPA 8270D	10-10-16	10-11-16	
2,3-Dichloroaniline	ND	0.039	EPA 8270D	10-10-16	10-11-16	
2,4,5-Trichlorophenol	ND	0.039	EPA 8270D	10-10-16	10-11-16	
2-Chloronaphthalene	ND	0.039	EPA 8270D	10-10-16	10-11-16	
2-Nitroaniline	ND	0.039	EPA 8270D	10-10-16	10-11-16	
1,4-Dinitrobenzene	ND	0.039	EPA 8270D	10-10-16	10-11-16	
Dimethylphthalate	ND	0.039	EPA 8270D	10-10-16	10-11-16	
1,3-Dinitrobenzene	ND	0.039	EPA 8270D	10-10-16	10-11-16	
2,6-Dinitrotoluene	ND	0.039	EPA 8270D	10-10-16	10-11-16	
1,2-Dinitrobenzene	ND	0.039	EPA 8270D	10-10-16	10-11-16	
Acenaphthylene	ND	0.0077	EPA 8270D/SIM	10-10-16	10-10-16	
3-Nitroaniline	ND	0.039	EPA 8270D	10-10-16	10-11-16	



Date of Report: October 17, 2016  
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 Project: 31008

**SEMIVOLATILES EPA 8270D/SIM**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-1-16-10</b>					
<b>Laboratory ID:</b>	<b>10-071-02</b>					
2,4-Dinitrophenol	ND	0.19	EPA 8270D	10-10-16	10-11-16	
Acenaphthene	ND	0.0077	EPA 8270D/SIM	10-10-16	10-10-16	
4-Nitrophenol	ND	0.039	EPA 8270D	10-10-16	10-11-16	
2,4-Dinitrotoluene	ND	0.039	EPA 8270D	10-10-16	10-11-16	
Dibenzofuran	ND	0.039	EPA 8270D	10-10-16	10-11-16	
2,3,5,6-Tetrachlorophenol	ND	0.039	EPA 8270D	10-10-16	10-11-16	
2,3,4,6-Tetrachlorophenol	ND	0.039	EPA 8270D	10-10-16	10-11-16	
Diethylphthalate	ND	0.19	EPA 8270D	10-10-16	10-11-16	
4-Chlorophenyl-phenylether	ND	0.039	EPA 8270D	10-10-16	10-11-16	
4-Nitroaniline	ND	0.039	EPA 8270D	10-10-16	10-11-16	
Fluorene	ND	0.0077	EPA 8270D/SIM	10-10-16	10-10-16	
4,6-Dinitro-2-methylphenol	ND	0.19	EPA 8270D	10-10-16	10-11-16	
n-Nitrosodiphenylamine	ND	0.039	EPA 8270D	10-10-16	10-11-16	
1,2-Diphenylhydrazine	ND	0.039	EPA 8270D	10-10-16	10-11-16	
4-Bromophenyl-phenylether	ND	0.039	EPA 8270D	10-10-16	10-11-16	
Hexachlorobenzene	ND	0.039	EPA 8270D	10-10-16	10-11-16	
Pentachlorophenol	ND	0.19	EPA 8270D	10-10-16	10-11-16	
Phenanthrene	ND	0.0077	EPA 8270D/SIM	10-10-16	10-10-16	
Anthracene	ND	0.0077	EPA 8270D/SIM	10-10-16	10-10-16	
Carbazole	ND	0.039	EPA 8270D	10-10-16	10-11-16	
Di-n-butylphthalate	ND	0.19	EPA 8270D	10-10-16	10-11-16	
Fluoranthene	ND	0.0077	EPA 8270D/SIM	10-10-16	10-10-16	
Benzidine	ND	0.39	EPA 8270D	10-10-16	10-11-16	
Pyrene	ND	0.0077	EPA 8270D/SIM	10-10-16	10-10-16	
Butylbenzylphthalate	ND	0.039	EPA 8270D	10-10-16	10-11-16	
bis-2-Ethylhexyladipate	ND	0.039	EPA 8270D	10-10-16	10-11-16	
3,3'-Dichlorobenzidine	ND	0.19	EPA 8270D	10-10-16	10-11-16	
Benzo[a]anthracene	ND	0.0077	EPA 8270D/SIM	10-10-16	10-10-16	
Chrysene	ND	0.0077	EPA 8270D/SIM	10-10-16	10-10-16	
bis(2-Ethylhexyl)phthalate	ND	0.039	EPA 8270D	10-10-16	10-11-16	
Di-n-octylphthalate	ND	0.039	EPA 8270D	10-10-16	10-11-16	
Benzo[b]fluoranthene	ND	0.0077	EPA 8270D/SIM	10-10-16	10-10-16	
Benzo(j,k)fluoranthene	ND	0.0077	EPA 8270D/SIM	10-10-16	10-10-16	
Benzo[a]pyrene	ND	0.0077	EPA 8270D/SIM	10-10-16	10-10-16	
Indeno[1,2,3-cd]pyrene	ND	0.0077	EPA 8270D/SIM	10-10-16	10-10-16	
Dibenz[a,h]anthracene	ND	0.0077	EPA 8270D/SIM	10-10-16	10-10-16	
Benzo[g,h,i]perylene	ND	0.0077	EPA 8270D/SIM	10-10-16	10-10-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorophenol</i>	<i>57</i>	<i>24 - 117</i>				
<i>Phenol-d6</i>	<i>60</i>	<i>30 - 120</i>				
<i>Nitrobenzene-d5</i>	<i>61</i>	<i>27 - 112</i>				
<i>2-Fluorobiphenyl</i>	<i>62</i>	<i>35 - 113</i>				
<i>2,4,6-Tribromophenol</i>	<i>65</i>	<i>21 - 120</i>				
<i>Terphenyl-d14</i>	<i>63</i>	<i>39 - 121</i>				



Date of Report: October 17, 2016  
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 Laboratory Reference: 1610-071  
 Project: 31008

**SEMIVOLATILES EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
 page 1 of 2

Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1010S1					
n-Nitrosodimethylamine	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Pyridine	ND	0.33	EPA 8270D	10-10-16	10-11-16	
Phenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Aniline	ND	0.17	EPA 8270D	10-10-16	10-11-16	
bis(2-Chloroethyl)ether	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2-Chlorophenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
1,3-Dichlorobenzene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
1,4-Dichlorobenzene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Benzyl alcohol	ND	0.17	EPA 8270D	10-10-16	10-11-16	
1,2-Dichlorobenzene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2-Methylphenol (o-Cresol)	ND	0.033	EPA 8270D	10-10-16	10-11-16	
bis(2-Chloroisopropyl)ether	ND	0.033	EPA 8270D	10-10-16	10-11-16	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.033	EPA 8270D	10-10-16	10-11-16	
n-Nitroso-di-n-propylamine	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Hexachloroethane	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Nitrobenzene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Isophorone	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2-Nitrophenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2,4-Dimethylphenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
bis(2-Chloroethoxy)methane	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2,4-Dichlorophenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
1,2,4-Trichlorobenzene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Naphthalene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
4-Chloroaniline	ND	0.17	EPA 8270D	10-10-16	10-11-16	
Hexachlorobutadiene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
4-Chloro-3-methylphenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2-Methylnaphthalene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
1-Methylnaphthalene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Hexachlorocyclopentadiene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2,4,6-Trichlorophenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2,3-Dichloroaniline	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2,4,5-Trichlorophenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2-Chloronaphthalene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2-Nitroaniline	ND	0.033	EPA 8270D	10-10-16	10-11-16	
1,4-Dinitrobenzene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Dimethylphthalate	ND	0.033	EPA 8270D	10-10-16	10-11-16	
1,3-Dinitrobenzene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2,6-Dinitrotoluene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
1,2-Dinitrobenzene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Acenaphthylene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
3-Nitroaniline	ND	0.033	EPA 8270D	10-10-16	10-11-16	



Date of Report: October 17, 2016  
 Samples Submitted: October 7, 2016  
 Laboratory Reference: 1610-071  
 Project: 31008

**SEMIVOLATILES EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1010S1					
2,4-Dinitrophenol	ND	0.17	EPA 8270D	10-10-16	10-11-16	
Acenaphthene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
4-Nitrophenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2,4-Dinitrotoluene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Dibenzofuran	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2,3,5,6-Tetrachlorophenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2,3,4,6-Tetrachlorophenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Diethylphthalate	ND	0.17	EPA 8270D	10-10-16	10-11-16	
4-Chlorophenyl-phenylether	ND	0.033	EPA 8270D	10-10-16	10-11-16	
4-Nitroaniline	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Fluorene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
4,6-Dinitro-2-methylphenol	ND	0.17	EPA 8270D	10-10-16	10-11-16	
n-Nitrosodiphenylamine	ND	0.033	EPA 8270D	10-10-16	10-11-16	
1,2-Diphenylhydrazine	ND	0.033	EPA 8270D	10-10-16	10-11-16	
4-Bromophenyl-phenylether	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Hexachlorobenzene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Pentachlorophenol	ND	0.17	EPA 8270D	10-10-16	10-11-16	
Phenanthrene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Anthracene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Carbazole	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Di-n-butylphthalate	ND	0.17	EPA 8270D	10-10-16	10-11-16	
Fluoranthene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Benzidine	ND	0.33	EPA 8270D	10-10-16	10-11-16	
Pyrene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Butylbenzylphthalate	ND	0.033	EPA 8270D	10-10-16	10-11-16	
bis-2-Ethylhexyladipate	ND	0.033	EPA 8270D	10-10-16	10-11-16	
3,3'-Dichlorobenzidine	ND	0.17	EPA 8270D	10-10-16	10-11-16	
Benzo[a]anthracene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Chrysene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
bis(2-Ethylhexyl)phthalate	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Di-n-octylphthalate	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Benzo[a]pyrene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Indeno[1,2,3-cd]pyrene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Benzo[g,h,i]perylene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorophenol</i>	<i>81</i>	<i>24 - 117</i>				
<i>Phenol-d6</i>	<i>83</i>	<i>30 - 120</i>				
<i>Nitrobenzene-d5</i>	<i>85</i>	<i>27 - 112</i>				
<i>2-Fluorobiphenyl</i>	<i>84</i>	<i>35 - 113</i>				
<i>2,4,6-Tribromophenol</i>	<i>85</i>	<i>21 - 120</i>				
<i>Terphenyl-d14</i>	<i>85</i>	<i>39 - 121</i>				



Date of Report: October 17, 2016  
 Samples Submitted: October 7, 2016  
 Laboratory Reference: 1610-071  
 Project: 31008

**SEMIVOLATILES EPA 8270D/SIM  
 MS/MSD QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result		Spike Level		Source	Percent		Recovery		RPD	RPD	Flags
	MS	MSD	MS	MSD	Result	Recovery	Recovery	Limits	RPD	Limit		
<b>MATRIX SPIKES</b>												
Laboratory ID:	10-080-03											
	MS	MSD	MS	MSD		MS	MSD					
Phenol	<b>0.914</b>	<b>0.997</b>	1.33	1.33	ND	69	75	31 - 108	9	36		
2-Chlorophenol	<b>0.940</b>	<b>1.03</b>	1.33	1.33	ND	71	77	38 - 103	9	38		
1,4-Dichlorobenzene	<b>0.465</b>	<b>0.502</b>	0.667	0.667	ND	70	75	25 - 101	8	40		
n-Nitroso-di-n-propylamine	<b>0.452</b>	<b>0.494</b>	0.667	0.667	ND	68	74	26 - 102	9	38		
1,2,4-Trichlorobenzene	<b>0.451</b>	<b>0.513</b>	0.667	0.667	ND	68	77	27 - 101	13	40		
4-Chloro-3-methylphenol	<b>0.914</b>	<b>1.02</b>	1.33	1.33	ND	69	77	42 - 106	11	29		
Acenaphthene	<b>0.460</b>	<b>0.513</b>	0.667	0.667	ND	69	77	42 - 103	11	30		
4-Nitrophenol	<b>0.927</b>	<b>1.02</b>	1.33	1.33	ND	70	77	25 - 125	10	29		
2,4-Dinitrotoluene	<b>0.466</b>	<b>0.507</b>	0.667	0.667	ND	70	76	45 - 107	8	30		
Pentachlorophenol	<b>1.03</b>	<b>1.15</b>	1.33	1.33	ND	77	86	30 - 103	11	31		
Pyrene	<b>0.478</b>	<b>0.532</b>	0.667	0.667	ND	72	80	50 - 118	11	28		
<i>Surrogate:</i>												
2-Fluorophenol						69	75	24 - 117				
Phenol-d6						71	77	30 - 120				
Nitrobenzene-d5						72	79	27 - 112				
2-Fluorobiphenyl						71	76	35 - 113				
2,4,6-Tribromophenol						75	81	21 - 120				
Terphenyl-d14						72	79	39 - 121				



Date of Report: October 17, 2016  
 Samples Submitted: October 7, 2016  
 Laboratory Reference: 1610-071  
 Project: 31008

**PCBs**  
**EPA 8082A**

Matrix: Soil  
 Units: mg/Kg (ppm)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>H-1-16-10</b>					
Laboratory ID:	10-071-02					
Aroclor 1016	<b>ND</b>	0.058	EPA 8082A	10-11-16	10-11-16	
Aroclor 1221	<b>ND</b>	0.058	EPA 8082A	10-11-16	10-11-16	
Aroclor 1232	<b>ND</b>	0.058	EPA 8082A	10-11-16	10-11-16	
Aroclor 1242	<b>ND</b>	0.058	EPA 8082A	10-11-16	10-11-16	
Aroclor 1248	<b>ND</b>	0.058	EPA 8082A	10-11-16	10-11-16	
Aroclor 1254	<b>ND</b>	0.058	EPA 8082A	10-11-16	10-11-16	
Aroclor 1260	<b>ND</b>	0.058	EPA 8082A	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>DCB</i>	<i>78</i>	<i>50-139</i>				



Date of Report: October 17, 2016  
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 Project: 31008

**PCBs EPA 8082A  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1011S1					
Aroclor 1016	ND	0.050	EPA 8082A	10-11-16	10-11-16	
Aroclor 1221	ND	0.050	EPA 8082A	10-11-16	10-11-16	
Aroclor 1232	ND	0.050	EPA 8082A	10-11-16	10-11-16	
Aroclor 1242	ND	0.050	EPA 8082A	10-11-16	10-11-16	
Aroclor 1248	ND	0.050	EPA 8082A	10-11-16	10-11-16	
Aroclor 1254	ND	0.050	EPA 8082A	10-11-16	10-11-16	
Aroclor 1260	ND	0.050	EPA 8082A	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
DCB	86		50-139			

Analyte	Result		Spike Level		Source Result	Percent Recovery		Recovery Limits	RPD	RPD Limit	Flags
<b>MATRIX SPIKES</b>											
Laboratory ID:	10-080-01										
	MS	MSD	MS	MSD		MS	MSD				
Aroclor 1260	0.389	0.387	0.500	0.500	ND	78	77	49-133	1	17	
<i>Surrogate:</i>											
DCB						80	81	50-139			



Date of Report: October 17, 2016  
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 Project: 31008

**TOTAL METALS**  
**EPA 6010C/6020A/7471B**

Matrix: Soil  
 Units: mg/kg (ppm)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>EPA Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
Lab ID:	10-071-02					
<b>Client ID:</b>	<b>H-1-16-10</b>					
Antimony	<b>ND</b>	5.8	6010C	10-13-16	10-13-16	
Arsenic	<b>ND</b>	12	6010C	10-13-16	10-13-16	
Beryllium	<b>ND</b>	0.58	6010C	10-13-16	10-13-16	
Cadmium	<b>ND</b>	0.58	6010C	10-13-16	10-13-16	
Chromium	<b>36</b>	0.58	6010C	10-13-16	10-13-16	
Copper	<b>8.4</b>	1.2	6010C	10-13-16	10-13-16	
Lead	<b>ND</b>	5.8	6010C	10-13-16	10-13-16	
Mercury	<b>ND</b>	0.29	7471B	10-11-16	10-11-16	
Nickel	<b>27</b>	2.9	6010C	10-13-16	10-13-16	
Selenium	<b>ND</b>	12	6010C	10-13-16	10-13-16	
Silver	<b>ND</b>	0.58	6010C	10-13-16	10-13-16	
Thallium	<b>ND</b>	1.4	6020A	10-13-16	10-17-16	
Zinc	<b>20</b>	2.9	6010C	10-13-16	10-13-16	



Date of Report: October 17, 2016  
 Samples Submitted: October 7, 2016  
 Laboratory Reference: 1610-071  
 Project: 31008

**TOTAL METALS  
 EPA 6010C/6020A/7471B  
 METHOD BLANK QUALITY CONTROL**

Date Extracted: 10-11&13-16  
 Date Analyzed: 10-11,13&17-16

Matrix: Soil  
 Units: mg/kg (ppm)

Lab ID: MB1013SH1&MB1011S1

Analyte	Method	Result	PQL
Antimony	6010C	ND	5.0
Arsenic	6010C	ND	10
Beryllium	6010C	ND	0.50
Cadmium	6010C	ND	0.50
Chromium	6010C	ND	0.50
Copper	6010C	ND	1.0
Lead	6010C	ND	5.0
Mercury	7471B	ND	0.25
Nickel	6010C	ND	2.5
Selenium	6010C	ND	10
Silver	6010C	ND	0.50
Thallium	6020A	ND	1.3
Zinc	6010C	ND	2.5



Date of Report: October 17, 2016  
 Samples Submitted: October 7, 2016  
 Laboratory Reference: 1610-071  
 Project: 31008

**TOTAL METALS  
 EPA 6010C/6020A/7471B  
 DUPLICATE QUALITY CONTROL**

Date Extracted: 10-11&13-16  
 Date Analyzed: 10-11,13&17-16

Matrix: Soil  
 Units: mg/kg (ppm)

Lab ID: 10-080-01

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Antimony	ND	ND	NA	5.0	
Arsenic	ND	ND	NA	10.0	
Beryllium	ND	ND	NA	0.50	
Cadmium	ND	ND	NA	0.50	
Chromium	25.9	24.3	6	0.50	
Copper	11.7	10.9	7	1.0	
Lead	ND	ND	NA	5.0	
Mercury	ND	ND	NA	0.25	
Nickel	30.4	29.2	4	2.5	
Selenium	ND	ND	NA	10	
Silver	ND	ND	NA	0.50	
Thallium	ND	ND	NA	1.3	
Zinc	24.0	22.0	9	2.5	



Date of Report: October 17, 2016  
 Samples Submitted: October 7, 2016  
 Laboratory Reference: 1610-071  
 Project: 31008

**TOTAL METALS  
 EPA 6010C/6020A/7471B  
 MS/MSD QUALITY CONTROL**

Date Extracted: 10-11&13-16  
 Date Analyzed: 10-11,13&17-16

Matrix: Soil  
 Units: mg/kg (ppm)

Lab ID: 10-080-01

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Antimony	100	<b>95.4</b>	95	<b>88.3</b>	88	8	
Arsenic	100	<b>101</b>	101	<b>94.3</b>	94	6	
Beryllium	50.0	<b>50.8</b>	102	<b>47.6</b>	95	7	
Cadmium	50.0	<b>49.3</b>	99	<b>47.5</b>	95	4	
Chromium	100	<b>128</b>	102	<b>119</b>	93	7	
Copper	50.0	<b>63.9</b>	105	<b>60.4</b>	97	6	
Lead	250	<b>238</b>	95	<b>232</b>	93	3	
Mercury	0.500	<b>0.499</b>	100	<b>0.546</b>	109	9	
Nickel	100	<b>127</b>	96	<b>120</b>	89	6	
Selenium	100	<b>104</b>	104	<b>99.4</b>	99	5	
Silver	25.0	<b>24.4</b>	97	<b>23.1</b>	92	5	
Thallium	50.0	<b>44.2</b>	88	<b>44.8</b>	90	1	
Zinc	100	<b>120</b>	96	<b>115</b>	91	4	



Date of Report: October 17, 2016  
Samples Submitted: October 7, 2016  
Laboratory Reference: 1610-071  
Project: 31008

**% MOISTURE**

Date Analyzed: 10-7-16

Client ID	Lab ID	% Moisture
H-1-16-10	10-071-02	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.
  - Z -
- ND - Not Detected at PQL  
 PQL - Practical Quantitation Limit  
 RPD - Relative Percent Difference





**MVA Onsite Environmental Inc.**  
 Analytical Laboratory Testing Services  
 14648 NE 95th Street • Redmond, WA 98052  
 Phone: (425) 883-3881 • www.onsite-env.com

# Chain of Custody

Turnaround Request  
 (in working days)  
 (Check One)

- Same Day  1 Day  
 2 Days  3 Days

Standard (7 Days)  
 (1 PH anal/sis 5 Days)

\_\_\_\_\_ (other)

Laboratory Number:

**10-071**

Company: **INVAEX**  
 Project Number: **31008**  
 Project Name: **SR-520 Phase II Investigation**  
 Project Manager: **Glenn Hayman**  
 Sampled by: **Jennifer Heise**

Lab ID	Sample Identification	Date Sampled	Time Sampled	Matrix	Number of Containers
1	H-1-16-5	161006	2152	soil	5
2	H-1-16-10	161006	2158	soil	5
3	H-1-16-15	161006	2204	soil	5
4	H-1-16-20	161006	2223	soil	5
5	H-1-16-25	161006	2236	soil	5
6	H-1-16-30	161006	2248	soil	5
7	H-1-16-45	161006	2330	soil	3
8	H-1-16-50	161006	2350	soil	3

Number of Containers	Analysis	Result
5	NWTPH-HCID	
5	NWTPH-Gx/BTEX	
5	NWTPH-Gx (Hold)	
5	NWTPH-Dx ( <input type="checkbox"/> Acid / SG Clean-up ) (Hold)	
5	Volatiles 8260C	
5	Halogenated Volatiles 8260C	
5	EDB EPA 8011 (Waters Only)	
5	Semivolatiles 8270D/SIM (with low-level PAHs)	
5	PAHs 8270D/SIM (low-level)	
5	PCBs 8082A	
5	Organochlorine Pesticides 8081B	
5	Organophosphorus Pesticides 8270D/SIM	
5	Chlorinated Acid Herbicides 8151A	
5	Total RCRA Metals	
5	Total MTCA Metals	
5	TCLP Metals	
5	HEM (oil and grease) 1664A	
5	PP13* - 6010	
5	% Moisture	

Signature	Company	Date	Time	Comments/Special Instructions
	INVAEX	10/7/16	1:05 pm	* Priority pollutant metals Hold all other samples until further notice
	OSE	10/7/16	1:05 pm	

Received \_\_\_\_\_  
 Relinquished \_\_\_\_\_  
 Received \_\_\_\_\_  
 Relinquished \_\_\_\_\_  
 Reviewed/Date \_\_\_\_\_

Data Package: Standard  Level III  Level IV   
 Chromatograms with final report  Electronic Data Deliverables (EDDs)



14648 NE 95<sup>th</sup> Street, Redmond, WA 98052 • (425) 883-3881

October 17, 2016

Glenn Hayman  
INNOVEX Environmental Mgt., Inc.  
16310 NE 80th St., Suite 300  
Redmond, WA 98052

Re: Analytical Data for Project 31008  
Laboratory Reference No. 1610-079

Dear Glenn:

Enclosed are the analytical results and associated quality control data for samples submitted on October 8, 2016.

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,

A handwritten signature in black ink, appearing to read "DB", with a long horizontal stroke extending to the right.

David Baumeister  
Project Manager

Enclosures



---

OnSite Environmental, Inc. 14648 NE 95<sup>th</sup> Street, Redmond, WA 98052 (425) 883-3881

This report pertains to the samples analyzed in accordance with the chain of custody, and is intended only for the use of the individual or company to whom it is addressed.

Date of Report: October 17, 2016  
Samples Submitted: October 8, 2016  
Laboratory Reference: 1610-079  
Project: 31008

### Case Narrative

Samples were collected on October 7, 2016 and received by the laboratory on October 8, 2016. 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.

#### Volatiles EPA 8260C 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.

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.



Date of Report: October 17, 2016  
 Samples Submitted: October 8, 2016  
 Laboratory Reference: 1610-079  
 Project: 31008

### NWTPH-HCID

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-2-16-13.5</b>					
Laboratory ID:	10-079-04					
Gasoline Range Organics	<b>ND</b>	29	NWTPH-HCID	10-10-16	10-10-16	
Diesel Range Organics	<b>ND</b>	71	NWTPH-HCID	10-10-16	10-10-16	
Lube Oil Range Organics	<b>ND</b>	140	NWTPH-HCID	10-10-16	10-10-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>94</i>	<i>50-150</i>				



Date of Report: October 17, 2016  
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**NWTPH-HCID  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1010S2					
Gasoline Range Organics	<b>ND</b>	20	NWTPH-HCID	10-10-16	10-10-16	
Diesel Range Organics	<b>ND</b>	50	NWTPH-HCID	10-10-16	10-10-16	
Lube Oil Range Organics	<b>ND</b>	100	NWTPH-HCID	10-10-16	10-10-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>84</i>	<i>50-150</i>				



Date of Report: October 17, 2016  
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 Project: 31008

**VOLATILES EPA 8260C**  
 page 1 of 2

Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>H-2-16-13.5</b>					
<b>Laboratory ID:</b>	<b>10-079-04</b>					
Dichlorodifluoromethane	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
Chloromethane	ND	0.0054	EPA 8260C	10-10-16	10-10-16	
Vinyl Chloride	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
Bromomethane	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
Chloroethane	ND	0.0054	EPA 8260C	10-10-16	10-10-16	
Trichlorofluoromethane	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
1,1-Dichloroethene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
Acetone	0.060	0.0054	EPA 8260C	10-10-16	10-10-16	
Iodomethane	ND	0.0054	EPA 8260C	10-10-16	10-10-16	
Carbon Disulfide	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
Methylene Chloride	ND	0.0054	EPA 8260C	10-10-16	10-10-16	
(trans) 1,2-Dichloroethene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
Methyl t-Butyl Ether	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
1,1-Dichloroethane	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
Vinyl Acetate	ND	0.0054	EPA 8260C	10-10-16	10-10-16	
2,2-Dichloropropane	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
(cis) 1,2-Dichloroethene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
2-Butanone	0.021	0.0054	EPA 8260C	10-10-16	10-10-16	
Bromochloromethane	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
Chloroform	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
1,1,1-Trichloroethane	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
Carbon Tetrachloride	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
1,1-Dichloropropene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
Benzene	0.0053	0.0011	EPA 8260C	10-10-16	10-10-16	
1,2-Dichloroethane	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
Trichloroethene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
1,2-Dichloropropane	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
Dibromomethane	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
Bromodichloromethane	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
2-Chloroethyl Vinyl Ether	ND	0.0054	EPA 8260C	10-10-16	10-10-16	
(cis) 1,3-Dichloropropene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
Methyl Isobutyl Ketone	ND	0.0054	EPA 8260C	10-10-16	10-10-16	
Toluene	ND	0.0054	EPA 8260C	10-10-16	10-10-16	
(trans) 1,3-Dichloropropene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	



Date of Report: October 17, 2016  
 Samples Submitted: October 8, 2016  
 Laboratory Reference: 1610-079  
 Project: 31008

**VOLATILES EPA 8260C**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-2-16-13.5</b>					
Laboratory ID:	10-079-04					
1,1,2-Trichloroethane	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
Tetrachloroethene	ND	0.0022	EPA 8260C	10-10-16	10-10-16	
1,3-Dichloropropane	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
2-Hexanone	ND	0.0054	EPA 8260C	10-10-16	10-10-16	
Dibromochloromethane	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
1,2-Dibromoethane	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
Chlorobenzene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
1,1,1,2-Tetrachloroethane	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
Ethylbenzene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
m,p-Xylene	ND	0.0022	EPA 8260C	10-10-16	10-10-16	
o-Xylene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
Styrene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
Bromoform	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
Isopropylbenzene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
Bromobenzene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
1,1,2,2-Tetrachloroethane	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
1,2,3-Trichloropropane	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
n-Propylbenzene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
2-Chlorotoluene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
4-Chlorotoluene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
1,3,5-Trimethylbenzene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
tert-Butylbenzene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
1,2,4-Trimethylbenzene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
sec-Butylbenzene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
1,3-Dichlorobenzene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
p-Isopropyltoluene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
1,4-Dichlorobenzene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
1,2-Dichlorobenzene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
n-Butylbenzene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
1,2-Dibromo-3-chloropropane	ND	0.0054	EPA 8260C	10-10-16	10-10-16	
1,2,4-Trichlorobenzene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
Hexachlorobutadiene	ND	0.0054	EPA 8260C	10-10-16	10-10-16	
Naphthalene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
1,2,3-Trichlorobenzene	ND	0.0011	EPA 8260C	10-10-16	10-10-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>97</i>	<i>76-131</i>				
<i>Toluene-d8</i>	<i>98</i>	<i>80-126</i>				
<i>4-Bromofluorobenzene</i>	<i>87</i>	<i>60-146</i>				



