

**Phase II Environmental Site Assessment**

A-1 Towing  
King County Tax Parcel 3223049048  
Des Moines, Washington

*for*  
**Washington State Department of Transportation  
c/o HNTB Corporation**

November 25, 2022



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File No. 0180-393-00

November 25, 2022

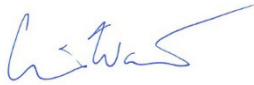
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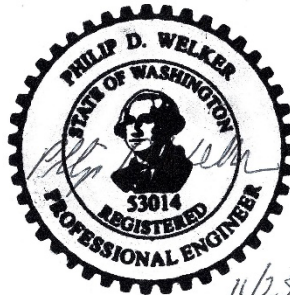
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## 1.0 INTRODUCTION AND BACKGROUND

This report describes the Phase II Environmental Site Assessment (ESA) activities conducted at the property located at King County Tax Parcel 3223049048 (the provided address of 18451 12<sup>th</sup> Avenue South in Des Moines, Washington does not map accurately) as part of the expansion of SR-509 (herein referred to as “site”). The property is owned by Washington State Department of Transportation (WSDOT) but is leased by a private party and WSDOT is evaluating environmental conditions. The approximate site location is shown in the attached Vicinity Map, Figure 1.

GeoEngineers, Inc. (GeoEngineers) reviewed a Phase II ESA prepared for the property (known as A-1 Towing and the Foreman Property) located at 18451 12<sup>th</sup> Avenue South in Des Moines, Washington by WSDOT, dated January 24, 2019. The Phase II ESA indicated that WSDOT purchased the property in 2000. A tenant (Mr. Foreman) conducted welding activities between 2000 and 2016, and the Washington State Patrol (WSP) used a portion of the property to dismantle seized vehicles in the summer of 2016. At the time of the report (January 2019), A-1 Towing leased the property as a towing yard, which included storage of vehicles throughout the site and repairing of vehicles within an existing garage structure. The Phase II report included a summary of the following activities:

- The Phase II ESA performed by WSDOT in 2016 following the termination of the lease to Mr. Foreman included soil sampling of areas with significant staining. Areas identified as Area A, B, C and D were sampled as part of the Phase II ESA. Results indicated the presence of lube oil and arsenic exceeding the Model Toxic Control Act (MTCA) Cleanup Levels (CULs) in Area B and C, respectively.
- In April 2017, WSDOT contracted Marine Vacuum Incorporated to clean up areas of the Site with trash, debris and soils with notable detectable analytes below MTCA Method A CULs. It is unclear from this statement as to what soils were actually removed and no additional details were provided.
- In January 2018, A-1 Towing graded Site soils including some that were contaminated. The soils were displaced from previously sampled areas A and B. A new stockpile was present following these activities just north of Area D.
- On March 15, 2018, WSDOT collected four additional soil samples from areas defined as Area A, B, and a new location. WSDOT collected a sample to determine if there was an indication of grading with known or suspected soil contamination, or where displaced/graded soils appeared to be consolidated into a stockpile generated by the lessee. Areas C and D did not have these indicators and did not warrant additional sampling. Sample results did not indicate results exceeding the MTCA Method A CULs.
- On September 20 and 21, 2018, Marine Vacuum Performed excavation activities to address the previously detected lube oil concentrations exceeding the MTCA Method A CULs and an area to remediate suspected contaminated materials in Area C and D. During site excavation activities soils were field screened using sight, olfactory and photoionization detector (PID) levels above 10 parts per million. Non-suspect soils were placed off to the side on a lined area and covered. On September 21, 2018, WSDOT collected two confirmation soil samples in locations where the soils had been removed. Laboratory analytical results were either non-detect or detected at concentrations less than MTCA Method A CULs. Following sample results, the stockpiled non-suspect soils were used to fill and grade the excavation areas.

- The January 24, 2019, Phase II ESA summary indicated that the known and suspected soils with contamination (approximately 8.4 cubic yards) were removed from the site and based on the analytical results no contamination remains on the property. WSDOT requested a no further action (NFA) determination from Washington State.
- Ecology responded to the request for a NFA determination indicating that additional information would be necessary including depth of samples, groundwater samples (if encountered) and more accurate sample location information (previous report had conflicting depictions).

The following sections describe our scope of services (prepared with approval by WSDOT), field activities, sampling results and opinions.

## 2.0 SCOPE OF SERVICES

Our general environmental scope of services was as follows:

- Prepared a Health and Safety Plan for use by GeoEngineers' personnel during the field activities.
- Coordinate the location of nine shallow proposed soil borings with one boring located at each historic background sample location SS-2, SS-3 and SS-4, which represent staining observed in Areas B, C and D, respectively (Area A does not appear to be of concern for additional assessment). Two additional borings were placed in each area based on the sample location map depicting rectangular areas of staining and or remedial excavations. At least one boring in each area will be advanced to groundwater or 15 feet below ground surface (bgs) and a groundwater monitoring well installed, if water is encountered. The additional borings will be placed in a manner to bracket the depicted area of concern. A total of nine shallow borings and three monitoring wells were proposed.
- Contact the Utility Notification Center public locating service. A private locating service will also be subcontracted to locate potential utilities.
- Retain a drilling subcontractor to advance nine direct push borings to a depth of 5 to 15 feet (bgs), dependent on field screening results, presence of groundwater and depth to native soils. If groundwater is encountered (reportedly as shallow as 3 feet bgs) at the former sample locations SS-2 through SS-4, a 2-inch-diameter polyvinyl chloride (PVC) casing monitoring well will be installed. Depths of the borings may be adjusted in an attempt to install a minimum of three wells to evaluate site specific groundwater flow direction. (Groundwater was not encountered during drilling activities and no wells were installed).
- Continuously sample push probes and visually classify and field screen (sheen, PID and odor) for potential laboratory analysis by a GeoEngineers representative. Visual classification of the soil samples is to be completed in general accordance with ASTM D 2488-90, the Standard Practice for Description and Identification of Soils.
- Collected two soil samples per boring for potential laboratory analysis based on field screening results and depth to groundwater.
- Submit the soil samples on a standard turnaround time (10 business days) for laboratory analysis of the following:
  - Gasoline-range petroleum hydrocarbons by Northwest Method NWTPH-Gx (Gx).

- Diesel- and lube-oil range petroleum hydrocarbons by Northwest Method NWTPH-Dx (Dx and Ox).
  - Resource Conservation and Recovery Act (RCRA) eight metals (arsenic, barium, cadmium, chromium, lead, mercury, selenium and silver) by United States Environmental Protection Agency (EPA) Method 6020.
  - Polycyclic aromatic hydrocarbons (PAHs) by EPA Method 8270.
  - Volatile organic compounds (VOCs) by EPA Method 8260C.
- Coordinate disposal of investigation-derived waste (IDW) at an appropriate off-site facility. Collect one composite soil sample directly from the drum or during drilling for waste profiling purposes.
  - Prepare a report including sample results tables, laboratory data, site figures and boring/sampling logs documenting results of the work performed at the site. The data will be tabulated and compared against MTCA CULs.
  - Upload data to the Ecology Environmental Information Management (EIM) system as required.

### **3.0 SUBSURFACE CONDITIONS**

Logs of borings are included in Appendix A. Soils were visually classified in general accordance with ASTM D 2488-94. Boring locations are shown in Exploration Locations, Figure 2.

#### **3.1. Subsurface Soil Conditions**

Surface conditions at the site consist primarily of fill gravel (from several inches up to 2 feet). Soils encountered during exploration activities consisted of sand, silt and varying gradations of sand and silt with occasional gravel to the maximum depth explored, 15 to 20 feet bgs.

#### **3.2. Subsurface Groundwater Conditions**

Soils observed during drilling were generally moist and groundwater was not encountered during drilling.

### **4.0 SOIL SAMPLING AND CHEMICAL ANALYTICAL RESULTS**

A GeoEngineers' field representative collected soil samples continuously from each boring for field screening and potential laboratory analysis. Samples were field screened for evidence of petroleum and/or VOC-related impact using visual, water sheen and headspace vapor screening methods using a PID. Soil samples were collected based on field screening data. The soil samples were submitted for laboratory analysis of VOCs, Gx, Dx, Ox, PAHs and/or RCRA eight metals to Pace National. Soil chemical analytical data are presented in Tables 1 and 2. Laboratory analytical reports are attached in Appendix B.

#### **4.1. Soil Sample Analytical Results**

As previously noted, three areas of concern with previous sample results were identified for additional assessment identified as areas B, C and D (no concerns were expressed by Ecology for the need to further explore Area A). The following sections summarize the results of the additional borings and soil sampling performed in each area. Approximate locations of previous sampling performed by others and the sampling performed during this assessment are shown on Figure 2.

#### 4.1.1. Area B

Area B is located off the southwest corner of the existing structure. Previous sample SS-2 indicated levels of arsenic (35 milligrams/kilogram [mg/kg]). Three borings (GEI-3, GEI-4 and MW-2) were performed within the identified Area B. Boring MW-2 was intended to be in the approximate location of previous sample SS-2 and planned to be installed as a monitoring well, but no groundwater was encountered to the maximum depth explored (15 feet bgs). Field screening the three boring locations did not indicate petroleum compounds or VOCs. One sample was analyzed from each boring. Soil sample results for the samples collected from each boring are as follows:

- Sample GEI-3 at 9-10 feet bgs was non-detect for Gx, Dx, Ox, VOCs and PAHs. Detected RCRA eight metals were at concentrations less than the MTCA Method A Cleanup Levels for Unrestricted Land Use.
- Sample GEI-4 at 2-3 feet bgs was non-detect for Gx, Dx and VOCs. Ox and PAHs and detected RCRA eight metals were at concentrations less than the MTCA Method A Cleanup Levels for Unrestricted Land Use.
- Sample MW-2 at 9-10 feet bgs was non-detect for Gx, Dx, Ox, VOCs and PAHs. Detected RCRA eight metals were at concentrations less than the MTCA Method A Cleanup Levels for Unrestricted Land Use.

#### 4.1.2. Area C

Area C is located off the east side of the of the existing structure. Previous sample SS-3 indicated levels of lube oil at 31,000 mg/kg. Three borings (GEI-1, GEI-2 and MW-1) were performed within the identified Area C. Boring MW-1 was intended to be in the approximate location of previous sample SS-3 and planned to be installed as a monitoring well, but no groundwater was encountered to the maximum depth explored (20 feet bgs at boring GEI-1). Field screening the three boring locations did not indicate petroleum compounds or VOCs. One sample was analyzed from each boring. Soil sample results for the samples collected from each boring are as follows:

- Sample GEI-1 at 9-10 feet bgs was non-detect for Gx, Dx, Ox, VOCs and PAHs. Detected RCRA eight metals were at concentrations less than the MTCA Method A Cleanup Levels for Unrestricted Land Use.
- Sample GEI-2 at 9-10 feet bgs was non-detect for Gx, Dx, Ox, VOCs and PAHs. Detected RCRA eight metals were at concentrations less than the MTCA Method A Cleanup Levels for Unrestricted Land Use.
- Sample MW-1 at 2-3 feet bgs was non-detect for Gx, Dx, Ox, VOCs and PAHs. Detected RCRA eight metals were at concentrations less than the MTCA Method A Cleanup Levels for Unrestricted Land Use.

#### 4.1.3. Area D

Area D is located off the north side of the existing structure. Previous observations indicated the presence of staining and groundwater in the exploration, which was not sampled. Three borings (GEI-4, GEI-6 and MW-3) were performed within the identified Area C. Boring MW-3 was intended to be in the approximate location of previous sample SS-4 and staining. MW-3 was planned to be installed as a monitoring well, but no groundwater was encountered to the maximum depth explored (15 feet bgs). Field screening the three boring locations did not indicate petroleum compounds or VOCs. One sample was analyzed from each boring. Soil sample results for the samples collected from each boring are as follows:

- Sample GEI-5 at 7-8 feet bgs was non-detect for Gx, Dx, Ox, VOCs and PAHs. Detected RCRA eight metals were at concentrations less than the MTCA Method A Cleanup Levels for Unrestricted Land Use.

- Sample GEI-6 at 7-8 feet bgs was non-detect for Gx, Dx, Ox, VOCs and PAHs. Detected RCRA eight metals were at concentrations less than the MTCA Method A Cleanup Levels for Unrestricted Land Use.
- Sample MW-3 at 2-3 feet bgs was non-detect for Gx, VOCs and PAHs. Detected Dx, Ox and RCRA eight metals were at concentrations less than the MTCA Method A Cleanup Levels for Unrestricted Land Use.

## 5.0 CONCLUSIONS

GeoEngineers performed additional sampling at three areas (Area B, C and D) located around the existing structure on the A-1 towing site to further evaluate the potential presence of contaminants related to past uses at the site. Results of the assessment did not identify concentrations of Gx, Dx, Ox, VOCs, PAHs and or RCRA eight metals at concentrations exceeding the MTCA Method A Cleanup Levels. Out of the nine soil samples analyzed, only two identified low concentrations of Dx, Ox and/or PAHs. Groundwater was not encountered at the maximum depth explored at the site of 15 to 20 feet bgs.

Based on the results of the assessment a No Further Action determination should be requested from Ecology.

## 6.0 LIMITATIONS

We have prepared this report for the exclusive use of WSDOT. Within the limitations of scope, schedule and budget, our services have been executed in accordance with generally accepted environmental science practices in this area at the time this report was prepared. The conclusions and opinions presented in this report are based on our professional knowledge, judgment, experience and discrete sample collection. Contaminants may be present in areas of the site not sampled or tested. No warranty or other conditions, express or implied, should be understood.

Please refer to Appendix C, Report Limitations and Guidelines for Use, for additional information regarding the use of this report.



**Table 1**

**Soil Chemical Analytical Results - Petroleum Hydrocarbons and VOCs  
A-1 Towing  
SeaTac, Washington**

Boring/ Well ID	Sample ID	Date Collected	Depth (Feet Below Ground Surface)	Field Screening Observation		NWTPH-Gx (mg/kg)	NWTPH-Dx (mg/kg)			VOCs EPA Method 8260 (mg/kg)	Anthracene	Acenaphthene	Acenaphthylene	Benzo (a) anthracene	Benzo (a) pyrene	Benzo (b) fluoranthene	Benzo(g,h,i)p erylene	Benzo (k) fluoranthene	Chrysene														
				Sheen	PID (ppm)		Gasoline	Diesel	Heavy Oil																								
GEI-1	GEI-1 (9-10')	10/25/2021	9-10	NS	<1	<3.01	<	4.4	<	10.9	ND	<	0.00654	<	0.00654	<	0.00654	<	0.00654	<	0.00654	<	0.00654	<	0.00654	<	0.00654	<	0.00654	<	0.00654	<	0.00654
GEI-2	GEI-2 (9-10')	10/25/2021	9-10	NS	<1	<2.99	<	4.34	<	10.80	ND	<	0.00650	<	0.00650	<	0.00650	<	0.00650	<	0.00650	<	0.00650	<	0.00650	<	0.00650	<	0.00650	<	0.00650	<	0.00650
GEI-3	GEI-3 (9-10')	10/25/2021	9-10	NS	<1	<3.23	<	4.50	<	11.2	ND	<	0.00675	<	0.00675	<	0.00675	<	0.00675	<	0.00675	<	0.00675	<	0.00675	<	0.00675	<	0.00675	<	0.00675	<	0.00675
GEI-4	GEI-4 (2-3')	10/25/2021	2-3	NS	<1	<2.97	<	21.8		<b>166</b>	ND	<	0.00655	<	0.00655	<	0.00655	<b>0.0298</b>	<b>0.0363</b>	<b>0.0500</b>	<b>0.0306</b>	<b>0.0150</b>	<b>0.0354</b>										
GEI-5	GEI-5 (7-8')	10/25/2021	7-8	NS	<1	<3.02	<	4.36	<	10.9	ND	<	0.00654	<	0.00654	<	0.00654	<	0.00654	<	0.00654	<	0.00654	<	0.00654	<	0.00654	<	0.00654	<	0.00654	<	0.00654
GEI-6	GEI-6 (7-8')	10/25/2021	7-8	NS	<1	<2.95	<	4.33	<	10.8	ND	<	0.00649	<	0.00649	<	0.00649	<	0.00649	<	0.00649	<	0.00649	<	0.00649	<	0.00649	<	0.00649	<	0.00649	<	0.00649
MW-1	MW-1 (2-3')	10/25/2021	2-3	NS	<1	<3.43	<	4.69	<	11.7	ND	<	0.00703	<	0.00703	<	0.00703	<	0.00703	<	0.00703	<	0.00703	<	0.00703	<	0.00703	<	0.00703	<	0.00703	<	0.00703
MW-2	MW-2 (9-10')	10/25/2021	9-10	NS	<1	<3.08	<	4.43	<	11.1	ND	<	0.00665	<	0.00665	<	0.00665	<	0.00665	<	0.00665	<	0.00665	<	0.00665	<	0.00665	<	0.00665	<	0.00665	<	0.00665
MW-3	MW-3 (2-3')	10/25/2021	2-3	NS	<1	<3.35		<b>5.55</b>		<b>24.3</b>	ND	<	0.00691	<	0.00691	<	0.00691	<	0.00691	<	0.00691	<	0.00691	<	0.00691	<	0.00691	<	0.00691	<	0.00691	<	0.00691
<b>MTCA Method A Cleanup Levels for Unrestricted Land Use</b>						800/1,000	2,000			Various	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-

Boring/ Well ID	Sample ID	Date Collected	Depth (Feet Below Ground Surface)	Field Screening Observation		NWTPH-Gx (mg/kg)	NWTPH-Dx (mg/kg)			VOCs EPA Method 8260 (mg/kg)									Total cPAHs <sup>1</sup> (mg/kg)
				Sheen	PID (ppm)	Gasoline	Diesel	Heavy Oil	Various	Dibenz(a,h)anthracene	Fluoranthene	Fluorene	Indeno (1,2,3-cd) pyrene	Naphthalene	Phenanthrene	Pyrene	1-Methyl-naphthalene	2-Methyl-naphthalene	TEQ
GEI-1	GEI-1 (9-10')	10/25/2021	9-10	NS	<1	<3.01	< 4.4	< 10.9	ND	< 0.00654	< 0.00654	< 0.00654	< 0.00654	< 0.0218	< 0.00654	< 0.00654	< 0.0218	< 0.0128	ND
GEI-2	GEI-2 (9-10')	10/25/2021	9-10	NS	<1	<2.99	< 4.34	< 10.80	ND	< 0.00650	< 0.00650	< 0.00650	< 0.00650	< 0.0217	< 0.00650	< 0.00650	< 0.0217	< 0.0217	ND
GEI-3	GEI-3 (9-10')	10/25/2021	9-10	NS	<1	<3.23	< 4.50	< 11.2	ND	< 0.00675	< 0.00675	< 0.00675	< 0.00675	< 0.0225	< 0.00675	< 0.00675	< 0.0225	< 0.0225	ND
GEI-4	GEI-4 (2-3')	10/25/2021	2-3	NS	<1	<2.97	< 22	<b>166</b>	ND	<b>0.00660</b>	<b>0.0622</b>	< 0.0066	<b>0.0325</b>	< 0.0218	<b>0.0244</b>	<b>0.0474</b>	< 0.0218	< 0.0218	<b>0.0223</b>
GEI-5	GEI-5 (7-8')	10/25/2021	7-8	NS	<1	<3.02	< 4.4	< 10.9	ND	< 0.00654	< 0.00654	< 0.00654	< 0.00654	< 0.02180	< 0.00654	< 0.00654	< 0.0218	< 0.0218	ND
GEI-6	GEI-6 (7-8')	10/25/2021	7-8	NS	<1	<2.95	< 4.3	< 10.8	ND	< 0.00649	< 0.00649	< 0.00649	< 0.00649	< 0.0216	< 0.00649	< 0.00649	< 0.0216	< 0.0216	ND
MW-1	MW-1 (2-3')	10/25/2021	2-3	NS	<1	<3.43	< 4.7	< 11.7	ND	< 0.00703	< 0.00703	< 0.00703	< 0.00703	< 0.02340	< 0.00703	< 0.00703	< 0.0234	< 0.0234	ND
MW-2	MW-2 (9-10')	10/25/2021	9-10	NS	<1	<3.08	< 4.43	< 11.1	ND	< 0.00665	< 0.00665	< 0.00665	< 0.00665	< 0.0222	< 0.00665	< 0.00665	< 0.0222	< 0.0222	ND
MW-3	MW-3 (2-3')	10/25/2021	2-3	NS	<1	<3.35	<b>5.6</b>	<b>24.3</b>	ND	< 0.00691	< 0.00691	< 0.00691	< 0.00691	< 0.023	< 0.00691	< 0.00691	< 0.023	< 0.023	ND
<b>MTCA Method A Cleanup Levels for Unrestricted Land Use</b>						800/1,000	2,000		Various		--	--	--	5	--	--	--	--	0.1

**Notes:**

<sup>1</sup>Carcinogenic Polycyclic Aromatic Hydrocarbons (cPAHs) analyzed by EPA Method 8270D/SIM. Total cPAHs calculated using the toxicity equivalency (TEQ) methodology specified in Washington Administrative Code (WAC) 173-340-780(8); cPAHs that were not detected were assigned half the value of the detection limit for these calculations. Not detected (ND) is indicated for samples which no cPAHs were detected.

<3.01 = analyte not detected at or less than the presented concentration

**Bold** = analyte was detected at a concentration above the laboratory reported detection limit

-- = not analyzed/no cleanup level

mg/kg = milligrams per kilogram; VOCs = volatile organic compounds

NS = no sheen; ppm = parts per million

PID = Photoionization Detector

ND = Not Detected

TEQ = toxic equivalency

EPA = Environmental Protection Agency

NWTPH = Northwest Total Petroleum Hydrocarbons

MTCA = Model Toxics Control Act

**Table 2**  
**Soil Chemical Analytical Results - Metals**  
**A-1 Towing**  
**SeaTac, Washington**

Boring/ Well ID	Sample ID	Date Collected	Depth (Feet Below Ground Surface)	Metals EPA Method 6010D/7471B (mg/kg)							
				Arsenic	Barium	Cadmium	Chromium	Lead	Mercury	Selenium	Silver
GEI-1	GEI-1 (9-10')	10/25/2021	9-10	<b>2.26</b>	<b>41.9</b>	< 1.09	<b>23.4</b>	< 2.2	< 0.0436	< 2.72	< 0.545
GEI-2	GEI-2 (9-10')	10/25/2021	9-10	<b>2.36</b>	<b>45.2</b>	< 1.08	<b>21.5</b>	< 2.17	< 0.0434	< 2.71	< 0.542
GEI-3	GEI-3 (9-10')	10/25/2021	9-10	<b>2.93</b>	<b>25.3</b>	< 1.12	<b>11.8</b>	< 2.25	< 0.0450	< 2.81	< 0.562
GEI-4	GEI-4 (2-3')	10/25/2021	2-3	<b>6.28</b>	<b>72.2</b>	< 1.09	<b>33.9</b>	<b>27.6</b>	<b>0.0731</b>	< 2.73	< 0.546
GEI-5	GEI-5 (7-8')	10/25/2021	7-8	<b>2.58</b>	<b>39.6</b>	< 1.09	<b>21.1</b>	< 2.18	< 0.0436	< 2.73	< 0.545
GEI-6	GEI-6 (7-8')	10/25/2021	7-8	<b>2.31</b>	<b>35.1</b>	< 1.08	<b>16.3</b>	< 2.16	< 0.0433	< 2.71	< 0.541
MW-1	MW-1 (2-3')	10/25/2021	2-3	<b>3.70</b>	<b>61.0</b>	< 1.17	<b>26.7</b>	< 2.34	< 0.0469	< 2.93	< 0.0586
MW- 2	MW-2 (9-10')	10/25/2021	9-10	<b>3.17</b>	<b>44.9</b>	< 1.11	<b>24.4</b>	< 2.22	< 0.0443	< 2.77	< 0.554
MW- 3	MW-3 (2-3')	10/25/2021	2-3	<b>7.77</b>	<b>58.8</b>	< 1.15	<b>41.6</b>	<b>3.16</b>	< 0.0460	< 2.88	< 0.575
<b>MTCA Method A Cleanup Levels for Unrestricted Land Use</b>				21	--	2	2,000	250	2	400	400
<b>MTCA Cleanup Level for the Protection of Groundwater</b>				--	1,600	--	--	--	--	5	14

**Notes:**

<1.09 = analyte not detected

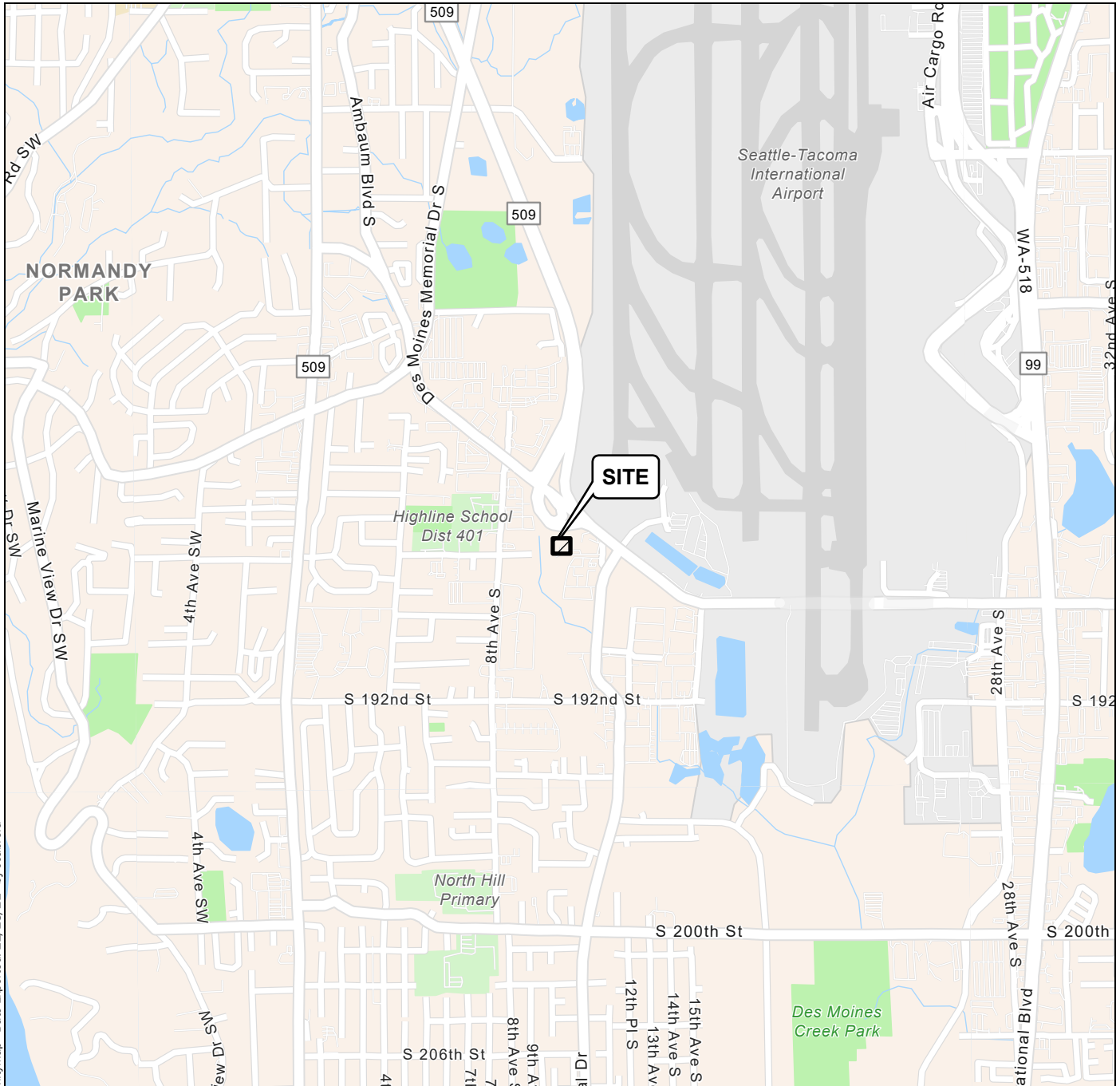
**Bold** = analyte was detected at a concentration above the laboratory method detection limit

mg/kg = milligrams per kilogram

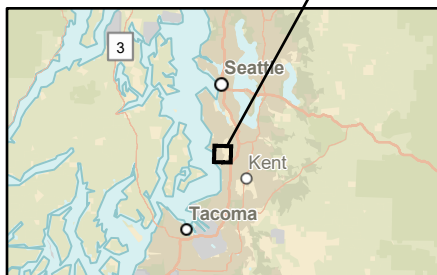
EPA = Environmental Protection Agency

MTCA = Model Toxics Control Act





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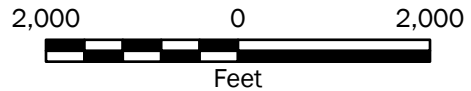
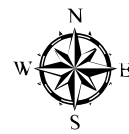


**Notes:**

1. The locations of all features shown are approximate.
2. This drawing is for information purposes. It is intended to assist in showing features discussed in an attached document. GeoEngineers, Inc. cannot guarantee the accuracy and content of electronic files. The master file is stored by GeoEngineers, Inc. and will serve as the official record of this communication.

Data Source: ESRI

Projection: NAD 1983 UTM Zone 10N



**Vicinity Map**

WSDOT - A-1 Towing  
SeaTac, Washington



**Figure 1**



P:\0\_0180393\GIS\0180393\Project\0180393\Project.aprx\018039300\_F02\_ExplorationLocations Date Exported: 12/01/21 by ccabrera






**Notes:**

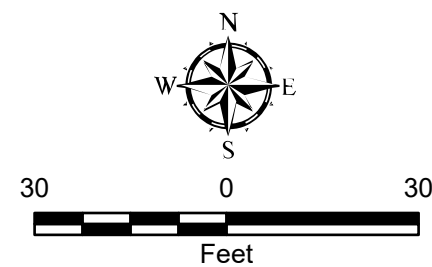
1. The locations of all features shown are approximate.
2. This drawing is for information purposes. It is intended to assist in showing features discussed in an attached document. GeoEngineers, Inc. cannot guarantee the accuracy and content of electronic files. The master file is stored by GeoEngineers, Inc. and will serve as the official record of this communication.


Data Source: ESRI Clarity

Projection: NAD 1983 UTM Zone 10N

**Legend**

-  Boring Number and Approximate Location
-  Monitoring Well Number and Approximate Location
-  Previous WSDOT Explorations (2016)
-  Drum Storage
-  Assessment Area



<b>Exploration Locations</b>	
WSDOT - A-1 Towing SeaTac, Washington	
	<b>Figure 2</b>



**APPENDIX A**  
**Boring Logs**

## SOIL CLASSIFICATION CHART

MAJOR DIVISIONS			SYMBOLS		TYPICAL DESCRIPTIONS
			GRAPH	LETTER	
COARSE GRAINED SOILS	GRAVEL AND GRAVELLY SOILS	CLEAN GRAVELS <small>(LITTLE OR NO FINES)</small>		<b>GW</b>	WELL-GRADED GRAVELS, GRAVEL - SAND MIXTURES
		GRAVELS WITH FINES <small>(APPRECIABLE AMOUNT OF FINES)</small>		<b>GP</b>	POORLY-GRADED GRAVELS, GRAVEL - SAND MIXTURES
		GRAVELS WITH FINES <small>(APPRECIABLE AMOUNT OF FINES)</small>		<b>GM</b>	SILTY GRAVELS, GRAVEL - SAND - SILT MIXTURES
	SAND AND SANDY SOILS	CLEAN SANDS <small>(LITTLE OR NO FINES)</small>		<b>SW</b>	WELL-GRADED SANDS, GRAVELLY SANDS
		SANDS WITH FINES <small>(APPRECIABLE AMOUNT OF FINES)</small>		<b>SP</b>	POORLY-GRADED SANDS, GRAVELLY SAND
		SANDS WITH FINES <small>(APPRECIABLE AMOUNT OF FINES)</small>		<b>SM</b>	SILTY SANDS, SAND - SILT MIXTURES
FINE GRAINED SOILS	SILTS AND CLAYS	LIQUID LIMIT LESS THAN 50		<b>ML</b>	INORGANIC SILTS, ROCK FLOUR, CLAYEY SILTS WITH SLIGHT PLASTICITY
				<b>CL</b>	INORGANIC CLAYS OF LOW TO MEDIUM PLASTICITY, GRAVELLY CLAYS, SANDY CLAYS, SILTY CLAYS, LEAN CLAYS
				<b>OL</b>	ORGANIC SILTS AND ORGANIC SILTY CLAYS OF LOW PLASTICITY
	SILTS AND CLAYS	LIQUID LIMIT GREATER THAN 50		<b>MH</b>	INORGANIC SILTS, MICACEOUS OR DIATOMACEOUS SILTY SOILS
				<b>CH</b>	INORGANIC CLAYS OF HIGH PLASTICITY
				<b>OH</b>	ORGANIC CLAYS AND SILTS OF MEDIUM TO HIGH PLASTICITY
HIGHLY ORGANIC SOILS				<b>PT</b>	PEAT, HUMUS, SWAMP SOILS WITH HIGH ORGANIC CONTENTS

NOTE: Multiple symbols are used to indicate borderline or dual soil classifications

### Sampler Symbol Descriptions

	2.4-inch I.D. split barrel
	Standard Penetration Test (SPT)
	Shelby tube
	Piston
	Direct-Push
	Bulk or grab
	Continuous Coring

Blowcount is recorded for driven samplers as the number of blows required to advance sampler 12 inches (or distance noted). See exploration log for hammer weight and drop.

"P" indicates sampler pushed using the weight of the drill rig.

"WOH" indicates sampler pushed using the weight of the hammer.

NOTE: The reader must refer to the discussion in the report text and the logs of explorations for a proper understanding of subsurface conditions. Descriptions on the logs apply only at the specific exploration locations and at the time the explorations were made; they are not warranted to be representative of subsurface conditions at other locations or times.

## ADDITIONAL MATERIAL SYMBOLS

SYMBOLS		TYPICAL DESCRIPTIONS
GRAPH	LETTER	
	<b>AC</b>	Asphalt Concrete
	<b>CC</b>	Cement Concrete
	<b>CR</b>	Crushed Rock/Quarry Spalls
	<b>SOD</b>	Sod/Forest Duff
	<b>TS</b>	Topsoil

### Groundwater Contact



Measured groundwater level in exploration, well, or piezometer



Measured free product in well or piezometer

### Graphic Log Contact

Distinct contact between soil strata

Approximate contact between soil strata

### Material Description Contact

Contact between geologic units

Contact between soil of the same geologic unit

### Laboratory / Field Tests

%F	Percent fines
%G	Percent gravel
AL	Atterberg limits
CA	Chemical analysis
CP	Laboratory compaction test
CS	Consolidation test
DD	Dry density
DS	Direct shear
HA	Hydrometer analysis
MC	Moisture content
MD	Moisture content and dry density
Mohs	Mohs hardness scale
OC	Organic content
PM	Permeability or hydraulic conductivity
PI	Plasticity index
PL	Point load test
PP	Pocket penetrometer
SA	Sieve analysis
TX	Triaxial compression
UC	Unconfined compression
VS	Vane shear

### Sheen Classification

NS	No Visible Sheen
SS	Slight Sheen
MS	Moderate Sheen
HS	Heavy Sheen

## Key to Exploration Logs

Start Drilled	10/25/2021	End	10/25/2021	Total Depth (ft)	20	Logged By	DC	Checked By	CJW	Driller	Holocene Drilling, Inc.	Drilling Method	Direct-push
Surface Elevation (ft) Vertical Datum	Undetermined					Hammer Data	Drilling Equipment Geoprobe 7822T						
Easting (X) Northing (Y)						System Datum	Groundwater not observed at time of exploration						
Notes:													

Elevation (feet)	FIELD DATA					Graphic Log	Group Classification	MATERIAL DESCRIPTION	Sheen	Headspace Vapor (ppm)	REMARKS
	Depth (feet)	Interval Recovered (in)	Blows/foot	Collected Sample	Sample Name Testing						
0	60					SP-SM	Dark brown silty fine to coarse sand with occasional gravel (loose, moist)	NS	<1		
						ML	Brown sandy silt				
						SM	Brown silty fine to medium sand				
						ML	Brown sandy silt				
5	60					SP-SM	Brown silty fine to medium sand with occasional gravel				
						ML	Brown sandy silt				
10	60					SP-SM	Brown silty fine to medium sand with occasional gravel				
						ML	Gray sandy silt (dense, moist)				
15	60										
20											

Note: See Figure A-1 for explanation of symbols.  
Coordinates Data Source: Horizontal approximated based on . Vertical approximated based on .

### Log of Boring GEI-1



Project: A-1 Towing  
Project Location: SeaTac, Washington  
Project Number: 0180-393-00

Figure A-2  
Sheet 1 of 1

Date: 11/18/21 Path: P:\04180393\GINT\018039300.GPJ DBL\Library\GEOENGINEERS\_DF\_STD\_US\_JUNE\_2017.GLB\GEB\_ENVIRONMENTAL\_STANDARD\_NO\_GW

Start Drilled	10/25/2021	End	10/25/2021	Total Depth (ft)	15	Logged By	DC	Checked By	CJW	Driller	Holocene Drilling, Inc.	Drilling Method	Direct-push
Surface Elevation (ft) Vertical Datum	Undetermined					Hammer Data	Geoprobe 7822T						
Easting (X) Northing (Y)						System Datum	Groundwater not observed at time of exploration						
Notes:													


Elevation (feet)	FIELD DATA					Graphic Log	Group Classification	MATERIAL DESCRIPTION	Sheen	Headspace Vapor (ppm)	REMARKS
	Depth (feet)	Interval Recovered (in)	Blows/foot	Collected Sample	Sample Name Testing						
0	60					SP	Brown silty fine to medium coarse sand (loose, moist)	NS	<1		
						ML	Brown sandy silt (loose, moist)				
5	60					SM	Brown silty fine to medium sand with occasional gravel (loose, moist)				
10	36										
15						SM	Gray silty fine to medium sand (loose, moist)				

GEI-2 (9-10)

GEI-2 (14-15)

Note: See Figure A-1 for explanation of symbols.  
Coordinates Data Source: Horizontal approximated based on . Vertical approximated based on .

Date: 11/18/21 Path: P:\01480393\GINT\018039300.GPJ DBLibrary\Library\GEOENGINEERS\_DF\_STD\_US\_JUNE\_2017.GLB\GEB\_ENVIRONMENTAL\_STANDARD\_NO\_GW

<b>Log of Boring GEI-2</b>		
	Project: A-1 Towing	Figure A-3 Sheet 1 of 1
	Project Location: SeaTac, Washington	
	Project Number: 0180-393-00	

Start Drilled	10/25/2021	End	10/25/2021	Total Depth (ft)	15	Logged By	DC	Checked By	CJW	Driller	Holocene Drilling, Inc.	Drilling Method	Direct-push	
Surface Elevation (ft) Vertical Datum	Undetermined					Hammer Data				Drilling Equipment	Geoprobe 7822T			
Easting (X) Northing (Y)						System Datum				Groundwater not observed at time of exploration				
Notes:														


Elevation (feet)	FIELD DATA					Graphic Log	Group Classification	MATERIAL DESCRIPTION	Sheen	Headspace Vapor (ppm)	REMARKS
	Depth (feet)	Interval Recovered (in)	Blows/foot	Collected Sample	Sample Name Testing						
0	36					DUF	Forest duff, sandy silt with organic matter and wood debris	NS	<1		
						SM	Brown silty fine to medium sand (loose, moist)				
						ML	Brown sandy silt (loose, moist)				
5	60					SM	Brown silty fine to medium sand with occasional gravel				
10	60										
						ML	Gray sandy silt (loose, moist)				
15											

GEI-3 (9-10)

GEI-3 (14-15)

Note: See Figure A-1 for explanation of symbols.  
Coordinates Data Source: Horizontal approximated based on . Vertical approximated based on .

Date: 11/18/21 Path: P:\04180393\GINT\018039300.GPJ DBLibrary\Library\GEOENGINEERS\_DF\_STD\_US\_JUNE\_2017.GLB\GEB\_ENVIRONMENTAL\_STANDARD\_NO\_GW

<b>Log of Boring GEI-3</b>	
	Project: A-1 Towing Project Location: SeaTac, Washington Project Number: 0180-393-00
Figure A-4 Sheet 1 of 1	

Start Drilled	10/25/2021	End	10/25/2021	Total Depth (ft)	15	Logged By	DC	Checked By	CJW	Driller	Holocene Drilling, Inc.	Drilling Method	Direct-push
Surface Elevation (ft) Vertical Datum	Undetermined					Hammer Data	Geoprobe 7822T						
Easting (X) Northing (Y)						System Datum	Groundwater not observed at time of exploration						
Notes:													

Elevation (feet)	FIELD DATA					Graphic Log	Group Classification	MATERIAL DESCRIPTION	Sheen	Headspace Vapor (ppm)	REMARKS
	Depth (feet)	Interval Recovered (in)	Blows/foot	Collected Sample	Sample Name Testing						
0	48					SP-SM	Brown silty fine to coarse sand with occasional gravel (loose, moist)	NS	<1		
						ML	Brown sandy silt				
5	60					SM	Brown silty fine to medium sand with occasional gravel				
10	60										
15						ML	Gray sandy silt (loose, moist)				

Note: See Figure A-1 for explanation of symbols.  
Coordinates Data Source: Horizontal approximated based on . Vertical approximated based on .

### Log of Boring GEI-4



Project: A-1 Towing  
Project Location: SeaTac, Washington  
Project Number: 0180-393-00

Figure A-5  
Sheet 1 of 1

Date: 11/18/21 Path: P:\0180393\GINT\018039300.GPJ DBLibrary\Library\GEOENGINEERS\_DF\_STD\_US\_JUNE\_2017.GLB\GEB\_ENVIRONMENTAL\_STANDARD\_NO\_GW

Start Drilled	10/26/2021	End	10/26/2021	Total Depth (ft)	16	Logged By	DC	Checked By	CJW	Driller	Holocene Drilling, Inc.	Drilling Method	Direct-push
Surface Elevation (ft) Vertical Datum	Undetermined					Hammer Data	Geoprobe 7822T						
Easting (X) Northing (Y)						System Datum	Groundwater not observed at time of exploration						
Notes:													

Elevation (feet)	FIELD DATA					Graphic Log	Group Classification	MATERIAL DESCRIPTION	Sheen	Headspace Vapor (ppm)	REMARKS
	Depth (feet)	Interval Recovered (in)	Blows/foot	Collected Sample	Sample Name Testing						
0								Vac truck used for 5 feet			
5	60					SP-SM	Brown silty fine to coarse sand with occasional gravel (loose, moist)	NS	<1		
				GEI-5 (7-8)							
10	60					ML	Gray sandy silt (loose, moist)				
				GEI-5 (14-15)							
15											

Note: See Figure A-1 for explanation of symbols.  
Coordinates Data Source: Horizontal approximated based on . Vertical approximated based on .

### Log of Boring GEI-5



Project: A-1 Towing  
Project Location: SeaTac, Washington  
Project Number: 0180-393-00

Figure A-6  
Sheet 1 of 1

Date: 11/18/21 Path: P:\0180393\GINT\018039300.GPJ DBLibrary\Library\GEOENGINEERS\_DF\_STD\_US\_JUNE\_2017.GLB\GEB\_ENVIRONMENTAL\_STANDARD\_NO\_GW

Start Drilled	10/26/2021	End	10/26/2021	Total Depth (ft)	15	Logged By	DC	Checked By	CJW	Driller	Holocene Drilling, Inc.	Drilling Method	Direct-push
Surface Elevation (ft) Vertical Datum	Undetermined					Hammer Data	Geoprobe 7822T						
Easting (X) Northing (Y)						System Datum	Groundwater not observed at time of exploration						
Notes:													

Elevation (feet)	FIELD DATA					Graphic Log	Group Classification	MATERIAL DESCRIPTION	Sheen	Headspace Vapor (ppm)	REMARKS
	Depth (feet)	Interval Recovered (in)	Blows/foot	Collected Sample	Sample Name Testing						
0								Vac truck used for 6 feet			
5	60								NS	<1	
					GEI-6 (7-8)		SP-SM	Brown silty fine to coarse sand (loose, moist)			
10	60										
					GEI-6 (14-15)		SM	Gray silty fine to medium sand (loose, moist)			
15											

Note: See Figure A-1 for explanation of symbols.  
Coordinates Data Source: Horizontal approximated based on . Vertical approximated based on .

### Log of Boring GEI-6

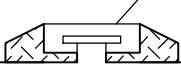


Project: A-1 Towing  
Project Location: SeaTac, Washington  
Project Number: 0180-393-00

Figure A-7  
Sheet 1 of 1

Date: 11/18/21 Path: P:\0180393\GINT\018039300.GPJ DBLlibrary\Library\GEOENGINEERS\_DF\_STD\_US\_JUNE\_2017.GLB\GEB\_ENVIRONMENTAL\_STANDARD\_NO\_GW

Start Drilled 10/25/2021	End 10/25/2021	Total Depth (ft)	15	Logged By Checked By	DC CJW	Driller Holocene Drilling, Inc.	Drilling Method	Direct-push
Hammer Data		Drilling Equipment			Geoprobe 7822T		A 2-in well was installed on 10/25/2021 to a depth of 15 ft.	
Surface Elevation (ft) Vertical Datum		Undetermined			Top of Casing Elevation (ft)		Groundwater Date Measured	
Easting (X) Northing (Y)		Horizontal Datum					Depth to Water (ft)	
							Elevation (ft)	
Notes:								


Elevation (feet)	FIELD DATA						MATERIAL DESCRIPTION	Sheen	Headspace Vapor (ppm)	WELL LOG
	Depth (feet)	Interval Recovered (in)	Blows/foot	Collected Sample	Sample Name Testing	Water Level				
0	60					SPSM	Brown silty fine to coarse sand (loose, moist)	NS	<1	 Steel surface monument
						ML	Brown sandy silt (loose, moist)			
5	60					SM	Brown silty fine to medium sand with occasional gravel (loose, moist)			
10	60									
15						SM	Gray silty fine to medium sand with occasional gravel (loose, moist)			

MW-1 (9-10)

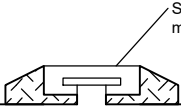
MW-1 (14-15)

Note: See Figure A-1 for explanation of symbols.  
Coordinates Data Source: Horizontal approximated based on . Vertical approximated based on .

Date: 11/18/21 Path: P:\01480393\GINT\018039300.GPJ DBLibrary\Library\GEOENGINEERS\_DF\_STD\_US\_JUNE\_2017.GLB\GEB\_ENVIRONMENTAL\_WELL

<b>Log of Boring MW-1</b>		
	Project: A-1 Towing	Figure A-8 Sheet 1 of 1
	Project Location: SeaTac, Washington	
	Project Number: 0180-393-00	

Start Drilled 10/25/2021	End 10/25/2021	Total Depth (ft)	15	Logged By Checked By	DC CJW	Driller Holocene Drilling, Inc.	Drilling Method	Direct-push
Hammer Data				Drilling Equipment Geoprobe 7822T			A 2-in well was installed on 10/26/2021 to a depth of 15 ft.	
Surface Elevation (ft) Vertical Datum				Undetermined			Top of Casing Elevation (ft)	
Easting (X) Northing (Y)				Horizontal Datum			Groundwater Date Measured	Depth to Water (ft) Elevation (ft)
Notes:								

Elevation (feet)	FIELD DATA						Group Classification	MATERIAL DESCRIPTION	Sheen	Headspace Vapor (ppm)	WELL LOG 
	Depth (feet)	Interval Recovered (in)	Blows/ foot	Collected Sample	Sample Name Testing	Water Level					
0	36					DUF	Forest duff	NS	<1		
						ML	Brown sandy silt (loose, moist)				
5	60					SM	Brown silty fine to medium sand with occasional gravel (loose, moist)				
10	60										
15						ML	Gray sandy silt (loose, moist)				

Note: See Figure A-1 for explanation of symbols.  
Coordinates Data Source: Horizontal approximated based on . Vertical approximated based on .

### Log of Boring MW-2

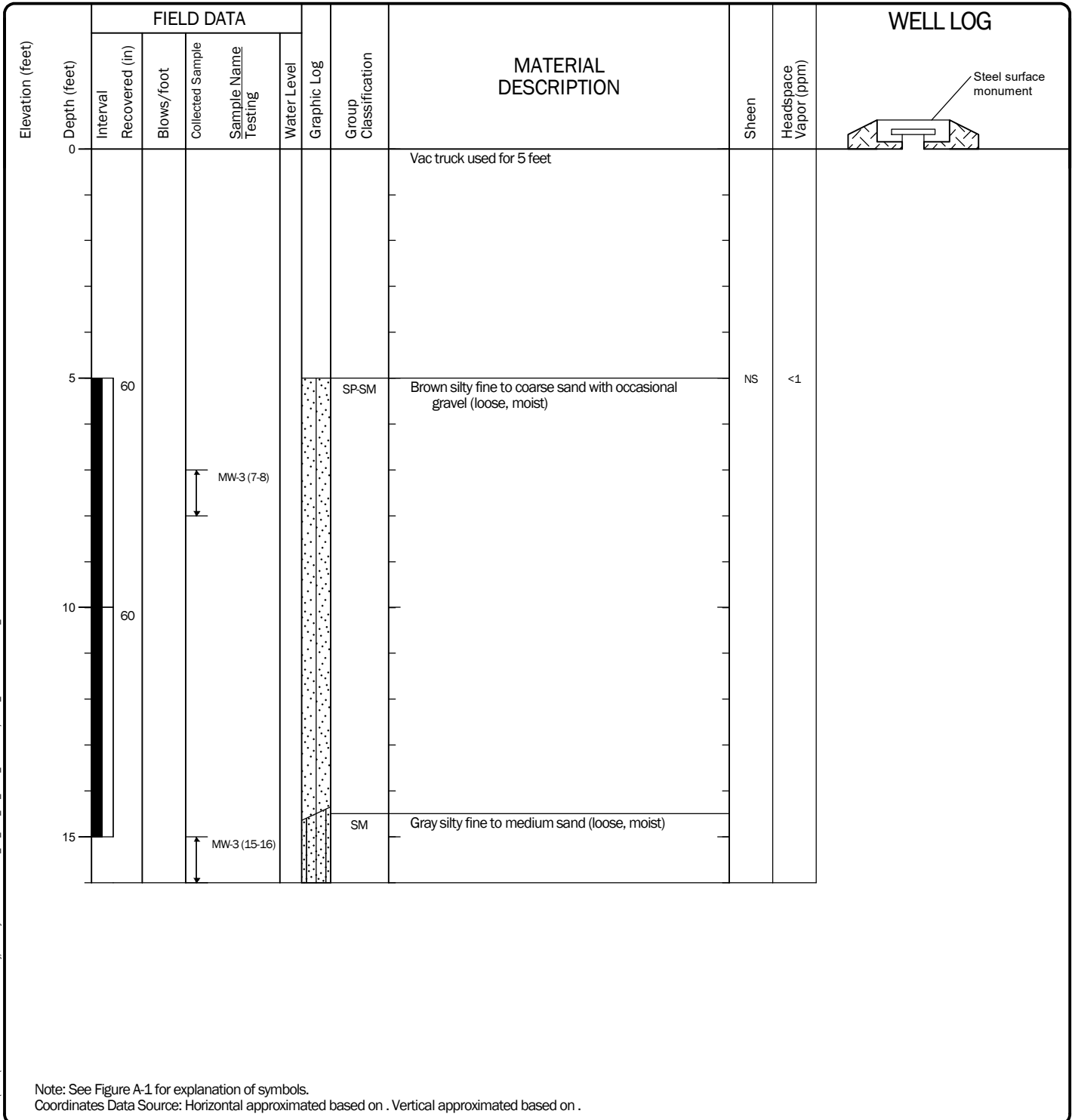


Project: A-1 Towing  
Project Location: SeaTac, Washington  
Project Number: 0180-393-00


Figure A-9  
Sheet 1 of 1

Date: 11/18/21 Path: P:\01480393\GINT\018039300.GPJ DBLibrary\Library\GEOENGINEERS\_DF\_STD\_US\_JUNE\_2017.GLB\GEB\_ENVIRONMENTAL\_WELL

Start Drilled 10/26/2021	End 10/26/2021	Total Depth (ft)	16	Logged By Checked By	DC CJW	Driller	Holocene Drilling, Inc.	Drilling Method	Direct-push
Hammer Data		Drilling Equipment			Geoprobe 7822T		A 2-in well was installed on 10/26/2021 to a depth of .16 ft.		
Surface Elevation (ft) Vertical Datum		Undetermined			Top of Casing Elevation (ft)		Groundwater Date Measured		
Easting (X) Northing (Y)		Horizontal Datum					Depth to Water (ft)		Elevation (ft)
Notes:									



Date: 11/18/21 Path: P:\01480393\GINT\018039300.GPJ DBLibrary\Library\GEOENGINEERS\_DF\_STD\_US\_JUNE\_2017.GLB\GEB\_ENVIRONMENTAL\_WELL

<b>Log of Boring MW-3</b>	
	Project: A-1 Towing Project Location: SeaTac, Washington Project Number: 0180-393-00
	Figure A-10 Sheet 1 of 1

**APPENDIX B**  
**Chemical Analytical Program**

## **APPENDIX B CHEMICAL ANALYTICAL PROGRAM**

### **Analytical Methods**

Chain-of-custody procedures were followed during the transport of the field samples to the analytical laboratory. The samples were held in cold storage pending extraction and/or analysis. The analytical results, analytical methods reference and laboratory quality control (QC) records are included in this appendix. The analytical results are also summarized in the text and tables of this report.

