



## **2017 Summary of investigations and remedial actions**

Former Kelly-Moore Manufacturing Facility  
5400–5580 Airport Way South  
Project # 0146970060 Kelly-Moore Paint Company, Inc.

Prepared for:

**Kelly-Moore Paint Company, Inc.**  
301 W Hurst Boulevard, Hurst, TX 76053

6/5/2018



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## **Prepared for:**

Kelly-Moore Paint Company, Inc.  
301 W Hurst Boulevard, Hurst, TX 76053

## **Prepared by:**

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**6/5/2018**

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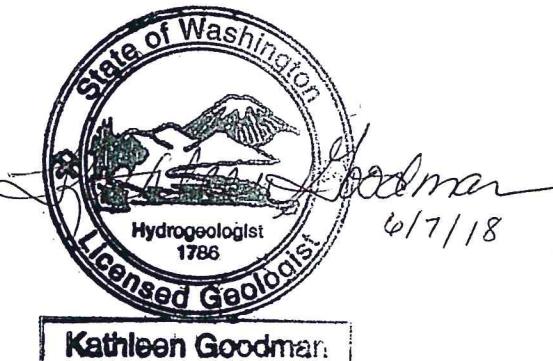
# wood.

**FORMER KELLY-MOORE MANUFACTURING FACILITY**  
2017 Summary of investigations and remedial actions  
5400-5580 Airport Way South  
Seattle, Washington

June 7, 2018  
Project No. 0146970060

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Licensed Geologist/Hydrogeologist #1786  
Expiration Date: September 6, 2018

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- Appendix F: Preliminary Screening Level Tables for Soil, Groundwater and Vapor

## 1.0 Introduction

Wood Environment & Infrastructure Solutions, Inc. (Wood), formerly Amec Foster Wheeler Environment & Infrastructure, Inc., prepared this report on behalf of Kelly-Moore Paint Company, Inc. (Kelly-Moore) for the former Kelly-Moore manufacturing facility located at 5400–5410 Airport Way South, in Seattle, Washington (the site) (Figure 1). This report provides a summary of groundwater samples collected since the Remedial Investigation, Feasibility Study, and Disproportionate Cost Analysis (RI/FS/DCA) was submitted to the Washington State Department of Ecology (Ecology) in March 2017 (Amec Foster Wheeler, 2017), as well as information regarding implementation of the remedial actions discussed in the RI/FS/DCA.

The RI/FS/DCA summarized the results of groundwater and soil sampling conducted since site activities began in 2011. Site activities conducted after March 2017 include dry and wet season groundwater sampling events, which were conducted in September 2017 and January 2018, respectively. Additionally, the soil vapor extraction (SVE) system was installed and started operating. While the air sparge system was also installed, it will not be operational until later in 2018. Details about the site activities conducted in 2017 are discussed below.

## 2.0 Groundwater sampling

Groundwater samples have been collected twice a year (during the wet and dry seasons) since June 2016. Tables 1 through 4 provide information on monitoring well construction, groundwater elevations, field parameters, and groundwater results.

### 2.1 Water level measurements and site hydrogeology

Monitoring well construction details are provided in Table 1. Depth-to-water measurements, top of casing (TOC) elevations, and groundwater elevations measured during the monitoring events are presented in Table 2. Groundwater depths were measured to the nearest 0.01 foot using a depth to water probe. Groundwater levels for each monitoring well were subtracted from the TOC elevations to determine groundwater elevations. TOC elevations were measured by Duane Hartman and Associates, of Seattle, Washington.

Groundwater elevation contours for water level measurements collected in September 2017 and January 2018 are presented on Figures 2 and 3. Water level measurements collected in September 2017 and January 2018 indicate that groundwater generally flows to the west-southwest. Groundwater elevations across the site vary seasonally, with higher groundwater elevations in the wet season and lower elevations in the dry season. The wet season/dry season range of elevations observed in 2017 was approximately 0.5 foot of elevation difference.

### 2.2 Groundwater sampling methodology

Groundwater samples were collected using low flow methodology (EPA, 1988) following the Additional Investigation Work Plan (Amec Foster Wheeler, 2016).

A peristaltic pump and new polyethylene tubing were used to collect groundwater from each well at a flow rate of approximately 100–200 milliliters per minute. Groundwater parameters were measured during purging using a YSI multi-parameter water quality meter and were recorded by hand on field data sheets (Appendix A). Parameters measured were turbidity, temperature, pH, dissolved oxygen, conductivity, and oxidation reduction potential (Table 3). Representative unfiltered groundwater samples were collected upon stabilization of the water quality parameters over the course of three consecutive measurements.

Groundwater sample containers were filled directly from the pump tubing and were immediately placed on ice. Samples were transported under chain-of-custody protocols to Onsite Environmental, Inc., in Redmond, Washington, for laboratory analyses. Each groundwater sample was analyzed for the following:

- Volatile organic compounds (VOCs) by U.S. Environmental Protection Agency (EPA) Method 8260C;
- Polycyclic aromatic hydrocarbons (PAHs) by EPA Method 8270D with selected ion monitoring for some compounds;
- Total metals (arsenic, barium, cadmium, chromium, lead, selenium and silver) by EPA Method 200.8 and total mercury by EPA Method 7470A;
- Total petroleum hydrocarbons (TPH) in the gasoline range (TPH-G) by Ecology method NWTPH Gx; and
- TPH in the diesel and motor oil ranges (TPH-D and TPH-O) by Ecology Method NWTPH-Dx.

Laboratory data packages and data validation memoranda are included in Appendix B.

## **3.0 Groundwater analytical results**

Groundwater results for commonly detected compounds are presented in Table 4, along with the results for detected compounds in sampling events conducted since 2011.

### **3.1 Total petroleum hydrocarbons**

The highest concentrations of TPH in the gasoline range have been observed in the groundwater from KMW-03, KMW-04, KMW-06, KMW-08, KMW-09, and KMW-10. Decreasing trends are observed in the groundwater samples collected from all of the monitoring wells except for from KMW-06, KMW-09, and KMW-10.

During the most recent sampling event conducted in January 2018, the concentrations of TPH in the gasoline range were below the Model Toxics Control Act (MTCA) Method A Cleanup Level of 800 micrograms per liter ( $\mu\text{g}/\text{L}$ ) (where benzene is present) except for KMW-04 and KMW-06.

TPH-D and TPH-O have been detected in the groundwater from all of the monitoring wells at least once since sampling began in 2011, except for KMW-07. During the recent sampling event, the concentrations of TPH-D exceeded the MTCA Method A Cleanup Level of 500  $\mu\text{g}/\text{L}$  in the groundwater collected from KMW-03R, KMW-04, KMW-06, KMW-09, and KMW-10.

### **3.2 Volatile organic compounds**

Groundwater samples were analyzed for the full list of VOC compounds, and benzene, toluene, ethylbenzene and xylenes (BTEX) compounds were the most frequently detected VOCs, predominantly in the central area of the site where high concentrations of TPH-G have been detected. The BTEX detections are most significant in the groundwater from KMW-04 and do not appear to have decreased in concentration over time, unlike the TPH-G concentrations in the groundwater from KMW-04. We expect to see concentrations decrease in groundwater with the startup of the SVE system.

Other VOC compounds detected in the groundwater are 1,2,4-trimethylbenzene and 1,3,5-trimethylbenzene, which were detected in the groundwater collected from KMW-04 and KMW-10.

Chlorinated VOCs were not detected in groundwater samples collected from the monitoring wells during the 2017 and 2018 sampling events, which is consistent with historical results.

### 3.3 Polycyclic aromatic hydrocarbons

During the August 2017 and January 2018 sampling events, PAHs were detected in the groundwater collected from KMW-04, KMW-06, KMW-08, and KMW-09. During the January 2018 sampling event, the concentrations of total PAHs in groundwater ranged from below detection (KMW-02R, KMW-03R, KMW-04, KMW-07, and KMW-10) to 0.095 µg/L detected in the groundwater sample collected from KMW-06.

### 3.4 Metals

Groundwater samples were analyzed for total arsenic, barium, cadmium, chromium, lead, mercury, selenium, and silver. The concentrations of metals in the groundwater samples were generally below detection, except for arsenic. During the January 2018 sampling event, arsenic concentrations detected in the groundwater from monitoring wells KMW-04 and KMW-10 only slightly exceed the Ecology background level of 5.0 µg/L, at concentrations of 7.6 µg/L and 5.7 µg/L, respectively.

## 4.0 Soil vapor extraction/air sparge system installation

Soil vapor extraction (SVE) and air sparging were selected to address past releases of hydrocarbons associated with former paint manufacturing activities at this site. SVE uses a vacuum to extract soil vapors from the subsurface while air sparging volatilizes hydrocarbons in the saturated zone. Both methods introduce oxygen into the subsurface, which also promotes aerobic biodegradation of residual hydrocarbons.

During redevelopment construction of the new warehouse on the eastern part of the site in 2015, a series of seven horizontal SVE wells (SVE-01 through SVE-07) were installed. After building construction was completed, a second set of five horizontal SVE wells (SVE-09 through SVE-13) were installed in the parking lot on the western side of the site. A set of five air sparge wells were installed between the western SVE wells. Figure 4 shows the location of the SVE horizontal wells at the site, and Figure 5 shows the location of the air sparge wells. Appendix C contains the SVE well construction logs and the air sparge boring logs. The air sparge wells were permitted as class 5 Underground Injection Control (UIC) wells and a copy of the UIC permit is included in Appendix D.

The SVE wells installed under the building were routed to a common manifold (referred to as the eastern manifold) located in a walkway between the north warehouse and the south warehouse. The SVE wells installed on the west side of the building were routed to the western manifold, which is located in a fenced-off area near the treatment equipment. Figure 6 shows the layout of the SVE and air sparge system based on the current equipment configuration.

The SVE blower and air sparge compressor were installed adjacent to the western manifold along with a catalytic thermal oxidizer (CATOX) unit. The CATOX is used to treat the extracted soil vapor, as well as volatized hydrocarbon sparged from the shallow groundwater that are recovered by the western SVE wells. The treatment system was permitted with the Puget Sound Clean Air Agency. Their approval letter and a copy of the Notice of Completion are contained in Appendix E.

Figure 7 is a process and instrumentation diagram showing the SVE and air sparge systems and the treatment equipment. This equipment was installed over a period of several months from July through October 2017, with additional time needed to replace polyvinyl chloride pipes used to deliver air to the air sparge wells. Polyvinyl chloride piping is not suitable for use in aboveground piping. A cellular auto-dialer also was installed that will notify Wood personnel if the CATOX system shuts down due to high water in the knockout pot, or if the CATOX unit stops operating. The SVE system started operation on October 31, 2017. After two months of operation we installed a 250-gallon polyethylene tote to provide additional

storage of knockout water produced by the SVE wells. The high water shut-off switch formerly installed on the knockout pot was reinstalled in the tote.

After startup, a set of soil vapor samples were collected from the SVE wells, and as well as the influent and effluent to the CATOX system. Table 5 presents the analytical data for these initial SVE samples. In addition to the analytical data from the laboratory, a flame ionization detector (FID) was used to measure the concentration of hydrocarbons in the SVE wells and the influent and effluent from the CATOX unit. Table 6 lists the FID readings collected through March 2018. The calculated removal efficiencies based on the influent and effluent FID readings are presented in Table 7. Based on the FID readings, the SVE removal efficiencies remain in compliance with the PSCAA permit requirements.