Date of Report: October 17, 2016  
 Samples Submitted: October 8, 2016  
 Laboratory Reference: 1610-079  
 Project: 31008

**VOLATILES by EPA 8260C**  
**METHOD BLANK QUALITY CONTROL**  
 page 1 of 2

Matrix: Soil  
 Units: mg/kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1010S1					
Dichlorodifluoromethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Chloromethane	ND	0.0050	EPA 8260C	10-10-16	10-10-16	
Vinyl Chloride	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Bromomethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Chloroethane	ND	0.0050	EPA 8260C	10-10-16	10-10-16	
Trichlorofluoromethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,1-Dichloroethene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Acetone	ND	0.0050	EPA 8260C	10-10-16	10-10-16	
Iodomethane	ND	0.0050	EPA 8260C	10-10-16	10-10-16	
Carbon Disulfide	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Methylene Chloride	ND	0.0050	EPA 8260C	10-10-16	10-10-16	
(trans) 1,2-Dichloroethene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Methyl t-Butyl Ether	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,1-Dichloroethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Vinyl Acetate	ND	0.0050	EPA 8260C	10-10-16	10-10-16	
2,2-Dichloropropane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
(cis) 1,2-Dichloroethene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
2-Butanone	ND	0.0050	EPA 8260C	10-10-16	10-10-16	
Bromochloromethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Chloroform	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,1,1-Trichloroethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Carbon Tetrachloride	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,1-Dichloropropene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Benzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,2-Dichloroethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Trichloroethene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,2-Dichloropropane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Dibromomethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Bromodichloromethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
2-Chloroethyl Vinyl Ether	ND	0.0050	EPA 8260C	10-10-16	10-10-16	
(cis) 1,3-Dichloropropene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Methyl Isobutyl Ketone	ND	0.0050	EPA 8260C	10-10-16	10-10-16	
Toluene	ND	0.0050	EPA 8260C	10-10-16	10-10-16	
(trans) 1,3-Dichloropropene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	



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 Laboratory Reference: 1610-079  
 Project: 31008

**VOLATILES by EPA 8260C**  
**METHOD BLANK QUALITY CONTROL**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1010S1					
1,1,2-Trichloroethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Tetrachloroethene	ND	0.0020	EPA 8260C	10-10-16	10-10-16	
1,3-Dichloropropane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
2-Hexanone	ND	0.0050	EPA 8260C	10-10-16	10-10-16	
Dibromochloromethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,2-Dibromoethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Chlorobenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,1,1,2-Tetrachloroethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Ethylbenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
m,p-Xylene	ND	0.0020	EPA 8260C	10-10-16	10-10-16	
o-Xylene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Styrene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Bromoform	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Isopropylbenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Bromobenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,1,2,2-Tetrachloroethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,2,3-Trichloropropane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
n-Propylbenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
2-Chlorotoluene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
4-Chlorotoluene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,3,5-Trimethylbenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
tert-Butylbenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,2,4-Trimethylbenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
sec-Butylbenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,3-Dichlorobenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
p-Isopropyltoluene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,4-Dichlorobenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,2-Dichlorobenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
n-Butylbenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,2-Dibromo-3-chloropropane	ND	0.0050	EPA 8260C	10-10-16	10-10-16	
1,2,4-Trichlorobenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Hexachlorobutadiene	ND	0.0050	EPA 8260C	10-10-16	10-10-16	
Naphthalene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,2,3-Trichlorobenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>98</i>	<i>76-131</i>				
<i>Toluene-d8</i>	<i>103</i>	<i>80-126</i>				
<i>4-Bromofluorobenzene</i>	<i>102</i>	<i>60-146</i>				



Date of Report: October 17, 2016  
 Samples Submitted: October 8, 2016  
 Laboratory Reference: 1610-079  
 Project: 31008

**VOLATILES by EPA 8260C  
 SB/SBD QUALITY CONTROL**

Matrix: Soil  
 Units: mg/kg

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD		Flags
					Recovery	Limits	RPD	Limit		
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB1010S1									
	SB	SBD	SB	SBD	SB	SBD				
1,1-Dichloroethene	<b>0.0457</b>	<b>0.0490</b>	0.0500	0.0500	91	98	68-126	7	15	
Benzene	<b>0.0461</b>	<b>0.0480</b>	0.0500	0.0500	92	96	70-121	4	15	
Trichloroethene	<b>0.0455</b>	<b>0.0476</b>	0.0500	0.0500	91	95	75-120	5	15	
Toluene	<b>0.0468</b>	<b>0.0495</b>	0.0500	0.0500	94	99	80-120	6	15	
Chlorobenzene	<b>0.0475</b>	<b>0.0490</b>	0.0500	0.0500	95	98	76-120	3	15	
<i>Surrogate:</i>										
Dibromofluoromethane					94	94	76-131			
Toluene-d8					94	100	80-126			
4-Bromofluorobenzene					96	96	60-146			



Date of Report: October 17, 2016  
 Samples Submitted: October 8, 2016  
 Laboratory Reference: 1610-079  
 Project: 31008

**SEMIVOLATILES EPA 8270D/SIM**  
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Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-2-16-13.5</b>					
<b>Laboratory ID:</b>	<b>10-079-04</b>					
n-Nitrosodimethylamine	ND	0.048	EPA 8270D	10-10-16	10-11-16	
Pyridine	ND	0.48	EPA 8270D	10-10-16	10-11-16	
Phenol	ND	0.048	EPA 8270D	10-10-16	10-11-16	
Aniline	ND	0.24	EPA 8270D	10-10-16	10-11-16	
bis(2-Chloroethyl)ether	ND	0.048	EPA 8270D	10-10-16	10-11-16	
2-Chlorophenol	ND	0.048	EPA 8270D	10-10-16	10-11-16	
1,3-Dichlorobenzene	ND	0.048	EPA 8270D	10-10-16	10-11-16	
1,4-Dichlorobenzene	ND	0.048	EPA 8270D	10-10-16	10-11-16	
Benzyl alcohol	ND	0.24	EPA 8270D	10-10-16	10-11-16	
1,2-Dichlorobenzene	ND	0.048	EPA 8270D	10-10-16	10-11-16	
2-Methylphenol (o-Cresol)	ND	0.048	EPA 8270D	10-10-16	10-11-16	
bis(2-Chloroisopropyl)ether	ND	0.048	EPA 8270D	10-10-16	10-11-16	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.048	EPA 8270D	10-10-16	10-11-16	
n-Nitroso-di-n-propylamine	ND	0.048	EPA 8270D	10-10-16	10-11-16	
Hexachloroethane	ND	0.048	EPA 8270D	10-10-16	10-11-16	
Nitrobenzene	ND	0.048	EPA 8270D	10-10-16	10-11-16	
Isophorone	ND	0.048	EPA 8270D	10-10-16	10-11-16	
2-Nitrophenol	ND	0.048	EPA 8270D	10-10-16	10-11-16	
2,4-Dimethylphenol	ND	0.048	EPA 8270D	10-10-16	10-11-16	
bis(2-Chloroethoxy)methane	ND	0.048	EPA 8270D	10-10-16	10-11-16	
2,4-Dichlorophenol	ND	0.048	EPA 8270D	10-10-16	10-11-16	
1,2,4-Trichlorobenzene	ND	0.048	EPA 8270D	10-10-16	10-11-16	
Naphthalene	0.0096	0.0095	EPA 8270D/SIM	10-10-16	10-10-16	
4-Chloroaniline	ND	0.24	EPA 8270D	10-10-16	10-11-16	
Hexachlorobutadiene	ND	0.048	EPA 8270D	10-10-16	10-11-16	
4-Chloro-3-methylphenol	ND	0.048	EPA 8270D	10-10-16	10-11-16	
2-Methylnaphthalene	ND	0.0095	EPA 8270D/SIM	10-10-16	10-10-16	
1-Methylnaphthalene	ND	0.0095	EPA 8270D/SIM	10-10-16	10-10-16	
Hexachlorocyclopentadiene	ND	0.048	EPA 8270D	10-10-16	10-11-16	
2,4,6-Trichlorophenol	ND	0.048	EPA 8270D	10-10-16	10-11-16	
2,3-Dichloroaniline	ND	0.048	EPA 8270D	10-10-16	10-11-16	
2,4,5-Trichlorophenol	ND	0.048	EPA 8270D	10-10-16	10-11-16	
2-Chloronaphthalene	ND	0.048	EPA 8270D	10-10-16	10-11-16	
2-Nitroaniline	ND	0.048	EPA 8270D	10-10-16	10-11-16	
1,4-Dinitrobenzene	ND	0.048	EPA 8270D	10-10-16	10-11-16	
Dimethylphthalate	ND	0.048	EPA 8270D	10-10-16	10-11-16	
1,3-Dinitrobenzene	ND	0.048	EPA 8270D	10-10-16	10-11-16	
2,6-Dinitrotoluene	ND	0.048	EPA 8270D	10-10-16	10-11-16	
1,2-Dinitrobenzene	ND	0.048	EPA 8270D	10-10-16	10-11-16	
Acenaphthylene	ND	0.0095	EPA 8270D/SIM	10-10-16	10-10-16	
3-Nitroaniline	ND	0.048	EPA 8270D	10-10-16	10-11-16	



Date of Report: October 17, 2016  
 Samples Submitted: October 8, 2016  
 Laboratory Reference: 1610-079  
 Project: 31008

**SEMIVOLATILES EPA 8270D/SIM**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-2-16-13.5</b>					
Laboratory ID:	10-079-04					
2,4-Dinitrophenol	ND	0.24	EPA 8270D	10-10-16	10-11-16	
Acenaphthene	ND	0.0095	EPA 8270D/SIM	10-10-16	10-10-16	
4-Nitrophenol	ND	0.048	EPA 8270D	10-10-16	10-11-16	
2,4-Dinitrotoluene	ND	0.048	EPA 8270D	10-10-16	10-11-16	
Dibenzofuran	ND	0.048	EPA 8270D	10-10-16	10-11-16	
2,3,5,6-Tetrachlorophenol	ND	0.048	EPA 8270D	10-10-16	10-11-16	
2,3,4,6-Tetrachlorophenol	ND	0.048	EPA 8270D	10-10-16	10-11-16	
Diethylphthalate	ND	0.24	EPA 8270D	10-10-16	10-11-16	
4-Chlorophenyl-phenylether	ND	0.048	EPA 8270D	10-10-16	10-11-16	
4-Nitroaniline	ND	0.048	EPA 8270D	10-10-16	10-11-16	
Fluorene	ND	0.0095	EPA 8270D/SIM	10-10-16	10-10-16	
4,6-Dinitro-2-methylphenol	ND	0.24	EPA 8270D	10-10-16	10-11-16	
n-Nitrosodiphenylamine	ND	0.048	EPA 8270D	10-10-16	10-11-16	
1,2-Diphenylhydrazine	ND	0.048	EPA 8270D	10-10-16	10-11-16	
4-Bromophenyl-phenylether	ND	0.048	EPA 8270D	10-10-16	10-11-16	
Hexachlorobenzene	ND	0.048	EPA 8270D	10-10-16	10-11-16	
Pentachlorophenol	ND	0.24	EPA 8270D	10-10-16	10-11-16	
Phenanthrene	ND	0.0095	EPA 8270D/SIM	10-10-16	10-10-16	
Anthracene	ND	0.0095	EPA 8270D/SIM	10-10-16	10-10-16	
Carbazole	ND	0.048	EPA 8270D	10-10-16	10-11-16	
Di-n-butylphthalate	ND	0.24	EPA 8270D	10-10-16	10-11-16	
Fluoranthene	ND	0.0095	EPA 8270D/SIM	10-10-16	10-10-16	
Benzidine	ND	0.48	EPA 8270D	10-10-16	10-11-16	
Pyrene	ND	0.0095	EPA 8270D/SIM	10-10-16	10-10-16	
Butylbenzylphthalate	ND	0.048	EPA 8270D	10-10-16	10-11-16	
bis-2-Ethylhexyladipate	ND	0.048	EPA 8270D	10-10-16	10-11-16	
3,3'-Dichlorobenzidine	ND	0.24	EPA 8270D	10-10-16	10-11-16	
Benzo[a]anthracene	ND	0.0095	EPA 8270D/SIM	10-10-16	10-10-16	
Chrysene	ND	0.0095	EPA 8270D/SIM	10-10-16	10-10-16	
bis(2-Ethylhexyl)phthalate	ND	0.048	EPA 8270D	10-10-16	10-11-16	
Di-n-octylphthalate	ND	0.048	EPA 8270D	10-10-16	10-11-16	
Benzo[b]fluoranthene	ND	0.0095	EPA 8270D/SIM	10-10-16	10-10-16	
Benzo(j,k)fluoranthene	ND	0.0095	EPA 8270D/SIM	10-10-16	10-10-16	
Benzo[a]pyrene	ND	0.0095	EPA 8270D/SIM	10-10-16	10-10-16	
Indeno[1,2,3-cd]pyrene	ND	0.0095	EPA 8270D/SIM	10-10-16	10-10-16	
Dibenz[a,h]anthracene	ND	0.0095	EPA 8270D/SIM	10-10-16	10-10-16	
Benzo[g,h,i]perylene	ND	0.0095	EPA 8270D/SIM	10-10-16	10-10-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	64	24 - 117				
Phenol-d6	66	30 - 120				
Nitrobenzene-d5	69	27 - 112				
2-Fluorobiphenyl	70	35 - 113				
2,4,6-Tribromophenol	75	21 - 120				
Terphenyl-d14	70	39 - 121				



Date of Report: October 17, 2016  
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 Project: 31008

**SEMIVOLATILES EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
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Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1010S1					
n-Nitrosodimethylamine	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Pyridine	ND	0.33	EPA 8270D	10-10-16	10-11-16	
Phenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Aniline	ND	0.17	EPA 8270D	10-10-16	10-11-16	
bis(2-Chloroethyl)ether	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2-Chlorophenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
1,3-Dichlorobenzene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
1,4-Dichlorobenzene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Benzyl alcohol	ND	0.17	EPA 8270D	10-10-16	10-11-16	
1,2-Dichlorobenzene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2-Methylphenol (o-Cresol)	ND	0.033	EPA 8270D	10-10-16	10-11-16	
bis(2-Chloroisopropyl)ether	ND	0.033	EPA 8270D	10-10-16	10-11-16	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.033	EPA 8270D	10-10-16	10-11-16	
n-Nitroso-di-n-propylamine	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Hexachloroethane	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Nitrobenzene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Isophorone	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2-Nitrophenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2,4-Dimethylphenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
bis(2-Chloroethoxy)methane	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2,4-Dichlorophenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
1,2,4-Trichlorobenzene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Naphthalene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
4-Chloroaniline	ND	0.17	EPA 8270D	10-10-16	10-11-16	
Hexachlorobutadiene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
4-Chloro-3-methylphenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2-Methylnaphthalene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
1-Methylnaphthalene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Hexachlorocyclopentadiene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2,4,6-Trichlorophenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2,3-Dichloroaniline	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2,4,5-Trichlorophenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2-Chloronaphthalene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2-Nitroaniline	ND	0.033	EPA 8270D	10-10-16	10-11-16	
1,4-Dinitrobenzene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Dimethylphthalate	ND	0.033	EPA 8270D	10-10-16	10-11-16	
1,3-Dinitrobenzene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2,6-Dinitrotoluene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
1,2-Dinitrobenzene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Acenaphthylene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
3-Nitroaniline	ND	0.033	EPA 8270D	10-10-16	10-11-16	



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**SEMIVOLATILES EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
 page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1010S1					
2,4-Dinitrophenol	ND	0.17	EPA 8270D	10-10-16	10-11-16	
Acenaphthene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
4-Nitrophenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2,4-Dinitrotoluene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Dibenzofuran	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2,3,5,6-Tetrachlorophenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2,3,4,6-Tetrachlorophenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Diethylphthalate	ND	0.17	EPA 8270D	10-10-16	10-11-16	
4-Chlorophenyl-phenylether	ND	0.033	EPA 8270D	10-10-16	10-11-16	
4-Nitroaniline	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Fluorene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
4,6-Dinitro-2-methylphenol	ND	0.17	EPA 8270D	10-10-16	10-11-16	
n-Nitrosodiphenylamine	ND	0.033	EPA 8270D	10-10-16	10-11-16	
1,2-Diphenylhydrazine	ND	0.033	EPA 8270D	10-10-16	10-11-16	
4-Bromophenyl-phenylether	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Hexachlorobenzene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Pentachlorophenol	ND	0.17	EPA 8270D	10-10-16	10-11-16	
Phenanthrene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Anthracene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Carbazole	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Di-n-butylphthalate	ND	0.17	EPA 8270D	10-10-16	10-11-16	
Fluoranthene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Benzidine	ND	0.33	EPA 8270D	10-10-16	10-11-16	
Pyrene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Butylbenzylphthalate	ND	0.033	EPA 8270D	10-10-16	10-11-16	
bis-2-Ethylhexyladipate	ND	0.033	EPA 8270D	10-10-16	10-11-16	
3,3'-Dichlorobenzidine	ND	0.17	EPA 8270D	10-10-16	10-11-16	
Benzo[a]anthracene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Chrysene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
bis(2-Ethylhexyl)phthalate	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Di-n-octylphthalate	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Benzo[a]pyrene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Indeno[1,2,3-cd]pyrene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Benzo[g,h,i]perylene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	81	24 - 117				
Phenol-d6	83	30 - 120				
Nitrobenzene-d5	85	27 - 112				
2-Fluorobiphenyl	84	35 - 113				
2,4,6-Tribromophenol	85	21 - 120				
Terphenyl-d14	85	39 - 121				



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**SEMIVOLATILES EPA 8270D/SIM  
 MS/MSD QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result		Spike Level		Source	Percent	Recovery	RPD		Flags
					Result	Recovery	Limits			
<b>MATRIX SPIKES</b>										
Laboratory ID:	10-080-03									
	MS	MSD	MS	MSD		MS	MSD			
Phenol	<b>0.914</b>	<b>0.997</b>	1.33	1.33	ND	69	75	31 - 108	9	36
2-Chlorophenol	<b>0.940</b>	<b>1.03</b>	1.33	1.33	ND	71	77	38 - 103	9	38
1,4-Dichlorobenzene	<b>0.465</b>	<b>0.502</b>	0.667	0.667	ND	70	75	25 - 101	8	40
n-Nitroso-di-n-propylamine	<b>0.452</b>	<b>0.494</b>	0.667	0.667	ND	68	74	26 - 102	9	38
1,2,4-Trichlorobenzene	<b>0.451</b>	<b>0.513</b>	0.667	0.667	ND	68	77	27 - 101	13	40
4-Chloro-3-methylphenol	<b>0.914</b>	<b>1.02</b>	1.33	1.33	ND	69	77	42 - 106	11	29
Acenaphthene	<b>0.460</b>	<b>0.513</b>	0.667	0.667	ND	69	77	42 - 103	11	30
4-Nitrophenol	<b>0.927</b>	<b>1.02</b>	1.33	1.33	ND	70	77	25 - 125	10	29
2,4-Dinitrotoluene	<b>0.466</b>	<b>0.507</b>	0.667	0.667	ND	70	76	45 - 107	8	30
Pentachlorophenol	<b>1.03</b>	<b>1.15</b>	1.33	1.33	ND	77	86	30 - 103	11	31
Pyrene	<b>0.478</b>	<b>0.532</b>	0.667	0.667	ND	72	80	50 - 118	11	28
<i>Surrogate:</i>										
<i>2-Fluorophenol</i>						69	75	24 - 117		
<i>Phenol-d6</i>						71	77	30 - 120		
<i>Nitrobenzene-d5</i>						72	79	27 - 112		
<i>2-Fluorobiphenyl</i>						71	76	35 - 113		
<i>2,4,6-Tribromophenol</i>						75	81	21 - 120		
<i>Terphenyl-d14</i>						72	79	39 - 121		



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**PCBs**  
**EPA 8082A**

Matrix: Soil  
 Units: mg/Kg (ppm)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>H-2-16-13.5</b>					
Laboratory ID:	10-079-04					
Aroclor 1016	<b>ND</b>	0.071	EPA 8082A	10-11-16	10-11-16	
Aroclor 1221	<b>ND</b>	0.071	EPA 8082A	10-11-16	10-11-16	
Aroclor 1232	<b>ND</b>	0.071	EPA 8082A	10-11-16	10-11-16	
Aroclor 1242	<b>ND</b>	0.071	EPA 8082A	10-11-16	10-11-16	
Aroclor 1248	<b>ND</b>	0.071	EPA 8082A	10-11-16	10-11-16	
Aroclor 1254	<b>ND</b>	0.071	EPA 8082A	10-11-16	10-11-16	
Aroclor 1260	<b>ND</b>	0.071	EPA 8082A	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>DCB</i>	<i>73</i>	<i>50-139</i>				



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**PCBs EPA 8082A  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1011S1					
Aroclor 1016	<b>ND</b>	0.050	EPA 8082A	10-11-16	10-11-16	
Aroclor 1221	<b>ND</b>	0.050	EPA 8082A	10-11-16	10-11-16	
Aroclor 1232	<b>ND</b>	0.050	EPA 8082A	10-11-16	10-11-16	
Aroclor 1242	<b>ND</b>	0.050	EPA 8082A	10-11-16	10-11-16	
Aroclor 1248	<b>ND</b>	0.050	EPA 8082A	10-11-16	10-11-16	
Aroclor 1254	<b>ND</b>	0.050	EPA 8082A	10-11-16	10-11-16	
Aroclor 1260	<b>ND</b>	0.050	EPA 8082A	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
<i>DCB</i>	86		50-139			

Analyte	Result		Spike Level		Source Result	Percent Recovery		Recovery Limits	RPD	RPD Limit	Flags
<b>MATRIX SPIKES</b>											
Laboratory ID:	10-080-01										
	MS	MSD	MS	MSD		MS	MSD				
Aroclor 1260	<b>0.389</b>	<b>0.387</b>	0.500	0.500	ND	<b>78</b>	<b>77</b>	49-133	1	17	
<i>Surrogate:</i>											
<i>DCB</i>						80	81	50-139			



Date of Report: October 17, 2016  
 Samples Submitted: October 8, 2016  
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**TOTAL METALS**  
**EPA 6010C/6020A/7471B**

Matrix: Soil  
 Units: mg/kg (ppm)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>EPA Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
Lab ID:	10-079-04					
<b>Client ID:</b>	<b>H-2-16-13.5</b>					
Antimony	<b>ND</b>	7.1	6010C	10-13-16	10-13-16	
Arsenic	<b>ND</b>	14	6010C	10-13-16	10-13-16	
Beryllium	<b>ND</b>	0.71	6010C	10-13-16	10-13-16	
Cadmium	<b>ND</b>	0.71	6010C	10-13-16	10-13-16	
Chromium	<b>37</b>	0.71	6010C	10-13-16	10-13-16	
Copper	<b>24</b>	1.4	6010C	10-13-16	10-13-16	
Lead	<b>11</b>	7.1	6010C	10-13-16	10-13-16	
Mercury	<b>ND</b>	0.36	7471B	10-11-16	10-11-16	
Nickel	<b>36</b>	3.6	6010C	10-13-16	10-13-16	
Selenium	<b>ND</b>	14	6010C	10-13-16	10-13-16	
Silver	<b>ND</b>	0.71	6010C	10-13-16	10-13-16	
Thallium	<b>ND</b>	1.8	6020A	10-13-16	10-17-16	
Zinc	<b>56</b>	3.6	6010C	10-13-16	10-13-16	



Date of Report: October 17, 2016  
 Samples Submitted: October 8, 2016  
 Laboratory Reference: 1610-079  
 Project: 31008

**TOTAL METALS  
 EPA 6010C/6020A/7471B  
 METHOD BLANK QUALITY CONTROL**

Date Extracted: 10-11&13-16  
 Date Analyzed: 10-11,13&17-16

Matrix: Soil  
 Units: mg/kg (ppm)

Lab ID: MB1013SH1&MB1011S1

Analyte	Method	Result	PQL
Antimony	6010C	ND	5.0
Arsenic	6010C	ND	10
Beryllium	6010C	ND	0.50
Cadmium	6010C	ND	0.50
Chromium	6010C	ND	0.50
Copper	6010C	ND	1.0
Lead	6010C	ND	5.0
Mercury	7471B	ND	0.25
Nickel	6010C	ND	2.5
Selenium	6010C	ND	10
Silver	6010C	ND	0.50
Thallium	6020A	ND	1.3
Zinc	6010C	ND	2.5



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**TOTAL METALS  
 EPA 6010C/6020A/7471B  
 DUPLICATE QUALITY CONTROL**

Date Extracted: 10-11&13-16  
 Date Analyzed: 10-11,13&17-16

Matrix: Soil  
 Units: mg/kg (ppm)

Lab ID: 10-080-01

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Antimony	ND	ND	NA	5.0	
Arsenic	ND	ND	NA	10.0	
Beryllium	ND	ND	NA	0.50	
Cadmium	ND	ND	NA	0.50	
Chromium	25.9	24.3	6	0.50	
Copper	11.7	10.9	7	1.0	
Lead	ND	ND	NA	5.0	
Mercury	ND	ND	NA	0.25	
Nickel	30.4	29.2	4	2.5	
Selenium	ND	ND	NA	10	
Silver	ND	ND	NA	0.50	
Thallium	ND	ND	NA	1.3	
Zinc	24.0	22.0	9	2.5	



Date of Report: October 17, 2016  
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 Project: 31008

**TOTAL METALS  
 EPA 6010C/6020A/7471B  
 MS/MSD QUALITY CONTROL**

Date Extracted: 10-11&13-16  
 Date Analyzed: 10-11,13&17-16

Matrix: Soil  
 Units: mg/kg (ppm)

Lab ID: 10-080-01

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Antimony	100	<b>95.4</b>	95	<b>88.3</b>	88	8	
Arsenic	100	<b>101</b>	101	<b>94.3</b>	94	6	
Beryllium	50.0	<b>50.8</b>	102	<b>47.6</b>	95	7	
Cadmium	50.0	<b>49.3</b>	99	<b>47.5</b>	95	4	
Chromium	100	<b>128</b>	102	<b>119</b>	93	7	
Copper	50.0	<b>63.9</b>	105	<b>60.4</b>	97	6	
Lead	250	<b>238</b>	95	<b>232</b>	93	3	
Mercury	0.500	<b>0.499</b>	100	<b>0.546</b>	109	9	
Nickel	100	<b>127</b>	96	<b>120</b>	89	6	
Selenium	100	<b>104</b>	104	<b>99.4</b>	99	5	
Silver	25.0	<b>24.4</b>	97	<b>23.1</b>	92	5	
Thallium	50.0	<b>44.2</b>	88	<b>44.8</b>	90	1	
Zinc	100	<b>120</b>	96	<b>115</b>	91	4	



Date of Report: October 17, 2016  
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Laboratory Reference: 1610-079  
Project: 31008

**% MOISTURE**

Date Analyzed: 10-10-16

Client ID	Lab ID	% Moisture
H-2-16-13.5	10-079-04	30





### 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.
  - Z -
- ND - Not Detected at PQL  
 PQL - Practical Quantitation Limit  
 RPD - Relative Percent Difference







14648 NE 95<sup>th</sup> Street, Redmond, WA 98052 • (425) 883-3881

October 17, 2016

Glenn Hayman  
INNOVEX Environmental Mgt., Inc.  
16310 NE 80th St., Suite 300  
Redmond, WA 98052

Re: Analytical Data for Project 31008  
Laboratory Reference No. 1610-080

Dear Glenn:

Enclosed are the analytical results and associated quality control data for samples submitted on October 8, 2016.

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,

A handwritten signature in black ink, appearing to read "DB", with a long horizontal flourish extending to the right.

David Baumeister  
Project Manager

Enclosures



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OnSite Environmental, Inc. 14648 NE 95<sup>th</sup> Street, Redmond, WA 98052 (425) 883-3881

This report pertains to the samples analyzed in accordance with the chain of custody, and is intended only for the use of the individual or company to whom it is addressed.

Date of Report: October 17, 2016  
Samples Submitted: October 8, 2016  
Laboratory Reference: 1610-080  
Project: 31008

### Case Narrative

Samples were collected on October 7, 2016 and received by the laboratory on October 8, 2016. 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.

#### Volatiles EPA 8260C 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.

All four Internal Standards did not meet acceptance criteria for samples H-3-16-6 and H-3-16-8.5. The samples were re-analyzed with similar results. Leaks in the sealed VOA environment caused by grit between the VOA lip and VOA cap septum have been shown to cause low internal standard recovery. The samples were consequently extracted from their respective 8-ounce jars, analyzed, and reported. Some loss of volatiles may have occurred, and common laboratory solvents Acetone and Methylene Chloride may have been introduced during sample preparation.

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.