### **Analytical Data Review**

The laboratory maintains an internal quality assurance program as documented in its laboratory quality assurance manual. The laboratory uses a combination of blanks, surrogate recoveries, duplicates, matrix spike recoveries, matrix spike duplicate recoveries, blank spike recoveries and blank spike duplicate recoveries to evaluate the validity of the analytical results. The laboratory also uses data quality goals for individual chemicals or groups of chemicals based on the long-term performance of the test methods. The data quality goals were included in the laboratory reports. The laboratory compared each group of samples with the existing data quality goals and noted any exceptions in the laboratory report. Data quality exceptions documented by the accredited laboratory were reviewed by GeoEngineers and are addressed in the data quality exception section of this appendix.

### **Analytical Data Review Summary**

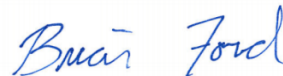
Quality controls were within required limits including duplicates. The data is of acceptable quality for their intended use.

**GeoEngineers- Portland, OR**

Sample Delivery Group: L1423266  
Samples Received: 10/27/2021  
Project Number: 0180-393-00  
Description: WSDOT-A-1 Towing

Report To: Cris Watkins  
4000 Kruse Way Place  
Bldg. 3, Suite 200  
Lake Oswego, OR 97035

Entire Report Reviewed By:



Brian Ford  
Project Manager

Results relate only to the items tested or calibrated and are reported as rounded values. This test report shall not be reproduced, except in full, without written approval of the laboratory. Where applicable, sampling conducted by Pace Analytical National is performed per guidance provided in laboratory standard operating procedures ENV-SOP-MTJL-0067 and ENV-SOP-MTJL-0068. Where sampling conducted by the customer, results relate to the accuracy of the information provided, and as the samples are received.

**Pace Analytical National**12065 Lebanon Rd Mount Juliet, TN 37122 615-758-5858 800-767-5859 [www.pacenational.com](http://www.pacenational.com)

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<b>Metals (ICPMS) by Method 6020B</b>	<b>35</b>
<b>Volatile Organic Compounds (GC) by Method NWTPHGX</b>	<b>36</b>
<b>Volatile Organic Compounds (GC/MS) by Method 8260D</b>	<b>38</b>
<b>Semi-Volatile Organic Compounds (GC) by Method NWTPHDX-NO SGT</b>	<b>42</b>
<b>Semi Volatile Organic Compounds (GC/MS) by Method 8270E-SIM</b>	<b>43</b>
<b>Gl: Glossary of Terms</b>	<b>45</b>
<b>Al: Accreditations &amp; Locations</b>	<b>46</b>
<b>Sc: Sample Chain of Custody</b>	<b>47</b>

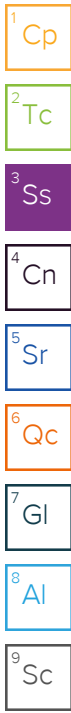
<sup>1</sup> Cp
<sup>2</sup> Tc
<sup>3</sup> Ss
<sup>4</sup> Cn
<sup>5</sup> Sr
<sup>6</sup> Qc
<sup>7</sup> Gl
<sup>8</sup> Al
<sup>9</sup> Sc

# SAMPLE SUMMARY

## GEI-1-9-10 L1423266-01 Solid

Collected by Dexter C.      Collected date/time 10/25/21 11:20      Received date/time 10/27/21 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Total Solids by Method 2540 G-2011	WG1765837	1	10/30/21 16:24	10/30/21 16:39	KDW	Mt. Juliet, TN
Mercury by Method 7471B	WG1766498	1	11/02/21 08:07	11/02/21 15:04	BMF	Mt. Juliet, TN
Metals (ICPMS) by Method 6020B	WG1765966	5	11/02/21 12:22	11/03/21 18:25	LD	Mt. Juliet, TN
Volatile Organic Compounds (GC) by Method NWTPHGX	WG1765532	25	10/25/21 11:20	10/29/21 11:17	JAH	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260D	WG1766323	1	10/25/21 11:20	10/31/21 03:56	JHH	Mt. Juliet, TN
Semi-Volatile Organic Compounds (GC) by Method NWTPHDX-NO SGT	WG1766948	1	11/02/21 03:07	11/03/21 07:15	JAS	Mt. Juliet, TN
Semi Volatile Organic Compounds (GC/MS) by Method 8270E-SIM	WG1765811	1	10/30/21 01:41	10/30/21 13:00	AAT	Mt. Juliet, TN



## GEI-2-9-10 L1423266-03 Solid

Collected by Dexter C.      Collected date/time 10/25/21 11:40      Received date/time 10/27/21 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Total Solids by Method 2540 G-2011	WG1765837	1	10/30/21 16:24	10/30/21 16:39	KDW	Mt. Juliet, TN
Mercury by Method 7471B	WG1766498	1	11/02/21 08:07	11/02/21 15:07	BMF	Mt. Juliet, TN
Metals (ICPMS) by Method 6020B	WG1765966	5	11/02/21 12:22	11/03/21 18:28	LD	Mt. Juliet, TN
Volatile Organic Compounds (GC) by Method NWTPHGX	WG1765532	25	10/25/21 11:40	10/29/21 11:38	JAH	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260D	WG1766323	1	10/25/21 11:40	10/31/21 04:15	JHH	Mt. Juliet, TN
Semi-Volatile Organic Compounds (GC) by Method NWTPHDX-NO SGT	WG1766948	1	11/02/21 03:07	11/03/21 06:11	JAS	Mt. Juliet, TN
Semi Volatile Organic Compounds (GC/MS) by Method 8270E-SIM	WG1765811	1	10/30/21 01:41	10/30/21 13:18	AAT	Mt. Juliet, TN

## GEI-3-9-10 L1423266-05 Solid

Collected by Dexter C.      Collected date/time 10/25/21 13:45      Received date/time 10/27/21 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Total Solids by Method 2540 G-2011	WG1765837	1	10/30/21 16:24	10/30/21 16:39	KDW	Mt. Juliet, TN
Mercury by Method 7471B	WG1766498	1	11/02/21 08:07	11/02/21 15:09	BMF	Mt. Juliet, TN
Metals (ICPMS) by Method 6020B	WG1765966	5	11/02/21 12:22	11/03/21 18:32	LD	Mt. Juliet, TN
Volatile Organic Compounds (GC) by Method NWTPHGX	WG1765532	25	10/25/21 13:45	10/29/21 12:00	JAH	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260D	WG1766323	1	10/25/21 13:45	10/31/21 04:34	JHH	Mt. Juliet, TN
Semi-Volatile Organic Compounds (GC) by Method NWTPHDX-NO SGT	WG1766948	1	11/02/21 03:07	11/03/21 06:24	JAS	Mt. Juliet, TN
Semi Volatile Organic Compounds (GC/MS) by Method 8270E-SIM	WG1765811	1	10/30/21 01:41	10/30/21 13:35	AAT	Mt. Juliet, TN

## MW-2-9-10 L1423266-11 Solid

Collected by Dexter C.      Collected date/time 10/25/21 14:00      Received date/time 10/27/21 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Total Solids by Method 2540 G-2011	WG1765837	1	10/30/21 16:24	10/30/21 16:39	KDW	Mt. Juliet, TN
Mercury by Method 7471B	WG1766498	1	11/02/21 08:07	11/02/21 15:12	BMF	Mt. Juliet, TN
Metals (ICPMS) by Method 6020B	WG1765966	5	11/02/21 12:22	11/03/21 18:35	LD	Mt. Juliet, TN
Volatile Organic Compounds (GC) by Method NWTPHGX	WG1765548	25	10/25/21 14:00	10/29/21 04:21	MGF	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260D	WG1766323	1	10/25/21 14:00	10/31/21 04:53	JHH	Mt. Juliet, TN
Semi-Volatile Organic Compounds (GC) by Method NWTPHDX-NO SGT	WG1766948	1	11/02/21 03:07	11/03/21 06:36	JAS	Mt. Juliet, TN
Semi Volatile Organic Compounds (GC/MS) by Method 8270E-SIM	WG1765811	1	10/30/21 01:41	10/30/21 13:52	AAT	Mt. Juliet, TN

## GEI-4-2-3 L1423266-13 Solid

Collected by Dexter C.      Collected date/time 10/26/21 10:00      Received date/time 10/27/21 09:30

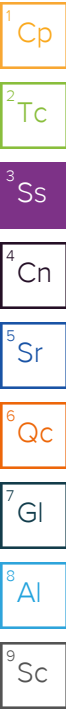
Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Total Solids by Method 2540 G-2011	WG1765837	1	10/30/21 16:24	10/30/21 16:39	KDW	Mt. Juliet, TN
Mercury by Method 7471B	WG1766498	1	11/02/21 08:07	11/02/21 15:15	BMF	Mt. Juliet, TN
Metals (ICPMS) by Method 6020B	WG1765966	5	11/02/21 12:22	11/03/21 18:39	LD	Mt. Juliet, TN
Volatile Organic Compounds (GC) by Method NWTPHGX	WG1765548	25	10/26/21 10:00	10/29/21 04:43	MGF	Mt. Juliet, TN

# SAMPLE SUMMARY

## GEI-4-2-3 L1423266-13 Solid

Collected by Dexter C.      Collected date/time 10/26/21 10:00      Received date/time 10/27/21 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Volatile Organic Compounds (GC/MS) by Method 8260D	WG1766323	1	10/26/21 10:00	10/31/21 05:12	JHH	Mt. Juliet, TN
Semi-Volatile Organic Compounds (GC) by Method NWTPHDX-NO SGT	WG1766948	5	11/02/21 03:07	11/03/21 15:22	JAS	Mt. Juliet, TN
Semi Volatile Organic Compounds (GC/MS) by Method 8270E-SIM	WG1765811	1	10/30/21 01:41	10/30/21 16:11	AAT	Mt. Juliet, TN



## MW-1-2-3 L1423266-14 Solid

Collected by Dexter C.      Collected date/time 10/26/21 10:20      Received date/time 10/27/21 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Total Solids by Method 2540 G-2011	WG1765837	1	10/30/21 16:24	10/30/21 16:39	KDW	Mt. Juliet, TN
Mercury by Method 7471B	WG1766498	1	11/02/21 08:07	11/02/21 15:22	BMF	Mt. Juliet, TN
Metals (ICPMS) by Method 6020B	WG1765966	5	11/02/21 12:22	11/03/21 18:42	LD	Mt. Juliet, TN
Volatile Organic Compounds (GC) by Method NWTPHGX	WG1765548	25	10/26/21 10:20	10/29/21 05:05	MGF	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260D	WG1766323	1	10/26/21 10:20	10/31/21 05:31	JHH	Mt. Juliet, TN
Semi-Volatile Organic Compounds (GC) by Method NWTPHDX-NO SGT	WG1766948	1	11/02/21 03:07	11/03/21 06:49	JAS	Mt. Juliet, TN
Semi Volatile Organic Compounds (GC/MS) by Method 8270E-SIM	WG1765811	1	10/30/21 01:41	10/30/21 14:27	AAT	Mt. Juliet, TN

## GEI-5-7-8 L1423266-15 Solid

Collected by Dexter C.      Collected date/time 10/26/21 10:40      Received date/time 10/27/21 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Total Solids by Method 2540 G-2011	WG1765837	1	10/30/21 16:24	10/30/21 16:39	KDW	Mt. Juliet, TN
Mercury by Method 7471B	WG1766498	1	11/02/21 08:07	11/02/21 15:25	BMF	Mt. Juliet, TN
Metals (ICPMS) by Method 6020B	WG1765966	5	11/02/21 12:22	11/03/21 18:46	LD	Mt. Juliet, TN
Volatile Organic Compounds (GC) by Method NWTPHGX	WG1765548	25	10/26/21 10:40	10/29/21 05:27	MGF	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260D	WG1766323	1	10/26/21 10:40	10/31/21 05:49	JHH	Mt. Juliet, TN
Semi-Volatile Organic Compounds (GC) by Method NWTPHDX-NO SGT	WG1766948	1	11/02/21 03:07	11/03/21 07:02	JAS	Mt. Juliet, TN
Semi Volatile Organic Compounds (GC/MS) by Method 8270E-SIM	WG1765811	1	10/30/21 01:41	10/30/21 14:44	AAT	Mt. Juliet, TN

## MW-3-2-3 L1423266-17 Solid

Collected by Dexter C.      Collected date/time 10/26/21 11:00      Received date/time 10/27/21 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Total Solids by Method 2540 G-2011	WG1765837	1	10/30/21 16:24	10/30/21 16:39	KDW	Mt. Juliet, TN
Mercury by Method 7471B	WG1766498	1	11/02/21 08:07	11/02/21 15:27	BMF	Mt. Juliet, TN
Metals (ICPMS) by Method 6020B	WG1765966	5	11/02/21 12:22	11/03/21 18:49	LD	Mt. Juliet, TN
Volatile Organic Compounds (GC) by Method NWTPHGX	WG1765548	25	10/26/21 11:00	10/29/21 05:49	MGF	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260D	WG1766323	1	10/26/21 11:00	10/31/21 06:08	JHH	Mt. Juliet, TN
Semi-Volatile Organic Compounds (GC) by Method NWTPHDX-NO SGT	WG1766948	1	11/02/21 03:07	11/03/21 07:54	JAS	Mt. Juliet, TN
Semi Volatile Organic Compounds (GC/MS) by Method 8270E-SIM	WG1765811	1	10/30/21 01:41	10/30/21 15:02	AAT	Mt. Juliet, TN

## GEI-6-7-8 L1423266-20 Solid

Collected by Dexter C.      Collected date/time 10/26/21 12:05      Received date/time 10/27/21 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Total Solids by Method 2540 G-2011	WG1765837	1	10/30/21 16:24	10/30/21 16:39	KDW	Mt. Juliet, TN
Mercury by Method 7471B	WG1766498	1	11/02/21 08:07	11/02/21 15:30	BMF	Mt. Juliet, TN
Metals (ICPMS) by Method 6020B	WG1765966	5	11/02/21 12:22	11/03/21 18:52	LD	Mt. Juliet, TN
Volatile Organic Compounds (GC) by Method NWTPHGX	WG1765548	25	10/26/21 12:05	10/29/21 06:11	MGF	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260D	WG1766323	1	10/26/21 12:05	10/31/21 06:27	JHH	Mt. Juliet, TN
Semi-Volatile Organic Compounds (GC) by Method NWTPHDX-NO SGT	WG1766948	1	11/02/21 03:07	11/03/21 07:41	JAS	Mt. Juliet, TN
Semi Volatile Organic Compounds (GC/MS) by Method 8270E-SIM	WG1765811	1	10/30/21 01:41	10/30/21 15:19	AAT	Mt. Juliet, TN

# CASE NARRATIVE

All sample aliquots were received at the correct temperature, in the proper containers, with the appropriate preservatives, and within method specified holding times, unless qualified or notated within the report. Where applicable, all MDL (LOD) and RDL (LOQ) values reported for environmental samples have been corrected for the dilution factor used in the analysis. All Method and Batch Quality Control are within established criteria except where addressed in this case narrative, a non-conformance form or properly qualified within the sample results. By my digital signature below, I affirm to the best of my knowledge, all problems/anomalies observed by the laboratory as having the potential to affect the quality of the data have been identified by the laboratory, and no information or data have been knowingly withheld that would affect the quality of the data.



Brian Ford  
Project Manager

- <sup>1</sup> Cp
- <sup>2</sup> Tc
- <sup>3</sup> Ss
- <sup>4</sup> Cn
- <sup>5</sup> Sr
- <sup>6</sup> Qc
- <sup>7</sup> Gl
- <sup>8</sup> Al
- <sup>9</sup> Sc

Total Solids by Method 2540 G-2011

Analyte	Result	Qualifier	Dilution	Analysis	Batch
Total Solids	91.8		1	10/30/2021 16:39	<a href="#">WG1765837</a>

Mercury by Method 7471B

Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
Mercury	ND		0.0436	1	11/02/2021 15:04	<a href="#">WG1766498</a>

Metals (ICPMS) by Method 6020B

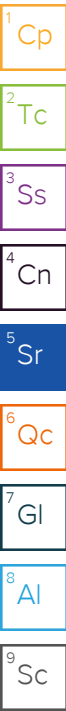
Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
Arsenic	2.26		1.09	5	11/03/2021 18:25	<a href="#">WG1765966</a>
Barium	41.9		2.72	5	11/03/2021 18:25	<a href="#">WG1765966</a>
Cadmium	ND		1.09	5	11/03/2021 18:25	<a href="#">WG1765966</a>
Chromium	23.4		5.45	5	11/03/2021 18:25	<a href="#">WG1765966</a>
Lead	ND		2.18	5	11/03/2021 18:25	<a href="#">WG1765966</a>
Selenium	ND		2.72	5	11/03/2021 18:25	<a href="#">WG1765966</a>
Silver	ND		0.545	5	11/03/2021 18:25	<a href="#">WG1765966</a>

Volatile Organic Compounds (GC) by Method NWTPHGX

Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
Gasoline Range Organics-NWTPH	ND		3.01	25	10/29/2021 11:17	<a href="#">WG1765532</a>
(S) a,a,a-Trifluorotoluene(FID)	119		77.0-120		10/29/2021 11:17	<a href="#">WG1765532</a>

Volatile Organic Compounds (GC/MS) by Method 8260D

Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
Acetone	ND		0.0607	1	10/31/2021 03:56	<a href="#">WG1766323</a>
Acrylonitrile	ND	<a href="#">C3</a>	0.0152	1	10/31/2021 03:56	<a href="#">WG1766323</a>
Benzene	ND		0.00121	1	10/31/2021 03:56	<a href="#">WG1766323</a>
Bromobenzene	ND		0.0152	1	10/31/2021 03:56	<a href="#">WG1766323</a>
Bromodichloromethane	ND		0.00303	1	10/31/2021 03:56	<a href="#">WG1766323</a>
Bromoform	ND		0.0303	1	10/31/2021 03:56	<a href="#">WG1766323</a>
Bromomethane	ND		0.0152	1	10/31/2021 03:56	<a href="#">WG1766323</a>
n-Butylbenzene	ND		0.0152	1	10/31/2021 03:56	<a href="#">WG1766323</a>
sec-Butylbenzene	ND		0.0152	1	10/31/2021 03:56	<a href="#">WG1766323</a>
tert-Butylbenzene	ND		0.00607	1	10/31/2021 03:56	<a href="#">WG1766323</a>
Carbon tetrachloride	ND		0.00607	1	10/31/2021 03:56	<a href="#">WG1766323</a>
Chlorobenzene	ND		0.00303	1	10/31/2021 03:56	<a href="#">WG1766323</a>
Chlorodibromomethane	ND		0.00303	1	10/31/2021 03:56	<a href="#">WG1766323</a>
Chloroethane	ND		0.00607	1	10/31/2021 03:56	<a href="#">WG1766323</a>
Chloroform	ND		0.00303	1	10/31/2021 03:56	<a href="#">WG1766323</a>
Chloromethane	ND		0.0152	1	10/31/2021 03:56	<a href="#">WG1766323</a>
2-Chlorotoluene	ND		0.00303	1	10/31/2021 03:56	<a href="#">WG1766323</a>
4-Chlorotoluene	ND		0.00607	1	10/31/2021 03:56	<a href="#">WG1766323</a>
1,2-Dibromo-3-Chloropropane	ND		0.0303	1	10/31/2021 03:56	<a href="#">WG1766323</a>
1,2-Dibromoethane	ND		0.00303	1	10/31/2021 03:56	<a href="#">WG1766323</a>
Dibromomethane	ND		0.00607	1	10/31/2021 03:56	<a href="#">WG1766323</a>
1,2-Dichlorobenzene	ND		0.00607	1	10/31/2021 03:56	<a href="#">WG1766323</a>
1,3-Dichlorobenzene	ND		0.00607	1	10/31/2021 03:56	<a href="#">WG1766323</a>
1,4-Dichlorobenzene	ND		0.00607	1	10/31/2021 03:56	<a href="#">WG1766323</a>
Dichlorodifluoromethane	ND		0.00303	1	10/31/2021 03:56	<a href="#">WG1766323</a>
1,1-Dichloroethane	ND		0.00303	1	10/31/2021 03:56	<a href="#">WG1766323</a>
1,2-Dichloroethane	ND		0.00303	1	10/31/2021 03:56	<a href="#">WG1766323</a>



Volatile Organic Compounds (GC/MS) by Method 8260D

Analyte	Result (dry) mg/kg	Qualifier	RDL (dry) mg/kg	Dilution	Analysis date / time	Batch
1,1-Dichloroethene	ND		0.00303	1	10/31/2021 03:56	WG1766323
cis-1,2-Dichloroethene	ND		0.00303	1	10/31/2021 03:56	WG1766323
trans-1,2-Dichloroethene	ND		0.00607	1	10/31/2021 03:56	WG1766323
1,2-Dichloropropane	ND		0.00607	1	10/31/2021 03:56	WG1766323
1,1-Dichloropropene	ND		0.00303	1	10/31/2021 03:56	WG1766323
1,3-Dichloropropane	ND		0.00607	1	10/31/2021 03:56	WG1766323
cis-1,3-Dichloropropene	ND		0.00303	1	10/31/2021 03:56	WG1766323
trans-1,3-Dichloropropene	ND		0.00607	1	10/31/2021 03:56	WG1766323
2,2-Dichloropropane	ND		0.00303	1	10/31/2021 03:56	WG1766323
Di-isopropyl ether	ND		0.00121	1	10/31/2021 03:56	WG1766323
Ethylbenzene	ND		0.00303	1	10/31/2021 03:56	WG1766323
Hexachloro-1,3-butadiene	ND		0.0303	1	10/31/2021 03:56	WG1766323
Isopropylbenzene	ND		0.00303	1	10/31/2021 03:56	WG1766323
p-Isopropyltoluene	ND		0.00607	1	10/31/2021 03:56	WG1766323
2-Butanone (MEK)	ND		0.121	1	10/31/2021 03:56	WG1766323
Methylene Chloride	ND		0.0303	1	10/31/2021 03:56	WG1766323
4-Methyl-2-pentanone (MIBK)	ND		0.0303	1	10/31/2021 03:56	WG1766323
Methyl tert-butyl ether	ND		0.00121	1	10/31/2021 03:56	WG1766323
Naphthalene	ND		0.0152	1	10/31/2021 03:56	WG1766323
n-Propylbenzene	ND		0.00607	1	10/31/2021 03:56	WG1766323
Styrene	ND		0.0152	1	10/31/2021 03:56	WG1766323
1,1,1,2-Tetrachloroethane	ND		0.00303	1	10/31/2021 03:56	WG1766323
1,1,2,2-Tetrachloroethane	ND		0.00303	1	10/31/2021 03:56	WG1766323
1,1,2-Trichlorotrifluoroethane	ND		0.00303	1	10/31/2021 03:56	WG1766323
Tetrachloroethene	ND		0.00303	1	10/31/2021 03:56	WG1766323
Toluene	ND		0.00607	1	10/31/2021 03:56	WG1766323
1,2,3-Trichlorobenzene	ND	C4	0.0152	1	10/31/2021 03:56	WG1766323
1,2,4-Trichlorobenzene	ND		0.0152	1	10/31/2021 03:56	WG1766323
1,1,1-Trichloroethane	ND		0.00303	1	10/31/2021 03:56	WG1766323
1,1,2-Trichloroethane	ND		0.00303	1	10/31/2021 03:56	WG1766323
Trichloroethene	ND		0.00121	1	10/31/2021 03:56	WG1766323
Trichlorofluoromethane	ND		0.00303	1	10/31/2021 03:56	WG1766323
1,2,3-Trichloropropane	ND		0.0152	1	10/31/2021 03:56	WG1766323
1,2,4-Trimethylbenzene	ND		0.00607	1	10/31/2021 03:56	WG1766323
1,2,3-Trimethylbenzene	ND		0.00607	1	10/31/2021 03:56	WG1766323
1,3,5-Trimethylbenzene	ND		0.00607	1	10/31/2021 03:56	WG1766323
Vinyl chloride	ND		0.00303	1	10/31/2021 03:56	WG1766323
Xylenes, Total	ND		0.00789	1	10/31/2021 03:56	WG1766323
(S) Toluene-d8	104		75.0-131		10/31/2021 03:56	WG1766323
(S) 4-Bromofluorobenzene	99.6		67.0-138		10/31/2021 03:56	WG1766323
(S) 1,2-Dichloroethane-d4	103		70.0-130		10/31/2021 03:56	WG1766323

- 1 Cp
- 2 Tc
- 3 Ss
- 4 Cn
- 5 Sr
- 6 Qc
- 7 Gl
- 8 Al
- 9 Sc

Semi-Volatile Organic Compounds (GC) by Method NWTPHDX-NO SGT

Analyte	Result (dry) mg/kg	Qualifier	RDL (dry) mg/kg	Dilution	Analysis date / time	Batch
Diesel Range Organics (DRO)	ND		4.36	1	11/03/2021 07:15	WG1766948
Residual Range Organics (RRO)	ND		10.9	1	11/03/2021 07:15	WG1766948
(S) o-Terphenyl	53.8		18.0-148		11/03/2021 07:15	WG1766948

Semi Volatile Organic Compounds (GC/MS) by Method 8270E-SIM

Analyte	Result (dry) mg/kg	Qualifier	RDL (dry) mg/kg	Dilution	Analysis date / time	Batch
Anthracene	ND		0.00654	1	10/30/2021 13:00	<a href="#">WG1765811</a>
Acenaphthene	ND		0.00654	1	10/30/2021 13:00	<a href="#">WG1765811</a>
Acenaphthylene	ND		0.00654	1	10/30/2021 13:00	<a href="#">WG1765811</a>
Benzo(a)anthracene	ND		0.00654	1	10/30/2021 13:00	<a href="#">WG1765811</a>
Benzo(a)pyrene	ND		0.00654	1	10/30/2021 13:00	<a href="#">WG1765811</a>
Benzo(b)fluoranthene	ND		0.00654	1	10/30/2021 13:00	<a href="#">WG1765811</a>
Benzo(g,h,i)perylene	ND		0.00654	1	10/30/2021 13:00	<a href="#">WG1765811</a>
Benzo(k)fluoranthene	ND		0.00654	1	10/30/2021 13:00	<a href="#">WG1765811</a>
Chrysene	ND		0.00654	1	10/30/2021 13:00	<a href="#">WG1765811</a>
Dibenz(a,h)anthracene	ND		0.00654	1	10/30/2021 13:00	<a href="#">WG1765811</a>
Fluoranthene	ND		0.00654	1	10/30/2021 13:00	<a href="#">WG1765811</a>
Fluorene	ND		0.00654	1	10/30/2021 13:00	<a href="#">WG1765811</a>
Indeno(1,2,3-cd)pyrene	ND		0.00654	1	10/30/2021 13:00	<a href="#">WG1765811</a>
Naphthalene	ND		0.0218	1	10/30/2021 13:00	<a href="#">WG1765811</a>
Phenanthrene	ND		0.00654	1	10/30/2021 13:00	<a href="#">WG1765811</a>
Pyrene	ND		0.00654	1	10/30/2021 13:00	<a href="#">WG1765811</a>
1-Methylnaphthalene	ND		0.0218	1	10/30/2021 13:00	<a href="#">WG1765811</a>
2-Methylnaphthalene	ND		0.0218	1	10/30/2021 13:00	<a href="#">WG1765811</a>
2-Chloronaphthalene	ND		0.0218	1	10/30/2021 13:00	<a href="#">WG1765811</a>
(S) Nitrobenzene-d5	85.9		14.0-149		10/30/2021 13:00	<a href="#">WG1765811</a>
(S) 2-Fluorobiphenyl	70.0		34.0-125		10/30/2021 13:00	<a href="#">WG1765811</a>
(S) p-Terphenyl-d14	110		23.0-120		10/30/2021 13:00	<a href="#">WG1765811</a>

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc

Total Solids by Method 2540 G-2011

Analyte	Result	Qualifier	Dilution	Analysis	Batch
Total Solids	92.3		1	10/30/2021 16:39	<a href="#">WG1765837</a>

Mercury by Method 7471B

Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
Mercury	ND		0.0434	1	11/02/2021 15:07	<a href="#">WG1766498</a>

Metals (ICPMS) by Method 6020B

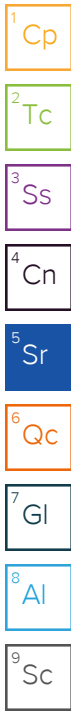
Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
Arsenic	2.36		1.08	5	11/03/2021 18:28	<a href="#">WG1765966</a>
Barium	45.2		2.71	5	11/03/2021 18:28	<a href="#">WG1765966</a>
Cadmium	ND		1.08	5	11/03/2021 18:28	<a href="#">WG1765966</a>
Chromium	21.5		5.42	5	11/03/2021 18:28	<a href="#">WG1765966</a>
Lead	ND		2.17	5	11/03/2021 18:28	<a href="#">WG1765966</a>
Selenium	ND		2.71	5	11/03/2021 18:28	<a href="#">WG1765966</a>
Silver	ND		0.542	5	11/03/2021 18:28	<a href="#">WG1765966</a>

Volatile Organic Compounds (GC) by Method NWTPHGX

Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
Gasoline Range Organics-NWTPH	ND		2.99	25	10/29/2021 11:38	<a href="#">WG1765532</a>
(S) a,a,a-Trifluorotoluene(FID)	119		77.0-120		10/29/2021 11:38	<a href="#">WG1765532</a>

Volatile Organic Compounds (GC/MS) by Method 8260D

Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
Acetone	ND		0.0594	1	10/31/2021 04:15	<a href="#">WG1766323</a>
Acrylonitrile	ND	<a href="#">C3</a>	0.0148	1	10/31/2021 04:15	<a href="#">WG1766323</a>
Benzene	ND		0.00119	1	10/31/2021 04:15	<a href="#">WG1766323</a>
Bromobenzene	ND		0.0148	1	10/31/2021 04:15	<a href="#">WG1766323</a>
Bromodichloromethane	ND		0.00297	1	10/31/2021 04:15	<a href="#">WG1766323</a>
Bromoform	ND		0.0297	1	10/31/2021 04:15	<a href="#">WG1766323</a>
Bromomethane	ND		0.0148	1	10/31/2021 04:15	<a href="#">WG1766323</a>
n-Butylbenzene	ND		0.0148	1	10/31/2021 04:15	<a href="#">WG1766323</a>
sec-Butylbenzene	ND		0.0148	1	10/31/2021 04:15	<a href="#">WG1766323</a>
tert-Butylbenzene	ND		0.00594	1	10/31/2021 04:15	<a href="#">WG1766323</a>
Carbon tetrachloride	ND		0.00594	1	10/31/2021 04:15	<a href="#">WG1766323</a>
Chlorobenzene	ND		0.00297	1	10/31/2021 04:15	<a href="#">WG1766323</a>
Chlorodibromomethane	ND		0.00297	1	10/31/2021 04:15	<a href="#">WG1766323</a>
Chloroethane	ND		0.00594	1	10/31/2021 04:15	<a href="#">WG1766323</a>
Chloroform	ND		0.00297	1	10/31/2021 04:15	<a href="#">WG1766323</a>
Chloromethane	ND		0.0148	1	10/31/2021 04:15	<a href="#">WG1766323</a>
2-Chlorotoluene	ND		0.00297	1	10/31/2021 04:15	<a href="#">WG1766323</a>
4-Chlorotoluene	ND		0.00594	1	10/31/2021 04:15	<a href="#">WG1766323</a>
1,2-Dibromo-3-Chloropropane	ND		0.0297	1	10/31/2021 04:15	<a href="#">WG1766323</a>
1,2-Dibromoethane	ND		0.00297	1	10/31/2021 04:15	<a href="#">WG1766323</a>
Dibromomethane	ND		0.00594	1	10/31/2021 04:15	<a href="#">WG1766323</a>
1,2-Dichlorobenzene	ND		0.00594	1	10/31/2021 04:15	<a href="#">WG1766323</a>
1,3-Dichlorobenzene	ND		0.00594	1	10/31/2021 04:15	<a href="#">WG1766323</a>
1,4-Dichlorobenzene	ND		0.00594	1	10/31/2021 04:15	<a href="#">WG1766323</a>
Dichlorodifluoromethane	ND		0.00297	1	10/31/2021 04:15	<a href="#">WG1766323</a>
1,1-Dichloroethane	ND		0.00297	1	10/31/2021 04:15	<a href="#">WG1766323</a>
1,2-Dichloroethane	ND		0.00297	1	10/31/2021 04:15	<a href="#">WG1766323</a>



Volatile Organic Compounds (GC/MS) by Method 8260D

Analyte	Result (dry) mg/kg	Qualifier	RDL (dry) mg/kg	Dilution	Analysis date / time	Batch
1,1-Dichloroethene	ND		0.00297	1	10/31/2021 04:15	WG1766323
cis-1,2-Dichloroethene	ND		0.00297	1	10/31/2021 04:15	WG1766323
trans-1,2-Dichloroethene	ND		0.00594	1	10/31/2021 04:15	WG1766323
1,2-Dichloropropane	ND		0.00594	1	10/31/2021 04:15	WG1766323
1,1-Dichloropropene	ND		0.00297	1	10/31/2021 04:15	WG1766323
1,3-Dichloropropane	ND		0.00594	1	10/31/2021 04:15	WG1766323
cis-1,3-Dichloropropene	ND		0.00297	1	10/31/2021 04:15	WG1766323
trans-1,3-Dichloropropene	ND		0.00594	1	10/31/2021 04:15	WG1766323
2,2-Dichloropropane	ND		0.00297	1	10/31/2021 04:15	WG1766323
Di-isopropyl ether	ND		0.00119	1	10/31/2021 04:15	WG1766323
Ethylbenzene	ND		0.00297	1	10/31/2021 04:15	WG1766323
Hexachloro-1,3-butadiene	ND		0.0297	1	10/31/2021 04:15	WG1766323
Isopropylbenzene	ND		0.00297	1	10/31/2021 04:15	WG1766323
p-Isopropyltoluene	ND		0.00594	1	10/31/2021 04:15	WG1766323
2-Butanone (MEK)	ND		0.119	1	10/31/2021 04:15	WG1766323
Methylene Chloride	ND		0.0297	1	10/31/2021 04:15	WG1766323
4-Methyl-2-pentanone (MIBK)	ND		0.0297	1	10/31/2021 04:15	WG1766323
Methyl tert-butyl ether	ND		0.00119	1	10/31/2021 04:15	WG1766323
Naphthalene	ND		0.0148	1	10/31/2021 04:15	WG1766323
n-Propylbenzene	ND		0.00594	1	10/31/2021 04:15	WG1766323
Styrene	ND		0.0148	1	10/31/2021 04:15	WG1766323
1,1,1,2-Tetrachloroethane	ND		0.00297	1	10/31/2021 04:15	WG1766323
1,1,2,2-Tetrachloroethane	ND		0.00297	1	10/31/2021 04:15	WG1766323
1,1,2-Trichlorotrifluoroethane	ND		0.00297	1	10/31/2021 04:15	WG1766323
Tetrachloroethene	ND		0.00297	1	10/31/2021 04:15	WG1766323
Toluene	ND		0.00594	1	10/31/2021 04:15	WG1766323
1,2,3-Trichlorobenzene	ND	C4	0.0148	1	10/31/2021 04:15	WG1766323
1,2,4-Trichlorobenzene	ND		0.0148	1	10/31/2021 04:15	WG1766323
1,1,1-Trichloroethane	ND		0.00297	1	10/31/2021 04:15	WG1766323
1,1,2-Trichloroethane	ND		0.00297	1	10/31/2021 04:15	WG1766323
Trichloroethene	ND		0.00119	1	10/31/2021 04:15	WG1766323
Trichlorofluoromethane	ND		0.00297	1	10/31/2021 04:15	WG1766323
1,2,3-Trichloropropane	ND		0.0148	1	10/31/2021 04:15	WG1766323
1,2,4-Trimethylbenzene	ND		0.00594	1	10/31/2021 04:15	WG1766323
1,2,3-Trimethylbenzene	ND		0.00594	1	10/31/2021 04:15	WG1766323
1,3,5-Trimethylbenzene	ND		0.00594	1	10/31/2021 04:15	WG1766323
Vinyl chloride	ND		0.00297	1	10/31/2021 04:15	WG1766323
Xylenes, Total	ND		0.00772	1	10/31/2021 04:15	WG1766323
(S) Toluene-d8	104		75.0-131		10/31/2021 04:15	WG1766323
(S) 4-Bromofluorobenzene	96.9		67.0-138		10/31/2021 04:15	WG1766323
(S) 1,2-Dichloroethane-d4	100		70.0-130		10/31/2021 04:15	WG1766323

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc

Semi-Volatile Organic Compounds (GC) by Method NWTPHDX-NO SGT

Analyte	Result (dry) mg/kg	Qualifier	RDL (dry) mg/kg	Dilution	Analysis date / time	Batch
Diesel Range Organics (DRO)	ND		4.34	1	11/03/2021 06:11	WG1766948
Residual Range Organics (RRO)	ND		10.8	1	11/03/2021 06:11	WG1766948
(S) o-Terphenyl	37.8		18.0-148		11/03/2021 06:11	WG1766948

Semi Volatile Organic Compounds (GC/MS) by Method 8270E-SIM

Analyte	Result (dry) mg/kg	Qualifier	RDL (dry) mg/kg	Dilution	Analysis date / time	Batch
Anthracene	ND		0.00650	1	10/30/2021 13:18	<a href="#">WG1765811</a>
Acenaphthene	ND		0.00650	1	10/30/2021 13:18	<a href="#">WG1765811</a>
Acenaphthylene	ND		0.00650	1	10/30/2021 13:18	<a href="#">WG1765811</a>
Benzo(a)anthracene	ND		0.00650	1	10/30/2021 13:18	<a href="#">WG1765811</a>
Benzo(a)pyrene	ND		0.00650	1	10/30/2021 13:18	<a href="#">WG1765811</a>
Benzo(b)fluoranthene	ND		0.00650	1	10/30/2021 13:18	<a href="#">WG1765811</a>
Benzo(g,h,i)perylene	ND		0.00650	1	10/30/2021 13:18	<a href="#">WG1765811</a>
Benzo(k)fluoranthene	ND		0.00650	1	10/30/2021 13:18	<a href="#">WG1765811</a>
Chrysene	ND		0.00650	1	10/30/2021 13:18	<a href="#">WG1765811</a>
Dibenz(a,h)anthracene	ND		0.00650	1	10/30/2021 13:18	<a href="#">WG1765811</a>
Fluoranthene	ND		0.00650	1	10/30/2021 13:18	<a href="#">WG1765811</a>
Fluorene	ND		0.00650	1	10/30/2021 13:18	<a href="#">WG1765811</a>
Indeno(1,2,3-cd)pyrene	ND		0.00650	1	10/30/2021 13:18	<a href="#">WG1765811</a>
Naphthalene	ND		0.0217	1	10/30/2021 13:18	<a href="#">WG1765811</a>
Phenanthrene	ND		0.00650	1	10/30/2021 13:18	<a href="#">WG1765811</a>
Pyrene	ND		0.00650	1	10/30/2021 13:18	<a href="#">WG1765811</a>
1-Methylnaphthalene	ND		0.0217	1	10/30/2021 13:18	<a href="#">WG1765811</a>
2-Methylnaphthalene	ND		0.0217	1	10/30/2021 13:18	<a href="#">WG1765811</a>
2-Chloronaphthalene	ND		0.0217	1	10/30/2021 13:18	<a href="#">WG1765811</a>
<i>(S)</i> Nitrobenzene-d5	98.0		14.0-149		10/30/2021 13:18	<a href="#">WG1765811</a>
<i>(S)</i> 2-Fluorobiphenyl	77.9		34.0-125		10/30/2021 13:18	<a href="#">WG1765811</a>
<i>(S)</i> p-Terphenyl-d14	96.1		23.0-120		10/30/2021 13:18	<a href="#">WG1765811</a>

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc

Total Solids by Method 2540 G-2011

Analyte	Result	Qualifier	Dilution	Analysis	Batch
	%			date / time	
Total Solids	89.0		1	10/30/2021 16:39	<a href="#">WG1765837</a>

Mercury by Method 7471B

Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
	mg/kg		mg/kg		date / time	
Mercury	ND		0.0450	1	11/02/2021 15:09	<a href="#">WG1766498</a>

Metals (ICPMS) by Method 6020B

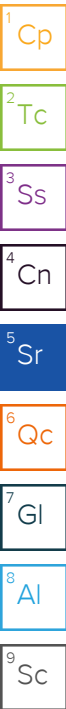
Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
	mg/kg		mg/kg		date / time	
Arsenic	2.93		1.12	5	11/03/2021 18:32	<a href="#">WG1765966</a>
Barium	25.3		2.81	5	11/03/2021 18:32	<a href="#">WG1765966</a>
Cadmium	ND		1.12	5	11/03/2021 18:32	<a href="#">WG1765966</a>
Chromium	11.8		5.62	5	11/03/2021 18:32	<a href="#">WG1765966</a>
Lead	ND		2.25	5	11/03/2021 18:32	<a href="#">WG1765966</a>
Selenium	ND		2.81	5	11/03/2021 18:32	<a href="#">WG1765966</a>
Silver	ND		0.562	5	11/03/2021 18:32	<a href="#">WG1765966</a>

Volatile Organic Compounds (GC) by Method NWTPHGX

Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
	mg/kg		mg/kg		date / time	
Gasoline Range Organics-NWTPH	ND		3.23	25	10/29/2021 12:00	<a href="#">WG1765532</a>
(S) a,a,a-Trifluorotoluene(FID)	119		77.0-120		10/29/2021 12:00	<a href="#">WG1765532</a>

Volatile Organic Compounds (GC/MS) by Method 8260D

Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
	mg/kg		mg/kg		date / time	
Acetone	ND		0.0646	1	10/31/2021 04:34	<a href="#">WG1766323</a>
Acrylonitrile	ND	<a href="#">C3</a>	0.0161	1	10/31/2021 04:34	<a href="#">WG1766323</a>
Benzene	ND		0.00129	1	10/31/2021 04:34	<a href="#">WG1766323</a>
Bromobenzene	ND		0.0161	1	10/31/2021 04:34	<a href="#">WG1766323</a>
Bromodichloromethane	ND		0.00323	1	10/31/2021 04:34	<a href="#">WG1766323</a>
Bromoform	ND		0.0323	1	10/31/2021 04:34	<a href="#">WG1766323</a>
Bromomethane	ND		0.0161	1	10/31/2021 04:34	<a href="#">WG1766323</a>
n-Butylbenzene	ND		0.0161	1	10/31/2021 04:34	<a href="#">WG1766323</a>
sec-Butylbenzene	ND		0.0161	1	10/31/2021 04:34	<a href="#">WG1766323</a>
tert-Butylbenzene	ND		0.00646	1	10/31/2021 04:34	<a href="#">WG1766323</a>
Carbon tetrachloride	ND		0.00646	1	10/31/2021 04:34	<a href="#">WG1766323</a>
Chlorobenzene	ND		0.00323	1	10/31/2021 04:34	<a href="#">WG1766323</a>
Chlorodibromomethane	ND		0.00323	1	10/31/2021 04:34	<a href="#">WG1766323</a>
Chloroethane	ND		0.00646	1	10/31/2021 04:34	<a href="#">WG1766323</a>
Chloroform	ND		0.00323	1	10/31/2021 04:34	<a href="#">WG1766323</a>
Chloromethane	ND		0.0161	1	10/31/2021 04:34	<a href="#">WG1766323</a>
2-Chlorotoluene	ND		0.00323	1	10/31/2021 04:34	<a href="#">WG1766323</a>
4-Chlorotoluene	ND		0.00646	1	10/31/2021 04:34	<a href="#">WG1766323</a>
1,2-Dibromo-3-Chloropropane	ND		0.0323	1	10/31/2021 04:34	<a href="#">WG1766323</a>
1,2-Dibromoethane	ND		0.00323	1	10/31/2021 04:34	<a href="#">WG1766323</a>
Dibromomethane	ND		0.00646	1	10/31/2021 04:34	<a href="#">WG1766323</a>
1,2-Dichlorobenzene	ND		0.00646	1	10/31/2021 04:34	<a href="#">WG1766323</a>
1,3-Dichlorobenzene	ND		0.00646	1	10/31/2021 04:34	<a href="#">WG1766323</a>
1,4-Dichlorobenzene	ND		0.00646	1	10/31/2021 04:34	<a href="#">WG1766323</a>
Dichlorodifluoromethane	ND		0.00323	1	10/31/2021 04:34	<a href="#">WG1766323</a>
1,1-Dichloroethane	ND		0.00323	1	10/31/2021 04:34	<a href="#">WG1766323</a>
1,2-Dichloroethane	ND		0.00323	1	10/31/2021 04:34	<a href="#">WG1766323</a>



Volatile Organic Compounds (GC/MS) by Method 8260D

Analyte	Result (dry) mg/kg	Qualifier	RDL (dry) mg/kg	Dilution	Analysis date / time	Batch
1,1-Dichloroethene	ND		0.00323	1	10/31/2021 04:34	WG1766323
cis-1,2-Dichloroethene	ND		0.00323	1	10/31/2021 04:34	WG1766323
trans-1,2-Dichloroethene	ND		0.00646	1	10/31/2021 04:34	WG1766323
1,2-Dichloropropane	ND		0.00646	1	10/31/2021 04:34	WG1766323
1,1-Dichloropropene	ND		0.00323	1	10/31/2021 04:34	WG1766323
1,3-Dichloropropane	ND		0.00646	1	10/31/2021 04:34	WG1766323
cis-1,3-Dichloropropene	ND		0.00323	1	10/31/2021 04:34	WG1766323
trans-1,3-Dichloropropene	ND		0.00646	1	10/31/2021 04:34	WG1766323
2,2-Dichloropropane	ND		0.00323	1	10/31/2021 04:34	WG1766323
Di-isopropyl ether	ND		0.00129	1	10/31/2021 04:34	WG1766323
Ethylbenzene	ND		0.00323	1	10/31/2021 04:34	WG1766323
Hexachloro-1,3-butadiene	ND		0.0323	1	10/31/2021 04:34	WG1766323
Isopropylbenzene	ND		0.00323	1	10/31/2021 04:34	WG1766323
p-Isopropyltoluene	ND		0.00646	1	10/31/2021 04:34	WG1766323
2-Butanone (MEK)	ND		0.129	1	10/31/2021 04:34	WG1766323
Methylene Chloride	ND		0.0323	1	10/31/2021 04:34	WG1766323
4-Methyl-2-pentanone (MIBK)	ND		0.0323	1	10/31/2021 04:34	WG1766323
Methyl tert-butyl ether	ND		0.00129	1	10/31/2021 04:34	WG1766323
Naphthalene	ND		0.0161	1	10/31/2021 04:34	WG1766323
n-Propylbenzene	ND		0.00646	1	10/31/2021 04:34	WG1766323
Styrene	ND		0.0161	1	10/31/2021 04:34	WG1766323
1,1,1,2-Tetrachloroethane	ND		0.00323	1	10/31/2021 04:34	WG1766323
1,1,2,2-Tetrachloroethane	ND		0.00323	1	10/31/2021 04:34	WG1766323
1,1,2-Trichlorotrifluoroethane	ND		0.00323	1	10/31/2021 04:34	WG1766323
Tetrachloroethene	ND		0.00323	1	10/31/2021 04:34	WG1766323
Toluene	ND		0.00646	1	10/31/2021 04:34	WG1766323
1,2,3-Trichlorobenzene	ND	C4	0.0161	1	10/31/2021 04:34	WG1766323
1,2,4-Trichlorobenzene	ND		0.0161	1	10/31/2021 04:34	WG1766323
1,1,1-Trichloroethane	ND		0.00323	1	10/31/2021 04:34	WG1766323
1,1,2-Trichloroethane	ND		0.00323	1	10/31/2021 04:34	WG1766323
Trichloroethene	ND		0.00129	1	10/31/2021 04:34	WG1766323
Trichlorofluoromethane	ND		0.00323	1	10/31/2021 04:34	WG1766323
1,2,3-Trichloropropane	ND		0.0161	1	10/31/2021 04:34	WG1766323
1,2,4-Trimethylbenzene	ND		0.00646	1	10/31/2021 04:34	WG1766323
1,2,3-Trimethylbenzene	ND		0.00646	1	10/31/2021 04:34	WG1766323
1,3,5-Trimethylbenzene	ND		0.00646	1	10/31/2021 04:34	WG1766323
Vinyl chloride	ND		0.00323	1	10/31/2021 04:34	WG1766323
Xylenes, Total	ND		0.00839	1	10/31/2021 04:34	WG1766323
(S) Toluene-d8	104		75.0-131		10/31/2021 04:34	WG1766323
(S) 4-Bromofluorobenzene	99.4		67.0-138		10/31/2021 04:34	WG1766323
(S) 1,2-Dichloroethane-d4	99.4		70.0-130		10/31/2021 04:34	WG1766323