As Table 6 shows, the SVE wells vary greatly in hydrocarbon content (Table 6). We are continuing to optimize the flow from the wells to maximize the concentration of hydrocarbons in the SVE influent. Once the SVE concentrations have decreased, the plan is to turn on the air sparge wells system, which will increase the hydrocarbon concentrations recovered by the SVE system. Eventually we may cycle the air sparge system to increase mixing in the subsurface.

## 5.0 Conclusions and recommendations

The following actions will be conducted before the end of 2018:

- Groundwater samples will be collected for the dry season sampling event in August 2018.
- Indoor air samples will be collected in order to assess baseline conditions prior to starting up the air sparge system. Soil vapor concentrations will be evaluated against results from baseline samples collected for the SVE system.
- Kelly Moore and Ecology would like to continue working together to take the necessary steps to eventually obtain "No Further Action" for the site. To that end, cleanup levels have been developed following the suggestions presented in our meeting on May 8, 2018. The cleanup levels are presented in Appendix F.

## 6.0 References

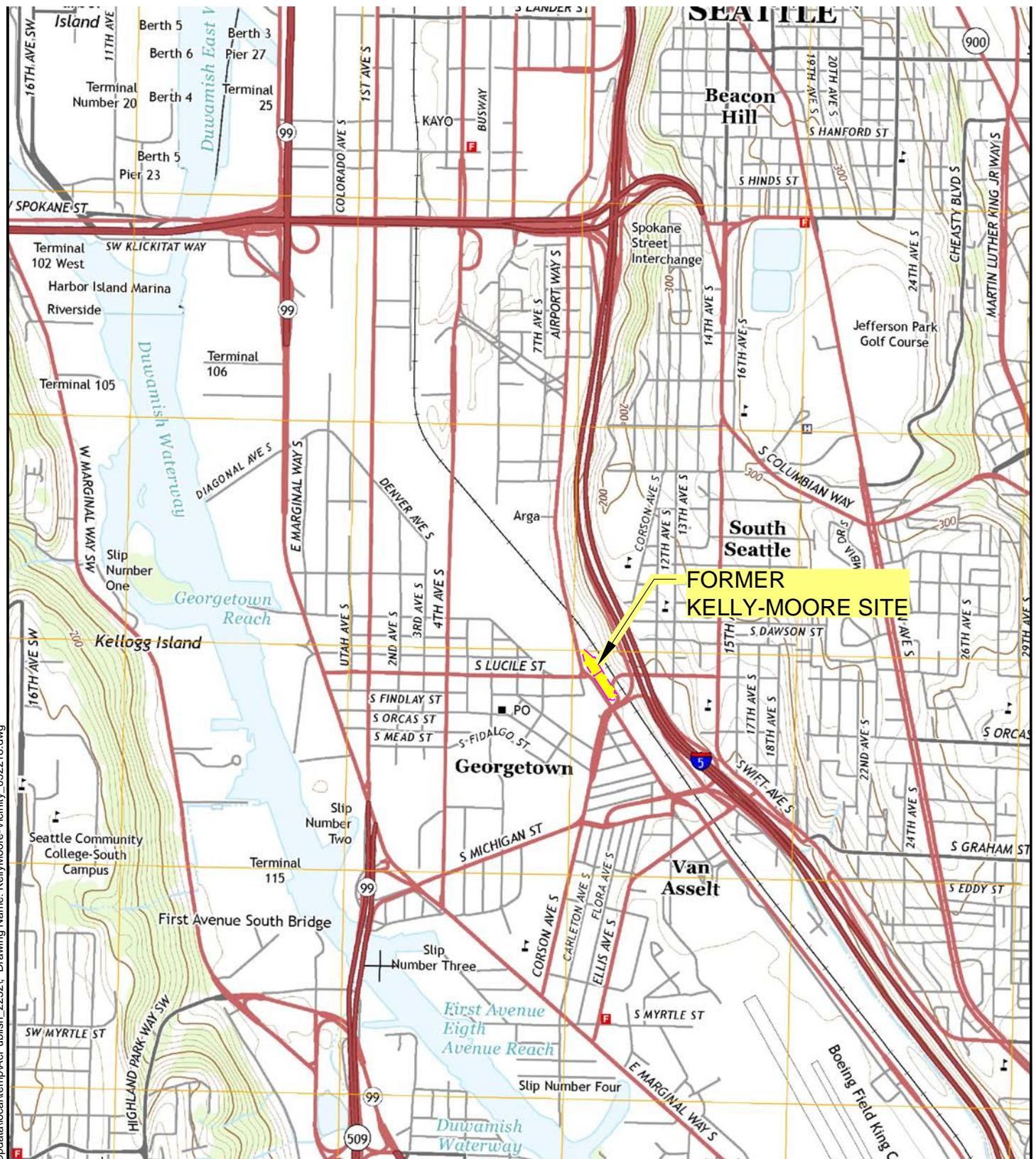
Amec Foster Wheeler Environment & Infrastructure, Inc. (Amec Foster Wheeler), 2017, Revised Remedial Investigation, Feasibility Study, and Disproportionate Cost Analysis, Former Kelly-Moore Manufacturing Facility, 5400–5580 Airport Way South, Seattle, Washington, March.

Amec Foster Wheeler, 2016, Additional Investigation Work Plan, Former Kelly-Moore Manufacturing Facility, 5400–5580 Airport Way South, Seattle, Washington, June.

U.S. Environmental Protection Agency (EPA), 1998, Groundwater Sampling Procedure Low Stress (Low Flow) Purging and Sampling, GW Sampling SOP, March 16.

**wood.**

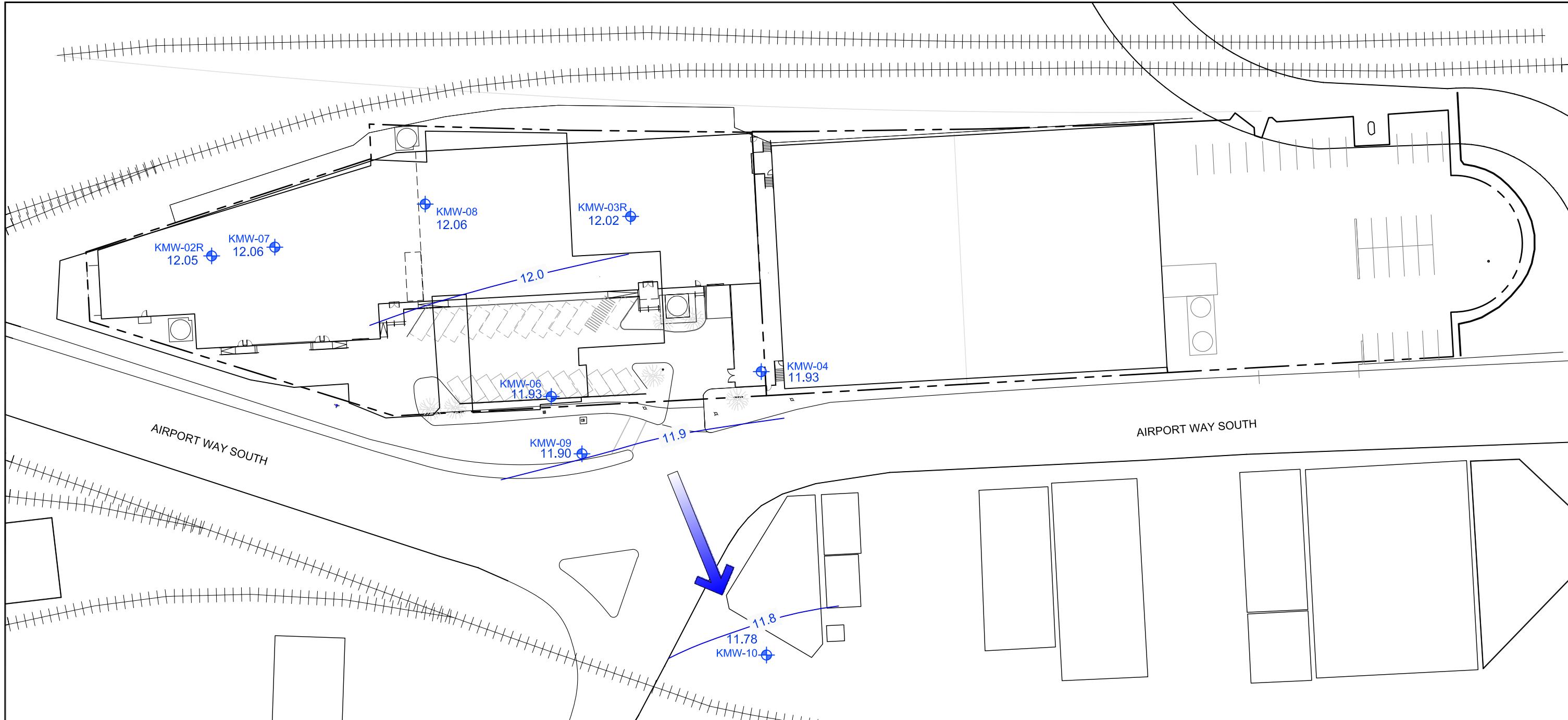
## **Figures**



**SITE LOCATION**  
Kelly-Moore Paint Company  
Seattle, Washington

By: APS	Date: 05/22/18	Project No.	14697
Wood Environment & Infrastructure Solutions, Inc.	Figure	1	

Plot Date: 05/22/18 - 4:57pm. Plotted by: adam.stenberg  
Drawing Path: C:\Users\adam.stenberg\appdata\local\temp\ACPublish\_2232\ Drawing Name: KellyMoore-Vicinity\_052218.dwg



#### EXPLANATION

DRAWN BY: APS CHECKED BY: B

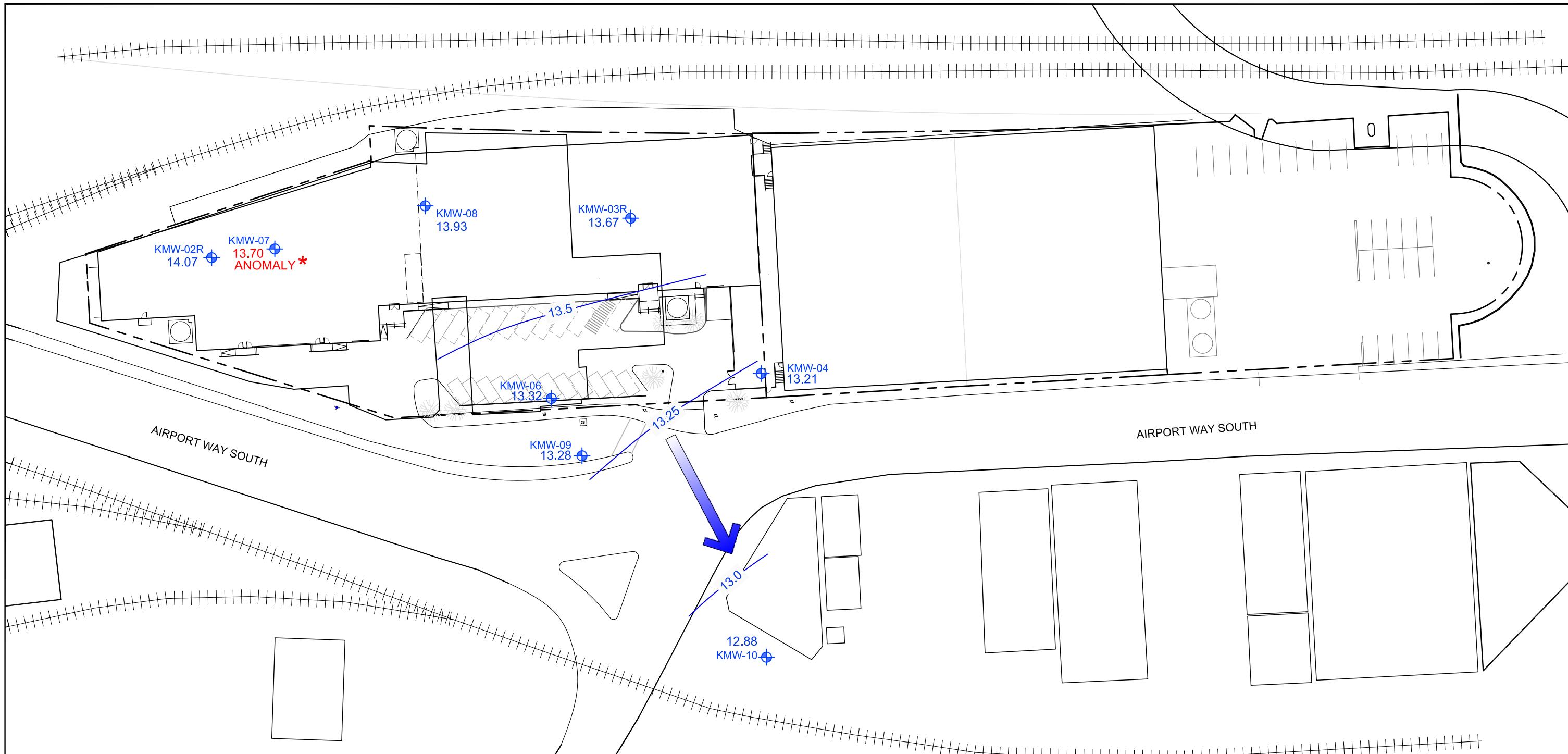
- PROPERTY LINE
- KMW-04 (11.93) MONITORING WELL LOCATION AND GROUNDWATER ELEVATION IN FEET BELOW MEAN SEA LEVEL (NAVD88-FT)
- 11.9 GROUNDWATER ELEVATION CONTOUR
- GROUNDWATER FLOW DIRECTION
- ||||| RAILROAD TRACK