Date of Report: October 17, 2016  
 Samples Submitted: October 8, 2016  
 Laboratory Reference: 1610-080  
 Project: 31008

### NWTPH-HCID

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-3-16-3</b>					
Laboratory ID:	10-080-01					
Gasoline Range Organics	<b>ND</b>	22	NWTPH-HCID	10-10-16	10-10-16	
Diesel Range Organics	<b>ND</b>	55	NWTPH-HCID	10-10-16	10-10-16	
Lube Oil Range Organics	<b>ND</b>	110	NWTPH-HCID	10-10-16	10-10-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	122	50-150				
<b>Client ID:</b>	<b>H-3-16-6</b>					
Laboratory ID:	10-080-02					
Gasoline Range Organics	<b>ND</b>	23	NWTPH-HCID	10-10-16	10-10-16	
Diesel Range Organics	<b>ND</b>	57	NWTPH-HCID	10-10-16	10-10-16	
Lube Oil Range Organics	<b>ND</b>	110	NWTPH-HCID	10-10-16	10-10-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	128	50-150				
<b>Client ID:</b>	<b>H-3-16-8.5</b>					
Laboratory ID:	10-080-03					
Gasoline Range Organics	<b>ND</b>	24	NWTPH-HCID	10-10-16	10-10-16	
Diesel Range Organics	<b>ND</b>	59	NWTPH-HCID	10-10-16	10-10-16	
Lube Oil Range Organics	<b>ND</b>	120	NWTPH-HCID	10-10-16	10-10-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	119	50-150				



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**NWTPH-HCID  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg (ppm)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>METHOD BLANK</b>						
Laboratory ID:	MB1010S2					
Gasoline Range Organics	<b>ND</b>	20	NWTPH-HCID	10-10-16	10-10-16	
Diesel Range Organics	<b>ND</b>	50	NWTPH-HCID	10-10-16	10-10-16	
Lube Oil Range Organics	<b>ND</b>	100	NWTPH-HCID	10-10-16	10-10-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>84</i>	<i>50-150</i>				



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Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>H-3-16-3</b>					
Laboratory ID:	10-080-01					
Dichlorodifluoromethane	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
Chloromethane	ND	0.0045	EPA 8260C	10-10-16	10-10-16	
Vinyl Chloride	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
Bromomethane	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
Chloroethane	ND	0.0045	EPA 8260C	10-10-16	10-10-16	
Trichlorofluoromethane	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
1,1-Dichloroethene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
Acetone	ND	0.0045	EPA 8260C	10-10-16	10-10-16	
Iodomethane	ND	0.0045	EPA 8260C	10-10-16	10-10-16	
Carbon Disulfide	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
Methylene Chloride	ND	0.0045	EPA 8260C	10-10-16	10-10-16	
(trans) 1,2-Dichloroethene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
Methyl t-Butyl Ether	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
1,1-Dichloroethane	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
Vinyl Acetate	ND	0.0045	EPA 8260C	10-10-16	10-10-16	
2,2-Dichloropropane	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
(cis) 1,2-Dichloroethene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
2-Butanone	ND	0.0045	EPA 8260C	10-10-16	10-10-16	
Bromochloromethane	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
Chloroform	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
1,1,1-Trichloroethane	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
Carbon Tetrachloride	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
1,1-Dichloropropene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
Benzene	0.0055	0.00090	EPA 8260C	10-10-16	10-10-16	
1,2-Dichloroethane	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
Trichloroethene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
1,2-Dichloropropane	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
Dibromomethane	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
Bromodichloromethane	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
2-Chloroethyl Vinyl Ether	ND	0.0045	EPA 8260C	10-10-16	10-10-16	
(cis) 1,3-Dichloropropene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
Methyl Isobutyl Ketone	ND	0.0045	EPA 8260C	10-10-16	10-10-16	
Toluene	ND	0.0045	EPA 8260C	10-10-16	10-10-16	
(trans) 1,3-Dichloropropene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	



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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-3-16-3</b>					
Laboratory ID:	10-080-01					
1,1,2-Trichloroethane	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
Tetrachloroethene	ND	0.0018	EPA 8260C	10-10-16	10-10-16	
1,3-Dichloropropane	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
2-Hexanone	ND	0.0045	EPA 8260C	10-10-16	10-10-16	
Dibromochloromethane	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
1,2-Dibromoethane	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
Chlorobenzene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
1,1,1,2-Tetrachloroethane	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
Ethylbenzene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
m,p-Xylene	ND	0.0018	EPA 8260C	10-10-16	10-10-16	
o-Xylene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
Styrene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
Bromoform	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
Isopropylbenzene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
Bromobenzene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
1,1,2,2-Tetrachloroethane	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
1,2,3-Trichloropropane	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
n-Propylbenzene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
2-Chlorotoluene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
4-Chlorotoluene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
1,3,5-Trimethylbenzene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
tert-Butylbenzene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
1,2,4-Trimethylbenzene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
sec-Butylbenzene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
1,3-Dichlorobenzene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
p-Isopropyltoluene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
1,4-Dichlorobenzene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
1,2-Dichlorobenzene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
n-Butylbenzene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
1,2-Dibromo-3-chloropropane	ND	0.0045	EPA 8260C	10-10-16	10-10-16	
1,2,4-Trichlorobenzene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
Hexachlorobutadiene	ND	0.0045	EPA 8260C	10-10-16	10-10-16	
Naphthalene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
1,2,3-Trichlorobenzene	ND	0.00090	EPA 8260C	10-10-16	10-10-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>99</i>	<i>76-131</i>				
<i>Toluene-d8</i>	<i>100</i>	<i>80-126</i>				
<i>4-Bromofluorobenzene</i>	<i>99</i>	<i>60-146</i>				



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Matrix: Soil  
 Units: mg/kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-3-16-6</b>					
Laboratory ID:	10-080-02					
Dichlorodifluoromethane	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
Chloromethane	ND	0.0057	EPA 8260C	10-11-16	10-11-16	
Vinyl Chloride	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
Bromomethane	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
Chloroethane	ND	0.0057	EPA 8260C	10-11-16	10-11-16	
Trichlorofluoromethane	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
1,1-Dichloroethene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
Acetone	0.023	0.0057	EPA 8260C	10-11-16	10-11-16	H
Iodomethane	ND	0.0057	EPA 8260C	10-11-16	10-11-16	
Carbon Disulfide	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
Methylene Chloride	0.053	0.0057	EPA 8260C	10-11-16	10-11-16	H
(trans) 1,2-Dichloroethene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
Methyl t-Butyl Ether	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
1,1-Dichloroethane	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
Vinyl Acetate	ND	0.0057	EPA 8260C	10-11-16	10-11-16	
2,2-Dichloropropane	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
(cis) 1,2-Dichloroethene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
2-Butanone	ND	0.0057	EPA 8260C	10-11-16	10-11-16	
Bromochloromethane	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
Chloroform	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
1,1,1-Trichloroethane	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
Carbon Tetrachloride	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
1,1-Dichloropropene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
Benzene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
1,2-Dichloroethane	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
Trichloroethene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
1,2-Dichloropropane	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
Dibromomethane	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
Bromodichloromethane	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
2-Chloroethyl Vinyl Ether	ND	0.0057	EPA 8260C	10-11-16	10-11-16	
(cis) 1,3-Dichloropropene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
Methyl Isobutyl Ketone	ND	0.0057	EPA 8260C	10-11-16	10-11-16	
Toluene	ND	0.0057	EPA 8260C	10-11-16	10-11-16	
(trans) 1,3-Dichloropropene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	



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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-3-16-6</b>					
<b>Laboratory ID:</b>	10-080-02					
1,1,2-Trichloroethane	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
Tetrachloroethene	ND	0.0023	EPA 8260C	10-11-16	10-11-16	
1,3-Dichloropropane	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
2-Hexanone	ND	0.0057	EPA 8260C	10-11-16	10-11-16	
Dibromochloromethane	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
1,2-Dibromoethane	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
Chlorobenzene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
1,1,1,2-Tetrachloroethane	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
Ethylbenzene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
m,p-Xylene	ND	0.0023	EPA 8260C	10-11-16	10-11-16	
o-Xylene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
Styrene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
Bromoform	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
Isopropylbenzene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
Bromobenzene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
1,1,2,2-Tetrachloroethane	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
1,2,3-Trichloropropane	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
n-Propylbenzene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
2-Chlorotoluene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
4-Chlorotoluene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
1,3,5-Trimethylbenzene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
tert-Butylbenzene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
1,2,4-Trimethylbenzene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
sec-Butylbenzene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
1,3-Dichlorobenzene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
p-Isopropyltoluene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
1,4-Dichlorobenzene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
1,2-Dichlorobenzene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
n-Butylbenzene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
1,2-Dibromo-3-chloropropane	ND	0.0057	EPA 8260C	10-11-16	10-11-16	
1,2,4-Trichlorobenzene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
Hexachlorobutadiene	ND	0.0057	EPA 8260C	10-11-16	10-11-16	
Naphthalene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
1,2,3-Trichlorobenzene	ND	0.0011	EPA 8260C	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>100</i>	<i>76-131</i>				
<i>Toluene-d8</i>	<i>96</i>	<i>80-126</i>				
<i>4-Bromofluorobenzene</i>	<i>100</i>	<i>60-146</i>				



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Matrix: Soil  
 Units: mg/kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-3-16-8.5</b>					
Laboratory ID:	10-080-03					
Dichlorodifluoromethane	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
Chloromethane	ND	0.0058	EPA 8260C	10-11-16	10-11-16	
Vinyl Chloride	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
Bromomethane	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
Chloroethane	ND	0.0058	EPA 8260C	10-11-16	10-11-16	
Trichlorofluoromethane	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
1,1-Dichloroethene	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
Acetone	ND	0.0058	EPA 8260C	10-11-16	10-11-16	
Iodomethane	ND	0.0058	EPA 8260C	10-11-16	10-11-16	
Carbon Disulfide	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
Methylene Chloride	0.022	0.0058	EPA 8260C	10-11-16	10-11-16	H
(trans) 1,2-Dichloroethene	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
Methyl t-Butyl Ether	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
1,1-Dichloroethane	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
Vinyl Acetate	ND	0.0058	EPA 8260C	10-11-16	10-11-16	
2,2-Dichloropropane	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
(cis) 1,2-Dichloroethene	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
2-Butanone	ND	0.0058	EPA 8260C	10-11-16	10-11-16	
Bromochloromethane	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
Chloroform	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
1,1,1-Trichloroethane	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
Carbon Tetrachloride	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
1,1-Dichloropropene	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
Benzene	0.038	0.0012	EPA 8260C	10-11-16	10-11-16	
1,2-Dichloroethane	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
Trichloroethene	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
1,2-Dichloropropane	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
Dibromomethane	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
Bromodichloromethane	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
2-Chloroethyl Vinyl Ether	ND	0.0058	EPA 8260C	10-11-16	10-11-16	
(cis) 1,3-Dichloropropene	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
Methyl Isobutyl Ketone	ND	0.0058	EPA 8260C	10-11-16	10-11-16	
Toluene	ND	0.0058	EPA 8260C	10-11-16	10-11-16	
(trans) 1,3-Dichloropropene	ND	0.0012	EPA 8260C	10-11-16	10-11-16	



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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-3-16-8.5</b>					
Laboratory ID:	10-080-03					
1,1,2-Trichloroethane	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
Tetrachloroethene	ND	0.0023	EPA 8260C	10-11-16	10-11-16	
1,3-Dichloropropane	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
2-Hexanone	ND	0.0058	EPA 8260C	10-11-16	10-11-16	
Dibromochloromethane	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
1,2-Dibromoethane	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
Chlorobenzene	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
1,1,1,2-Tetrachloroethane	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
Ethylbenzene	0.0050	0.0012	EPA 8260C	10-11-16	10-11-16	
m,p-Xylene	0.014	0.0023	EPA 8260C	10-11-16	10-11-16	
o-Xylene	0.0026	0.0012	EPA 8260C	10-11-16	10-11-16	
Styrene	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
Bromoform	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
Isopropylbenzene	0.0018	0.0012	EPA 8260C	10-11-16	10-11-16	
Bromobenzene	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
1,1,2,2-Tetrachloroethane	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
1,2,3-Trichloropropane	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
n-Propylbenzene	0.0032	0.0012	EPA 8260C	10-11-16	10-11-16	
2-Chlorotoluene	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
4-Chlorotoluene	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
1,3,5-Trimethylbenzene	0.0015	0.0012	EPA 8260C	10-11-16	10-11-16	
tert-Butylbenzene	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
1,2,4-Trimethylbenzene	0.0018	0.0012	EPA 8260C	10-11-16	10-11-16	
sec-Butylbenzene	0.0013	0.0012	EPA 8260C	10-11-16	10-11-16	
1,3-Dichlorobenzene	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
p-Isopropyltoluene	0.0020	0.0012	EPA 8260C	10-11-16	10-11-16	
1,4-Dichlorobenzene	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
1,2-Dichlorobenzene	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
n-Butylbenzene	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
1,2-Dibromo-3-chloropropane	ND	0.0058	EPA 8260C	10-11-16	10-11-16	
1,2,4-Trichlorobenzene	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
Hexachlorobutadiene	ND	0.0058	EPA 8260C	10-11-16	10-11-16	
Naphthalene	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
1,2,3-Trichlorobenzene	ND	0.0012	EPA 8260C	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>87</i>	<i>76-131</i>				
<i>Toluene-d8</i>	<i>89</i>	<i>80-126</i>				
<i>4-Bromofluorobenzene</i>	<i>83</i>	<i>60-146</i>				



Date of Report: October 17, 2016  
 Samples Submitted: October 8, 2016  
 Laboratory Reference: 1610-080  
 Project: 31008

**VOLATILES by EPA 8260C**  
**METHOD BLANK QUALITY CONTROL**  
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Matrix: Soil  
 Units: mg/kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1010S1					
Dichlorodifluoromethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Chloromethane	ND	0.0050	EPA 8260C	10-10-16	10-10-16	
Vinyl Chloride	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Bromomethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Chloroethane	ND	0.0050	EPA 8260C	10-10-16	10-10-16	
Trichlorofluoromethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,1-Dichloroethene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Acetone	ND	0.0050	EPA 8260C	10-10-16	10-10-16	
Iodomethane	ND	0.0050	EPA 8260C	10-10-16	10-10-16	
Carbon Disulfide	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Methylene Chloride	ND	0.0050	EPA 8260C	10-10-16	10-10-16	
(trans) 1,2-Dichloroethene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Methyl t-Butyl Ether	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,1-Dichloroethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Vinyl Acetate	ND	0.0050	EPA 8260C	10-10-16	10-10-16	
2,2-Dichloropropane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
(cis) 1,2-Dichloroethene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
2-Butanone	ND	0.0050	EPA 8260C	10-10-16	10-10-16	
Bromochloromethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Chloroform	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,1,1-Trichloroethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Carbon Tetrachloride	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,1-Dichloropropene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Benzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,2-Dichloroethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Trichloroethene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,2-Dichloropropane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Dibromomethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Bromodichloromethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
2-Chloroethyl Vinyl Ether	ND	0.0050	EPA 8260C	10-10-16	10-10-16	
(cis) 1,3-Dichloropropene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Methyl Isobutyl Ketone	ND	0.0050	EPA 8260C	10-10-16	10-10-16	
Toluene	ND	0.0050	EPA 8260C	10-10-16	10-10-16	
(trans) 1,3-Dichloropropene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	



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**VOLATILES by EPA 8260C**  
**METHOD BLANK QUALITY CONTROL**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:		MB1010S1				
1,1,2-Trichloroethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Tetrachloroethene	ND	0.0020	EPA 8260C	10-10-16	10-10-16	
1,3-Dichloropropane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
2-Hexanone	ND	0.0050	EPA 8260C	10-10-16	10-10-16	
Dibromochloromethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,2-Dibromoethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Chlorobenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,1,1,2-Tetrachloroethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Ethylbenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
m,p-Xylene	ND	0.0020	EPA 8260C	10-10-16	10-10-16	
o-Xylene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Styrene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Bromoform	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Isopropylbenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Bromobenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,1,2,2-Tetrachloroethane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,2,3-Trichloropropane	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
n-Propylbenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
2-Chlorotoluene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
4-Chlorotoluene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,3,5-Trimethylbenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
tert-Butylbenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,2,4-Trimethylbenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
sec-Butylbenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,3-Dichlorobenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
p-Isopropyltoluene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,4-Dichlorobenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,2-Dichlorobenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
n-Butylbenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,2-Dibromo-3-chloropropane	ND	0.0050	EPA 8260C	10-10-16	10-10-16	
1,2,4-Trichlorobenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
Hexachlorobutadiene	ND	0.0050	EPA 8260C	10-10-16	10-10-16	
Naphthalene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
1,2,3-Trichlorobenzene	ND	0.0010	EPA 8260C	10-10-16	10-10-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>98</i>	<i>76-131</i>				
<i>Toluene-d8</i>	<i>103</i>	<i>80-126</i>				
<i>4-Bromofluorobenzene</i>	<i>102</i>	<i>60-146</i>				



Date of Report: October 17, 2016  
 Samples Submitted: October 8, 2016  
 Laboratory Reference: 1610-080  
 Project: 31008

**VOLATILES by EPA 8260C**  
**METHOD BLANK QUALITY CONTROL**  
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Matrix: Soil  
 Units: mg/kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1011S1					
Dichlorodifluoromethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Chloromethane	ND	0.0050	EPA 8260C	10-11-16	10-11-16	
Vinyl Chloride	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Bromomethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Chloroethane	ND	0.0050	EPA 8260C	10-11-16	10-11-16	
Trichlorofluoromethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,1-Dichloroethene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Acetone	ND	0.0050	EPA 8260C	10-11-16	10-11-16	
Iodomethane	ND	0.0050	EPA 8260C	10-11-16	10-11-16	
Carbon Disulfide	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Methylene Chloride	ND	0.0050	EPA 8260C	10-11-16	10-11-16	
(trans) 1,2-Dichloroethene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Methyl t-Butyl Ether	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,1-Dichloroethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Vinyl Acetate	ND	0.0050	EPA 8260C	10-11-16	10-11-16	
2,2-Dichloropropane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
(cis) 1,2-Dichloroethene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
2-Butanone	ND	0.0050	EPA 8260C	10-11-16	10-11-16	
Bromochloromethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Chloroform	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,1,1-Trichloroethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Carbon Tetrachloride	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,1-Dichloropropene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Benzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,2-Dichloroethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Trichloroethene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,2-Dichloropropane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Dibromomethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Bromodichloromethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
2-Chloroethyl Vinyl Ether	ND	0.0050	EPA 8260C	10-11-16	10-11-16	
(cis) 1,3-Dichloropropene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Methyl Isobutyl Ketone	ND	0.0050	EPA 8260C	10-11-16	10-11-16	
Toluene	ND	0.0050	EPA 8260C	10-11-16	10-11-16	
(trans) 1,3-Dichloropropene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	



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**VOLATILES by EPA 8260C**  
**METHOD BLANK QUALITY CONTROL**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1011S1					
1,1,2-Trichloroethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Tetrachloroethene	ND	0.0020	EPA 8260C	10-11-16	10-11-16	
1,3-Dichloropropane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
2-Hexanone	ND	0.0050	EPA 8260C	10-11-16	10-11-16	
Dibromochloromethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,2-Dibromoethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Chlorobenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,1,1,2-Tetrachloroethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Ethylbenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
m,p-Xylene	ND	0.0020	EPA 8260C	10-11-16	10-11-16	
o-Xylene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Styrene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Bromoform	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Isopropylbenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Bromobenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,1,2,2-Tetrachloroethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,2,3-Trichloropropane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
n-Propylbenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
2-Chlorotoluene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
4-Chlorotoluene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,3,5-Trimethylbenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
tert-Butylbenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,2,4-Trimethylbenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
sec-Butylbenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,3-Dichlorobenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
p-Isopropyltoluene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,4-Dichlorobenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,2-Dichlorobenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
n-Butylbenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,2-Dibromo-3-chloropropane	ND	0.0050	EPA 8260C	10-11-16	10-11-16	
1,2,4-Trichlorobenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Hexachlorobutadiene	ND	0.0050	EPA 8260C	10-11-16	10-11-16	
Naphthalene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,2,3-Trichlorobenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>104</i>	<i>76-131</i>				
<i>Toluene-d8</i>	<i>107</i>	<i>80-126</i>				
<i>4-Bromofluorobenzene</i>	<i>102</i>	<i>60-146</i>				



Date of Report: October 17, 2016  
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 Laboratory Reference: 1610-080  
 Project: 31008

**VOLATILES by EPA 8260C  
 SB/SBD QUALITY CONTROL**

Matrix: Soil  
 Units: mg/kg

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD		Flags
					Recovery	Limits	RPD	Limit		
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB1010S1									
	SB	SBD	SB	SBD	SB	SBD				
1,1-Dichloroethene	<b>0.0457</b>	<b>0.0490</b>	0.0500	0.0500	91	98	68-126	7	15	
Benzene	<b>0.0461</b>	<b>0.0480</b>	0.0500	0.0500	92	96	70-121	4	15	
Trichloroethene	<b>0.0455</b>	<b>0.0476</b>	0.0500	0.0500	91	95	75-120	5	15	
Toluene	<b>0.0468</b>	<b>0.0495</b>	0.0500	0.0500	94	99	80-120	6	15	
Chlorobenzene	<b>0.0475</b>	<b>0.0490</b>	0.0500	0.0500	95	98	76-120	3	15	
<i>Surrogate:</i>										
Dibromofluoromethane					94	94	76-131			
Toluene-d8					94	100	80-126			
4-Bromofluorobenzene					96	96	60-146			



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**VOLATILES by EPA 8260C  
 SB/SBD QUALITY CONTROL**

Matrix: Soil  
 Units: mg/kg

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD		Flags
					Recovery	Limits	RPD	Limit		
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB1011S1									
	SB	SBD	SB	SBD	SB	SBD				
1,1-Dichloroethene	<b>0.0472</b>	<b>0.0496</b>	0.0500	0.0500	94	99	68-126	5	15	
Benzene	<b>0.0473</b>	<b>0.0487</b>	0.0500	0.0500	95	97	70-121	3	15	
Trichloroethene	<b>0.0440</b>	<b>0.0462</b>	0.0500	0.0500	88	92	75-120	5	15	
Toluene	<b>0.0459</b>	<b>0.0487</b>	0.0500	0.0500	92	97	80-120	6	15	
Chlorobenzene	<b>0.0474</b>	<b>0.0478</b>	0.0500	0.0500	95	96	76-120	1	15	
<i>Surrogate:</i>										
Dibromofluoromethane					98	99	76-131			
Toluene-d8					98	100	80-126			
4-Bromofluorobenzene					100	101	60-146			



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**SEMIVOLATILES EPA 8270D/SIM**  
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Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-3-16-3</b>					
<b>Laboratory ID:</b>	<b>10-080-01</b>					
n-Nitrosodimethylamine	ND	0.037	EPA 8270D	10-10-16	10-12-16	
Pyridine	ND	0.37	EPA 8270D	10-10-16	10-12-16	
Phenol	ND	0.037	EPA 8270D	10-10-16	10-12-16	
Aniline	ND	0.18	EPA 8270D	10-10-16	10-12-16	
bis(2-Chloroethyl)ether	ND	0.037	EPA 8270D	10-10-16	10-12-16	
2-Chlorophenol	ND	0.037	EPA 8270D	10-10-16	10-12-16	
1,3-Dichlorobenzene	ND	0.037	EPA 8270D	10-10-16	10-12-16	
1,4-Dichlorobenzene	ND	0.037	EPA 8270D	10-10-16	10-12-16	
Benzyl alcohol	ND	0.18	EPA 8270D	10-10-16	10-12-16	
1,2-Dichlorobenzene	ND	0.037	EPA 8270D	10-10-16	10-12-16	
2-Methylphenol (o-Cresol)	ND	0.037	EPA 8270D	10-10-16	10-12-16	
bis(2-Chloroisopropyl)ether	ND	0.037	EPA 8270D	10-10-16	10-12-16	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.037	EPA 8270D	10-10-16	10-12-16	
n-Nitroso-di-n-propylamine	ND	0.037	EPA 8270D	10-10-16	10-12-16	
Hexachloroethane	ND	0.037	EPA 8270D	10-10-16	10-12-16	
Nitrobenzene	ND	0.037	EPA 8270D	10-10-16	10-12-16	
Isophorone	ND	0.037	EPA 8270D	10-10-16	10-12-16	
2-Nitrophenol	ND	0.037	EPA 8270D	10-10-16	10-12-16	
2,4-Dimethylphenol	ND	0.037	EPA 8270D	10-10-16	10-12-16	
bis(2-Chloroethoxy)methane	ND	0.037	EPA 8270D	10-10-16	10-12-16	
2,4-Dichlorophenol	ND	0.037	EPA 8270D	10-10-16	10-12-16	
1,2,4-Trichlorobenzene	ND	0.037	EPA 8270D	10-10-16	10-12-16	
Naphthalene	ND	0.0073	EPA 8270D/SIM	10-10-16	10-10-16	
4-Chloroaniline	ND	0.18	EPA 8270D	10-10-16	10-12-16	
Hexachlorobutadiene	ND	0.037	EPA 8270D	10-10-16	10-12-16	
4-Chloro-3-methylphenol	ND	0.037	EPA 8270D	10-10-16	10-12-16	
2-Methylnaphthalene	ND	0.0073	EPA 8270D/SIM	10-10-16	10-10-16	
1-Methylnaphthalene	ND	0.0073	EPA 8270D/SIM	10-10-16	10-10-16	
Hexachlorocyclopentadiene	ND	0.037	EPA 8270D	10-10-16	10-12-16	
2,4,6-Trichlorophenol	ND	0.037	EPA 8270D	10-10-16	10-12-16	
2,3-Dichloroaniline	ND	0.037	EPA 8270D	10-10-16	10-12-16	
2,4,5-Trichlorophenol	ND	0.037	EPA 8270D	10-10-16	10-12-16	
2-Chloronaphthalene	ND	0.037	EPA 8270D	10-10-16	10-12-16	
2-Nitroaniline	ND	0.037	EPA 8270D	10-10-16	10-12-16	
1,4-Dinitrobenzene	ND	0.037	EPA 8270D	10-10-16	10-12-16	
Dimethylphthalate	ND	0.037	EPA 8270D	10-10-16	10-12-16	
1,3-Dinitrobenzene	ND	0.037	EPA 8270D	10-10-16	10-12-16	
2,6-Dinitrotoluene	ND	0.037	EPA 8270D	10-10-16	10-12-16	
1,2-Dinitrobenzene	ND	0.037	EPA 8270D	10-10-16	10-12-16	
Acenaphthylene	ND	0.0073	EPA 8270D/SIM	10-10-16	10-10-16	
3-Nitroaniline	ND	0.037	EPA 8270D	10-10-16	10-12-16	