1 Cp  
2 Tc  
3 Ss  
4 Cn  
5 Sr  
6 Qc  
7 Gl  
8 Al  
9 Sc

Semi-Volatile Organic Compounds (GC) by Method NWTPHDX-NO SGT

Analyte	Result (dry) mg/kg	Qualifier	RDL (dry) mg/kg	Dilution	Analysis date / time	Batch
Diesel Range Organics (DRO)	ND		4.50	1	11/03/2021 06:24	WG1766948
Residual Range Organics (RRO)	ND		11.2	1	11/03/2021 06:24	WG1766948
(S) o-Terphenyl	45.3		18.0-148		11/03/2021 06:24	WG1766948

Semi Volatile Organic Compounds (GC/MS) by Method 8270E-SIM

Analyte	Result (dry) mg/kg	Qualifier	RDL (dry) mg/kg	Dilution	Analysis date / time	Batch
Anthracene	ND		0.00675	1	10/30/2021 13:35	<a href="#">WG1765811</a>
Acenaphthene	ND		0.00675	1	10/30/2021 13:35	<a href="#">WG1765811</a>
Acenaphthylene	ND		0.00675	1	10/30/2021 13:35	<a href="#">WG1765811</a>
Benzo(a)anthracene	ND		0.00675	1	10/30/2021 13:35	<a href="#">WG1765811</a>
Benzo(a)pyrene	ND		0.00675	1	10/30/2021 13:35	<a href="#">WG1765811</a>
Benzo(b)fluoranthene	ND		0.00675	1	10/30/2021 13:35	<a href="#">WG1765811</a>
Benzo(g,h,i)perylene	ND		0.00675	1	10/30/2021 13:35	<a href="#">WG1765811</a>
Benzo(k)fluoranthene	ND		0.00675	1	10/30/2021 13:35	<a href="#">WG1765811</a>
Chrysene	ND		0.00675	1	10/30/2021 13:35	<a href="#">WG1765811</a>
Dibenz(a,h)anthracene	ND		0.00675	1	10/30/2021 13:35	<a href="#">WG1765811</a>
Fluoranthene	ND		0.00675	1	10/30/2021 13:35	<a href="#">WG1765811</a>
Fluorene	ND		0.00675	1	10/30/2021 13:35	<a href="#">WG1765811</a>
Indeno(1,2,3-cd)pyrene	ND		0.00675	1	10/30/2021 13:35	<a href="#">WG1765811</a>
Naphthalene	ND		0.0225	1	10/30/2021 13:35	<a href="#">WG1765811</a>
Phenanthrene	ND		0.00675	1	10/30/2021 13:35	<a href="#">WG1765811</a>
Pyrene	ND		0.00675	1	10/30/2021 13:35	<a href="#">WG1765811</a>
1-Methylnaphthalene	ND		0.0225	1	10/30/2021 13:35	<a href="#">WG1765811</a>
2-Methylnaphthalene	ND		0.0225	1	10/30/2021 13:35	<a href="#">WG1765811</a>
2-Chloronaphthalene	ND		0.0225	1	10/30/2021 13:35	<a href="#">WG1765811</a>
(S) Nitrobenzene-d5	93.3		14.0-149		10/30/2021 13:35	<a href="#">WG1765811</a>
(S) 2-Fluorobiphenyl	75.9		34.0-125		10/30/2021 13:35	<a href="#">WG1765811</a>
(S) p-Terphenyl-d14	96.6		23.0-120		10/30/2021 13:35	<a href="#">WG1765811</a>

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc

Total Solids by Method 2540 G-2011

Analyte	Result	Qualifier	Dilution	Analysis	Batch
	%			date / time	
Total Solids	90.2		1	10/30/2021 16:39	<a href="#">WG1765837</a>

Mercury by Method 7471B

Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
	mg/kg		mg/kg		date / time	
Mercury	ND		0.0443	1	11/02/2021 15:12	<a href="#">WG1766498</a>

Metals (ICPMS) by Method 6020B

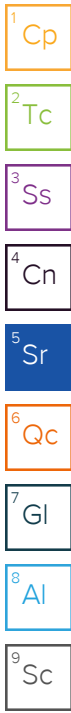
Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
	mg/kg		mg/kg		date / time	
Arsenic	3.17		1.11	5	11/03/2021 18:35	<a href="#">WG1765966</a>
Barium	44.9		2.77	5	11/03/2021 18:35	<a href="#">WG1765966</a>
Cadmium	ND		1.11	5	11/03/2021 18:35	<a href="#">WG1765966</a>
Chromium	24.4		5.54	5	11/03/2021 18:35	<a href="#">WG1765966</a>
Lead	ND		2.22	5	11/03/2021 18:35	<a href="#">WG1765966</a>
Selenium	ND		2.77	5	11/03/2021 18:35	<a href="#">WG1765966</a>
Silver	ND		0.554	5	11/03/2021 18:35	<a href="#">WG1765966</a>

Volatile Organic Compounds (GC) by Method NWTPHGX

Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
	mg/kg		mg/kg		date / time	
Gasoline Range Organics-NWTPH	ND		3.08	25	10/29/2021 04:21	<a href="#">WG1765548</a>
(S) a,a,a-Trifluorotoluene(FID)	91.8		77.0-120		10/29/2021 04:21	<a href="#">WG1765548</a>

Volatile Organic Compounds (GC/MS) by Method 8260D

Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
	mg/kg		mg/kg		date / time	
Acetone	ND		0.0616	1	10/31/2021 04:53	<a href="#">WG1766323</a>
Acrylonitrile	ND	<a href="#">C3</a>	0.0154	1	10/31/2021 04:53	<a href="#">WG1766323</a>
Benzene	ND		0.00123	1	10/31/2021 04:53	<a href="#">WG1766323</a>
Bromobenzene	ND		0.0154	1	10/31/2021 04:53	<a href="#">WG1766323</a>
Bromodichloromethane	ND		0.00308	1	10/31/2021 04:53	<a href="#">WG1766323</a>
Bromoform	ND		0.0308	1	10/31/2021 04:53	<a href="#">WG1766323</a>
Bromomethane	ND		0.0154	1	10/31/2021 04:53	<a href="#">WG1766323</a>
n-Butylbenzene	ND		0.0154	1	10/31/2021 04:53	<a href="#">WG1766323</a>
sec-Butylbenzene	ND		0.0154	1	10/31/2021 04:53	<a href="#">WG1766323</a>
tert-Butylbenzene	ND		0.00616	1	10/31/2021 04:53	<a href="#">WG1766323</a>
Carbon tetrachloride	ND		0.00616	1	10/31/2021 04:53	<a href="#">WG1766323</a>
Chlorobenzene	ND		0.00308	1	10/31/2021 04:53	<a href="#">WG1766323</a>
Chlorodibromomethane	ND		0.00308	1	10/31/2021 04:53	<a href="#">WG1766323</a>
Chloroethane	ND		0.00616	1	10/31/2021 04:53	<a href="#">WG1766323</a>
Chloroform	ND		0.00308	1	10/31/2021 04:53	<a href="#">WG1766323</a>
Chloromethane	ND		0.0154	1	10/31/2021 04:53	<a href="#">WG1766323</a>
2-Chlorotoluene	ND		0.00308	1	10/31/2021 04:53	<a href="#">WG1766323</a>
4-Chlorotoluene	ND		0.00616	1	10/31/2021 04:53	<a href="#">WG1766323</a>
1,2-Dibromo-3-Chloropropane	ND		0.0308	1	10/31/2021 04:53	<a href="#">WG1766323</a>
1,2-Dibromoethane	ND		0.00308	1	10/31/2021 04:53	<a href="#">WG1766323</a>
Dibromomethane	ND		0.00616	1	10/31/2021 04:53	<a href="#">WG1766323</a>
1,2-Dichlorobenzene	ND		0.00616	1	10/31/2021 04:53	<a href="#">WG1766323</a>
1,3-Dichlorobenzene	ND		0.00616	1	10/31/2021 04:53	<a href="#">WG1766323</a>
1,4-Dichlorobenzene	ND		0.00616	1	10/31/2021 04:53	<a href="#">WG1766323</a>
Dichlorodifluoromethane	ND		0.00308	1	10/31/2021 04:53	<a href="#">WG1766323</a>
1,1-Dichloroethane	ND		0.00308	1	10/31/2021 04:53	<a href="#">WG1766323</a>
1,2-Dichloroethane	ND		0.00308	1	10/31/2021 04:53	<a href="#">WG1766323</a>



Volatile Organic Compounds (GC/MS) by Method 8260D

Analyte	Result (dry) mg/kg	Qualifier	RDL (dry) mg/kg	Dilution	Analysis date / time	Batch
1,1-Dichloroethene	ND		0.00308	1	10/31/2021 04:53	WG1766323
cis-1,2-Dichloroethene	ND		0.00308	1	10/31/2021 04:53	WG1766323
trans-1,2-Dichloroethene	ND		0.00616	1	10/31/2021 04:53	WG1766323
1,2-Dichloropropane	ND		0.00616	1	10/31/2021 04:53	WG1766323
1,1-Dichloropropene	ND		0.00308	1	10/31/2021 04:53	WG1766323
1,3-Dichloropropane	ND		0.00616	1	10/31/2021 04:53	WG1766323
cis-1,3-Dichloropropene	ND		0.00308	1	10/31/2021 04:53	WG1766323
trans-1,3-Dichloropropene	ND		0.00616	1	10/31/2021 04:53	WG1766323
2,2-Dichloropropane	ND		0.00308	1	10/31/2021 04:53	WG1766323
Di-isopropyl ether	ND		0.00123	1	10/31/2021 04:53	WG1766323
Ethylbenzene	ND		0.00308	1	10/31/2021 04:53	WG1766323
Hexachloro-1,3-butadiene	ND		0.0308	1	10/31/2021 04:53	WG1766323
Isopropylbenzene	ND		0.00308	1	10/31/2021 04:53	WG1766323
p-Isopropyltoluene	ND		0.00616	1	10/31/2021 04:53	WG1766323
2-Butanone (MEK)	ND		0.123	1	10/31/2021 04:53	WG1766323
Methylene Chloride	ND		0.0308	1	10/31/2021 04:53	WG1766323
4-Methyl-2-pentanone (MIBK)	ND		0.0308	1	10/31/2021 04:53	WG1766323
Methyl tert-butyl ether	ND		0.00123	1	10/31/2021 04:53	WG1766323
Naphthalene	ND		0.0154	1	10/31/2021 04:53	WG1766323
n-Propylbenzene	ND		0.00616	1	10/31/2021 04:53	WG1766323
Styrene	ND		0.0154	1	10/31/2021 04:53	WG1766323
1,1,1,2-Tetrachloroethane	ND		0.00308	1	10/31/2021 04:53	WG1766323
1,1,2,2-Tetrachloroethane	ND		0.00308	1	10/31/2021 04:53	WG1766323
1,1,2-Trichlorotrifluoroethane	ND		0.00308	1	10/31/2021 04:53	WG1766323
Tetrachloroethene	ND		0.00308	1	10/31/2021 04:53	WG1766323
Toluene	ND		0.00616	1	10/31/2021 04:53	WG1766323
1,2,3-Trichlorobenzene	ND	C4	0.0154	1	10/31/2021 04:53	WG1766323
1,2,4-Trichlorobenzene	ND		0.0154	1	10/31/2021 04:53	WG1766323
1,1,1-Trichloroethane	ND		0.00308	1	10/31/2021 04:53	WG1766323
1,1,2-Trichloroethane	ND		0.00308	1	10/31/2021 04:53	WG1766323
Trichloroethene	ND		0.00123	1	10/31/2021 04:53	WG1766323
Trichlorofluoromethane	ND		0.00308	1	10/31/2021 04:53	WG1766323
1,2,3-Trichloropropane	ND		0.0154	1	10/31/2021 04:53	WG1766323
1,2,4-Trimethylbenzene	ND		0.00616	1	10/31/2021 04:53	WG1766323
1,2,3-Trimethylbenzene	ND		0.00616	1	10/31/2021 04:53	WG1766323
1,3,5-Trimethylbenzene	ND		0.00616	1	10/31/2021 04:53	WG1766323
Vinyl chloride	ND		0.00308	1	10/31/2021 04:53	WG1766323
Xylenes, Total	ND		0.00801	1	10/31/2021 04:53	WG1766323
(S) Toluene-d8	104		75.0-131		10/31/2021 04:53	WG1766323
(S) 4-Bromofluorobenzene	100		67.0-138		10/31/2021 04:53	WG1766323
(S) 1,2-Dichloroethane-d4	101		70.0-130		10/31/2021 04:53	WG1766323

1 Cp  
2 Tc  
3 Ss  
4 Cn  
5 Sr  
6 Qc  
7 Gl  
8 Al  
9 Sc

Semi-Volatile Organic Compounds (GC) by Method NWTPHDX-NO SGT

Analyte	Result (dry) mg/kg	Qualifier	RDL (dry) mg/kg	Dilution	Analysis date / time	Batch
Diesel Range Organics (DRO)	ND		4.43	1	11/03/2021 06:36	WG1766948
Residual Range Organics (RRO)	ND		11.1	1	11/03/2021 06:36	WG1766948
(S) o-Terphenyl	45.8		18.0-148		11/03/2021 06:36	WG1766948

Semi Volatile Organic Compounds (GC/MS) by Method 8270E-SIM

Analyte	Result (dry) mg/kg	Qualifier	RDL (dry) mg/kg	Dilution	Analysis date / time	Batch
Anthracene	ND		0.00665	1	10/30/2021 13:52	<a href="#">WG1765811</a>
Acenaphthene	ND		0.00665	1	10/30/2021 13:52	<a href="#">WG1765811</a>
Acenaphthylene	ND		0.00665	1	10/30/2021 13:52	<a href="#">WG1765811</a>
Benzo(a)anthracene	ND		0.00665	1	10/30/2021 13:52	<a href="#">WG1765811</a>
Benzo(a)pyrene	ND		0.00665	1	10/30/2021 13:52	<a href="#">WG1765811</a>
Benzo(b)fluoranthene	ND		0.00665	1	10/30/2021 13:52	<a href="#">WG1765811</a>
Benzo(g,h,i)perylene	ND		0.00665	1	10/30/2021 13:52	<a href="#">WG1765811</a>
Benzo(k)fluoranthene	ND		0.00665	1	10/30/2021 13:52	<a href="#">WG1765811</a>
Chrysene	ND		0.00665	1	10/30/2021 13:52	<a href="#">WG1765811</a>
Dibenz(a,h)anthracene	ND		0.00665	1	10/30/2021 13:52	<a href="#">WG1765811</a>
Fluoranthene	ND		0.00665	1	10/30/2021 13:52	<a href="#">WG1765811</a>
Fluorene	ND		0.00665	1	10/30/2021 13:52	<a href="#">WG1765811</a>
Indeno(1,2,3-cd)pyrene	ND		0.00665	1	10/30/2021 13:52	<a href="#">WG1765811</a>
Naphthalene	ND		0.0222	1	10/30/2021 13:52	<a href="#">WG1765811</a>
Phenanthrene	ND		0.00665	1	10/30/2021 13:52	<a href="#">WG1765811</a>
Pyrene	ND		0.00665	1	10/30/2021 13:52	<a href="#">WG1765811</a>
1-Methylnaphthalene	ND		0.0222	1	10/30/2021 13:52	<a href="#">WG1765811</a>
2-Methylnaphthalene	ND		0.0222	1	10/30/2021 13:52	<a href="#">WG1765811</a>
2-Chloronaphthalene	ND		0.0222	1	10/30/2021 13:52	<a href="#">WG1765811</a>
<i>(S)</i> Nitrobenzene-d5	94.9		14.0-149		10/30/2021 13:52	<a href="#">WG1765811</a>
<i>(S)</i> 2-Fluorobiphenyl	81.7		34.0-125		10/30/2021 13:52	<a href="#">WG1765811</a>
<i>(S)</i> p-Terphenyl-d14	95.1		23.0-120		10/30/2021 13:52	<a href="#">WG1765811</a>

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc

Total Solids by Method 2540 G-2011

Analyte	Result	Qualifier	Dilution	Analysis	Batch
Total Solids	91.6		1	10/30/2021 16:39	<a href="#">WG1765837</a>

Mercury by Method 7471B

Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
Mercury	0.0731		0.0437	1	11/02/2021 15:15	<a href="#">WG1766498</a>

Metals (ICPMS) by Method 6020B

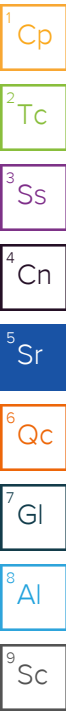
Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
Arsenic	6.28		1.09	5	11/03/2021 18:39	<a href="#">WG1765966</a>
Barium	72.2		2.73	5	11/03/2021 18:39	<a href="#">WG1765966</a>
Cadmium	ND		1.09	5	11/03/2021 18:39	<a href="#">WG1765966</a>
Chromium	33.9		5.46	5	11/03/2021 18:39	<a href="#">WG1765966</a>
Lead	27.6		2.18	5	11/03/2021 18:39	<a href="#">WG1765966</a>
Selenium	ND		2.73	5	11/03/2021 18:39	<a href="#">WG1765966</a>
Silver	ND		0.546	5	11/03/2021 18:39	<a href="#">WG1765966</a>

Volatile Organic Compounds (GC) by Method NWTPHGX

Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
Gasoline Range Organics-NWTPH	ND		2.97	25	10/29/2021 04:43	<a href="#">WG1765548</a>
(S) a,a,a-Trifluorotoluene(FID)	94.7		77.0-120		10/29/2021 04:43	<a href="#">WG1765548</a>

Volatile Organic Compounds (GC/MS) by Method 8260D

Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
Acetone	ND		0.0594	1	10/31/2021 05:12	<a href="#">WG1766323</a>
Acrylonitrile	ND	<a href="#">C3</a>	0.0148	1	10/31/2021 05:12	<a href="#">WG1766323</a>
Benzene	ND		0.00119	1	10/31/2021 05:12	<a href="#">WG1766323</a>
Bromobenzene	ND		0.0148	1	10/31/2021 05:12	<a href="#">WG1766323</a>
Bromodichloromethane	ND		0.00297	1	10/31/2021 05:12	<a href="#">WG1766323</a>
Bromoform	ND		0.0297	1	10/31/2021 05:12	<a href="#">WG1766323</a>
Bromomethane	ND		0.0148	1	10/31/2021 05:12	<a href="#">WG1766323</a>
n-Butylbenzene	ND		0.0148	1	10/31/2021 05:12	<a href="#">WG1766323</a>
sec-Butylbenzene	ND		0.0148	1	10/31/2021 05:12	<a href="#">WG1766323</a>
tert-Butylbenzene	ND		0.00594	1	10/31/2021 05:12	<a href="#">WG1766323</a>
Carbon tetrachloride	ND		0.00594	1	10/31/2021 05:12	<a href="#">WG1766323</a>
Chlorobenzene	ND		0.00297	1	10/31/2021 05:12	<a href="#">WG1766323</a>
Chlorodibromomethane	ND		0.00297	1	10/31/2021 05:12	<a href="#">WG1766323</a>
Chloroethane	ND		0.00594	1	10/31/2021 05:12	<a href="#">WG1766323</a>
Chloroform	ND		0.00297	1	10/31/2021 05:12	<a href="#">WG1766323</a>
Chloromethane	ND		0.0148	1	10/31/2021 05:12	<a href="#">WG1766323</a>
2-Chlorotoluene	ND		0.00297	1	10/31/2021 05:12	<a href="#">WG1766323</a>
4-Chlorotoluene	ND		0.00594	1	10/31/2021 05:12	<a href="#">WG1766323</a>
1,2-Dibromo-3-Chloropropane	ND		0.0297	1	10/31/2021 05:12	<a href="#">WG1766323</a>
1,2-Dibromoethane	ND		0.00297	1	10/31/2021 05:12	<a href="#">WG1766323</a>
Dibromomethane	ND		0.00594	1	10/31/2021 05:12	<a href="#">WG1766323</a>
1,2-Dichlorobenzene	ND		0.00594	1	10/31/2021 05:12	<a href="#">WG1766323</a>
1,3-Dichlorobenzene	ND		0.00594	1	10/31/2021 05:12	<a href="#">WG1766323</a>
1,4-Dichlorobenzene	ND		0.00594	1	10/31/2021 05:12	<a href="#">WG1766323</a>
Dichlorodifluoromethane	ND		0.00297	1	10/31/2021 05:12	<a href="#">WG1766323</a>
1,1-Dichloroethane	ND		0.00297	1	10/31/2021 05:12	<a href="#">WG1766323</a>
1,2-Dichloroethane	ND		0.00297	1	10/31/2021 05:12	<a href="#">WG1766323</a>



Volatile Organic Compounds (GC/MS) by Method 8260D

Analyte	Result (dry) mg/kg	Qualifier	RDL (dry) mg/kg	Dilution	Analysis date / time	Batch
1,1-Dichloroethene	ND		0.00297	1	10/31/2021 05:12	WG1766323
cis-1,2-Dichloroethene	ND		0.00297	1	10/31/2021 05:12	WG1766323
trans-1,2-Dichloroethene	ND		0.00594	1	10/31/2021 05:12	WG1766323
1,2-Dichloropropane	ND		0.00594	1	10/31/2021 05:12	WG1766323
1,1-Dichloropropene	ND		0.00297	1	10/31/2021 05:12	WG1766323
1,3-Dichloropropane	ND		0.00594	1	10/31/2021 05:12	WG1766323
cis-1,3-Dichloropropene	ND		0.00297	1	10/31/2021 05:12	WG1766323
trans-1,3-Dichloropropene	ND		0.00594	1	10/31/2021 05:12	WG1766323
2,2-Dichloropropane	ND		0.00297	1	10/31/2021 05:12	WG1766323
Di-isopropyl ether	ND		0.00119	1	10/31/2021 05:12	WG1766323
Ethylbenzene	ND		0.00297	1	10/31/2021 05:12	WG1766323
Hexachloro-1,3-butadiene	ND		0.0297	1	10/31/2021 05:12	WG1766323
Isopropylbenzene	ND		0.00297	1	10/31/2021 05:12	WG1766323
p-Isopropyltoluene	ND		0.00594	1	10/31/2021 05:12	WG1766323
2-Butanone (MEK)	ND		0.119	1	10/31/2021 05:12	WG1766323
Methylene Chloride	ND		0.0297	1	10/31/2021 05:12	WG1766323
4-Methyl-2-pentanone (MIBK)	ND		0.0297	1	10/31/2021 05:12	WG1766323
Methyl tert-butyl ether	ND		0.00119	1	10/31/2021 05:12	WG1766323
Naphthalene	ND		0.0148	1	10/31/2021 05:12	WG1766323
n-Propylbenzene	ND		0.00594	1	10/31/2021 05:12	WG1766323
Styrene	ND		0.0148	1	10/31/2021 05:12	WG1766323
1,1,1,2-Tetrachloroethane	ND		0.00297	1	10/31/2021 05:12	WG1766323
1,1,2,2-Tetrachloroethane	ND		0.00297	1	10/31/2021 05:12	WG1766323
1,1,2-Trichlorotrifluoroethane	ND		0.00297	1	10/31/2021 05:12	WG1766323
Tetrachloroethene	ND		0.00297	1	10/31/2021 05:12	WG1766323
Toluene	ND		0.00594	1	10/31/2021 05:12	WG1766323
1,2,3-Trichlorobenzene	ND	C4	0.0148	1	10/31/2021 05:12	WG1766323
1,2,4-Trichlorobenzene	ND		0.0148	1	10/31/2021 05:12	WG1766323
1,1,1-Trichloroethane	ND		0.00297	1	10/31/2021 05:12	WG1766323
1,1,2-Trichloroethane	ND		0.00297	1	10/31/2021 05:12	WG1766323
Trichloroethene	ND		0.00119	1	10/31/2021 05:12	WG1766323
Trichlorofluoromethane	ND		0.00297	1	10/31/2021 05:12	WG1766323
1,2,3-Trichloropropane	ND		0.0148	1	10/31/2021 05:12	WG1766323
1,2,4-Trimethylbenzene	ND		0.00594	1	10/31/2021 05:12	WG1766323
1,2,3-Trimethylbenzene	ND		0.00594	1	10/31/2021 05:12	WG1766323
1,3,5-Trimethylbenzene	ND		0.00594	1	10/31/2021 05:12	WG1766323
Vinyl chloride	ND		0.00297	1	10/31/2021 05:12	WG1766323
Xylenes, Total	ND		0.00772	1	10/31/2021 05:12	WG1766323
(S) Toluene-d8	104		75.0-131		10/31/2021 05:12	WG1766323
(S) 4-Bromofluorobenzene	101		67.0-138		10/31/2021 05:12	WG1766323
(S) 1,2-Dichloroethane-d4	105		70.0-130		10/31/2021 05:12	WG1766323

1  
Cp

2  
Tc

3  
Ss

4  
Cn

5  
Sr

6  
Qc

7  
Gl

8  
Al

9  
Sc

Semi-Volatile Organic Compounds (GC) by Method NWTPHDX-NO SGT

Analyte	Result (dry) mg/kg	Qualifier	RDL (dry) mg/kg	Dilution	Analysis date / time	Batch
Diesel Range Organics (DRO)	ND		21.8	5	11/03/2021 15:22	WG1766948
Residual Range Organics (RRO)	166		54.6	5	11/03/2021 15:22	WG1766948
(S) o-Terphenyl	60.1		18.0-148		11/03/2021 15:22	WG1766948

Sample Narrative:

L1423266-13 WG1766948: Cannot run at lower dilution due to viscosity of extract

Semi Volatile Organic Compounds (GC/MS) by Method 8270E-SIM

Analyte	Result (dry) mg/kg	Qualifier	RDL (dry) mg/kg	Dilution	Analysis date / time	Batch
Anthracene	ND		0.00655	1	10/30/2021 16:11	<a href="#">WG1765811</a>
Acenaphthene	ND		0.00655	1	10/30/2021 16:11	<a href="#">WG1765811</a>
Acenaphthylene	ND		0.00655	1	10/30/2021 16:11	<a href="#">WG1765811</a>
Benzo(a)anthracene	0.0298		0.00655	1	10/30/2021 16:11	<a href="#">WG1765811</a>
Benzo(a)pyrene	0.0363		0.00655	1	10/30/2021 16:11	<a href="#">WG1765811</a>
Benzo(b)fluoranthene	0.0500		0.00655	1	10/30/2021 16:11	<a href="#">WG1765811</a>
Benzo(g,h,i)perylene	0.0306		0.00655	1	10/30/2021 16:11	<a href="#">WG1765811</a>
Benzo(k)fluoranthene	0.0150		0.00655	1	10/30/2021 16:11	<a href="#">WG1765811</a>
Chrysene	0.0354		0.00655	1	10/30/2021 16:11	<a href="#">WG1765811</a>
Dibenz(a,h)anthracene	0.00660		0.00655	1	10/30/2021 16:11	<a href="#">WG1765811</a>
Fluoranthene	0.0622		0.00655	1	10/30/2021 16:11	<a href="#">WG1765811</a>
Fluorene	ND		0.00655	1	10/30/2021 16:11	<a href="#">WG1765811</a>
Indeno(1,2,3-cd)pyrene	0.0325		0.00655	1	10/30/2021 16:11	<a href="#">WG1765811</a>
Naphthalene	ND		0.0218	1	10/30/2021 16:11	<a href="#">WG1765811</a>
Phenanthrene	0.0244		0.00655	1	10/30/2021 16:11	<a href="#">WG1765811</a>
Pyrene	0.0474		0.00655	1	10/30/2021 16:11	<a href="#">WG1765811</a>
1-Methylnaphthalene	ND		0.0218	1	10/30/2021 16:11	<a href="#">WG1765811</a>
2-Methylnaphthalene	ND		0.0218	1	10/30/2021 16:11	<a href="#">WG1765811</a>
2-Chloronaphthalene	ND		0.0218	1	10/30/2021 16:11	<a href="#">WG1765811</a>
(S) Nitrobenzene-d5	113		14.0-149		10/30/2021 16:11	<a href="#">WG1765811</a>
(S) 2-Fluorobiphenyl	81.7		34.0-125		10/30/2021 16:11	<a href="#">WG1765811</a>
(S) p-Terphenyl-d14	99.3		23.0-120		10/30/2021 16:11	<a href="#">WG1765811</a>

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc

Total Solids by Method 2540 G-2011

Analyte	Result	Qualifier	Dilution	Analysis	Batch
	%			date / time	
Total Solids	85.4		1	10/30/2021 16:39	<a href="#">WG1765837</a>

Mercury by Method 7471B

Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
	mg/kg		mg/kg		date / time	
Mercury	ND		0.0469	1	11/02/2021 15:22	<a href="#">WG1766498</a>

Metals (ICPMS) by Method 6020B

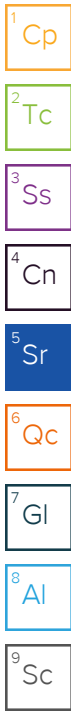
Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
	mg/kg		mg/kg		date / time	
Arsenic	3.70		1.17	5	11/03/2021 18:42	<a href="#">WG1765966</a>
Barium	61.0		2.93	5	11/03/2021 18:42	<a href="#">WG1765966</a>
Cadmium	ND		1.17	5	11/03/2021 18:42	<a href="#">WG1765966</a>
Chromium	26.7		5.86	5	11/03/2021 18:42	<a href="#">WG1765966</a>
Lead	ND		2.34	5	11/03/2021 18:42	<a href="#">WG1765966</a>
Selenium	ND		2.93	5	11/03/2021 18:42	<a href="#">WG1765966</a>
Silver	ND		0.586	5	11/03/2021 18:42	<a href="#">WG1765966</a>

Volatile Organic Compounds (GC) by Method NWTPHGX

Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
	mg/kg		mg/kg		date / time	
Gasoline Range Organics-NWTPH	ND		3.43	25	10/29/2021 05:05	<a href="#">WG1765548</a>
(S) a,a,a-Trifluorotoluene(FID)	88.3		77.0-120		10/29/2021 05:05	<a href="#">WG1765548</a>

Volatile Organic Compounds (GC/MS) by Method 8260D

Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
	mg/kg		mg/kg		date / time	
Acetone	ND		0.0686	1	10/31/2021 05:31	<a href="#">WG1766323</a>
Acrylonitrile	ND	<a href="#">C3</a>	0.0171	1	10/31/2021 05:31	<a href="#">WG1766323</a>
Benzene	ND		0.00137	1	10/31/2021 05:31	<a href="#">WG1766323</a>
Bromobenzene	ND		0.0171	1	10/31/2021 05:31	<a href="#">WG1766323</a>
Bromodichloromethane	ND		0.00343	1	10/31/2021 05:31	<a href="#">WG1766323</a>
Bromoform	ND		0.0343	1	10/31/2021 05:31	<a href="#">WG1766323</a>
Bromomethane	ND		0.0171	1	10/31/2021 05:31	<a href="#">WG1766323</a>
n-Butylbenzene	ND		0.0171	1	10/31/2021 05:31	<a href="#">WG1766323</a>
sec-Butylbenzene	ND		0.0171	1	10/31/2021 05:31	<a href="#">WG1766323</a>
tert-Butylbenzene	ND		0.00686	1	10/31/2021 05:31	<a href="#">WG1766323</a>
Carbon tetrachloride	ND		0.00686	1	10/31/2021 05:31	<a href="#">WG1766323</a>
Chlorobenzene	ND		0.00343	1	10/31/2021 05:31	<a href="#">WG1766323</a>
Chlorodibromomethane	ND		0.00343	1	10/31/2021 05:31	<a href="#">WG1766323</a>
Chloroethane	ND		0.00686	1	10/31/2021 05:31	<a href="#">WG1766323</a>
Chloroform	ND		0.00343	1	10/31/2021 05:31	<a href="#">WG1766323</a>
Chloromethane	ND		0.0171	1	10/31/2021 05:31	<a href="#">WG1766323</a>
2-Chlorotoluene	ND		0.00343	1	10/31/2021 05:31	<a href="#">WG1766323</a>
4-Chlorotoluene	ND		0.00686	1	10/31/2021 05:31	<a href="#">WG1766323</a>
1,2-Dibromo-3-Chloropropane	ND		0.0343	1	10/31/2021 05:31	<a href="#">WG1766323</a>
1,2-Dibromoethane	ND		0.00343	1	10/31/2021 05:31	<a href="#">WG1766323</a>
Dibromomethane	ND		0.00686	1	10/31/2021 05:31	<a href="#">WG1766323</a>
1,2-Dichlorobenzene	ND		0.00686	1	10/31/2021 05:31	<a href="#">WG1766323</a>
1,3-Dichlorobenzene	ND		0.00686	1	10/31/2021 05:31	<a href="#">WG1766323</a>
1,4-Dichlorobenzene	ND		0.00686	1	10/31/2021 05:31	<a href="#">WG1766323</a>
Dichlorodifluoromethane	ND		0.00343	1	10/31/2021 05:31	<a href="#">WG1766323</a>
1,1-Dichloroethane	ND		0.00343	1	10/31/2021 05:31	<a href="#">WG1766323</a>
1,2-Dichloroethane	ND		0.00343	1	10/31/2021 05:31	<a href="#">WG1766323</a>



Volatile Organic Compounds (GC/MS) by Method 8260D

Analyte	Result (dry) mg/kg	Qualifier	RDL (dry) mg/kg	Dilution	Analysis date / time	Batch
1,1-Dichloroethene	ND		0.00343	1	10/31/2021 05:31	WG1766323
cis-1,2-Dichloroethene	ND		0.00343	1	10/31/2021 05:31	WG1766323
trans-1,2-Dichloroethene	ND		0.00686	1	10/31/2021 05:31	WG1766323
1,2-Dichloropropane	ND		0.00686	1	10/31/2021 05:31	WG1766323
1,1-Dichloropropene	ND		0.00343	1	10/31/2021 05:31	WG1766323
1,3-Dichloropropane	ND		0.00686	1	10/31/2021 05:31	WG1766323
cis-1,3-Dichloropropene	ND		0.00343	1	10/31/2021 05:31	WG1766323
trans-1,3-Dichloropropene	ND		0.00686	1	10/31/2021 05:31	WG1766323
2,2-Dichloropropane	ND		0.00343	1	10/31/2021 05:31	WG1766323
Di-isopropyl ether	ND		0.00137	1	10/31/2021 05:31	WG1766323
Ethylbenzene	ND		0.00343	1	10/31/2021 05:31	WG1766323
Hexachloro-1,3-butadiene	ND		0.0343	1	10/31/2021 05:31	WG1766323
Isopropylbenzene	ND		0.00343	1	10/31/2021 05:31	WG1766323
p-Isopropyltoluene	ND		0.00686	1	10/31/2021 05:31	WG1766323
2-Butanone (MEK)	ND		0.137	1	10/31/2021 05:31	WG1766323
Methylene Chloride	ND		0.0343	1	10/31/2021 05:31	WG1766323
4-Methyl-2-pentanone (MIBK)	ND		0.0343	1	10/31/2021 05:31	WG1766323
Methyl tert-butyl ether	ND		0.00137	1	10/31/2021 05:31	WG1766323
Naphthalene	ND		0.0171	1	10/31/2021 05:31	WG1766323
n-Propylbenzene	ND		0.00686	1	10/31/2021 05:31	WG1766323
Styrene	ND		0.0171	1	10/31/2021 05:31	WG1766323
1,1,1,2-Tetrachloroethane	ND		0.00343	1	10/31/2021 05:31	WG1766323
1,1,2,2-Tetrachloroethane	ND		0.00343	1	10/31/2021 05:31	WG1766323
1,1,2-Trichlorotrifluoroethane	ND		0.00343	1	10/31/2021 05:31	WG1766323
Tetrachloroethene	ND		0.00343	1	10/31/2021 05:31	WG1766323
Toluene	ND		0.00686	1	10/31/2021 05:31	WG1766323
1,2,3-Trichlorobenzene	ND	C4	0.0171	1	10/31/2021 05:31	WG1766323
1,2,4-Trichlorobenzene	ND		0.0171	1	10/31/2021 05:31	WG1766323
1,1,1-Trichloroethane	ND		0.00343	1	10/31/2021 05:31	WG1766323
1,1,2-Trichloroethane	ND		0.00343	1	10/31/2021 05:31	WG1766323
Trichloroethene	ND		0.00137	1	10/31/2021 05:31	WG1766323
Trichlorofluoromethane	ND		0.00343	1	10/31/2021 05:31	WG1766323
1,2,3-Trichloropropane	ND		0.0171	1	10/31/2021 05:31	WG1766323
1,2,4-Trimethylbenzene	ND		0.00686	1	10/31/2021 05:31	WG1766323
1,2,3-Trimethylbenzene	ND		0.00686	1	10/31/2021 05:31	WG1766323
1,3,5-Trimethylbenzene	ND		0.00686	1	10/31/2021 05:31	WG1766323
Vinyl chloride	ND		0.00343	1	10/31/2021 05:31	WG1766323
Xylenes, Total	ND		0.00892	1	10/31/2021 05:31	WG1766323
(S) Toluene-d8	105		75.0-131		10/31/2021 05:31	WG1766323
(S) 4-Bromofluorobenzene	100		67.0-138		10/31/2021 05:31	WG1766323
(S) 1,2-Dichloroethane-d4	101		70.0-130		10/31/2021 05:31	WG1766323

- 1 Cp
- 2 Tc
- 3 Ss
- 4 Cn
- 5 Sr
- 6 Qc
- 7 Gl
- 8 Al
- 9 Sc

Semi-Volatile Organic Compounds (GC) by Method NWTPHDX-NO SGT

Analyte	Result (dry) mg/kg	Qualifier	RDL (dry) mg/kg	Dilution	Analysis date / time	Batch
Diesel Range Organics (DRO)	ND		4.69	1	11/03/2021 06:49	WG1766948
Residual Range Organics (RRO)	ND		11.7	1	11/03/2021 06:49	WG1766948
(S) o-Terphenyl	32.5		18.0-148		11/03/2021 06:49	WG1766948

Semi Volatile Organic Compounds (GC/MS) by Method 8270E-SIM

Analyte	Result (dry) mg/kg	Qualifier	RDL (dry) mg/kg	Dilution	Analysis date / time	Batch
Anthracene	ND		0.00703	1	10/30/2021 14:27	<a href="#">WG1765811</a>
Acenaphthene	ND		0.00703	1	10/30/2021 14:27	<a href="#">WG1765811</a>
Acenaphthylene	ND		0.00703	1	10/30/2021 14:27	<a href="#">WG1765811</a>
Benzo(a)anthracene	ND		0.00703	1	10/30/2021 14:27	<a href="#">WG1765811</a>
Benzo(a)pyrene	ND		0.00703	1	10/30/2021 14:27	<a href="#">WG1765811</a>
Benzo(b)fluoranthene	ND		0.00703	1	10/30/2021 14:27	<a href="#">WG1765811</a>
Benzo(g,h,i)perylene	ND		0.00703	1	10/30/2021 14:27	<a href="#">WG1765811</a>
Benzo(k)fluoranthene	ND		0.00703	1	10/30/2021 14:27	<a href="#">WG1765811</a>
Chrysene	ND		0.00703	1	10/30/2021 14:27	<a href="#">WG1765811</a>
Dibenz(a,h)anthracene	ND		0.00703	1	10/30/2021 14:27	<a href="#">WG1765811</a>
Fluoranthene	ND		0.00703	1	10/30/2021 14:27	<a href="#">WG1765811</a>
Fluorene	ND		0.00703	1	10/30/2021 14:27	<a href="#">WG1765811</a>
Indeno(1,2,3-cd)pyrene	ND		0.00703	1	10/30/2021 14:27	<a href="#">WG1765811</a>
Naphthalene	ND		0.0234	1	10/30/2021 14:27	<a href="#">WG1765811</a>
Phenanthrene	ND		0.00703	1	10/30/2021 14:27	<a href="#">WG1765811</a>
Pyrene	ND		0.00703	1	10/30/2021 14:27	<a href="#">WG1765811</a>
1-Methylnaphthalene	ND		0.0234	1	10/30/2021 14:27	<a href="#">WG1765811</a>
2-Methylnaphthalene	ND		0.0234	1	10/30/2021 14:27	<a href="#">WG1765811</a>
2-Chloronaphthalene	ND		0.0234	1	10/30/2021 14:27	<a href="#">WG1765811</a>
<i>(S)</i> Nitrobenzene-d5	85.6		14.0-149		10/30/2021 14:27	<a href="#">WG1765811</a>
<i>(S)</i> 2-Fluorobiphenyl	75.7		34.0-125		10/30/2021 14:27	<a href="#">WG1765811</a>
<i>(S)</i> p-Terphenyl-d14	89.7		23.0-120		10/30/2021 14:27	<a href="#">WG1765811</a>

1  
Cp

2  
Tc

3  
Ss

4  
Cn

5  
Sr

6  
Qc

7  
Gl

8  
Al

9  
Sc

Total Solids by Method 2540 G-2011

Analyte	Result	Qualifier	Dilution	Analysis	Batch
Total Solids	91.7		1	10/30/2021 16:39	<a href="#">WG1765837</a>

Mercury by Method 7471B

Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
Mercury	ND		0.0436	1	11/02/2021 15:25	<a href="#">WG1766498</a>

Metals (ICPMS) by Method 6020B

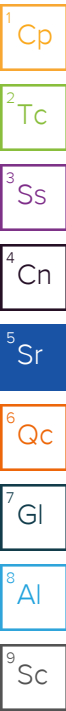
Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
Arsenic	2.58		1.09	5	11/03/2021 18:46	<a href="#">WG1765966</a>
Barium	39.6		2.73	5	11/03/2021 18:46	<a href="#">WG1765966</a>
Cadmium	ND		1.09	5	11/03/2021 18:46	<a href="#">WG1765966</a>
Chromium	21.1		5.45	5	11/03/2021 18:46	<a href="#">WG1765966</a>
Lead	ND		2.18	5	11/03/2021 18:46	<a href="#">WG1765966</a>
Selenium	ND		2.73	5	11/03/2021 18:46	<a href="#">WG1765966</a>
Silver	ND		0.545	5	11/03/2021 18:46	<a href="#">WG1765966</a>

Volatile Organic Compounds (GC) by Method NWTPHGX

Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
Gasoline Range Organics-NWTPH	ND		3.02	25	10/29/2021 05:27	<a href="#">WG1765548</a>
(S) a,a,a-Trifluorotoluene(FID)	94.5		77.0-120		10/29/2021 05:27	<a href="#">WG1765548</a>

Volatile Organic Compounds (GC/MS) by Method 8260D

Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
Acetone	ND		0.0604	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Acrylonitrile	ND	C3	0.0151	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Benzene	ND		0.00121	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Bromobenzene	ND		0.0151	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Bromodichloromethane	ND		0.00302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Bromoform	ND		0.0302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Bromomethane	ND		0.0151	1	10/31/2021 05:49	<a href="#">WG1766323</a>
n-Butylbenzene	ND		0.0151	1	10/31/2021 05:49	<a href="#">WG1766323</a>
sec-Butylbenzene	ND		0.0151	1	10/31/2021 05:49	<a href="#">WG1766323</a>
tert-Butylbenzene	ND		0.00604	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Carbon tetrachloride	ND		0.00604	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Chlorobenzene	ND		0.00302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Chlorodibromomethane	ND		0.00302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Chloroethane	ND		0.00604	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Chloroform	ND		0.00302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Chloromethane	ND		0.0151	1	10/31/2021 05:49	<a href="#">WG1766323</a>
2-Chlorotoluene	ND		0.00302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
4-Chlorotoluene	ND		0.00604	1	10/31/2021 05:49	<a href="#">WG1766323</a>
1,2-Dibromo-3-Chloropropane	ND		0.0302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
1,2-Dibromoethane	ND		0.00302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Dibromomethane	ND		0.00604	1	10/31/2021 05:49	<a href="#">WG1766323</a>
1,2-Dichlorobenzene	ND		0.00604	1	10/31/2021 05:49	<a href="#">WG1766323</a>
1,3-Dichlorobenzene	ND		0.00604	1	10/31/2021 05:49	<a href="#">WG1766323</a>
1,4-Dichlorobenzene	ND		0.00604	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Dichlorodifluoromethane	ND		0.00302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
1,1-Dichloroethane	ND		0.00302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
1,2-Dichloroethane	ND		0.00302	1	10/31/2021 05:49	<a href="#">WG1766323</a>



Volatile Organic Compounds (GC/MS) by Method 8260D

Analyte	Result (dry) mg/kg	Qualifier	RDL (dry) mg/kg	Dilution	Analysis date / time	Batch
1,1-Dichloroethene	ND		0.00302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
cis-1,2-Dichloroethene	ND		0.00302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
trans-1,2-Dichloroethene	ND		0.00604	1	10/31/2021 05:49	<a href="#">WG1766323</a>
1,2-Dichloropropane	ND		0.00604	1	10/31/2021 05:49	<a href="#">WG1766323</a>
1,1-Dichloropropene	ND		0.00302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
1,3-Dichloropropane	ND		0.00604	1	10/31/2021 05:49	<a href="#">WG1766323</a>
cis-1,3-Dichloropropene	ND		0.00302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
trans-1,3-Dichloropropene	ND		0.00604	1	10/31/2021 05:49	<a href="#">WG1766323</a>
2,2-Dichloropropane	ND		0.00302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Di-isopropyl ether	ND		0.00121	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Ethylbenzene	ND		0.00302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Hexachloro-1,3-butadiene	ND		0.0302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Isopropylbenzene	ND		0.00302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
p-Isopropyltoluene	ND		0.00604	1	10/31/2021 05:49	<a href="#">WG1766323</a>
2-Butanone (MEK)	ND		0.121	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Methylene Chloride	ND		0.0302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
4-Methyl-2-pentanone (MIBK)	ND		0.0302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Methyl tert-butyl ether	ND		0.00121	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Naphthalene	ND		0.0151	1	10/31/2021 05:49	<a href="#">WG1766323</a>
n-Propylbenzene	ND		0.00604	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Styrene	ND		0.0151	1	10/31/2021 05:49	<a href="#">WG1766323</a>
1,1,1,2-Tetrachloroethane	ND		0.00302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
1,1,2,2-Tetrachloroethane	ND		0.00302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
1,1,2-Trichlorotrifluoroethane	ND		0.00302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Tetrachloroethene	ND		0.00302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Toluene	ND		0.00604	1	10/31/2021 05:49	<a href="#">WG1766323</a>
1,2,3-Trichlorobenzene	ND	<a href="#">C4</a>	0.0151	1	10/31/2021 05:49	<a href="#">WG1766323</a>
1,2,4-Trichlorobenzene	ND		0.0151	1	10/31/2021 05:49	<a href="#">WG1766323</a>
1,1,1-Trichloroethane	ND		0.00302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
1,1,2-Trichloroethane	ND		0.00302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Trichloroethene	ND		0.00121	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Trichlorofluoromethane	ND		0.00302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
1,2,3-Trichloropropane	ND		0.0151	1	10/31/2021 05:49	<a href="#">WG1766323</a>
1,2,4-Trimethylbenzene	ND		0.00604	1	10/31/2021 05:49	<a href="#">WG1766323</a>
1,2,3-Trimethylbenzene	ND		0.00604	1	10/31/2021 05:49	<a href="#">WG1766323</a>
1,3,5-Trimethylbenzene	ND		0.00604	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Vinyl chloride	ND		0.00302	1	10/31/2021 05:49	<a href="#">WG1766323</a>
Xylenes, Total	ND		0.00785	1	10/31/2021 05:49	<a href="#">WG1766323</a>
(S) Toluene-d8	104		75.0-131		10/31/2021 05:49	<a href="#">WG1766323</a>
(S) 4-Bromofluorobenzene	100		67.0-138		10/31/2021 05:49	<a href="#">WG1766323</a>
(S) 1,2-Dichloroethane-d4	101		70.0-130		10/31/2021 05:49	<a href="#">WG1766323</a>

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc

Semi-Volatile Organic Compounds (GC) by Method NWTPHDX-NO SGT

Analyte	Result (dry) mg/kg	Qualifier	RDL (dry) mg/kg	Dilution	Analysis date / time	Batch
Diesel Range Organics (DRO)	ND		4.36	1	11/03/2021 07:02	<a href="#">WG1766948</a>
Residual Range Organics (RRO)	ND		10.9	1	11/03/2021 07:02	<a href="#">WG1766948</a>
(S) o-Terphenyl	31.1		18.0-148		11/03/2021 07:02	<a href="#">WG1766948</a>

Semi Volatile Organic Compounds (GC/MS) by Method 8270E-SIM

Analyte	Result (dry) mg/kg	Qualifier	RDL (dry) mg/kg	Dilution	Analysis date / time	Batch
Anthracene	ND		0.00654	1	10/30/2021 14:44	<a href="#">WG1765811</a>
Acenaphthene	ND		0.00654	1	10/30/2021 14:44	<a href="#">WG1765811</a>
Acenaphthylene	ND		0.00654	1	10/30/2021 14:44	<a href="#">WG1765811</a>
Benzo(a)anthracene	ND		0.00654	1	10/30/2021 14:44	<a href="#">WG1765811</a>
Benzo(a)pyrene	ND		0.00654	1	10/30/2021 14:44	<a href="#">WG1765811</a>
Benzo(b)fluoranthene	ND		0.00654	1	10/30/2021 14:44	<a href="#">WG1765811</a>
Benzo(g,h,i)perylene	ND		0.00654	1	10/30/2021 14:44	<a href="#">WG1765811</a>
Benzo(k)fluoranthene	ND		0.00654	1	10/30/2021 14:44	<a href="#">WG1765811</a>
Chrysene	ND		0.00654	1	10/30/2021 14:44	<a href="#">WG1765811</a>
Dibenz(a,h)anthracene	ND		0.00654	1	10/30/2021 14:44	<a href="#">WG1765811</a>
Fluoranthene	ND		0.00654	1	10/30/2021 14:44	<a href="#">WG1765811</a>
Fluorene	ND		0.00654	1	10/30/2021 14:44	<a href="#">WG1765811</a>
Indeno(1,2,3-cd)pyrene	ND		0.00654	1	10/30/2021 14:44	<a href="#">WG1765811</a>
Naphthalene	ND		0.0218	1	10/30/2021 14:44	<a href="#">WG1765811</a>
Phenanthrene	ND		0.00654	1	10/30/2021 14:44	<a href="#">WG1765811</a>
Pyrene	ND		0.00654	1	10/30/2021 14:44	<a href="#">WG1765811</a>
1-Methylnaphthalene	ND		0.0218	1	10/30/2021 14:44	<a href="#">WG1765811</a>
2-Methylnaphthalene	ND		0.0218	1	10/30/2021 14:44	<a href="#">WG1765811</a>
2-Chloronaphthalene	ND		0.0218	1	10/30/2021 14:44	<a href="#">WG1765811</a>
(S) Nitrobenzene-d5	86.5		14.0-149		10/30/2021 14:44	<a href="#">WG1765811</a>
(S) 2-Fluorobiphenyl	71.6		34.0-125		10/30/2021 14:44	<a href="#">WG1765811</a>
(S) p-Terphenyl-d14	97.4		23.0-120		10/30/2021 14:44	<a href="#">WG1765811</a>

1  
Cp

2  
Tc

3  
Ss

4  
Cn

5  
Sr

6  
Qc

7  
Gl

8  
Al

9  
Sc

Total Solids by Method 2540 G-2011

Analyte	Result	Qualifier	Dilution	Analysis	Batch
	%			date / time	
Total Solids	86.9		1	10/30/2021 16:39	<a href="#">WG1765837</a>