0 30 60  
APPROXIMATE SCALE IN FEET

AUGUST 2017 DRY SEASON  
GROUNDWATER CONTOURS  
Kelly-Moore Paint Company  
Seattle, Washington

**wood.**

By: APS	Date: 05/21/18	Project No. 14697
Wood Environment & Infrastructure Solutions, Inc.	Figure	2



#### EXPLANATION

DRAWN BY: APS CHECKED BY:

KMW-04 13.21  
MONITORING WELL LOCATION AND GROUNDWATER ELEVATION IN FEET BELOW MEAN SEA LEVEL (NAVD88-FT)

\* ANOMALOUS WATER LEVEL NOT USED FOR CONTOURING

13.5 GROUNDWATER ELEVATION CONTOUR

GROUNDWATER FLOW DIRECTION

RAILROAD TRACK

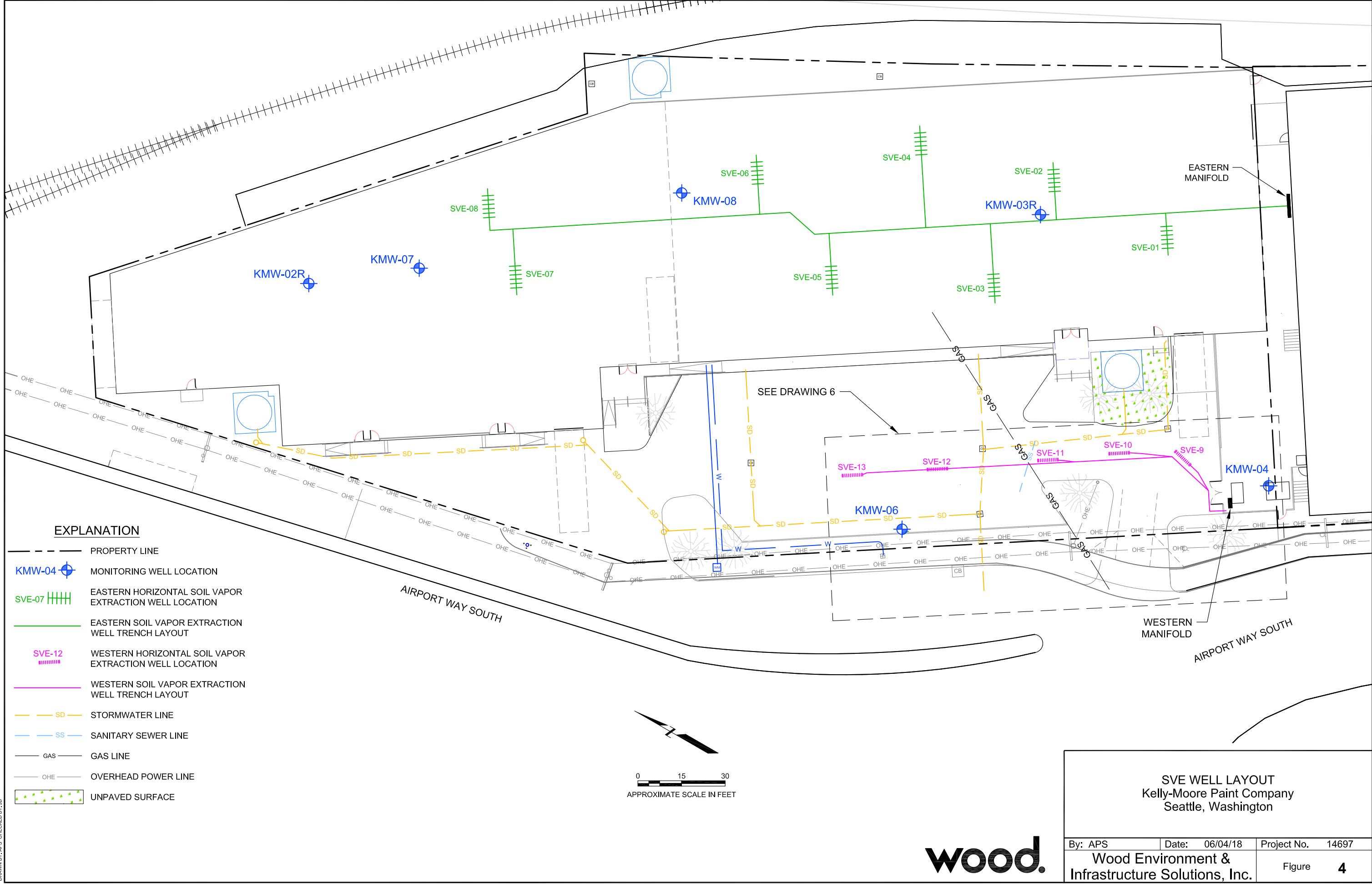
0 30 60  
APPROXIMATE SCALE IN FEET

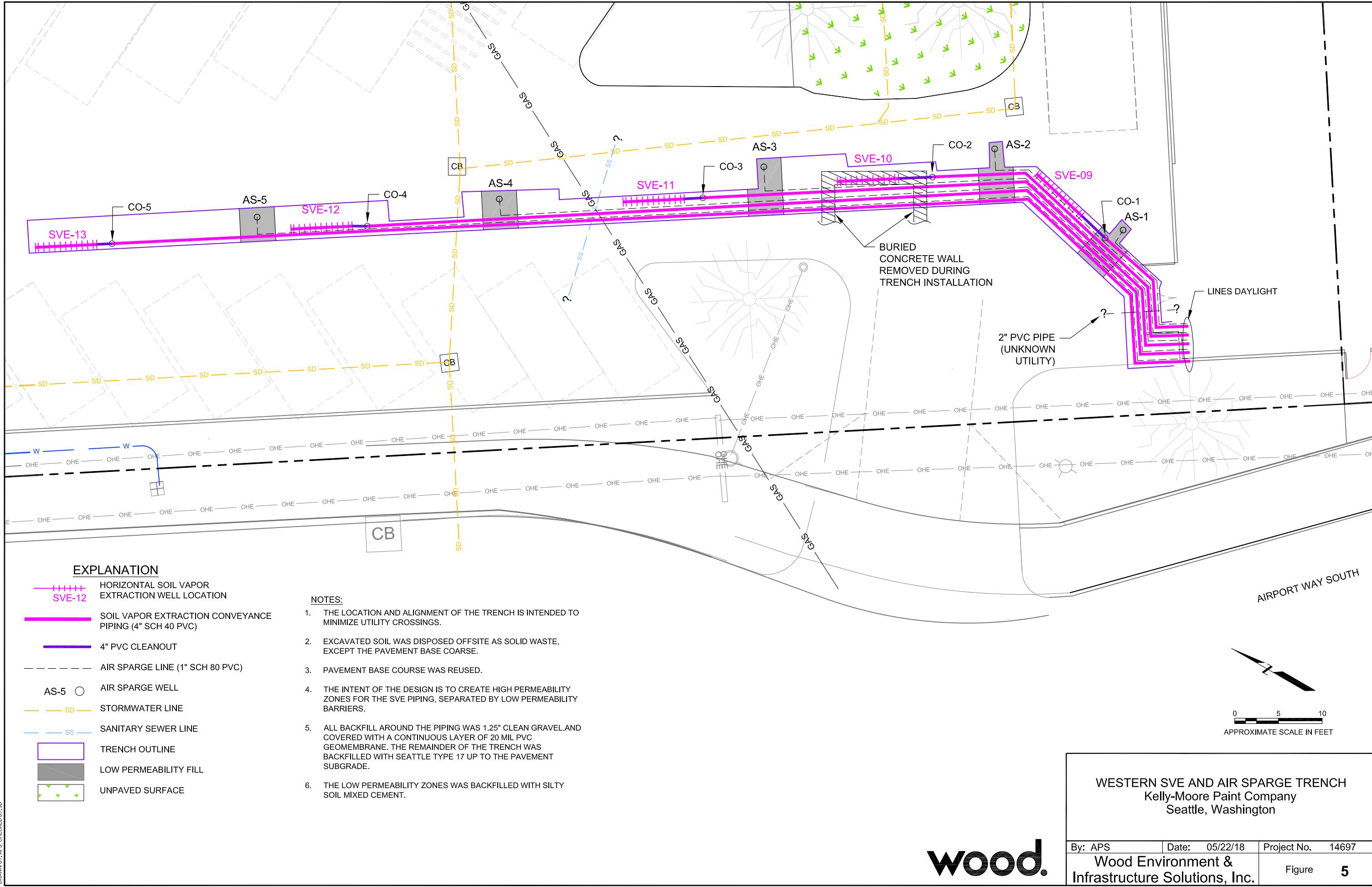
S. LUCILE ST.