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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-3-16-3</b>					
<b>Laboratory ID:</b>	<b>10-080-01</b>					
2,4-Dinitrophenol	ND	0.18	EPA 8270D	10-10-16	10-12-16	
Acenaphthene	ND	0.0073	EPA 8270D/SIM	10-10-16	10-10-16	
4-Nitrophenol	ND	0.037	EPA 8270D	10-10-16	10-12-16	
2,4-Dinitrotoluene	ND	0.037	EPA 8270D	10-10-16	10-12-16	
Dibenzofuran	ND	0.037	EPA 8270D	10-10-16	10-12-16	
2,3,5,6-Tetrachlorophenol	ND	0.037	EPA 8270D	10-10-16	10-12-16	
2,3,4,6-Tetrachlorophenol	ND	0.037	EPA 8270D	10-10-16	10-12-16	
Diethylphthalate	ND	0.18	EPA 8270D	10-10-16	10-12-16	
4-Chlorophenyl-phenylether	ND	0.037	EPA 8270D	10-10-16	10-12-16	
4-Nitroaniline	ND	0.037	EPA 8270D	10-10-16	10-12-16	
Fluorene	ND	0.0073	EPA 8270D/SIM	10-10-16	10-10-16	
4,6-Dinitro-2-methylphenol	ND	0.18	EPA 8270D	10-10-16	10-12-16	
n-Nitrosodiphenylamine	ND	0.037	EPA 8270D	10-10-16	10-12-16	
1,2-Diphenylhydrazine	ND	0.037	EPA 8270D	10-10-16	10-12-16	
4-Bromophenyl-phenylether	ND	0.037	EPA 8270D	10-10-16	10-12-16	
Hexachlorobenzene	ND	0.037	EPA 8270D	10-10-16	10-12-16	
Pentachlorophenol	ND	0.18	EPA 8270D	10-10-16	10-12-16	
Phenanthrene	ND	0.0073	EPA 8270D/SIM	10-10-16	10-10-16	
Anthracene	ND	0.0073	EPA 8270D/SIM	10-10-16	10-10-16	
Carbazole	ND	0.037	EPA 8270D	10-10-16	10-12-16	
Di-n-butylphthalate	ND	0.18	EPA 8270D	10-10-16	10-12-16	
Fluoranthene	ND	0.0073	EPA 8270D/SIM	10-10-16	10-10-16	
Benzidine	ND	0.37	EPA 8270D	10-10-16	10-12-16	
Pyrene	ND	0.0073	EPA 8270D/SIM	10-10-16	10-10-16	
Butylbenzylphthalate	ND	0.037	EPA 8270D	10-10-16	10-12-16	
bis-2-Ethylhexyladipate	ND	0.037	EPA 8270D	10-10-16	10-12-16	
3,3'-Dichlorobenzidine	ND	0.18	EPA 8270D	10-10-16	10-12-16	
Benzo[a]anthracene	ND	0.0073	EPA 8270D/SIM	10-10-16	10-10-16	
Chrysene	ND	0.0073	EPA 8270D/SIM	10-10-16	10-10-16	
bis(2-Ethylhexyl)phthalate	ND	0.037	EPA 8270D	10-10-16	10-12-16	
Di-n-octylphthalate	ND	0.037	EPA 8270D	10-10-16	10-12-16	
Benzo[b]fluoranthene	ND	0.0073	EPA 8270D/SIM	10-10-16	10-10-16	
Benzo(j,k)fluoranthene	ND	0.0073	EPA 8270D/SIM	10-10-16	10-10-16	
Benzo[a]pyrene	ND	0.0073	EPA 8270D/SIM	10-10-16	10-10-16	
Indeno[1,2,3-cd]pyrene	ND	0.0073	EPA 8270D/SIM	10-10-16	10-10-16	
Dibenz[a,h]anthracene	ND	0.0073	EPA 8270D/SIM	10-10-16	10-10-16	
Benzo[g,h,i]perylene	ND	0.0073	EPA 8270D/SIM	10-10-16	10-10-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorophenol</i>	<i>57</i>	<i>24 - 117</i>				
<i>Phenol-d6</i>	<i>61</i>	<i>30 - 120</i>				
<i>Nitrobenzene-d5</i>	<i>61</i>	<i>27 - 112</i>				
<i>2-Fluorobiphenyl</i>	<i>62</i>	<i>35 - 113</i>				
<i>2,4,6-Tribromophenol</i>	<i>67</i>	<i>21 - 120</i>				
<i>Terphenyl-d14</i>	<i>65</i>	<i>39 - 121</i>				



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Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-3-16-6</b>					
<b>Laboratory ID:</b>	<b>10-080-02</b>					
n-Nitrosodimethylamine	ND	0.038	EPA 8270D	10-10-16	10-11-16	
Pyridine	ND	0.38	EPA 8270D	10-10-16	10-11-16	
Phenol	ND	0.038	EPA 8270D	10-10-16	10-11-16	
Aniline	ND	0.19	EPA 8270D	10-10-16	10-11-16	
bis(2-Chloroethyl)ether	ND	0.038	EPA 8270D	10-10-16	10-11-16	
2-Chlorophenol	ND	0.038	EPA 8270D	10-10-16	10-11-16	
1,3-Dichlorobenzene	ND	0.038	EPA 8270D	10-10-16	10-11-16	
1,4-Dichlorobenzene	ND	0.038	EPA 8270D	10-10-16	10-11-16	
Benzyl alcohol	ND	0.19	EPA 8270D	10-10-16	10-11-16	
1,2-Dichlorobenzene	ND	0.038	EPA 8270D	10-10-16	10-11-16	
2-Methylphenol (o-Cresol)	ND	0.038	EPA 8270D	10-10-16	10-11-16	
bis(2-Chloroisopropyl)ether	ND	0.038	EPA 8270D	10-10-16	10-11-16	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.038	EPA 8270D	10-10-16	10-11-16	
n-Nitroso-di-n-propylamine	ND	0.038	EPA 8270D	10-10-16	10-11-16	
Hexachloroethane	ND	0.038	EPA 8270D	10-10-16	10-11-16	
Nitrobenzene	ND	0.038	EPA 8270D	10-10-16	10-11-16	
Isophorone	ND	0.038	EPA 8270D	10-10-16	10-11-16	
2-Nitrophenol	ND	0.038	EPA 8270D	10-10-16	10-11-16	
2,4-Dimethylphenol	ND	0.038	EPA 8270D	10-10-16	10-11-16	
bis(2-Chloroethoxy)methane	ND	0.038	EPA 8270D	10-10-16	10-11-16	
2,4-Dichlorophenol	ND	0.038	EPA 8270D	10-10-16	10-11-16	
1,2,4-Trichlorobenzene	ND	0.038	EPA 8270D	10-10-16	10-11-16	
Naphthalene	ND	0.0076	EPA 8270D/SIM	10-10-16	10-10-16	
4-Chloroaniline	ND	0.19	EPA 8270D	10-10-16	10-11-16	
Hexachlorobutadiene	ND	0.038	EPA 8270D	10-10-16	10-11-16	
4-Chloro-3-methylphenol	ND	0.038	EPA 8270D	10-10-16	10-11-16	
2-Methylnaphthalene	ND	0.0076	EPA 8270D/SIM	10-10-16	10-10-16	
1-Methylnaphthalene	ND	0.0076	EPA 8270D/SIM	10-10-16	10-10-16	
Hexachlorocyclopentadiene	ND	0.038	EPA 8270D	10-10-16	10-11-16	
2,4,6-Trichlorophenol	ND	0.038	EPA 8270D	10-10-16	10-11-16	
2,3-Dichloroaniline	ND	0.038	EPA 8270D	10-10-16	10-11-16	
2,4,5-Trichlorophenol	ND	0.038	EPA 8270D	10-10-16	10-11-16	
2-Chloronaphthalene	ND	0.038	EPA 8270D	10-10-16	10-11-16	
2-Nitroaniline	ND	0.038	EPA 8270D	10-10-16	10-11-16	
1,4-Dinitrobenzene	ND	0.038	EPA 8270D	10-10-16	10-11-16	
Dimethylphthalate	ND	0.038	EPA 8270D	10-10-16	10-11-16	
1,3-Dinitrobenzene	ND	0.038	EPA 8270D	10-10-16	10-11-16	
2,6-Dinitrotoluene	ND	0.038	EPA 8270D	10-10-16	10-11-16	
1,2-Dinitrobenzene	ND	0.038	EPA 8270D	10-10-16	10-11-16	
Acenaphthylene	ND	0.0076	EPA 8270D/SIM	10-10-16	10-10-16	
3-Nitroaniline	ND	0.038	EPA 8270D	10-10-16	10-11-16	



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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-3-16-6</b>					
Laboratory ID:	10-080-02					
2,4-Dinitrophenol	ND	0.19	EPA 8270D	10-10-16	10-11-16	
Acenaphthene	ND	0.0076	EPA 8270D/SIM	10-10-16	10-10-16	
4-Nitrophenol	ND	0.038	EPA 8270D	10-10-16	10-11-16	
2,4-Dinitrotoluene	ND	0.038	EPA 8270D	10-10-16	10-11-16	
Dibenzofuran	ND	0.038	EPA 8270D	10-10-16	10-11-16	
2,3,5,6-Tetrachlorophenol	ND	0.038	EPA 8270D	10-10-16	10-11-16	
2,3,4,6-Tetrachlorophenol	ND	0.038	EPA 8270D	10-10-16	10-11-16	
Diethylphthalate	ND	0.19	EPA 8270D	10-10-16	10-11-16	
4-Chlorophenyl-phenylether	ND	0.038	EPA 8270D	10-10-16	10-11-16	
4-Nitroaniline	ND	0.038	EPA 8270D	10-10-16	10-11-16	
Fluorene	ND	0.0076	EPA 8270D/SIM	10-10-16	10-10-16	
4,6-Dinitro-2-methylphenol	ND	0.19	EPA 8270D	10-10-16	10-11-16	
n-Nitrosodiphenylamine	ND	0.038	EPA 8270D	10-10-16	10-11-16	
1,2-Diphenylhydrazine	ND	0.038	EPA 8270D	10-10-16	10-11-16	
4-Bromophenyl-phenylether	ND	0.038	EPA 8270D	10-10-16	10-11-16	
Hexachlorobenzene	ND	0.038	EPA 8270D	10-10-16	10-11-16	
Pentachlorophenol	ND	0.19	EPA 8270D	10-10-16	10-11-16	
Phenanthrene	ND	0.0076	EPA 8270D/SIM	10-10-16	10-10-16	
Anthracene	ND	0.0076	EPA 8270D/SIM	10-10-16	10-10-16	
Carbazole	ND	0.038	EPA 8270D	10-10-16	10-11-16	
Di-n-butylphthalate	ND	0.19	EPA 8270D	10-10-16	10-11-16	
Fluoranthene	ND	0.0076	EPA 8270D/SIM	10-10-16	10-10-16	
Benzidine	ND	0.38	EPA 8270D	10-10-16	10-11-16	
Pyrene	ND	0.0076	EPA 8270D/SIM	10-10-16	10-10-16	
Butylbenzylphthalate	ND	0.038	EPA 8270D	10-10-16	10-11-16	
bis-2-Ethylhexyladipate	ND	0.038	EPA 8270D	10-10-16	10-11-16	
3,3'-Dichlorobenzidine	ND	0.19	EPA 8270D	10-10-16	10-11-16	
Benzo[a]anthracene	ND	0.0076	EPA 8270D/SIM	10-10-16	10-10-16	
Chrysene	ND	0.0076	EPA 8270D/SIM	10-10-16	10-10-16	
bis(2-Ethylhexyl)phthalate	ND	0.038	EPA 8270D	10-10-16	10-11-16	
Di-n-octylphthalate	ND	0.038	EPA 8270D	10-10-16	10-11-16	
Benzo[b]fluoranthene	ND	0.0076	EPA 8270D/SIM	10-10-16	10-10-16	
Benzo(j,k)fluoranthene	ND	0.0076	EPA 8270D/SIM	10-10-16	10-10-16	
Benzo[a]pyrene	ND	0.0076	EPA 8270D/SIM	10-10-16	10-10-16	
Indeno[1,2,3-cd]pyrene	ND	0.0076	EPA 8270D/SIM	10-10-16	10-10-16	
Dibenz[a,h]anthracene	ND	0.0076	EPA 8270D/SIM	10-10-16	10-10-16	
Benzo[g,h,i]perylene	ND	0.0076	EPA 8270D/SIM	10-10-16	10-10-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	71	24 - 117				
Phenol-d6	75	30 - 120				
Nitrobenzene-d5	77	27 - 112				
2-Fluorobiphenyl	76	35 - 113				
2,4,6-Tribromophenol	79	21 - 120				
Terphenyl-d14	78	39 - 121				



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Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-3-16-8.5</b>					
<b>Laboratory ID:</b>	<b>10-080-03</b>					
n-Nitrosodimethylamine	ND	0.039	EPA 8270D	10-10-16	10-12-16	
Pyridine	ND	0.39	EPA 8270D	10-10-16	10-12-16	
Phenol	ND	0.039	EPA 8270D	10-10-16	10-12-16	
Aniline	ND	0.20	EPA 8270D	10-10-16	10-12-16	
bis(2-Chloroethyl)ether	ND	0.039	EPA 8270D	10-10-16	10-12-16	
2-Chlorophenol	ND	0.039	EPA 8270D	10-10-16	10-12-16	
1,3-Dichlorobenzene	ND	0.039	EPA 8270D	10-10-16	10-12-16	
1,4-Dichlorobenzene	ND	0.039	EPA 8270D	10-10-16	10-12-16	
Benzyl alcohol	ND	0.20	EPA 8270D	10-10-16	10-12-16	
1,2-Dichlorobenzene	ND	0.039	EPA 8270D	10-10-16	10-12-16	
2-Methylphenol (o-Cresol)	ND	0.039	EPA 8270D	10-10-16	10-12-16	
bis(2-Chloroisopropyl)ether	ND	0.039	EPA 8270D	10-10-16	10-12-16	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.039	EPA 8270D	10-10-16	10-12-16	
n-Nitroso-di-n-propylamine	ND	0.039	EPA 8270D	10-10-16	10-12-16	
Hexachloroethane	ND	0.039	EPA 8270D	10-10-16	10-12-16	
Nitrobenzene	ND	0.039	EPA 8270D	10-10-16	10-12-16	
Isophorone	ND	0.039	EPA 8270D	10-10-16	10-12-16	
2-Nitrophenol	ND	0.039	EPA 8270D	10-10-16	10-12-16	
2,4-Dimethylphenol	ND	0.039	EPA 8270D	10-10-16	10-12-16	
bis(2-Chloroethoxy)methane	ND	0.039	EPA 8270D	10-10-16	10-12-16	
2,4-Dichlorophenol	ND	0.039	EPA 8270D	10-10-16	10-12-16	
1,2,4-Trichlorobenzene	ND	0.039	EPA 8270D	10-10-16	10-12-16	
Naphthalene	ND	0.0078	EPA 8270D/SIM	10-10-16	10-10-16	
4-Chloroaniline	ND	0.20	EPA 8270D	10-10-16	10-12-16	
Hexachlorobutadiene	ND	0.039	EPA 8270D	10-10-16	10-12-16	
4-Chloro-3-methylphenol	ND	0.039	EPA 8270D	10-10-16	10-12-16	
2-Methylnaphthalene	0.018	0.0078	EPA 8270D/SIM	10-10-16	10-10-16	
1-Methylnaphthalene	0.013	0.0078	EPA 8270D/SIM	10-10-16	10-10-16	
Hexachlorocyclopentadiene	ND	0.039	EPA 8270D	10-10-16	10-12-16	
2,4,6-Trichlorophenol	ND	0.039	EPA 8270D	10-10-16	10-12-16	
2,3-Dichloroaniline	ND	0.039	EPA 8270D	10-10-16	10-12-16	
2,4,5-Trichlorophenol	ND	0.039	EPA 8270D	10-10-16	10-12-16	
2-Chloronaphthalene	ND	0.039	EPA 8270D	10-10-16	10-12-16	
2-Nitroaniline	ND	0.039	EPA 8270D	10-10-16	10-12-16	
1,4-Dinitrobenzene	ND	0.039	EPA 8270D	10-10-16	10-12-16	
Dimethylphthalate	ND	0.039	EPA 8270D	10-10-16	10-12-16	
1,3-Dinitrobenzene	ND	0.039	EPA 8270D	10-10-16	10-12-16	
2,6-Dinitrotoluene	ND	0.039	EPA 8270D	10-10-16	10-12-16	
1,2-Dinitrobenzene	ND	0.039	EPA 8270D	10-10-16	10-12-16	
Acenaphthylene	ND	0.0078	EPA 8270D/SIM	10-10-16	10-10-16	
3-Nitroaniline	ND	0.039	EPA 8270D	10-10-16	10-12-16	



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 Laboratory Reference: 1610-080  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-3-16-8.5</b>					
<b>Laboratory ID:</b>	10-080-03					
2,4-Dinitrophenol	ND	0.20	EPA 8270D	10-10-16	10-12-16	
Acenaphthene	ND	0.0078	EPA 8270D/SIM	10-10-16	10-10-16	
4-Nitrophenol	ND	0.039	EPA 8270D	10-10-16	10-12-16	
2,4-Dinitrotoluene	ND	0.039	EPA 8270D	10-10-16	10-12-16	
Dibenzofuran	ND	0.039	EPA 8270D	10-10-16	10-12-16	
2,3,5,6-Tetrachlorophenol	ND	0.039	EPA 8270D	10-10-16	10-12-16	
2,3,4,6-Tetrachlorophenol	ND	0.039	EPA 8270D	10-10-16	10-12-16	
Diethylphthalate	ND	0.20	EPA 8270D	10-10-16	10-12-16	
4-Chlorophenyl-phenylether	ND	0.039	EPA 8270D	10-10-16	10-12-16	
4-Nitroaniline	ND	0.039	EPA 8270D	10-10-16	10-12-16	
Fluorene	ND	0.0078	EPA 8270D/SIM	10-10-16	10-10-16	
4,6-Dinitro-2-methylphenol	ND	0.20	EPA 8270D	10-10-16	10-12-16	
n-Nitrosodiphenylamine	ND	0.039	EPA 8270D	10-10-16	10-12-16	
1,2-Diphenylhydrazine	ND	0.039	EPA 8270D	10-10-16	10-12-16	
4-Bromophenyl-phenylether	ND	0.039	EPA 8270D	10-10-16	10-12-16	
Hexachlorobenzene	ND	0.039	EPA 8270D	10-10-16	10-12-16	
Pentachlorophenol	ND	0.20	EPA 8270D	10-10-16	10-12-16	
Phenanthrene	ND	0.0078	EPA 8270D/SIM	10-10-16	10-10-16	
Anthracene	ND	0.0078	EPA 8270D/SIM	10-10-16	10-10-16	
Carbazole	ND	0.039	EPA 8270D	10-10-16	10-12-16	
Di-n-butylphthalate	ND	0.20	EPA 8270D	10-10-16	10-12-16	
Fluoranthene	ND	0.0078	EPA 8270D/SIM	10-10-16	10-10-16	
Benzidine	ND	0.39	EPA 8270D	10-10-16	10-12-16	
Pyrene	ND	0.0078	EPA 8270D/SIM	10-10-16	10-10-16	
Butylbenzylphthalate	ND	0.039	EPA 8270D	10-10-16	10-12-16	
bis-2-Ethylhexyladipate	ND	0.039	EPA 8270D	10-10-16	10-12-16	
3,3'-Dichlorobenzidine	ND	0.20	EPA 8270D	10-10-16	10-12-16	
Benzo[a]anthracene	ND	0.0078	EPA 8270D/SIM	10-10-16	10-10-16	
Chrysene	ND	0.0078	EPA 8270D/SIM	10-10-16	10-10-16	
bis(2-Ethylhexyl)phthalate	ND	0.039	EPA 8270D	10-10-16	10-12-16	
Di-n-octylphthalate	ND	0.039	EPA 8270D	10-10-16	10-12-16	
Benzo[b]fluoranthene	ND	0.0078	EPA 8270D/SIM	10-10-16	10-10-16	
Benzo(j,k)fluoranthene	ND	0.0078	EPA 8270D/SIM	10-10-16	10-10-16	
Benzo[a]pyrene	ND	0.0078	EPA 8270D/SIM	10-10-16	10-10-16	
Indeno[1,2,3-cd]pyrene	ND	0.0078	EPA 8270D/SIM	10-10-16	10-10-16	
Dibenz[a,h]anthracene	ND	0.0078	EPA 8270D/SIM	10-10-16	10-10-16	
Benzo[g,h,i]perylene	ND	0.0078	EPA 8270D/SIM	10-10-16	10-10-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorophenol</i>	<i>69</i>	<i>24 - 117</i>				
<i>Phenol-d6</i>	<i>71</i>	<i>30 - 120</i>				
<i>Nitrobenzene-d5</i>	<i>70</i>	<i>27 - 112</i>				
<i>2-Fluorobiphenyl</i>	<i>71</i>	<i>35 - 113</i>				
<i>2,4,6-Tribromophenol</i>	<i>77</i>	<i>21 - 120</i>				
<i>Terphenyl-d14</i>	<i>73</i>	<i>39 - 121</i>				



Date of Report: October 17, 2016  
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 Project: 31008

**SEMIVOLATILES EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
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Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1010S1					
n-Nitrosodimethylamine	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Pyridine	ND	0.33	EPA 8270D	10-10-16	10-11-16	
Phenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Aniline	ND	0.17	EPA 8270D	10-10-16	10-11-16	
bis(2-Chloroethyl)ether	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2-Chlorophenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
1,3-Dichlorobenzene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
1,4-Dichlorobenzene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Benzyl alcohol	ND	0.17	EPA 8270D	10-10-16	10-11-16	
1,2-Dichlorobenzene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2-Methylphenol (o-Cresol)	ND	0.033	EPA 8270D	10-10-16	10-11-16	
bis(2-Chloroisopropyl)ether	ND	0.033	EPA 8270D	10-10-16	10-11-16	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.033	EPA 8270D	10-10-16	10-11-16	
n-Nitroso-di-n-propylamine	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Hexachloroethane	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Nitrobenzene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Isophorone	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2-Nitrophenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2,4-Dimethylphenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
bis(2-Chloroethoxy)methane	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2,4-Dichlorophenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
1,2,4-Trichlorobenzene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Naphthalene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
4-Chloroaniline	ND	0.17	EPA 8270D	10-10-16	10-11-16	
Hexachlorobutadiene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
4-Chloro-3-methylphenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2-Methylnaphthalene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
1-Methylnaphthalene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Hexachlorocyclopentadiene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2,4,6-Trichlorophenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2,3-Dichloroaniline	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2,4,5-Trichlorophenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2-Chloronaphthalene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2-Nitroaniline	ND	0.033	EPA 8270D	10-10-16	10-11-16	
1,4-Dinitrobenzene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Dimethylphthalate	ND	0.033	EPA 8270D	10-10-16	10-11-16	
1,3-Dinitrobenzene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2,6-Dinitrotoluene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
1,2-Dinitrobenzene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Acenaphthylene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
3-Nitroaniline	ND	0.033	EPA 8270D	10-10-16	10-11-16	



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**SEMIVOLATILES EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1010S1					
2,4-Dinitrophenol	ND	0.17	EPA 8270D	10-10-16	10-11-16	
Acenaphthene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
4-Nitrophenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2,4-Dinitrotoluene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Dibenzofuran	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2,3,5,6-Tetrachlorophenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
2,3,4,6-Tetrachlorophenol	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Diethylphthalate	ND	0.17	EPA 8270D	10-10-16	10-11-16	
4-Chlorophenyl-phenylether	ND	0.033	EPA 8270D	10-10-16	10-11-16	
4-Nitroaniline	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Fluorene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
4,6-Dinitro-2-methylphenol	ND	0.17	EPA 8270D	10-10-16	10-11-16	
n-Nitrosodiphenylamine	ND	0.033	EPA 8270D	10-10-16	10-11-16	
1,2-Diphenylhydrazine	ND	0.033	EPA 8270D	10-10-16	10-11-16	
4-Bromophenyl-phenylether	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Hexachlorobenzene	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Pentachlorophenol	ND	0.17	EPA 8270D	10-10-16	10-11-16	
Phenanthrene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Anthracene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Carbazole	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Di-n-butylphthalate	ND	0.17	EPA 8270D	10-10-16	10-11-16	
Fluoranthene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Benzidine	ND	0.33	EPA 8270D	10-10-16	10-11-16	
Pyrene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Butylbenzylphthalate	ND	0.033	EPA 8270D	10-10-16	10-11-16	
bis-2-Ethylhexyladipate	ND	0.033	EPA 8270D	10-10-16	10-11-16	
3,3'-Dichlorobenzidine	ND	0.17	EPA 8270D	10-10-16	10-11-16	
Benzo[a]anthracene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Chrysene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
bis(2-Ethylhexyl)phthalate	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Di-n-octylphthalate	ND	0.033	EPA 8270D	10-10-16	10-11-16	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Benzo[a]pyrene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Indeno[1,2,3-cd]pyrene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
Benzo[g,h,i]perylene	ND	0.0067	EPA 8270D/SIM	10-10-16	10-10-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorophenol</i>	<i>81</i>	<i>24 - 117</i>				
<i>Phenol-d6</i>	<i>83</i>	<i>30 - 120</i>				
<i>Nitrobenzene-d5</i>	<i>85</i>	<i>27 - 112</i>				
<i>2-Fluorobiphenyl</i>	<i>84</i>	<i>35 - 113</i>				
<i>2,4,6-Tribromophenol</i>	<i>85</i>	<i>21 - 120</i>				
<i>Terphenyl-d14</i>	<i>85</i>	<i>39 - 121</i>				



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**SEMIVOLATILES EPA 8270D/SIM  
 MS/MSD QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result		Spike Level		Source	Percent		Recovery	RPD	RPD	Flags
	MS	MSD	MS	MSD	Result	Recovery	Limits	RPD	Limit		
<b>MATRIX SPIKES</b>											
Laboratory ID:	10-080-03										
	MS	MSD	MS	MSD		MS	MSD				
Phenol	<b>0.914</b>	<b>0.997</b>	1.33	1.33	ND	69	75	31 - 108	9	36	
2-Chlorophenol	<b>0.940</b>	<b>1.03</b>	1.33	1.33	ND	71	77	38 - 103	9	38	
1,4-Dichlorobenzene	<b>0.465</b>	<b>0.502</b>	0.667	0.667	ND	70	75	25 - 101	8	40	
n-Nitroso-di-n-propylamine	<b>0.452</b>	<b>0.494</b>	0.667	0.667	ND	68	74	26 - 102	9	38	
1,2,4-Trichlorobenzene	<b>0.451</b>	<b>0.513</b>	0.667	0.667	ND	68	77	27 - 101	13	40	
4-Chloro-3-methylphenol	<b>0.914</b>	<b>1.02</b>	1.33	1.33	ND	69	77	42 - 106	11	29	
Acenaphthene	<b>0.460</b>	<b>0.513</b>	0.667	0.667	ND	69	77	42 - 103	11	30	
4-Nitrophenol	<b>0.927</b>	<b>1.02</b>	1.33	1.33	ND	70	77	25 - 125	10	29	
2,4-Dinitrotoluene	<b>0.466</b>	<b>0.507</b>	0.667	0.667	ND	70	76	45 - 107	8	30	
Pentachlorophenol	<b>1.03</b>	<b>1.15</b>	1.33	1.33	ND	77	86	30 - 103	11	31	
Pyrene	<b>0.478</b>	<b>0.532</b>	0.667	0.667	ND	72	80	50 - 118	11	28	
<i>Surrogate:</i>											
2-Fluorophenol						69	75	24 - 117			
Phenol-d6						71	77	30 - 120			
Nitrobenzene-d5						72	79	27 - 112			
2-Fluorobiphenyl						71	76	35 - 113			
2,4,6-Tribromophenol						75	81	21 - 120			
Terphenyl-d14						72	79	39 - 121			



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**PCBs**  
**EPA 8082A**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-3-16-3</b>					
Laboratory ID:	10-080-01					
Aroclor 1016	ND	0.055	EPA 8082A	10-11-16	10-11-16	
Aroclor 1221	ND	0.055	EPA 8082A	10-11-16	10-11-16	
Aroclor 1232	ND	0.055	EPA 8082A	10-11-16	10-11-16	
Aroclor 1242	ND	0.055	EPA 8082A	10-11-16	10-11-16	
Aroclor 1248	ND	0.055	EPA 8082A	10-11-16	10-11-16	
Aroclor 1254	ND	0.055	EPA 8082A	10-11-16	10-11-16	
Aroclor 1260	ND	0.055	EPA 8082A	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	82	50-139				
<b>Client ID:</b>	<b>H-3-16-6</b>					
Laboratory ID:	10-080-02					
Aroclor 1016	ND	0.057	EPA 8082A	10-11-16	10-11-16	
Aroclor 1221	ND	0.057	EPA 8082A	10-11-16	10-11-16	
Aroclor 1232	ND	0.057	EPA 8082A	10-11-16	10-11-16	
Aroclor 1242	ND	0.057	EPA 8082A	10-11-16	10-11-16	
Aroclor 1248	ND	0.057	EPA 8082A	10-11-16	10-11-16	
Aroclor 1254	ND	0.057	EPA 8082A	10-11-16	10-11-16	
Aroclor 1260	ND	0.057	EPA 8082A	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	77	50-139				
<b>Client ID:</b>	<b>H-3-16-8.5</b>					
Laboratory ID:	10-080-03					
Aroclor 1016	ND	0.059	EPA 8082A	10-11-16	10-11-16	
Aroclor 1221	ND	0.059	EPA 8082A	10-11-16	10-11-16	
Aroclor 1232	ND	0.059	EPA 8082A	10-11-16	10-11-16	
Aroclor 1242	ND	0.059	EPA 8082A	10-11-16	10-11-16	
Aroclor 1248	ND	0.059	EPA 8082A	10-11-16	10-11-16	
Aroclor 1254	ND	0.059	EPA 8082A	10-11-16	10-11-16	
Aroclor 1260	ND	0.059	EPA 8082A	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	70	50-139				



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**PCBs EPA 8082A  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1011S1					
Aroclor 1016	ND	0.050	EPA 8082A	10-11-16	10-11-16	
Aroclor 1221	ND	0.050	EPA 8082A	10-11-16	10-11-16	
Aroclor 1232	ND	0.050	EPA 8082A	10-11-16	10-11-16	
Aroclor 1242	ND	0.050	EPA 8082A	10-11-16	10-11-16	
Aroclor 1248	ND	0.050	EPA 8082A	10-11-16	10-11-16	
Aroclor 1254	ND	0.050	EPA 8082A	10-11-16	10-11-16	
Aroclor 1260	ND	0.050	EPA 8082A	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
DCB	86		50-139			