Mercury by Method 7471B

Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
	mg/kg		mg/kg		date / time	
Mercury	ND		0.0460	1	11/02/2021 15:27	<a href="#">WG1766498</a>

Metals (ICPMS) by Method 6020B

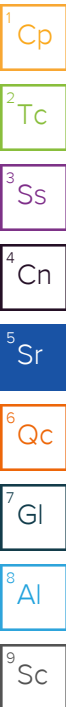
Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
	mg/kg		mg/kg		date / time	
Arsenic	7.77		1.15	5	11/03/2021 18:49	<a href="#">WG1765966</a>
Barium	58.8		2.88	5	11/03/2021 18:49	<a href="#">WG1765966</a>
Cadmium	ND		1.15	5	11/03/2021 18:49	<a href="#">WG1765966</a>
Chromium	41.6		5.75	5	11/03/2021 18:49	<a href="#">WG1765966</a>
Lead	3.16		2.30	5	11/03/2021 18:49	<a href="#">WG1765966</a>
Selenium	ND		2.88	5	11/03/2021 18:49	<a href="#">WG1765966</a>
Silver	ND		0.575	5	11/03/2021 18:49	<a href="#">WG1765966</a>

Volatile Organic Compounds (GC) by Method NWTPHGX

Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
	mg/kg		mg/kg		date / time	
Gasoline Range Organics-NWTPH	ND		3.35	25	10/29/2021 05:49	<a href="#">WG1765548</a>
(S) a,a,a-Trifluorotoluene(FID)	94.3		77.0-120		10/29/2021 05:49	<a href="#">WG1765548</a>

Volatile Organic Compounds (GC/MS) by Method 8260D

Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
	mg/kg		mg/kg		date / time	
Acetone	ND		0.0669	1	10/31/2021 06:08	<a href="#">WG1766323</a>
Acrylonitrile	ND	<a href="#">C3</a>	0.0167	1	10/31/2021 06:08	<a href="#">WG1766323</a>
Benzene	ND		0.00134	1	10/31/2021 06:08	<a href="#">WG1766323</a>
Bromobenzene	ND		0.0167	1	10/31/2021 06:08	<a href="#">WG1766323</a>
Bromodichloromethane	ND		0.00335	1	10/31/2021 06:08	<a href="#">WG1766323</a>
Bromoform	ND		0.0335	1	10/31/2021 06:08	<a href="#">WG1766323</a>
Bromomethane	ND		0.0167	1	10/31/2021 06:08	<a href="#">WG1766323</a>
n-Butylbenzene	ND		0.0167	1	10/31/2021 06:08	<a href="#">WG1766323</a>
sec-Butylbenzene	ND		0.0167	1	10/31/2021 06:08	<a href="#">WG1766323</a>
tert-Butylbenzene	ND		0.00669	1	10/31/2021 06:08	<a href="#">WG1766323</a>
Carbon tetrachloride	ND		0.00669	1	10/31/2021 06:08	<a href="#">WG1766323</a>
Chlorobenzene	ND		0.00335	1	10/31/2021 06:08	<a href="#">WG1766323</a>
Chlorodibromomethane	ND		0.00335	1	10/31/2021 06:08	<a href="#">WG1766323</a>
Chloroethane	ND		0.00669	1	10/31/2021 06:08	<a href="#">WG1766323</a>
Chloroform	ND		0.00335	1	10/31/2021 06:08	<a href="#">WG1766323</a>
Chloromethane	ND		0.0167	1	10/31/2021 06:08	<a href="#">WG1766323</a>
2-Chlorotoluene	ND		0.00335	1	10/31/2021 06:08	<a href="#">WG1766323</a>
4-Chlorotoluene	ND		0.00669	1	10/31/2021 06:08	<a href="#">WG1766323</a>
1,2-Dibromo-3-Chloropropane	ND		0.0335	1	10/31/2021 06:08	<a href="#">WG1766323</a>
1,2-Dibromoethane	ND		0.00335	1	10/31/2021 06:08	<a href="#">WG1766323</a>
Dibromomethane	ND		0.00669	1	10/31/2021 06:08	<a href="#">WG1766323</a>
1,2-Dichlorobenzene	ND		0.00669	1	10/31/2021 06:08	<a href="#">WG1766323</a>
1,3-Dichlorobenzene	ND		0.00669	1	10/31/2021 06:08	<a href="#">WG1766323</a>
1,4-Dichlorobenzene	ND		0.00669	1	10/31/2021 06:08	<a href="#">WG1766323</a>
Dichlorodifluoromethane	ND		0.00335	1	10/31/2021 06:08	<a href="#">WG1766323</a>
1,1-Dichloroethane	ND		0.00335	1	10/31/2021 06:08	<a href="#">WG1766323</a>
1,2-Dichloroethane	ND		0.00335	1	10/31/2021 06:08	<a href="#">WG1766323</a>



Volatile Organic Compounds (GC/MS) by Method 8260D

Analyte	Result (dry) mg/kg	Qualifier	RDL (dry) mg/kg	Dilution	Analysis date / time	Batch
1,1-Dichloroethene	ND		0.00335	1	10/31/2021 06:08	WG1766323
cis-1,2-Dichloroethene	ND		0.00335	1	10/31/2021 06:08	WG1766323
trans-1,2-Dichloroethene	ND		0.00669	1	10/31/2021 06:08	WG1766323
1,2-Dichloropropane	ND		0.00669	1	10/31/2021 06:08	WG1766323
1,1-Dichloropropene	ND		0.00335	1	10/31/2021 06:08	WG1766323
1,3-Dichloropropane	ND		0.00669	1	10/31/2021 06:08	WG1766323
cis-1,3-Dichloropropene	ND		0.00335	1	10/31/2021 06:08	WG1766323
trans-1,3-Dichloropropene	ND		0.00669	1	10/31/2021 06:08	WG1766323
2,2-Dichloropropane	ND		0.00335	1	10/31/2021 06:08	WG1766323
Di-isopropyl ether	ND		0.00134	1	10/31/2021 06:08	WG1766323
Ethylbenzene	ND		0.00335	1	10/31/2021 06:08	WG1766323
Hexachloro-1,3-butadiene	ND		0.0335	1	10/31/2021 06:08	WG1766323
Isopropylbenzene	ND		0.00335	1	10/31/2021 06:08	WG1766323
p-Isopropyltoluene	ND		0.00669	1	10/31/2021 06:08	WG1766323
2-Butanone (MEK)	ND		0.134	1	10/31/2021 06:08	WG1766323
Methylene Chloride	ND		0.0335	1	10/31/2021 06:08	WG1766323
4-Methyl-2-pentanone (MIBK)	ND		0.0335	1	10/31/2021 06:08	WG1766323
Methyl tert-butyl ether	ND		0.00134	1	10/31/2021 06:08	WG1766323
Naphthalene	ND		0.0167	1	10/31/2021 06:08	WG1766323
n-Propylbenzene	ND		0.00669	1	10/31/2021 06:08	WG1766323
Styrene	ND		0.0167	1	10/31/2021 06:08	WG1766323
1,1,1,2-Tetrachloroethane	ND		0.00335	1	10/31/2021 06:08	WG1766323
1,1,2,2-Tetrachloroethane	ND		0.00335	1	10/31/2021 06:08	WG1766323
1,1,2-Trichlorotrifluoroethane	ND		0.00335	1	10/31/2021 06:08	WG1766323
Tetrachloroethene	ND		0.00335	1	10/31/2021 06:08	WG1766323
Toluene	ND		0.00669	1	10/31/2021 06:08	WG1766323
1,2,3-Trichlorobenzene	ND	C4	0.0167	1	10/31/2021 06:08	WG1766323
1,2,4-Trichlorobenzene	ND		0.0167	1	10/31/2021 06:08	WG1766323
1,1,1-Trichloroethane	ND		0.00335	1	10/31/2021 06:08	WG1766323
1,1,2-Trichloroethane	ND		0.00335	1	10/31/2021 06:08	WG1766323
Trichloroethene	ND		0.00134	1	10/31/2021 06:08	WG1766323
Trichlorofluoromethane	ND		0.00335	1	10/31/2021 06:08	WG1766323
1,2,3-Trichloropropane	ND		0.0167	1	10/31/2021 06:08	WG1766323
1,2,4-Trimethylbenzene	ND		0.00669	1	10/31/2021 06:08	WG1766323
1,2,3-Trimethylbenzene	ND		0.00669	1	10/31/2021 06:08	WG1766323
1,3,5-Trimethylbenzene	ND		0.00669	1	10/31/2021 06:08	WG1766323
Vinyl chloride	ND		0.00335	1	10/31/2021 06:08	WG1766323
Xylenes, Total	ND		0.00870	1	10/31/2021 06:08	WG1766323
(S) Toluene-d8	102		75.0-131		10/31/2021 06:08	WG1766323
(S) 4-Bromofluorobenzene	101		67.0-138		10/31/2021 06:08	WG1766323
(S) 1,2-Dichloroethane-d4	107		70.0-130		10/31/2021 06:08	WG1766323

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc

Semi-Volatile Organic Compounds (GC) by Method NWTPHDX-NO SGT

Analyte	Result (dry) mg/kg	Qualifier	RDL (dry) mg/kg	Dilution	Analysis date / time	Batch
Diesel Range Organics (DRO)	5.55		4.60	1	11/03/2021 07:54	WG1766948
Residual Range Organics (RRO)	24.3		11.5	1	11/03/2021 07:54	WG1766948
(S) o-Terphenyl	61.6		18.0-148		11/03/2021 07:54	WG1766948

Semi Volatile Organic Compounds (GC/MS) by Method 8270E-SIM

Analyte	Result (dry) mg/kg	Qualifier	RDL (dry) mg/kg	Dilution	Analysis date / time	Batch
Anthracene	ND		0.00691	1	10/30/2021 15:02	<a href="#">WG1765811</a>
Acenaphthene	ND		0.00691	1	10/30/2021 15:02	<a href="#">WG1765811</a>
Acenaphthylene	ND		0.00691	1	10/30/2021 15:02	<a href="#">WG1765811</a>
Benzo(a)anthracene	ND		0.00691	1	10/30/2021 15:02	<a href="#">WG1765811</a>
Benzo(a)pyrene	ND		0.00691	1	10/30/2021 15:02	<a href="#">WG1765811</a>
Benzo(b)fluoranthene	ND		0.00691	1	10/30/2021 15:02	<a href="#">WG1765811</a>
Benzo(g,h,i)perylene	ND		0.00691	1	10/30/2021 15:02	<a href="#">WG1765811</a>
Benzo(k)fluoranthene	ND		0.00691	1	10/30/2021 15:02	<a href="#">WG1765811</a>
Chrysene	ND		0.00691	1	10/30/2021 15:02	<a href="#">WG1765811</a>
Dibenz(a,h)anthracene	ND		0.00691	1	10/30/2021 15:02	<a href="#">WG1765811</a>
Fluoranthene	ND		0.00691	1	10/30/2021 15:02	<a href="#">WG1765811</a>
Fluorene	ND		0.00691	1	10/30/2021 15:02	<a href="#">WG1765811</a>
Indeno(1,2,3-cd)pyrene	ND		0.00691	1	10/30/2021 15:02	<a href="#">WG1765811</a>
Naphthalene	ND		0.0230	1	10/30/2021 15:02	<a href="#">WG1765811</a>
Phenanthrene	ND		0.00691	1	10/30/2021 15:02	<a href="#">WG1765811</a>
Pyrene	ND		0.00691	1	10/30/2021 15:02	<a href="#">WG1765811</a>
1-Methylnaphthalene	ND		0.0230	1	10/30/2021 15:02	<a href="#">WG1765811</a>
2-Methylnaphthalene	ND		0.0230	1	10/30/2021 15:02	<a href="#">WG1765811</a>
2-Chloronaphthalene	ND		0.0230	1	10/30/2021 15:02	<a href="#">WG1765811</a>
(S) Nitrobenzene-d5	77.8		14.0-149		10/30/2021 15:02	<a href="#">WG1765811</a>
(S) 2-Fluorobiphenyl	62.6		34.0-125		10/30/2021 15:02	<a href="#">WG1765811</a>
(S) p-Terphenyl-d14	80.0		23.0-120		10/30/2021 15:02	<a href="#">WG1765811</a>

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc

Total Solids by Method 2540 G-2011

Analyte	Result	Qualifier	Dilution	Analysis	Batch
	%			date / time	
Total Solids	92.4		1	10/30/2021 16:39	<a href="#">WG1765837</a>

Mercury by Method 7471B

Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
	mg/kg		mg/kg		date / time	
Mercury	ND		0.0433	1	11/02/2021 15:30	<a href="#">WG1766498</a>

Metals (ICPMS) by Method 6020B

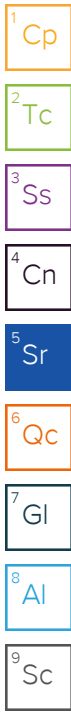
Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
	mg/kg		mg/kg		date / time	
Arsenic	2.31		1.08	5	11/03/2021 18:52	<a href="#">WG1765966</a>
Barium	35.1		2.71	5	11/03/2021 18:52	<a href="#">WG1765966</a>
Cadmium	ND		1.08	5	11/03/2021 18:52	<a href="#">WG1765966</a>
Chromium	16.3		5.41	5	11/03/2021 18:52	<a href="#">WG1765966</a>
Lead	ND		2.16	5	11/03/2021 18:52	<a href="#">WG1765966</a>
Selenium	ND		2.71	5	11/03/2021 18:52	<a href="#">WG1765966</a>
Silver	ND		0.541	5	11/03/2021 18:52	<a href="#">WG1765966</a>

Volatile Organic Compounds (GC) by Method NWTPHGX

Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
	mg/kg		mg/kg		date / time	
Gasoline Range Organics-NWTPH	ND		2.95	25	10/29/2021 06:11	<a href="#">WG1765548</a>
(S) a,a,a-Trifluorotoluene(FID)	94.3		77.0-120		10/29/2021 06:11	<a href="#">WG1765548</a>

Volatile Organic Compounds (GC/MS) by Method 8260D

Analyte	Result (dry)	Qualifier	RDL (dry)	Dilution	Analysis	Batch
	mg/kg		mg/kg		date / time	
Acetone	ND		0.0590	1	10/31/2021 06:27	<a href="#">WG1766323</a>
Acrylonitrile	ND	<a href="#">C3</a>	0.0147	1	10/31/2021 06:27	<a href="#">WG1766323</a>
Benzene	ND		0.00118	1	10/31/2021 06:27	<a href="#">WG1766323</a>
Bromobenzene	ND		0.0147	1	10/31/2021 06:27	<a href="#">WG1766323</a>
Bromodichloromethane	ND		0.00295	1	10/31/2021 06:27	<a href="#">WG1766323</a>
Bromoform	ND		0.0295	1	10/31/2021 06:27	<a href="#">WG1766323</a>
Bromomethane	ND		0.0147	1	10/31/2021 06:27	<a href="#">WG1766323</a>
n-Butylbenzene	ND		0.0147	1	10/31/2021 06:27	<a href="#">WG1766323</a>
sec-Butylbenzene	ND		0.0147	1	10/31/2021 06:27	<a href="#">WG1766323</a>
tert-Butylbenzene	ND		0.00590	1	10/31/2021 06:27	<a href="#">WG1766323</a>
Carbon tetrachloride	ND		0.00590	1	10/31/2021 06:27	<a href="#">WG1766323</a>
Chlorobenzene	ND		0.00295	1	10/31/2021 06:27	<a href="#">WG1766323</a>
Chlorodibromomethane	ND		0.00295	1	10/31/2021 06:27	<a href="#">WG1766323</a>
Chloroethane	ND		0.00590	1	10/31/2021 06:27	<a href="#">WG1766323</a>
Chloroform	ND		0.00295	1	10/31/2021 06:27	<a href="#">WG1766323</a>
Chloromethane	ND		0.0147	1	10/31/2021 06:27	<a href="#">WG1766323</a>
2-Chlorotoluene	ND		0.00295	1	10/31/2021 06:27	<a href="#">WG1766323</a>
4-Chlorotoluene	ND		0.00590	1	10/31/2021 06:27	<a href="#">WG1766323</a>
1,2-Dibromo-3-Chloropropane	ND		0.0295	1	10/31/2021 06:27	<a href="#">WG1766323</a>
1,2-Dibromoethane	ND		0.00295	1	10/31/2021 06:27	<a href="#">WG1766323</a>
Dibromomethane	ND		0.00590	1	10/31/2021 06:27	<a href="#">WG1766323</a>
1,2-Dichlorobenzene	ND		0.00590	1	10/31/2021 06:27	<a href="#">WG1766323</a>
1,3-Dichlorobenzene	ND		0.00590	1	10/31/2021 06:27	<a href="#">WG1766323</a>
1,4-Dichlorobenzene	ND		0.00590	1	10/31/2021 06:27	<a href="#">WG1766323</a>
Dichlorodifluoromethane	ND		0.00295	1	10/31/2021 06:27	<a href="#">WG1766323</a>
1,1-Dichloroethane	ND		0.00295	1	10/31/2021 06:27	<a href="#">WG1766323</a>
1,2-Dichloroethane	ND		0.00295	1	10/31/2021 06:27	<a href="#">WG1766323</a>



Volatile Organic Compounds (GC/MS) by Method 8260D

Analyte	Result (dry) mg/kg	Qualifier	RDL (dry) mg/kg	Dilution	Analysis date / time	Batch
1,1-Dichloroethene	ND		0.00295	1	10/31/2021 06:27	WG1766323
cis-1,2-Dichloroethene	ND		0.00295	1	10/31/2021 06:27	WG1766323
trans-1,2-Dichloroethene	ND		0.00590	1	10/31/2021 06:27	WG1766323
1,2-Dichloropropane	ND		0.00590	1	10/31/2021 06:27	WG1766323
1,1-Dichloropropene	ND		0.00295	1	10/31/2021 06:27	WG1766323
1,3-Dichloropropane	ND		0.00590	1	10/31/2021 06:27	WG1766323
cis-1,3-Dichloropropene	ND		0.00295	1	10/31/2021 06:27	WG1766323
trans-1,3-Dichloropropene	ND		0.00590	1	10/31/2021 06:27	WG1766323
2,2-Dichloropropane	ND		0.00295	1	10/31/2021 06:27	WG1766323
Di-isopropyl ether	ND		0.00118	1	10/31/2021 06:27	WG1766323
Ethylbenzene	ND		0.00295	1	10/31/2021 06:27	WG1766323
Hexachloro-1,3-butadiene	ND		0.0295	1	10/31/2021 06:27	WG1766323
Isopropylbenzene	ND		0.00295	1	10/31/2021 06:27	WG1766323
p-Isopropyltoluene	ND		0.00590	1	10/31/2021 06:27	WG1766323
2-Butanone (MEK)	ND		0.118	1	10/31/2021 06:27	WG1766323
Methylene Chloride	ND		0.0295	1	10/31/2021 06:27	WG1766323
4-Methyl-2-pentanone (MIBK)	ND		0.0295	1	10/31/2021 06:27	WG1766323
Methyl tert-butyl ether	ND		0.00118	1	10/31/2021 06:27	WG1766323
Naphthalene	ND		0.0147	1	10/31/2021 06:27	WG1766323
n-Propylbenzene	ND		0.00590	1	10/31/2021 06:27	WG1766323
Styrene	ND		0.0147	1	10/31/2021 06:27	WG1766323
1,1,1,2-Tetrachloroethane	ND		0.00295	1	10/31/2021 06:27	WG1766323
1,1,2,2-Tetrachloroethane	ND		0.00295	1	10/31/2021 06:27	WG1766323
1,1,2-Trichlorotrifluoroethane	ND		0.00295	1	10/31/2021 06:27	WG1766323
Tetrachloroethene	ND		0.00295	1	10/31/2021 06:27	WG1766323
Toluene	ND		0.00590	1	10/31/2021 06:27	WG1766323
1,2,3-Trichlorobenzene	ND	C4	0.0147	1	10/31/2021 06:27	WG1766323
1,2,4-Trichlorobenzene	ND		0.0147	1	10/31/2021 06:27	WG1766323
1,1,1-Trichloroethane	ND		0.00295	1	10/31/2021 06:27	WG1766323
1,1,2-Trichloroethane	ND		0.00295	1	10/31/2021 06:27	WG1766323
Trichloroethene	ND		0.00118	1	10/31/2021 06:27	WG1766323
Trichlorofluoromethane	ND		0.00295	1	10/31/2021 06:27	WG1766323
1,2,3-Trichloropropane	ND		0.0147	1	10/31/2021 06:27	WG1766323
1,2,4-Trimethylbenzene	ND		0.00590	1	10/31/2021 06:27	WG1766323
1,2,3-Trimethylbenzene	ND		0.00590	1	10/31/2021 06:27	WG1766323
1,3,5-Trimethylbenzene	ND		0.00590	1	10/31/2021 06:27	WG1766323
Vinyl chloride	ND		0.00295	1	10/31/2021 06:27	WG1766323
Xylenes, Total	ND		0.00767	1	10/31/2021 06:27	WG1766323
(S) Toluene-d8	103		75.0-131		10/31/2021 06:27	WG1766323
(S) 4-Bromofluorobenzene	96.5		67.0-138		10/31/2021 06:27	WG1766323
(S) 1,2-Dichloroethane-d4	101		70.0-130		10/31/2021 06:27	WG1766323

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc

Semi-Volatile Organic Compounds (GC) by Method NWTPHDX-NO SGT

Analyte	Result (dry) mg/kg	Qualifier	RDL (dry) mg/kg	Dilution	Analysis date / time	Batch
Diesel Range Organics (DRO)	ND		4.33	1	11/03/2021 07:41	WG1766948
Residual Range Organics (RRO)	ND		10.8	1	11/03/2021 07:41	WG1766948
(S) o-Terphenyl	48.8		18.0-148		11/03/2021 07:41	WG1766948

Semi Volatile Organic Compounds (GC/MS) by Method 8270E-SIM

Analyte	Result (dry) mg/kg	Qualifier	RDL (dry) mg/kg	Dilution	Analysis date / time	Batch
Anthracene	ND		0.00649	1	10/30/2021 15:19	<a href="#">WG1765811</a>
Acenaphthene	ND		0.00649	1	10/30/2021 15:19	<a href="#">WG1765811</a>
Acenaphthylene	ND		0.00649	1	10/30/2021 15:19	<a href="#">WG1765811</a>
Benzo(a)anthracene	ND		0.00649	1	10/30/2021 15:19	<a href="#">WG1765811</a>
Benzo(a)pyrene	ND		0.00649	1	10/30/2021 15:19	<a href="#">WG1765811</a>
Benzo(b)fluoranthene	ND		0.00649	1	10/30/2021 15:19	<a href="#">WG1765811</a>
Benzo(g,h,i)perylene	ND		0.00649	1	10/30/2021 15:19	<a href="#">WG1765811</a>
Benzo(k)fluoranthene	ND		0.00649	1	10/30/2021 15:19	<a href="#">WG1765811</a>
Chrysene	ND		0.00649	1	10/30/2021 15:19	<a href="#">WG1765811</a>
Dibenz(a,h)anthracene	ND		0.00649	1	10/30/2021 15:19	<a href="#">WG1765811</a>
Fluoranthene	ND		0.00649	1	10/30/2021 15:19	<a href="#">WG1765811</a>
Fluorene	ND		0.00649	1	10/30/2021 15:19	<a href="#">WG1765811</a>
Indeno(1,2,3-cd)pyrene	ND		0.00649	1	10/30/2021 15:19	<a href="#">WG1765811</a>
Naphthalene	ND		0.0216	1	10/30/2021 15:19	<a href="#">WG1765811</a>
Phenanthrene	ND		0.00649	1	10/30/2021 15:19	<a href="#">WG1765811</a>
Pyrene	ND		0.00649	1	10/30/2021 15:19	<a href="#">WG1765811</a>
1-Methylnaphthalene	ND		0.0216	1	10/30/2021 15:19	<a href="#">WG1765811</a>
2-Methylnaphthalene	ND		0.0216	1	10/30/2021 15:19	<a href="#">WG1765811</a>
2-Chloronaphthalene	ND		0.0216	1	10/30/2021 15:19	<a href="#">WG1765811</a>
(S) Nitrobenzene-d5	99.3		14.0-149		10/30/2021 15:19	<a href="#">WG1765811</a>
(S) 2-Fluorobiphenyl	86.3		34.0-125		10/30/2021 15:19	<a href="#">WG1765811</a>
(S) p-Terphenyl-d14	121	<a href="#">J1</a>	23.0-120		10/30/2021 15:19	<a href="#">WG1765811</a>

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc

Method Blank (MB)

(MB) R3723681-1 10/30/21 16:39

Analyte	MB Result	MB Qualifier	MB MDL	MB RDL
	%		%	%
Total Solids	0.00100			

<sup>1</sup>Cp

<sup>2</sup>Tc

<sup>3</sup>Ss

L1423266-05 Original Sample (OS) • Duplicate (DUP)

(OS) L1423266-05 10/30/21 16:39 • (DUP) R3723681-3 10/30/21 16:39

Analyte	Original Result	DUP Result	Dilution	DUP RPD	DUP Qualifier	DUP RPD Limits
	%	%		%		%
Total Solids	89.0	89.2	1	0.237		10

<sup>4</sup>Cn

<sup>5</sup>Sr

Laboratory Control Sample (LCS)

(LCS) R3723681-2 10/30/21 16:39

Analyte	Spike Amount	LCS Result	LCS Rec.	Rec. Limits	LCS Qualifier
	%	%	%	%	
Total Solids	50.0	50.0	100	85.0-115	

<sup>6</sup>Qc

<sup>7</sup>Gl

<sup>8</sup>Al

<sup>9</sup>Sc

Method Blank (MB)

(MB) R3724558-1 11/02/21 14:52

Analyte	MB Result	MB Qualifier	MB MDL	MB RDL
Mercury	U		0.0180	0.0400

1 Cp

2 Tc

3 Ss

Laboratory Control Sample (LCS)

(LCS) R3724558-2 11/02/21 14:54

Analyte	Spike Amount	LCS Result	LCS Rec.	Rec. Limits	LCS Qualifier
Mercury	0.500	0.509	102	80.0-120	

4 Cn

5 Sr

L1423622-05 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1423622-05 11/02/21 14:57 • (MS) R3724558-3 11/02/21 14:59 • (MSD) R3724558-4 11/02/21 15:02

Analyte	Spike Amount (dry)	Original Result (dry)	MS Result (dry)	MSD Result (dry)	MS Rec.	MSD Rec.	Dilution	Rec. Limits	MS Qualifier	MSD Qualifier	RPD	RPD Limits
Mercury	0.669	ND	0.790	0.736	111	103	1	75.0-125			7.13	20

6 Qc

7 Gl

8 Al

9 Sc

Method Blank (MB)

(MB) R3725162-1 11/03/21 17:52

Analyte	MB Result	MB Qualifier	MB MDL	MB RDL
	mg/kg		mg/kg	mg/kg
Arsenic	U		0.100	1.00
Barium	U		0.152	2.50
Cadmium	U		0.0855	1.00
Chromium	U		0.297	5.00
Lead	U		0.0990	2.00
Selenium	0.208	J	0.180	2.50
Silver	U		0.0865	0.500

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc

Laboratory Control Sample (LCS)

(LCS) R3725162-2 11/03/21 17:55

Analyte	Spike Amount	LCS Result	LCS Rec.	Rec. Limits	LCS Qualifier
	mg/kg	mg/kg	%	%	
Arsenic	100	91.7	91.7	80.0-120	
Barium	100	90.8	90.8	80.0-120	
Cadmium	100	101	101	80.0-120	
Chromium	100	96.7	96.7	80.0-120	
Lead	100	93.1	93.1	80.0-120	
Selenium	100	98.2	98.2	80.0-120	
Silver	20.0	19.7	98.4	80.0-120	

L1423646-29 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1423646-29 11/03/21 17:59 • (MS) R3725162-5 11/03/21 18:09 • (MSD) R3725162-6 11/03/21 18:12

Analyte	Spike Amount	Original Result	MS Result	MSD Result	MS Rec.	MSD Rec.	Dilution	Rec. Limits	MS Qualifier	MSD Qualifier	RPD	RPD Limits
	mg/kg	mg/kg	mg/kg	mg/kg	%	%		%			%	%
Arsenic	100	29.6	112	116	82.6	85.9	5	75.0-125			2.90	20
Barium	100	382	476	502	94.4	120	5	75.0-125			5.33	20
Cadmium	100	ND	91.5	98.6	90.9	98.0	5	75.0-125			7.48	20
Chromium	100	26.3	113	115	86.4	88.8	5	75.0-125			2.09	20
Lead	100	19.4	106	110	86.2	91.0	5	75.0-125			4.46	20
Selenium	100	ND	91.3	98.2	90.2	97.1	5	75.0-125			7.27	20
Silver	20.0	ND	17.6	18.6	88.2	93.0	5	75.0-125			5.36	20

Method Blank (MB)

(MB) R3725739-3 10/29/21 04:36

Analyte	MB Result	MB Qualifier	MB MDL	MB RDL
TPHG C6 - C12	U		0.848	2.50
(S) a,a,a-Trifluorotoluene(FID)	117			77.0-120

1 Cp

2 Tc

3 Ss

4 Cn

Laboratory Control Sample (LCS)

(LCS) R3725739-2 10/29/21 03:47

Analyte	Spike Amount	LCS Result	LCS Rec.	Rec. Limits	LCS Qualifier
TPHG C6 - C12	5.50	6.17	112	71.0-124	
(S) a,a,a-Trifluorotoluene(FID)			104	77.0-120	

5 Sr

6 Qc

7 Gl

L1423058-01 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1423058-01 10/29/21 08:23 • (MS) R3725739-6 10/29/21 13:48 • (MSD) R3725739-7 10/29/21 14:10

Analyte	Spike Amount (dry)	Original Result (dry)	MS Result (dry)	MSD Result (dry)	MS Rec.	MSD Rec.	Dilution	Rec. Limits	MS Qualifier	MSD Qualifier	RPD	RPD Limits
Gasoline Range Organics-NWTPH	162	ND	199	188	123	117	25	50.0-150			5.36	27
(S) a,a,a-Trifluorotoluene(FID)					110	109		77.0-120				

8 Al

9 Sc

Method Blank (MB)

(MB) R3723041-2 10/29/21 03:49

Analyte	MB Result mg/kg	MB Qualifier	MB MDL mg/kg	MB RDL mg/kg
Gasoline Range Organics-NWTPH	U		0.848	2.50
(S) a,a,a-Trifluorotoluene(FID)	93.8			77.0-120

Laboratory Control Sample (LCS)

(LCS) R3723041-1 10/29/21 03:05

Analyte	Spike Amount mg/kg	LCS Result mg/kg	LCS Rec. %	Rec. Limits %	LCS Qualifier
Gasoline Range Organics-NWTPH	5.50	5.75	105	71.0-124	
(S) a,a,a-Trifluorotoluene(FID)			100	77.0-120	

- 1 Cp
- 2 Tc
- 3 Ss
- 4 Cn
- 5 Sr
- 6 Qc
- 7 Gl
- 8 Al
- 9 Sc

Method Blank (MB)

(MB) R3725864-2 10/31/21 00:10

Analyte	MB Result mg/kg	MB Qualifier	MB MDL mg/kg	MB RDL mg/kg
Acetone	U		0.0365	0.0500
Acrylonitrile	U		0.00361	0.0125
Benzene	U		0.000467	0.00100
Bromobenzene	U		0.000900	0.0125
Bromodichloromethane	U		0.000725	0.00250
Bromoform	U		0.00117	0.0250
Bromomethane	U		0.00197	0.0125
n-Butylbenzene	U		0.00525	0.0125
sec-Butylbenzene	U		0.00288	0.0125
tert-Butylbenzene	U		0.00195	0.00500
Carbon tetrachloride	U		0.000898	0.00500
Chlorobenzene	U		0.000210	0.00250
Chlorodibromomethane	U		0.000612	0.00250
Chloroethane	U		0.00170	0.00500
Chloroform	U		0.00103	0.00250
Chloromethane	U		0.00435	0.0125
2-Chlorotoluene	U		0.000865	0.00250
4-Chlorotoluene	U		0.000450	0.00500
1,2-Dibromo-3-Chloropropane	U		0.00390	0.0250
1,2-Dibromoethane	U		0.000648	0.00250
Dibromomethane	U		0.000750	0.00500
1,2-Dichlorobenzene	U		0.000425	0.00500
1,3-Dichlorobenzene	U		0.000600	0.00500
1,4-Dichlorobenzene	U		0.000700	0.00500
Dichlorodifluoromethane	U		0.00161	0.00250
1,1-Dichloroethane	U		0.000491	0.00250
1,2-Dichloroethane	U		0.000649	0.00250
1,1-Dichloroethene	U		0.000606	0.00250
cis-1,2-Dichloroethene	U		0.000734	0.00250
trans-1,2-Dichloroethene	U		0.00104	0.00500
1,2-Dichloropropane	U		0.00142	0.00500
1,1-Dichloropropene	U		0.000809	0.00250
1,3-Dichloropropane	U		0.000501	0.00500
cis-1,3-Dichloropropene	U		0.000757	0.00250
trans-1,3-Dichloropropene	U		0.00114	0.00500
2,2-Dichloropropane	U		0.00138	0.00250
Di-isopropyl ether	U		0.000410	0.00100
Ethylbenzene	U		0.000737	0.00250
Hexachloro-1,3-butadiene	U		0.00600	0.0250
Isopropylbenzene	U		0.000425	0.00250

<sup>1</sup>Cp

<sup>2</sup>Tc

<sup>3</sup>Ss

<sup>4</sup>Cn

<sup>5</sup>Sr

<sup>6</sup>Qc

<sup>7</sup>Gl

<sup>8</sup>Al

<sup>9</sup>Sc

Method Blank (MB)

(MB) R3725864-2 10/31/21 00:10

Analyte	MB Result mg/kg	MB Qualifier	MB MDL mg/kg	MB RDL mg/kg
p-Isopropyltoluene	U		0.00255	0.00500
2-Butanone (MEK)	0.0821	U	0.0635	0.100
Methylene Chloride	U		0.00664	0.0250
4-Methyl-2-pentanone (MIBK)	U		0.00228	0.0250
Methyl tert-butyl ether	U		0.000350	0.00100
Naphthalene	U		0.00488	0.0125
n-Propylbenzene	U		0.000950	0.00500
Styrene	U		0.000229	0.0125
1,1,1,2-Tetrachloroethane	U		0.000948	0.00250
1,1,2,2-Tetrachloroethane	U		0.000695	0.00250
Tetrachloroethene	U		0.000896	0.00250
Toluene	U		0.00130	0.00500
1,1,2-Trichlorotrifluoroethane	U		0.000754	0.00250
1,2,3-Trichlorobenzene	U		0.00733	0.0125
1,2,4-Trichlorobenzene	U		0.00440	0.0125
1,1,1-Trichloroethane	U		0.000923	0.00250
1,1,2-Trichloroethane	U		0.000597	0.00250
Trichloroethene	U		0.000584	0.00100
Trichlorofluoromethane	U		0.000827	0.00250
1,2,3-Trichloropropane	U		0.00162	0.0125
1,2,3-Trimethylbenzene	U		0.00158	0.00500
1,2,4-Trimethylbenzene	U		0.00158	0.00500
1,3,5-Trimethylbenzene	U		0.00200	0.00500
Vinyl chloride	U		0.00116	0.00250
Xylenes, Total	U		0.000880	0.00650
(S) Toluene-d8	103			75.0-131
(S) 4-Bromofluorobenzene	100			67.0-138
(S) 1,2-Dichloroethane-d4	107			70.0-130

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc

Laboratory Control Sample (LCS)

(LCS) R3725864-1 10/30/21 23:13

Analyte	Spike Amount mg/kg	LCS Result mg/kg	LCS Rec. %	Rec. Limits %	LCS Qualifier
Acetone	0.625	0.554	88.6	10.0-160	
Acrylonitrile	0.625	0.518	82.9	45.0-153	
Benzene	0.125	0.112	89.6	70.0-123	
Bromobenzene	0.125	0.125	100	73.0-121	
Bromodichloromethane	0.125	0.115	92.0	73.0-121	

Laboratory Control Sample (LCS)

(LCS) R3725864-1 10/30/21 23:13

Analyte	Spike Amount mg/kg	LCS Result mg/kg	LCS Rec. %	Rec. Limits %	<u>LCS Qualifier</u>
Bromoform	0.125	0.112	89.6	64.0-132	
Bromomethane	0.125	0.114	91.2	56.0-147	
n-Butylbenzene	0.125	0.109	87.2	68.0-135	
sec-Butylbenzene	0.125	0.108	86.4	74.0-130	
tert-Butylbenzene	0.125	0.111	88.8	75.0-127	
Carbon tetrachloride	0.125	0.113	90.4	66.0-128	
Chlorobenzene	0.125	0.115	92.0	76.0-128	
Chlorodibromomethane	0.125	0.110	88.0	74.0-127	
Chloroethane	0.125	0.121	96.8	61.0-134	
Chloroform	0.125	0.114	91.2	72.0-123	
Chloromethane	0.125	0.133	106	51.0-138	
2-Chlorotoluene	0.125	0.117	93.6	75.0-124	
4-Chlorotoluene	0.125	0.114	91.2	75.0-124	
1,2-Dibromo-3-Chloropropane	0.125	0.104	83.2	59.0-130	
1,2-Dibromoethane	0.125	0.117	93.6	74.0-128	
Dibromomethane	0.125	0.125	100	75.0-122	
1,2-Dichlorobenzene	0.125	0.115	92.0	76.0-124	
1,3-Dichlorobenzene	0.125	0.114	91.2	76.0-125	
1,4-Dichlorobenzene	0.125	0.112	89.6	77.0-121	
Dichlorodifluoromethane	0.125	0.107	85.6	43.0-156	
1,1-Dichloroethane	0.125	0.115	92.0	70.0-127	
1,2-Dichloroethane	0.125	0.128	102	65.0-131	
1,1-Dichloroethene	0.125	0.110	88.0	65.0-131	
cis-1,2-Dichloroethene	0.125	0.113	90.4	73.0-125	
trans-1,2-Dichloroethene	0.125	0.105	84.0	71.0-125	
1,2-Dichloropropane	0.125	0.117	93.6	74.0-125	
1,1-Dichloropropene	0.125	0.120	96.0	73.0-125	
1,3-Dichloropropane	0.125	0.120	96.0	80.0-125	
cis-1,3-Dichloropropene	0.125	0.121	96.8	76.0-127	
trans-1,3-Dichloropropene	0.125	0.124	99.2	73.0-127	
2,2-Dichloropropane	0.125	0.0921	73.7	59.0-135	
Di-isopropyl ether	0.125	0.116	92.8	60.0-136	
Ethylbenzene	0.125	0.107	85.6	74.0-126	
Hexachloro-1,3-butadiene	0.125	0.145	116	57.0-150	
Isopropylbenzene	0.125	0.106	84.8	72.0-127	
p-Isopropyltoluene	0.125	0.101	80.8	72.0-133	
2-Butanone (MEK)	0.625	0.531	85.0	30.0-160	
Methylene Chloride	0.125	0.111	88.8	68.0-123	
4-Methyl-2-pentanone (MIBK)	0.625	0.603	96.5	56.0-143	
Methyl tert-butyl ether	0.125	0.111	88.8	66.0-132	

<sup>1</sup>Cp

<sup>2</sup>Tc

<sup>3</sup>Ss

<sup>4</sup>Cn

<sup>5</sup>Sr

<sup>6</sup>Qc

<sup>7</sup>Gl

<sup>8</sup>Al

<sup>9</sup>Sc

Laboratory Control Sample (LCS)

(LCS) R3725864-1 10/30/21 23:13

Analyte	Spike Amount mg/kg	LCS Result mg/kg	LCS Rec. %	Rec. Limits %	<u>LCS Qualifier</u>
Naphthalene	0.125	0.116	92.8	59.0-130	
n-Propylbenzene	0.125	0.111	88.8	74.0-126	
Styrene	0.125	0.107	85.6	72.0-127	
1,1,1,2-Tetrachloroethane	0.125	0.104	83.2	74.0-129	
1,1,2,2-Tetrachloroethane	0.125	0.109	87.2	68.0-128	
Tetrachloroethene	0.125	0.123	98.4	70.0-136	
Toluene	0.125	0.113	90.4	75.0-121	
1,1,2-Trichlorotrifluoroethane	0.125	0.122	97.6	61.0-139	
1,2,3-Trichlorobenzene	0.125	0.126	101	59.0-139	
1,2,4-Trichlorobenzene	0.125	0.128	102	62.0-137	
1,1,1-Trichloroethane	0.125	0.107	85.6	69.0-126	
1,1,2-Trichloroethane	0.125	0.120	96.0	78.0-123	
Trichloroethene	0.125	0.126	101	76.0-126	
Trichlorofluoromethane	0.125	0.113	90.4	61.0-142	
1,2,3-Trichloropropane	0.125	0.113	90.4	67.0-129	
1,2,3-Trimethylbenzene	0.125	0.105	84.0	74.0-124	
1,2,4-Trimethylbenzene	0.125	0.103	82.4	70.0-126	
1,3,5-Trimethylbenzene	0.125	0.106	84.8	73.0-127	
Vinyl chloride	0.125	0.125	100	63.0-134	
Xylenes, Total	0.375	0.325	86.7	72.0-127	
<i>(S) Toluene-d8</i>			103	75.0-131	
<i>(S) 4-Bromofluorobenzene</i>			101	67.0-138	
<i>(S) 1,2-Dichloroethane-d4</i>			108	70.0-130	

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc

Method Blank (MB)

(MB) R3724440-1 11/02/21 07:41

Analyte	MB Result mg/kg	MB Qualifier	MB MDL mg/kg	MB RDL mg/kg
Diesel Range Organics (DRO)	U		1.33	4.00
Residual Range Organics (RRO)	U		3.33	10.0
<i>(S) o-Terphenyl</i>	53.8			18.0-148

Laboratory Control Sample (LCS)

(LCS) R3724440-2 11/02/21 07:54

Analyte	Spike Amount mg/kg	LCS Result mg/kg	LCS Rec. %	Rec. Limits %	LCS Qualifier
Diesel Range Organics (DRO)	50.0	32.2	64.4	50.0-150	
<i>(S) o-Terphenyl</i>			70.6	18.0-148	

L1423415-01 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1423415-01 11/02/21 16:57 • (MS) R3724440-3 11/02/21 17:10 • (MSD) R3724440-4 11/02/21 17:22

Analyte	Spike Amount mg/kg	Original Result mg/kg	MS Result mg/kg	MSD Result mg/kg	MS Rec. %	MSD Rec. %	Dilution	Rec. Limits %	MS Qualifier	MSD Qualifier	RPD %	RPD Limits %
Diesel Range Organics (DRO)	47.8	ND	33.5	42.3	53.9	71.3	5	50.0-150		J3	23.2	20
<i>(S) o-Terphenyl</i>					129	62.5		18.0-148				

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

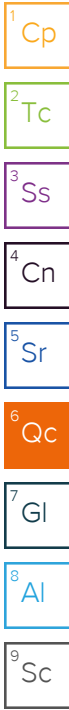
8 Al

9 Sc

Method Blank (MB)

(MB) R3723791-2 10/30/21 10:59

Analyte	MB Result mg/kg	MB Qualifier	MB MDL mg/kg	MB RDL mg/kg
Anthracene	U		0.00230	0.00600
Acenaphthene	U		0.00209	0.00600
Acenaphthylene	U		0.00216	0.00600
Benzo(a)anthracene	U		0.00173	0.00600
Benzo(a)pyrene	U		0.00179	0.00600
Benzo(b)fluoranthene	U		0.00153	0.00600
Benzo(g,h,i)perylene	U		0.00177	0.00600
Benzo(k)fluoranthene	U		0.00215	0.00600
Chrysene	U		0.00232	0.00600
Dibenz(a,h)anthracene	U		0.00172	0.00600
Fluoranthene	U		0.00227	0.00600
Fluorene	U		0.00205	0.00600
Indeno(1,2,3-cd)pyrene	U		0.00181	0.00600
Naphthalene	U		0.00408	0.0200
Phenanthrene	U		0.00231	0.00600
Pyrene	U		0.00200	0.00600
1-Methylnaphthalene	U		0.00449	0.0200
2-Methylnaphthalene	U		0.00427	0.0200
2-Chloronaphthalene	U		0.00466	0.0200
(S) Nitrobenzene-d5	105			14.0-149
(S) 2-Fluorobiphenyl	87.4			34.0-125
(S) p-Terphenyl-d14	109			23.0-120



Laboratory Control Sample (LCS)

(LCS) R3723791-1 10/30/21 10:42

Analyte	Spike Amount mg/kg	LCS Result mg/kg	LCS Rec. %	Rec. Limits %	LCS Qualifier
Anthracene	0.0800	0.0587	73.4	50.0-126	
Acenaphthene	0.0800	0.0587	73.4	50.0-120	
Acenaphthylene	0.0800	0.0573	71.6	50.0-120	
Benzo(a)anthracene	0.0800	0.0587	73.4	45.0-120	
Benzo(a)pyrene	0.0800	0.0532	66.5	42.0-120	
Benzo(b)fluoranthene	0.0800	0.0611	76.4	42.0-121	
Benzo(g,h,i)perylene	0.0800	0.0611	76.4	45.0-125	
Benzo(k)fluoranthene	0.0800	0.0601	75.1	49.0-125	
Chrysene	0.0800	0.0615	76.9	49.0-122	
Dibenz(a,h)anthracene	0.0800	0.0610	76.3	47.0-125	
Fluoranthene	0.0800	0.0596	74.5	49.0-129	

Laboratory Control Sample (LCS)

(LCS) R3723791-1 10/30/21 10:42

Analyte	Spike Amount mg/kg	LCS Result mg/kg	LCS Rec. %	Rec. Limits %	LCS Qualifier
Fluorene	0.0800	0.0620	77.5	49.0-120	
Indeno(1,2,3-cd)pyrene	0.0800	0.0590	73.8	46.0-125	
Naphthalene	0.0800	0.0582	72.8	50.0-120	
Phenanthrene	0.0800	0.0627	78.4	47.0-120	
Pyrene	0.0800	0.0593	74.1	43.0-123	
1-Methylnaphthalene	0.0800	0.0548	68.5	51.0-121	
2-Methylnaphthalene	0.0800	0.0538	67.3	50.0-120	
2-Chloronaphthalene	0.0800	0.0555	69.4	50.0-120	
(S) Nitrobenzene-d5			107	14.0-149	
(S) 2-Fluorobiphenyl			79.1	34.0-125	
(S) p-Terphenyl-d14			103	23.0-120	

L1422174-01 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1422174-01 10/30/21 16:46 • (MS) R3723791-3 10/30/21 17:04 • (MSD) R3723791-4 10/30/21 17:21

Analyte	Spike Amount mg/kg	Original Result mg/kg	MS Result mg/kg	MSD Result mg/kg	MS Rec. %	MSD Rec. %	Dilution	Rec. Limits %	MS Qualifier	MSD Qualifier	RPD %	RPD Limits %
Anthracene	0.0760	0.00867	0.0599	0.0670	67.4	76.8	1	10.0-145			11.2	30
Acenaphthene	0.0760	ND	0.0537	0.0618	63.3	73.9	1	14.0-127			14.0	27
Acenaphthylene	0.0760	ND	0.0555	0.0625	73.0	82.2	1	21.0-124			11.9	25
Benzo(a)anthracene	0.0760	0.0208	0.0739	0.0757	69.9	72.2	1	10.0-139			2.41	30
Benzo(a)pyrene	0.0760	0.0134	0.0567	0.0634	57.0	65.8	1	10.0-141			11.2	31
Benzo(b)fluoranthene	0.0760	0.0366	0.0795	0.0783	56.4	54.9	1	10.0-140			1.52	36
Benzo(g,h,i)perylene	0.0760	0.0259	0.0604	0.0733	45.4	62.4	1	10.0-140			19.3	33
Benzo(k)fluoranthene	0.0760	0.00795	0.0550	0.0627	61.9	72.0	1	10.0-137			13.1	31
Chrysene	0.0760	0.0322	0.0986	0.0974	87.4	85.8	1	10.0-145			1.22	30
Dibenz(a,h)anthracene	0.0760	ND	0.0479	0.0656	56.6	79.9	1	10.0-132	J3		31.2	31
Fluoranthene	0.0760	0.0456	0.0859	0.0880	53.0	55.8	1	10.0-153			2.42	33
Fluorene	0.0760	ND	0.0620	0.0633	81.6	83.3	1	11.0-130			2.08	29
Indeno(1,2,3-cd)pyrene	0.0760	0.0145	0.0627	0.0731	63.4	77.1	1	10.0-137			15.3	32
Naphthalene	0.0760	0.0545	0.0884	0.0944	44.6	52.5	1	10.0-135			6.56	27
Phenanthrene	0.0760	0.137	0.160	0.145	30.3	10.5	1	10.0-144			9.84	31
Pyrene	0.0760	0.0469	0.0944	0.0809	62.5	44.7	1	10.0-148			15.4	35
1-Methylnaphthalene	0.0760	0.0965	0.119	0.123	29.6	34.9	1	10.0-142			3.31	28
2-Methylnaphthalene	0.0760	0.121	0.143	0.138	28.9	22.4	1	10.0-137			3.56	28
2-Chloronaphthalene	0.0760	ND	0.0497	0.0601	65.4	79.1	1	29.0-120			18.9	24
(S) Nitrobenzene-d5					114	128		14.0-149				
(S) 2-Fluorobiphenyl					81.2	84.2		34.0-125				
(S) p-Terphenyl-d14					98.9	104		23.0-120				

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc

# GLOSSARY OF TERMS

## Guide to Reading and Understanding Your Laboratory Report

The information below is designed to better explain the various terms used in your report of analytical results from the Laboratory. This is not intended as a comprehensive explanation, and if you have additional questions please contact your project representative.

Results Disclaimer - Information that may be provided by the customer, and contained within this report, include Permit Limits, Project Name, Sample ID, Sample Matrix, Sample Preservation, Field Blanks, Field Spikes, Field Duplicates, On-Site Data, Sampling Collection Dates/Times, and Sampling Location. Results relate to the accuracy of this information provided, and as the samples are received.