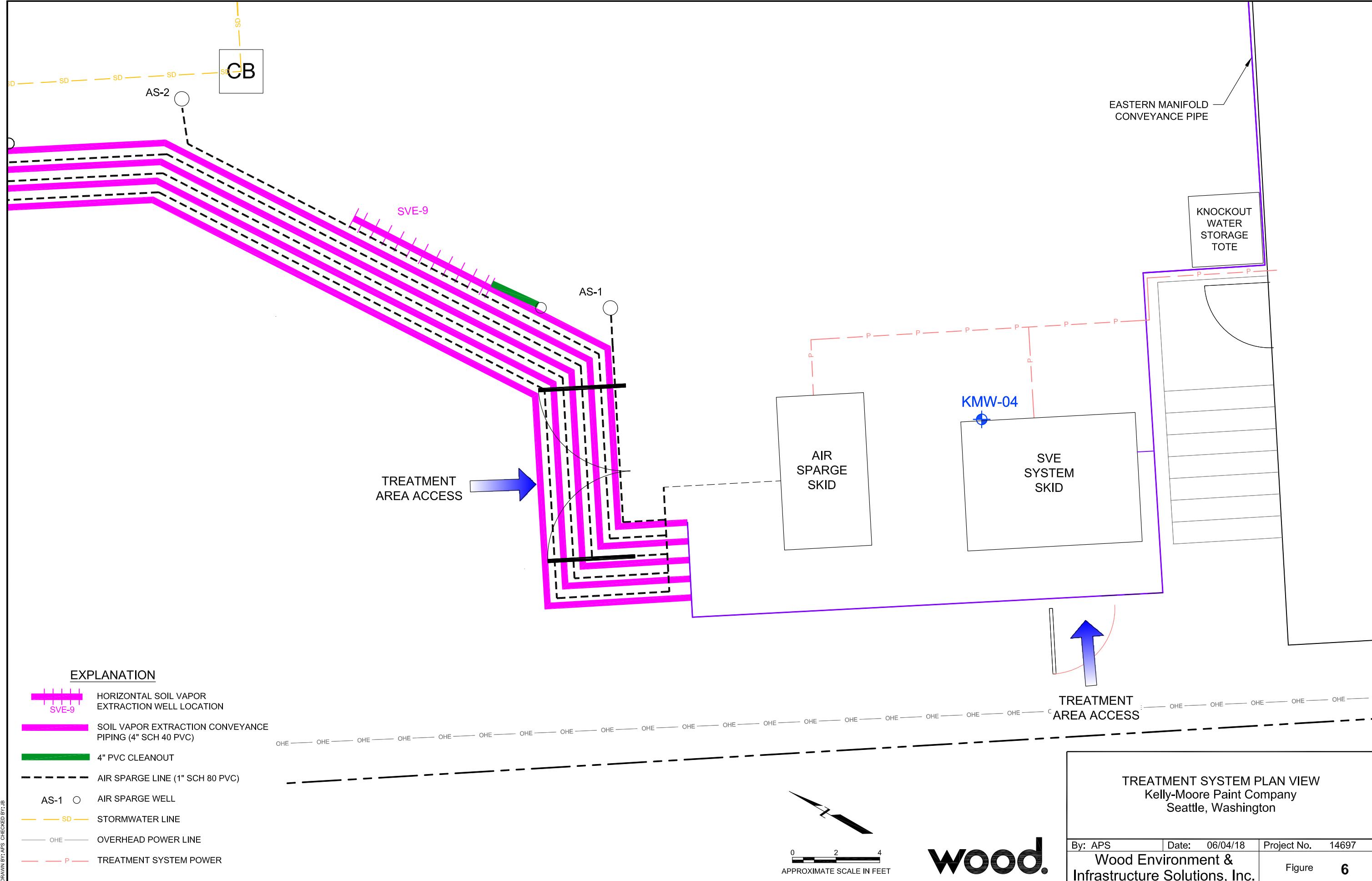
JANUARY 2018 WET SEASON  
GROUNDWATER CONTOURS  
Kelly-Moore Paint Company  
Seattle, Washington

**wood.**

By: APS	Date: 05/21/18	Project No. 14697
Wood Environment & Infrastructure Solutions, Inc.	Figure 3	







DRAWN BY: APS CHECKED BY: JB

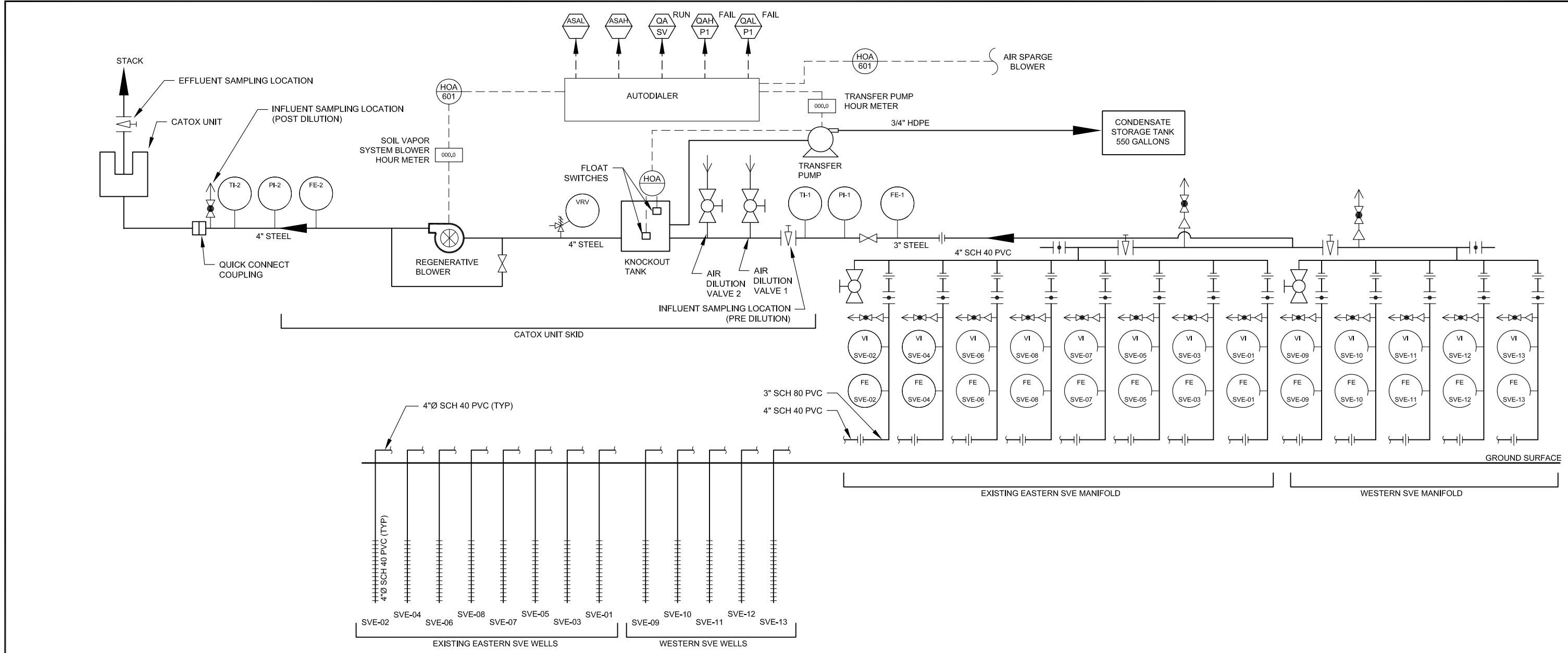
S:\14697\016\_2018-RIFS\KellyMoore-WesternSVETrench\_052218.dwg - Figure6-treatmentLayout - Jun. 04, 2018 12:24pm - adam.stenback

APPROXIMATE SCALE IN FEET

**wood.**

**TREATMENT SYSTEM PLAN VIEW**  
Kelly-Moore Paint Company  
Seattle, Washington

By: APS	Date: 06/04/18	Project No. 14697
Wood Environment & Infrastructure Solutions, Inc.	Figure	6



**wood.**

## **Tables**

**TABLE 1: MONITORING WELL SURVEY DATA**  
 Former Kelly-Moore Manufacturing Facility, Seattle, Washington

Well Name	WCS North Zone NAD 83(91) <sup>3</sup>		Ground Surface Elevation	Top of Casing Elevation
	Northing	Easting		
KMW-02R <sup>1</sup>	205743.9	1273010.4	22.01	21.63
KMW-03R <sup>1</sup>	205538.1	1273156.6	21.99	21.54
KMW-04 <sup>1</sup>	205423.6	1273115.0	18.90	18.56
KMW-06 <sup>1</sup>	205525.2	1273039.2	20.16	19.80
KMW-07 <sup>1</sup>	205713.7	1273034.0	22.00	21.63
KMW-08 <sup>1</sup>	205648.5	1273101.3	22.03	21.65
KMW-09 <sup>2</sup>	205508.9	1273025.5	18.60	18.14
KMW-10 <sup>2</sup>	205336.2	1272955.0	20.84	20.39

Notes:

1. Survey completed on June 30, 2016, by Duane Hartman and Associates.
2. Survey completed on December 13, 2016, by Duane Hartman and Associates.
3. Coordinate System and Zone: Washington State Plane, North Zone Coordinates.  
 Horizontal Datum: NAD 83(91), North Zone, US feet.  
 Vertical Datum: NAVD88, US feet.

Abbreviations:

- NAD = North American Datum  
 NAVD88 = North American Vertical Datum of 1988  
 WCS = Washington Coordinate System

**TABLE 2: GROUNDWATER ELEVATIONS**  
 Former Kelly-Moore Manufacturing Facility, Seattle, Washington

Well ID	TOC Elevation (feet) <sup>1</sup>	Date	Depth to Water (feet below TOC)	Groundwater Elevation (feet) <sup>1</sup>
KMW-02R	21.63 <sup>2</sup>	8/31/2017	9.58	12.05
		1/26/2018	7.56	14.07
KMW-03R	21.54 <sup>2</sup>	8/31/2017	9.52	12.02
		1/26/2018	7.87	13.67
KMW-04	18.56 <sup>2</sup>	8/31/2017	6.63	11.93
		1/26/2018	5.35	13.21
KMW-06	19.80 <sup>2</sup>	8/31/2017	7.87	11.93
		1/26/2018	6.48	13.32
KMW-07	21.63 <sup>2</sup>	8/31/2017	9.57	12.06
		1/26/2018	7.93	13.70
KMW-08	21.65 <sup>2</sup>	8/31/2017	9.59	12.06
		1/26/2018	7.72	13.93
KMW-09	18.14 <sup>3</sup>	8/31/2017	6.24	11.90
		1/26/2018	4.86	13.28
KMW-10	20.39 <sup>3</sup>	8/31/2017	8.61	11.78
		1/26/2018	7.51	12.88

Notes:

1. Elevations in feet above mean sea level.
2. Surveys conducted on June 30, 2016. Vertical Datum: NAVD 88. Bench Mark: City of Seattle 2" Brass disk 3805-3801 located at the SW corner of Airport Way South and Corson Ave South, Elevation 18.532 feet
3. Surveys conducted on December 13, 2016. Same datum as June 30, 2016.