Analyte	Result		Spike Level		Source Result	Percent Recovery		Recovery Limits	RPD	RPD Limit	Flags
<b>MATRIX SPIKES</b>											
Laboratory ID:	10-080-01										
	MS	MSD	MS	MSD		MS	MSD				
Aroclor 1260	0.389	0.387	0.500	0.500	ND	78	77	49-133	1	17	
<i>Surrogate:</i>											
DCB						80	81	50-139			



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 Project: 31008

**TOTAL METALS**  
**EPA 6010C/6020A/7471B**

Matrix: Soil  
 Units: mg/kg (ppm)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>EPA Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
Lab ID:	10-080-01					
<b>Client ID:</b>	<b>H-3-16-3</b>					
Antimony	<b>ND</b>	5.5	6010C	10-13-16	10-13-16	
Arsenic	<b>ND</b>	11	6010C	10-13-16	10-13-16	
Beryllium	<b>ND</b>	0.55	6010C	10-13-16	10-13-16	
Cadmium	<b>ND</b>	0.55	6010C	10-13-16	10-13-16	
Chromium	<b>28</b>	0.55	6010C	10-13-16	10-13-16	
Copper	<b>13</b>	1.1	6010C	10-13-16	10-13-16	
Lead	<b>ND</b>	5.5	6010C	10-13-16	10-13-16	
Mercury	<b>ND</b>	0.28	7471B	10-11-16	10-11-16	
Nickel	<b>33</b>	2.8	6010C	10-13-16	10-13-16	
Selenium	<b>ND</b>	11	6010C	10-13-16	10-13-16	
Silver	<b>ND</b>	0.55	6010C	10-13-16	10-13-16	
Thallium	<b>ND</b>	1.4	6020A	10-13-16	10-17-16	
Zinc	<b>26</b>	2.8	6010C	10-13-16	10-13-16	



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**TOTAL METALS**  
**EPA 6010C/6020A/7471B**

Matrix: Soil  
 Units: mg/kg (ppm)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>EPA Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
Lab ID:	10-080-02					
Client ID:	H-3-16-6					
Antimony	ND	5.7	6010C	10-13-16	10-13-16	
Arsenic	ND	11	6010C	10-13-16	10-13-16	
Beryllium	ND	0.57	6010C	10-13-16	10-13-16	
Cadmium	ND	0.57	6010C	10-13-16	10-13-16	
Chromium	27	0.57	6010C	10-13-16	10-13-16	
Copper	11	1.1	6010C	10-13-16	10-13-16	
Lead	ND	5.7	6010C	10-13-16	10-13-16	
Mercury	ND	0.29	7471B	10-11-16	10-11-16	
Nickel	30	2.9	6010C	10-13-16	10-13-16	
Selenium	ND	11	6010C	10-13-16	10-13-16	
Silver	ND	0.57	6010C	10-13-16	10-13-16	
Thallium	ND	1.4	6020A	10-13-16	10-17-16	
Zinc	24	2.9	6010C	10-13-16	10-13-16	



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 Laboratory Reference: 1610-080  
 Project: 31008

**TOTAL METALS**  
**EPA 6010C/6020A/7471B**

Matrix: Soil  
 Units: mg/kg (ppm)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>EPA Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
Lab ID:	10-080-03					
Client ID:	H-3-16-8.5					
Antimony	ND	5.9	6010C	10-13-16	10-13-16	
Arsenic	ND	12	6010C	10-13-16	10-13-16	
Beryllium	ND	0.59	6010C	10-13-16	10-13-16	
Cadmium	ND	0.59	6010C	10-13-16	10-13-16	
Chromium	29	0.59	6010C	10-13-16	10-13-16	
Copper	8.6	1.2	6010C	10-13-16	10-13-16	
Lead	ND	5.9	6010C	10-13-16	10-13-16	
Mercury	ND	0.29	7471B	10-11-16	10-11-16	
Nickel	24	2.9	6010C	10-13-16	10-13-16	
Selenium	ND	12	6010C	10-13-16	10-13-16	
Silver	ND	0.59	6010C	10-13-16	10-13-16	
Thallium	ND	1.5	6020A	10-13-16	10-17-16	
Zinc	26	2.9	6010C	10-13-16	10-13-16	



Date of Report: October 17, 2016  
 Samples Submitted: October 8, 2016  
 Laboratory Reference: 1610-080  
 Project: 31008

**TOTAL METALS  
 EPA 6010C/6020A/7471B  
 METHOD BLANK QUALITY CONTROL**

Date Extracted: 10-11&13-16  
 Date Analyzed: 10-11,13&17-16

Matrix: Soil  
 Units: mg/kg (ppm)

Lab ID: MB1013SH1&MB1011S1

Analyte	Method	Result	PQL
Antimony	6010C	ND	5.0
Arsenic	6010C	ND	10
Beryllium	6010C	ND	0.50
Cadmium	6010C	ND	0.50
Chromium	6010C	ND	0.50
Copper	6010C	ND	1.0
Lead	6010C	ND	5.0
Mercury	7471B	ND	0.25
Nickel	6010C	ND	2.5
Selenium	6010C	ND	10
Silver	6010C	ND	0.50
Thallium	6020A	ND	1.3
Zinc	6010C	ND	2.5



Date of Report: October 17, 2016  
 Samples Submitted: October 8, 2016  
 Laboratory Reference: 1610-080  
 Project: 31008

**TOTAL METALS  
 EPA 6010C/6020A/7471B  
 DUPLICATE QUALITY CONTROL**

Date Extracted: 10-11&13-16  
 Date Analyzed: 10-11,13&17-16

Matrix: Soil  
 Units: mg/kg (ppm)

Lab ID: 10-080-01

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Antimony	ND	ND	NA	5.0	
Arsenic	ND	ND	NA	10.0	
Beryllium	ND	ND	NA	0.50	
Cadmium	ND	ND	NA	0.50	
Chromium	25.9	24.3	6	0.50	
Copper	11.7	10.9	7	1.0	
Lead	ND	ND	NA	5.0	
Mercury	ND	ND	NA	0.25	
Nickel	30.4	29.2	4	2.5	
Selenium	ND	ND	NA	10	
Silver	ND	ND	NA	0.50	
Thallium	ND	ND	NA	1.3	
Zinc	24.0	22.0	9	2.5	



Date of Report: October 17, 2016  
 Samples Submitted: October 8, 2016  
 Laboratory Reference: 1610-080  
 Project: 31008

**TOTAL METALS  
 EPA 6010C/6020A/7471B  
 MS/MSD QUALITY CONTROL**

Date Extracted: 10-11&13-16  
 Date Analyzed: 10-11,13&17-16

Matrix: Soil  
 Units: mg/kg (ppm)

Lab ID: 10-080-01

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Antimony	100	<b>95.4</b>	95	<b>88.3</b>	88	8	
Arsenic	100	<b>101</b>	101	<b>94.3</b>	94	6	
Beryllium	50.0	<b>50.8</b>	102	<b>47.6</b>	95	7	
Cadmium	50.0	<b>49.3</b>	99	<b>47.5</b>	95	4	
Chromium	100	<b>128</b>	102	<b>119</b>	93	7	
Copper	50.0	<b>63.9</b>	105	<b>60.4</b>	97	6	
Lead	250	<b>238</b>	95	<b>232</b>	93	3	
Mercury	0.500	<b>0.499</b>	100	<b>0.546</b>	109	9	
Nickel	100	<b>127</b>	96	<b>120</b>	89	6	
Selenium	100	<b>104</b>	104	<b>99.4</b>	99	5	
Silver	25.0	<b>24.4</b>	97	<b>23.1</b>	92	5	
Thallium	50.0	<b>44.2</b>	88	<b>44.8</b>	90	1	
Zinc	100	<b>120</b>	96	<b>115</b>	91	4	



Date of Report: October 17, 2016  
Samples Submitted: October 8, 2016  
Laboratory Reference: 1610-080  
Project: 31008

### % MOISTURE

Date Analyzed: 10-10-16

Client ID	Lab ID	% Moisture
H-3-16-3	10-080-01	9
H-3-16-6	10-080-02	12
H-3-16-8.5	10-080-03	15





### 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.
  - Z -
- ND - Not Detected at PQL  
 PQL - Practical Quantitation Limit  
 RPD - Relative Percent Difference





# Onsite Environmental Inc.

Analytical Laboratory Testing Services  
14648 NE 95th Street • Redmond, WA 98052  
Phone: (425) 883-3881 • www.onsite-env.com

## Chain of Custody

### Turnaround Request (in working days)

(Check One)

- Same Day  1 Day
- 2 Days  3 Days
- Standard (7 Days)  
(TPH analysis 5 Days)
- \_\_\_\_\_ (other)

### Laboratory Number:

10-080

Company: INNOVEX

Project Number: 31008

Project Name: SR520

Project Manager: Glenn Hayman

Sampled by: Andrew J. Sinden

Lab ID	Sample Identification	Date Sampled	Time Sampled	Matrix	Number of Containers																						
1	H-3-16-3	10/7/16	2325	S	5	NWTPH-HCID	NWTPH-Gx/BTEX	NWTPH-Gx*	NWTPH-Dx ( <input type="checkbox"/> Acid / SG Clean-up )	<input checked="" type="checkbox"/>	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	PP-13 metals 6010	% Moisture		
2	H-3-16-6		2330																								
3	H-3-16-8.5		2335																								
4	H-3-16-13.5		2350																								
5	H-3-16-17.4		2450																								

*Signature* 10/7/16

Signature \_\_\_\_\_ Company \_\_\_\_\_ Date \_\_\_\_\_ Time \_\_\_\_\_

Comments/Special Instructions  
\* Follow up w/ appropriate analysis based on HCTD

Relinquished  
Received  
Relinquished  
Received  
Relinquished  
Received  
Reviewed/Date

Signature  
Company  
Date  
Time  
Comments/Special Instructions  
Data Package: Standard  Level III  Level IV   
Chromatograms with final report  Electronic Data Deliverables (EDDs)



14648 NE 95<sup>th</sup> Street, Redmond, WA 98052 • (425) 883-3881

October 18, 2016

Glenn Hayman  
INNOVEX Environmental Mgt., Inc.  
16310 NE 80th St., Suite 300  
Redmond, WA 98052

Re: Analytical Data for Project 31008  
Laboratory Reference No. 1610-081

Dear Glenn:

Enclosed are the analytical results and associated quality control data for samples submitted on October 8, 2016.

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,

A handwritten signature in black ink, appearing to read "D. Baumeister", with a long horizontal flourish extending to the right.

David Baumeister  
Project Manager

Enclosures



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OnSite Environmental, Inc. 14648 NE 95<sup>th</sup> Street, Redmond, WA 98052 (425) 883-3881

This report pertains to the samples analyzed in accordance with the chain of custody, and is intended only for the use of the individual or company to whom it is addressed.

Date of Report: October 18, 2016  
Samples Submitted: October 8, 2016  
Laboratory Reference: 1610-081  
Project: 31008

### Case Narrative

Samples were collected on October 7, 2016 and received by the laboratory on October 8, 2016. 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.

#### PCBs EPA 8082A Analysis

Due to limited sample volume, H-3-16 was extracted from a 500 mL poly bottle.

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.



Date of Report: October 18, 2016  
 Samples Submitted: October 8, 2016  
 Laboratory Reference: 1610-081  
 Project: 31008

### NWTPH-HCID

Matrix: Water  
 Units: mg/L (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-3-16</b>					
Laboratory ID:	10-081-01					
Gasoline Range Organics	<b>ND</b>	0.11	NWTPH-HCID	10-12-16	10-12-16	
Diesel Range Organics	<b>ND</b>	0.29	NWTPH-HCID	10-12-16	10-12-16	
Lube Oil Range Organics	<b>ND</b>	0.46	NWTPH-HCID	10-12-16	10-12-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>112</i>	<i>50-150</i>				



Date of Report: October 18, 2016  
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**NWTPH-HCID  
 QUALITY CONTROL**

Matrix: Water  
 Units: mg/L (ppm)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>METHOD BLANK</b>						
Laboratory ID:	MB1012W1					
Gasoline Range Organics	<b>ND</b>	0.10	NWTPH-HCID	10-12-16	10-12-16	
Diesel Range Organics	<b>ND</b>	0.25	NWTPH-HCID	10-12-16	10-12-16	
Lube Oil Range Organics	<b>ND</b>	0.40	NWTPH-HCID	10-12-16	10-12-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>94</i>	<i>50-150</i>				



Date of Report: October 18, 2016  
 Samples Submitted: October 8, 2016  
 Laboratory Reference: 1610-081  
 Project: 31008

**VOLATILES EPA 8260C**  
 page 1 of 2

Matrix: Water  
 Units: ug/L

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>H-3-16</b>					
Laboratory ID:	10-081-01					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Chloromethane	ND	1.0	EPA 8260C	10-11-16	10-11-16	
Vinyl Chloride	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Bromomethane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Chloroethane	ND	1.0	EPA 8260C	10-11-16	10-11-16	
Trichlorofluoromethane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,1-Dichloroethene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Acetone	6.5	5.0	EPA 8260C	10-11-16	10-11-16	
Iodomethane	ND	1.0	EPA 8260C	10-11-16	10-11-16	
Carbon Disulfide	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Methylene Chloride	ND	1.0	EPA 8260C	10-11-16	10-11-16	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,1-Dichloroethane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Vinyl Acetate	ND	1.0	EPA 8260C	10-11-16	10-11-16	
2,2-Dichloropropane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
2-Butanone	ND	5.0	EPA 8260C	10-11-16	10-11-16	
Bromochloromethane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Chloroform	8.3	0.20	EPA 8260C	10-11-16	10-11-16	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Carbon Tetrachloride	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,1-Dichloropropene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Benzene	7.4	0.20	EPA 8260C	10-11-16	10-11-16	
1,2-Dichloroethane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Trichloroethene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,2-Dichloropropane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Dibromomethane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Bromodichloromethane	1.0	0.20	EPA 8260C	10-11-16	10-11-16	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	10-11-16	10-11-16	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	10-11-16	10-11-16	
Toluene	ND	1.0	EPA 8260C	10-11-16	10-11-16	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	10-11-16	10-11-16	



Date of Report: October 18, 2016  
 Samples Submitted: October 8, 2016  
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 Project: 31008

**VOLATILES EPA 8260C**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-3-16</b>					
Laboratory ID:	10-081-01					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Tetrachloroethene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,3-Dichloropropane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
2-Hexanone	ND	2.0	EPA 8260C	10-11-16	10-11-16	
Dibromochloromethane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,2-Dibromoethane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Chlorobenzene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Ethylbenzene	0.70	0.20	EPA 8260C	10-11-16	10-11-16	
m,p-Xylene	2.1	0.40	EPA 8260C	10-11-16	10-11-16	
o-Xylene	0.67	0.20	EPA 8260C	10-11-16	10-11-16	
Styrene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Bromoform	ND	1.0	EPA 8260C	10-11-16	10-11-16	
Isopropylbenzene	0.25	0.20	EPA 8260C	10-11-16	10-11-16	
Bromobenzene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
n-Propylbenzene	0.37	0.20	EPA 8260C	10-11-16	10-11-16	
2-Chlorotoluene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
4-Chlorotoluene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,3,5-Trimethylbenzene	0.30	0.20	EPA 8260C	10-11-16	10-11-16	
tert-Butylbenzene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,2,4-Trimethylbenzene	0.44	0.20	EPA 8260C	10-11-16	10-11-16	
sec-Butylbenzene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
p-Isopropyltoluene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
n-Butylbenzene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	10-11-16	10-11-16	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Hexachlorobutadiene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Naphthalene	ND	1.0	EPA 8260C	10-11-16	10-11-16	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>104</i>	<i>71-131</i>				
<i>Toluene-d8</i>	<i>102</i>	<i>80-127</i>				
<i>4-Bromofluorobenzene</i>	<i>100</i>	<i>80-125</i>				



Date of Report: October 18, 2016  
 Samples Submitted: October 8, 2016  
 Laboratory Reference: 1610-081  
 Project: 31008

**VOLATILES by EPA 8260C**  
**METHOD BLANK QUALITY CONTROL**  
 page 1 of 2

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1011W2					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Chloromethane	ND	1.0	EPA 8260C	10-11-16	10-11-16	
Vinyl Chloride	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Bromomethane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Chloroethane	ND	1.0	EPA 8260C	10-11-16	10-11-16	
Trichlorofluoromethane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,1-Dichloroethene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Acetone	ND	5.0	EPA 8260C	10-11-16	10-11-16	
Iodomethane	ND	1.0	EPA 8260C	10-11-16	10-11-16	
Carbon Disulfide	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Methylene Chloride	ND	1.0	EPA 8260C	10-11-16	10-11-16	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,1-Dichloroethane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Vinyl Acetate	ND	1.0	EPA 8260C	10-11-16	10-11-16	
2,2-Dichloropropane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
2-Butanone	ND	5.0	EPA 8260C	10-11-16	10-11-16	
Bromochloromethane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Chloroform	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Carbon Tetrachloride	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,1-Dichloropropene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Benzene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,2-Dichloroethane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Trichloroethene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,2-Dichloropropane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Dibromomethane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Bromodichloromethane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	10-11-16	10-11-16	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	10-11-16	10-11-16	
Toluene	ND	1.0	EPA 8260C	10-11-16	10-11-16	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	10-11-16	10-11-16	



Date of Report: October 18, 2016  
 Samples Submitted: October 8, 2016  
 Laboratory Reference: 1610-081  
 Project: 31008

**VOLATILES by EPA 8260C**  
**METHOD BLANK QUALITY CONTROL**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:		MB1011W2				
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Tetrachloroethene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,3-Dichloropropane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
2-Hexanone	ND	2.0	EPA 8260C	10-11-16	10-11-16	
Dibromochloromethane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,2-Dibromoethane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Chlorobenzene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Ethylbenzene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
m,p-Xylene	ND	0.40	EPA 8260C	10-11-16	10-11-16	
o-Xylene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Styrene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Bromoform	ND	1.0	EPA 8260C	10-11-16	10-11-16	
Isopropylbenzene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Bromobenzene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	10-11-16	10-11-16	
n-Propylbenzene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
2-Chlorotoluene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
4-Chlorotoluene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
tert-Butylbenzene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
sec-Butylbenzene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
p-Isopropyltoluene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
n-Butylbenzene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	10-11-16	10-11-16	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Hexachlorobutadiene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
Naphthalene	ND	1.0	EPA 8260C	10-11-16	10-11-16	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>102</i>	<i>71-131</i>				
<i>Toluene-d8</i>	<i>99</i>	<i>80-127</i>				
<i>4-Bromofluorobenzene</i>	<i>95</i>	<i>80-125</i>				



Date of Report: October 18, 2016  
 Samples Submitted: October 8, 2016  
 Laboratory Reference: 1610-081  
 Project: 31008

**VOLATILES by EPA 8260C  
 SB/SBD QUALITY CONTROL**

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD		Flags
					SB	SBD	Limits	RPD	Limit	
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB1011W2									
	SB	SBD	SB	SBD	SB	SBD				
1,1-Dichloroethene	9.76	9.64	10.0	10.0	98	96	62-132	1	20	
Benzene	9.37	9.66	10.0	10.0	94	97	75-121	3	15	
Trichloroethene	8.72	8.92	10.0	10.0	87	89	65-115	2	15	
Toluene	8.96	9.43	10.0	10.0	90	94	78-120	5	15	
Chlorobenzene	8.98	9.32	10.0	10.0	90	93	77-118	4	15	
<i>Surrogate:</i>										
Dibromofluoromethane					104	102	71-131			
Toluene-d8					100	100	80-127			
4-Bromofluorobenzene					97	98	80-125			



Date of Report: October 18, 2016  
 Samples Submitted: October 8, 2016  
 Laboratory Reference: 1610-081  
 Project: 31008

**SEMIVOLATILES EPA 8270D/SIM**  
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Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-3-16</b>					
<b>Laboratory ID:</b>	<b>10-081-01</b>					
n-Nitrosodimethylamine	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Pyridine	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Phenol	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Aniline	ND	5.0	EPA 8270D	10-13-16	10-13-16	
bis(2-Chloroethyl)ether	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2-Chlorophenol	ND	1.0	EPA 8270D	10-13-16	10-13-16	
1,3-Dichlorobenzene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
1,4-Dichlorobenzene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Benzyl alcohol	ND	1.0	EPA 8270D	10-13-16	10-13-16	
1,2-Dichlorobenzene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2-Methylphenol (o-Cresol)	ND	1.0	EPA 8270D	10-13-16	10-13-16	
bis(2-Chloroisopropyl)ether	ND	1.0	EPA 8270D	10-13-16	10-13-16	
(3+4)-Methylphenol (m,p-Cresol)	ND	1.0	EPA 8270D	10-13-16	10-13-16	
n-Nitroso-di-n-propylamine	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Hexachloroethane	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Nitrobenzene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Isophorone	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2-Nitrophenol	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2,4-Dimethylphenol	ND	1.0	EPA 8270D	10-13-16	10-13-16	
bis(2-Chloroethoxy)methane	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2,4-Dichlorophenol	ND	1.0	EPA 8270D	10-13-16	10-13-16	
1,2,4-Trichlorobenzene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Naphthalene	0.25	0.10	EPA 8270D/SIM	10-13-16	10-13-16	
4-Chloroaniline	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Hexachlorobutadiene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
4-Chloro-3-methylphenol	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2-Methylnaphthalene	0.23	0.10	EPA 8270D/SIM	10-13-16	10-13-16	
1-Methylnaphthalene	0.13	0.10	EPA 8270D/SIM	10-13-16	10-13-16	
Hexachlorocyclopentadiene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2,4,6-Trichlorophenol	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2,3-Dichloroaniline	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2,4,5-Trichlorophenol	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2-Chloronaphthalene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2-Nitroaniline	ND	1.0	EPA 8270D	10-13-16	10-13-16	
1,4-Dinitrobenzene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Dimethylphthalate	ND	1.0	EPA 8270D	10-13-16	10-13-16	
1,3-Dinitrobenzene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2,6-Dinitrotoluene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
1,2-Dinitrobenzene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Acenaphthylene	ND	0.10	EPA 8270D/SIM	10-13-16	10-13-16	
3-Nitroaniline	ND	1.0	EPA 8270D	10-13-16	10-13-16	



Date of Report: October 18, 2016  
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**SEMIVOLATILES EPA 8270D/SIM**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-3-16</b>					
Laboratory ID:	10-081-01					
2,4-Dinitrophenol	ND	5.0	EPA 8270D	10-13-16	10-13-16	
Acenaphthene	ND	0.10	EPA 8270D/SIM	10-13-16	10-13-16	
4-Nitrophenol	ND	5.0	EPA 8270D	10-13-16	10-13-16	
2,4-Dinitrotoluene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Dibenzofuran	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2,3,5,6-Tetrachlorophenol	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2,3,4,6-Tetrachlorophenol	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Diethylphthalate	1.3	1.0	EPA 8270D	10-13-16	10-13-16	
4-Chlorophenyl-phenylether	ND	1.0	EPA 8270D	10-13-16	10-13-16	
4-Nitroaniline	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Fluorene	ND	0.10	EPA 8270D/SIM	10-13-16	10-13-16	
4,6-Dinitro-2-methylphenol	ND	5.0	EPA 8270D	10-13-16	10-13-16	
n-Nitrosodiphenylamine	ND	1.0	EPA 8270D	10-13-16	10-13-16	
1,2-Diphenylhydrazine	ND	1.0	EPA 8270D	10-13-16	10-13-16	
4-Bromophenyl-phenylether	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Hexachlorobenzene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Pentachlorophenol	ND	5.0	EPA 8270D	10-13-16	10-13-16	
Phenanthrene	0.11	0.10	EPA 8270D/SIM	10-13-16	10-13-16	
Anthracene	ND	0.10	EPA 8270D/SIM	10-13-16	10-13-16	
Carbazole	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Di-n-butylphthalate	9.3	1.0	EPA 8270D	10-13-16	10-13-16	
Fluoranthene	ND	0.10	EPA 8270D/SIM	10-13-16	10-13-16	
Benzidine	ND	5.0	EPA 8270D	10-13-16	10-13-16	
Pyrene	ND	0.10	EPA 8270D/SIM	10-13-16	10-13-16	
Butylbenzylphthalate	ND	1.0	EPA 8270D	10-13-16	10-13-16	
bis-2-Ethylhexyladipate	ND	1.0	EPA 8270D	10-13-16	10-13-16	
3,3'-Dichlorobenzidine	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Benzo[a]anthracene	0.024	0.010	EPA 8270D/SIM	10-13-16	10-13-16	
Chrysene	0.012	0.010	EPA 8270D/SIM	10-13-16	10-13-16	
bis(2-Ethylhexyl)phthalate	5.7	1.0	EPA 8270D	10-13-16	10-13-16	
Di-n-octylphthalate	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Benzo[b]fluoranthene	0.082	0.010	EPA 8270D/SIM	10-13-16	10-13-16	
Benzo(j,k)fluoranthene	0.023	0.010	EPA 8270D/SIM	10-13-16	10-13-16	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	10-13-16	10-13-16	
Indeno[1,2,3-cd]pyrene	ND	0.010	EPA 8270D/SIM	10-13-16	10-13-16	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	10-13-16	10-13-16	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270D/SIM	10-13-16	10-13-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	37	19 - 87				
Phenol-d6	39	10 - 83				
Nitrobenzene-d5	65	35 - 112				
2-Fluorobiphenyl	59	45 - 112				
2,4,6-Tribromophenol	39	37 - 115				
Terphenyl-d14	61	49 - 126				



Date of Report: October 18, 2016  
 Samples Submitted: October 8, 2016  
 Laboratory Reference: 1610-081  
 Project: 31008

**SEMIVOLATILES EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
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Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1013W1					
n-Nitrosodimethylamine	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Pyridine	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Phenol	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Aniline	ND	5.0	EPA 8270D	10-13-16	10-13-16	
bis(2-Chloroethyl)ether	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2-Chlorophenol	ND	1.0	EPA 8270D	10-13-16	10-13-16	
1,3-Dichlorobenzene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
1,4-Dichlorobenzene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Benzyl alcohol	ND	1.0	EPA 8270D	10-13-16	10-13-16	
1,2-Dichlorobenzene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2-Methylphenol (o-Cresol)	ND	1.0	EPA 8270D	10-13-16	10-13-16	
bis(2-Chloroisopropyl)ether	ND	1.0	EPA 8270D	10-13-16	10-13-16	
(3+4)-Methylphenol (m,p-Cresol)	ND	1.0	EPA 8270D	10-13-16	10-13-16	
n-Nitroso-di-n-propylamine	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Hexachloroethane	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Nitrobenzene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Isophorone	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2-Nitrophenol	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2,4-Dimethylphenol	ND	1.0	EPA 8270D	10-13-16	10-13-16	
bis(2-Chloroethoxy)methane	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2,4-Dichlorophenol	ND	1.0	EPA 8270D	10-13-16	10-13-16	
1,2,4-Trichlorobenzene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Naphthalene	ND	0.10	EPA 8270D/SIM	10-13-16	10-13-16	
4-Chloroaniline	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Hexachlorobutadiene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
4-Chloro-3-methylphenol	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2-Methylnaphthalene	ND	0.10	EPA 8270D/SIM	10-13-16	10-13-16	
1-Methylnaphthalene	ND	0.10	EPA 8270D/SIM	10-13-16	10-13-16	
Hexachlorocyclopentadiene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2,4,6-Trichlorophenol	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2,3-Dichloroaniline	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2,4,5-Trichlorophenol	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2-Chloronaphthalene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2-Nitroaniline	ND	1.0	EPA 8270D	10-13-16	10-13-16	
1,4-Dinitrobenzene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Dimethylphthalate	ND	1.0	EPA 8270D	10-13-16	10-13-16	
1,3-Dinitrobenzene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2,6-Dinitrotoluene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
1,2-Dinitrobenzene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Acenaphthylene	ND	0.10	EPA 8270D/SIM	10-13-16	10-13-16	
3-Nitroaniline	ND	1.0	EPA 8270D	10-13-16	10-13-16	



Date of Report: October 18, 2016  
 Samples Submitted: October 8, 2016  
 Laboratory Reference: 1610-081  
 Project: 31008