### Abbreviations and Definitions

(dry)	Results are reported based on the dry weight of the sample. [this will only be present on a dry report basis for soils].
MDL	Method Detection Limit.
ND	Not detected at the Reporting Limit (or MDL where applicable).
RDL	Reported Detection Limit.
RDL (dry)	Reported Detection Limit.
Rec.	Recovery.
RPD	Relative Percent Difference.
SDG	Sample Delivery Group.
(S)	Surrogate (Surrogate Standard) - Analytes added to every blank, sample, Laboratory Control Sample/Duplicate and Matrix Spike/Duplicate; used to evaluate analytical efficiency by measuring recovery. Surrogates are not expected to be detected in all environmental media.
U	Not detected at the Reporting Limit (or MDL where applicable).
Analyte	The name of the particular compound or analysis performed. Some Analyses and Methods will have multiple analytes reported.
Dilution	If the sample matrix contains an interfering material, the sample preparation volume or weight values differ from the standard, or if concentrations of analytes in the sample are higher than the highest limit of concentration that the laboratory can accurately report, the sample may be diluted for analysis. If a value different than 1 is used in this field, the result reported has already been corrected for this factor.
Limits	These are the target % recovery ranges or % difference value that the laboratory has historically determined as normal for the method and analyte being reported. Successful QC Sample analysis will target all analytes recovered or duplicated within these ranges.
Original Sample	The non-spiked sample in the prep batch used to determine the Relative Percent Difference (RPD) from a quality control sample. The Original Sample may not be included within the reported SDG.
Qualifier	This column provides a letter and/or number designation that corresponds to additional information concerning the result reported. If a Qualifier is present, a definition per Qualifier is provided within the Glossary and Definitions page and potentially a discussion of possible implications of the Qualifier in the Case Narrative if applicable.
Result	The actual analytical final result (corrected for any sample specific characteristics) reported for your sample. If there was no measurable result returned for a specific analyte, the result in this column may state "ND" (Not Detected) or "BDL" (Below Detectable Levels). The information in the results column should always be accompanied by either an MDL (Method Detection Limit) or RDL (Reporting Detection Limit) that defines the lowest value that the laboratory could detect or report for this analyte.
Uncertainty (Radiochemistry)	Confidence level of 2 sigma.
Case Narrative (Cn)	A brief discussion about the included sample results, including a discussion of any non-conformances to protocol observed either at sample receipt by the laboratory from the field or during the analytical process. If present, there will be a section in the Case Narrative to discuss the meaning of any data qualifiers used in the report.
Quality Control Summary (Qc)	This section of the report includes the results of the laboratory quality control analyses required by procedure or analytical methods to assist in evaluating the validity of the results reported for your samples. These analyses are not being performed on your samples typically, but on laboratory generated material.
Sample Chain of Custody (Sc)	This is the document created in the field when your samples were initially collected. This is used to verify the time and date of collection, the person collecting the samples, and the analyses that the laboratory is requested to perform. This chain of custody also documents all persons (excluding commercial shippers) that have had control or possession of the samples from the time of collection until delivery to the laboratory for analysis.
Sample Results (Sr)	This section of your report will provide the results of all testing performed on your samples. These results are provided by sample ID and are separated by the analyses performed on each sample. The header line of each analysis section for each sample will provide the name and method number for the analysis reported.
Sample Summary (Ss)	This section of the Analytical Report defines the specific analyses performed for each sample ID, including the dates and times of preparation and/or analysis.

Qualifier	Description
C3	The reported concentration is an estimate. The continuing calibration standard associated with this data responded low. Method sensitivity check is acceptable.
C4	The reported concentration is an estimate. The continuing calibration standard associated with this data responded low. Data is likely to show a low bias concerning the result.
J	The identification of the analyte is acceptable; the reported value is an estimate.
J1	Surrogate recovery limits have been exceeded; values are outside upper control limits.
J3	The associated batch QC was outside the established quality control range for precision.

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc

# ACCREDITATIONS & LOCATIONS

## Pace Analytical National 12065 Lebanon Rd Mount Juliet, TN 37122

Alabama	40660	Nebraska	NE-OS-15-05
Alaska	17-026	Nevada	TN000032021-1
Arizona	AZ0612	New Hampshire	2975
Arkansas	88-0469	New Jersey–NELAP	TN002
California	2932	New Mexico <sup>1</sup>	TN00003
Colorado	TN00003	New York	11742
Connecticut	PH-0197	North Carolina	Env375
Florida	E87487	North Carolina <sup>1</sup>	DW21704
Georgia	NELAP	North Carolina <sup>3</sup>	41
Georgia <sup>1</sup>	923	North Dakota	R-140
Idaho	TN00003	Ohio–VAP	CL0069
Illinois	200008	Oklahoma	9915
Indiana	C-TN-01	Oregon	TN200002
Iowa	364	Pennsylvania	68-02979
Kansas	E-10277	Rhode Island	LA000356
Kentucky <sup>1,6</sup>	KY90010	South Carolina	84004002
Kentucky <sup>2</sup>	16	South Dakota	n/a
Louisiana	AI30792	Tennessee <sup>1,4</sup>	2006
Louisiana	LA018	Texas	T104704245-20-18
Maine	TN00003	Texas <sup>5</sup>	LAB0152
Maryland	324	Utah	TN000032021-11
Massachusetts	M-TN003	Vermont	VT2006
Michigan	9958	Virginia	110033
Minnesota	047-999-395	Washington	C847
Mississippi	TN00003	West Virginia	233
Missouri	340	Wisconsin	998093910
Montana	CERT0086	Wyoming	A2LA
A2LA – ISO 17025	1461.01	AIHA-LAP,LLC EMLAP	100789
A2LA – ISO 17025 <sup>5</sup>	1461.02	DOD	1461.01
Canada	1461.01	USDA	P330-15-00234
EPA–Crypto	TN00003		

<sup>1</sup> Drinking Water <sup>2</sup> Underground Storage Tanks <sup>3</sup> Aquatic Toxicity <sup>4</sup> Chemical/Microbiological <sup>5</sup> Mold <sup>6</sup> Wastewater n/a Accreditation not applicable

\* Not all certifications held by the laboratory are applicable to the results reported in the attached report.

\* Accreditation is only applicable to the test methods specified on each scope of accreditation held by Pace Analytical.

<sup>1</sup> Cp

<sup>2</sup> Tc

<sup>3</sup> Ss

<sup>4</sup> Cn

<sup>5</sup> Sr

<sup>6</sup> Qc

<sup>7</sup> Gl

<sup>8</sup> Al

<sup>9</sup> Sc

Company/Address: **GeoEngineers- Portland, OR**  
 4000 Kruse Way Place  
 Bldg. 3, Suite 200  
 Lake Oswego, OR 97035

Billing Information:  
 Accounts Payable (Marlee Johnston)  
 17425 NE Union Hill Rd, Suite 250  
 Redmond, WA 98052

Report to: **Cris Watkins**  
 Email To: [cwatkins@geoengineers.com](mailto:cwatkins@geoengineers.com)

Project Description: \_\_\_\_\_ City/State Collected: **Seeta, WA** Please Circle: PT MT CT ET

Phone: **503-603-6661** Client Project #: **0180-393-00** Lab Project #: **GEOENGPOR-018039300**

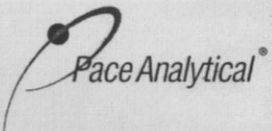
Collected by (print): **Dexter Chem** Site/Facility ID #: \_\_\_\_\_ P.O. #: \_\_\_\_\_

Collected by (signature): *[Signature]* **Rush?** (Lab MUST Be Notified)  
 \_\_\_ Same Day \_\_\_ Five Day \_\_\_ Next Day \_\_\_ 5 Day (Rad Only)  
 \_\_\_ Two Day \_\_\_ **X** 10 Day (Rad Only) \_\_\_ Three Day

Immediately Packed on Ice N \_\_\_ Y **X** Date Results Needed: \_\_\_\_\_ No. of Cntrs: \_\_\_\_\_

Sample ID	Comp/Grab	Matrix *	Depth	Date	Time	Cntrs	NWTPHDX no silica 8ozClr-NoPres	NWTPHDX w/ silica 8ozClr-NoPres	NWTPHGX 40mlAmb/MeOH10ml/Syr	PAHs 8270ESIM 8ozClr-NoPres	RCRA8 Metals 6020 4ozClr-NoPres	VOCs 8260D 40mlAmb/MeOH10ml/Syr	Hold
GEL-1-9-10		SS	9-10	10/25/21	1120	3							X
GEL-1-14-15		SS	14-15		1125								X
GEL-2-9-10		SS	9-10		1140								X
GEL-2-14-15		SS	14-15		1145								X
GEL-3-9-10		SS	9-10		BK5								X
GEL-3-14-15		SS	14-15		1230								X
GEL-4-9-10		SS	9-10		1430								X
GEL-4-14-15		SS	14-15		1430								X
MW-1-9-10		SS	9-10		1225								X
MW-1-14-15		SS	14-15		1230								X

Chain of Custody Page 1 of 3



12065 Lebanon Rd Mount Juliet, TN 37122  
 Submitting a sample via this chain of custody constitutes acknowledgment and acceptance of the Pace Terms and Conditions found at: <https://info.pacelabs.com/hubfs/pas-standard-terms.pdf>

SDG # **1423266**  
**J226**

Acctnum: **GEOENGPOR**  
 Template: **T197917**  
 Prelogin: **P882161**  
 PM: **110 - Brian Ford**  
 PB: \_\_\_\_\_  
 Shipped Via: \_\_\_\_\_

Remarks	Sample # (lab only)
	-01
	-02
	-03
	-04
	-05
	-06
	-07
	-08
	-09
	-10

\* Matrix: SS - Soil AIR - Air F - Filter  
 GW - Groundwater B - Bioassay  
 WW - WasteWater  
 DW - Drinking Water  
 OT - Other \_\_\_\_\_

Remarks: \_\_\_\_\_

Samples returned via: \_\_\_ UPS \_\_\_ FedEx \_\_\_ Courier \_\_\_\_\_ Tracking # **5217 3314 1922**

pH \_\_\_\_\_ Temp \_\_\_\_\_  
 Flow \_\_\_\_\_ Other \_\_\_\_\_

**Sample Receipt Checklist**

COC Seal Present/Intact:	NP	Y	N
COC Signed/Accurate:		Y	N
Bottles arrive intact:		Y	N
Correct bottles used:		Y	N
Sufficient volume sent:		Y	N
If Applicable			
VOA Zero Headspace:		Y	N
Preservation Correct/Checked:		Y	N
RAD Screen <0.5 mR/hr:		Y	N

Relinquished by: (Signature) <i>[Signature]</i>	Date: <b>10/26/21</b>	Time: <b>1500</b>	Received by: (Signature) <i>[Signature]</i>	Trip Blank Received: <input checked="" type="checkbox"/> Yes / <input type="checkbox"/> No <input checked="" type="checkbox"/> HCL / MeOH <input type="checkbox"/> TBR
Relinquished by: (Signature) _____	Date: _____	Time: _____	Received by: (Signature) _____	Temp: <b>17.0</b> °C Bottles Received: <b>17 63</b>
Relinquished by: (Signature) _____	Date: _____	Time: _____	Received for lab by: (Signature) <i>[Signature]</i>	Date: <b>10/27/21</b> Time: <b>930</b>

Hold: \_\_\_\_\_ Condition: **NCF / OK**


Company Address: **GeoEngineers- Portland, OR**  
**4000 Kruse Way Place**  
**Bldg. 3, Suite 200**  
**Lake Oswego, OR 97035**

Billing Information:  
**Accounts Payable (Marlee Johnston)**  
**17425 NE Union Hill Rd, Suite 250**  
**Redmond, WA 98052**

Report to:  
**Cris Watkins**

Email To: **cwatkins@geoengineers.com**

Chain of Custody Page **2** of **3**



12065 Lebanon Rd Mount Juliet, TN 37122  
 Submitting a sample via this chain of custody constitutes acknowledgment and acceptance of the Pace Terms and Conditions found at:  
<https://info.pacelabs.com/hubs/pas-standard-terms.pdf>

Project Description: \_\_\_\_\_ City/State Collected: **Seattle, WA** Please Circle: PT MT CT ET

Phone: **503-603-6661** Client Project #: **0180-393-00** Lab Project #: **GEOENGPOR-018039300**

Collected by (print): \_\_\_\_\_ Site/Facility ID # \_\_\_\_\_ P.O. # \_\_\_\_\_

Collected by (signature): *Dexter Chen* **Rush?** (Lab MUST Be Notified)  
 Same Day  Five Day  
 Next Day  5 Day (Rad Only)  
 Two Day  10 Day (Rad Only)  
 Three Day

Immediately Packed on Ice N  Y  Quote # \_\_\_\_\_ Date Results Needed \_\_\_\_\_ No. of Cntrs \_\_\_\_\_

Analysis / Container / Preservative	Pres Chk
NWTPHDX no silica 8ozClr-NoPres	
NWTPHDX w/ silica 8ozClr-NoPres	
NWTPHGX 40mlAmb/MeOH10ml/Syr	
PAHs 8270ESIM 8ozClr-NoPres	
RCRA8 Metals 6020 4ozClr-NoPres	
VOCs 8260D 40mlAmb/MeOH10ml/Syr	

SDG #: **1423268**

Table #: \_\_\_\_\_

Acctnum: **GEOENGPOR**

Template: **T197917**

Prelogin: **P882161**

PM: **110 - Brian Ford**

PB: \_\_\_\_\_

Shipped Via: \_\_\_\_\_

Sample ID	Comp/Grab	Matrix *	Depth	Date	Time	No. of Cntrs
<b>MW-2-9-10</b>		<b>SS</b>	<b>9-10</b>	<b>10/25/21</b>	<b>1400</b>	<b>3</b>
<b>MW-2-14-15</b>		<b>SS</b>	<b>14-15</b>	<b>10/25/21</b>	<b>1405</b>	
<b>GE1-4-2-3</b>		<b>SS</b>	<b>2-3</b>	<b>10/26/21</b>	<b>1000</b>	
<b>MW-1-2-3</b>		<b>SS</b>	<b>2-3</b>		<b>1020</b>	
<b>GE1-5-7-8</b>		<b>SS</b>	<b>7-8</b>		<b>1040</b>	
<b>GE1-5-15-16</b>		<b>SS</b>	<b>15-16</b>		<b>1045</b>	
<b>MW-3-2-3</b>		<b>SS</b>	<b>2-3</b>		<b>1100</b>	
<b>MW-3-7-8</b>		<b>SS</b>	<b>7-8</b>		<b>1105</b>	
<b>MW-3-14-15</b>		<b>SS</b>	<b>14-15</b>		<b>1110</b>	
<b>GE1-6-7-8</b>		<b>SS</b>	<b>7-8</b>		<b>1205</b>	

\* Matrix: **SS - Soil AIR - Air F - Filter**  
**GW - Groundwater B - Bioassay**  
**WW - WasteWater**  
**DW - Drinking Water**  
**OT - Other**

Remarks: \_\_\_\_\_

Samples returned via:  UPS  FedEx  Courier Tracking # \_\_\_\_\_

Sample Receipt Checklist

COC Seal Present/Intact:  Y  N

COC Signed/Accurate:  Y  N

Bottles arrive intact:  Y  N

Correct bottles used:  Y  N

Sufficient volume sent:  Y  N

If Applicable

VOA Zero Headspace:  Y  N

Preservation Correct/Checked:  Y  N

RAD Screen <0.5 mR/hr:  Y  N

Relinquished by: (Signature) *[Signature]* Date: **10/26/21** Time: **1520** Received by: (Signature) \_\_\_\_\_ Trip Blank Received: Yes  No  TBR

Relinquished by: (Signature) \_\_\_\_\_ Date: \_\_\_\_\_ Time: \_\_\_\_\_ Received by: (Signature) \_\_\_\_\_ Temp: **17.0** °C Bottles Received: **63**

Relinquished by: (Signature) \_\_\_\_\_ Date: \_\_\_\_\_ Time: \_\_\_\_\_ Received for lab by: (Signature) *[Signature]* Date: **10/27/21** Time: **930** Hold: \_\_\_\_\_ Condition: **NCF / OK**

Company Name/Address: **GeoEngineers- Portland, OR**  
**4000 Kruse Way Place**  
**Bldg. 3, Suite 200**  
**Lake Oswego, OR 97035**

Billing Information:  
**Accounts Payable (Marlee Johnston)**  
**17425 NE Union Hill Rd, Suite 250**  
**Redmond, WA 98052**

Report to:  
**Cris Watkins**

Email To: **cwatkins@geoengineers.com**

City/State Collected: **Seattle, WA**

Please Circle: **PT MT CT ET**

Project Description: \_\_\_\_\_

Phone: **503-603-6661**

Client Project # **0180-393-00**

Lab Project # **GEOENGPOR-018039300**

Collected by (print): **Deborah Chan**

Site/Facility ID # \_\_\_\_\_

P.O. # \_\_\_\_\_

Collected by (signature): \_\_\_\_\_

**Rush?** (Lab MUST Be Notified)  
 \_\_\_ Same Day \_\_\_ Five Day  
 \_\_\_ Next Day \_\_\_ 5 Day (Rad Only)  
 \_\_\_ Two Day  10 Day (Rad Only)  
 \_\_\_ Three Day

Date Results Needed \_\_\_\_\_

Immediately Packed on Ice N \_\_\_ Y \_\_\_

No. of Cntrs \_\_\_\_\_

Sample ID	Comp/Grab	Matrix *	Depth	Date	Time	No. of Cntrs	NWTPHDX no silica 8ozClr-NoPres	NWTPHDX w/ silica 8ozClr-NoPres	NWTPHGX 40mlAmb/MeOH10ml/Syr	PAHs 8270ESIM 8ozClr-NoPres	RCRA8 Metals 6020 4ozClr-NoPres	VOCs 8260D 40mlAmb/MeOH10ml/Syr	Chain of Custody
<b>201-6-14-15</b>		<b>SS</b>	<b>14-15</b>	<b>10/26/21</b>	<b>1210</b>	<b>3</b>							Pace Analytical 12065 Lebanon Rd Mount Juliet, TN 37122 Submitting a sample via this chain of custody constitutes acknowledgment and acceptance of the Pace Terms and Conditions found at: <a href="https://info.pacelabs.com/hubs/pas-standard-terms.pdf">https://info.pacelabs.com/hubs/pas-standard-terms.pdf</a> SDG # <b>1923266</b> Table # _____ Acctnum: <b>GEOENGPOR</b> Template: <b>T197917</b> Prelogin: <b>P882161</b> PM: <b>110 - Brian Ford</b> PB: _____ Shipped Via: _____ Remarks   Sample # (lab only) _____   <b>-21</b>

\* Matrix: **SS - Soil AIR - Air F - Filter**  
**GW - Groundwater B - Bioassay**  
**WW - WasteWater**  
**DW - Drinking Water**  
**OT - Other \_\_\_\_\_**

Remarks: \_\_\_\_\_

Samples returned via:  UPS  FedEx  Courier \_\_\_\_\_

Tracking # \_\_\_\_\_

Relinquished by: (Signature) \_\_\_\_\_ Date: **10/26/21** Time: **1500**

Received by: (Signature) \_\_\_\_\_

Trip Blank Received:  Yes  No  
 HCL/MeOH TBR

Temp: **17.0** °C Bottles Received: **63**

Relinquished by: (Signature) \_\_\_\_\_ Date: \_\_\_\_\_ Time: \_\_\_\_\_

Received for lab by: (Signature) \_\_\_\_\_ Date: **10/27/21** Time: **930**

Hold: \_\_\_\_\_ Condition: **NCF / OK**

Sample Receipt Checklist:  
 COC Seal Present/Intact:  Y  N  
 COC Signed/Accurate:  Y  N  
 Bottles arrive intact:  Y  N  
 Correct bottles used:  Y  N  
 Sufficient volume sent:  Y  N  
 If Applicable  
 VOA Zero Headspace:  Y  N  
 Preservation Correct/Checked:  Y  N  
 RAD Screen <0.5 mR/hr:  Y  N

## Brian Ford

---

**From:** Cris J. Watkins <cwatkins@geoengineers.com>  
**Sent:** Thursday, October 28, 2021 5:54 PM  
**To:** Brian Ford  
**Cc:** Phil D. Welker  
**Subject:** RE: updated analysis request FW: L1423266 Proj 0180-393-00 Hold Samples Received

CAUTION: This email originated from outside Pace Analytical. Do not click links or open attachments unless you recognize the sender and know the content is safe.

Also – please update the project name to WSDOT-A-1 Towing. Thanks,

**Cris J. Watkins**  
**Senior Environmental Scientist | GeoEngineers, Inc.**  
Telephone: 503.603.6687  
Mobile: 971.235.6147

---

**From:** Cris J. Watkins  
**Sent:** Thursday, October 28, 2021 3:53 PM  
**To:** Brian Ford <Brian.Ford@pacelabs.com>  
**Cc:** Phil D. Welker <pwelker@geoengineers.com>  
**Subject:** updated analysis request FW: L1423266 Proj 0180-393-00 Hold Samples Received

Hey Brian,

Can you run the following samples on a 5-day TAT:

GEI-1-9-10  
GEI-2-9-10  
GEI-3-9-10  
GEI-4-2-3  
GEI-5-7-8  
GEI-6-7-8  
MW-1-2-3  
MW-2-9-10  
MW-3-2-3

We will run Dx-no silica, Gx, PAHs, RCRA 8 and VOCs.

Thanks,

**Cris J. Watkins**  
**Senior Environmental Scientist | GeoEngineers, Inc.**  
Telephone: 503.603.6687  
Mobile: 971.235.6147

---

**From:** Brian Ford <[Brian.Ford@pacelabs.com](mailto:Brian.Ford@pacelabs.com)>  
**Sent:** Thursday, October 28, 2021 8:22 AM  
**To:** Cris J. Watkins <[cwatkins@geoengineers.com](mailto:cwatkins@geoengineers.com)>  
**Subject:** L1423266 Proj 0180-393-00 Hold Samples Received

[EXTERNAL]

Thanks,



**Brian Ford**

*Project Manager 2 / Pace National*

12065 Lebanon Road | Mt. Juliet, TN 37122

Office: 615.773.9772

[brian.ford@pacelabs.com](mailto:brian.ford@pacelabs.com)

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 Please consider the environment before printing this email

Confidentiality: This message is confidential and intended solely for use of the individual or entity to whom it is addressed. If you are not the person for whom this message is intended, please delete it and notify me immediately, and please do not copy or send this message to anyone else.

**APPENDIX C**  
**Report Limitations and Guidelines for Use**

## **APPENDIX C**

### **REPORT LIMITATIONS AND GUIDELINES FOR USE<sup>1</sup>**

This appendix provides information to help you manage your risks with respect to the use of this report.

#### **Read These Provisions Closely**

Some clients, design professionals, and contractors may not recognize that the geoscience practices (geotechnical engineering, geology and environmental science) are far less exact than other engineering and natural science disciplines. This lack of understanding can create unrealistic expectations that could lead to disappointments, claims and disputes. GeoEngineers includes these explanatory “limitations” provisions in our reports to help reduce such risks. Please confer with GeoEngineers if you are unclear how these “Report Limitations and Guidelines for Use” apply to your project or site.

#### **Environmental Services Are Performed for Specific Purposes, Persons and Projects**

This report has been prepared for the exclusive use of the WSDOT, their authorized agents and regulatory agencies. This report is not intended for use by others, and the information contained herein is not applicable to other sites.

GeoEngineers structures our services to meet the specific needs of our clients. For example, an environmental site assessment or remedial action study conducted for a property owner may not fulfill the needs of a prospective purchaser of the same property. Because each environmental study is unique, each environmental report is unique, prepared solely for the specific client and project site. No one except WSDOT should rely on this plan without first conferring with GeoEngineers. This report should not be applied for any purpose or project except the one originally contemplated.

#### **This Environmental Report Is Based on a Unique Set of Project-Specific Factors**

This report applies to the King County Tax Parcel 3223049048 project located in Des Moines, Washington. GeoEngineers considered a number of unique, project-specific factors when establishing the scope of services for this project and report. Unless GeoEngineers specifically indicates otherwise, do not rely on this report if it was:

- Not prepared for you,
- Not prepared for your project,
- Not prepared for the specific site explored, or
- Completed before important project changes were made.

If important changes are made after the date of this remedial action plan, GeoEngineers should be given the opportunity to review our interpretations and recommendations and provide written modifications or confirmation, as appropriate.

---

<sup>1</sup> Developed based on material provided by Geoprofessional Business Association (GBA), Professional Firms Practicing in the GeoSciences, [www.geoprofesional.org](http://www.geoprofesional.org).

### **Reliance Conditions for Third Parties**

No third party may rely on the product of our services unless GeoEngineers agrees in advance, and in writing to such reliance. This is to provide our firm with reasonable protection against open-ended liability claims by third parties with whom there would otherwise be no contractual limits to their actions.

### **Environmental Regulations Are Always Evolving**

Some substances may be present in the site vicinity in quantities or under conditions that may have led, or may lead, to contamination of the subject site, but are not included in current local, state or federal regulatory definitions of hazardous substances or do not otherwise present current potential liability. GeoEngineers cannot be responsible if the standards for appropriate inquiry, or regulatory definitions of hazardous substance, change or if more stringent environmental standards are developed in the future.

### **Subsurface Conditions Can Change**

This report is based on conditions that existed at the time our site studies were performed. The findings and conclusions of this report may be affected by the passage of time, by manmade events such as construction on or adjacent to the site, by new releases of hazardous substances, or by natural events such as floods, earthquakes, slope instability or groundwater fluctuations. Always contact GeoEngineers before applying this report to determine if it is still applicable.

### **Soil and Groundwater End Use**

The cleanup levels referenced in this report are site- and situation-specific. The cleanup levels may not be applicable for other sites or for other on-site uses of the affected media (soil and/or groundwater). Note that hazardous substances may be present in some of the Site soil and/or groundwater at detectable concentrations that are less than the referenced cleanup levels. GeoEngineers should be contacted prior to the export of soil or groundwater from the subject Site or reuse of the affected media on Site to evaluate the potential for associated environmental liabilities. We cannot be responsible for potential environmental liability arising out of the transfer of soil and/or groundwater from the subject site to another location or its reuse on site in instances that we were not aware of or could not control.

### **Biological Pollutants**

GeoEngineers' Scope of Work specifically excludes the investigation, detection, prevention or assessment of the presence of Biological Pollutants. Accordingly, this report does not include any interpretations, recommendations, findings, or conclusions regarding the detecting, assessing, preventing or abating of Biological Pollutants and no conclusions or inferences should be drawn regarding Biological Pollutants, as they may relate to this project. The term "Biological Pollutants" includes, but is not limited to, molds, fungi, spores, bacteria, and viruses, and/or any of their byproducts.

If the Client desires these specialized services, they should be obtained from a consultant who offers services in this specialized field.

### **Do Not Redraw the Exploration Logs**

Environmental scientists prepare final boring and testing logs based upon their interpretation of field logs and laboratory data. To prevent errors or omissions, the logs included in an environmental report should never be redrawn for inclusion in other design drawings. Only photographic or electronic reproduction is acceptable, but recognize that separating logs from the report can elevate risk.

### **Geotechnical, Geologic and Environmental Reports Should Not Be Interchanged**

The equipment, techniques and personnel used to perform an environmental study differ significantly from those used to perform a geotechnical or geologic study and vice versa. For that reason, a geotechnical engineering or geologic report does not usually relate any environmental findings, conclusions or recommendations; e.g., about the likelihood of encountering underground storage tanks or regulated contaminants. Similarly, environmental reports are not used to address geotechnical or geologic concerns regarding a specific project.

### **Most Environmental Findings Are Professional Opinions**

Our interpretations of subsurface conditions are based on field observations and chemical analytical data from the sampling locations at the site documented in past reports. Site exploration identifies subsurface conditions only at those points where subsurface tests are conducted or samples are taken. GeoEngineers reviewed field and laboratory data and then applied our professional judgment to render an opinion about subsurface conditions throughout the site. Actual subsurface conditions may differ – sometimes significantly – from those indicated in this report. There is always a potential that areas of contamination exist in portions of the site that were not sampled or tested during this or previous studies. Our remedial action plan, conclusions and interpretations should not be construed as a warranty of the subsurface conditions.



**Phase II Environmental Site Assessment, Soil Sampling Report  
State Route (SR) 509  
Foreman Property 18451 12th Avenue South**

Prepared for

**Washington State Department of Transportation  
SR 509 Project Office**

Prepared by

**Headquarters Hazardous Materials and Solid Waste Program  
Environmental Services Office**

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## Acronyms and Abbreviations

ASTM E 1903-11	American Society for Testing and Materials Standard Practice for Environmental Site Assessments: Phase II Environmental Assessment Process, Designation: E 1903-11
bgs	Below ground surface
CUL	Cleanup Level
Ecology	Washington State Department of Ecology
ESA	Environmental Site Assessment
ESO	Environmental Services Office
HazMat Program	Hazardous Materials and Solid Waste Program
HQ	Headquarters
MTCA	Model Toxics Control Act
NFA	No Further Action
NWR	Northwest Region
RCRA	Resource Conservation and Recovery Act
WAC	Washington Administrative Code
WSDOT	Washington State Department of Transportation

## **1.0 INTRODUCTION**

The Washington State Department of Transportation (WSDOT) Headquarters (HQ) Hazardous Materials and Solid Waste (HazMat) Program prepared this Phase II Environmental Site Assessment (ESA) Report to provide a summary of information associated with soil investigation and excavation activities within existing unconstructed WSDOT right-of-way (ROW) at 18451 12th Avenue South in Des Moines, WA (Site) (See Appendix 1).. The Site is associated with the WSDOT 509 Project Office in support of the future construction of the State Route (SR) 509 Roadway Project.

The Phase II ESA soil sampling activities assesses soil conditions within the property and confirms the presence or absence of contaminated soil after cleanup activities. A contractor, Marine Vacuum Incorporated (Marine Vac.) remediated two areas of confirmed and suspected contaminated materials.

HazMat conducted this Phase II ESA in general accordance with the American Society for Testing and Materials Standard Practice for Environmental Site Assessments (ASTM): Phase II Environmental Site Assessment Process, Designation: E 1903-11 (ASTM 1903-1) and the WSDOT Environmental Manual (EM) 31-11, Chapter 447.

## **1.1 SITE INFORMATION**

The following sections provide information pertaining to the background, physical, geological and hydrogeological setting of the site.

### **1.1.1 Physical Setting and Historical Background**

WSDOT purchased this Site as part of the future SR 509 Project, and leased the property to tenants since 2000. The Site is located near the intersection of Des Moines Memorial Drive and South 188 Street and the intersection of SR 509 and South 188 Street. The previous tenant, Mr. Foreman, conducted welding activities on this property from 2000 through 2016. In the summer of 2016, WSDOT allowed the Washington State Patrol (WSP) to use a portion of the property to dismantle seized vehicles. Currently, A-1 Towing leases the property as a towing yard, which includes parking vehicles throughout the site and repairing the vehicles within the garage structure.

### **1.2.1 Geologic and Hydrogeologic Setting**

Based upon review available information near the Site vicinity, soils from surface to six feet (ft.) below ground surface (bgs) are defined as poorly graded sand, medium dense, olive gray, moist and homogenous. Soils from six to 10 ft. are defined as poorly graded sand, medium dense, olive gray, moist, and homogeneous. Soils from 10 to 15 ft. are defined as silty sand, dense, grayish brown, moist, and homogenous.

Groundwater depth varied across the Site; although, it was encountered as shallow as three ft. bgs in select locations.

## 2.0 BACKGROUND INFORMATION

The following provides a chronological order of the relevant events that led up to the cleanup of contaminated soils on the Site. The background describes two separate sample events prior to the contaminated soil remedial activities and confirmation sampling.

In October 2016, WSDOT Northwest Region (NWR) Real Estate Services (RES) requested the WSDOT NWR Hazardous Materials Specialist to perform a limited Phase II Environmental Site Assessment (ESA) after MR. Foreman's lease terminated. WSDOT NWR sampled four locations (A, B, C, & D) (See Appendix 1) of suspected contaminated soil identified on the property. NWR HazMat identified two relatively large heavily stained areas, numerous small stains located on the north end of the site, and abandoned solid waste debris and trash scattered throughout. WSDOT collected a total of four representative samples (SS1 in Area A, SS2 in Area B, SS3 in Area C, and SS4 in Area D). WSDOT collected samples in areas where large stains were evident (A and C), numerous small stains (D), unusual odors (A-D), and solid wastes like trash and debris were abandoned (B and D). The soil samples were analyzed using the following methods:

- Northwest Total Petroleum Hydrocarbons - Method NWTPH – Benzene, Toluene, Ethylbenzene, and Xylene (BTEX)
- Diesel and Heavy Oil - Method NWTPH (Dx)
- SVOCs - EPA Method 8270
- RCRA 8 Metals - EPA Method 6010

Analytical results of one the four soil samples indicated that that lube range petroleum hydrocarbons were detected at a concentration greater than current Model Toxics Control Act (MTCA) Method A cleanup level (CUL) of 2,000 milligrams per kilograms (SS-3 31,000 mg/kg). WSDOT considered this asphalt. The analytical results of another sample indicated that arsenic were detected at a concentration greater the current MTCA Method A CUL of 20 milligrams per kilograms (SS-2 35 mg/kg). The remaining analytes sampled were either non-detect or below MTCA Method A cleanup levels. Based upon the sample results, field remediation of contaminated soils was needed in area B (SS-2) and C (SS-3).

In April 2017, WSDOT contracted Marine Vacuum Incorporated (Marine Vac.) to cleanup areas on the Site with trash, debris, and soils with notable detectable analytes below MTCA Method A CULs.

In January 2018 A-1 Towing, the lessee, graded and displaced Site soils. Some soils were contaminated. The soils were displaced from previously sampled areas A and B. There also was a new stockpile located just north of area D.

On March 15, 2018, WSDOT Headquarters (HQ) HazMat Program collected four soil samples (SS1a, SS2a, SS3a, and SS4a) on the Site in areas previously defined as A, B, and in a new location. The HazMat Program collected a sample if there was an indication of grading with known or suspected soil contamination, or where displaced/graded soils appeared to be consolidated into a stock pile generated by the lessee. Areas C and D did not have these indicators and did not warrant additional sampling. The stock pile location is just north of area D and was estimated to contain approximately 25 cubic yards of material.

The analytical results from the March 2018 investigative sampling, individual chemical constituents detected in each soil sample were analyzed using:

- NWTPH-Gx,
- NWTPH-Dx,
- Volatiles EPA Method 8260C,
- Semi-Volatiles Method 8270D,
- Total metals EPA Method 6010D.

No sample results were above MTCA Method A cleanup standards.

### **3.0 FIELD INVESTIGATION**

On September 21, 2018, HQ HazMat collected two soil confirmation samples (Fcl1 and Fcl2) using a pre-cleaned stainless steel spoon. HQ HazMat decontaminated the stainless steel spoon with a phosphate free cleaner (Alconox) and rinsed it with distilled water between the confirmation sample locations. HQ HazMat collected the two confirmation samples in areas remediated where:

- Soils were previously identified in concentrations that exceeded MTCA Method A cleanup levels (CULs) for lube oil; and
- Numerous stains were located throughout the area.

The Fcl1 confirmation sample was collected at the bottom from the center of where the crushed asphalt was removed.

The depth of the excavation trench for the 2nd confirmation sample could not safely be accessed. Therefore, the Fcl2 confirmation sample was collected from soil excavated with a backhoe bucket taken from the center of the bottom of the excavation.

### **4.0 ANALYTICAL SAMPLING RESULTS**

The analytical data was reviewed and deemed acceptable. The individual chemical constituents in each soil sample was evaluated against the MTCA Method A cleanup levels for unrestricted land use. The sample results were below MTCA Method A cleanup levels. All chemical results from each of the soil sample events are presented in its entirety in Appendix 3. All of the soil samples were delivered to OnSite Environmental Incorporated located in Redmond, WA.

The analytical results from the September 2018 confirmation samples, individual constituents detected in each soil sample was evaluated against:

- NWTPH-Dx
- Semi-Volatiles EPA Method 8260C
- MTCA Metals

The laboratory analytical results from the two confirmation samples were either non-detect or at concentrations below MTCA Method A CULs.

## **5.0 FIELD ACTIVITIES**

On September 20 and 21, 2018, Marine Vac. conducted excavation activities to address lube oil concentrations exceeding MTCA Method A CULs. Additionally, Marine Vac. excavated an area to remediate suspected contaminated materials (See Appendix 1 areas C and D). During site cleanup activities, excavated soils were segregated based on field screening results. Soils included in the direct haul to Marine Vac. for disposal as contaminated included soils based up sight, olfactory indicators, and/or had Photoionization Indication Detector (PID) levels above 10 ppm. The non-suspect soils were placed off to the side in a designated area lined and covered with 6 mil polyethylene sheeting. On September 24, 2018, Marine Vac. filled and graded the excavation areas with the existing non-suspect soils.

A total of about 8.4 cubic yards of suspected and/or confirmed contaminated material was excavated and transported for disposal by Marine Vac. Waste disposal receipts are located in Appendix 4.

No groundwater was encountered during cleanup activities.

## **6.0 SUMMARY AND CONCLUSIONS**

WSDOT completed soil investigation and excavation activities within the Site. WSDOT investigated the soils on the Site at the end of an existing long term lease and also after the new lessee displaced and graded soils. Confirmed and suspect contamination existed on the site. Approximately 8.4 cubic yards of confirmed and suspected contaminated soil was disposed. Based on the laboratory analytical results from the confirmation sample locations, it was determined that no contamination remains on the property.

Based on the cleanup activities and the laboratory analytical results, it appears that the subsequent environmental investigation removed the confirmed or suspected contamination; therefore, with no immediate threat to human health or the environment no additional environmental investigation appears warranted at this time. WSDOT recommends Washington State Department of Ecology to accept this Site as a No Further Action.

**APPENDIX 1 – VICINITY MAP and SAMPLE POINTS & BACKGROUND**

VICINITY MAP WITH SAMPLE POINTS



BACKGROUND AREA



○ Sample locations

## **APPENDIX 2 SUMMARY TABLES**

BACKGROUND CONDITION

Sample Collected (10/2016)	Gasoline mg/kg	Diesel mg/kg	Lube Oil mg/kg	Benzene mg/kg	Chromium mg/kg	Arsenic mg/kg
SS 1	ND	ND	1900	ND	48	ND
SS 2	ND	ND	220	0.022 (0.03)	46	<b>35</b>
SS 3	ND	ND	<b>31,000</b>	ND	22	ND
SS 4 MTCA Method A Cleanup levels mg/kg	ND 30 mg/kg with Benzene 100 mg/kg without Benzene	ND 2000 mg/kg	60 2000 mg/kg	ND 0.03	42 100 mg/kg(replace d by values for III (2000) and VI (19))	ND 20 mg/kg

MARCH 2018 INVESTIGATION

Sample Collected (10/2016)	Gasoline mg/kg	Diesel mg/kg	Lube Oil mg/kg	Benzene mg/kg	Chromium mg/kg	Arsenic mg/kg
SS 1	ND	ND	82	ND	41	ND
SS 2	ND	58	370	ND	52	ND
SS 3 (STOCKPILE)	ND	ND	470	ND	47	ND
SS 4 (STOCKPILE)	ND	ND	260	ND	60	ND
MTCA Method A Cleanup levels mg/kg	30 mg/kg with Benzene 100 mg/kg without Benzene	2000 mg/kg	2000 mg/kg	0.03 mg/kg	100 mg/kg(replace d by values for III (2000) and VI (19))	5 mg/kg

CONFIRMATION SAMPLE

<u>Sample Collected</u> 8/21/2018	<u>Lube Oil</u>
<u>FC1-1</u> <u>(Baseline ss3)</u>	<u>590</u>
<u>FC1-2</u> <u>(Baseline ss4)</u>	<u>ND</u>
<u>MTCA</u> <u>Method A</u> <u>cleanup level</u> <u>mg/kg</u>	<u>2000 mg/kg</u>

**APPENDIX 3 LABORATORY REPORTS**



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

November 8, 2016

Patrick Svoboda  
WSDOT  
15700 Dayton Avenue North  
NB82-138  
P.O. Box 330310  
Seattle, WA 98133-9710

Re: Analytical Data for Project SR 509 - 12th Street  
Laboratory Reference No. 1610-295

Dear Pat:

Enclosed are the analytical results and associated quality control data for samples submitted on October 26, 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: November 8, 2016  
Samples Submitted: October 26, 2016  
Laboratory Reference: 1610-295  
Project: SR 509 - 12th Street

### Case Narrative

Samples were collected on October 25, 2016 and received by the laboratory on October 26, 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 within 48 hours of sample collection. They were stored in a freezer at between -7°C and -20°C until extraction or analysis.

#### Semivolatiles EPA 8270D/SIM Analysis

Sample SS 4 had one surrogate recovery out of control limits. This was due to a large co-eluting non-target analyte peak, tentatively identified as 2-butoxy ethanol by spectral library search.

The Spike Blank/Spike Blank Duplicate had one recovery slightly above control limits, due to a small upward bias in the instruments' calibration and the sample was non-detect for this analyte. No further action was taken.

**Please note that any other QA/QC issues associated with these extractions and analyses will be indicated with a footnote reference and discussed in detail on the Data Qualifier page.**



Date of Report: November 8, 2016  
 Samples Submitted: October 26, 2016  
 Laboratory Reference: 1610-295  
 Project: SR 509 - 12th Street

### NWTPH-Gx/BTEX

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>Soil Sample 1</b>					
Laboratory ID:	10-295-01					
Benzene	ND	0.020	EPA 8021B	10-28-16	10-31-16	
Toluene	ND	0.051	EPA 8021B	10-28-16	10-31-16	
Ethyl Benzene	ND	0.051	EPA 8021B	10-28-16	10-31-16	
m,p-Xylene	ND	0.051	EPA 8021B	10-28-16	10-31-16	
o-Xylene	ND	0.051	EPA 8021B	10-28-16	10-31-16	
Gasoline	ND	5.1	NWTPH-Gx	10-28-16	10-31-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	91	63-124				
<b>Client ID:</b>	<b>Soil Sample 2</b>					
Laboratory ID:	10-295-02					
Benzene	0.022	0.020	EPA 8021B	10-28-16	11-1-16	
Toluene	ND	0.094	EPA 8021B	10-28-16	11-1-16	
Ethyl Benzene	ND	0.094	EPA 8021B	10-28-16	11-1-16	
m,p-Xylene	ND	0.094	EPA 8021B	10-28-16	11-1-16	
o-Xylene	ND	0.094	EPA 8021B	10-28-16	11-1-16	
Gasoline	ND	9.4	NWTPH-Gx	10-28-16	11-1-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	80	63-124				
<b>Client ID:</b>	<b>Soil Sample 3</b>					
Laboratory ID:	10-295-03					
Benzene	ND	0.020	EPA 8021B	10-28-16	11-1-16	
Toluene	ND	0.048	EPA 8021B	10-28-16	11-1-16	
Ethyl Benzene	ND	0.048	EPA 8021B	10-28-16	11-1-16	
m,p-Xylene	ND	0.048	EPA 8021B	10-28-16	11-1-16	
o-Xylene	ND	0.048	EPA 8021B	10-28-16	11-1-16	
Gasoline	ND	4.8	NWTPH-Gx	10-28-16	11-1-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	86	63-124				



Date of Report: November 8, 2016  
 Samples Submitted: October 26, 2016  
 Laboratory Reference: 1610-295  
 Project: SR 509 - 12th Street

**NWTPH-Gx/BTEX**

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>SS 4</b>					
Laboratory ID:	10-295-04					
Benzene	<b>ND</b>	0.020	EPA 8021B	10-28-16	11-1-16	
Toluene	<b>ND</b>	0.055	EPA 8021B	10-28-16	11-1-16	
Ethyl Benzene	<b>ND</b>	0.055	EPA 8021B	10-28-16	11-1-16	
m,p-Xylene	<b>ND</b>	0.055	EPA 8021B	10-28-16	11-1-16	
o-Xylene	<b>ND</b>	0.055	EPA 8021B	10-28-16	11-1-16	
Gasoline	<b>ND</b>	5.5	NWTPH-Gx	10-28-16	11-1-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	<i>89</i>	<i>63-124</i>				



Date of Report: November 8, 2016  
 Samples Submitted: October 26, 2016  
 Laboratory Reference: 1610-295  
 Project: SR 509 - 12th Street

**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:	MB1028S1					
Benzene	ND	0.020	EPA 8021B	10-28-16	10-28-16	
Toluene	ND	0.050	EPA 8021B	10-28-16	10-28-16	
Ethyl Benzene	ND	0.050	EPA 8021B	10-28-16	10-28-16	
m,p-Xylene	ND	0.050	EPA 8021B	10-28-16	10-28-16	
o-Xylene	ND	0.050	EPA 8021B	10-28-16	10-28-16	
Gasoline	ND	5.0	NWTPH-Gx	10-28-16	10-28-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	90	63-124				

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>								
Laboratory ID:	10-293-16							
	ORIG	DUP						
Benzene	ND	ND	NA	NA	NA	NA	30	
Toluene	ND	ND	NA	NA	NA	NA	30	
Ethyl Benzene	ND	ND	NA	NA	NA	NA	30	
m,p-Xylene	ND	ND	NA	NA	NA	NA	30	
o-Xylene	ND	ND	NA	NA	NA	NA	30	
Gasoline	ND	ND	NA	NA	NA	NA	30	
<i>Surrogate:</i>								
<i>Fluorobenzene</i>				100	95	63-124		

**SPIKE BLANKS**

Laboratory ID:	SB1028S1								
	SB	SBD	SB	SBD	SB	SBD			
Benzene	0.929	0.914	1.00	1.00	93	91	70-124	2	12
Toluene	0.935	0.927	1.00	1.00	94	93	73-119	1	12
Ethyl Benzene	0.953	0.942	1.00	1.00	95	94	74-117	1	12
m,p-Xylene	0.905	0.907	1.00	1.00	91	91	75-117	0	13
o-Xylene	0.943	0.932	1.00	1.00	94	93	75-116	1	12
<i>Surrogate:</i>									
<i>Fluorobenzene</i>					95	91	63-124		



Date of Report: November 8, 2016  
 Samples Submitted: October 26, 2016  
 Laboratory Reference: 1610-295  
 Project: SR 509 - 12th Street

### NWTPH-Dx

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>Soil Sample 1</b>					
Laboratory ID:	10-295-01					
Diesel Range Organics	<b>ND</b>	550	NWTPH-Dx	10-28-16	10-28-16	U1
Lube Oil	<b>1900</b>	56	NWTPH-Dx	10-28-16	10-28-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	84	50-150				
<b>Client ID:</b>	<b>Soil Sample 2</b>					
Laboratory ID:	10-295-02					
Diesel Range Organics	<b>ND</b>	46	NWTPH-Dx	10-28-16	10-28-16	U1
Lube Oil Range Organics	<b>220</b>	70	NWTPH-Dx	10-28-16	10-28-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	107	50-150				
<b>Client ID:</b>	<b>Soil Sample 3</b>					
Laboratory ID:	10-295-03					
Diesel Range Organics	<b>ND</b>	6000	NWTPH-Dx	10-28-16	10-31-16	U1
Lube Oil	<b>31000</b>	1100	NWTPH-Dx	10-28-16	10-31-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	---	50-150				S
<b>Client ID:</b>	<b>SS 4</b>					
Laboratory ID:	10-295-04					
Diesel Range Organics	<b>ND</b>	30	NWTPH-Dx	10-28-16	10-28-16	
Lube Oil Range Organics	<b>60</b>	59	NWTPH-Dx	10-28-16	10-28-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	99	50-150				



Date of Report: November 8, 2016  
 Samples Submitted: October 26, 2016  
 Laboratory Reference: 1610-295  
 Project: SR 509 - 12th Street

**NWTPH-Dx  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1028S1					
Diesel Range Organics	<b>ND</b>	25	NWTPH-Dx	10-28-16	10-28-16	
Lube Oil Range Organics	<b>ND</b>	50	NWTPH-Dx	10-28-16	10-28-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	97	50-150				

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>								
Laboratory ID:	10-295-01							
	ORIG	DUP						
Diesel Range	<b>ND</b>	<b>ND</b>	NA	NA	NA	NA	NA	U1
Lube Oil	<b>1670</b>	<b>1670</b>	NA	NA	NA	0	NA	
<i>Surrogate:</i>								
<i>o-Terphenyl</i>			84	100	50-150			



Date of Report: November 8, 2016  
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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>Soil Sample 1</b>					
<b>Laboratory ID:</b>	10-295-01					
n-Nitrosodimethylamine	ND	0.037	EPA 8270D	11-2-16	11-3-16	
Pyridine	ND	0.37	EPA 8270D	11-2-16	11-3-16	
Phenol	ND	0.037	EPA 8270D	11-2-16	11-3-16	
Aniline	ND	0.19	EPA 8270D	11-2-16	11-3-16	
bis(2-Chloroethyl)ether	ND	0.037	EPA 8270D	11-2-16	11-3-16	
2-Chlorophenol	ND	0.037	EPA 8270D	11-2-16	11-3-16	
1,3-Dichlorobenzene	ND	0.037	EPA 8270D	11-2-16	11-3-16	
1,4-Dichlorobenzene	ND	0.037	EPA 8270D	11-2-16	11-3-16	
Benzyl alcohol	ND	0.19	EPA 8270D	11-2-16	11-3-16	
1,2-Dichlorobenzene	ND	0.037	EPA 8270D	11-2-16	11-3-16	
2-Methylphenol (o-Cresol)	ND	0.037	EPA 8270D	11-2-16	11-3-16	
bis(2-Chloroisopropyl)ether	ND	0.037	EPA 8270D	11-2-16	11-3-16	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.037	EPA 8270D	11-2-16	11-3-16	
n-Nitroso-di-n-propylamine	ND	0.037	EPA 8270D	11-2-16	11-3-16	
Hexachloroethane	ND	0.037	EPA 8270D	11-2-16	11-3-16	
Nitrobenzene	ND	0.037	EPA 8270D	11-2-16	11-3-16	
Isophorone	ND	0.037	EPA 8270D	11-2-16	11-3-16	
2-Nitrophenol	ND	0.037	EPA 8270D	11-2-16	11-3-16	
2,4-Dimethylphenol	ND	0.037	EPA 8270D	11-2-16	11-3-16	
bis(2-Chloroethoxy)methane	ND	0.037	EPA 8270D	11-2-16	11-3-16	
2,4-Dichlorophenol	ND	0.037	EPA 8270D	11-2-16	11-3-16	
1,2,4-Trichlorobenzene	ND	0.037	EPA 8270D	11-2-16	11-3-16	
Naphthalene	ND	0.0075	EPA 8270D/SIM	11-2-16	11-3-16	
4-Chloroaniline	ND	0.19	EPA 8270D	11-2-16	11-3-16	
Hexachlorobutadiene	ND	0.037	EPA 8270D	11-2-16	11-3-16	
4-Chloro-3-methylphenol	ND	0.037	EPA 8270D	11-2-16	11-3-16	
2-Methylnaphthalene	<b>0.0099</b>	0.0075	EPA 8270D/SIM	11-2-16	11-3-16	
1-Methylnaphthalene	ND	0.0075	EPA 8270D/SIM	11-2-16	11-3-16	
Hexachlorocyclopentadiene	ND	0.037	EPA 8270D	11-2-16	11-3-16	
2,4,6-Trichlorophenol	ND	0.037	EPA 8270D	11-2-16	11-3-16	
2,3-Dichloroaniline	ND	0.037	EPA 8270D	11-2-16	11-3-16	
2,4,5-Trichlorophenol	ND	0.037	EPA 8270D	11-2-16	11-3-16	
2-Chloronaphthalene	ND	0.037	EPA 8270D	11-2-16	11-3-16	
2-Nitroaniline	ND	0.037	EPA 8270D	11-2-16	11-3-16	
1,4-Dinitrobenzene	ND	0.037	EPA 8270D	11-2-16	11-3-16	
Dimethylphthalate	ND	0.037	EPA 8270D	11-2-16	11-3-16	
1,3-Dinitrobenzene	ND	0.037	EPA 8270D	11-2-16	11-3-16	
2,6-Dinitrotoluene	ND	0.037	EPA 8270D	11-2-16	11-3-16	
1,2-Dinitrobenzene	ND	0.037	EPA 8270D	11-2-16	11-3-16	
Acenaphthylene	ND	0.0075	EPA 8270D/SIM	11-2-16	11-3-16	
3-Nitroaniline	ND	0.037	EPA 8270D	11-2-16	11-3-16	