Abbreviations:

NAVD88 = North American Vertical Datum of 1988  
 TOC = top of casing

**TABLE 3: GROUNDWATER PARAMETERS**  
 Former Kelly-Moore Manufacturing Facility, Seattle, Washington

<b>Monitoring Well</b>	<b>Date</b>	<b>pH</b>	<b>SC</b>	<b>ORP</b>	<b>DO</b>
			(ms/cm)	(mv)	(mg/L)
KMW-02R	8/31/2017	5.89	0.175	142.9	0.21
	1/26/2018	5.99	0.199	150.9	0.28
KMW-03R	8/31/2017	7.07	0.477	-117.2	0.15
	1/26/2018	7.27	0.454	-102.2	0.19
KMW-04	8/31/2017	6.31	0.485	-92.0	0.07
	1/25/2018	6.40	0.276	-40.0	0.58
KMW-06	8/31/2017	6.35	0.453	-90.3	0.10
	1/24/2018	6.56	0.314	-91.4	0.24
KMW-07	8/31/2017	6.02	0.283	56.2	0.15
	1/26/2018	6.32	0.280	56.1	0.32
KMW-08	8/31/2017	6.15	0.177	1.90	0.10
	1/26/2018	5.98	0.526	32.9	0.50
KMW-09	8/31/2017	6.32	0.415	-95.1	0.21
	1/24/2018	6.56	0.396	-79.5	0.40
KMW-10	8/31/2017	6.21	0.567	-86.3	0.15
	1/25/2018	6.46	0.656	-69.4	0.28

Abbreviations

DO = dissolved oxygen

mg/L = milligrams per liter

ms/cm = millisiemens per centimeter

mv = millivolts

ORP = oxidation reduction potential

SC = specific conductivity

**TABLE 4: GROUNDWATER ANALYTICAL RESULTS<sup>1, 2</sup>**  
 Former Kelly-Moore Manufacturing Facility, Seattle, Washington

**TABLE 4: GROUNDWATER ANALYTICAL RESULTS<sup>1, 2</sup>**  
Former Kelly-Moore Manufacturing Facility, Seattle, Washington

Analyte	KMW-04										KMW-05				KMW-06				KMW-07			
	3/28/2011	8/4/2011	8/4/2011 (D)	6/7/2013	3/10/2015	6/30/2016 (D)	11/11/2016	8/31/2017	1/25/2018	3/28/2011	8/4/2011	6/7/2013	6/30/2016	11/11/2016	8/31/2017	1/24/2018	7/1/2016	11/10/2016	9/1/2017	1/26/2018		
<b>Total Metals (µg/L)</b>																						
Arsenic	<b>12</b>	<b>12</b>	<b>10</b>	--	--	<b>12</b>	<b>12</b>	<b>20</b>	<b>14</b>	<b>7.6</b>	3.3 U	3.3 U	--	3.5	3.4	3.3 U	3.3 U	3.3 U	3.3 U	3.3 U	3.3 U	
Chromium	11 U	--	--	--	--	11 U	11 U	11 U	11 U	11 U	--	--	11 U	11 U	11 U	11 U	11 U	11 U	11 U	11 U	11 U	
Copper	--	11 U	11 U	--	--	--	--	--	--	--	--	--	11 U	--	--	--	--	--	--	--	--	
Lead	1.1 U	1.1 U	1.1 U	--	--	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	--	--	1.1 U	1.1 U	--	3.7	1.1	2.7	7.3	1.1 U	1.1 U	
Mercury	--	--	--	--	--	0.5 U	0.5 U	0.50 U	0.50 U	0.50 U	--	--	--	0.5 U	0.50 U	0.50 U	0.5 U	0.50 U	0.50 U	0.50 U	0.50 U	
Nickel	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	
Zinc	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	
<b>Polycyclic Aromatic Hydrocarbons (µg/L)</b>																						
Benzo(a)anthracene	0.012 U	0.0095 U	0.0095 U	--	--	<b>0.011</b>	<b>0.014</b>	0.0094 U	0.0096 U	0.010 U	0.0095 U	0.0095 U	--	<b>0.047</b>	<b>0.013</b>	<b>0.041</b>	<b>0.055</b>	0.0095 U	0.0095 U	0.0099 U	0.011 U	
Benzo(a)pyrene	0.012 U	0.0095 U	0.0095 U	--	--	0.0095 U	<b>0.015 J</b>	0.0094 U	0.0096 U	0.010 U	0.0095 U	0.0095 U	--	<b>0.038</b>	<b>0.022</b>	<b>0.033</b>	<b>0.071</b>	0.0095 U	0.0095 U	0.0099 U	0.011 U	
Benzo(b)fluoranthene	0.012 U	0.0095 U	0.0095 U	--	--	0.0095 U	<b>0.022 J</b>	0.0094 U	<b>0.0120</b>	0.010 U	0.0095 U	0.0095 U	--	<b>0.047</b>	<b>0.021</b>	<b>0.034</b>	<b>0.082</b>	0.0095 U	0.0095 U	0.0099 U	0.011 U	
Benzo(ghi)perylene	0.012 U	0.0095 U	0.0095 U	--	--	0.0095 U	<b>0.023 J</b>	0.0094 U	0.0096 U	0.010 U	0.0095 U	0.0095 U	--	<b>0.041</b>	<b>0.024</b>	<b>0.025</b>	<b>0.056</b>	0.0095 U	0.0095 U	0.0099 U	0.011 U	
Benzo(j,k)fluoranthene	0.012 U	0.0095 U	0.0095 U	--	--	0.0095 U	0.0095 U	0.0094 U	0.0096 U	0.010 U	0.0095 U	0.0095 U	--	<b>0.018</b>	<b>0.031</b>	<b>0.018</b>	<b>0.034</b>	0.0095 U	0.0095 U	0.0099 U	0.011 U	
Chrysene	0.012 U	0.0095 U	0.0095 U	--	--	0.0095 U	0.0095 U	0.0094 U	<b>0.0100</b>	0.010 U	0.0095 U	0.0095 U	--	<b>0.035</b>	<b>0.028</b>	<b>0.032</b>	<b>0.062</b>	0.0095 U	0.0095 U	0.0099 U	0.011 U	
Dibenz(a,h)anthracene	0.012 U	0.0095 U	0.0095 U	--	--	0.0095 U	0.0095 U	0.0094 U	0.0096 U	0.010 U	0.0095 U	0.0095 U	--	0.0095 U	0.0095 U	0.0095 U	<b>0.0110</b>	0.0095 U	0.0095 U	0.0099 U	0.011 U	
Indeno(1,2,3-cd)pyrene	0.012 U	0.0095 U	0.0095 U	--	--	0.0095 U	<b>0.016 J</b>	0.0094 U	0.0096 U	0.010 U	0.0095 U	0.0095 U	--	<b>0.028</b>	<b>0.023</b>	<b>0.023</b>	<b>0.054</b>	0.0095 U	0.0095 U	0.0099 U	0.011 U	
Total cPAHs	0.009 U	0.007 U	0.007 U	--	--	<b>0.008</b>	<b>0.020 J</b>	0.007 U	<b>0.008</b>	0.008 U	0.007 U	0.007 U	--	<b>0.050</b>	<b>0.030</b>	<b>0.045</b>	<b>0.095</b>	0.007 U	0.007 U	0.007 U	0.008 U	
<b>Polychlorinated Biphenyls (µg/L)</b>																						
Aroclor 1016	--	--	--	--	--	0.048 U	0.047 U	--	--	--	--	--	--	0.047 U	--	--	--	0.047 U	--	--	--	
Aroclor 1221	--	--	--	--	--	0.048 U	0.047 U	--	--	--	--	--	--	0.047 U	--	--	--	0.047 U	--	--	--	
Aroclor 1232	--	--	--	--	--	0.048 U	0.047 U	--	--	--	--	--	--	0.047 U	--	--	--	0.047 U	--	--	--	
Aroclor 1242	--	--	--	--	--	0.048 U	0.047 U	--	--	--	--	--	--	0.047 U	--	--	--	0.047 U	--	--	--	
Aroclor 1248	--	--	--	--	--	0.048 U	0.047 U	--	--	--	--	--	--	0.047 U	--	--	--	0.047 U	--	--	--	
Aroclor 1254	--	--	--	--	--	0.048 U	0.047 U	--	--	--	--	--	--	0.047 U	--	--	--	0.047 U	--	--	--	
Aroclor 1260	--	--	--	--	--	0.048 U	0.047 U	--	--	--	--	--	--	0.047 U	--	--	--	0.047 U	--	--	--	
<b>Total Petroleum Hydrocarbons (µg/L)</b>																						
Gasoline Range Organics	<b>75,000</b>	<b>55,000</b>	<b>50,000</b>	<b>48,000</b>	<b>27,000</b>	<b>27,000</b>	<b>27,000</b>	<b>63,000</b>	<b>8,000</b>	<b>9,000</b>	100 U	100 U	100 U	<b>2,700</b>	<b>850</b>	<b>1,600</b>	<b>1,300</b>	500 U	100 U	100 U	100 U	
Diesel Range Organics	--	--	--	--	--	<b>3,000 J</b>	<b>2,700 J</b>	6,400 U	<b>1,600</b>	<b>1,700</b>	--	--	--	<b>5,400 J</b>	<b>3,500</b>	<b>4,400</b>	<b>4,200</b>	260 U	260 U	280 U	260 U	
Lube Oil	--	--	--	--	--	<b>510</b>	<b>870</b>	410 U	440 U	410 U	--	--	--	<b>1,500 J</b>	<b>1,200</b>	<b>1,600</b>	<b>600</b>	410 U	420 U	450 U	410 U	
<b>Volatile Organic Compounds (µg/L)</b>																						
1,2,4-Trimethylbenzene	--	--	--	<b>77</b>	--	<b>44 J</b>	<b>65 J</b>	<b>160</b>	<b>97</b>	<b>110</b>	--	--	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	
1,3,5-Trimethylbenzene	--	--	--	20 U	--	20 U	<b>20</b>	<b>52</b>	<b>30</b>	50 U	--	--	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	
Acetone	--	--	--	500 U	--</td																	

**TABLE 4: GROUNDWATER ANALYTICAL RESULTS<sup>1, 2</sup>**  
Former Kelly-Moore Manufacturing Facility, Seattle, Washington