**SEMIVOLATILES EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1013W1					
2,4-Dinitrophenol	ND	5.0	EPA 8270D	10-13-16	10-13-16	
Acenaphthene	ND	0.10	EPA 8270D/SIM	10-13-16	10-13-16	
4-Nitrophenol	ND	5.0	EPA 8270D	10-13-16	10-13-16	
2,4-Dinitrotoluene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Dibenzofuran	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2,3,5,6-Tetrachlorophenol	ND	1.0	EPA 8270D	10-13-16	10-13-16	
2,3,4,6-Tetrachlorophenol	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Diethylphthalate	ND	1.0	EPA 8270D	10-13-16	10-13-16	
4-Chlorophenyl-phenylether	ND	1.0	EPA 8270D	10-13-16	10-13-16	
4-Nitroaniline	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Fluorene	ND	0.10	EPA 8270D/SIM	10-13-16	10-13-16	
4,6-Dinitro-2-methylphenol	ND	5.0	EPA 8270D	10-13-16	10-13-16	
n-Nitrosodiphenylamine	ND	1.0	EPA 8270D	10-13-16	10-13-16	
1,2-Diphenylhydrazine	ND	1.0	EPA 8270D	10-13-16	10-13-16	
4-Bromophenyl-phenylether	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Hexachlorobenzene	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Pentachlorophenol	ND	5.0	EPA 8270D	10-13-16	10-13-16	
Phenanthrene	ND	0.10	EPA 8270D/SIM	10-13-16	10-13-16	
Anthracene	ND	0.10	EPA 8270D/SIM	10-13-16	10-13-16	
Carbazole	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Di-n-butylphthalate	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Fluoranthene	ND	0.10	EPA 8270D/SIM	10-13-16	10-13-16	
Benzidine	ND	5.0	EPA 8270D	10-13-16	10-13-16	
Pyrene	ND	0.10	EPA 8270D/SIM	10-13-16	10-13-16	
Butylbenzylphthalate	ND	1.0	EPA 8270D	10-13-16	10-13-16	
bis-2-Ethylhexyladipate	ND	1.0	EPA 8270D	10-13-16	10-13-16	
3,3'-Dichlorobenzidine	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	10-13-16	10-13-16	
Chrysene	ND	0.010	EPA 8270D/SIM	10-13-16	10-13-16	
bis(2-Ethylhexyl)phthalate	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Di-n-octylphthalate	ND	1.0	EPA 8270D	10-13-16	10-13-16	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	10-13-16	10-13-16	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	10-13-16	10-13-16	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	10-13-16	10-13-16	
Indeno[1,2,3-cd]pyrene	ND	0.010	EPA 8270D/SIM	10-13-16	10-13-16	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	10-13-16	10-13-16	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270D/SIM	10-13-16	10-13-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	50	19 - 87				
Phenol-d6	38	10 - 83				
Nitrobenzene-d5	79	35 - 112				
2-Fluorobiphenyl	75	45 - 112				
2,4,6-Tribromophenol	85	37 - 115				
Terphenyl-d14	85	49 - 126				



Date of Report: October 18, 2016  
 Samples Submitted: October 8, 2016  
 Laboratory Reference: 1610-081  
 Project: 31008

**SEMIVOLATILES EPA 8270D/SIM  
 SB/SBD QUALITY CONTROL**

Matrix: Water  
 Units: ug/L

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD	RPD	Flags
					Recovery	Limits	Limits	Limit		
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB1013W1									
	SB	SBD	SB	SBD	SB	SBD				
Phenol	17.1	17.4	40.0	40.0	43	44	25 - 70	2	32	
2-Chlorophenol	30.8	31.8	40.0	40.0	77	80	55 - 99	3	27	
1,4-Dichlorobenzene	14.5	14.5	20.0	20.0	73	73	48 - 93	0	30	
n-Nitroso-di-n-propylamine	15.3	15.4	20.0	20.0	77	77	47 - 108	1	26	
1,2,4-Trichlorobenzene	14.8	15.1	20.0	20.0	74	76	52 - 94	2	24	
4-Chloro-3-methylphenol	32.1	32.7	40.0	40.0	80	82	67 - 108	2	16	
Acenaphthene	16.2	15.5	20.0	20.0	81	78	50 - 113	4	17	
4-Nitrophenol	16.9	16.4	40.0	40.0	42	41	29 - 78	3	37	
2,4-Dinitrotoluene	14.8	14.2	20.0	20.0	74	71	64 - 107	4	19	
Pentachlorophenol	26.5	27.3	40.0	40.0	66	68	35 - 116	3	25	
Pyrene	17.0	16.8	20.0	20.0	85	84	61 - 112	1	15	
<i>Surrogate:</i>										
2-Fluorophenol					56	57	19 - 87			
Phenol-d6					43	44	10 - 83			
Nitrobenzene-d5					84	86	35 - 112			
2-Fluorobiphenyl					80	79	45 - 112			
2,4,6-Tribromophenol					86	87	37 - 115			
Terphenyl-d14					89	86	49 - 126			



Date of Report: October 18, 2016  
 Samples Submitted: October 8, 2016  
 Laboratory Reference: 1610-081  
 Project: 31008

**PCBs**  
**EPA 8082A**

Matrix: Water  
 Units: ug/L (ppb)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>H-3-16</b>					
Laboratory ID:	10-081-01					
Aroclor 1016	<b>ND</b>	0.045	EPA 8082A	10-10-16	10-10-16	
Aroclor 1221	<b>ND</b>	0.045	EPA 8082A	10-10-16	10-10-16	
Aroclor 1232	<b>ND</b>	0.045	EPA 8082A	10-10-16	10-10-16	
Aroclor 1242	<b>ND</b>	0.045	EPA 8082A	10-10-16	10-10-16	
Aroclor 1248	<b>ND</b>	0.045	EPA 8082A	10-10-16	10-10-16	
Aroclor 1254	<b>ND</b>	0.045	EPA 8082A	10-10-16	10-10-16	
Aroclor 1260	<b>ND</b>	0.045	EPA 8082A	10-10-16	10-10-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	79	38-137				



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**PCBs EPA 8082A  
 QUALITY CONTROL**

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1010W1					
Aroclor 1016	<b>ND</b>	0.020	EPA 8082A	10-10-16	10-10-16	
Aroclor 1221	<b>ND</b>	0.020	EPA 8082A	10-10-16	10-10-16	
Aroclor 1232	<b>ND</b>	0.020	EPA 8082A	10-10-16	10-10-16	
Aroclor 1242	<b>ND</b>	0.020	EPA 8082A	10-10-16	10-10-16	
Aroclor 1248	<b>ND</b>	0.020	EPA 8082A	10-10-16	10-10-16	
Aroclor 1254	<b>ND</b>	0.020	EPA 8082A	10-10-16	10-10-16	
Aroclor 1260	<b>ND</b>	0.020	EPA 8082A	10-10-16	10-10-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
DCB	84		38-137			

Analyte	Result		Spike Level		Source Result	Percent Recovery		Recovery Limits	RPD	RPD Limit	Flags
<b>SPIKE BLANKS</b>											
Laboratory ID:	SB1010W1										
	SB	SBD	SB	SBD		SB	SBD				
Aroclor 1260	<b>0.382</b>	<b>0.406</b>	0.500	0.500	N/A	<b>76</b>	<b>81</b>	68-114	6	12	
<i>Surrogate:</i>											
DCB						74	86	38-137			



Date of Report: October 18, 2016  
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**DISSOLVED METALS**  
**EPA 200.8/7470A**

Matrix: Water  
 Units: ug/L (ppb)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>EPA Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
Lab ID:	10-081-01					
<b>Client ID:</b>	<b>H-3-16</b>					
Antimony	<b>18</b>	5.0	200.8	10-10-16	10-12-16	
Arsenic	<b>3.3</b>	3.0	200.8	10-10-16	10-12-16	
Beryllium	<b>ND</b>	10	200.8	10-10-16	10-12-16	
Cadmium	<b>ND</b>	4.0	200.8	10-10-16	10-12-16	
Chromium	<b>ND</b>	10	200.8	10-10-16	10-12-16	
Copper	<b>ND</b>	10	200.8	10-10-16	10-12-16	
Lead	<b>2.5</b>	1.0	200.8	10-10-16	10-12-16	
Mercury	<b>ND</b>	0.50	7470A	10-10-16	10-11-16	
Nickel	<b>ND</b>	20	200.8	10-10-16	10-12-16	
Selenium	<b>ND</b>	5.0	200.8	10-10-16	10-12-16	
Silver	<b>ND</b>	10	200.8	10-10-16	10-12-16	
Thallium	<b>ND</b>	5.0	200.8	10-10-16	10-12-16	
Zinc	<b>ND</b>	25	200.8	10-10-16	10-12-16	



Date of Report: October 18, 2016  
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 Project: 31008

**DISSOLVED METALS  
 EPA 200.8  
 METHOD BLANK QUALITY CONTROL**

Date Filtered: 10-10-16  
 Date Analyzed: 10-12-16

Matrix: Water  
 Units: ug/L (ppb)

Lab ID: MB1010F1

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



Date of Report: October 18, 2016  
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Laboratory Reference: 1610-081  
Project: 31008

**DISSOLVED MERCURY  
EPA 7470A  
METHOD BLANK QUALITY CONTROL**

Date Filtered: 10-10-16

Date Analyzed: 10-11-16

Matrix: Water

Units: ug/L (ppb)

Lab ID: MB1010F1

Analyte	Method	Result	PQL
Mercury	7470A	<b>ND</b>	0.50



Date of Report: October 18, 2016  
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 Laboratory Reference: 1610-081  
 Project: 31008

**DISSOLVED METALS  
 EPA 200.8  
 DUPLICATE QUALITY CONTROL**

Date Filtered: 10-10-16  
 Date Analyzed: 10-12-16

Matrix: Water  
 Units: ug/L (ppb)

Lab ID: 10-081-01

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Antimony	17.7	16.3	8	5.0	
Arsenic	3.33	3.59	8	3.0	
Beryllium	ND	ND	NA	10	
Cadmium	ND	ND	NA	4.0	
Chromium	ND	ND	NA	10	
Copper	ND	ND	NA	10	
Lead	2.46	2.47	0	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	ND	ND	NA	25	



Date of Report: October 18, 2016  
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**DISSOLVED MERCURY  
EPA 7470A  
DUPLICATE QUALITY CONTROL**

Date Filtered: 10-10-16

Date Analyzed: 10-11-16

Matrix: Water

Units: ug/L (ppb)

Lab ID: 10-051-01

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Mercury	<b>ND</b>	<b>ND</b>	NA	0.50	



Date of Report: October 18, 2016  
 Samples Submitted: October 8, 2016  
 Laboratory Reference: 1610-081  
 Project: 31008

**DISSOLVED METALS  
 EPA 200.8  
 MS/MSD QUALITY CONTROL**

Date Filtered: 10-10-16

Date Analyzed: 10-12-16

Matrix: Water

Units: ug/L (ppb)

Lab ID: 10-081-01

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Antimony	200	<b>213</b>	98	<b>231</b>	107	8	
Arsenic	200	<b>207</b>	102	<b>224</b>	110	8	
Beryllium	200	<b>200</b>	100	<b>217</b>	109	9	
Cadmium	200	<b>199</b>	99	<b>213</b>	107	7	
Chromium	200	<b>193</b>	96	<b>207</b>	104	7	
Copper	200	<b>194</b>	97	<b>208</b>	104	7	
Lead	200	<b>189</b>	93	<b>205</b>	101	8	
Nickel	200	<b>188</b>	94	<b>203</b>	102	8	
Selenium	200	<b>202</b>	101	<b>230</b>	115	13	
Silver	200	<b>182</b>	91	<b>202</b>	101	10	
Thallium	200	<b>183</b>	92	<b>197</b>	98	7	
Zinc	200	<b>211</b>	106	<b>228</b>	114	8	



Date of Report: October 18, 2016  
 Samples Submitted: October 8, 2016  
 Laboratory Reference: 1610-081  
 Project: 31008

**DISSOLVED MERCURY  
 EPA 7470A  
 MS/MSD QUALITY CONTROL**

Date Filtered: 10-10-16

Date Analyzed: 10-11-16

Matrix: Water

Units: ug/L (ppb)

Lab ID: 10-051-01

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Mercury	12.5	<b>13.5</b>	108	<b>12.7</b>	102	6	





### 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.
  - Z -
- ND - Not Detected at PQL  
 PQL - Practical Quantitation Limit  
 RPD - Relative Percent Difference







14648 NE 95<sup>th</sup> Street, Redmond, WA 98052 • (425) 883-3881

October 19, 2016

Glenn Hayman  
INNOVEX Environmental Mgt., Inc.  
16310 NE 80th St., Suite 300  
Redmond, WA 98052

Re: Analytical Data for Project 31008  
Laboratory Reference No. 1610-084

Dear Glenn:

Enclosed are the analytical results and associated quality control data for samples submitted on October 10, 2016.

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,

A handwritten signature in black ink, appearing to read "DB", with a long horizontal flourish extending to the right.

David Baumeister  
Project Manager

Enclosures



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OnSite Environmental, Inc. 14648 NE 95<sup>th</sup> Street, Redmond, WA 98052 (425) 883-3881

This report pertains to the samples analyzed in accordance with the chain of custody, and is intended only for the use of the individual or company to whom it is addressed.

Date of Report: October 19, 2016  
Samples Submitted: October 10, 2016  
Laboratory Reference: 1610-084  
Project: 31008

### Case Narrative

Samples were collected on October 8, 2016 and received by the laboratory on October 10, 2016. 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/BTEX and Volatiles EPA 8260C Analysis

Per EPA method 5035A, samples were received by the laboratory in pre-weighed 40 ml VOA vials preserved with either Methanol or Sodium Bisulfate.

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.



Date of Report: October 19, 2016  
 Samples Submitted: October 10, 2016  
 Laboratory Reference: 1610-084  
 Project: 31008

### NWTPH-HCID

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-5-16-3</b>					
Laboratory ID:	10-084-01					
Gasoline Range Organics	ND	23	NWTPH-HCID	10-11-16	10-11-16	
Diesel Range Organics	ND	57	NWTPH-HCID	10-11-16	10-11-16	
Lube Oil Range Organics	ND	120	NWTPH-HCID	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	119	50-150				
<b>Client ID:</b>	<b>H-5-16-6</b>					
Laboratory ID:	10-084-02					
Gasoline Range Organics	ND	22	NWTPH-HCID	10-11-16	10-11-16	
Diesel Range Organics	ND	56	NWTPH-HCID	10-11-16	10-11-16	
Lube Oil Range Organics	ND	110	NWTPH-HCID	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	116	50-150				
<b>Client ID:</b>	<b>H-5-16-8.5</b>					
Laboratory ID:	10-084-03					
Gasoline Range Organics	ND	29	NWTPH-HCID	10-11-16	10-11-16	
Diesel Range Organics	ND	72	NWTPH-HCID	10-11-16	10-11-16	
Lube Oil Range Organics	ND	140	NWTPH-HCID	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	111	50-150				
<b>Client ID:</b>	<b>H-5-16-11</b>					
Laboratory ID:	10-084-04					
Gasoline Range Organics	ND	29	NWTPH-HCID	10-11-16	10-11-16	
Diesel Range Organics	ND	73	NWTPH-HCID	10-11-16	10-11-16	
Lube Oil Range Organics	ND	150	NWTPH-HCID	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	106	50-150				
<b>Client ID:</b>	<b>H-5-16-13.5</b>					
Laboratory ID:	10-084-05					
Gasoline Range Organics	ND	28	NWTPH-HCID	10-11-16	10-11-16	
Diesel Range Organics	ND	69	NWTPH-HCID	10-11-16	10-11-16	
Lube Oil Range Organics	ND	140	NWTPH-HCID	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	111	50-150				



Date of Report: October 19, 2016  
 Samples Submitted: October 10, 2016  
 Laboratory Reference: 1610-084  
 Project: 31008

### NWTPH-HCID

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-5-16-16</b>					
Laboratory ID:	10-084-06					
Gasoline Range Organics	<b>ND</b>	23	NWTPH-HCID	10-11-16	10-11-16	
Diesel Range Organics	<b>ND</b>	58	NWTPH-HCID	10-11-16	10-11-16	
Lube Oil Range Organics	<b>ND</b>	120	NWTPH-HCID	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	116	50-150				

<b>Client ID:</b>	<b>H-5-16-18.5</b>					
Laboratory ID:	10-084-07					
Gasoline Range Organics	<b>ND</b>	23	NWTPH-HCID	10-11-16	10-11-16	
Diesel Range Organics	<b>ND</b>	57	NWTPH-HCID	10-11-16	10-11-16	
Lube Oil Range Organics	<b>ND</b>	110	NWTPH-HCID	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	114	50-150				

<b>Client ID:</b>	<b>H-4-16-3</b>					
Laboratory ID:	10-084-12					
Gasoline Range Organics	<b>ND</b>	22	NWTPH-HCID	10-11-16	10-11-16	
Diesel Range Organics	<b>ND</b>	55	NWTPH-HCID	10-11-16	10-11-16	
Lube Oil Range Organics	<b>ND</b>	110	NWTPH-HCID	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	108	50-150				

<b>Client ID:</b>	<b>H-4-16-6</b>					
Laboratory ID:	10-084-13					
Gasoline Range Organics	<b>ND</b>	23	NWTPH-HCID	10-11-16	10-11-16	
Diesel Range Organics	<b>ND</b>	57	NWTPH-HCID	10-11-16	10-11-16	
Lube Oil Range Organics	<b>ND</b>	110	NWTPH-HCID	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	106	50-150				

<b>Client ID:</b>	<b>H-4-16-8.5</b>					
Laboratory ID:	10-084-14					
Gasoline Range Organics	<b>ND</b>	27	NWTPH-HCID	10-11-16	10-11-16	
Diesel Range Organics	<b>ND</b>	67	NWTPH-HCID	10-11-16	10-11-16	
Lube Oil Range Organics	<b>ND</b>	130	NWTPH-HCID	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	104	50-150				



Date of Report: October 19, 2016  
 Samples Submitted: October 10, 2016  
 Laboratory Reference: 1610-084  
 Project: 31008

### NWTPH-HCID

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-4-16-11</b>					
Laboratory ID:	10-084-15					
Gasoline Range Organics	<b>ND</b>	26	NWTPH-HCID	10-11-16	10-11-16	
Diesel Range Organics	<b>ND</b>	66	NWTPH-HCID	10-11-16	10-11-16	
Lube Oil Range Organics	<b>ND</b>	130	NWTPH-HCID	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>89</i>	<i>50-150</i>				
<b>Client ID:</b>	<b>H-4-16-16</b>					
Laboratory ID:	10-084-16					
Gasoline Range Organics	<b>Detected</b>	23	NWTPH-HCID	10-11-16	10-11-16	
Diesel Range Organics	<b>ND</b>	56	NWTPH-HCID	10-11-16	10-11-16	
Lube Oil Range Organics	<b>ND</b>	110	NWTPH-HCID	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>111</i>	<i>50-150</i>				
<b>Client ID:</b>	<b>H-4-16-18.5</b>					
Laboratory ID:	10-084-17					
Gasoline Range Organics	<b>Detected</b>	23	NWTPH-HCID	10-11-16	10-11-16	
Diesel Range Organics	<b>ND</b>	110	NWTPH-HCID	10-11-16	10-11-16	U1
Lube Oil Range Organics	<b>ND</b>	120	NWTPH-HCID	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>112</i>	<i>50-150</i>				
<b>Client ID:</b>	<b>H-4-16-19.9</b>					
Laboratory ID:	10-084-18					
Gasoline Range Organics	<b>ND</b>	23	NWTPH-HCID	10-11-16	10-11-16	
Diesel Range Organics	<b>ND</b>	58	NWTPH-HCID	10-11-16	10-11-16	
Lube Oil Range Organics	<b>ND</b>	120	NWTPH-HCID	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>91</i>	<i>50-150</i>				
<b>Client ID:</b>	<b>H-4-16-25.4</b>					
Laboratory ID:	10-084-19					
Gasoline Range Organics	<b>ND</b>	24	NWTPH-HCID	10-11-16	10-11-16	
Diesel Range Organics	<b>ND</b>	60	NWTPH-HCID	10-11-16	10-11-16	
Lube Oil Range Organics	<b>ND</b>	120	NWTPH-HCID	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>108</i>	<i>50-150</i>				



Date of Report: October 19, 2016  
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 Project: 31008

**NWTPH-HCID  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg (ppm)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>METHOD BLANK</b>						
Laboratory ID:	MB1011S1					
Gasoline Range Organics	<b>ND</b>	20	NWTPH-HCID	10-11-16	10-11-16	
Diesel Range Organics	<b>ND</b>	50	NWTPH-HCID	10-11-16	10-11-16	
Lube Oil Range Organics	<b>ND</b>	100	NWTPH-HCID	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>90</i>	<i>50-150</i>				



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 Project: 31008

### NWTPH-Gx/BTEX

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-4-16-16</b>					
Laboratory ID:	10-084-16					
Gasoline	<b>69</b>	23	NWTPH-Gx	10-17-16	10-17-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	<i>88</i>	<i>63-124</i>				
<b>Client ID:</b>	<b>H-4-16-18.5</b>					
Laboratory ID:	10-084-17					
Benzene	<b>0.13</b>	0.020	EPA 8021B	10-17-16	10-18-16	
Toluene	<b>0.074</b>	0.058	EPA 8021B	10-17-16	10-18-16	
Ethyl Benzene	<b>0.76</b>	0.058	EPA 8021B	10-17-16	10-18-16	
m,p-Xylene	<b>1.9</b>	0.058	EPA 8021B	10-17-16	10-18-16	
o-Xylene	<b>0.38</b>	0.058	EPA 8021B	10-17-16	10-18-16	
Gasoline	<b>30</b>	5.8	NWTPH-Gx	10-17-16	10-18-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	<i>111</i>	<i>63-124</i>				



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 Project: 31008

**NWTPH-Gx/BTEX  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1017S1					
Benzene	ND	0.020	EPA 8021B	10-17-16	10-17-16	
Toluene	ND	0.050	EPA 8021B	10-17-16	10-17-16	
Ethyl Benzene	ND	0.050	EPA 8021B	10-17-16	10-17-16	
m,p-Xylene	ND	0.050	EPA 8021B	10-17-16	10-17-16	
o-Xylene	ND	0.050	EPA 8021B	10-17-16	10-17-16	
Gasoline	ND	5.0	NWTPH-Gx	10-17-16	10-17-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	91	63-124				

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>								
Laboratory ID:	10-084-17							
	ORIG	DUP						
Benzene	0.111	0.112	NA	NA	NA	NA	1	30
Toluene	0.0634	0.0510	NA	NA	NA	NA	22	30
Ethyl Benzene	0.652	0.645	NA	NA	NA	NA	1	30
m,p-Xylene	1.65	1.62	NA	NA	NA	NA	2	30
o-Xylene	0.326	0.316	NA	NA	NA	NA	3	30
Gasoline	25.2	23.3	NA	NA	NA	NA	8	30
<i>Surrogate:</i>								
<i>Fluorobenzene</i>				111	111	63-124		

**SPIKE BLANKS**

Laboratory ID:	SB	SBD	SB	SBD	SB	SBD			
SB1017S1									
Benzene	0.914	0.902	1.00	1.00	91	90	70-124	1	12
Toluene	0.965	0.910	1.00	1.00	97	91	73-119	6	12
Ethyl Benzene	0.934	0.911	1.00	1.00	93	91	74-117	2	12
m,p-Xylene	0.949	0.885	1.00	1.00	95	89	75-117	7	13
o-Xylene	0.929	0.907	1.00	1.00	93	91	75-116	2	12
<i>Surrogate:</i>									
<i>Fluorobenzene</i>					89	89	63-124		



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**VOLATILES EPA 8260C**  
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Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>H-5-16-13.5</b>					
Laboratory ID:	10-084-05					
Dichlorodifluoromethane	ND	0.083	EPA 8260C	10-11-16	10-11-16	
Chloromethane	ND	0.41	EPA 8260C	10-11-16	10-11-16	
Vinyl Chloride	ND	0.083	EPA 8260C	10-11-16	10-11-16	
Bromomethane	ND	0.083	EPA 8260C	10-11-16	10-11-16	
Chloroethane	ND	0.41	EPA 8260C	10-11-16	10-11-16	
Trichlorofluoromethane	ND	0.083	EPA 8260C	10-11-16	10-11-16	
1,1-Dichloroethene	ND	0.083	EPA 8260C	10-11-16	10-11-16	
Acetone	ND	0.41	EPA 8260C	10-11-16	10-11-16	
Iodomethane	ND	0.41	EPA 8260C	10-11-16	10-11-16	
Carbon Disulfide	ND	0.083	EPA 8260C	10-11-16	10-11-16	
Methylene Chloride	ND	0.41	EPA 8260C	10-11-16	10-11-16	
(trans) 1,2-Dichloroethene	ND	0.083	EPA 8260C	10-11-16	10-11-16	
Methyl t-Butyl Ether	ND	0.083	EPA 8260C	10-11-16	10-11-16	
1,1-Dichloroethane	ND	0.083	EPA 8260C	10-11-16	10-11-16	
Vinyl Acetate	ND	0.41	EPA 8260C	10-11-16	10-11-16	
2,2-Dichloropropane	ND	0.083	EPA 8260C	10-11-16	10-11-16	
(cis) 1,2-Dichloroethene	ND	0.083	EPA 8260C	10-11-16	10-11-16	
2-Butanone	ND	0.41	EPA 8260C	10-11-16	10-11-16	
Bromochloromethane	ND	0.083	EPA 8260C	10-11-16	10-11-16	
Chloroform	ND	0.083	EPA 8260C	10-11-16	10-11-16	
1,1,1-Trichloroethane	ND	0.083	EPA 8260C	10-11-16	10-11-16	
Carbon Tetrachloride	ND	0.083	EPA 8260C	10-11-16	10-11-16	
1,1-Dichloropropene	ND	0.083	EPA 8260C	10-11-16	10-11-16	
Benzene	ND	0.083	EPA 8260C	10-11-16	10-11-16	
1,2-Dichloroethane	ND	0.083	EPA 8260C	10-11-16	10-11-16	
Trichloroethene	ND	0.083	EPA 8260C	10-11-16	10-11-16	
1,2-Dichloropropane	ND	0.083	EPA 8260C	10-11-16	10-11-16	
Dibromomethane	ND	0.083	EPA 8260C	10-11-16	10-11-16	
Bromodichloromethane	ND	0.083	EPA 8260C	10-11-16	10-11-16	
2-Chloroethyl Vinyl Ether	ND	0.41	EPA 8260C	10-11-16	10-11-16	
(cis) 1,3-Dichloropropene	ND	0.083	EPA 8260C	10-11-16	10-11-16	
Methyl Isobutyl Ketone	ND	0.41	EPA 8260C	10-11-16	10-11-16	
Toluene	ND	0.41	EPA 8260C	10-11-16	10-11-16	
(trans) 1,3-Dichloropropene	ND	0.083	EPA 8260C	10-11-16	10-11-16	



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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-5-16-13.5</b>					
<b>Laboratory ID:</b>	10-084-05					
1,1,2-Trichloroethane	ND	0.083	EPA 8260C	10-11-16	10-11-16	
Tetrachloroethene	ND	0.083	EPA 8260C	10-11-16	10-11-16	
1,3-Dichloropropane	ND	0.083	EPA 8260C	10-11-16	10-11-16	
2-Hexanone	ND	0.41	EPA 8260C	10-11-16	10-11-16	
Dibromochloromethane	ND	0.083	EPA 8260C	10-11-16	10-11-16	
1,2-Dibromoethane	ND	0.083	EPA 8260C	10-11-16	10-11-16	
Chlorobenzene	ND	0.083	EPA 8260C	10-11-16	10-11-16	
1,1,1,2-Tetrachloroethane	ND	0.083	EPA 8260C	10-11-16	10-11-16	
Ethylbenzene	0.089	0.083	EPA 8260C	10-11-16	10-11-16	
m,p-Xylene	ND	0.17	EPA 8260C	10-11-16	10-11-16	
o-Xylene	ND	0.083	EPA 8260C	10-11-16	10-11-16	
Styrene	ND	0.083	EPA 8260C	10-11-16	10-11-16	
Bromoform	ND	0.083	EPA 8260C	10-11-16	10-11-16	
Isopropylbenzene	0.19	0.083	EPA 8260C	10-11-16	10-11-16	
Bromobenzene	ND	0.083	EPA 8260C	10-11-16	10-11-16	
1,1,2,2-Tetrachloroethane	ND	0.083	EPA 8260C	10-11-16	10-11-16	
1,2,3-Trichloropropane	ND	0.083	EPA 8260C	10-11-16	10-11-16	
n-Propylbenzene	1.2	0.083	EPA 8260C	10-11-16	10-11-16	
2-Chlorotoluene	ND	0.083	EPA 8260C	10-11-16	10-11-16	
4-Chlorotoluene	ND	0.083	EPA 8260C	10-11-16	10-11-16	
1,3,5-Trimethylbenzene	0.15	0.083	EPA 8260C	10-11-16	10-11-16	
tert-Butylbenzene	ND	0.083	EPA 8260C	10-11-16	10-11-16	
1,2,4-Trimethylbenzene	ND	0.083	EPA 8260C	10-11-16	10-11-16	
sec-Butylbenzene	0.51	0.083	EPA 8260C	10-11-16	10-11-16	
1,3-Dichlorobenzene	ND	0.083	EPA 8260C	10-11-16	10-11-16	
p-Isopropyltoluene	0.32	0.083	EPA 8260C	10-11-16	10-11-16	
1,4-Dichlorobenzene	ND	0.083	EPA 8260C	10-11-16	10-11-16	
1,2-Dichlorobenzene	ND	0.083	EPA 8260C	10-11-16	10-11-16	
n-Butylbenzene	2.0	0.083	EPA 8260C	10-11-16	10-11-16	
1,2-Dibromo-3-chloropropane	ND	0.41	EPA 8260C	10-11-16	10-11-16	
1,2,4-Trichlorobenzene	ND	0.083	EPA 8260C	10-11-16	10-11-16	
Hexachlorobutadiene	ND	0.41	EPA 8260C	10-11-16	10-11-16	
Naphthalene	ND	0.083	EPA 8260C	10-11-16	10-11-16	
1,2,3-Trichlorobenzene	ND	0.083	EPA 8260C	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>93</i>	<i>76-131</i>				
<i>Toluene-d8</i>	<i>102</i>	<i>80-126</i>				
<i>4-Bromofluorobenzene</i>	<i>96</i>	<i>60-146</i>				