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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>Soil Sample 1</b>					
Laboratory ID:	10-295-01					
2,4-Dinitrophenol	ND	0.19	EPA 8270D	11-2-16	11-3-16	
Acenaphthene	ND	0.0075	EPA 8270D/SIM	11-2-16	11-3-16	
4-Nitrophenol	ND	0.037	EPA 8270D	11-2-16	11-3-16	
2,4-Dinitrotoluene	ND	0.037	EPA 8270D	11-2-16	11-3-16	
Dibenzofuran	ND	0.037	EPA 8270D	11-2-16	11-3-16	
2,3,5,6-Tetrachlorophenol	ND	0.037	EPA 8270D	11-2-16	11-3-16	
2,3,4,6-Tetrachlorophenol	ND	0.037	EPA 8270D	11-2-16	11-3-16	
Diethylphthalate	ND	0.19	EPA 8270D	11-2-16	11-3-16	
4-Chlorophenyl-phenylether	ND	0.037	EPA 8270D	11-2-16	11-3-16	
4-Nitroaniline	ND	0.037	EPA 8270D	11-2-16	11-3-16	
Fluorene	<b>0.0088</b>	0.0075	EPA 8270D/SIM	11-2-16	11-3-16	
4,6-Dinitro-2-methylphenol	ND	0.19	EPA 8270D	11-2-16	11-3-16	
n-Nitrosodiphenylamine	ND	0.037	EPA 8270D	11-2-16	11-3-16	
1,2-Diphenylhydrazine	ND	0.037	EPA 8270D	11-2-16	11-3-16	
4-Bromophenyl-phenylether	ND	0.037	EPA 8270D	11-2-16	11-3-16	
Hexachlorobenzene	ND	0.037	EPA 8270D	11-2-16	11-3-16	
Pentachlorophenol	ND	0.19	EPA 8270D	11-2-16	11-3-16	
Phenanthrene	<b>0.014</b>	0.0075	EPA 8270D/SIM	11-2-16	11-3-16	
Anthracene	ND	0.0075	EPA 8270D/SIM	11-2-16	11-3-16	
Carbazole	ND	0.037	EPA 8270D	11-2-16	11-3-16	
Di-n-butylphthalate	ND	0.19	EPA 8270D	11-2-16	11-3-16	
Fluoranthene	<b>0.011</b>	0.0075	EPA 8270D/SIM	11-2-16	11-3-16	
Benzidine	ND	0.37	EPA 8270D	11-2-16	11-3-16	
Pyrene	<b>0.018</b>	0.0075	EPA 8270D/SIM	11-2-16	11-3-16	
Butylbenzylphthalate	<b>0.053</b>	0.037	EPA 8270D	11-2-16	11-3-16	
bis-2-Ethylhexyladipate	ND	0.037	EPA 8270D	11-2-16	11-3-16	
3,3'-Dichlorobenzidine	ND	0.19	EPA 8270D	11-2-16	11-3-16	
Benzo[a]anthracene	ND	0.015	EPA 8270D/SIM	11-2-16	11-4-16	
Chrysene	ND	0.015	EPA 8270D/SIM	11-2-16	11-4-16	
bis(2-Ethylhexyl)phthalate	<b>0.26</b>	0.037	EPA 8270D	11-2-16	11-3-16	
Di-n-octylphthalate	ND	0.037	EPA 8270D	11-2-16	11-3-16	
Benzo[b]fluoranthene	ND	0.015	EPA 8270D/SIM	11-2-16	11-4-16	
Benzo(j,k)fluoranthene	ND	0.015	EPA 8270D/SIM	11-2-16	11-4-16	
Benzo[a]pyrene	ND	0.015	EPA 8270D/SIM	11-2-16	11-4-16	
Indeno[1,2,3-cd]pyrene	ND	0.015	EPA 8270D/SIM	11-2-16	11-4-16	
Dibenz[a,h]anthracene	ND	0.015	EPA 8270D/SIM	11-2-16	11-4-16	
Benzo[g,h,i]perylene	ND	0.015	EPA 8270D/SIM	11-2-16	11-4-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorophenol</i>	<i>51</i>	<i>18 - 109</i>				
<i>Phenol-d6</i>	<i>63</i>	<i>25 - 111</i>				
<i>Nitrobenzene-d5</i>	<i>64</i>	<i>22 - 113</i>				
<i>2-Fluorobiphenyl</i>	<i>82</i>	<i>30 - 114</i>				
<i>2,4,6-Tribromophenol</i>	<i>88</i>	<i>22 - 116</i>				
<i>Terphenyl-d14</i>	<i>78</i>	<i>33 - 114</i>				



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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>Soil Sample 2</b>					
<b>Laboratory ID:</b>	10-295-02					
n-Nitrosodimethylamine	ND	0.046	EPA 8270D	11-2-16	11-3-16	
Pyridine	ND	0.46	EPA 8270D	11-2-16	11-3-16	
Phenol	ND	0.046	EPA 8270D	11-2-16	11-3-16	
Aniline	ND	0.23	EPA 8270D	11-2-16	11-3-16	
bis(2-Chloroethyl)ether	ND	0.046	EPA 8270D	11-2-16	11-3-16	
2-Chlorophenol	ND	0.046	EPA 8270D	11-2-16	11-3-16	
1,3-Dichlorobenzene	ND	0.046	EPA 8270D	11-2-16	11-3-16	
1,4-Dichlorobenzene	ND	0.046	EPA 8270D	11-2-16	11-3-16	
Benzyl alcohol	<b>0.30</b>	0.23	EPA 8270D	11-2-16	11-3-16	
1,2-Dichlorobenzene	ND	0.046	EPA 8270D	11-2-16	11-3-16	
2-Methylphenol (o-Cresol)	ND	0.046	EPA 8270D	11-2-16	11-3-16	
bis(2-Chloroisopropyl)ether	ND	0.046	EPA 8270D	11-2-16	11-3-16	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.046	EPA 8270D	11-2-16	11-3-16	
n-Nitroso-di-n-propylamine	ND	0.046	EPA 8270D	11-2-16	11-3-16	
Hexachloroethane	ND	0.046	EPA 8270D	11-2-16	11-3-16	
Nitrobenzene	ND	0.046	EPA 8270D	11-2-16	11-3-16	
Isophorone	ND	0.046	EPA 8270D	11-2-16	11-3-16	
2-Nitrophenol	ND	0.046	EPA 8270D	11-2-16	11-3-16	
2,4-Dimethylphenol	ND	0.046	EPA 8270D	11-2-16	11-3-16	
bis(2-Chloroethoxy)methane	ND	0.046	EPA 8270D	11-2-16	11-3-16	
2,4-Dichlorophenol	ND	0.046	EPA 8270D	11-2-16	11-3-16	
1,2,4-Trichlorobenzene	ND	0.046	EPA 8270D	11-2-16	11-3-16	
Naphthalene	<b>0.044</b>	0.0093	EPA 8270D/SIM	11-2-16	11-3-16	
4-Chloroaniline	ND	0.23	EPA 8270D	11-2-16	11-3-16	
Hexachlorobutadiene	ND	0.046	EPA 8270D	11-2-16	11-3-16	
4-Chloro-3-methylphenol	ND	0.046	EPA 8270D	11-2-16	11-3-16	
2-Methylnaphthalene	<b>0.026</b>	0.0093	EPA 8270D/SIM	11-2-16	11-3-16	
1-Methylnaphthalene	<b>0.014</b>	0.0093	EPA 8270D/SIM	11-2-16	11-3-16	
Hexachlorocyclopentadiene	ND	0.046	EPA 8270D	11-2-16	11-3-16	
2,4,6-Trichlorophenol	ND	0.046	EPA 8270D	11-2-16	11-3-16	
2,3-Dichloroaniline	ND	0.046	EPA 8270D	11-2-16	11-3-16	
2,4,5-Trichlorophenol	ND	0.046	EPA 8270D	11-2-16	11-3-16	
2-Chloronaphthalene	ND	0.046	EPA 8270D	11-2-16	11-3-16	
2-Nitroaniline	ND	0.046	EPA 8270D	11-2-16	11-3-16	
1,4-Dinitrobenzene	ND	0.046	EPA 8270D	11-2-16	11-3-16	
Dimethylphthalate	ND	0.046	EPA 8270D	11-2-16	11-3-16	
1,3-Dinitrobenzene	ND	0.046	EPA 8270D	11-2-16	11-3-16	
2,6-Dinitrotoluene	ND	0.046	EPA 8270D	11-2-16	11-3-16	
1,2-Dinitrobenzene	ND	0.046	EPA 8270D	11-2-16	11-3-16	
Acenaphthylene	ND	0.0093	EPA 8270D/SIM	11-2-16	11-3-16	
3-Nitroaniline	ND	0.046	EPA 8270D	11-2-16	11-3-16	



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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>Soil Sample 2</b>					
Laboratory ID:	10-295-02					
2,4-Dinitrophenol	ND	0.23	EPA 8270D	11-2-16	11-3-16	
Acenaphthene	ND	0.0093	EPA 8270D/SIM	11-2-16	11-3-16	
4-Nitrophenol	ND	0.046	EPA 8270D	11-2-16	11-3-16	
2,4-Dinitrotoluene	ND	0.046	EPA 8270D	11-2-16	11-3-16	
Dibenzofuran	ND	0.046	EPA 8270D	11-2-16	11-3-16	
2,3,5,6-Tetrachlorophenol	ND	0.046	EPA 8270D	11-2-16	11-3-16	
2,3,4,6-Tetrachlorophenol	ND	0.046	EPA 8270D	11-2-16	11-3-16	
Diethylphthalate	ND	0.23	EPA 8270D	11-2-16	11-3-16	
4-Chlorophenyl-phenylether	ND	0.046	EPA 8270D	11-2-16	11-3-16	
4-Nitroaniline	ND	0.046	EPA 8270D	11-2-16	11-3-16	
Fluorene	ND	0.0093	EPA 8270D/SIM	11-2-16	11-3-16	
4,6-Dinitro-2-methylphenol	ND	0.23	EPA 8270D	11-2-16	11-3-16	
n-Nitrosodiphenylamine	ND	0.046	EPA 8270D	11-2-16	11-3-16	
1,2-Diphenylhydrazine	ND	0.046	EPA 8270D	11-2-16	11-3-16	
4-Bromophenyl-phenylether	ND	0.046	EPA 8270D	11-2-16	11-3-16	
Hexachlorobenzene	ND	0.046	EPA 8270D	11-2-16	11-3-16	
Pentachlorophenol	ND	0.23	EPA 8270D	11-2-16	11-3-16	
Phenanthrene	0.012	0.0093	EPA 8270D/SIM	11-2-16	11-3-16	
Anthracene	ND	0.0093	EPA 8270D/SIM	11-2-16	11-3-16	
Carbazole	ND	0.046	EPA 8270D	11-2-16	11-3-16	
Di-n-butylphthalate	ND	0.23	EPA 8270D	11-2-16	11-3-16	
Fluoranthene	0.0098	0.0093	EPA 8270D/SIM	11-2-16	11-3-16	
Benzidine	ND	0.46	EPA 8270D	11-2-16	11-3-16	
Pyrene	ND	0.0093	EPA 8270D/SIM	11-2-16	11-3-16	
Butylbenzylphthalate	ND	0.046	EPA 8270D	11-2-16	11-3-16	
bis-2-Ethylhexyladipate	ND	0.046	EPA 8270D	11-2-16	11-3-16	
3,3'-Dichlorobenzidine	ND	0.23	EPA 8270D	11-2-16	11-3-16	
Benzo[a]anthracene	ND	0.0093	EPA 8270D/SIM	11-2-16	11-3-16	
Chrysene	ND	0.0093	EPA 8270D/SIM	11-2-16	11-3-16	
bis(2-Ethylhexyl)phthalate	0.059	0.046	EPA 8270D	11-2-16	11-3-16	
Di-n-octylphthalate	ND	0.046	EPA 8270D	11-2-16	11-3-16	
Benzo[b]fluoranthene	ND	0.0093	EPA 8270D/SIM	11-2-16	11-3-16	
Benzo(j,k)fluoranthene	ND	0.0093	EPA 8270D/SIM	11-2-16	11-3-16	
Benzo[a]pyrene	ND	0.0093	EPA 8270D/SIM	11-2-16	11-3-16	
Indeno[1,2,3-cd]pyrene	ND	0.0093	EPA 8270D/SIM	11-2-16	11-3-16	
Dibenz[a,h]anthracene	ND	0.0093	EPA 8270D/SIM	11-2-16	11-3-16	
Benzo[g,h,i]perylene	ND	0.0093	EPA 8270D/SIM	11-2-16	11-3-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	28	18 - 109				
Phenol-d6	29	25 - 111				
Nitrobenzene-d5	49	22 - 113				
2-Fluorobiphenyl	57	30 - 114				
2,4,6-Tribromophenol	51	22 - 116				
Terphenyl-d14	57	33 - 114				



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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>Soil Sample 3</b>					
<b>Laboratory ID:</b>	10-295-03					
n-Nitrosodimethylamine	ND	0.18	EPA 8270D	11-2-16	11-4-16	
Pyridine	ND	1.8	EPA 8270D	11-2-16	11-4-16	
Phenol	ND	0.18	EPA 8270D	11-2-16	11-4-16	
Aniline	ND	0.92	EPA 8270D	11-2-16	11-4-16	
bis(2-Chloroethyl)ether	ND	0.18	EPA 8270D	11-2-16	11-4-16	
2-Chlorophenol	ND	0.18	EPA 8270D	11-2-16	11-4-16	
1,3-Dichlorobenzene	ND	0.18	EPA 8270D	11-2-16	11-4-16	
1,4-Dichlorobenzene	ND	0.18	EPA 8270D	11-2-16	11-4-16	
Benzyl alcohol	ND	0.92	EPA 8270D	11-2-16	11-4-16	
1,2-Dichlorobenzene	ND	0.18	EPA 8270D	11-2-16	11-4-16	
2-Methylphenol (o-Cresol)	ND	0.18	EPA 8270D	11-2-16	11-4-16	
bis(2-Chloroisopropyl)ether	ND	0.18	EPA 8270D	11-2-16	11-4-16	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.18	EPA 8270D	11-2-16	11-4-16	
n-Nitroso-di-n-propylamine	ND	0.18	EPA 8270D	11-2-16	11-4-16	
Hexachloroethane	ND	0.18	EPA 8270D	11-2-16	11-4-16	
Nitrobenzene	ND	0.18	EPA 8270D	11-2-16	11-4-16	
Isophorone	ND	0.18	EPA 8270D	11-2-16	11-4-16	
2-Nitrophenol	ND	0.18	EPA 8270D	11-2-16	11-4-16	
2,4-Dimethylphenol	ND	0.18	EPA 8270D	11-2-16	11-4-16	
bis(2-Chloroethoxy)methane	ND	0.18	EPA 8270D	11-2-16	11-4-16	
2,4-Dichlorophenol	ND	0.18	EPA 8270D	11-2-16	11-4-16	
1,2,4-Trichlorobenzene	ND	0.18	EPA 8270D	11-2-16	11-4-16	
Naphthalene	0.011	0.0074	EPA 8270D/SIM	11-2-16	11-3-16	
4-Chloroaniline	ND	0.92	EPA 8270D	11-2-16	11-4-16	
Hexachlorobutadiene	ND	0.18	EPA 8270D	11-2-16	11-4-16	
4-Chloro-3-methylphenol	ND	0.18	EPA 8270D	11-2-16	11-4-16	
2-Methylnaphthalene	0.042	0.0074	EPA 8270D/SIM	11-2-16	11-3-16	
1-Methylnaphthalene	0.030	0.0074	EPA 8270D/SIM	11-2-16	11-3-16	
Hexachlorocyclopentadiene	ND	0.18	EPA 8270D	11-2-16	11-4-16	
2,4,6-Trichlorophenol	ND	0.18	EPA 8270D	11-2-16	11-4-16	
2,3-Dichloroaniline	ND	0.18	EPA 8270D	11-2-16	11-4-16	
2,4,5-Trichlorophenol	ND	0.18	EPA 8270D	11-2-16	11-4-16	
2-Chloronaphthalene	ND	0.18	EPA 8270D	11-2-16	11-4-16	
2-Nitroaniline	ND	0.18	EPA 8270D	11-2-16	11-4-16	
1,4-Dinitrobenzene	ND	0.18	EPA 8270D	11-2-16	11-4-16	
Dimethylphthalate	ND	0.18	EPA 8270D	11-2-16	11-4-16	
1,3-Dinitrobenzene	ND	0.18	EPA 8270D	11-2-16	11-4-16	
2,6-Dinitrotoluene	ND	0.18	EPA 8270D	11-2-16	11-4-16	
1,2-Dinitrobenzene	ND	0.18	EPA 8270D	11-2-16	11-4-16	
Acenaphthylene	ND	0.15	EPA 8270D/SIM	11-2-16	11-4-16	
3-Nitroaniline	ND	0.18	EPA 8270D	11-2-16	11-4-16	



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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>Soil Sample 3</b>					
Laboratory ID:	10-295-03					
2,4-Dinitrophenol	ND	0.92	EPA 8270D	11-2-16	11-4-16	
Acenaphthene	ND	0.15	EPA 8270D/SIM	11-2-16	11-4-16	
4-Nitrophenol	ND	0.18	EPA 8270D	11-2-16	11-4-16	
2,4-Dinitrotoluene	ND	0.18	EPA 8270D	11-2-16	11-4-16	
Dibenzofuran	ND	0.18	EPA 8270D	11-2-16	11-4-16	
2,3,5,6-Tetrachlorophenol	ND	0.18	EPA 8270D	11-2-16	11-4-16	
2,3,4,6-Tetrachlorophenol	ND	0.18	EPA 8270D	11-2-16	11-4-16	
Diethylphthalate	ND	0.92	EPA 8270D	11-2-16	11-4-16	
4-Chlorophenyl-phenylether	ND	0.18	EPA 8270D	11-2-16	11-4-16	
4-Nitroaniline	ND	0.18	EPA 8270D	11-2-16	11-4-16	
Fluorene	0.15	0.15	EPA 8270D/SIM	11-2-16	11-4-16	
4,6-Dinitro-2-methylphenol	ND	0.92	EPA 8270D	11-2-16	11-4-16	
n-Nitrosodiphenylamine	ND	0.18	EPA 8270D	11-2-16	11-4-16	
1,2-Diphenylhydrazine	ND	0.18	EPA 8270D	11-2-16	11-4-16	
4-Bromophenyl-phenylether	ND	0.18	EPA 8270D	11-2-16	11-4-16	
Hexachlorobenzene	ND	0.18	EPA 8270D	11-2-16	11-4-16	
Pentachlorophenol	ND	0.92	EPA 8270D	11-2-16	11-4-16	
Phenanthrene	ND	0.15	EPA 8270D/SIM	11-2-16	11-4-16	
Anthracene	ND	0.15	EPA 8270D/SIM	11-2-16	11-4-16	
Carbazole	ND	0.18	EPA 8270D	11-2-16	11-4-16	
Di-n-butylphthalate	ND	0.92	EPA 8270D	11-2-16	11-4-16	
Fluoranthene	0.18	0.15	EPA 8270D/SIM	11-2-16	11-4-16	
Benzidine	ND	1.8	EPA 8270D	11-2-16	11-4-16	
Pyrene	0.30	0.15	EPA 8270D/SIM	11-2-16	11-4-16	
Butylbenzylphthalate	ND	0.18	EPA 8270D	11-2-16	11-4-16	
bis-2-Ethylhexyladipate	ND	0.18	EPA 8270D	11-2-16	11-4-16	
3,3'-Dichlorobenzidine	ND	0.92	EPA 8270D	11-2-16	11-4-16	
Benzo[a]anthracene	0.21	0.15	EPA 8270D/SIM	11-2-16	11-4-16	
Chrysene	ND	0.15	EPA 8270D/SIM	11-2-16	11-4-16	
bis(2-Ethylhexyl)phthalate	ND	0.18	EPA 8270D	11-2-16	11-4-16	
Di-n-octylphthalate	ND	0.18	EPA 8270D	11-2-16	11-4-16	
Benzo[b]fluoranthene	ND	0.15	EPA 8270D/SIM	11-2-16	11-4-16	
Benzo(j,k)fluoranthene	ND	0.15	EPA 8270D/SIM	11-2-16	11-4-16	
Benzo[a]pyrene	0.12	0.037	EPA 8270D/SIM	11-2-16	11-4-16	
Indeno[1,2,3-cd]pyrene	0.078	0.037	EPA 8270D/SIM	11-2-16	11-4-16	
Dibenz[a,h]anthracene	ND	0.037	EPA 8270D/SIM	11-2-16	11-4-16	
Benzo[g,h,i]perylene	0.15	0.15	EPA 8270D/SIM	11-2-16	11-4-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	53	18 - 109				
Phenol-d6	62	25 - 111				
Nitrobenzene-d5	67	22 - 113				
2-Fluorobiphenyl	68	30 - 114				
2,4,6-Tribromophenol	76	22 - 116				
Terphenyl-d14	38	33 - 114				



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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>SS 4</b>					
<b>Laboratory ID:</b>	<b>10-295-04</b>					
n-Nitrosodimethylamine	ND	0.040	EPA 8270D	11-2-16	11-7-16	
Pyridine	ND	0.40	EPA 8270D	11-2-16	11-7-16	
Phenol	ND	0.040	EPA 8270D	11-2-16	11-7-16	
Aniline	ND	0.20	EPA 8270D	11-2-16	11-7-16	
bis(2-Chloroethyl)ether	ND	0.040	EPA 8270D	11-2-16	11-7-16	
2-Chlorophenol	ND	0.040	EPA 8270D	11-2-16	11-7-16	
1,3-Dichlorobenzene	ND	0.040	EPA 8270D	11-2-16	11-7-16	
1,4-Dichlorobenzene	ND	0.040	EPA 8270D	11-2-16	11-7-16	
Benzyl alcohol	ND	0.20	EPA 8270D	11-2-16	11-7-16	
1,2-Dichlorobenzene	ND	0.040	EPA 8270D	11-2-16	11-7-16	
2-Methylphenol (o-Cresol)	ND	0.040	EPA 8270D	11-2-16	11-7-16	
bis(2-Chloroisopropyl)ether	ND	0.040	EPA 8270D	11-2-16	11-7-16	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.040	EPA 8270D	11-2-16	11-7-16	
n-Nitroso-di-n-propylamine	ND	0.040	EPA 8270D	11-2-16	11-7-16	
Hexachloroethane	ND	0.040	EPA 8270D	11-2-16	11-7-16	
Nitrobenzene	ND	0.040	EPA 8270D	11-2-16	11-7-16	
Isophorone	ND	0.040	EPA 8270D	11-2-16	11-7-16	
2-Nitrophenol	ND	0.040	EPA 8270D	11-2-16	11-7-16	
2,4-Dimethylphenol	ND	0.040	EPA 8270D	11-2-16	11-7-16	
bis(2-Chloroethoxy)methane	ND	0.040	EPA 8270D	11-2-16	11-7-16	
2,4-Dichlorophenol	ND	0.040	EPA 8270D	11-2-16	11-7-16	
1,2,4-Trichlorobenzene	ND	0.040	EPA 8270D	11-2-16	11-7-16	
Naphthalene	ND	0.0079	EPA 8270D/SIM	11-2-16	11-7-16	
4-Chloroaniline	ND	0.20	EPA 8270D	11-2-16	11-7-16	
Hexachlorobutadiene	ND	0.040	EPA 8270D	11-2-16	11-7-16	
4-Chloro-3-methylphenol	ND	0.040	EPA 8270D	11-2-16	11-7-16	
2-Methylnaphthalene	ND	0.0079	EPA 8270D/SIM	11-2-16	11-7-16	
1-Methylnaphthalene	ND	0.0079	EPA 8270D/SIM	11-2-16	11-7-16	
Hexachlorocyclopentadiene	ND	0.040	EPA 8270D	11-2-16	11-7-16	
2,4,6-Trichlorophenol	ND	0.040	EPA 8270D	11-2-16	11-7-16	
2,3-Dichloroaniline	ND	0.040	EPA 8270D	11-2-16	11-7-16	
2,4,5-Trichlorophenol	ND	0.040	EPA 8270D	11-2-16	11-7-16	
2-Chloronaphthalene	ND	0.040	EPA 8270D	11-2-16	11-7-16	
2-Nitroaniline	ND	0.040	EPA 8270D	11-2-16	11-7-16	
1,4-Dinitrobenzene	ND	0.040	EPA 8270D	11-2-16	11-7-16	
Dimethylphthalate	ND	0.040	EPA 8270D	11-2-16	11-7-16	
1,3-Dinitrobenzene	ND	0.040	EPA 8270D	11-2-16	11-7-16	
2,6-Dinitrotoluene	ND	0.040	EPA 8270D	11-2-16	11-7-16	
1,2-Dinitrobenzene	ND	0.040	EPA 8270D	11-2-16	11-7-16	
Acenaphthylene	ND	0.0079	EPA 8270D/SIM	11-2-16	11-7-16	
3-Nitroaniline	ND	0.040	EPA 8270D	11-2-16	11-7-16	



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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>SS 4</b>					
Laboratory ID:	10-295-04					
2,4-Dinitrophenol	ND	0.20	EPA 8270D	11-2-16	11-7-16	
Acenaphthene	ND	0.0079	EPA 8270D/SIM	11-2-16	11-7-16	
4-Nitrophenol	ND	0.040	EPA 8270D	11-2-16	11-7-16	
2,4-Dinitrotoluene	ND	0.040	EPA 8270D	11-2-16	11-7-16	
Dibenzofuran	ND	0.040	EPA 8270D	11-2-16	11-7-16	
2,3,5,6-Tetrachlorophenol	ND	0.040	EPA 8270D	11-2-16	11-7-16	
2,3,4,6-Tetrachlorophenol	ND	0.040	EPA 8270D	11-2-16	11-7-16	
Diethylphthalate	ND	0.20	EPA 8270D	11-2-16	11-7-16	
4-Chlorophenyl-phenylether	ND	0.040	EPA 8270D	11-2-16	11-7-16	
4-Nitroaniline	ND	0.040	EPA 8270D	11-2-16	11-7-16	
Fluorene	ND	0.0079	EPA 8270D/SIM	11-2-16	11-7-16	
4,6-Dinitro-2-methylphenol	ND	0.20	EPA 8270D	11-2-16	11-7-16	
n-Nitrosodiphenylamine	ND	0.040	EPA 8270D	11-2-16	11-7-16	
1,2-Diphenylhydrazine	ND	0.040	EPA 8270D	11-2-16	11-7-16	
4-Bromophenyl-phenylether	ND	0.040	EPA 8270D	11-2-16	11-7-16	
Hexachlorobenzene	ND	0.040	EPA 8270D	11-2-16	11-7-16	
Pentachlorophenol	ND	0.20	EPA 8270D	11-2-16	11-7-16	
Phenanthrene	ND	0.0079	EPA 8270D/SIM	11-2-16	11-7-16	
Anthracene	ND	0.0079	EPA 8270D/SIM	11-2-16	11-7-16	
Carbazole	ND	0.040	EPA 8270D	11-2-16	11-7-16	
Di-n-butylphthalate	ND	0.20	EPA 8270D	11-2-16	11-7-16	
Fluoranthene	ND	0.0079	EPA 8270D/SIM	11-2-16	11-7-16	
Benzidine	ND	0.40	EPA 8270D	11-2-16	11-7-16	
Pyrene	ND	0.0079	EPA 8270D/SIM	11-2-16	11-7-16	
Butylbenzylphthalate	ND	0.040	EPA 8270D	11-2-16	11-7-16	
bis-2-Ethylhexyladipate	ND	0.040	EPA 8270D	11-2-16	11-7-16	
3,3'-Dichlorobenzidine	ND	0.20	EPA 8270D	11-2-16	11-7-16	
Benzo[a]anthracene	ND	0.0079	EPA 8270D/SIM	11-2-16	11-7-16	
Chrysene	ND	0.0079	EPA 8270D/SIM	11-2-16	11-7-16	
bis(2-Ethylhexyl)phthalate	ND	0.040	EPA 8270D	11-2-16	11-7-16	
Di-n-octylphthalate	ND	0.040	EPA 8270D	11-2-16	11-7-16	
Benzo[b]fluoranthene	ND	0.0079	EPA 8270D/SIM	11-2-16	11-7-16	
Benzo(j,k)fluoranthene	ND	0.0079	EPA 8270D/SIM	11-2-16	11-7-16	
Benzo[a]pyrene	ND	0.0079	EPA 8270D/SIM	11-2-16	11-7-16	
Indeno[1,2,3-cd]pyrene	ND	0.0079	EPA 8270D/SIM	11-2-16	11-7-16	
Dibenz[a,h]anthracene	ND	0.0079	EPA 8270D/SIM	11-2-16	11-7-16	
Benzo[g,h,i]perylene	ND	0.0079	EPA 8270D/SIM	11-2-16	11-7-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	1.6	18 - 109				Q
Phenol-d6	75	25 - 111				
Nitrobenzene-d5	78	22 - 113				
2-Fluorobiphenyl	83	30 - 114				
2,4,6-Tribromophenol	92	22 - 116				
Terphenyl-d14	85	33 - 114				



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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:	MB1102S1					
n-Nitrosodimethylamine	ND	0.033	EPA 8270D	11-2-16	11-3-16	
Pyridine	ND	0.33	EPA 8270D	11-2-16	11-3-16	
Phenol	ND	0.033	EPA 8270D	11-2-16	11-3-16	
Aniline	ND	0.17	EPA 8270D	11-2-16	11-3-16	
bis(2-Chloroethyl)ether	ND	0.033	EPA 8270D	11-2-16	11-3-16	
2-Chlorophenol	ND	0.033	EPA 8270D	11-2-16	11-3-16	
1,3-Dichlorobenzene	ND	0.033	EPA 8270D	11-2-16	11-3-16	
1,4-Dichlorobenzene	ND	0.033	EPA 8270D	11-2-16	11-3-16	
Benzyl alcohol	ND	0.17	EPA 8270D	11-2-16	11-3-16	
1,2-Dichlorobenzene	ND	0.033	EPA 8270D	11-2-16	11-3-16	
2-Methylphenol (o-Cresol)	ND	0.033	EPA 8270D	11-2-16	11-3-16	
bis(2-Chloroisopropyl)ether	ND	0.033	EPA 8270D	11-2-16	11-3-16	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.033	EPA 8270D	11-2-16	11-3-16	
n-Nitroso-di-n-propylamine	ND	0.033	EPA 8270D	11-2-16	11-3-16	
Hexachloroethane	ND	0.033	EPA 8270D	11-2-16	11-3-16	
Nitrobenzene	ND	0.033	EPA 8270D	11-2-16	11-3-16	
Isophorone	ND	0.033	EPA 8270D	11-2-16	11-3-16	
2-Nitrophenol	ND	0.033	EPA 8270D	11-2-16	11-3-16	
2,4-Dimethylphenol	ND	0.033	EPA 8270D	11-2-16	11-3-16	
bis(2-Chloroethoxy)methane	ND	0.033	EPA 8270D	11-2-16	11-3-16	
2,4-Dichlorophenol	ND	0.033	EPA 8270D	11-2-16	11-3-16	
1,2,4-Trichlorobenzene	ND	0.033	EPA 8270D	11-2-16	11-3-16	
Naphthalene	ND	0.0067	EPA 8270D/SIM	11-2-16	11-3-16	
4-Chloroaniline	ND	0.17	EPA 8270D	11-2-16	11-3-16	
Hexachlorobutadiene	ND	0.033	EPA 8270D	11-2-16	11-3-16	
4-Chloro-3-methylphenol	ND	0.033	EPA 8270D	11-2-16	11-3-16	
2-Methylnaphthalene	ND	0.0067	EPA 8270D/SIM	11-2-16	11-3-16	
1-Methylnaphthalene	ND	0.0067	EPA 8270D/SIM	11-2-16	11-3-16	
Hexachlorocyclopentadiene	ND	0.033	EPA 8270D	11-2-16	11-3-16	
2,4,6-Trichlorophenol	ND	0.033	EPA 8270D	11-2-16	11-3-16	
2,3-Dichloroaniline	ND	0.033	EPA 8270D	11-2-16	11-3-16	
2,4,5-Trichlorophenol	ND	0.033	EPA 8270D	11-2-16	11-3-16	
2-Chloronaphthalene	ND	0.033	EPA 8270D	11-2-16	11-3-16	
2-Nitroaniline	ND	0.033	EPA 8270D	11-2-16	11-3-16	
1,4-Dinitrobenzene	ND	0.033	EPA 8270D	11-2-16	11-3-16	
Dimethylphthalate	ND	0.033	EPA 8270D	11-2-16	11-3-16	
1,3-Dinitrobenzene	ND	0.033	EPA 8270D	11-2-16	11-3-16	
2,6-Dinitrotoluene	ND	0.033	EPA 8270D	11-2-16	11-3-16	
1,2-Dinitrobenzene	ND	0.033	EPA 8270D	11-2-16	11-3-16	
Acenaphthylene	ND	0.0067	EPA 8270D/SIM	11-2-16	11-3-16	
3-Nitroaniline	ND	0.033	EPA 8270D	11-2-16	11-3-16	



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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:	MB1102S1					
2,4-Dinitrophenol	ND	0.17	EPA 8270D	11-2-16	11-3-16	
Acenaphthene	ND	0.0067	EPA 8270D/SIM	11-2-16	11-3-16	
4-Nitrophenol	ND	0.033	EPA 8270D	11-2-16	11-3-16	
2,4-Dinitrotoluene	ND	0.033	EPA 8270D	11-2-16	11-3-16	
Dibenzofuran	ND	0.033	EPA 8270D	11-2-16	11-3-16	
2,3,5,6-Tetrachlorophenol	ND	0.033	EPA 8270D	11-2-16	11-3-16	
2,3,4,6-Tetrachlorophenol	ND	0.033	EPA 8270D	11-2-16	11-3-16	
Diethylphthalate	ND	0.17	EPA 8270D	11-2-16	11-3-16	
4-Chlorophenyl-phenylether	ND	0.033	EPA 8270D	11-2-16	11-3-16	
4-Nitroaniline	ND	0.033	EPA 8270D	11-2-16	11-3-16	
Fluorene	ND	0.0067	EPA 8270D/SIM	11-2-16	11-3-16	
4,6-Dinitro-2-methylphenol	ND	0.17	EPA 8270D	11-2-16	11-3-16	
n-Nitrosodiphenylamine	ND	0.033	EPA 8270D	11-2-16	11-3-16	
1,2-Diphenylhydrazine	ND	0.033	EPA 8270D	11-2-16	11-3-16	
4-Bromophenyl-phenylether	ND	0.033	EPA 8270D	11-2-16	11-3-16	
Hexachlorobenzene	ND	0.033	EPA 8270D	11-2-16	11-3-16	
Pentachlorophenol	ND	0.17	EPA 8270D	11-2-16	11-3-16	
Phenanthrene	ND	0.0067	EPA 8270D/SIM	11-2-16	11-3-16	
Anthracene	ND	0.0067	EPA 8270D/SIM	11-2-16	11-3-16	
Carbazole	ND	0.033	EPA 8270D	11-2-16	11-3-16	
Di-n-butylphthalate	ND	0.17	EPA 8270D	11-2-16	11-3-16	
Fluoranthene	ND	0.0067	EPA 8270D/SIM	11-2-16	11-3-16	
Benzidine	ND	0.33	EPA 8270D	11-2-16	11-3-16	
Pyrene	ND	0.0067	EPA 8270D/SIM	11-2-16	11-3-16	
Butylbenzylphthalate	ND	0.033	EPA 8270D	11-2-16	11-3-16	
bis-2-Ethylhexyladipate	ND	0.033	EPA 8270D	11-2-16	11-3-16	
3,3'-Dichlorobenzidine	ND	0.17	EPA 8270D	11-2-16	11-3-16	
Benzo[a]anthracene	ND	0.0067	EPA 8270D/SIM	11-2-16	11-3-16	
Chrysene	ND	0.0067	EPA 8270D/SIM	11-2-16	11-3-16	
bis(2-Ethylhexyl)phthalate	ND	0.033	EPA 8270D	11-2-16	11-3-16	
Di-n-octylphthalate	ND	0.033	EPA 8270D	11-2-16	11-3-16	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270D/SIM	11-2-16	11-3-16	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270D/SIM	11-2-16	11-3-16	
Benzo[a]pyrene	ND	0.0067	EPA 8270D/SIM	11-2-16	11-3-16	
Indeno[1,2,3-cd]pyrene	ND	0.0067	EPA 8270D/SIM	11-2-16	11-3-16	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270D/SIM	11-2-16	11-3-16	
Benzo[g,h,i]perylene	ND	0.0067	EPA 8270D/SIM	11-2-16	11-3-16	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorophenol</i>	<i>69</i>	<i>18 - 109</i>				
<i>Phenol-d6</i>	<i>70</i>	<i>25 - 111</i>				
<i>Nitrobenzene-d5</i>	<i>72</i>	<i>22 - 113</i>				
<i>2-Fluorobiphenyl</i>	<i>74</i>	<i>30 - 114</i>				
<i>2,4,6-Tribromophenol</i>	<i>76</i>	<i>22 - 116</i>				
<i>Terphenyl-d14</i>	<i>81</i>	<i>33 - 114</i>				



Date of Report: November 8, 2016  
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**SEMIVOLATILES EPA 8270D/SIM  
 SB/SBD QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD	RPD	Flags
					SB	SBD	Limits		Limit	
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB1102S1									
	SB	SBD	SB	SBD	SB	SBD				
Phenol	1.07	1.02	1.33	1.33	80	77	37 - 107	5	39	
2-Chlorophenol	1.08	1.01	1.33	1.33	81	76	36 - 107	7	40	
1,4-Dichlorobenzene	0.503	0.485	0.667	0.667	75	73	28 - 108	4	41	
n-Nitroso-di-n-propylamine	0.488	0.465	0.667	0.667	73	70	28 - 113	5	33	
1,2,4-Trichlorobenzene	0.505	0.489	0.667	0.667	76	73	33 - 106	3	38	
4-Chloro-3-methylphenol	1.15	1.11	1.33	1.33	86	83	52 - 106	4	30	
Acenaphthene	0.525	0.510	0.667	0.667	79	76	52 - 90	3	30	
4-Nitrophenol	1.44	1.48	1.33	1.33	108	111	30 - 109	3	32	I
2,4-Dinitrotoluene	0.576	0.569	0.667	0.667	86	85	50 - 101	1	32	
Pentachlorophenol	0.981	0.855	1.33	1.33	74	64	21 - 114	14	40	
Pyrene	0.590	0.592	0.667	0.667	88	89	52 - 104	0	30	
<i>Surrogate:</i>										
2-Fluorophenol					79	75	18 - 109			
Phenol-d6					84	78	25 - 111			
Nitrobenzene-d5					81	76	22 - 113			
2-Fluorobiphenyl					87	80	30 - 114			
2,4,6-Tribromophenol					90	94	22 - 116			
Terphenyl-d14					95	92	33 - 114			



Date of Report: November 8, 2016  
 Samples Submitted: October 26, 2016  
 Laboratory Reference: 1610-295  
 Project: SR 509 - 12th Street

**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:	10-295-01					
<b>Client ID:</b>	<b>Soil Sample 1</b>					
Arsenic	<b>ND</b>	11	6010C	10-27-16	10-27-16	
Barium	<b>79</b>	2.8	6010C	10-27-16	10-27-16	
Cadmium	<b>ND</b>	0.56	6010C	10-27-16	10-27-16	
Chromium	<b>48</b>	0.56	6010C	10-27-16	10-27-16	
Lead	<b>6.2</b>	5.6	6010C	10-27-16	10-27-16	
Mercury	<b>ND</b>	0.28	7471B	10-27-16	10-27-16	
Selenium	<b>ND</b>	11	6010C	10-27-16	10-27-16	
Silver	<b>ND</b>	1.1	6010C	10-27-16	10-27-16	

Lab ID:	10-295-02					
<b>Client ID:</b>	<b>Soil Sample 2</b>					
Arsenic	<b>35</b>	14	6010C	10-27-16	10-27-16	
Barium	<b>170</b>	3.5	6010C	10-27-16	10-27-16	
Cadmium	<b>0.77</b>	0.70	6010C	10-27-16	10-27-16	
Chromium	<b>45</b>	0.70	6010C	10-27-16	10-27-16	
Lead	<b>37</b>	7.0	6010C	10-27-16	10-27-16	
Mercury	<b>ND</b>	0.35	7471B	10-27-16	10-27-16	
Selenium	<b>ND</b>	14	6010C	10-27-16	10-27-16	
Silver	<b>ND</b>	1.4	6010C	10-27-16	10-27-16	



Date of Report: November 8, 2016  
 Samples Submitted: October 26, 2016  
 Laboratory Reference: 1610-295  
 Project: SR 509 - 12th Street

**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:	10-295-03					
<b>Client ID:</b>	<b>Soil Sample 3</b>					
Arsenic	ND	11	6010C	10-27-16	10-27-16	
Barium	55	2.8	6010C	10-27-16	10-27-16	
Cadmium	ND	0.55	6010C	10-27-16	10-27-16	
Chromium	22	0.55	6010C	10-27-16	10-27-16	
Lead	45	5.5	6010C	10-27-16	10-27-16	
Mercury	ND	0.28	7471B	10-27-16	10-27-16	
Selenium	ND	11	6010C	10-27-16	10-27-16	
Silver	ND	1.1	6010C	10-27-16	10-27-16	

Lab ID:	10-295-04					
<b>Client ID:</b>	<b>SS 4</b>					
Arsenic	ND	12	6010C	10-27-16	10-27-16	
Barium	73	3.0	6010C	10-27-16	10-27-16	
Cadmium	ND	0.59	6010C	10-27-16	10-27-16	
Chromium	42	0.59	6010C	10-27-16	10-27-16	
Lead	ND	5.9	6010C	10-27-16	10-27-16	
Mercury	ND	0.30	7471B	10-27-16	10-27-16	
Selenium	ND	12	6010C	10-27-16	10-27-16	
Silver	ND	1.2	6010C	10-27-16	10-27-16	



Date of Report: November 8, 2016  
 Samples Submitted: October 26, 2016  
 Laboratory Reference: 1610-295  
 Project: SR 509 - 12th Street

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

Date Extracted: 10-27-16

Date Analyzed: 10-27-16

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: MB1027SM1,MB1027SM2&MB1027S1

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: November 8, 2016  
 Samples Submitted: October 26, 2016  
 Laboratory Reference: 1610-295  
 Project: SR 509 - 12th Street

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

Date Extracted: 10-27-16

Date Analyzed: 10-27-16

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 10-289-02

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Arsenic	ND	ND	NA	10	
Barium	47.1	46.5	1	2.5	
Cadmium	ND	ND	NA	0.50	
Chromium	50.1	48.5	3	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: November 8, 2016  
 Samples Submitted: October 26, 2016  
 Laboratory Reference: 1610-295  
 Project: SR 509 - 12th Street

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

Date Extracted: 10-27-16

Date Analyzed: 10-27-16

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 10-289-02

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Arsenic	100	<b>92.3</b>	92	<b>91.8</b>	92	0	
Barium	100	<b>148</b>	101	<b>142</b>	95	4	
Cadmium	50.0	<b>48.7</b>	97	<b>48.3</b>	97	1	
Chromium	100	<b>135</b>	85	<b>137</b>	86	1	
Lead	250	<b>265</b>	106	<b>250</b>	100	6	
Mercury	0.500	<b>0.535</b>	107	<b>0.542</b>	108	1	
Selenium	100	<b>92.3</b>	92	<b>91.9</b>	92	0	
Silver	25.0	<b>20.8</b>	83	<b>21.0</b>	84	1	



Date of Report: November 8, 2016  
Samples Submitted: October 26, 2016  
Laboratory Reference: 1610-295  
Project: SR 509 - 12th Street

### % MOISTURE

Date Analyzed: 10-27-16

Client ID	Lab ID	% Moisture
Soil Sample 1	10-295-01	11
Soil Sample 2	10-295-02	28
Soil Sample 3	10-295-03	10
SS 4	10-295-04	16





### 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) (TPH analysis 5 Days)  
 (other) \_\_\_\_\_

Laboratory Number: **10-295**

10-295

Company: **WSDOT**  
 Project Number:  
 Project Name: **SR509 - 12th St**  
 Project Manager: **PAT SYOBODA**  
 Sampled by: **WSDOT**

Lab ID	Sample Identification	Date Sampled	Time Sampled	Matrix	Number of Containers	Laboratory Tests										% Moisture							
1	Soil Sample 1	10/26/16	3:10	Soil	3	NWTPH-HCID	NWTPH-Gx/BTEX	NWTPH-Gx	NWTPH-Dx	Volatiles 8260C	Halogenated Volatiles 8260C	Volatiles 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		
2	Soil Sample 2	10/26/16	3:00	Soil	1																		X
3	Soil Sample 3	10/26/16	3:30	Soil	1																		X
4	SS 4	10/26/16	3:45	SS	1																		X

Signature	Company	Date	Time	Comments/Special Instructions
<i>Pat Syoboda</i>	WSDOT	10/26/16	8:15A	
<i>Marie Sylvestre</i>	SPEDDY	10-26-16	8:15A	
<i>Marie Sylvestre</i>	SPEDDY	10-26-16	11:10A	
<i>[Signature]</i>	ORRE	10/26/16	1110	



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

March 27, 2018

Patrick Svoboda  
WSDOT  
15700 Dayton Avenue North  
NB82-138  
P.O. Box 330310  
Seattle, WA 98133-9710

Re: Analytical Data for Project SR509 Foreman Property  
Laboratory Reference No. 1803-152

Dear Pat:

Enclosed are the analytical results and associated quality control data for samples submitted on March 16, 2018.

The standard policy of OnSite Environmental, Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

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: March 27, 2018  
Samples Submitted: March 16, 2018  
Laboratory Reference: 1803-152  
Project: SR509 Foreman Property

### Case Narrative

Samples were collected on March 15, 2018 and received by the laboratory on March 16, 2018. They were maintained at the laboratory at a temperature of 2°C to 6°C.

Please note that any and all soil sample results are reported on a dry-weight basis, unless otherwise noted below.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

#### Semivolatiles EPA 8270D/SIM Analysis

The spike blank duplicate had one recovery slightly above control limits. The samples were non-detect for this analyte, further action was taken.