Analyte	KMW-08				KMW-09				KMW-10				1/25/2018 (D)
	7/1/2016	11/10/2016	9/1/2017	1/26/2018	11/11/2016	8/31/2017	1/24/2018	11/11/2016	8/31/2017	1/25/2018			
<b>Total Metals (µg/L)</b>													
Arsenic	3.3 U	3.3 U	3.3 U	<b>9.1</b>	<b>10</b>	<b>6.8</b>	<b>5.7</b>						
Chromium	11 U	11 U	11 U	11 U	11 U	11 U	11 U						
Copper	--	--	--	--	--	--	--	--	--	--	--		
Lead	1.1 U	1.1 U	<b>3.0</b>	1.1 U	1.1 U	<b>1.1</b>	1.1 U						
Mercury	0.5 U	0.50 U	0.50 U	0.50 U	0.50 U	0.50 U	0.50 U	0.50 U	0.50 U	0.50 U	0.50 U		
Nickel	--	--	--	--	--	--	--	--	--	--	--		
Zinc	--	--	--	--	--	--	--	--	--	--	--		
<b>Polycyclic Aromatic Hydrocarbons (µg/L)</b>													
Benzo(a)anthracene	<b>0.086</b>	<b>0.015</b>	<b>0.012</b>	<b>0.018</b>	0.0098 U	0.0098 U	<b>0.020</b>	0.0098 U	0.0094 U	0.011 U	0.011 U		
Benzo(a)pyrene	<b>0.11</b>	<b>0.013</b>	0.010 U	<b>0.015</b>	0.0098 U	0.0098 U	<b>0.021</b>	0.0098 U	0.0094 U	0.011 U	0.011 U		
Benzo(b)fluoranthene	<b>0.12</b>	<b>0.020</b>	0.010 U	<b>0.018</b>	0.0098 U	0.0098 U	<b>0.017</b>	0.0098 U	0.0094 U	0.011 U	0.011 U		
Benzo(ghi)perylene	<b>0.1</b>	0.0095 U	0.010 U	0.012 U	0.0098 U	0.0098 U	<b>0.011</b>	0.0098 U	0.0094 U	0.011 U	0.011 U		
Benzo(j,k)fluoranthene	<b>0.046</b>	0.0095 U	0.010 U	0.012 U	0.0098 U	0.0098 U	<b>0.015</b>	0.0098 U	0.0094 U	0.011 U	0.011 U		
Chrysene	<b>0.09</b>	<b>0.042</b>	<b>0.012</b>	<b>0.028</b>	0.0098 U	0.0098 U	<b>0.018</b>	0.0098 U	0.0094 U	0.011 U	0.011 U		
Dibenz(a,h)anthracene	<b>0.024</b>	0.0095 U	0.010 U	0.012 U	0.0098 U	0.0098 U	0.011 U	0.0098 U	0.0094 U	0.011 U	0.011 U		
Indeno(1,2,3-cd)pyrene	<b>0.063</b>	0.0095 U	0.010 U	0.012 U	0.0098 U	0.0098 U	<b>0.0130</b>	0.0098 U	0.0094 U	0.011 U	0.011 U		
Total cPAHs	<b>0.14</b>	<b>0.018</b>	<b>0.008</b>	<b>0.021</b>	0.007 U	0.007 U	<b>0.028</b>	0.007 U	0.007 U	0.008 U	0.008 U		
<b>Polychlorinated Biphenyls (µg/L)</b>													
Aroclor 1016	0.047 U	--	--	--	--	--	--	--	--	--	--		
Aroclor 1221	0.047 U	--	--	--	--	--	--	--	--	--	--		
Aroclor 1232	0.047 U	--	--	--	--	--	--	--	--	--	--		
Aroclor 1242	0.047 U	--	--	--	--	--	--	--	--	--	--		
Aroclor 1248	0.047 U	--	--	--	--	--	--	--	--	--	--		
Aroclor 1254	0.047 U	--	--	--	--	--	--	--	--	--	--		
Aroclor 1260	0.047 U	--	--	--	--	--	--	--	--	--	--		
<b>Total Petroleum Hydrocarbons (µg/L)</b>													
Gasoline Range Organics	<b>1,000</b>	<b>400</b>	<b>130</b>	<b>120</b>	<b>370</b>	<b>360</b>	<b>760</b>	<b>110</b>	<b>3,400</b>	<b>270</b>	<b>260</b>		
Diesel Range Organics	<b>770 J</b>	370 U	300 U	<b>450</b>	<b>1,700</b>	<b>2300</b>	<b>3100</b>	1,300 U	<b>1,800</b>	<b>2,300</b>	<b>2,300</b>		
Lube Oil	410 U	410 U	480 U	410 U	<b>660</b>	<b>810</b>	<b>690</b>	420 U	430 U	410 U	410 U		
<b>Volatile Organic Compounds (µg/L)</b>													
1,2,4-Trimethylbenzene	0.20 U	0.20 U	0.20 U	0.20 U	0.2 U	0.20 U	0.20 U	<b>3.7</b>	<b>53</b>	<b>2.7</b>	<b>5.4</b>		
1,3,5-Trimethylbenzene	0.20 U	0.20 U	0.20 U	0.20 U	0.2 U	0.20 U	0.20 U	<b>0.38</b>	<b>23</b>	<b>0.9</b>	<b>1.7</b>		
Acetone	12 U	<b>10</b>	5.0 U	5.0 U	5.0 U	5.0 U	5.0 U	5.0 U	100 U	5.0 U	5.0 U		
Benzene	0.20 U	<b>0.20</b>	0.20 U	0.20 U	0.2 U	0.20 U	0.20 U	<b>0.7</b>	<b>8.2</b>	0.20 U	0.20 U		
Ethylbenzene	0.20 U	<b>0.31</b>	<b>0.36</b>	0.20 U	<b>3.1</b>	0.20 U	<b>0.44</b>	<b>1.6</b>	<b>810</b>	<b>14</b>	<b>30</b>		
m,p-Xylene	0.4 U	<b>0.76</b>	<b>0.69</b>	0.40 U	<b>0.51</b>	0.40 U	0.40 U	<b>11</b>	<b>1100</b>	<b>28</b>	<b>65</b>		
Naphthalene	<b>1.5</b>	1.3 U	1.0 U	1.0 U	1.3 U	1.0 U	1.0 U	1.3 U	20 U	1.0 U	1.0 U		
o-Xylene	0.20 U	<b>0.34</b>	0.20 U	0.20 U	0.2 U	0.20 U	0.20 U	<b>0.29</b>	<b>22</b>	<b>0.42</b>	<b>0.58</b>		
Toluene	1.0 U	1.0 U	1.0 U	1.0 U	20 U	<b>1.0</b>	<b>1.0</b>						
Trichloroethene	0.20 U	0.20 U	0.20 U	0.20 U	0.2 U	0.20 U	0.20 U	0.2 U	4.0 U	0.20 U	0.20 U		
Vinyl Chloride	0.20 U	0.20 U	0.20 U	0.20 U	0.2 U	0.20 U	0.20 U	0.2 U	4.0 U	0.20 U	0.20 U		

**Notes:**

1. Data qualifiers are as follows:  
U = The analyte was not detected at the reporting limit indicated.  
J = The value is an estimate.  
UJ = The analyte was not detected at the estimated reporting limit indicated.
2. Bold values indicate detections.
3. KMW-02 was abandoned by backfilling with bentonite on February 4, 2015, and was replaced June 28, 2016.
4. KMW-03 was destroyed during building demolition on June 3, 2015, and was replaced June 27, 2016.

**Abbreviations:**

- = not analyzed
- µg/L = micrograms per liter
- cPAHS = carcinogenic polycyclic aromatic hydrocarbons
- (D) = duplicate sample collected
- mg/L - milligram per liter
- R = replaced

**TABLE 5: SOIL VAPOR EXTRACTION SYSTEM ANALYTICAL SUMMARY<sup>1, 2</sup>**

Former Kelly-Moore Manufacturing Facility, Seattle, Washington

Sample	Date	Benzene (ppmv)	TPH as Hexane (ppmv)
<b>Western SVE Wells</b>			
SVE-09	11/7/17	<b>0.069</b>	<b>310</b>
SVE-10	11/7/17	<b>0.53</b>	<b>820 J</b>
SVE-11	11/7/17	<b>0.069</b>	<b>220</b>
SVE-12	11/7/17	<b>0.44</b>	<b>1,400 J</b>
SVE-13	11/7/17	<b>0.23</b>	<b>600 J</b>
<b>Eastern SVE Wells</b>			
SVE-02	11/7/17	<0.03	<b>3.4</b>
SVE-04	11/7/17	<0.03	<b>310</b>
SVE-06	11/7/17	<b>0.041</b>	<b>280</b>
SVE-08	11/7/17	<0.03	<b>65</b>
SVE-07	11/7/17	<0.03	<b>82</b>
SVE-05	11/7/17	<b>0.50</b>	<b>2,200 J</b>
SVE-03	11/7/17	<b>1.1</b>	<b>1,900 J</b>
SVE-01	11/7/17	<b>0.14</b>	<b>450</b>
<b>Influent</b>	11/7/17	<b>0.18</b>	<b>650 J</b>
<b>Effluent</b>	11/7/17	<0.03	<b>28</b>

Notes:

1. **Bold** numbers indicate results above the reporting limits.
2. J = the result is estimated because the concentration exceeded the calibration range of the instrument.

Abbreviations:

ppmv = parts per million volume

TPH = total petroleum hydrocarbons

**TABLE 6: FLAME IONIZATION DETECTOR READINGS**  
 Former Kelly-Moore Manufacturing Facility, Seattle, Washington

Date	FID Readings <sup>1</sup> (ppmv)				
	12/13/2017	1/16/2018	2/13/2018	3/21/2018	4/24/2018
<b>Western SVE Wells</b>					
West Manifold	19.4	NM	NM	NM	NM
SVE-09	NM	0	53.5	53.5	7.6
SVE-10	54.5	55.5	258.1	258.1	14.4
SVE-11	NM	6.7	0.7	0.7	6.2
SVE-12	30.7	10.7	65.2	65.2	14.8
SVE-13	11.7	4.1	27.6	27.6	7.4
<b>Eastern SVE Wells</b>					
East Manifold	106.5	NM	NM	NM	NM
SVE-02	NM	0	0.4	0	0.1
SVE-04	NM	16.6	8.3	0	0.5
SVE-06	NM	52.5	24.5	20.8	13.9
SVE-08	NM	1.9	3.5	1.2	6.3
SVE-07	NM	4.7	4.2	24.8	7.5
SVE-05	1,105	828.9	483.7	1,811	43
SVE-03	446.8	246.7	85.9	622.8	21.8
SVE-01	11.0	6.3	2.7	2.1	5.8
<b>Influent</b>	85.6	72.3	16.4	45.2	6.8
<b>Effluent</b>	1.2	0.2	0.6	0.7	0.2

Notes:

1. FID calibrated to 100 parts per million hexane.

Abbreviations:

- FID = flame ionization detector  
 NM = not measured  
 ppmv = parts per million by volume  
 SVE = soil vapor extraction

**TABLE 7: SOIL VAPOR EXTRACTION SYSTEM REMOVAL EFFICIENCIES**

Former Kelly-Moore Manufacturing Facility, Seattle, Washington

<b>Date</b>	<b>Influent</b>	<b>Effluent</b>	<b>Removal Efficiency<sup>1</sup></b>
	<b>(ppm)</b>	<b>(ppm)</b>	<b>(percent)</b>
11/7/17	358.0	0.0	100.0%
11/15/17	547.5	3.7	99.3%
11/21/17	123.7	1.9	98.5%
12/5/17	59.6	0.2	99.7%
12/13/17	85.6	1.2	98.6%
1/16/18	72.3	0.2	99.7%
2/13/18	16.4	0.6	96.3%
3/21/18	45.0	0.7	98.4%
4/24/18	6.8	0.2	97.1%

Notes:

1. Removal efficiencies exceed the Puget Sound Clean Air Agency 95 percent standard for removal efficiency.

Abbreviations:

ppm - parts per million

**wood.**

## **Appendix A**



## **GROUNDWATER SAMPLING LOG**

### **Low Flow Sampling**

MONITORING WELL/PIEZOMETER NUMBER KMW-02R  
Project Name: kelly - Moore  
Project Number: 14697.00008 Date: 9/1/17  
Location: Seattle WA Weather Conditions: Sunny 60's  
Sampler: BH Wind Speed/Direction: N/A

## WELL INFORMATION

Casing Diameter (in): 2 Groundwater Elevation (ft): \_\_\_\_\_  
Top of Casing Elevation (ft): \_\_\_\_\_ Depth of Well Casing (ft): \_\_\_\_\_  
Initial Depth to Water (ft): 8.58 Actual Purge Volume (gal): 2  
Wellhead Condition: Needs Rock

## PURGING MEASUREMENTS

Sample ID No.: KMW-02R-9-1-17  
Water Level Ind. Model & No.: Sensus Interface Model 122  
ORP/DO Meter Model & No.: XSI PRO DSS  
Purge Equipment Used: Peristaltic pump w/ dedicated tubing  
Sampling Equipment Used: SAA  
Purge Start Time: 0853  
Purge Completion Time: 1035  
Average Purge Rate (mL/min): 700  
Analytical Lab: Onsite  
Sample Collection Time: 0935  
Purging Method: SAA  
Sample Containers Used: Lab provided  
Chemical Analyses: See COE  
Other Field Observations: Purge water has strong odor -



## **GROUNDWATER SAMPLING LOG**

### **Low Flow Sampling**

**MONITORING WELL/P**  
Project Name: Kelly Moore  
Project Number: 14697.00008  
Location: Seattle WA  
Sampler: RK

Casing Diameter (in): 2  
Top of Casing Elevation (ft): 953  
Initial Depth to Water (ft): 953  
Wellhead Condition: needs rock

Date: 9-1-17  
Weather Conditions: Sunny 70's  
Wind Speed/Direction: NA

## WELL INFORMATION

Casing Diameter (in): 2  
Top of Casing Elevation (ft): 953  
Initial Depth to Water (ft): 953  
Wellhead Condition: needs rock

**Groundwater Elevation (ft):** \_\_\_\_\_  
**Depth of Well Casing (ft):** \_\_\_\_\_  
**Actual Purge Volume (gal):** 2