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 Project: 31008

**VOLATILES EPA 8260C**  
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Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>H-4-16-16</b>					
<b>Laboratory ID:</b>	<b>10-084-16</b>					
Dichlorodifluoromethane	ND	0.051	EPA 8260C	10-11-16	10-11-16	
Chloromethane	ND	0.25	EPA 8260C	10-11-16	10-11-16	
Vinyl Chloride	ND	0.051	EPA 8260C	10-11-16	10-11-16	
Bromomethane	ND	0.051	EPA 8260C	10-11-16	10-11-16	
Chloroethane	ND	0.25	EPA 8260C	10-11-16	10-11-16	
Trichlorofluoromethane	ND	0.051	EPA 8260C	10-11-16	10-11-16	
1,1-Dichloroethene	ND	0.051	EPA 8260C	10-11-16	10-11-16	
Acetone	ND	0.25	EPA 8260C	10-11-16	10-11-16	
Iodomethane	ND	0.25	EPA 8260C	10-11-16	10-11-16	
Carbon Disulfide	ND	0.051	EPA 8260C	10-11-16	10-11-16	
Methylene Chloride	ND	0.25	EPA 8260C	10-11-16	10-11-16	
(trans) 1,2-Dichloroethene	ND	0.051	EPA 8260C	10-11-16	10-11-16	
Methyl t-Butyl Ether	ND	0.051	EPA 8260C	10-11-16	10-11-16	
1,1-Dichloroethane	ND	0.051	EPA 8260C	10-11-16	10-11-16	
Vinyl Acetate	ND	0.25	EPA 8260C	10-11-16	10-11-16	
2,2-Dichloropropane	ND	0.051	EPA 8260C	10-11-16	10-11-16	
(cis) 1,2-Dichloroethene	ND	0.051	EPA 8260C	10-11-16	10-11-16	
2-Butanone	ND	0.25	EPA 8260C	10-11-16	10-11-16	
Bromochloromethane	ND	0.051	EPA 8260C	10-11-16	10-11-16	
Chloroform	ND	0.051	EPA 8260C	10-11-16	10-11-16	
1,1,1-Trichloroethane	ND	0.051	EPA 8260C	10-11-16	10-11-16	
Carbon Tetrachloride	ND	0.051	EPA 8260C	10-11-16	10-11-16	
1,1-Dichloropropene	ND	0.051	EPA 8260C	10-11-16	10-11-16	
Benzene	ND	0.051	EPA 8260C	10-11-16	10-11-16	
1,2-Dichloroethane	ND	0.051	EPA 8260C	10-11-16	10-11-16	
Trichloroethene	ND	0.051	EPA 8260C	10-11-16	10-11-16	
1,2-Dichloropropane	ND	0.051	EPA 8260C	10-11-16	10-11-16	
Dibromomethane	ND	0.051	EPA 8260C	10-11-16	10-11-16	
Bromodichloromethane	ND	0.051	EPA 8260C	10-11-16	10-11-16	
2-Chloroethyl Vinyl Ether	ND	0.25	EPA 8260C	10-11-16	10-11-16	
(cis) 1,3-Dichloropropene	ND	0.051	EPA 8260C	10-11-16	10-11-16	
Methyl Isobutyl Ketone	ND	0.25	EPA 8260C	10-11-16	10-11-16	
Toluene	ND	0.25	EPA 8260C	10-11-16	10-11-16	
(trans) 1,3-Dichloropropene	ND	0.051	EPA 8260C	10-11-16	10-11-16	



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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-4-16-16</b>					
Laboratory ID:	10-084-16					
1,1,2-Trichloroethane	ND	0.051	EPA 8260C	10-11-16	10-11-16	
Tetrachloroethene	ND	0.051	EPA 8260C	10-11-16	10-11-16	
1,3-Dichloropropane	ND	0.051	EPA 8260C	10-11-16	10-11-16	
2-Hexanone	ND	0.25	EPA 8260C	10-11-16	10-11-16	
Dibromochloromethane	ND	0.051	EPA 8260C	10-11-16	10-11-16	
1,2-Dibromoethane	ND	0.051	EPA 8260C	10-11-16	10-11-16	
Chlorobenzene	ND	0.051	EPA 8260C	10-11-16	10-11-16	
1,1,1,2-Tetrachloroethane	ND	0.051	EPA 8260C	10-11-16	10-11-16	
Ethylbenzene	0.55	0.051	EPA 8260C	10-11-16	10-11-16	
m,p-Xylene	1.4	0.10	EPA 8260C	10-11-16	10-11-16	
o-Xylene	0.49	0.051	EPA 8260C	10-11-16	10-11-16	
Styrene	ND	0.051	EPA 8260C	10-11-16	10-11-16	
Bromoform	ND	0.051	EPA 8260C	10-11-16	10-11-16	
Isopropylbenzene	0.092	0.051	EPA 8260C	10-11-16	10-11-16	
Bromobenzene	ND	0.051	EPA 8260C	10-11-16	10-11-16	
1,1,2,2-Tetrachloroethane	ND	0.051	EPA 8260C	10-11-16	10-11-16	
1,2,3-Trichloropropane	ND	0.051	EPA 8260C	10-11-16	10-11-16	
n-Propylbenzene	0.36	0.051	EPA 8260C	10-11-16	10-11-16	
2-Chlorotoluene	ND	0.051	EPA 8260C	10-11-16	10-11-16	
4-Chlorotoluene	ND	0.051	EPA 8260C	10-11-16	10-11-16	
1,3,5-Trimethylbenzene	0.56	0.051	EPA 8260C	10-11-16	10-11-16	
tert-Butylbenzene	ND	0.051	EPA 8260C	10-11-16	10-11-16	
1,2,4-Trimethylbenzene	1.8	0.051	EPA 8260C	10-11-16	10-11-16	
sec-Butylbenzene	0.064	0.051	EPA 8260C	10-11-16	10-11-16	
1,3-Dichlorobenzene	ND	0.051	EPA 8260C	10-11-16	10-11-16	
p-Isopropyltoluene	0.053	0.051	EPA 8260C	10-11-16	10-11-16	
1,4-Dichlorobenzene	ND	0.051	EPA 8260C	10-11-16	10-11-16	
1,2-Dichlorobenzene	ND	0.051	EPA 8260C	10-11-16	10-11-16	
n-Butylbenzene	0.29	0.051	EPA 8260C	10-11-16	10-11-16	
1,2-Dibromo-3-chloropropane	ND	0.25	EPA 8260C	10-11-16	10-11-16	
1,2,4-Trichlorobenzene	ND	0.051	EPA 8260C	10-11-16	10-11-16	
Hexachlorobutadiene	ND	0.25	EPA 8260C	10-11-16	10-11-16	
Naphthalene	0.64	0.051	EPA 8260C	10-11-16	10-11-16	
1,2,3-Trichlorobenzene	ND	0.051	EPA 8260C	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>92</i>	<i>76-131</i>				
<i>Toluene-d8</i>	<i>95</i>	<i>80-126</i>				
<i>4-Bromofluorobenzene</i>	<i>96</i>	<i>60-146</i>				



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**VOLATILES EPA 8260C**  
**METHOD BLANK QUALITY CONTROL**

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Matrix: Soil  
 Units: mg/kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1011S2					
Dichlorodifluoromethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Chloromethane	ND	0.0050	EPA 8260C	10-11-16	10-11-16	
Vinyl Chloride	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Bromomethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Chloroethane	ND	0.0050	EPA 8260C	10-11-16	10-11-16	
Trichlorofluoromethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,1-Dichloroethene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Acetone	ND	0.0050	EPA 8260C	10-11-16	10-11-16	
Iodomethane	ND	0.0050	EPA 8260C	10-11-16	10-11-16	
Carbon Disulfide	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Methylene Chloride	ND	0.0050	EPA 8260C	10-11-16	10-11-16	
(trans) 1,2-Dichloroethene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Methyl t-Butyl Ether	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,1-Dichloroethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Vinyl Acetate	ND	0.0050	EPA 8260C	10-11-16	10-11-16	
2,2-Dichloropropane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
(cis) 1,2-Dichloroethene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
2-Butanone	ND	0.0050	EPA 8260C	10-11-16	10-11-16	
Bromochloromethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Chloroform	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,1,1-Trichloroethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Carbon Tetrachloride	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,1-Dichloropropene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Benzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,2-Dichloroethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Trichloroethene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,2-Dichloropropane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Dibromomethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Bromodichloromethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
2-Chloroethyl Vinyl Ether	ND	0.0050	EPA 8260C	10-11-16	10-11-16	
(cis) 1,3-Dichloropropene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Methyl Isobutyl Ketone	ND	0.0050	EPA 8260C	10-11-16	10-11-16	
Toluene	ND	0.0050	EPA 8260C	10-11-16	10-11-16	
(trans) 1,3-Dichloropropene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	



Date of Report: October 19, 2016  
 Samples Submitted: October 10, 2016  
 Laboratory Reference: 1610-084  
 Project: 31008

**VOLATILES EPA 8260C**  
**METHOD BLANK QUALITY CONTROL**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1011S2					
1,1,2-Trichloroethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Tetrachloroethene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,3-Dichloropropane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
2-Hexanone	ND	0.0050	EPA 8260C	10-11-16	10-11-16	
Dibromochloromethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,2-Dibromoethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Chlorobenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,1,1,2-Tetrachloroethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Ethylbenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
m,p-Xylene	ND	0.0020	EPA 8260C	10-11-16	10-11-16	
o-Xylene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Styrene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Bromoform	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Isopropylbenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Bromobenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,1,2,2-Tetrachloroethane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,2,3-Trichloropropane	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
n-Propylbenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
2-Chlorotoluene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
4-Chlorotoluene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,3,5-Trimethylbenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
tert-Butylbenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,2,4-Trimethylbenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
sec-Butylbenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,3-Dichlorobenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
p-Isopropyltoluene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,4-Dichlorobenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,2-Dichlorobenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
n-Butylbenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,2-Dibromo-3-chloropropane	ND	0.0050	EPA 8260C	10-11-16	10-11-16	
1,2,4-Trichlorobenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
Hexachlorobutadiene	ND	0.0050	EPA 8260C	10-11-16	10-11-16	
Naphthalene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
1,2,3-Trichlorobenzene	ND	0.0010	EPA 8260C	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>100</i>	<i>76-131</i>				
<i>Toluene-d8</i>	<i>103</i>	<i>80-126</i>				
<i>4-Bromofluorobenzene</i>	<i>103</i>	<i>60-146</i>				



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**VOLATILES EPA 8260C  
 SB/SBD QUALITY CONTROL**

Matrix: Soil  
 Units: mg/kg

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD		Flags
					Recovery	Limits	RPD	Limit		
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB1011S1									
	SB	SBD	SB	SBD	SB	SBD				
1,1-Dichloroethene	<b>0.0472</b>	<b>0.0496</b>	0.0500	0.0500	94	99	68-126	5	15	
Benzene	<b>0.0473</b>	<b>0.0487</b>	0.0500	0.0500	95	97	70-121	3	15	
Trichloroethene	<b>0.0440</b>	<b>0.0462</b>	0.0500	0.0500	88	92	75-120	5	15	
Toluene	<b>0.0459</b>	<b>0.0487</b>	0.0500	0.0500	92	97	80-120	6	15	
Chlorobenzene	<b>0.0474</b>	<b>0.0478</b>	0.0500	0.0500	95	96	76-120	1	15	
<i>Surrogate:</i>										
Dibromofluoromethane					98	99	76-131			
Toluene-d8					98	100	80-126			
4-Bromofluorobenzene					100	101	60-146			



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 Project: 31008

**SEMIVOLATILES EPA 8270D/SIM**

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Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-5-16-13.5</b>					
Laboratory ID:	10-084-05					
n-Nitrosodimethylamine	ND	0.046	EPA 8270D	10-12-16	10-13-16	
Pyridine	ND	0.46	EPA 8270D	10-12-16	10-13-16	
Phenol	ND	0.046	EPA 8270D	10-12-16	10-13-16	
Aniline	ND	0.23	EPA 8270D	10-12-16	10-13-16	
bis(2-Chloroethyl)ether	ND	0.046	EPA 8270D	10-12-16	10-13-16	
2-Chlorophenol	ND	0.046	EPA 8270D	10-12-16	10-13-16	
1,3-Dichlorobenzene	ND	0.046	EPA 8270D	10-12-16	10-13-16	
1,4-Dichlorobenzene	ND	0.046	EPA 8270D	10-12-16	10-13-16	
Benzyl alcohol	ND	0.23	EPA 8270D	10-12-16	10-13-16	
1,2-Dichlorobenzene	ND	0.046	EPA 8270D	10-12-16	10-13-16	
2-Methylphenol (o-Cresol)	ND	0.046	EPA 8270D	10-12-16	10-13-16	
bis(2-Chloroisopropyl)ether	ND	0.046	EPA 8270D	10-12-16	10-13-16	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.046	EPA 8270D	10-12-16	10-13-16	
n-Nitroso-di-n-propylamine	ND	0.046	EPA 8270D	10-12-16	10-13-16	
Hexachloroethane	ND	0.046	EPA 8270D	10-12-16	10-13-16	
Nitrobenzene	ND	0.046	EPA 8270D	10-12-16	10-13-16	
Isophorone	ND	0.046	EPA 8270D	10-12-16	10-13-16	
2-Nitrophenol	ND	0.046	EPA 8270D	10-12-16	10-13-16	
2,4-Dimethylphenol	ND	0.046	EPA 8270D	10-12-16	10-13-16	
bis(2-Chloroethoxy)methane	ND	0.046	EPA 8270D	10-12-16	10-13-16	
2,4-Dichlorophenol	ND	0.046	EPA 8270D	10-12-16	10-13-16	
1,2,4-Trichlorobenzene	ND	0.046	EPA 8270D	10-12-16	10-13-16	
Naphthalene	<b>0.039</b>	0.0092	EPA 8270D/SIM	10-12-16	10-13-16	
4-Chloroaniline	ND	0.23	EPA 8270D	10-12-16	10-13-16	
Hexachlorobutadiene	ND	0.046	EPA 8270D	10-12-16	10-13-16	
4-Chloro-3-methylphenol	ND	0.046	EPA 8270D	10-12-16	10-13-16	
2-Methylnaphthalene	<b>0.35</b>	0.046	EPA 8270D	10-12-16	10-13-16	
1-Methylnaphthalene	<b>0.037</b>	0.0092	EPA 8270D/SIM	10-12-16	10-13-16	
Hexachlorocyclopentadiene	ND	0.046	EPA 8270D	10-12-16	10-13-16	
2,4,6-Trichlorophenol	ND	0.046	EPA 8270D	10-12-16	10-13-16	
2,3-Dichloroaniline	ND	0.046	EPA 8270D	10-12-16	10-13-16	
2,4,5-Trichlorophenol	ND	0.046	EPA 8270D	10-12-16	10-13-16	
2-Chloronaphthalene	ND	0.046	EPA 8270D	10-12-16	10-13-16	
2-Nitroaniline	ND	0.046	EPA 8270D	10-12-16	10-13-16	
1,4-Dinitrobenzene	ND	0.046	EPA 8270D	10-12-16	10-13-16	
Dimethylphthalate	ND	0.046	EPA 8270D	10-12-16	10-13-16	
1,3-Dinitrobenzene	ND	0.046	EPA 8270D	10-12-16	10-13-16	
2,6-Dinitrotoluene	ND	0.046	EPA 8270D	10-12-16	10-13-16	
1,2-Dinitrobenzene	ND	0.046	EPA 8270D	10-12-16	10-13-16	
Acenaphthylene	ND	0.0092	EPA 8270D/SIM	10-12-16	10-13-16	
3-Nitroaniline	ND	0.046	EPA 8270D	10-12-16	10-13-16	



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**SEMIVOLATILES EPA 8270D/SIM**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-5-16-13.5</b>					
Laboratory ID:	10-084-05					
2,4-Dinitrophenol	ND	0.23	EPA 8270D	10-12-16	10-13-16	
Acenaphthene	ND	0.0092	EPA 8270D/SIM	10-12-16	10-13-16	
4-Nitrophenol	ND	0.046	EPA 8270D	10-12-16	10-13-16	
2,4-Dinitrotoluene	ND	0.046	EPA 8270D	10-12-16	10-13-16	
Dibenzofuran	ND	0.046	EPA 8270D	10-12-16	10-13-16	
2,3,5,6-Tetrachlorophenol	ND	0.046	EPA 8270D	10-12-16	10-13-16	
2,3,4,6-Tetrachlorophenol	ND	0.046	EPA 8270D	10-12-16	10-13-16	
Diethylphthalate	ND	0.23	EPA 8270D	10-12-16	10-13-16	
4-Chlorophenyl-phenylether	ND	0.046	EPA 8270D	10-12-16	10-13-16	
4-Nitroaniline	ND	0.046	EPA 8270D	10-12-16	10-13-16	
Fluorene	ND	0.0092	EPA 8270D/SIM	10-12-16	10-13-16	
4,6-Dinitro-2-methylphenol	ND	0.23	EPA 8270D	10-12-16	10-13-16	
n-Nitrosodiphenylamine	ND	0.046	EPA 8270D	10-12-16	10-13-16	
1,2-Diphenylhydrazine	ND	0.046	EPA 8270D	10-12-16	10-13-16	
4-Bromophenyl-phenylether	ND	0.046	EPA 8270D	10-12-16	10-13-16	
Hexachlorobenzene	ND	0.046	EPA 8270D	10-12-16	10-13-16	
Pentachlorophenol	ND	0.23	EPA 8270D	10-12-16	10-13-16	
Phenanthrene	ND	0.0092	EPA 8270D/SIM	10-12-16	10-13-16	
Anthracene	ND	0.0092	EPA 8270D/SIM	10-12-16	10-13-16	
Carbazole	ND	0.046	EPA 8270D	10-12-16	10-13-16	
Di-n-butylphthalate	ND	0.23	EPA 8270D	10-12-16	10-13-16	
Fluoranthene	ND	0.0092	EPA 8270D/SIM	10-12-16	10-13-16	
Benzidine	ND	0.46	EPA 8270D	10-12-16	10-13-16	
Pyrene	ND	0.0092	EPA 8270D/SIM	10-12-16	10-13-16	
Butylbenzylphthalate	ND	0.046	EPA 8270D	10-12-16	10-13-16	
bis-2-Ethylhexyladipate	ND	0.046	EPA 8270D	10-12-16	10-13-16	
3,3'-Dichlorobenzidine	ND	0.23	EPA 8270D	10-12-16	10-13-16	
Benzo[a]anthracene	ND	0.0092	EPA 8270D/SIM	10-12-16	10-13-16	
Chrysene	ND	0.0092	EPA 8270D/SIM	10-12-16	10-13-16	
bis(2-Ethylhexyl)phthalate	ND	0.046	EPA 8270D	10-12-16	10-13-16	
Di-n-octylphthalate	ND	0.046	EPA 8270D	10-12-16	10-13-16	
Benzo[b]fluoranthene	ND	0.0092	EPA 8270D/SIM	10-12-16	10-13-16	
Benzo(j,k)fluoranthene	ND	0.0092	EPA 8270D/SIM	10-12-16	10-13-16	
Benzo[a]pyrene	ND	0.0092	EPA 8270D/SIM	10-12-16	10-13-16	
Indeno[1,2,3-cd]pyrene	ND	0.0092	EPA 8270D/SIM	10-12-16	10-13-16	
Dibenz[a,h]anthracene	ND	0.0092	EPA 8270D/SIM	10-12-16	10-13-16	
Benzo[g,h,i]perylene	ND	0.0092	EPA 8270D/SIM	10-12-16	10-13-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	62	24 - 117				
Phenol-d6	65	30 - 120				
Nitrobenzene-d5	67	27 - 112				
2-Fluorobiphenyl	66	35 - 113				
2,4,6-Tribromophenol	68	21 - 120				
Terphenyl-d14	70	39 - 121				



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**SEMIVOLATILES EPA 8270D/SIM**  
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Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-4-16-16</b>					
<b>Laboratory ID:</b>	<b>10-084-16</b>					
n-Nitrosodimethylamine	ND	0.037	EPA 8270D	10-12-16	10-13-16	
Pyridine	ND	0.37	EPA 8270D	10-12-16	10-13-16	
Phenol	ND	0.037	EPA 8270D	10-12-16	10-13-16	
Aniline	ND	0.19	EPA 8270D	10-12-16	10-13-16	
bis(2-Chloroethyl)ether	ND	0.037	EPA 8270D	10-12-16	10-13-16	
2-Chlorophenol	ND	0.037	EPA 8270D	10-12-16	10-13-16	
1,3-Dichlorobenzene	ND	0.037	EPA 8270D	10-12-16	10-13-16	
1,4-Dichlorobenzene	ND	0.037	EPA 8270D	10-12-16	10-13-16	
Benzyl alcohol	ND	0.19	EPA 8270D	10-12-16	10-13-16	
1,2-Dichlorobenzene	ND	0.037	EPA 8270D	10-12-16	10-13-16	
2-Methylphenol (o-Cresol)	ND	0.037	EPA 8270D	10-12-16	10-13-16	
bis(2-Chloroisopropyl)ether	ND	0.037	EPA 8270D	10-12-16	10-13-16	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.037	EPA 8270D	10-12-16	10-13-16	
n-Nitroso-di-n-propylamine	ND	0.037	EPA 8270D	10-12-16	10-13-16	
Hexachloroethane	ND	0.037	EPA 8270D	10-12-16	10-13-16	
Nitrobenzene	ND	0.037	EPA 8270D	10-12-16	10-13-16	
Isophorone	ND	0.037	EPA 8270D	10-12-16	10-13-16	
2-Nitrophenol	ND	0.037	EPA 8270D	10-12-16	10-13-16	
2,4-Dimethylphenol	ND	0.037	EPA 8270D	10-12-16	10-13-16	
bis(2-Chloroethoxy)methane	ND	0.037	EPA 8270D	10-12-16	10-13-16	
2,4-Dichlorophenol	ND	0.037	EPA 8270D	10-12-16	10-13-16	
1,2,4-Trichlorobenzene	ND	0.037	EPA 8270D	10-12-16	10-13-16	
Naphthalene	0.59	0.037	EPA 8270D	10-12-16	10-13-16	
4-Chloroaniline	ND	0.19	EPA 8270D	10-12-16	10-13-16	
Hexachlorobutadiene	ND	0.037	EPA 8270D	10-12-16	10-13-16	
4-Chloro-3-methylphenol	ND	0.037	EPA 8270D	10-12-16	10-13-16	
2-Methylnaphthalene	0.74	0.037	EPA 8270D	10-12-16	10-13-16	
1-Methylnaphthalene	0.37	0.037	EPA 8270D	10-12-16	10-13-16	
Hexachlorocyclopentadiene	ND	0.037	EPA 8270D	10-12-16	10-13-16	
2,4,6-Trichlorophenol	ND	0.037	EPA 8270D	10-12-16	10-13-16	
2,3-Dichloroaniline	ND	0.037	EPA 8270D	10-12-16	10-13-16	
2,4,5-Trichlorophenol	ND	0.037	EPA 8270D	10-12-16	10-13-16	
2-Chloronaphthalene	ND	0.037	EPA 8270D	10-12-16	10-13-16	
2-Nitroaniline	ND	0.037	EPA 8270D	10-12-16	10-13-16	
1,4-Dinitrobenzene	ND	0.037	EPA 8270D	10-12-16	10-13-16	
Dimethylphthalate	ND	0.037	EPA 8270D	10-12-16	10-13-16	
1,3-Dinitrobenzene	ND	0.037	EPA 8270D	10-12-16	10-13-16	
2,6-Dinitrotoluene	ND	0.037	EPA 8270D	10-12-16	10-13-16	
1,2-Dinitrobenzene	ND	0.037	EPA 8270D	10-12-16	10-13-16	
Acenaphthylene	ND	0.0075	EPA 8270D/SIM	10-12-16	10-13-16	
3-Nitroaniline	ND	0.037	EPA 8270D	10-12-16	10-13-16	



Date of Report: October 19, 2016  
 Samples Submitted: October 10, 2016  
 Laboratory Reference: 1610-084  
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**SEMIVOLATILES EPA 8270D/SIM**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-4-16-16</b>					
<b>Laboratory ID:</b>	<b>10-084-16</b>					
2,4-Dinitrophenol	ND	0.19	EPA 8270D	10-12-16	10-13-16	
Acenaphthene	ND	0.0075	EPA 8270D/SIM	10-12-16	10-13-16	
4-Nitrophenol	ND	0.037	EPA 8270D	10-12-16	10-13-16	
2,4-Dinitrotoluene	ND	0.037	EPA 8270D	10-12-16	10-13-16	
Dibenzofuran	ND	0.037	EPA 8270D	10-12-16	10-13-16	
2,3,5,6-Tetrachlorophenol	ND	0.037	EPA 8270D	10-12-16	10-13-16	
2,3,4,6-Tetrachlorophenol	ND	0.037	EPA 8270D	10-12-16	10-13-16	
Diethylphthalate	ND	0.19	EPA 8270D	10-12-16	10-13-16	
4-Chlorophenyl-phenylether	ND	0.037	EPA 8270D	10-12-16	10-13-16	
4-Nitroaniline	ND	0.037	EPA 8270D	10-12-16	10-13-16	
Fluorene	ND	0.0075	EPA 8270D/SIM	10-12-16	10-13-16	
4,6-Dinitro-2-methylphenol	ND	0.19	EPA 8270D	10-12-16	10-13-16	
n-Nitrosodiphenylamine	ND	0.037	EPA 8270D	10-12-16	10-13-16	
1,2-Diphenylhydrazine	ND	0.037	EPA 8270D	10-12-16	10-13-16	
4-Bromophenyl-phenylether	ND	0.037	EPA 8270D	10-12-16	10-13-16	
Hexachlorobenzene	ND	0.037	EPA 8270D	10-12-16	10-13-16	
Pentachlorophenol	ND	0.19	EPA 8270D	10-12-16	10-13-16	
Phenanthrene	ND	0.0075	EPA 8270D/SIM	10-12-16	10-13-16	
Anthracene	ND	0.0075	EPA 8270D/SIM	10-12-16	10-13-16	
Carbazole	ND	0.037	EPA 8270D	10-12-16	10-13-16	
Di-n-butylphthalate	ND	0.19	EPA 8270D	10-12-16	10-13-16	
Fluoranthene	ND	0.0075	EPA 8270D/SIM	10-12-16	10-13-16	
Benzidine	ND	0.37	EPA 8270D	10-12-16	10-13-16	
Pyrene	ND	0.0075	EPA 8270D/SIM	10-12-16	10-13-16	
Butylbenzylphthalate	ND	0.037	EPA 8270D	10-12-16	10-13-16	
bis-2-Ethylhexyladipate	ND	0.037	EPA 8270D	10-12-16	10-13-16	
3,3'-Dichlorobenzidine	ND	0.19	EPA 8270D	10-12-16	10-13-16	
Benzo[a]anthracene	ND	0.0075	EPA 8270D/SIM	10-12-16	10-13-16	
Chrysene	ND	0.0075	EPA 8270D/SIM	10-12-16	10-13-16	
bis(2-Ethylhexyl)phthalate	ND	0.037	EPA 8270D	10-12-16	10-13-16	
Di-n-octylphthalate	ND	0.037	EPA 8270D	10-12-16	10-13-16	
Benzo[b]fluoranthene	ND	0.0075	EPA 8270D/SIM	10-12-16	10-13-16	
Benzo(j,k)fluoranthene	ND	0.0075	EPA 8270D/SIM	10-12-16	10-13-16	
Benzo[a]pyrene	ND	0.0075	EPA 8270D/SIM	10-12-16	10-13-16	
Indeno[1,2,3-cd]pyrene	ND	0.0075	EPA 8270D/SIM	10-12-16	10-13-16	
Dibenz[a,h]anthracene	ND	0.0075	EPA 8270D/SIM	10-12-16	10-13-16	
Benzo[g,h,i]perylene	ND	0.0075	EPA 8270D/SIM	10-12-16	10-13-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorophenol</i>	<i>67</i>	<i>24 - 117</i>				
<i>Phenol-d6</i>	<i>69</i>	<i>30 - 120</i>				
<i>Nitrobenzene-d5</i>	<i>71</i>	<i>27 - 112</i>				
<i>2-Fluorobiphenyl</i>	<i>70</i>	<i>35 - 113</i>				
<i>2,4,6-Tribromophenol</i>	<i>74</i>	<i>21 - 120</i>				
<i>Terphenyl-d14</i>	<i>76</i>	<i>39 - 121</i>				