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 27, 2018  
 Samples Submitted: March 16, 2018  
 Laboratory Reference: 1803-152  
 Project: SR509 Foreman Property

### NWTPH-Gx

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SS1 N A'</b>					
Laboratory ID:	03-152-01					
Gasoline	<b>ND</b>	7.3	NWTPH-Gx	3-19-18	3-19-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	101	66-130				
<b>Client ID:</b>	<b>SS2</b>					
Laboratory ID:	03-152-02					
Gasoline	<b>ND</b>	12	NWTPH-Gx	3-19-18	3-19-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	97	66-130				
<b>Client ID:</b>	<b>SS3</b>					
Laboratory ID:	03-152-03					
Gasoline	<b>ND</b>	5.4	NWTPH-Gx	3-19-18	3-19-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	95	66-130				
<b>Client ID:</b>	<b>SS4</b>					
Laboratory ID:	03-152-04					
Gasoline	<b>ND</b>	10	NWTPH-Gx	3-19-18	3-19-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	97	66-130				



Date of Report: March 27, 2018  
 Samples Submitted: March 16, 2018  
 Laboratory Reference: 1803-152  
 Project: SR509 Foreman Property

**NWTPH-Gx  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB0319S1					
Gasoline	<b>ND</b>	5.0	NWTPH-Gx	3-19-18	3-19-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	86	66-130				

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>								
Laboratory ID:	03-145-01							
	ORIG	DUP						
Gasoline	<b>ND</b>	<b>ND</b>	NA	NA	NA	NA	NA	30
<i>Surrogate:</i>								
<i>Fluorobenzene</i>				95	97	66-130		



Date of Report: March 27, 2018  
 Samples Submitted: March 16, 2018  
 Laboratory Reference: 1803-152  
 Project: SR509 Foreman Property

### NWTPH-Dx

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SS1 N A'</b>					
Laboratory ID:	03-152-01					
Diesel Range Organics	<b>ND</b>	35	NWTPH-Dx	3-19-18	3-19-18	
Lube Oil Range Organics	<b>82</b>	71	NWTPH-Dx	3-19-18	3-19-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	92	50-150				
<b>Client ID:</b>	<b>SS2</b>					
Laboratory ID:	03-152-02					
Diesel Range Organics	<b>58</b>	34	NWTPH-Dx	3-19-18	3-19-18	N
Lube Oil Range Organics	<b>370</b>	68	NWTPH-Dx	3-19-18	3-19-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	86	50-150				
<b>Client ID:</b>	<b>SS3</b>					
Laboratory ID:	03-152-03					
Diesel Range Organics	<b>ND</b>	60	NWTPH-Dx	3-19-18	3-20-18	
Lube Oil	<b>470</b>	120	NWTPH-Dx	3-19-18	3-20-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	89	50-150				
<b>Client ID:</b>	<b>SS4</b>					
Laboratory ID:	03-152-04					
Diesel Range Organics	<b>ND</b>	33	NWTPH-Dx	3-19-18	3-19-18	U1
Lube Oil	<b>260</b>	61	NWTPH-Dx	3-19-18	3-19-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	79	50-150				



Date of Report: March 27, 2018  
 Samples Submitted: March 16, 2018  
 Laboratory Reference: 1803-152  
 Project: SR509 Foreman Property

**NWTPH-Dx  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB0319S1					
Diesel Range Organics	<b>ND</b>	25	NWTPH-Dx	3-19-18	3-19-18	
Lube Oil Range Organics	<b>ND</b>	50	NWTPH-Dx	3-19-18	3-19-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	97	50-150				

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>								
Laboratory ID:	03-145-01							
	ORIG	DUP						
Diesel Range Organics	<b>64.2</b>	<b>42.0</b>	NA	NA	NA	NA	42	NA
Lube Oil Range Organics	<b>188</b>	<b>137</b>	NA	NA	NA	NA	31	NA
<i>Surrogate:</i>								
<i>o-Terphenyl</i>				88	81	50-150		



Date of Report: March 27, 2018  
 Samples Submitted: March 16, 2018  
 Laboratory Reference: 1803-152  
 Project: SR509 Foreman Property

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

Matrix: Soil  
 Units: mg/kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SS1 N A'</b>					
Laboratory ID:	03-152-01					
Dichlorodifluoromethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Chloromethane	ND	0.0054	EPA 8260C	3-19-18	3-19-18	
Vinyl Chloride	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Bromomethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Chloroethane	ND	0.0054	EPA 8260C	3-19-18	3-19-18	
Trichlorofluoromethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,1-Dichloroethene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Acetone	0.023	0.0054	EPA 8260C	3-19-18	3-19-18	
Iodomethane	ND	0.0054	EPA 8260C	3-19-18	3-19-18	
Carbon Disulfide	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Methylene Chloride	ND	0.0054	EPA 8260C	3-19-18	3-19-18	
(trans) 1,2-Dichloroethene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Methyl t-Butyl Ether	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,1-Dichloroethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Vinyl Acetate	ND	0.0054	EPA 8260C	3-19-18	3-19-18	
2,2-Dichloropropane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
(cis) 1,2-Dichloroethene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
2-Butanone	ND	0.0054	EPA 8260C	3-19-18	3-19-18	
Bromochloromethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Chloroform	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,1,1-Trichloroethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Carbon Tetrachloride	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,1-Dichloropropene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Benzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,2-Dichloroethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Trichloroethene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,2-Dichloropropane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Dibromomethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Bromodichloromethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
2-Chloroethyl Vinyl Ether	ND	0.0054	EPA 8260C	3-19-18	3-19-18	
(cis) 1,3-Dichloropropene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Methyl Isobutyl Ketone	ND	0.0054	EPA 8260C	3-19-18	3-19-18	
Toluene	ND	0.0054	EPA 8260C	3-19-18	3-19-18	
(trans) 1,3-Dichloropropene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	



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**VOLATILES EPA 8260C**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SS1 N A'</b>					
Laboratory ID:	03-152-01					
1,1,2-Trichloroethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Tetrachloroethene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,3-Dichloropropane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
2-Hexanone	ND	0.0054	EPA 8260C	3-19-18	3-19-18	
Dibromochloromethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,2-Dibromoethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Chlorobenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,1,1,2-Tetrachloroethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Ethylbenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
m,p-Xylene	ND	0.0022	EPA 8260C	3-19-18	3-19-18	
o-Xylene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Styrene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Bromoform	ND	0.0054	EPA 8260C	3-19-18	3-19-18	
Isopropylbenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Bromobenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,1,2,2-Tetrachloroethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,2,3-Trichloropropane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
n-Propylbenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
2-Chlorotoluene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
4-Chlorotoluene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,3,5-Trimethylbenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
tert-Butylbenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,2,4-Trimethylbenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
sec-Butylbenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,3-Dichlorobenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
p-Isopropyltoluene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,4-Dichlorobenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,2-Dichlorobenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
n-Butylbenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,2-Dibromo-3-chloropropane	ND	0.0054	EPA 8260C	3-19-18	3-19-18	
1,2,4-Trichlorobenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Hexachlorobutadiene	ND	0.0054	EPA 8260C	3-19-18	3-19-18	
Naphthalene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,2,3-Trichlorobenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>120</i>	<i>75-131</i>				
<i>Toluene-d8</i>	<i>119</i>	<i>83-130</i>				
<i>4-Bromofluorobenzene</i>	<i>115</i>	<i>78-130</i>				



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

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SS2</b>					
Laboratory ID:	03-152-02					
Dichlorodifluoromethane	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
Chloromethane	ND	0.0063	EPA 8260C	3-19-18	3-19-18	
Vinyl Chloride	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
Bromomethane	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
Chloroethane	ND	0.0063	EPA 8260C	3-19-18	3-19-18	
Trichlorofluoromethane	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
1,1-Dichloroethene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
Acetone	0.0075	0.0063	EPA 8260C	3-19-18	3-19-18	
Iodomethane	ND	0.0063	EPA 8260C	3-19-18	3-19-18	
Carbon Disulfide	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
Methylene Chloride	ND	0.0063	EPA 8260C	3-19-18	3-19-18	
(trans) 1,2-Dichloroethene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
Methyl t-Butyl Ether	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
1,1-Dichloroethane	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
Vinyl Acetate	ND	0.0063	EPA 8260C	3-19-18	3-19-18	
2,2-Dichloropropane	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
(cis) 1,2-Dichloroethene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
2-Butanone	ND	0.0063	EPA 8260C	3-19-18	3-19-18	
Bromochloromethane	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
Chloroform	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
1,1,1-Trichloroethane	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
Carbon Tetrachloride	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
1,1-Dichloropropene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
Benzene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
1,2-Dichloroethane	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
Trichloroethene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
1,2-Dichloropropane	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
Dibromomethane	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
Bromodichloromethane	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
2-Chloroethyl Vinyl Ether	ND	0.0063	EPA 8260C	3-19-18	3-19-18	
(cis) 1,3-Dichloropropene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
Methyl Isobutyl Ketone	ND	0.0063	EPA 8260C	3-19-18	3-19-18	
Toluene	ND	0.0063	EPA 8260C	3-19-18	3-19-18	
(trans) 1,3-Dichloropropene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	



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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SS2</b>					
Laboratory ID:	03-152-02					
1,1,2-Trichloroethane	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
Tetrachloroethene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
1,3-Dichloropropane	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
2-Hexanone	ND	0.0063	EPA 8260C	3-19-18	3-19-18	
Dibromochloromethane	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
1,2-Dibromoethane	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
Chlorobenzene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
1,1,1,2-Tetrachloroethane	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
Ethylbenzene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
m,p-Xylene	ND	0.0025	EPA 8260C	3-19-18	3-19-18	
o-Xylene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
Styrene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
Bromoform	ND	0.0063	EPA 8260C	3-19-18	3-19-18	
Isopropylbenzene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
Bromobenzene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
1,1,2,2-Tetrachloroethane	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
1,2,3-Trichloropropane	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
n-Propylbenzene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
2-Chlorotoluene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
4-Chlorotoluene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
1,3,5-Trimethylbenzene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
tert-Butylbenzene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
1,2,4-Trimethylbenzene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
sec-Butylbenzene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
1,3-Dichlorobenzene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
p-Isopropyltoluene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
1,4-Dichlorobenzene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
1,2-Dichlorobenzene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
n-Butylbenzene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
1,2-Dibromo-3-chloropropane	ND	0.0063	EPA 8260C	3-19-18	3-19-18	
1,2,4-Trichlorobenzene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
Hexachlorobutadiene	ND	0.0063	EPA 8260C	3-19-18	3-19-18	
Naphthalene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
1,2,3-Trichlorobenzene	ND	0.0013	EPA 8260C	3-19-18	3-19-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>107</i>	<i>75-131</i>				
<i>Toluene-d8</i>	<i>104</i>	<i>83-130</i>				
<i>4-Bromofluorobenzene</i>	<i>100</i>	<i>78-130</i>				



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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>SS3</b>					
Laboratory ID:	03-152-03					
Dichlorodifluoromethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Chloromethane	ND	0.0055	EPA 8260C	3-19-18	3-19-18	
Vinyl Chloride	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Bromomethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Chloroethane	ND	0.0055	EPA 8260C	3-19-18	3-19-18	
Trichlorofluoromethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,1-Dichloroethene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Acetone	ND	0.0055	EPA 8260C	3-19-18	3-19-18	
Iodomethane	ND	0.0055	EPA 8260C	3-19-18	3-19-18	
Carbon Disulfide	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Methylene Chloride	ND	0.0055	EPA 8260C	3-19-18	3-19-18	
(trans) 1,2-Dichloroethene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Methyl t-Butyl Ether	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,1-Dichloroethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Vinyl Acetate	ND	0.0055	EPA 8260C	3-19-18	3-19-18	
2,2-Dichloropropane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
(cis) 1,2-Dichloroethene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
2-Butanone	ND	0.0055	EPA 8260C	3-19-18	3-19-18	
Bromochloromethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Chloroform	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,1,1-Trichloroethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Carbon Tetrachloride	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,1-Dichloropropene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Benzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,2-Dichloroethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Trichloroethene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,2-Dichloropropane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Dibromomethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Bromodichloromethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
2-Chloroethyl Vinyl Ether	ND	0.0055	EPA 8260C	3-19-18	3-19-18	
(cis) 1,3-Dichloropropene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Methyl Isobutyl Ketone	ND	0.0055	EPA 8260C	3-19-18	3-19-18	
Toluene	ND	0.0055	EPA 8260C	3-19-18	3-19-18	
(trans) 1,3-Dichloropropene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	



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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SS3</b>					
Laboratory ID:	03-152-03					
1,1,2-Trichloroethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Tetrachloroethene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,3-Dichloropropane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
2-Hexanone	ND	0.0055	EPA 8260C	3-19-18	3-19-18	
Dibromochloromethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,2-Dibromoethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Chlorobenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,1,1,2-Tetrachloroethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Ethylbenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
m,p-Xylene	ND	0.0022	EPA 8260C	3-19-18	3-19-18	
o-Xylene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Styrene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Bromoform	ND	0.0055	EPA 8260C	3-19-18	3-19-18	
Isopropylbenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Bromobenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,1,2,2-Tetrachloroethane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,2,3-Trichloropropane	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
n-Propylbenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
2-Chlorotoluene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
4-Chlorotoluene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,3,5-Trimethylbenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
tert-Butylbenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,2,4-Trimethylbenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
sec-Butylbenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,3-Dichlorobenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
p-Isopropyltoluene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,4-Dichlorobenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,2-Dichlorobenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
n-Butylbenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,2-Dibromo-3-chloropropane	ND	0.0055	EPA 8260C	3-19-18	3-19-18	
1,2,4-Trichlorobenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
Hexachlorobutadiene	ND	0.0055	EPA 8260C	3-19-18	3-19-18	
Naphthalene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
1,2,3-Trichlorobenzene	ND	0.0011	EPA 8260C	3-19-18	3-19-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>123</i>	<i>75-131</i>				
<i>Toluene-d8</i>	<i>119</i>	<i>83-130</i>				
<i>4-Bromofluorobenzene</i>	<i>115</i>	<i>78-130</i>				



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 Laboratory Reference: 1803-152  
 Project: SR509 Foreman Property

**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>SS4</b>					
Laboratory ID:	03-152-04					
Dichlorodifluoromethane	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
Chloromethane	ND	0.0081	EPA 8260C	3-19-18	3-19-18	
Vinyl Chloride	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
Bromomethane	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
Chloroethane	ND	0.0081	EPA 8260C	3-19-18	3-19-18	
Trichlorofluoromethane	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
1,1-Dichloroethene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
Acetone	ND	0.0081	EPA 8260C	3-19-18	3-19-18	
Iodomethane	ND	0.0081	EPA 8260C	3-19-18	3-19-18	
Carbon Disulfide	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
Methylene Chloride	ND	0.0081	EPA 8260C	3-19-18	3-19-18	
(trans) 1,2-Dichloroethene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
Methyl t-Butyl Ether	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
1,1-Dichloroethane	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
Vinyl Acetate	ND	0.0081	EPA 8260C	3-19-18	3-19-18	
2,2-Dichloropropane	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
(cis) 1,2-Dichloroethene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
2-Butanone	ND	0.0081	EPA 8260C	3-19-18	3-19-18	
Bromochloromethane	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
Chloroform	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
1,1,1-Trichloroethane	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
Carbon Tetrachloride	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
1,1-Dichloropropene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
Benzene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
1,2-Dichloroethane	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
Trichloroethene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
1,2-Dichloropropane	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
Dibromomethane	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
Bromodichloromethane	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
2-Chloroethyl Vinyl Ether	ND	0.0081	EPA 8260C	3-19-18	3-19-18	
(cis) 1,3-Dichloropropene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
Methyl Isobutyl Ketone	ND	0.0081	EPA 8260C	3-19-18	3-19-18	
Toluene	ND	0.0081	EPA 8260C	3-19-18	3-19-18	
(trans) 1,3-Dichloropropene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	



Date of Report: March 27, 2018  
 Samples Submitted: March 16, 2018  
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**VOLATILES EPA 8260C**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SS4</b>					
Laboratory ID:	03-152-04					
1,1,2-Trichloroethane	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
Tetrachloroethene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
1,3-Dichloropropane	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
2-Hexanone	ND	0.0081	EPA 8260C	3-19-18	3-19-18	
Dibromochloromethane	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
1,2-Dibromoethane	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
Chlorobenzene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
1,1,1,2-Tetrachloroethane	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
Ethylbenzene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
m,p-Xylene	ND	0.0032	EPA 8260C	3-19-18	3-19-18	
o-Xylene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
Styrene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
Bromoform	ND	0.0081	EPA 8260C	3-19-18	3-19-18	
Isopropylbenzene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
Bromobenzene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
1,1,2,2-Tetrachloroethane	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
1,2,3-Trichloropropane	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
n-Propylbenzene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
2-Chlorotoluene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
4-Chlorotoluene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
1,3,5-Trimethylbenzene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
tert-Butylbenzene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
1,2,4-Trimethylbenzene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
sec-Butylbenzene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
1,3-Dichlorobenzene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
p-Isopropyltoluene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
1,4-Dichlorobenzene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
1,2-Dichlorobenzene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
n-Butylbenzene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
1,2-Dibromo-3-chloropropane	ND	0.0081	EPA 8260C	3-19-18	3-19-18	
1,2,4-Trichlorobenzene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
Hexachlorobutadiene	ND	0.0081	EPA 8260C	3-19-18	3-19-18	
Naphthalene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
1,2,3-Trichlorobenzene	ND	0.0016	EPA 8260C	3-19-18	3-19-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>119</i>	<i>75-131</i>				
<i>Toluene-d8</i>	<i>121</i>	<i>83-130</i>				
<i>4-Bromofluorobenzene</i>	<i>118</i>	<i>78-130</i>				



Date of Report: March 27, 2018  
 Samples Submitted: March 16, 2018  
 Laboratory Reference: 1803-152  
 Project: SR509 Foreman Property

**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:	MB0319S1					
Dichlorodifluoromethane	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
Chloromethane	ND	0.0050	EPA 8260C	3-19-18	3-19-18	
Vinyl Chloride	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
Bromomethane	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
Chloroethane	ND	0.0050	EPA 8260C	3-19-18	3-19-18	
Trichlorofluoromethane	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
1,1-Dichloroethene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
Acetone	ND	0.0050	EPA 8260C	3-19-18	3-19-18	
Iodomethane	ND	0.0050	EPA 8260C	3-19-18	3-19-18	
Carbon Disulfide	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
Methylene Chloride	ND	0.0050	EPA 8260C	3-19-18	3-19-18	
(trans) 1,2-Dichloroethene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
Methyl t-Butyl Ether	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
1,1-Dichloroethane	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
Vinyl Acetate	ND	0.0050	EPA 8260C	3-19-18	3-19-18	
2,2-Dichloropropane	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
(cis) 1,2-Dichloroethene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
2-Butanone	ND	0.0050	EPA 8260C	3-19-18	3-19-18	
Bromochloromethane	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
Chloroform	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
1,1,1-Trichloroethane	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
Carbon Tetrachloride	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
1,1-Dichloropropene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
Benzene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
1,2-Dichloroethane	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
Trichloroethene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
1,2-Dichloropropane	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
Dibromomethane	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
Bromodichloromethane	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
2-Chloroethyl Vinyl Ether	ND	0.0050	EPA 8260C	3-19-18	3-19-18	
(cis) 1,3-Dichloropropene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
Methyl Isobutyl Ketone	ND	0.0050	EPA 8260C	3-19-18	3-19-18	
Toluene	ND	0.0050	EPA 8260C	3-19-18	3-19-18	
(trans) 1,3-Dichloropropene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	



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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:		MB0319S1				
1,1,2-Trichloroethane	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
Tetrachloroethene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
1,3-Dichloropropane	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
2-Hexanone	ND	0.0050	EPA 8260C	3-19-18	3-19-18	
Dibromochloromethane	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
1,2-Dibromoethane	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
Chlorobenzene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
1,1,1,2-Tetrachloroethane	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
Ethylbenzene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
m,p-Xylene	ND	0.0020	EPA 8260C	3-19-18	3-19-18	
o-Xylene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
Styrene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
Bromoform	ND	0.0050	EPA 8260C	3-19-18	3-19-18	
Isopropylbenzene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
Bromobenzene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
1,1,2,2-Tetrachloroethane	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
1,2,3-Trichloropropane	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
n-Propylbenzene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
2-Chlorotoluene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
4-Chlorotoluene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
1,3,5-Trimethylbenzene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
tert-Butylbenzene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
1,2,4-Trimethylbenzene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
sec-Butylbenzene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
1,3-Dichlorobenzene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
p-Isopropyltoluene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
1,4-Dichlorobenzene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
1,2-Dichlorobenzene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
n-Butylbenzene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
1,2-Dibromo-3-chloropropane	ND	0.0050	EPA 8260C	3-19-18	3-19-18	
1,2,4-Trichlorobenzene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
Hexachlorobutadiene	ND	0.0050	EPA 8260C	3-19-18	3-19-18	
Naphthalene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
1,2,3-Trichlorobenzene	ND	0.0010	EPA 8260C	3-19-18	3-19-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>103</i>	<i>75-131</i>				
<i>Toluene-d8</i>	<i>102</i>	<i>83-130</i>				
<i>4-Bromofluorobenzene</i>	<i>103</i>	<i>78-130</i>				



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 Samples Submitted: March 16, 2018  
 Laboratory Reference: 1803-152  
 Project: SR509 Foreman Property

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

Matrix: Soil  
 Units: mg/kg

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD	RPD	Flags
					SB	SBD	Limits	RPD	Limit	
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB0319S1									
	SB	SBD	SB	SBD	SB	SBD				
1,1-Dichloroethene	<b>0.0518</b>	<b>0.0521</b>	0.0500	0.0500	104	104	58-126	1	20	
Benzene	<b>0.0538</b>	<b>0.0549</b>	0.0500	0.0500	108	110	72-122	2	19	
Trichloroethene	<b>0.0503</b>	<b>0.0497</b>	0.0500	0.0500	101	99	75-120	1	20	
Toluene	<b>0.0506</b>	<b>0.0503</b>	0.0500	0.0500	101	101	78-123	1	19	
Chlorobenzene	<b>0.0453</b>	<b>0.0456</b>	0.0500	0.0500	91	91	75-120	1	18	
<i>Surrogate:</i>										
<i>Dibromofluoromethane</i>					<i>109</i>	<i>103</i>	<i>75-131</i>			
<i>Toluene-d8</i>					<i>109</i>	<i>99</i>	<i>83-130</i>			
<i>4-Bromofluorobenzene</i>					<i>105</i>	<i>99</i>	<i>78-130</i>			



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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>SS1 N A'</b>					
<b>Laboratory ID:</b>	<b>03-152-01</b>					
n-Nitrosodimethylamine	ND	0.047	EPA 8270D	3-21-18	3-22-18	
Pyridine	ND	0.47	EPA 8270D	3-21-18	3-22-18	
Phenol	ND	0.047	EPA 8270D	3-21-18	3-22-18	
Aniline	ND	0.24	EPA 8270D	3-21-18	3-22-18	
bis(2-Chloroethyl)ether	ND	0.047	EPA 8270D	3-21-18	3-22-18	
2-Chlorophenol	ND	0.047	EPA 8270D	3-21-18	3-22-18	
1,3-Dichlorobenzene	ND	0.047	EPA 8270D	3-21-18	3-22-18	
1,4-Dichlorobenzene	ND	0.047	EPA 8270D	3-21-18	3-22-18	
Benzyl alcohol	ND	0.24	EPA 8270D	3-21-18	3-22-18	
1,2-Dichlorobenzene	ND	0.047	EPA 8270D	3-21-18	3-22-18	
2-Methylphenol (o-Cresol)	ND	0.047	EPA 8270D	3-21-18	3-22-18	
bis(2-Chloroisopropyl)ether	ND	0.047	EPA 8270D	3-21-18	3-22-18	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.047	EPA 8270D	3-21-18	3-22-18	
n-Nitroso-di-n-propylamine	ND	0.047	EPA 8270D	3-21-18	3-22-18	
Hexachloroethane	ND	0.047	EPA 8270D	3-21-18	3-22-18	
Nitrobenzene	ND	0.047	EPA 8270D	3-21-18	3-22-18	
Isophorone	ND	0.047	EPA 8270D	3-21-18	3-22-18	
2-Nitrophenol	ND	0.047	EPA 8270D	3-21-18	3-22-18	
2,4-Dimethylphenol	ND	0.047	EPA 8270D	3-21-18	3-22-18	
bis(2-Chloroethoxy)methane	ND	0.047	EPA 8270D	3-21-18	3-22-18	
2,4-Dichlorophenol	ND	0.047	EPA 8270D	3-21-18	3-22-18	
1,2,4-Trichlorobenzene	ND	0.047	EPA 8270D	3-21-18	3-22-18	
Naphthalene	ND	0.0094	EPA 8270D/SIM	3-21-18	3-22-18	
4-Chloroaniline	ND	0.24	EPA 8270D	3-21-18	3-22-18	
Hexachlorobutadiene	ND	0.047	EPA 8270D	3-21-18	3-22-18	
4-Chloro-3-methylphenol	ND	0.047	EPA 8270D	3-21-18	3-22-18	
2-Methylnaphthalene	ND	0.0094	EPA 8270D/SIM	3-21-18	3-22-18	
1-Methylnaphthalene	ND	0.0094	EPA 8270D/SIM	3-21-18	3-22-18	
Hexachlorocyclopentadiene	ND	0.047	EPA 8270D	3-21-18	3-22-18	
2,4,6-Trichlorophenol	ND	0.047	EPA 8270D	3-21-18	3-22-18	
2,3-Dichloroaniline	ND	0.047	EPA 8270D	3-21-18	3-22-18	
2,4,5-Trichlorophenol	ND	0.047	EPA 8270D	3-21-18	3-22-18	
2-Chloronaphthalene	ND	0.047	EPA 8270D	3-21-18	3-22-18	
2-Nitroaniline	ND	0.047	EPA 8270D	3-21-18	3-22-18	
1,4-Dinitrobenzene	ND	0.047	EPA 8270D	3-21-18	3-22-18	
Dimethylphthalate	ND	0.047	EPA 8270D	3-21-18	3-22-18	
1,3-Dinitrobenzene	ND	0.047	EPA 8270D	3-21-18	3-22-18	
2,6-Dinitrotoluene	ND	0.047	EPA 8270D	3-21-18	3-22-18	
1,2-Dinitrobenzene	ND	0.047	EPA 8270D	3-21-18	3-22-18	
Acenaphthylene	ND	0.0094	EPA 8270D/SIM	3-21-18	3-22-18	
3-Nitroaniline	ND	0.047	EPA 8270D	3-21-18	3-22-18	



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 Project: SR509 Foreman Property

**SEMIVOLATILES EPA 8270D/SIM**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SS1 N A'</b>					
Laboratory ID:	03-152-01					
2,4-Dinitrophenol	ND	0.24	EPA 8270D	3-21-18	3-22-18	
Acenaphthene	ND	0.0094	EPA 8270D/SIM	3-21-18	3-22-18	
4-Nitrophenol	ND	0.047	EPA 8270D	3-21-18	3-22-18	
2,4-Dinitrotoluene	ND	0.047	EPA 8270D	3-21-18	3-22-18	
Dibenzofuran	ND	0.047	EPA 8270D	3-21-18	3-22-18	
2,3,5,6-Tetrachlorophenol	ND	0.047	EPA 8270D	3-21-18	3-22-18	
2,3,4,6-Tetrachlorophenol	ND	0.047	EPA 8270D	3-21-18	3-22-18	
Diethylphthalate	ND	0.24	EPA 8270D	3-21-18	3-22-18	
4-Chlorophenyl-phenylether	ND	0.047	EPA 8270D	3-21-18	3-22-18	
4-Nitroaniline	ND	0.047	EPA 8270D	3-21-18	3-22-18	
Fluorene	ND	0.0094	EPA 8270D/SIM	3-21-18	3-22-18	
4,6-Dinitro-2-methylphenol	ND	0.24	EPA 8270D	3-21-18	3-22-18	
n-Nitrosodiphenylamine	ND	0.047	EPA 8270D	3-21-18	3-22-18	
1,2-Diphenylhydrazine	ND	0.047	EPA 8270D	3-21-18	3-22-18	
4-Bromophenyl-phenylether	ND	0.047	EPA 8270D	3-21-18	3-22-18	
Hexachlorobenzene	ND	0.047	EPA 8270D	3-21-18	3-22-18	
Pentachlorophenol	ND	0.24	EPA 8270D	3-21-18	3-22-18	
Phenanthrene	ND	0.0094	EPA 8270D/SIM	3-21-18	3-22-18	
Anthracene	ND	0.0094	EPA 8270D/SIM	3-21-18	3-22-18	
Carbazole	ND	0.047	EPA 8270D	3-21-18	3-22-18	
Di-n-butylphthalate	ND	0.24	EPA 8270D	3-21-18	3-22-18	
Fluoranthene	ND	0.0094	EPA 8270D/SIM	3-21-18	3-22-18	
Benzidine	ND	0.47	EPA 8270D	3-21-18	3-22-18	
Pyrene	ND	0.0094	EPA 8270D/SIM	3-21-18	3-22-18	
Butylbenzylphthalate	ND	0.24	EPA 8270D	3-21-18	3-22-18	
bis-2-Ethylhexyladipate	ND	0.24	EPA 8270D	3-21-18	3-22-18	
3,3'-Dichlorobenzidine	ND	0.24	EPA 8270D	3-21-18	3-22-18	
Benzo[a]anthracene	ND	0.0094	EPA 8270D/SIM	3-21-18	3-22-18	
Chrysene	ND	0.0094	EPA 8270D/SIM	3-21-18	3-22-18	
bis(2-Ethylhexyl)phthalate	ND	0.24	EPA 8270D	3-21-18	3-22-18	
Di-n-octylphthalate	ND	0.24	EPA 8270D	3-21-18	3-22-18	
Benzo[b]fluoranthene	ND	0.0094	EPA 8270D/SIM	3-21-18	3-22-18	
Benzo(j,k)fluoranthene	ND	0.0094	EPA 8270D/SIM	3-21-18	3-22-18	
Benzo[a]pyrene	ND	0.0094	EPA 8270D/SIM	3-21-18	3-22-18	
Indeno[1,2,3-cd]pyrene	ND	0.0094	EPA 8270D/SIM	3-21-18	3-22-18	
Dibenz[a,h]anthracene	ND	0.0094	EPA 8270D/SIM	3-21-18	3-22-18	
Benzo[g,h,i]perylene	ND	0.0094	EPA 8270D/SIM	3-21-18	3-22-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	68	18 - 113				
Phenol-d6	71	19 - 119				
Nitrobenzene-d5	71	19 - 119				
2-Fluorobiphenyl	78	33 - 109				
2,4,6-Tribromophenol	80	19 - 121				
Terphenyl-d14	74	30 - 116				



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 Samples Submitted: March 16, 2018  
 Laboratory Reference: 1803-152  
 Project: SR509 Foreman Property

**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>SS2</b>					
<b>Laboratory ID:</b>	<b>03-152-02</b>					
n-Nitrosodimethylamine	ND	0.046	EPA 8270D	3-21-18	3-22-18	
Pyridine	ND	0.46	EPA 8270D	3-21-18	3-22-18	
Phenol	ND	0.046	EPA 8270D	3-21-18	3-22-18	
Aniline	ND	0.23	EPA 8270D	3-21-18	3-22-18	
bis(2-Chloroethyl)ether	ND	0.046	EPA 8270D	3-21-18	3-22-18	
2-Chlorophenol	ND	0.046	EPA 8270D	3-21-18	3-22-18	
1,3-Dichlorobenzene	ND	0.046	EPA 8270D	3-21-18	3-22-18	
1,4-Dichlorobenzene	ND	0.046	EPA 8270D	3-21-18	3-22-18	
Benzyl alcohol	2.9	1.1	EPA 8270D	3-21-18	3-23-18	
1,2-Dichlorobenzene	ND	0.046	EPA 8270D	3-21-18	3-22-18	
2-Methylphenol (o-Cresol)	ND	0.046	EPA 8270D	3-21-18	3-22-18	
bis(2-Chloroisopropyl)ether	ND	0.046	EPA 8270D	3-21-18	3-22-18	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.046	EPA 8270D	3-21-18	3-22-18	
n-Nitroso-di-n-propylamine	ND	0.046	EPA 8270D	3-21-18	3-22-18	
Hexachloroethane	ND	0.046	EPA 8270D	3-21-18	3-22-18	
Nitrobenzene	ND	0.046	EPA 8270D	3-21-18	3-22-18	
Isophorone	ND	0.046	EPA 8270D	3-21-18	3-22-18	
2-Nitrophenol	ND	0.046	EPA 8270D	3-21-18	3-22-18	
2,4-Dimethylphenol	ND	0.046	EPA 8270D	3-21-18	3-22-18	
bis(2-Chloroethoxy)methane	ND	0.046	EPA 8270D	3-21-18	3-22-18	
2,4-Dichlorophenol	ND	0.046	EPA 8270D	3-21-18	3-22-18	
1,2,4-Trichlorobenzene	ND	0.046	EPA 8270D	3-21-18	3-22-18	
Naphthalene	0.016	0.0091	EPA 8270D/SIM	3-21-18	3-22-18	
4-Chloroaniline	ND	0.23	EPA 8270D	3-21-18	3-22-18	
Hexachlorobutadiene	ND	0.046	EPA 8270D	3-21-18	3-22-18	
4-Chloro-3-methylphenol	ND	0.046	EPA 8270D	3-21-18	3-22-18	
2-Methylnaphthalene	ND	0.0091	EPA 8270D/SIM	3-21-18	3-22-18	
1-Methylnaphthalene	ND	0.0091	EPA 8270D/SIM	3-21-18	3-22-18	
Hexachlorocyclopentadiene	ND	0.046	EPA 8270D	3-21-18	3-22-18	
2,4,6-Trichlorophenol	ND	0.046	EPA 8270D	3-21-18	3-22-18	
2,3-Dichloroaniline	ND	0.046	EPA 8270D	3-21-18	3-22-18	
2,4,5-Trichlorophenol	ND	0.046	EPA 8270D	3-21-18	3-22-18	
2-Chloronaphthalene	ND	0.046	EPA 8270D	3-21-18	3-22-18	
2-Nitroaniline	ND	0.046	EPA 8270D	3-21-18	3-22-18	
1,4-Dinitrobenzene	ND	0.046	EPA 8270D	3-21-18	3-22-18	
Dimethylphthalate	ND	0.046	EPA 8270D	3-21-18	3-22-18	
1,3-Dinitrobenzene	ND	0.046	EPA 8270D	3-21-18	3-22-18	
2,6-Dinitrotoluene	ND	0.046	EPA 8270D	3-21-18	3-22-18	
1,2-Dinitrobenzene	ND	0.046	EPA 8270D	3-21-18	3-22-18	
Acenaphthylene	ND	0.0091	EPA 8270D/SIM	3-21-18	3-22-18	
3-Nitroaniline	ND	0.046	EPA 8270D	3-21-18	3-22-18	



Date of Report: March 27, 2018  
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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>SS2</b>					
Laboratory ID:	03-152-02					
2,4-Dinitrophenol	ND	0.23	EPA 8270D	3-21-18	3-22-18	
Acenaphthene	ND	0.0091	EPA 8270D/SIM	3-21-18	3-22-18	
4-Nitrophenol	ND	0.046	EPA 8270D	3-21-18	3-22-18	
2,4-Dinitrotoluene	ND	0.046	EPA 8270D	3-21-18	3-22-18	
Dibenzofuran	ND	0.046	EPA 8270D	3-21-18	3-22-18	
2,3,5,6-Tetrachlorophenol	ND	0.046	EPA 8270D	3-21-18	3-22-18	
2,3,4,6-Tetrachlorophenol	ND	0.046	EPA 8270D	3-21-18	3-22-18	
Diethylphthalate	ND	0.23	EPA 8270D	3-21-18	3-22-18	
4-Chlorophenyl-phenylether	ND	0.046	EPA 8270D	3-21-18	3-22-18	
4-Nitroaniline	ND	0.046	EPA 8270D	3-21-18	3-22-18	
Fluorene	ND	0.0091	EPA 8270D/SIM	3-21-18	3-22-18	
4,6-Dinitro-2-methylphenol	ND	0.23	EPA 8270D	3-21-18	3-22-18	
n-Nitrosodiphenylamine	ND	0.046	EPA 8270D	3-21-18	3-22-18	
1,2-Diphenylhydrazine	ND	0.046	EPA 8270D	3-21-18	3-22-18	
4-Bromophenyl-phenylether	ND	0.046	EPA 8270D	3-21-18	3-22-18	
Hexachlorobenzene	ND	0.046	EPA 8270D	3-21-18	3-22-18	
Pentachlorophenol	ND	0.23	EPA 8270D	3-21-18	3-22-18	
Phenanthrene	0.014	0.0091	EPA 8270D/SIM	3-21-18	3-22-18	
Anthracene	ND	0.0091	EPA 8270D/SIM	3-21-18	3-22-18	
Carbazole	ND	0.046	EPA 8270D	3-21-18	3-22-18	
Di-n-butylphthalate	ND	0.23	EPA 8270D	3-21-18	3-22-18	
Fluoranthene	0.028	0.0091	EPA 8270D/SIM	3-21-18	3-22-18	
Benzidine	ND	0.46	EPA 8270D	3-21-18	3-22-18	
Pyrene	0.023	0.0091	EPA 8270D/SIM	3-21-18	3-22-18	
Butylbenzylphthalate	ND	0.23	EPA 8270D	3-21-18	3-22-18	
bis-2-Ethylhexyladipate	ND	0.23	EPA 8270D	3-21-18	3-22-18	
3,3'-Dichlorobenzidine	ND	0.23	EPA 8270D	3-21-18	3-22-18	
Benzo[a]anthracene	ND	0.0091	EPA 8270D/SIM	3-21-18	3-22-18	
Chrysene	0.022	0.0091	EPA 8270D/SIM	3-21-18	3-22-18	
bis(2-Ethylhexyl)phthalate	ND	0.23	EPA 8270D	3-21-18	3-22-18	
Di-n-octylphthalate	ND	0.23	EPA 8270D	3-21-18	3-22-18	
Benzo[b]fluoranthene	0.028	0.0091	EPA 8270D/SIM	3-21-18	3-22-18	
Benzo(j,k)fluoranthene	ND	0.0091	EPA 8270D/SIM	3-21-18	3-22-18	
Benzo[a]pyrene	0.012	0.0091	EPA 8270D/SIM	3-21-18	3-22-18	
Indeno[1,2,3-cd]pyrene	0.012	0.0091	EPA 8270D/SIM	3-21-18	3-22-18	
Dibenz[a,h]anthracene	ND	0.0091	EPA 8270D/SIM	3-21-18	3-22-18	
Benzo[g,h,i]perylene	0.020	0.0091	EPA 8270D/SIM	3-21-18	3-22-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	61	18 - 113				
Phenol-d6	67	19 - 119				
Nitrobenzene-d5	67	19 - 119				
2-Fluorobiphenyl	77	33 - 109				
2,4,6-Tribromophenol	79	19 - 121				
Terphenyl-d14	76	30 - 116				



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Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SS3</b>					
<b>Laboratory ID:</b>	<b>03-152-03</b>					
n-Nitrosodimethylamine	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Pyridine	ND	0.40	EPA 8270D	3-21-18	3-22-18	
Phenol	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Aniline	ND	0.20	EPA 8270D	3-21-18	3-22-18	
bis(2-Chloroethyl)ether	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2-Chlorophenol	ND	0.040	EPA 8270D	3-21-18	3-22-18	
1,3-Dichlorobenzene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
1,4-Dichlorobenzene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Benzyl alcohol	ND	0.20	EPA 8270D	3-21-18	3-22-18	
1,2-Dichlorobenzene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2-Methylphenol (o-Cresol)	ND	0.040	EPA 8270D	3-21-18	3-22-18	
bis(2-Chloroisopropyl)ether	ND	0.040	EPA 8270D	3-21-18	3-22-18	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.040	EPA 8270D	3-21-18	3-22-18	
n-Nitroso-di-n-propylamine	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Hexachloroethane	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Nitrobenzene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Isophorone	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2-Nitrophenol	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2,4-Dimethylphenol	ND	0.040	EPA 8270D	3-21-18	3-22-18	
bis(2-Chloroethoxy)methane	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2,4-Dichlorophenol	ND	0.040	EPA 8270D	3-21-18	3-22-18	
1,2,4-Trichlorobenzene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Naphthalene	ND	0.0080	EPA 8270D/SIM	3-21-18	3-22-18	
4-Chloroaniline	ND	0.20	EPA 8270D	3-21-18	3-22-18	
Hexachlorobutadiene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
4-Chloro-3-methylphenol	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2-Methylnaphthalene	ND	0.0080	EPA 8270D/SIM	3-21-18	3-22-18	
1-Methylnaphthalene	ND	0.0080	EPA 8270D/SIM	3-21-18	3-22-18	
Hexachlorocyclopentadiene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2,4,6-Trichlorophenol	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2,3-Dichloroaniline	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2,4,5-Trichlorophenol	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2-Chloronaphthalene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2-Nitroaniline	ND	0.040	EPA 8270D	3-21-18	3-22-18	
1,4-Dinitrobenzene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Dimethylphthalate	0.12	0.040	EPA 8270D	3-21-18	3-22-18	
1,3-Dinitrobenzene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2,6-Dinitrotoluene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
1,2-Dinitrobenzene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Acenaphthylene	ND	0.0080	EPA 8270D/SIM	3-21-18	3-22-18	
3-Nitroaniline	ND	0.040	EPA 8270D	3-21-18	3-22-18	



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 Laboratory Reference: 1803-152  
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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>SS3</b>					
Laboratory ID:	03-152-03					
2,4-Dinitrophenol	ND	0.20	EPA 8270D	3-21-18	3-22-18	
Acenaphthene	ND	0.0080	EPA 8270D/SIM	3-21-18	3-22-18	
4-Nitrophenol	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2,4-Dinitrotoluene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Dibenzofuran	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2,3,5,6-Tetrachlorophenol	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2,3,4,6-Tetrachlorophenol	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Diethylphthalate	ND	0.20	EPA 8270D	3-21-18	3-22-18	
4-Chlorophenyl-phenylether	ND	0.040	EPA 8270D	3-21-18	3-22-18	
4-Nitroaniline	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Fluorene	ND	0.0080	EPA 8270D/SIM	3-21-18	3-22-18	
4,6-Dinitro-2-methylphenol	ND	0.20	EPA 8270D	3-21-18	3-22-18	
n-Nitrosodiphenylamine	ND	0.040	EPA 8270D	3-21-18	3-22-18	
1,2-Diphenylhydrazine	ND	0.040	EPA 8270D	3-21-18	3-22-18	
4-Bromophenyl-phenylether	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Hexachlorobenzene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Pentachlorophenol	ND	0.20	EPA 8270D	3-21-18	3-22-18	
Phenanthrene	ND	0.0080	EPA 8270D/SIM	3-21-18	3-22-18	
Anthracene	ND	0.0080	EPA 8270D/SIM	3-21-18	3-22-18	
Carbazole	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Di-n-butylphthalate	ND	0.20	EPA 8270D	3-21-18	3-22-18	
Fluoranthene	0.013	0.0080	EPA 8270D/SIM	3-21-18	3-22-18	
Benzidine	ND	0.40	EPA 8270D	3-21-18	3-22-18	
Pyrene	0.012	0.0080	EPA 8270D/SIM	3-21-18	3-22-18	
Butylbenzylphthalate	0.51	0.20	EPA 8270D	3-21-18	3-22-18	
bis-2-Ethylhexyladipate	ND	0.20	EPA 8270D	3-21-18	3-22-18	
3,3'-Dichlorobenzidine	ND	0.20	EPA 8270D	3-21-18	3-22-18	
Benzo[a]anthracene	ND	0.0080	EPA 8270D/SIM	3-21-18	3-22-18	
Chrysene	0.0092	0.0080	EPA 8270D/SIM	3-21-18	3-22-18	
bis(2-Ethylhexyl)phthalate	0.25	0.20	EPA 8270D	3-21-18	3-22-18	
Di-n-octylphthalate	ND	0.20	EPA 8270D	3-21-18	3-22-18	
Benzo[b]fluoranthene	0.011	0.0080	EPA 8270D/SIM	3-21-18	3-22-18	
Benzo(j,k)fluoranthene	ND	0.0080	EPA 8270D/SIM	3-21-18	3-22-18	
Benzo[a]pyrene	ND	0.0080	EPA 8270D/SIM	3-21-18	3-22-18	
Indeno[1,2,3-cd]pyrene	ND	0.0080	EPA 8270D/SIM	3-21-18	3-22-18	
Dibenz[a,h]anthracene	ND	0.0080	EPA 8270D/SIM	3-21-18	3-22-18	
Benzo[g,h,i]perylene	0.014	0.0080	EPA 8270D/SIM	3-21-18	3-22-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	55	18 - 113				
Phenol-d6	68	19 - 119				
Nitrobenzene-d5	61	19 - 119				
2-Fluorobiphenyl	76	33 - 109				
2,4,6-Tribromophenol	80	19 - 121				
Terphenyl-d14	77	30 - 116				



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Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SS4</b>					
<b>Laboratory ID:</b>	<b>03-152-04</b>					
n-Nitrosodimethylamine	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Pyridine	ND	0.40	EPA 8270D	3-21-18	3-22-18	
Phenol	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Aniline	ND	0.20	EPA 8270D	3-21-18	3-22-18	
bis(2-Chloroethyl)ether	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2-Chlorophenol	ND	0.040	EPA 8270D	3-21-18	3-22-18	
1,3-Dichlorobenzene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
1,4-Dichlorobenzene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Benzyl alcohol	ND	0.20	EPA 8270D	3-21-18	3-22-18	
1,2-Dichlorobenzene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2-Methylphenol (o-Cresol)	ND	0.040	EPA 8270D	3-21-18	3-22-18	
bis(2-Chloroisopropyl)ether	ND	0.040	EPA 8270D	3-21-18	3-22-18	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.040	EPA 8270D	3-21-18	3-22-18	
n-Nitroso-di-n-propylamine	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Hexachloroethane	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Nitrobenzene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Isophorone	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2-Nitrophenol	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2,4-Dimethylphenol	ND	0.040	EPA 8270D	3-21-18	3-22-18	
bis(2-Chloroethoxy)methane	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2,4-Dichlorophenol	ND	0.040	EPA 8270D	3-21-18	3-22-18	
1,2,4-Trichlorobenzene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Naphthalene	ND	0.0081	EPA 8270D/SIM	3-21-18	3-22-18	
4-Chloroaniline	ND	0.20	EPA 8270D	3-21-18	3-22-18	
Hexachlorobutadiene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
4-Chloro-3-methylphenol	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2-Methylnaphthalene	ND	0.0081	EPA 8270D/SIM	3-21-18	3-22-18	
1-Methylnaphthalene	ND	0.0081	EPA 8270D/SIM	3-21-18	3-22-18	
Hexachlorocyclopentadiene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2,4,6-Trichlorophenol	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2,3-Dichloroaniline	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2,4,5-Trichlorophenol	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2-Chloronaphthalene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2-Nitroaniline	ND	0.040	EPA 8270D	3-21-18	3-22-18	
1,4-Dinitrobenzene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Dimethylphthalate	ND	0.040	EPA 8270D	3-21-18	3-22-18	
1,3-Dinitrobenzene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2,6-Dinitrotoluene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
1,2-Dinitrobenzene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Acenaphthylene	ND	0.0081	EPA 8270D/SIM	3-21-18	3-22-18	
3-Nitroaniline	ND	0.040	EPA 8270D	3-21-18	3-22-18	



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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SS4</b>					
Laboratory ID:	03-152-04					
2,4-Dinitrophenol	ND	0.20	EPA 8270D	3-21-18	3-22-18	
Acenaphthene	ND	0.0081	EPA 8270D/SIM	3-21-18	3-22-18	
4-Nitrophenol	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2,4-Dinitrotoluene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Dibenzofuran	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2,3,5,6-Tetrachlorophenol	ND	0.040	EPA 8270D	3-21-18	3-22-18	
2,3,4,6-Tetrachlorophenol	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Diethylphthalate	ND	0.20	EPA 8270D	3-21-18	3-22-18	
4-Chlorophenyl-phenylether	ND	0.040	EPA 8270D	3-21-18	3-22-18	
4-Nitroaniline	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Fluorene	ND	0.0081	EPA 8270D/SIM	3-21-18	3-22-18	
4,6-Dinitro-2-methylphenol	ND	0.20	EPA 8270D	3-21-18	3-22-18	
n-Nitrosodiphenylamine	ND	0.040	EPA 8270D	3-21-18	3-22-18	
1,2-Diphenylhydrazine	ND	0.040	EPA 8270D	3-21-18	3-22-18	
4-Bromophenyl-phenylether	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Hexachlorobenzene	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Pentachlorophenol	ND	0.20	EPA 8270D	3-21-18	3-22-18	
Phenanthrene	<b>0.0097</b>	0.0081	EPA 8270D/SIM	3-21-18	3-22-18	
Anthracene	ND	0.0081	EPA 8270D/SIM	3-21-18	3-22-18	
Carbazole	ND	0.040	EPA 8270D	3-21-18	3-22-18	
Di-n-butylphthalate	ND	0.20	EPA 8270D	3-21-18	3-22-18	
Fluoranthene	<b>0.012</b>	0.0081	EPA 8270D/SIM	3-21-18	3-22-18	
Benzidine	ND	0.40	EPA 8270D	3-21-18	3-22-18	
Pyrene	<b>0.013</b>	0.0081	EPA 8270D/SIM	3-21-18	3-22-18	
Butylbenzylphthalate	ND	0.20	EPA 8270D	3-21-18	3-22-18	
bis-2-Ethylhexyladipate	ND	0.20	EPA 8270D	3-21-18	3-22-18	
3,3'-Dichlorobenzidine	ND	0.20	EPA 8270D	3-21-18	3-22-18	
Benzo[a]anthracene	ND	0.0081	EPA 8270D/SIM	3-21-18	3-22-18	
Chrysene	<b>0.0084</b>	0.0081	EPA 8270D/SIM	3-21-18	3-22-18	
bis(2-Ethylhexyl)phthalate	ND	0.20	EPA 8270D	3-21-18	3-22-18	
Di-n-octylphthalate	ND	0.20	EPA 8270D	3-21-18	3-22-18	
Benzo[b]fluoranthene	<b>0.011</b>	0.0081	EPA 8270D/SIM	3-21-18	3-22-18	
Benzo(j,k)fluoranthene	ND	0.0081	EPA 8270D/SIM	3-21-18	3-22-18	
Benzo[a]pyrene	ND	0.0081	EPA 8270D/SIM	3-21-18	3-22-18	
Indeno[1,2,3-cd]pyrene	ND	0.0081	EPA 8270D/SIM	3-21-18	3-22-18	
Dibenz[a,h]anthracene	ND	0.0081	EPA 8270D/SIM	3-21-18	3-22-18	
Benzo[g,h,i]perylene	ND	0.0081	EPA 8270D/SIM	3-21-18	3-22-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	63	18 - 113				
Phenol-d6	67	19 - 119				
Nitrobenzene-d5	68	19 - 119				
2-Fluorobiphenyl	75	33 - 109				
2,4,6-Tribromophenol	74	19 - 121				
Terphenyl-d14	71	30 - 116				



Date of Report: March 27, 2018  
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 Project: SR509 Foreman Property

**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:	MB0321S1					
n-Nitrosodimethylamine	ND	0.033	EPA 8270D	3-21-18	3-21-18	
Pyridine	ND	0.33	EPA 8270D	3-21-18	3-21-18	
Phenol	ND	0.033	EPA 8270D	3-21-18	3-21-18	
Aniline	ND	0.17	EPA 8270D	3-21-18	3-21-18	
bis(2-Chloroethyl)ether	ND	0.033	EPA 8270D	3-21-18	3-21-18	
2-Chlorophenol	ND	0.033	EPA 8270D	3-21-18	3-21-18	
1,3-Dichlorobenzene	ND	0.033	EPA 8270D	3-21-18	3-21-18	
1,4-Dichlorobenzene	ND	0.033	EPA 8270D	3-21-18	3-21-18	
Benzyl alcohol	ND	0.17	EPA 8270D	3-21-18	3-21-18	
1,2-Dichlorobenzene	ND	0.033	EPA 8270D	3-21-18	3-21-18	
2-Methylphenol (o-Cresol)	ND	0.033	EPA 8270D	3-21-18	3-21-18	
bis(2-Chloroisopropyl)ether	ND	0.033	EPA 8270D	3-21-18	3-21-18	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.033	EPA 8270D	3-21-18	3-21-18	
n-Nitroso-di-n-propylamine	ND	0.033	EPA 8270D	3-21-18	3-21-18	
Hexachloroethane	ND	0.033	EPA 8270D	3-21-18	3-21-18	
Nitrobenzene	ND	0.033	EPA 8270D	3-21-18	3-21-18	
Isophorone	ND	0.033	EPA 8270D	3-21-18	3-21-18	
2-Nitrophenol	ND	0.033	EPA 8270D	3-21-18	3-21-18	
2,4-Dimethylphenol	ND	0.033	EPA 8270D	3-21-18	3-21-18	
bis(2-Chloroethoxy)methane	ND	0.033	EPA 8270D	3-21-18	3-21-18	
2,4-Dichlorophenol	ND	0.033	EPA 8270D	3-21-18	3-21-18	
1,2,4-Trichlorobenzene	ND	0.033	EPA 8270D	3-21-18	3-21-18	
Naphthalene	ND	0.0067	EPA 8270D/SIM	3-21-18	3-21-18	
4-Chloroaniline	ND	0.17	EPA 8270D	3-21-18	3-21-18	
Hexachlorobutadiene	ND	0.033	EPA 8270D	3-21-18	3-21-18	
4-Chloro-3-methylphenol	ND	0.033	EPA 8270D	3-21-18	3-21-18	
2-Methylnaphthalene	ND	0.0067	EPA 8270D/SIM	3-21-18	3-21-18	
1-Methylnaphthalene	ND	0.0067	EPA 8270D/SIM	3-21-18	3-21-18	
Hexachlorocyclopentadiene	ND	0.033	EPA 8270D	3-21-18	3-21-18	
2,4,6-Trichlorophenol	ND	0.033	EPA 8270D	3-21-18	3-21-18	
2,3-Dichloroaniline	ND	0.033	EPA 8270D	3-21-18	3-21-18	
2,4,5-Trichlorophenol	ND	0.033	EPA 8270D	3-21-18	3-21-18	
2-Chloronaphthalene	ND	0.033	EPA 8270D	3-21-18	3-21-18	
2-Nitroaniline	ND	0.033	EPA 8270D	3-21-18	3-21-18	
1,4-Dinitrobenzene	ND	0.033	EPA 8270D	3-21-18	3-21-18	
Dimethylphthalate	ND	0.033	EPA 8270D	3-21-18	3-21-18	
1,3-Dinitrobenzene	ND	0.033	EPA 8270D	3-21-18	3-21-18	
2,6-Dinitrotoluene	ND	0.033	EPA 8270D	3-21-18	3-21-18	
1,2-Dinitrobenzene	ND	0.033	EPA 8270D	3-21-18	3-21-18	
Acenaphthylene	ND	0.0067	EPA 8270D/SIM	3-21-18	3-21-18	
3-Nitroaniline	ND	0.033	EPA 8270D	3-21-18	3-21-18	



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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:	MB0321S1					
2,4-Dinitrophenol	ND	0.17	EPA 8270D	3-21-18	3-21-18	
Acenaphthene	ND	0.0067	EPA 8270D/SIM	3-21-18	3-21-18	
4-Nitrophenol	ND	0.033	EPA 8270D	3-21-18	3-21-18	
2,4-Dinitrotoluene	ND	0.033	EPA 8270D	3-21-18	3-21-18	
Dibenzofuran	ND	0.033	EPA 8270D	3-21-18	3-21-18	
2,3,5,6-Tetrachlorophenol	ND	0.033	EPA 8270D	3-21-18	3-21-18	
2,3,4,6-Tetrachlorophenol	ND	0.033	EPA 8270D	3-21-18	3-21-18	
Diethylphthalate	ND	0.17	EPA 8270D	3-21-18	3-21-18	
4-Chlorophenyl-phenylether	ND	0.033	EPA 8270D	3-21-18	3-21-18	
4-Nitroaniline	ND	0.033	EPA 8270D	3-21-18	3-21-18	
Fluorene	ND	0.0067	EPA 8270D/SIM	3-21-18	3-21-18	
4,6-Dinitro-2-methylphenol	ND	0.17	EPA 8270D	3-21-18	3-21-18	
n-Nitrosodiphenylamine	ND	0.033	EPA 8270D	3-21-18	3-21-18	
1,2-Diphenylhydrazine	ND	0.033	EPA 8270D	3-21-18	3-21-18	
4-Bromophenyl-phenylether	ND	0.033	EPA 8270D	3-21-18	3-21-18	
Hexachlorobenzene	ND	0.033	EPA 8270D	3-21-18	3-21-18	
Pentachlorophenol	ND	0.17	EPA 8270D	3-21-18	3-21-18	
Phenanthrene	ND	0.0067	EPA 8270D/SIM	3-21-18	3-21-18	
Anthracene	ND	0.0067	EPA 8270D/SIM	3-21-18	3-21-18	
Carbazole	ND	0.033	EPA 8270D	3-21-18	3-21-18	
Di-n-butylphthalate	ND	0.17	EPA 8270D	3-21-18	3-21-18	
Fluoranthene	ND	0.0067	EPA 8270D/SIM	3-21-18	3-21-18	
Benzidine	ND	0.33	EPA 8270D	3-21-18	3-21-18	
Pyrene	ND	0.0067	EPA 8270D/SIM	3-21-18	3-21-18	
Butylbenzylphthalate	ND	0.17	EPA 8270D	3-21-18	3-21-18	
bis-2-Ethylhexyladipate	ND	0.17	EPA 8270D	3-21-18	3-21-18	
3,3'-Dichlorobenzidine	ND	0.17	EPA 8270D	3-21-18	3-21-18	
Benzo[a]anthracene	ND	0.0067	EPA 8270D/SIM	3-21-18	3-21-18	
Chrysene	ND	0.0067	EPA 8270D/SIM	3-21-18	3-21-18	
bis(2-Ethylhexyl)phthalate	ND	0.17	EPA 8270D	3-21-18	3-21-18	
Di-n-octylphthalate	ND	0.17	EPA 8270D	3-21-18	3-21-18	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270D/SIM	3-21-18	3-21-18	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270D/SIM	3-21-18	3-21-18	
Benzo[a]pyrene	ND	0.0067	EPA 8270D/SIM	3-21-18	3-21-18	
Indeno[1,2,3-cd]pyrene	ND	0.0067	EPA 8270D/SIM	3-21-18	3-21-18	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270D/SIM	3-21-18	3-21-18	
Benzo[g,h,i]perylene	ND	0.0067	EPA 8270D/SIM	3-21-18	3-21-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	85	18 - 113				
Phenol-d6	86	19 - 119				
Nitrobenzene-d5	85	19 - 119				
2-Fluorobiphenyl	89	33 - 109				
2,4,6-Tribromophenol	85	19 - 121				
Terphenyl-d14	81	30 - 116				



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 Laboratory Reference: 1803-152  
 Project: SR509 Foreman Property

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

Matrix: Soil  
 Units: mg/Kg

Analyte	Result		Spike Level		Percent Recovery		Recovery Limits	RPD	RPD Limit	Flags
	SB	SBD	SB	SBD	SB	SBD				
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB0321S1									
Phenol	1.11	1.15	1.33	1.33	83	86	39 - 109	4	36	
2-Chlorophenol	1.18	1.28	1.33	1.33	89	96	42 - 105	8	35	
1,4-Dichlorobenzene	0.561	0.613	0.667	0.667	84	92	31 - 103	9	37	
n-Nitroso-di-n-propylamine	0.598	0.615	0.667	0.667	90	92	36 - 104	3	34	
1,2,4-Trichlorobenzene	0.601	0.618	0.667	0.667	90	93	32 - 104	3	38	
4-Chloro-3-methylphenol	1.17	1.21	1.33	1.33	88	91	48 - 107	3	31	
Acenaphthene	0.554	0.597	0.667	0.667	83	90	38 - 102	7	33	
4-Nitrophenol	1.21	1.23	1.33	1.33	91	92	27 - 121	2	35	
2,4-Dinitrotoluene	0.551	0.582	0.667	0.667	83	87	36 - 103	5	34	
Pentachlorophenol	1.43	1.53	1.33	1.33	108	115	21 - 114	7	37	I
Pyrene	0.585	0.619	0.667	0.667	88	93	46 - 108	6	31	
<i>Surrogate:</i>										
2-Fluorophenol					86	90	18 - 113			
Phenol-d6					87	90	19 - 119			
Nitrobenzene-d5					84	87	19 - 119			
2-Fluorobiphenyl					85	86	33 - 109			
2,4,6-Tribromophenol					90	92	19 - 121			
Terphenyl-d14					83	86	30 - 116			



Date of Report: March 27, 2018  
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 Project: SR509 Foreman Property

**TOTAL METALS  
 EPA 6010D/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:	03-152-01					
<b>Client ID:</b>	<b>SS1 N A'</b>					
Arsenic	ND	14	6010D	3-21-18	3-21-18	
Barium	84	3.5	6010D	3-21-18	3-21-18	
Cadmium	ND	0.71	6010D	3-21-18	3-21-18	
Chromium	41	0.71	6010D	3-21-18	3-21-18	
Lead	ND	7.1	6010D	3-21-18	3-21-18	
Mercury	ND	0.35	7471B	3-19-18	3-19-18	
Selenium	ND	14	6010D	3-21-18	3-21-18	
Silver	ND	1.4	6010D	3-21-18	3-21-18	

Lab ID:	03-152-02					
<b>Client ID:</b>	<b>SS2</b>					
Arsenic	ND	14	6010D	3-21-18	3-21-18	
Barium	110	3.4	6010D	3-21-18	3-21-18	
Cadmium	ND	0.68	6010D	3-21-18	3-21-18	
Chromium	52	0.68	6010D	3-21-18	3-21-18	
Lead	53	6.8	6010D	3-21-18	3-21-18	
Mercury	ND	0.34	7471B	3-19-18	3-19-18	
Selenium	ND	14	6010D	3-21-18	3-21-18	
Silver	ND	1.4	6010D	3-21-18	3-21-18	



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**TOTAL METALS  
 EPA 6010D/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:	03-152-03					
<b>Client ID:</b>	<b>SS3</b>					
Arsenic	<b>ND</b>	12	6010D	3-21-18	3-21-18	
Barium	<b>61</b>	3.0	6010D	3-21-18	3-21-18	
Cadmium	<b>ND</b>	0.60	6010D	3-21-18	3-21-18	
Chromium	<b>47</b>	0.60	6010D	3-21-18	3-21-18	
Lead	<b>27</b>	6.0	6010D	3-21-18	3-21-18	
Mercury	<b>ND</b>	0.30	7471B	3-19-18	3-19-18	
Selenium	<b>ND</b>	12	6010D	3-21-18	3-21-18	
Silver	<b>ND</b>	1.2	6010D	3-21-18	3-21-18	

Lab ID:	03-152-04					
<b>Client ID:</b>	<b>SS4</b>					
Arsenic	<b>ND</b>	12	6010D	3-21-18	3-21-18	
Barium	<b>160</b>	3.0	6010D	3-21-18	3-21-18	
Cadmium	<b>ND</b>	0.61	6010D	3-21-18	3-21-18	
Chromium	<b>60</b>	0.61	6010D	3-21-18	3-21-18	
Lead	<b>23</b>	6.1	6010D	3-21-18	3-21-18	
Mercury	<b>ND</b>	0.30	7471B	3-19-18	3-19-18	
Selenium	<b>ND</b>	12	6010D	3-21-18	3-21-18	
Silver	<b>ND</b>	1.2	6010D	3-21-18	3-21-18	



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 Project: SR509 Foreman Property

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

Date Extracted: 3-21-18  
 Date Analyzed: 3-21-18  
 Matrix: Soil  
 Units: mg/kg (ppm)  
 Lab ID: MB0321SM1

Analyte	Method	Result	PQL
Arsenic	6010D	ND	10
Barium	6010D	ND	2.5
Cadmium	6010D	ND	0.50
Chromium	6010D	ND	0.50
Lead	6010D	ND	5.0
Selenium	6010D	ND	10
Silver	6010D	ND	1.0



Date of Report: March 27, 2018  
Samples Submitted: March 16, 2018  
Laboratory Reference: 1803-152  
Project: SR509 Foreman Property

**TOTAL MERCURY  
EPA 7471B  
METHOD BLANK QUALITY CONTROL**

Date Extracted: 3-19-18  
Date Analyzed: 3-19-18  
  
Matrix: Soil  
Units: mg/kg (ppm)  
  
Lab ID: MB0319S1

Analyte	Method	Result	PQL
Mercury	7471B	<b>ND</b>	0.25



Date of Report: March 27, 2018  
 Samples Submitted: March 16, 2018  
 Laboratory Reference: 1803-152  
 Project: SR509 Foreman Property

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

Date Extracted: 3-21-18

Date Analyzed: 3-21-18

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-145-02

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Arsenic	ND	ND	NA	10	
Barium	83.0	68.1	20	2.5	
Cadmium	ND	ND	NA	0.50	
Chromium	30.1	36.9	20	0.50	
Lead	57.3	49.0	16	5.0	
Selenium	ND	ND	NA	10	
Silver	ND	ND	NA	1.0	



Date of Report: March 27, 2018  
Samples Submitted: March 16, 2018  
Laboratory Reference: 1803-152  
Project: SR509 Foreman Property

**TOTAL MERCURY  
EPA 7471B  
DUPLICATE QUALITY CONTROL**

Date Extracted: 3-19-18

Date Analyzed: 3-19-18

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-145-01

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Mercury	<b>ND</b>	<b>ND</b>	NA	0.25	



Date of Report: March 27, 2018  
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 Laboratory Reference: 1803-152  
 Project: SR509 Foreman Property

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

Date Extracted: 3-21-18

Date Analyzed: 3-21-18

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-145-02

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Arsenic	100	<b>95.4</b>	95	<b>90.7</b>	91	5	
Barium	100	<b>175</b>	92	<b>181</b>	98	4	
Cadmium	50.0	<b>46.4</b>	93	<b>45.9</b>	92	1	
Chromium	100	<b>123</b>	93	<b>122</b>	92	1	
Lead	250	<b>291</b>	93	<b>290</b>	93	0	
Selenium	100	<b>97.3</b>	97	<b>94.8</b>	95	3	
Silver	25.0	<b>19.3</b>	77	<b>18.8</b>	75	3	



Date of Report: March 27, 2018  
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**TOTAL MERCURY  
 EPA 7471B  
 MS/MSD QUALITY CONTROL**

Date Extracted: 3-19-18

Date Analyzed: 3-19-18

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-145-01

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Mercury	0.500	<b>0.510</b>	102	<b>0.490</b>	98	4	



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Project: SR509 Foreman Property

### % MOISTURE

Date Analyzed: 3-19-18

Client ID	Lab ID	% Moisture
SS1 N A'	03-152-01	29
SS2	03-152-02	27
SS3	03-152-03	16
SS4	03-152-04	18





### Data Qualifiers and Abbreviations

- A - Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
  - B - The analyte indicated was also found in the blank sample.
  - C - The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
  - E - The value reported exceeds the quantitation range and is an estimate.
  - F - Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
  - H - The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
  - I - Compound recovery is outside of the control limits.
  - J - The value reported was below the practical quantitation limit. The value is an estimate.
  - K - Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
  - L - The RPD is outside of the control limits.
  - M - Hydrocarbons in the gasoline range are impacting the diesel range result.
  - M1 - Hydrocarbons in the gasoline range (toluene-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)  
(TPH analysis 5 Days)

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

**Laboratory Number:**

**03-152**

Company: **WSDOT**  
Project Number:  
Project Name: **SR509 Foreman Property**  
Project Manager: **PATRICK SVOBODA**  
Sampled by:

Lab ID      Sample Identification      Date Sampled      Time Sampled      Matrix

1	SS1NPA	3/15/18	12:30	S
2	SS2	3/15/18	1:00	S
3	SS3	3/15/18	2:00	S
4	SS4	3/15/18	2:15	S

**Number of Containers**

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

**Signature**

**Company**

**Date**

**Time**

**Comments/Special Instructions**

Relinquished	<i>[Signature]</i>	WSDOT	3/16/18	8:30	
Received	<i>[Signature]</i>	Alpha	3/16/18	8:30	
Relinquished	<i>[Signature]</i>	Alpha	3/16/18	11:23	
Received	<i>[Signature]</i>				
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

September 24, 2018

Patrick Svoboda  
WSDOT  
15700 Dayton Avenue North  
NB82-138  
P.O. Box 330310  
Seattle, WA 98133-9710

Re: Analytical Data for Project Foreman  
Laboratory Reference No. 1809-229

Dear Pat:

Enclosed are the analytical results and associated quality control data for samples submitted on September 21, 2018.

The standard policy of OnSite Environmental, Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

A handwritten signature in black ink, appearing to read "DB", with a long horizontal flourish 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: September 24, 2018  
Samples Submitted: September 21, 2018  
Laboratory Reference: 1809-229  
Project: Foreman

### Case Narrative

Samples were collected on September 20 and 21, 2018 and received by the laboratory on September 21, 2018. They were maintained at the laboratory at a temperature of 2°C to 6°C.

Please note that any and all soil sample results are reported on a dry-weight basis, unless otherwise noted below.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.





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

September 24, 2018

Patrick Svoboda  
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15700 Dayton Avenue North  
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Re: Analytical Data for Project Foreman  
Laboratory Reference No. 1809-229

Dear Pat:

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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



---

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

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Project: Foreman

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Samples were collected on September 20 and 21, 2018 and received by the laboratory on September 21, 2018. They were maintained at the laboratory at a temperature of 2°C to 6°C.

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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: September 24, 2018  
 Samples Submitted: September 21, 2018  
 Laboratory Reference: 1809-229  
 Project: Foreman

**DIESEL AND HEAVY OIL RANGE ORGANICS  
 NWTPH-Dx**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>FC1-1</b>					
Laboratory ID:	09-229-01					
Diesel Range Organics	<b>ND</b>	41	NWTPH-Dx	9-21-18	9-21-18	U1
Lube Oil	<b>590</b>	57	NWTPH-Dx	9-21-18	9-21-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>88</i>	<i>50-150</i>				

<b>Client ID:</b>	<b>FC1-2</b>					
Laboratory ID:	09-229-02					
Diesel Range Organics	<b>ND</b>	27	NWTPH-Dx	9-21-18	9-21-18	
Lube Oil Range Organics	<b>ND</b>	54	NWTPH-Dx	9-21-18	9-21-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>81</i>	<i>50-150</i>				



Date of Report: September 24, 2018  
 Samples Submitted: September 21, 2018  
 Laboratory Reference: 1809-229  
 Project: Foreman

**DIESEL AND HEAVY OIL RANGE ORGANICS  
 NWTPH-Dx  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB0921S3					
Diesel Range Organics	<b>ND</b>	25	NWTPH-Dx	9-21-18	9-21-18	
Lube Oil Range Organics	<b>ND</b>	50	NWTPH-Dx	9-21-18	9-21-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	88	50-150				

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>								
Laboratory ID:	09-217-01							
	ORIG	DUP						
Diesel Range	<b>ND</b>	<b>ND</b>	NA	NA	NA	NA	NA	NA
Lube Oil Range Organics	<b>178</b>	<b>164</b>	NA	NA	NA	NA	8	NA
<i>Surrogate:</i>								
<i>o-Terphenyl</i>				82	80	50-150		



Date of Report: September 24, 2018  
 Samples Submitted: September 21, 2018  
 Laboratory Reference: 1809-229  
 Project: Foreman

**SEMIVOLATILE ORGANICS 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>FC1-1</b>					
<b>Laboratory ID:</b>	<b>09-229-01</b>					
n-Nitrosodimethylamine	ND	0.38	EPA 8270D	9-21-18	9-21-18	
Pyridine	ND	3.8	EPA 8270D	9-21-18	9-21-18	
Phenol	ND	0.38	EPA 8270D	9-21-18	9-21-18	
Aniline	ND	1.9	EPA 8270D	9-21-18	9-21-18	
bis(2-Chloroethyl)ether	ND	0.38	EPA 8270D	9-21-18	9-21-18	
2-Chlorophenol	ND	0.38	EPA 8270D	9-21-18	9-21-18	
1,3-Dichlorobenzene	ND	0.38	EPA 8270D	9-21-18	9-21-18	
1,4-Dichlorobenzene	ND	0.38	EPA 8270D	9-21-18	9-21-18	
Benzyl alcohol	ND	1.9	EPA 8270D	9-21-18	9-21-18	
1,2-Dichlorobenzene	ND	0.38	EPA 8270D	9-21-18	9-21-18	
2-Methylphenol (o-Cresol)	ND	0.38	EPA 8270D	9-21-18	9-21-18	
bis(2-Chloroisopropyl)ether	ND	0.38	EPA 8270D	9-21-18	9-21-18	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.38	EPA 8270D	9-21-18	9-21-18	
n-Nitroso-di-n-propylamine	ND	0.38	EPA 8270D	9-21-18	9-21-18	
Hexachloroethane	ND	0.38	EPA 8270D	9-21-18	9-21-18	
Nitrobenzene	ND	0.38	EPA 8270D	9-21-18	9-21-18	
Isophorone	ND	0.38	EPA 8270D	9-21-18	9-21-18	
2-Nitrophenol	ND	0.38	EPA 8270D	9-21-18	9-21-18	
2,4-Dimethylphenol	ND	0.38	EPA 8270D	9-21-18	9-21-18	
bis(2-Chloroethoxy)methane	ND	0.38	EPA 8270D	9-21-18	9-21-18	
2,4-Dichlorophenol	ND	0.38	EPA 8270D	9-21-18	9-21-18	
1,2,4-Trichlorobenzene	ND	0.38	EPA 8270D	9-21-18	9-21-18	
Naphthalene	ND	0.015	EPA 8270D/SIM	9-21-18	9-21-18	
4-Chloroaniline	ND	1.9	EPA 8270D	9-21-18	9-21-18	
Hexachlorobutadiene	ND	0.38	EPA 8270D	9-21-18	9-21-18	
4-Chloro-3-methylphenol	ND	0.38	EPA 8270D	9-21-18	9-21-18	
2-Methylnaphthalene	ND	0.015	EPA 8270D/SIM	9-21-18	9-21-18	
1-Methylnaphthalene	ND	0.015	EPA 8270D/SIM	9-21-18	9-21-18	
Hexachlorocyclopentadiene	ND	0.38	EPA 8270D	9-21-18	9-21-18	
2,4,6-Trichlorophenol	ND	0.38	EPA 8270D	9-21-18	9-21-18	
2,3-Dichloroaniline	ND	0.38	EPA 8270D	9-21-18	9-21-18	
2,4,5-Trichlorophenol	ND	0.38	EPA 8270D	9-21-18	9-21-18	
2-Chloronaphthalene	ND	0.38	EPA 8270D	9-21-18	9-21-18	
2-Nitroaniline	ND	0.38	EPA 8270D	9-21-18	9-21-18	
1,4-Dinitrobenzene	ND	0.38	EPA 8270D	9-21-18	9-21-18	
Dimethylphthalate	ND	0.38	EPA 8270D	9-21-18	9-21-18	
1,3-Dinitrobenzene	ND	0.38	EPA 8270D	9-21-18	9-21-18	
2,6-Dinitrotoluene	ND	0.38	EPA 8270D	9-21-18	9-21-18	
1,2-Dinitrobenzene	ND	0.38	EPA 8270D	9-21-18	9-21-18	
Acenaphthylene	ND	0.015	EPA 8270D/SIM	9-21-18	9-21-18	
3-Nitroaniline	ND	0.38	EPA 8270D	9-21-18	9-21-18	



Date of Report: September 24, 2018  
 Samples Submitted: September 21, 2018  
 Laboratory Reference: 1809-229  
 Project: Foreman

**SEMIVOLATILE ORGANICS EPA 8270D/SIM**  
 page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>FC1-1</b>					
Laboratory ID:	09-229-01					
2,4-Dinitrophenol	ND	1.9	EPA 8270D	9-21-18	9-21-18	
Acenaphthene	ND	0.015	EPA 8270D/SIM	9-21-18	9-21-18	
4-Nitrophenol	ND	0.38	EPA 8270D	9-21-18	9-21-18	
2,4-Dinitrotoluene	ND	0.38	EPA 8270D	9-21-18	9-21-18	
Dibenzofuran	ND	0.38	EPA 8270D	9-21-18	9-21-18	
2,3,5,6-Tetrachlorophenol	ND	0.38	EPA 8270D	9-21-18	9-21-18	
2,3,4,6-Tetrachlorophenol	ND	0.38	EPA 8270D	9-21-18	9-21-18	
Diethylphthalate	ND	1.9	EPA 8270D	9-21-18	9-21-18	
4-Chlorophenyl-phenylether	ND	0.38	EPA 8270D	9-21-18	9-21-18	
4-Nitroaniline	ND	0.38	EPA 8270D	9-21-18	9-21-18	
Fluorene	ND	0.015	EPA 8270D/SIM	9-21-18	9-21-18	
4,6-Dinitro-2-methylphenol	ND	1.9	EPA 8270D	9-21-18	9-21-18	
n-Nitrosodiphenylamine	ND	0.38	EPA 8270D	9-21-18	9-21-18	
1,2-Diphenylhydrazine	ND	0.38	EPA 8270D	9-21-18	9-21-18	
4-Bromophenyl-phenylether	ND	0.38	EPA 8270D	9-21-18	9-21-18	
Hexachlorobenzene	ND	0.38	EPA 8270D	9-21-18	9-21-18	
Pentachlorophenol	ND	1.9	EPA 8270D	9-21-18	9-21-18	
Phenanthrene	ND	0.015	EPA 8270D/SIM	9-21-18	9-21-18	
Anthracene	ND	0.015	EPA 8270D/SIM	9-21-18	9-21-18	
Carbazole	ND	0.38	EPA 8270D	9-21-18	9-21-18	
Di-n-butylphthalate	ND	1.9	EPA 8270D	9-21-18	9-21-18	
Fluoranthene	ND	0.015	EPA 8270D/SIM	9-21-18	9-21-18	
Benzidine	ND	3.8	EPA 8270D	9-21-18	9-21-18	
Pyrene	ND	0.015	EPA 8270D/SIM	9-21-18	9-21-18	
Butylbenzylphthalate	ND	1.9	EPA 8270D	9-21-18	9-21-18	
bis-2-Ethylhexyladipate	ND	1.9	EPA 8270D	9-21-18	9-21-18	
3,3'-Dichlorobenzidine	ND	1.9	EPA 8270D	9-21-18	9-21-18	
Benzo[a]anthracene	ND	0.015	EPA 8270D/SIM	9-21-18	9-21-18	
Chrysene	ND	0.015	EPA 8270D/SIM	9-21-18	9-21-18	
bis(2-Ethylhexyl)phthalate	ND	1.9	EPA 8270D	9-21-18	9-21-18	
Di-n-octylphthalate	ND	1.9	EPA 8270D	9-21-18	9-21-18	
Benzo[b]fluoranthene	ND	0.015	EPA 8270D/SIM	9-21-18	9-21-18	
Benzo(j,k)fluoranthene	ND	0.015	EPA 8270D/SIM	9-21-18	9-21-18	
Benzo[a]pyrene	ND	0.015	EPA 8270D/SIM	9-21-18	9-21-18	
Indeno[1,2,3-cd]pyrene	ND	0.015	EPA 8270D/SIM	9-21-18	9-21-18	
Dibenz[a,h]anthracene	ND	0.015	EPA 8270D/SIM	9-21-18	9-21-18	
Benzo[g,h,i]perylene	ND	0.015	EPA 8270D/SIM	9-21-18	9-21-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	38	19 - 103				
Phenol-d6	48	30 - 103				
Nitrobenzene-d5	44	27 - 105				
2-Fluorobiphenyl	56	36 - 102				
2,4,6-Tribromophenol	55	33 - 110				
Terphenyl-d14	64	38 - 108				



Date of Report: September 24, 2018  
 Samples Submitted: September 21, 2018  
 Laboratory Reference: 1809-229  
 Project: Foreman

**SEMIVOLATILE ORGANICS 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>FC1-2</b>					
<b>Laboratory ID:</b>	<b>09-229-02</b>					
n-Nitrosodimethylamine	ND	0.036	EPA 8270D	9-21-18	9-21-18	
Pyridine	ND	0.36	EPA 8270D	9-21-18	9-21-18	
Phenol	ND	0.036	EPA 8270D	9-21-18	9-21-18	
Aniline	ND	0.18	EPA 8270D	9-21-18	9-21-18	
bis(2-Chloroethyl)ether	ND	0.036	EPA 8270D	9-21-18	9-21-18	
2-Chlorophenol	ND	0.036	EPA 8270D	9-21-18	9-21-18	
1,3-Dichlorobenzene	ND	0.036	EPA 8270D	9-21-18	9-21-18	
1,4-Dichlorobenzene	ND	0.036	EPA 8270D	9-21-18	9-21-18	
Benzyl alcohol	ND	0.18	EPA 8270D	9-21-18	9-21-18	
1,2-Dichlorobenzene	ND	0.036	EPA 8270D	9-21-18	9-21-18	
2-Methylphenol (o-Cresol)	ND	0.036	EPA 8270D	9-21-18	9-21-18	
bis(2-Chloroisopropyl)ether	ND	0.036	EPA 8270D	9-21-18	9-21-18	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.036	EPA 8270D	9-21-18	9-21-18	
n-Nitroso-di-n-propylamine	ND	0.036	EPA 8270D	9-21-18	9-21-18	
Hexachloroethane	ND	0.036	EPA 8270D	9-21-18	9-21-18	
Nitrobenzene	ND	0.036	EPA 8270D	9-21-18	9-21-18	
Isophorone	ND	0.036	EPA 8270D	9-21-18	9-21-18	
2-Nitrophenol	ND	0.036	EPA 8270D	9-21-18	9-21-18	
2,4-Dimethylphenol	ND	0.036	EPA 8270D	9-21-18	9-21-18	
bis(2-Chloroethoxy)methane	ND	0.036	EPA 8270D	9-21-18	9-21-18	
2,4-Dichlorophenol	ND	0.036	EPA 8270D	9-21-18	9-21-18	
1,2,4-Trichlorobenzene	ND	0.036	EPA 8270D	9-21-18	9-21-18	
Naphthalene	ND	0.0072	EPA 8270D/SIM	9-21-18	9-21-18	
4-Chloroaniline	ND	0.18	EPA 8270D	9-21-18	9-21-18	
Hexachlorobutadiene	ND	0.036	EPA 8270D	9-21-18	9-21-18	
4-Chloro-3-methylphenol	ND	0.036	EPA 8270D	9-21-18	9-21-18	
2-Methylnaphthalene	ND	0.0072	EPA 8270D/SIM	9-21-18	9-21-18	
1-Methylnaphthalene	ND	0.0072	EPA 8270D/SIM	9-21-18	9-21-18	
Hexachlorocyclopentadiene	ND	0.036	EPA 8270D	9-21-18	9-21-18	
2,4,6-Trichlorophenol	ND	0.036	EPA 8270D	9-21-18	9-21-18	
2,3-Dichloroaniline	ND	0.036	EPA 8270D	9-21-18	9-21-18	
2,4,5-Trichlorophenol	ND	0.036	EPA 8270D	9-21-18	9-21-18	
2-Chloronaphthalene	ND	0.036	EPA 8270D	9-21-18	9-21-18	
2-Nitroaniline	ND	0.036	EPA 8270D	9-21-18	9-21-18	
1,4-Dinitrobenzene	ND	0.036	EPA 8270D	9-21-18	9-21-18	
Dimethylphthalate	ND	0.036	EPA 8270D	9-21-18	9-21-18	
1,3-Dinitrobenzene	ND	0.036	EPA 8270D	9-21-18	9-21-18	
2,6-Dinitrotoluene	ND	0.036	EPA 8270D	9-21-18	9-21-18	
1,2-Dinitrobenzene	ND	0.036	EPA 8270D	9-21-18	9-21-18	
Acenaphthylene	ND	0.0072	EPA 8270D/SIM	9-21-18	9-21-18	
3-Nitroaniline	ND	0.036	EPA 8270D	9-21-18	9-21-18	



Date of Report: September 24, 2018  
 Samples Submitted: September 21, 2018  
 Laboratory Reference: 1809-229  
 Project: Foreman

**SEMIVOLATILE ORGANICS EPA 8270D/SIM**  
 page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>FC1-2</b>					
Laboratory ID:	09-229-02					
2,4-Dinitrophenol	ND	0.18	EPA 8270D	9-21-18	9-21-18	
Acenaphthene	ND	0.0072	EPA 8270D/SIM	9-21-18	9-21-18	
4-Nitrophenol	ND	0.036	EPA 8270D	9-21-18	9-21-18	
2,4-Dinitrotoluene	ND	0.036	EPA 8270D	9-21-18	9-21-18	
Dibenzofuran	ND	0.036	EPA 8270D	9-21-18	9-21-18	
2,3,5,6-Tetrachlorophenol	ND	0.036	EPA 8270D	9-21-18	9-21-18	
2,3,4,6-Tetrachlorophenol	ND	0.036	EPA 8270D	9-21-18	9-21-18	
Diethylphthalate	ND	0.18	EPA 8270D	9-21-18	9-21-18	
4-Chlorophenyl-phenylether	ND	0.036	EPA 8270D	9-21-18	9-21-18	
4-Nitroaniline	ND	0.036	EPA 8270D	9-21-18	9-21-18	
Fluorene	ND	0.0072	EPA 8270D/SIM	9-21-18	9-21-18	
4,6-Dinitro-2-methylphenol	ND	0.18	EPA 8270D	9-21-18	9-21-18	
n-Nitrosodiphenylamine	ND	0.036	EPA 8270D	9-21-18	9-21-18	
1,2-Diphenylhydrazine	ND	0.036	EPA 8270D	9-21-18	9-21-18	
4-Bromophenyl-phenylether	ND	0.036	EPA 8270D	9-21-18	9-21-18	
Hexachlorobenzene	ND	0.036	EPA 8270D	9-21-18	9-21-18	
Pentachlorophenol	ND	0.18	EPA 8270D	9-21-18	9-21-18	
Phenanthrene	ND	0.0072	EPA 8270D/SIM	9-21-18	9-21-18	
Anthracene	ND	0.0072	EPA 8270D/SIM	9-21-18	9-21-18	
Carbazole	ND	0.036	EPA 8270D	9-21-18	9-21-18	
Di-n-butylphthalate	ND	0.18	EPA 8270D	9-21-18	9-21-18	
Fluoranthene	ND	0.0072	EPA 8270D/SIM	9-21-18	9-21-18	
Benzidine	ND	0.36	EPA 8270D	9-21-18	9-21-18	
Pyrene	ND	0.0072	EPA 8270D/SIM	9-21-18	9-21-18	
Butylbenzylphthalate	ND	0.18	EPA 8270D	9-21-18	9-21-18	
bis-2-Ethylhexyladipate	ND	0.18	EPA 8270D	9-21-18	9-21-18	
3,3'-Dichlorobenzidine	ND	0.18	EPA 8270D	9-21-18	9-21-18	
Benzo[a]anthracene	ND	0.0072	EPA 8270D/SIM	9-21-18	9-21-18	
Chrysene	ND	0.0072	EPA 8270D/SIM	9-21-18	9-21-18	
bis(2-Ethylhexyl)phthalate	ND	0.18	EPA 8270D	9-21-18	9-21-18	
Di-n-octylphthalate	ND	0.18	EPA 8270D	9-21-18	9-21-18	
Benzo[b]fluoranthene	ND	0.0072	EPA 8270D/SIM	9-21-18	9-21-18	
Benzo(j,k)fluoranthene	ND	0.0072	EPA 8270D/SIM	9-21-18	9-21-18	
Benzo[a]pyrene	ND	0.0072	EPA 8270D/SIM	9-21-18	9-21-18	
Indeno[1,2,3-cd]pyrene	ND	0.0072	EPA 8270D/SIM	9-21-18	9-21-18	
Dibenz[a,h]anthracene	ND	0.0072	EPA 8270D/SIM	9-21-18	9-21-18	
Benzo[g,h,i]perylene	ND	0.0072	EPA 8270D/SIM	9-21-18	9-21-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	41	19 - 103				
Phenol-d6	47	30 - 103				
Nitrobenzene-d5	43	27 - 105				
2-Fluorobiphenyl	52	36 - 102				
2,4,6-Tribromophenol	78	33 - 110				
Terphenyl-d14	79	38 - 108				



Date of Report: September 24, 2018  
 Samples Submitted: September 21, 2018  
 Laboratory Reference: 1809-229  
 Project: Foreman

**SEMIVOLATILE ORGANICS 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:	MB0921S1					
n-Nitrosodimethylamine	ND	0.033	EPA 8270D	9-21-18	9-21-18	
Pyridine	ND	0.33	EPA 8270D	9-21-18	9-21-18	
Phenol	ND	0.033	EPA 8270D	9-21-18	9-21-18	
Aniline	ND	0.17	EPA 8270D	9-21-18	9-21-18	
bis(2-Chloroethyl)ether	ND	0.033	EPA 8270D	9-21-18	9-21-18	
2-Chlorophenol	ND	0.033	EPA 8270D	9-21-18	9-21-18	
1,3-Dichlorobenzene	ND	0.033	EPA 8270D	9-21-18	9-21-18	
1,4-Dichlorobenzene	ND	0.033	EPA 8270D	9-21-18	9-21-18	
Benzyl alcohol	ND	0.17	EPA 8270D	9-21-18	9-21-18	
1,2-Dichlorobenzene	ND	0.033	EPA 8270D	9-21-18	9-21-18	
2-Methylphenol (o-Cresol)	ND	0.033	EPA 8270D	9-21-18	9-21-18	
bis(2-Chloroisopropyl)ether	ND	0.033	EPA 8270D	9-21-18	9-21-18	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.033	EPA 8270D	9-21-18	9-21-18	
n-Nitroso-di-n-propylamine	ND	0.033	EPA 8270D	9-21-18	9-21-18	
Hexachloroethane	ND	0.033	EPA 8270D	9-21-18	9-21-18	
Nitrobenzene	ND	0.033	EPA 8270D	9-21-18	9-21-18	
Isophorone	ND	0.033	EPA 8270D	9-21-18	9-21-18	
2-Nitrophenol	ND	0.033	EPA 8270D	9-21-18	9-21-18	
2,4-Dimethylphenol	ND	0.033	EPA 8270D	9-21-18	9-21-18	
bis(2-Chloroethoxy)methane	ND	0.033	EPA 8270D	9-21-18	9-21-18	
2,4-Dichlorophenol	ND	0.033	EPA 8270D	9-21-18	9-21-18	
1,2,4-Trichlorobenzene	ND	0.033	EPA 8270D	9-21-18	9-21-18	
Naphthalene	ND	0.0067	EPA 8270D/SIM	9-21-18	9-21-18	
4-Chloroaniline	ND	0.17	EPA 8270D	9-21-18	9-21-18	
Hexachlorobutadiene	ND	0.033	EPA 8270D	9-21-18	9-21-18	
4-Chloro-3-methylphenol	ND	0.033	EPA 8270D	9-21-18	9-21-18	
2-Methylnaphthalene	ND	0.0067	EPA 8270D/SIM	9-21-18	9-21-18	
1-Methylnaphthalene	ND	0.0067	EPA 8270D/SIM	9-21-18	9-21-18	
Hexachlorocyclopentadiene	ND	0.033	EPA 8270D	9-21-18	9-21-18	
2,4,6-Trichlorophenol	ND	0.033	EPA 8270D	9-21-18	9-21-18	
2,3-Dichloroaniline	ND	0.033	EPA 8270D	9-21-18	9-21-18	
2,4,5-Trichlorophenol	ND	0.033	EPA 8270D	9-21-18	9-21-18	
2-Chloronaphthalene	ND	0.033	EPA 8270D	9-21-18	9-21-18	
2-Nitroaniline	ND	0.033	EPA 8270D	9-21-18	9-21-18	
1,4-Dinitrobenzene	ND	0.033	EPA 8270D	9-21-18	9-21-18	
Dimethylphthalate	ND	0.033	EPA 8270D	9-21-18	9-21-18	
1,3-Dinitrobenzene	ND	0.033	EPA 8270D	9-21-18	9-21-18	
2,6-Dinitrotoluene	ND	0.033	EPA 8270D	9-21-18	9-21-18	
1,2-Dinitrobenzene	ND	0.033	EPA 8270D	9-21-18	9-21-18	
Acenaphthylene	ND	0.0067	EPA 8270D/SIM	9-21-18	9-21-18	
3-Nitroaniline	ND	0.033	EPA 8270D	9-21-18	9-21-18	



Date of Report: September 24, 2018  
 Samples Submitted: September 21, 2018  
 Laboratory Reference: 1809-229  
 Project: Foreman

**SEMIVOLATILE ORGANICS EPA 8270D/SIM  
 METHOD BLANK QUALITY CONTROL**

page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0921S1					
2,4-Dinitrophenol	ND	0.17	EPA 8270D	9-21-18	9-21-18	
Acenaphthene	ND	0.0067	EPA 8270D/SIM	9-21-18	9-21-18	
4-Nitrophenol	ND	0.033	EPA 8270D	9-21-18	9-21-18	
2,4-Dinitrotoluene	ND	0.033	EPA 8270D	9-21-18	9-21-18	
Dibenzofuran	ND	0.033	EPA 8270D	9-21-18	9-21-18	
2,3,5,6-Tetrachlorophenol	ND	0.033	EPA 8270D	9-21-18	9-21-18	
2,3,4,6-Tetrachlorophenol	ND	0.033	EPA 8270D	9-21-18	9-21-18	
Diethylphthalate	ND	0.17	EPA 8270D	9-21-18	9-21-18	
4-Chlorophenyl-phenylether	ND	0.033	EPA 8270D	9-21-18	9-21-18	
4-Nitroaniline	ND	0.033	EPA 8270D	9-21-18	9-21-18	
Fluorene	ND	0.0067	EPA 8270D/SIM	9-21-18	9-21-18	
4,6-Dinitro-2-methylphenol	ND	0.17	EPA 8270D	9-21-18	9-21-18	
n-Nitrosodiphenylamine	ND	0.033	EPA 8270D	9-21-18	9-21-18	
1,2-Diphenylhydrazine	ND	0.033	EPA 8270D	9-21-18	9-21-18	
4-Bromophenyl-phenylether	ND	0.033	EPA 8270D	9-21-18	9-21-18	
Hexachlorobenzene	ND	0.033	EPA 8270D	9-21-18	9-21-18	
Pentachlorophenol	ND	0.17	EPA 8270D	9-21-18	9-21-18	
Phenanthrene	ND	0.0067	EPA 8270D/SIM	9-21-18	9-21-18	
Anthracene	ND	0.0067	EPA 8270D/SIM	9-21-18	9-21-18	
Carbazole	ND	0.033	EPA 8270D	9-21-18	9-21-18	
Di-n-butylphthalate	ND	0.17	EPA 8270D	9-21-18	9-21-18	
Fluoranthene	ND	0.0067	EPA 8270D/SIM	9-21-18	9-21-18	
Benzidine	ND	0.33	EPA 8270D	9-21-18	9-21-18	
Pyrene	ND	0.0067	EPA 8270D/SIM	9-21-18	9-21-18	
Butylbenzylphthalate	ND	0.17	EPA 8270D	9-21-18	9-21-18	
bis-2-Ethylhexyladipate	ND	0.17	EPA 8270D	9-21-18	9-21-18	
3,3'-Dichlorobenzidine	ND	0.17	EPA 8270D	9-21-18	9-21-18	
Benzo[a]anthracene	ND	0.0067	EPA 8270D/SIM	9-21-18	9-21-18	
Chrysene	ND	0.0067	EPA 8270D/SIM	9-21-18	9-21-18	
bis(2-Ethylhexyl)phthalate	ND	0.17	EPA 8270D	9-21-18	9-21-18	
Di-n-octylphthalate	ND	0.17	EPA 8270D	9-21-18	9-21-18	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270D/SIM	9-21-18	9-21-18	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270D/SIM	9-21-18	9-21-18	
Benzo[a]pyrene	ND	0.0067	EPA 8270D/SIM	9-21-18	9-21-18	
Indeno[1,2,3-cd]pyrene	ND	0.0067	EPA 8270D/SIM	9-21-18	9-21-18	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270D/SIM	9-21-18	9-21-18	
Benzo[g,h,i]perylene	ND	0.0067	EPA 8270D/SIM	9-21-18	9-21-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	42	19 - 103				
Phenol-d6	47	30 - 103				
Nitrobenzene-d5	43	27 - 105				
2-Fluorobiphenyl	50	36 - 102				
2,4,6-Tribromophenol	70	33 - 110				
Terphenyl-d14	69	38 - 108				



Date of Report: September 24, 2018  
 Samples Submitted: September 21, 2018  
 Laboratory Reference: 1809-229  
 Project: Foreman

**SEMIVOLATILE ORGANICS EPA 8270D/SIM  
 SB/SBD QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result		Spike Level		Percent Recovery		Recovery Limits	RPD	RPD Limit	Flags
	SB	SBD	SB	SBD	SB	SBD				
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB0921S1									
Phenol	<b>0.782</b>	<b>0.896</b>	1.33	1.33	59	67	45 - 94	14	29	
2-Chlorophenol	<b>0.770</b>	<b>0.879</b>	1.33	1.33	58	66	46 - 94	13	33	
1,4-Dichlorobenzene	<b>0.360</b>	<b>0.429</b>	0.667	0.667	54	64	42 - 91	17	37	
n-Nitroso-di-n-propylamine	<b>0.382</b>	<b>0.479</b>	0.667	0.667	57	72	45 - 100	23	26	
1,2,4-Trichlorobenzene	<b>0.396</b>	<b>0.454</b>	0.667	0.667	59	68	45 - 100	14	32	
4-Chloro-3-methylphenol	<b>0.950</b>	<b>1.01</b>	1.33	1.33	71	76	55 - 97	6	21	
Acenaphthene	<b>0.444</b>	<b>0.497</b>	0.667	0.667	67	75	48 - 91	11	21	
4-Nitrophenol	<b>1.04</b>	<b>1.11</b>	1.33	1.33	78	83	53 - 102	7	20	
2,4-Dinitrotoluene	<b>0.496</b>	<b>0.524</b>	0.667	0.667	74	79	47 - 96	5	19	
Pentachlorophenol	<b>1.32</b>	<b>1.35</b>	1.33	1.33	99	102	35 - 125	2	26	
Pyrene	<b>0.540</b>	<b>0.545</b>	0.667	0.667	81	82	55 - 110	1	17	
<i>Surrogate:</i>										
2-Fluorophenol					48	55	19 - 103			
Phenol-d6					55	61	30 - 103			
Nitrobenzene-d5					51	57	27 - 105			
2-Fluorobiphenyl					56	64	36 - 102			
2,4,6-Tribromophenol					76	74	33 - 110			
Terphenyl-d14					71	71	38 - 108			



Date of Report: September 24, 2018  
 Samples Submitted: September 21, 2018  
 Laboratory Reference: 1809-229  
 Project: Foreman

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

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>FC1-1</b>					
Laboratory ID:	09-229-01					
Arsenic	<b>ND</b>	11	EPA 6010D	9-21-18	9-21-18	
Cadmium	<b>ND</b>	0.57	EPA 6010D	9-21-18	9-21-18	
Chromium	<b>64</b>	0.57	EPA 6010D	9-21-18	9-21-18	
Lead	<b>27</b>	5.7	EPA 6010D	9-21-18	9-21-18	
Mercury	<b>ND</b>	0.28	EPA 7471B	9-21-18	9-21-18	

<b>Client ID:</b>	<b>FC1-2</b>					
Laboratory ID:	09-229-02					
Arsenic	<b>ND</b>	11	EPA 6010D	9-21-18	9-21-18	
Cadmium	<b>ND</b>	0.54	EPA 6010D	9-21-18	9-21-18	
Chromium	<b>11</b>	0.54	EPA 6010D	9-21-18	9-21-18	
Lead	<b>ND</b>	5.4	EPA 6010D	9-21-18	9-21-18	
Mercury	<b>ND</b>	0.27	EPA 7471B	9-21-18	9-21-18	



Date of Report: September 24, 2018  
 Samples Submitted: September 21, 2018  
 Laboratory Reference: 1809-229  
 Project: Foreman

**TOTAL METALS  
 EPA 6010D/7471B  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB0921SM2					
Arsenic	ND	10	EPA 6010D	9-21-18	9-21-18	
Cadmium	ND	0.50	EPA 6010D	9-21-18	9-21-18	
Chromium	ND	0.50	EPA 6010D	9-21-18	9-21-18	
Lead	ND	5.0	EPA 6010D	9-21-18	9-21-18	

Laboratory ID:	MB0921S1					
Mercury	ND	0.25	EPA 7471B	9-21-18	9-21-18	

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>								
Laboratory ID:	09-138-01							
	ORIG	DUP						
Arsenic	14.6	14.8	NA	NA	NA	NA	2	20
Cadmium	ND	ND	NA	NA	NA	NA	NA	20
Chromium	43.8	50.1	NA	NA	NA	NA	13	20
Lead	52.9	55.3	NA	NA	NA	NA	5	20

Laboratory ID:	09-213-01							
Mercury	ND	ND	NA	NA	NA	NA	NA	20

**MATRIX SPIKES**

Laboratory ID:	09-138-01									
	MS	MSD	MS	MSD	MS	MSD				
Arsenic	106	105	100	100	14.6	92	91	75-125	1	20
Cadmium	48.0	47.1	50.0	50.0	ND	96	94	75-125	2	20
Chromium	145	147	100	100	43.8	101	103	75-125	1	20
Lead	296	287	250	250	52.9	97	94	75-125	3	20

Laboratory ID:	09-213-01									
Mercury	0.566	0.586	0.500	0.500	0.0548	102	106	80-120	3	20



Date of Report: September 24, 2018  
Samples Submitted: September 21, 2018  
Laboratory Reference: 1809-229  
Project: Foreman

### % MOISTURE

Date Analyzed: 9-21-18

Client ID	Lab ID	% Moisture
FC1-1	09-229-01	12
FC1-2	09-229-02	8





### 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





**Mn Onsite Environmental Inc.**  
 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) 24 hours

Same Day  1 Day

2 Days  3 Days

Standard (7 Days)  
 (TPH analysis 5 Days)

(other)

Laboratory Number: **09-229**

Company: WSDOT  
 Project Number:  
 Project Name: Foreman  
 Project Manager: PAT SUBORDA  
 Sampled by: WSDOT

Lab ID	Sample Identification	Date Sampled	Time Sampled	Matrix	Number of Containers
1	FCI-1	9/20	14:15	S	6
2	FCI-2	9/21	11:00	S	6

Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX	NWTPH-Gx	NWTPH-Dx	Volatiles 8260B	Halogenated Volatiles 8260B	Semivolatiles 8270D/SIM (with low-level PAHs)	PAHs 8270D/SIM (low-level)	PCBs 8082	Organochlorine Pesticides 8081A	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA (MTCA) Metals (circle one)	TCLP Metals	HEM (oil and grease) 1664	% Moisture
6				✓	✓		✓						✓			X
6				✓												X

Signature	Company	Date	Time	Comments/Special Instructions
<u>[Signature]</u>	<u>WSDOT</u>	<u>9/21/18</u>	<u>11:10</u>	<u>Call when you receive these</u>
<u>[Signature]</u>	<u>AC PHH</u>	<u>9/21/18</u>	<u>11:10</u>	
<u>[Signature]</u>	<u>AC PHH</u>	<u>9/21/18</u>	<u>12:39</u>	
<u>[Signature]</u>	<u>ORE</u>	<u>9/21/18</u>	<u>12:39</u>	

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

Reviewed/Date \_\_\_\_\_

Chromatograms with final report

**APPENDIX 4 DISPOSAL TICKETS**



19 + cel

Load

401187

Phone (206) 621.9777

Certified Weight Ticket

Company \_\_\_\_\_

Charge \_\_\_\_\_ Cash

Truck # \_\_\_\_\_

Account # \_\_\_\_\_

Trailer # \_\_\_\_\_

Card # \_\_\_\_\_

Weighmaster [Signature]

Signature [Signature]

16:53:47

20 Sep 2018

Weigh fee: 10.00

3215 4TH AVE S

Axle 1:	3240 lb
Axle 2:	5620 lb
Axle 3:	5660 lb
Total:	14520 lb



2<sup>nd</sup> Load

401252

Phone (206) 621.9777

Certified Weight Ticket

Company  MARINE VACUME SERVICE Charge  Cash

Truck #  \_\_\_\_\_

Account # \_\_\_\_\_

Trailer # \_\_\_\_\_

Card # \_\_\_\_\_

Weighmaster \_\_\_\_\_ *[Signature]*

Signature  R.B. ALLEN \_\_\_\_\_

08:44:47 22 Sep 2018

Weigh fee: 10.00  
3215 4TH AVE S

Axle 1: 10680 lb  
Axle 2: 7540 lb  
Total: 18220 lb

*V I S E*



Light weight

52  
401254

Phone (206) 621.9777

Certified Weight Ticket

Company  MARINE VACUUM SERVICE

Charge \_\_\_\_\_

Cash

Truck #  \_\_\_\_\_

Account # \_\_\_\_\_

Trailer # \_\_\_\_\_

Card # \_\_\_\_\_

Weighmaster \_\_\_\_\_ 

Signature  R B ALLEN

10:13:36

22 Sep 2018

Weigh fee: 10.00

3215 4TH AVE S



Axle 1: 3620 lb  
Axle 2: 7640 lb  
Total: 11260 lb





Light weight

52  
401254

Phone (206) 621.9777

Certified Weight Ticket

Company  MARINE VACUUM SERVICE

Charge \_\_\_\_\_ Cash

Truck #  \_\_\_\_\_

Account # \_\_\_\_\_

Trailer # \_\_\_\_\_

Card # \_\_\_\_\_

Weighmaster \_\_\_\_\_ *[Signature]*

Signature  R B ALLEN

10:13:36 22 Sep 2018

Weigh fee: 10.00  
3215 4TH AVE S

102



*[Handwritten mark]*

Axle 1: 3620 lb  
Axle 2: 7640 lb  
Total: 11260 lb



**APPENDIX 5 PHOTOS**



SS-1 during investigation



SS-2 during investigation



SS-3 during investigation



SS-4 during investigation



SS-4 During investigation



General trash and debris scattered around the property



Grading and soil displacement on property



Soil displacement near SS-1



Soil displacement near SS-2



SS-2 Sample near soil displacement



SS3 of stockpile



SS4 of stockpile



Cleanup of SS-3 Area and confirmation sample area



Part of Cleanup of SS-4 area and confirmation sample taken from this area



Backfill of front area



Backfill of side area