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**SEMIVOLATILES EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
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Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1012S1					
n-Nitrosodimethylamine	ND	0.033	EPA 8270D	10-12-16	10-12-16	
Pyridine	ND	0.33	EPA 8270D	10-12-16	10-12-16	
Phenol	ND	0.033	EPA 8270D	10-12-16	10-12-16	
Aniline	ND	0.17	EPA 8270D	10-12-16	10-12-16	
bis(2-Chloroethyl)ether	ND	0.033	EPA 8270D	10-12-16	10-12-16	
2-Chlorophenol	ND	0.033	EPA 8270D	10-12-16	10-12-16	
1,3-Dichlorobenzene	ND	0.033	EPA 8270D	10-12-16	10-12-16	
1,4-Dichlorobenzene	ND	0.033	EPA 8270D	10-12-16	10-12-16	
Benzyl alcohol	ND	0.17	EPA 8270D	10-12-16	10-12-16	
1,2-Dichlorobenzene	ND	0.033	EPA 8270D	10-12-16	10-12-16	
2-Methylphenol (o-Cresol)	ND	0.033	EPA 8270D	10-12-16	10-12-16	
bis(2-Chloroisopropyl)ether	ND	0.033	EPA 8270D	10-12-16	10-12-16	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.033	EPA 8270D	10-12-16	10-12-16	
n-Nitroso-di-n-propylamine	ND	0.033	EPA 8270D	10-12-16	10-12-16	
Hexachloroethane	ND	0.033	EPA 8270D	10-12-16	10-12-16	
Nitrobenzene	ND	0.033	EPA 8270D	10-12-16	10-12-16	
Isophorone	ND	0.033	EPA 8270D	10-12-16	10-12-16	
2-Nitrophenol	ND	0.033	EPA 8270D	10-12-16	10-12-16	
2,4-Dimethylphenol	ND	0.033	EPA 8270D	10-12-16	10-12-16	
bis(2-Chloroethoxy)methane	ND	0.033	EPA 8270D	10-12-16	10-12-16	
2,4-Dichlorophenol	ND	0.033	EPA 8270D	10-12-16	10-12-16	
1,2,4-Trichlorobenzene	ND	0.033	EPA 8270D	10-12-16	10-12-16	
Naphthalene	ND	0.0067	EPA 8270D/SIM	10-12-16	10-13-16	
4-Chloroaniline	ND	0.17	EPA 8270D	10-12-16	10-12-16	
Hexachlorobutadiene	ND	0.033	EPA 8270D	10-12-16	10-12-16	
4-Chloro-3-methylphenol	ND	0.033	EPA 8270D	10-12-16	10-12-16	
2-Methylnaphthalene	ND	0.0067	EPA 8270D/SIM	10-12-16	10-13-16	
1-Methylnaphthalene	ND	0.0067	EPA 8270D/SIM	10-12-16	10-13-16	
Hexachlorocyclopentadiene	ND	0.033	EPA 8270D	10-12-16	10-12-16	
2,4,6-Trichlorophenol	ND	0.033	EPA 8270D	10-12-16	10-12-16	
2,3-Dichloroaniline	ND	0.033	EPA 8270D	10-12-16	10-12-16	
2,4,5-Trichlorophenol	ND	0.033	EPA 8270D	10-12-16	10-12-16	
2-Chloronaphthalene	ND	0.033	EPA 8270D	10-12-16	10-12-16	
2-Nitroaniline	ND	0.033	EPA 8270D	10-12-16	10-12-16	
1,4-Dinitrobenzene	ND	0.033	EPA 8270D	10-12-16	10-12-16	
Dimethylphthalate	ND	0.033	EPA 8270D	10-12-16	10-12-16	
1,3-Dinitrobenzene	ND	0.033	EPA 8270D	10-12-16	10-12-16	
2,6-Dinitrotoluene	ND	0.033	EPA 8270D	10-12-16	10-12-16	
1,2-Dinitrobenzene	ND	0.033	EPA 8270D	10-12-16	10-12-16	
Acenaphthylene	ND	0.0067	EPA 8270D/SIM	10-12-16	10-13-16	
3-Nitroaniline	ND	0.033	EPA 8270D	10-12-16	10-12-16	



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**SEMIVOLATILES EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1012S1					
2,4-Dinitrophenol	ND	0.17	EPA 8270D	10-12-16	10-12-16	
Acenaphthene	ND	0.0067	EPA 8270D/SIM	10-12-16	10-13-16	
4-Nitrophenol	ND	0.033	EPA 8270D	10-12-16	10-12-16	
2,4-Dinitrotoluene	ND	0.033	EPA 8270D	10-12-16	10-12-16	
Dibenzofuran	ND	0.033	EPA 8270D	10-12-16	10-12-16	
2,3,5,6-Tetrachlorophenol	ND	0.033	EPA 8270D	10-12-16	10-12-16	
2,3,4,6-Tetrachlorophenol	ND	0.033	EPA 8270D	10-12-16	10-12-16	
Diethylphthalate	ND	0.17	EPA 8270D	10-12-16	10-12-16	
4-Chlorophenyl-phenylether	ND	0.033	EPA 8270D	10-12-16	10-12-16	
4-Nitroaniline	ND	0.033	EPA 8270D	10-12-16	10-12-16	
Fluorene	ND	0.0067	EPA 8270D/SIM	10-12-16	10-13-16	
4,6-Dinitro-2-methylphenol	ND	0.17	EPA 8270D	10-12-16	10-12-16	
n-Nitrosodiphenylamine	ND	0.033	EPA 8270D	10-12-16	10-12-16	
1,2-Diphenylhydrazine	ND	0.033	EPA 8270D	10-12-16	10-12-16	
4-Bromophenyl-phenylether	ND	0.033	EPA 8270D	10-12-16	10-12-16	
Hexachlorobenzene	ND	0.033	EPA 8270D	10-12-16	10-12-16	
Pentachlorophenol	ND	0.17	EPA 8270D	10-12-16	10-12-16	
Phenanthrene	ND	0.0067	EPA 8270D/SIM	10-12-16	10-13-16	
Anthracene	ND	0.0067	EPA 8270D/SIM	10-12-16	10-13-16	
Carbazole	ND	0.033	EPA 8270D	10-12-16	10-12-16	
Di-n-butylphthalate	ND	0.17	EPA 8270D	10-12-16	10-12-16	
Fluoranthene	ND	0.0067	EPA 8270D/SIM	10-12-16	10-13-16	
Benzidine	ND	0.33	EPA 8270D	10-12-16	10-12-16	
Pyrene	ND	0.0067	EPA 8270D/SIM	10-12-16	10-13-16	
Butylbenzylphthalate	ND	0.033	EPA 8270D	10-12-16	10-12-16	
bis-2-Ethylhexyladipate	ND	0.033	EPA 8270D	10-12-16	10-12-16	
3,3'-Dichlorobenzidine	ND	0.17	EPA 8270D	10-12-16	10-12-16	
Benzo[a]anthracene	ND	0.0067	EPA 8270D/SIM	10-12-16	10-13-16	
Chrysene	ND	0.0067	EPA 8270D/SIM	10-12-16	10-13-16	
bis(2-Ethylhexyl)phthalate	ND	0.033	EPA 8270D	10-12-16	10-12-16	
Di-n-octylphthalate	ND	0.033	EPA 8270D	10-12-16	10-12-16	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270D/SIM	10-12-16	10-13-16	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270D/SIM	10-12-16	10-13-16	
Benzo[a]pyrene	ND	0.0067	EPA 8270D/SIM	10-12-16	10-13-16	
Indeno[1,2,3-cd]pyrene	ND	0.0067	EPA 8270D/SIM	10-12-16	10-13-16	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270D/SIM	10-12-16	10-13-16	
Benzo[g,h,i]perylene	ND	0.0067	EPA 8270D/SIM	10-12-16	10-13-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	89	24 - 117				
Phenol-d6	94	30 - 120				
Nitrobenzene-d5	96	27 - 112				
2-Fluorobiphenyl	89	35 - 113				
2,4,6-Tribromophenol	97	21 - 120				
Terphenyl-d14	93	39 - 121				



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 Project: 31008

**SEMIVOLATILES EPA 8270D/SIM  
 MS/MSD QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result		Spike Level		Source	Percent		Recovery	RPD	RPD	Flags
	MS	MSD	MS	MSD	Result	Recovery	Limits	RPD	Limit		
<b>MATRIX SPIKES</b>											
Laboratory ID:	10-015-01										
	MS	MSD	MS	MSD		MS	MSD				
Phenol	<b>0.679</b>	<b>0.658</b>	1.33	1.33	ND	51	49	31 - 108	3	36	
2-Chlorophenol	<b>0.671</b>	<b>0.659</b>	1.33	1.33	ND	50	50	38 - 103	2	38	
1,4-Dichlorobenzene	<b>0.325</b>	<b>0.329</b>	0.667	0.667	ND	49	49	25 - 101	1	40	
n-Nitroso-di-n-propylamine	<b>0.322</b>	<b>0.324</b>	0.667	0.667	ND	48	49	26 - 102	1	38	
1,2,4-Trichlorobenzene	<b>0.325</b>	<b>0.328</b>	0.667	0.667	ND	49	49	27 - 101	1	40	
4-Chloro-3-methylphenol	<b>0.765</b>	<b>0.802</b>	1.33	1.33	ND	58	60	42 - 106	5	29	
Acenaphthene	<b>0.385</b>	<b>0.404</b>	0.667	0.667	ND	58	61	42 - 103	5	30	
4-Nitrophenol	<b>0.664</b>	<b>0.713</b>	1.33	1.33	ND	50	54	25 - 125	7	29	
2,4-Dinitrotoluene	<b>0.403</b>	<b>0.422</b>	0.667	0.667	ND	60	63	45 - 107	5	30	
Pentachlorophenol	<b>0.629</b>	<b>0.625</b>	1.33	1.33	ND	47	47	30 - 103	1	31	
Pyrene	<b>0.488</b>	<b>0.500</b>	0.667	0.667	ND	73	75	50 - 118	2	28	
<i>Surrogate:</i>											
2-Fluorophenol						47	48	24 - 117			
Phenol-d6						52	50	30 - 120			
Nitrobenzene-d5						54	54	27 - 112			
2-Fluorobiphenyl						53	55	35 - 113			
2,4,6-Tribromophenol						69	74	21 - 120			
Terphenyl-d14						73	75	39 - 121			



Date of Report: October 19, 2016  
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 Laboratory Reference: 1610-084  
 Project: 31008

**PCBs**  
**EPA 8082A**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-5-16-13.5</b>					
Laboratory ID:	10-084-05					
Aroclor 1016	<b>ND</b>	0.069	EPA 8082A	10-11-16	10-11-16	
Aroclor 1221	<b>ND</b>	0.069	EPA 8082A	10-11-16	10-11-16	
Aroclor 1232	<b>ND</b>	0.069	EPA 8082A	10-11-16	10-11-16	
Aroclor 1242	<b>ND</b>	0.069	EPA 8082A	10-11-16	10-11-16	
Aroclor 1248	<b>ND</b>	0.069	EPA 8082A	10-11-16	10-11-16	
Aroclor 1254	<b>ND</b>	0.069	EPA 8082A	10-11-16	10-11-16	
Aroclor 1260	<b>ND</b>	0.069	EPA 8082A	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>DCB</i>	<i>70</i>	<i>50-139</i>				
<b>Client ID:</b>	<b>H-4-16-16</b>					
Laboratory ID:	10-084-16					
Aroclor 1016	<b>ND</b>	0.056	EPA 8082A	10-11-16	10-11-16	
Aroclor 1221	<b>ND</b>	0.056	EPA 8082A	10-11-16	10-11-16	
Aroclor 1232	<b>ND</b>	0.056	EPA 8082A	10-11-16	10-11-16	
Aroclor 1242	<b>ND</b>	0.056	EPA 8082A	10-11-16	10-11-16	
Aroclor 1248	<b>ND</b>	0.056	EPA 8082A	10-11-16	10-11-16	
Aroclor 1254	<b>ND</b>	0.056	EPA 8082A	10-11-16	10-11-16	
Aroclor 1260	<b>ND</b>	0.056	EPA 8082A	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>DCB</i>	<i>85</i>	<i>50-139</i>				



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 Project: 31008

**PCBs EPA 8082A  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1011S1					
Aroclor 1016	ND	0.050	EPA 8082A	10-11-16	10-11-16	
Aroclor 1221	ND	0.050	EPA 8082A	10-11-16	10-11-16	
Aroclor 1232	ND	0.050	EPA 8082A	10-11-16	10-11-16	
Aroclor 1242	ND	0.050	EPA 8082A	10-11-16	10-11-16	
Aroclor 1248	ND	0.050	EPA 8082A	10-11-16	10-11-16	
Aroclor 1254	ND	0.050	EPA 8082A	10-11-16	10-11-16	
Aroclor 1260	ND	0.050	EPA 8082A	10-11-16	10-11-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
DCB	86		50-139			

Analyte	Result		Spike Level		Source Result	Percent Recovery		Recovery Limits	RPD	RPD Limit	Flags
<b>MATRIX SPIKES</b>											
Laboratory ID:	10-080-01										
	MS	MSD	MS	MSD		MS	MSD				
Aroclor 1260	0.389	0.387	0.500	0.500	ND	78	77	49-133	1	17	
<i>Surrogate:</i>											
DCB						80	81	50-139			



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 Project: 31008

**TOTAL METALS**  
**EPA 6010C/6020A/7471B**

Matrix: Soil  
 Units: mg/kg (ppm)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>EPA Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
Lab ID:	10-084-05					
<b>Client ID:</b>	<b>H-5-16-13.5</b>					
Antimony	<b>ND</b>	6.9	6010C	10-13-16	10-13-16	
Arsenic	<b>ND</b>	14	6010C	10-13-16	10-13-16	
Beryllium	<b>ND</b>	0.69	6010C	10-13-16	10-13-16	
Cadmium	<b>ND</b>	0.69	6010C	10-13-16	10-13-16	
Chromium	<b>64</b>	0.69	6010C	10-13-16	10-13-16	
Copper	<b>46</b>	1.4	6010C	10-13-16	10-13-16	
Lead	<b>ND</b>	6.9	6010C	10-13-16	10-13-16	
Mercury	<b>ND</b>	0.35	7471B	10-11-16	10-11-16	
Nickel	<b>69</b>	3.5	6010C	10-13-16	10-13-16	
Selenium	<b>ND</b>	14	6010C	10-13-16	10-13-16	
Silver	<b>ND</b>	0.69	6010C	10-13-16	10-13-16	
Thallium	<b>ND</b>	1.7	6020A	10-13-16	10-17-16	
Zinc	<b>69</b>	3.5	6010C	10-13-16	10-13-16	



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**TOTAL METALS**  
**EPA 6010C/6020A/7471B**

Matrix: Soil  
 Units: mg/kg (ppm)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>EPA Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
Lab ID:	10-084-16					
Client ID:	H-4-16-16					
Antimony	ND	5.6	6010C	10-13-16	10-13-16	
Arsenic	ND	11	6010C	10-13-16	10-13-16	
Beryllium	ND	0.56	6010C	10-13-16	10-13-16	
Cadmium	ND	0.56	6010C	10-13-16	10-13-16	
Chromium	28	0.56	6010C	10-13-16	10-13-16	
Copper	11	1.1	6010C	10-13-16	10-13-16	
Lead	ND	5.6	6010C	10-13-16	10-13-16	
Mercury	ND	0.28	7471B	10-11-16	10-11-16	
Nickel	30	2.8	6010C	10-13-16	10-13-16	
Selenium	ND	11	6010C	10-13-16	10-13-16	
Silver	ND	0.56	6010C	10-13-16	10-13-16	
Thallium	ND	1.4	6020A	10-13-16	10-17-16	
Zinc	24	2.8	6010C	10-13-16	10-13-16	



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 Project: 31008

**TOTAL METALS  
 EPA 6010C/6020A/7471B  
 METHOD BLANK QUALITY CONTROL**

Date Extracted: 10-11&13-16  
 Date Analyzed: 10-11,13&17-16

Matrix: Soil  
 Units: mg/kg (ppm)

Lab ID: MB1013SH1&MB1011S1

Analyte	Method	Result	PQL
Antimony	6010C	ND	5.0
Arsenic	6010C	ND	10
Beryllium	6010C	ND	0.50
Cadmium	6010C	ND	0.50
Chromium	6010C	ND	0.50
Copper	6010C	ND	1.0
Lead	6010C	ND	5.0
Mercury	7471B	ND	0.25
Nickel	6010C	ND	2.5
Selenium	6010C	ND	10
Silver	6010C	ND	0.50
Thallium	6020A	ND	1.3
Zinc	6010C	ND	2.5



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**TOTAL METALS  
 EPA 6010C/6020A/7471B  
 DUPLICATE QUALITY CONTROL**

Date Extracted: 10-11&13-16  
 Date Analyzed: 10-11,13&17-16

Matrix: Soil  
 Units: mg/kg (ppm)

Lab ID: 10-080-01

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Antimony	ND	ND	NA	5.0	
Arsenic	ND	ND	NA	10.0	
Beryllium	ND	ND	NA	0.50	
Cadmium	ND	ND	NA	0.50	
Chromium	25.9	24.3	6	0.50	
Copper	11.7	10.9	7	1.0	
Lead	ND	ND	NA	5.0	
Mercury	ND	ND	NA	0.25	
Nickel	30.4	29.2	4	2.5	
Selenium	ND	ND	NA	10	
Silver	ND	ND	NA	0.50	
Thallium	ND	ND	NA	1.3	
Zinc	24.0	22.0	9	2.5	



Date of Report: October 19, 2016  
 Samples Submitted: October 10, 2016  
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 Project: 31008

**TOTAL METALS  
 EPA 6010C/6020A/7471B  
 MS/MSD QUALITY CONTROL**

Date Extracted: 10-11&13-16  
 Date Analyzed: 10-11,13&17-16

Matrix: Soil  
 Units: mg/kg (ppm)

Lab ID: 10-080-01

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Antimony	100	<b>95.4</b>	95	<b>88.3</b>	88	8	
Arsenic	100	<b>101</b>	101	<b>94.3</b>	94	6	
Beryllium	50.0	<b>50.8</b>	102	<b>47.6</b>	95	7	
Cadmium	50.0	<b>49.3</b>	99	<b>47.5</b>	95	4	
Chromium	100	<b>128</b>	102	<b>119</b>	93	7	
Copper	50.0	<b>63.9</b>	105	<b>60.4</b>	97	6	
Lead	250	<b>238</b>	95	<b>232</b>	93	3	
Mercury	0.500	<b>0.499</b>	100	<b>0.546</b>	109	9	
Nickel	100	<b>127</b>	96	<b>120</b>	89	6	
Selenium	100	<b>104</b>	104	<b>99.4</b>	99	5	
Silver	25.0	<b>24.4</b>	97	<b>23.1</b>	92	5	
Thallium	50.0	<b>44.2</b>	88	<b>44.8</b>	90	1	
Zinc	100	<b>120</b>	96	<b>115</b>	91	4	



Date of Report: October 19, 2016  
Samples Submitted: October 10, 2016  
Laboratory Reference: 1610-084  
Project: 31008

### % MOISTURE

Date Analyzed: 10-12-16

Client ID	Lab ID	% Moisture
H-5-16-3	10-084-01	13
H-5-16-6	10-084-02	10
H-5-16-8.5	10-084-03	31
H-5-16-11	10-084-04	31
H-5-16-13.5	10-084-05	28
H-5-16-16	10-084-06	13
H-5-16-18.5	10-084-07	12
H-4-16-3	10-084-12	9
H-4-16-6	10-084-13	12
H-4-16-8.5	10-084-14	26
H-4-16-11	10-084-15	24
H-4-16-16	10-084-16	11
H-4-16-18.5	10-084-17	15
H-4-16-19.9	10-084-18	13
H-4-16-25.4	10-084-19	17





### 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.
  - Z -
- ND - Not Detected at PQL  
 PQL - Practical Quantitation Limit  
 RPD - Relative Percent Difference









14648 NE 95<sup>th</sup> Street, Redmond, WA 98052 • (425) 883-3881

October 28, 2016

Glenn Hayman  
INNOVEX Environmental Mgt., Inc.  
16310 NE 80th St., Suite 300  
Redmond, WA 98052

Re: Analytical Data for Project 31008  
Laboratory Reference No. 1610-084B

Dear Glenn:

Enclosed are the analytical results and associated quality control data for samples submitted on October 10, 2016.

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,

A handwritten signature in black ink, appearing to read "DB", with a long horizontal stroke extending to the right.

David Baumeister  
Project Manager

Enclosures



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OnSite Environmental, Inc. 14648 NE 95<sup>th</sup> Street, Redmond, WA 98052 (425) 883-3881

This report pertains to the samples analyzed in accordance with the chain of custody, and is intended only for the use of the individual or company to whom it is addressed.

Date of Report: October 28, 2016  
Samples Submitted: October 10, 2016  
Laboratory Reference: 1610-084B  
Project: 31008

### Case Narrative

Samples were collected on October 8, 2016 and received by the laboratory on October 10, 2016. 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/BTEX Analysis

Per EPA method 5035A, samples were received by the laboratory in pre-weighed 40 ml VOA vials preserved with either Methanol or Sodium Bisulfate.

Samples H-4-16-3, H-4-16-6, H-4-16-8.5, H-4-16-11, H-4-16-19.9 and H-4-16-25.4 were extracted and analyzed outside the holding time. Some loss of volatiles might have occurred.

The surrogate recovery was above the upper control limit for sample H-4-16-6. The recovery was confirmed by reanalysis.

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.



Date of Report: October 28, 2016  
 Samples Submitted: October 10, 2016  
 Laboratory Reference: 1610-084B  
 Project: 31008

### NWTPH-Gx/BTEX

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-4-16-3</b>					
Laboratory ID:	10-084-12					
Benzene	ND	0.020	EPA 8021B	10-27-16	10-27-16	
Toluene	ND	0.055	EPA 8021B	10-27-16	10-27-16	
Ethyl Benzene	ND	0.055	EPA 8021B	10-27-16	10-27-16	
m,p-Xylene	ND	0.055	EPA 8021B	10-27-16	10-27-16	
o-Xylene	ND	0.055	EPA 8021B	10-27-16	10-27-16	
Gasoline	ND	5.5	NWTPH-Gx	10-27-16	10-27-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	109	63-124				
<b>Client ID:</b>	<b>H-4-16-6</b>					
Laboratory ID:	10-084-13					
Benzene	0.024	0.020	EPA 8021B	10-27-16	10-27-16	
Toluene	ND	0.068	EPA 8021B	10-27-16	10-27-16	
Ethyl Benzene	ND	0.068	EPA 8021B	10-27-16	10-27-16	
m,p-Xylene	ND	0.068	EPA 8021B	10-27-16	10-27-16	
o-Xylene	ND	0.068	EPA 8021B	10-27-16	10-27-16	
Gasoline	ND	6.8	NWTPH-Gx	10-27-16	10-27-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	147	63-124				Q
<b>Client ID:</b>	<b>H-4-16-8.5</b>					
Laboratory ID:	10-084-14					
Benzene	0.045	0.020	EPA 8021B	10-27-16	10-27-16	
Toluene	ND	0.077	EPA 8021B	10-27-16	10-27-16	
Ethyl Benzene	ND	0.077	EPA 8021B	10-27-16	10-27-16	
m,p-Xylene	ND	0.077	EPA 8021B	10-27-16	10-27-16	
o-Xylene	ND	0.077	EPA 8021B	10-27-16	10-27-16	
Gasoline	ND	7.7	NWTPH-Gx	10-27-16	10-27-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	103	63-124				



Date of Report: October 28, 2016  
 Samples Submitted: October 10, 2016  
 Laboratory Reference: 1610-084B  
 Project: 31008

### NWTPH-Gx/BTEX

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>H-4-16-11</b>					
Laboratory ID:	10-084-15					
Benzene	<b>0.026</b>	0.020	EPA 8021B	10-27-16	10-27-16	
Toluene	<b>ND</b>	0.069	EPA 8021B	10-27-16	10-27-16	
Ethyl Benzene	<b>ND</b>	0.069	EPA 8021B	10-27-16	10-27-16	
m,p-Xylene	<b>ND</b>	0.069	EPA 8021B	10-27-16	10-27-16	
o-Xylene	<b>ND</b>	0.069	EPA 8021B	10-27-16	10-27-16	
Gasoline	<b>ND</b>	6.9	NWTPH-Gx	10-27-16	10-27-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	<i>111</i>	<i>63-124</i>				
<b>Client ID:</b>	<b>H-4-16-19.9</b>					
Laboratory ID:	10-084-18					
Benzene	<b>0.35</b>	0.020	EPA 8021B	10-27-16	10-27-16	
Toluene	<b>0.090</b>	0.053	EPA 8021B	10-27-16	10-27-16	
Ethyl Benzene	<b>1.4</b>	0.053	EPA 8021B	10-27-16	10-27-16	
m,p-Xylene	<b>2.2</b>	0.053	EPA 8021B	10-27-16	10-27-16	
o-Xylene	<b>0.59</b>	0.053	EPA 8021B	10-27-16	10-27-16	
Gasoline	<b>99</b>	5.3	NWTPH-Gx	10-27-16	10-27-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	<i>100</i>	<i>63-124</i>				
<b>Client ID:</b>	<b>H-4-16-25.4</b>					
Laboratory ID:	10-084-19					
Benzene	<b>0.092</b>	0.020	EPA 8021B	10-27-16	10-27-16	
Toluene	<b>0.064</b>	0.060	EPA 8021B	10-27-16	10-27-16	
Ethyl Benzene	<b>ND</b>	0.060	EPA 8021B	10-27-16	10-27-16	
m,p-Xylene	<b>0.088</b>	0.060	EPA 8021B	10-27-16	10-27-16	
o-Xylene	<b>ND</b>	0.060	EPA 8021B	10-27-16	10-27-16	
Gasoline	<b>ND</b>	6.0	NWTPH-Gx	10-27-16	10-27-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	<i>105</i>	<i>63-124</i>				



Date of Report: October 28, 2016  
 Samples Submitted: October 10, 2016  
 Laboratory Reference: 1610-084B  
 Project: 31008

**NWTPH-Gx/BTEX  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1027S2					
Benzene	ND	0.020	EPA 8021B	10-27-16	10-27-16	
Toluene	ND	0.050	EPA 8021B	10-27-16	10-27-16	
Ethyl Benzene	ND	0.050	EPA 8021B	10-27-16	10-27-16	
m,p-Xylene	ND	0.050	EPA 8021B	10-27-16	10-27-16	
o-Xylene	ND	0.050	EPA 8021B	10-27-16	10-27-16	
Gasoline	ND	5.0	NWTPH-Gx	10-27-16	10-27-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	92	63-124				

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>								
Laboratory ID:	10-084-19							
	ORIG	DUP						
Benzene	0.0766	0.0696	NA	NA	NA	NA	10	30
Toluene	0.0534	ND	NA	NA	NA	NA	NA	30
Ethyl Benzene	ND	ND	NA	NA	NA	NA	NA	30
m,p-Xylene	0.0734	0.0621	NA	NA	NA	NA	17	30
o-Xylene	ND	ND	NA	NA	NA	NA	NA	30
Gasoline	ND	ND	NA	NA	NA	NA	NA	30
<i>Surrogate:</i>								
<i>Fluorobenzene</i>				105	105	63-124		

**SPIKE BLANKS**

Laboratory ID:	SB	SBD	SB	SBD	SB	SBD			
SB1027S1									
Benzene	0.934	0.951	1.00	1.00	93	95	70-124	2	12
Toluene	0.932	0.957	1.00	1.00	93	96	73-119	3	12
Ethyl Benzene	0.952	0.972	1.00	1.00	95	97	74-117	2	12
m,p-Xylene	0.900	0.923	1.00	1.00	90	92	75-117	3	13
o-Xylene	0.935	0.951	1.00	1.00	94	95	75-116	2	12
<i>Surrogate:</i>									
<i>Fluorobenzene</i>					98	93	63-124		





### Data Qualifiers and Abbreviations

- A - Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
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  - 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.
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  - 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.
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  - 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.
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- ND - Not Detected at PQL  
 PQL - Practical Quantitation Limit  
 RPD - Relative Percent Difference