## PURGING MEASUREMENTS

Sample ID No.: KMW-03R:9.1.17  
Water Level Ind. Model & No.: Salinst Interface Model 12C  
ORP/DO Meter Model & No.: YSI PRO PDS  
Purge Equipment Used: Peri pump w/ dedicated tubing  
Sampling Equipment Used: SPA

Purge Start Time: 1405  
Purge Completion Time: 1500  
Average Purge Rate (mL/min): 200  
Analytical Lab: Ousler

Sample Collection Time: 1440  
Purging Method: SAT  
Sample Containers Used: Lab provided  
Chemical Analyses: See CCR

Other Field Observations: Strong exuvia



## **GROUNDWATER SAMPLING LOG**

### **Low Flow Sampling**

Project Name: Kelly-Moore MONITORING WELL/PIEZOMETER NUMBER KMW-09

Project Number: 14697.00008  
Location: Seattle, WA  
Sampler: BTT

Date: 8-31-17  
Weather Conditions: Sunny 70's

Wind Speed/Direction: 5 mph - west

## **WELL INFORMATION**

Casing Diameter (in): 2  
Top of Casing Elevation (ft): \_\_\_\_\_  
Initial Depth to Water (ft): 6.63  
Wellhead Condition: Flooded

Groundwater Elevation (ft): \_\_\_\_\_  
Depth of Well Casing (ft): \_\_\_\_\_  
Actual Purge Volume (gal): 2.5

## PURGING MEASUREMENTS

Sample ID No.: EHW-09.8-31-17

Water Level Ind. Model & No.: Solust Interface Model 022

**ORP/DO Meter Model & No.:** YSI Pro DSS

**Purge Equipment Used:** Perf pump w/ dedicated tubing

### **Sampling Equipment Used:**

**Purge Start Time:**

**Purge Completion Time:**

### Average Purge Rate (mL/min)

Analytical Lab: Onsite

**Sample Collection Time:**

#### **Purging Method:**

**Sample Containers Used:**

Chemical Analyses: *See CAC*

**Other Field Observations:** WC meter periodically produced sustained tone (prolonged). Range of DOPUS of possible product  $\leq 0.01$  @ 1.65'. Pump water has sheen





## **GROUNDWATER SAMPLING LOG**

### **Low Flow Sampling**

MONITORING WELL/PIEZOMETER NUMBER KMW-07

**Project Name:** Kelly-Moore

Project Number: 14697.0008

Location: Gatlinburg

Sampler: BTH

Date: 9.1.17

Weather Conditions: Sunny 70's

Wind Speed/Direction: NA

## WELL INFORMATION

**Casing Diameter (in):** 2

### **Groundwater Elevation (ft):**

**Top of Casing Elevation (ft):**

**Depth of Well Casing (ft):**

Initial Depth to Water (ft): 9.58

Actual Purge Volume (gal): 7.5

Wellhead Condition: Needs repair

## PURGING MEASUREMENTS

Sample ID No.: KMW-07.9.1.17

Water Level Ind. Model & No.: Solinst Interface Model 122

**ORP/DO Meter Model & No.:** VSI PRO PPS

**Purge Equipment Used:** *No air purges w/ de-activated turbine*

### **Sampling Equipment Used:**

**Purge Start Time:**

Purge Start Time: 105  
Purge Completion Time: 1330

**Sample Collection Time:** 1150

#### Purging Method:

Sample Containers Used: Cabs provided

Chemical Analyses: See ~~the~~

Other Field Observations: Sustained tone (Product) : Range of depth of possible product 50.0' @ 9.64' range water has strong odor

# GROUNDWATER SAMPLING LOG

## Low Flow Sampling

### MONITORING WELL/PIEZOMETER NUMBER KMW-08

Project Name: kelly moore  
Project Number: 14697.00008  
Location: Seattle WA  
Sampler: BTT

Date: 9.1.17  
Weather Conditions: Sunny 70° S  
Wind Speed/Direction: NA

#### WELL INFORMATION

Casing Diameter (in): 2  
Top of Casing Elevation (ft):   
Initial Depth to Water (ft): 9.60  
Wellhead Condition: needs tock  
Groundwater Elevation (ft):   
Depth of Well Casing (ft):   
Actual Purge Volume (gal): 5

#### PURGING MEASUREMENTS

WL (ft btoc)	Time	pH (std. units)	SC (ms/cm)	Temp. (°C)	ORP (mv)	DO (mg/L)	Turbidity (NTUs)	Notes
9.72	1242	6.21	0.177	14.2	-3.9	0.066	55.2	
9.72	1247	6.14	0.173	14.3	11.9	0.24	271.4	
9.72	1252	6.11	0.172	14.2	17.3	0.123	443.2	
9.72	1257	6.12	0.174	14.2	15.1	0.19	607.2	
9.70	1302	6.13	0.174	14.4	11.9	0.16	653.0	
9.70	1307	6.14	0.175	14.3	8.5	0.14	543.2	
9.70	1310	6.14	0.176	14.3	7.9	0.14	634.2	
9.70	1313	6.14	0.176	14.3	7.5	0.113	720.5	
9.70	1316	6.15	0.176	14.3	5.7	0.12	625.7	
9.70	1319	6.15	0.177	14.3	4.4	0.12	696.7	
9.70	1322	6.18	0.177	14.3	2.5	0.11	695.7	
9.70	BTT	9.1.17						
9.71	1325	6.15	0.177	14.3	2.9	0.11	765.0	
9.71	1328	6.15	0.177	14.3	2.4	0.11	795.2	

Sample ID No.: KMW-08-9.1.17

Water Level Ind. Model & No.: Satrist Interface Model 12C

ORP/DO Meter Model & No.: XSI PRO DPS

Purge Equipment Used: Peri pump w/ dechlorinated tubing

Sampling Equipment Used: SMA

Purge Start Time: 1240

Purge Completion Time: 1400

Average Purge Rate (mL/min): 200

Analytical Lab: Oncile

Sample Collection Time: 1340

Purging Method: SMA

Sample Containers Used: Lab provided

Chemical Analyses: See doc

Other Field Observations: Strong odor





## **GROUNDWATER SAMPLING LOG**

### **Low Flow Sampling**

**Project Name:** Kelly-Moore      **MONITORING WELL/PIEZOMETER NUMBER** KMW-09

Project Number: 14697,00008  
Location: Seattle WA  
Sampler: BH

Date: 8-31-17  
Weather Conditions: overcast 70°s  
Wind Speed/Direction: 9 mph / East

## WELL INFORMATION

Casing Diameter (in): 2  
Top of Casing Elevation (ft): \_\_\_\_\_  
Initial Depth to Water (ft): 612.3  
Wellhead Condition: OK - Secure

**Groundwater Elevation (ft):** \_\_\_\_\_  
**Depth of Well Casing (ft):** \_\_\_\_\_  
**Actual Purge Volume (gal):** \_\_\_\_\_ / 1.5

## PURGING MEASUREMENTS

Sample ID No.: KMW-09-8.31.17

Water Level Ind. Model & No.: Selbst Interface Modell 122

**ORP/DO Meter Model & No.:**

#### **Purge Equipment Used:**

#### **Large Equipment Used:**

### **Sampling Equipment Used:**

#### Purge Start Time: Burns Consultation

#### Purge Completion Time:

**Average Purge Rate (mL/min):**

**Analytical Lab:** Cust #

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**Sample Collection Time:** 1350

#### **Purging Method:**

Sample Containers Used: Lab provided

**Chemical Analyses:**

#### **Other Field Observations:**







## GROUNDWATER SAMPLING LOG

### Low Flow Sampling

**MONITORING WELL/PIEZOMETER NUMBER** KMW-03R

**Project Name:** \_\_\_\_\_ Date: 1/26/08

**Project Number:** 014b910060 Weather Conditions: 40s overcast

**Location:** Seattle, WA

**Sampler:** P. Blaett Wind Speed/Direction: -

### WELL INFORMATION

**Casing Diameter (in):** 2" **Groundwater Elevation (ft):** \_\_\_\_\_

**Top of Casing Elevation (ft):** \_\_\_\_\_

**Initial Depth to Water (ft):** 7.89 **Depth of Well Casing (ft):** \_\_\_\_\_

**Wellhead Condition:** \_\_\_\_\_ **Actual Purge Volume (gal):** 8.5 gal

### PURGING MEASUREMENTS

WL (ft btoc)	Time	pH (std. units)	SC (ms/cm)	Temp. (°C)	ORP (mv)	DO (mg/L)	Turbidity (NTUs)	Notes
7.91	0858	6.92	13.9	17.9	+56.0	3.71	8.4	
7.95	0901	7.10	243.2	17.5	8.0	0.79	7.0	
7.99	0904	7.16	299.9	17.4	-9.9	0.65	4.9	
7.99	0907	7.21	332.8	17.3	-3.5	0.48	4.4	
7.99	0910	7.21	455.5	17.0	-61.1	0.35	6.3	
7.99	0913	7.26	447.7	16.9	-79.3	0.29	7.9	
7.99	0916	7.26	449.0	16.8	-84.0	0.29	11.2	
7.99	0919	7.26	451.4	16.9	-86.7	0.29	13.3	
8.00	0922	7.26	453.5	16.8	-90.6	0.27	21.0	
8.00	0925	7.26	454.0	16.8	-93.1	0.26	29.2	
8.00	0928	7.26	454.8	16.8	-95.1	0.25	31.0	
8.00	0931	7.26	455.6	16.8	-96.5	0.24	43.0	
8.00	0934	7.26	456.0	16.8	-98.2	0.24	59.6	
8.00	0937	7.26	455.7	16.8	-100.0	0.23	75.7	

**Sample ID No.:** KMW-03R-012618

→ See pg. 2

**Water Level Ind. Model & No.:** \_\_\_\_\_

**ORP/DO Meter Model & No.:** \_\_\_\_\_

**Purge Equipment Used:** \_\_\_\_\_

**Sampling Equipment Used:** \_\_\_\_\_

**Purge Start Time:** 0855 **Sample Collection Time:** 0945

**Purge Completion Time:** 1008 **Purging Method:** low flow

**Average Purge Rate (mL/min):** 160 mL/min **Sample Containers Used:** \_\_\_\_\_

**Analytical Lab:** CGS/inside environment **Chemical Analyses:** \_\_\_\_\_

**Other Field Observations:** suction in well, when removed reddish brown  
black residue, some suction in well

**Final DTW = 8.04**

## **GROUNDWATER SAMPLING LOG**

### **Low Flow Sampling**



MONITORING WELL/PIEZOMETER NUMBER KMW-037

**Project Name:** \_\_\_\_\_

**Project Number:** 123 **Date:** \_\_\_\_\_ **Weather Conditions:** \_\_\_\_\_

Location: *[Signature]* Father Conditions: \_\_\_\_\_

**Sampler:** \_\_\_\_\_ **Wind Speed/Direction:** \_\_\_\_\_

Date: \_\_\_\_\_

**Weather Conditions:** \_\_\_\_\_

Wind Speed/Direction:

## WELL INFORMATION

**Casing Diameter (in):** \_\_\_\_\_ **Groundwater Elevation (ft):** \_\_\_\_\_

**Top of Casing Elevation (ft):** \_\_\_\_\_ **Depth of Well Casing (ft):** \_\_\_\_\_

**Initial Depth to Water (ft):** \_\_\_\_\_ **Actual Purge Volume (gal):** \_\_\_\_\_

**Wellhead Condition:** \_\_\_\_\_

## PURGING MEASUREMENTS

**Sample ID No.:**

**Water Level Ind. Model & No.:**

**ORP/DO Meter Model & No.:** \_\_\_\_\_

**Purge Equipment Used:** \_\_\_\_\_

**Sampling Equipment Used:** \_\_\_\_\_

**Purge Start Time:** \_\_\_\_\_ **Sample Collection Time:** \_\_\_\_\_

**Purge Completion Time:** \_\_\_\_\_ **Purging Method:** \_\_\_\_\_

**Average Purge Rate (mL/min):** \_\_\_\_\_ **Sample Containers Used:** \_\_\_\_\_

Analytical Lab: \_\_\_\_\_ Chemical Analyses: \_\_\_\_\_

**Sample Collection Time:** \_\_\_\_\_

## Purging Method:

**Sample Containers Used:** \_\_\_\_\_

**Chemical Analyses:** \_\_\_\_\_

**Other Field Observations:**



**GROUNDWATER SAMPLING LOG**  
**Low Flow Sampling**

amec  
foster  
wheeler

**MONITORING WELL/PIEZOMETER NUMBER** KMW-04

Project Name: Former Kelly Moore Date: 1/25/08  
Project Number: 0146970060 Weather Conditions: rainy 40°  
Location: Seattle, WA  
Sampler: K. Black Wind Speed/Direction: /

**WELL INFORMATION**

Casing Diameter (in): 2 1/4 Groundwater Elevation (ft): \_\_\_\_\_  
Top of Casing Elevation (ft): \_\_\_\_\_ Depth of Well Casing (ft): \_\_\_\_\_  
Initial Depth to Water (ft): 5.36 Actual Purge Volume (gal): 2  
Wellhead Condition: \_\_\_\_\_

**PURGING MEASUREMENTS**

WL (ft btoc)	Time	pH (std. units)	SC (ms/cm)	Temp. (°C)	ORP (mv)	DO (mg/L)	Turbidity (NTUs)	Notes
5.36	1143	6.73	188.2	11.4	-2.4	1.99	41.7	
5.37	1146	6.48	201.4	11.4	-30.6	0.65	31.5	
5.37	1149	6.47	218.6	11.5	-27.9	0.53	28.7	
5.37	1152	6.45	229.6	11.5	-32.2	0.51	27.8	
5.37	1155	6.42	249.6	11.5	-36.5	0.46	25.8	
5.37	1158	6.41	264.3	11.5	-39.6	0.43	24.4	
5.37	1201	6.42	262.2	11.4	-40.2	0.60	23.8	
5.37	1204	6.41	269.9	11.3	-41.0	0.56	23.6	
5.37	1207	6.40	274.6	11.4	-40.6	0.56	19.3	
5.38	1210	6.40	278.3	11.4	-41.7	0.57	18.9	
5.38	1213	6.40	276.3	11.4	-40.0	0.58	18.3	

Sample ID No.: KMW-04 - 012518

Water Level Ind. Model & No.: \_\_\_\_\_

ORP/DO Meter Model & No.: \_\_\_\_\_

Purge Equipment Used: \_\_\_\_\_

Sampling Equipment Used: \_\_\_\_\_

Purge Start Time: 1140

Sample Collection Time: 1215

Purge Completion Time: 1240

Purging Method: low flow

Average Purge Rate (mL/min): 200 mL/min

Sample Containers Used: \_\_\_\_\_

Analytical Lab: \_\_\_\_\_

Chemical Analyses: \_\_\_\_\_

Other Field Observations: tubing present when open well

black residue on tubing when opened





## **GROUNDWATER SAMPLING LOG**

### **Low Flow Sampling**



MONITORING WELL/PIEZOMETER NUMBER KMW-08

Project Name: Telluride Moore Date: 1/26/18  
Project Number: 0146970060 Weather Conditions: 40°s rainy  
Location: Seattle, WA  
Sampler: K. Black Wind Speed/Direction: SW 10-15 mph

## WELL INFORMATION

Casing Diameter (in): 2      Groundwater Elevation (ft): \_\_\_\_\_  
Top of Casing Elevation (ft): \_\_\_\_\_      Depth of Well Casing (ft): \_\_\_\_\_  
Initial Depth to Water (ft): 7.77      Actual Purge Volume (gal): 1.5  
Wellhead Condition:

## PURGING MEASUREMENTS

Sample ID No.: KMW-08-P12618

Water Level Ind. Model & No.:

ORP/DO Meter Model & No.:

#### Purge Equipment Used:

#### **Sampling Equipment Used:**

**Purge Start Time:**

Purge Start Time:

Average Purge Rate (mL/min): 15 (P) mL/min

Analytical Lab: *present environmental*

Sample Collection Time: 11:00

#### **Sample Collection Purging Method:**

#### **Purging Method: Sample Containers Used:**

#### **Sample Containers & Chemical Analyses:**

**Other Field Observations:** tubercles in well when open; when removed  
orange brown in <sup>inner</sup> side on tubercles

Final DTW = 7.79

## **GROUNDWATER SAMPLING LOG**

### **Low Flow Sampling**



MONITORING WELL/PIEZOMETER NUMBER MW-09

Project Name: Former Kelly Moore Date: 1/24/18  
Project Number: 01469-0060 Weather Conditions: 50% sunny  
Location: Seattle, WA  
Sampler: \_\_\_\_\_ Wind Speed/Direction: /

## WELL INFORMATION

Casing Diameter (in): 2      Groundwater Elevation (ft): \_\_\_\_\_  
Top of Casing Elevation (ft): 1      Depth of Well Casing (ft): \_\_\_\_\_  
Initial Depth to Water (ft): 4,910      Actual Purge Volume (gal): 1 gal  
Wellhead Condition: \_\_\_\_\_

## PURGING MEASUREMENTS

Sample ID No.: KMW-09-012419

**Water Level Ind. Model & No.:** \_\_\_\_\_

**ORP/DO Meter Model & No.:**

**Purge Equipment Used:** \_\_\_\_\_

**Sampling Equipment Used:** The D-113-1-51-

Sampling Equipment Used: Hand Brew - #16

Purge Start Time: 1420 Sample Collection Time: 1445

Purge Completion Time: \_\_\_\_\_ Purging Method: \_\_\_\_\_

Average Purge Rate (mL/min): 110 mL/min      Sample Containers Used: \_\_\_\_\_

**Analytical Lab:** \_\_\_\_\_ **Chemical Analyses:** \_\_\_\_\_

20. What is the name of the author of the book?

**Other Field Observations:** tubing left in well from previous sampler

~~black. Residue) or until whom removed~~

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Digitized by srujanika@gmail.com

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Digitized by srujanika@gmail.com



**wood.**

## **Appendix B**



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

September 14, 2017

Crystal Thimsen  
AMEC Foster Wheeler  
Environment and Infrastructure, Inc.  
One Union Square  
600 University Street, Suite 600  
Seattle, WA 98101

Re: Analytical Data for Project 14697.00008  
Laboratory Reference No. 1709-019

Dear Crystal:

Enclosed are the analytical results and associated quality control data for samples submitted on September 1, 2017.

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

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

Sincerely,

A handwritten signature in black ink, appearing to read "DB" followed by a cursive surname.

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 14, 2017  
Samples Submitted: September 1, 2017  
Laboratory Reference: 1709-019  
Project: 14697.00008

### Case Narrative

Samples were collected on August 31 and September 1, 2017 and received by the laboratory on September 1, 2017. They were maintained at the laboratory at a temperature of 2°C to 6°C.

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

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

#### PAHs EPA 8270D/SIM Analysis

Sample KMW-09-8-31-17 had one surrogate recovery out of control limits. This is within allowance of our standard operating procedure as long as the recovery is above 10%.

#### Dissolved Metals by EPA 200.8/7470A Analysis

Samples KMW-02R-9-1-17, KMW-07-9-1-17, and KMW-03R-9-1-17 were refiltered.

**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.**



---

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Date of Report: September 14, 2017  
 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

### NWTPH-Gx

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-10-8-31-17</b>					
Laboratory ID:	09-019-01					
Gasoline	<b>3400</b>	100	NWTPH-Gx	9-7-17	9-7-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	90	61-118				
<b>Client ID:</b>	<b>KMW-09-8-31-17</b>					
Laboratory ID:	09-019-02					
Gasoline	<b>360</b>	100	NWTPH-Gx	9-7-17	9-7-17	O
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	82	61-118				
<b>Client ID:</b>	<b>KMW-06-8-31-17</b>					
Laboratory ID:	09-019-03					
Gasoline	<b>1600</b>	100	NWTPH-Gx	9-7-17	9-7-17	O
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	94	61-118				
<b>Client ID:</b>	<b>KMW-04-8-31-17</b>					
Laboratory ID:	09-019-04					
Gasoline	<b>8000</b>	100	NWTPH-Gx	9-7-17	9-7-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	92	61-118				
<b>Client ID:</b>	<b>KMW-02R-9-1-17</b>					
Laboratory ID:	09-019-05					
Gasoline	<b>ND</b>	100	NWTPH-Gx	9-7-17	9-7-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	88	61-118				
<b>Client ID:</b>	<b>KMW-07-9-1-17</b>					
Laboratory ID:	09-019-06					
Gasoline	<b>ND</b>	100	NWTPH-Gx	9-7-17	9-7-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	94	61-118				



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Date of Report: September 14, 2017  
 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

### NWTPH-Gx

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-08-9-1-17</b>					
Laboratory ID:	09-019-07					
Gasoline	<b>130</b>	100	NWTPH-Gx	9-13-17	9-13-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	89	61-118				
<b>Client ID:</b>	<b>KMW-03R-9-1-17</b>					
Laboratory ID:	09-019-08					
Gasoline	<b>270</b>	100	NWTPH-Gx	9-13-17	9-13-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	85	61-118				
<b>Client ID:</b>	<b>TRIP BLANK</b>					
Laboratory ID:	09-019-09					
Gasoline	<b>ND</b>	100	NWTPH-Gx	9-7-17	9-7-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	97	61-118				



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Date of Report: September 14, 2017  
 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

**NWTPH-Gx**  
**QUALITY CONTROL**

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB0907W1					
Gasoline	ND	100	NWTPH-Gx	9-7-17	9-7-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	89	61-118				
Laboratory ID:	MB0913W1					
Gasoline	ND	100	NWTPH-Gx	9-13-17	9-13-17	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	97	61-118				
Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD RPD Limit Flags
<b>DUPLICATE</b>						
Laboratory ID:	09-019-05					
	ORIG	DUP				
Gasoline	ND	ND	NA	NA	NA	NA 30
Surrogate:						
Fluorobenzene				88 103	61-118	



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Date of Report: September 14, 2017  
 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

### NWTPH-Dx

Matrix: Water  
 Units: mg/L (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-10-8-31-17</b>					
Laboratory ID:	09-019-01					
Diesel Range Organics	<b>1.8</b>	0.27	NWTPH-Dx	9-8-17	9-8-17	M
Lube Oil Range Organics	<b>ND</b>	0.43	NWTPH-Dx	9-8-17	9-8-17	
Surrogate: <i>o-Terphenyl</i>	Percent Recovery 91	Control Limits 50-150				
<b>Client ID:</b>	<b>KMW-09-8-31-17</b>					
Laboratory ID:	09-019-02					
Diesel Range Organics	<b>2.3</b>	0.27	NWTPH-Dx	9-8-17	9-8-17	
Lube Oil Range Organics	<b>0.81</b>	0.43	NWTPH-Dx	9-8-17	9-8-17	N1
Surrogate: <i>o-Terphenyl</i>	Percent Recovery 85	Control Limits 50-150				
<b>Client ID:</b>	<b>KMW-06-8-31-17</b>					
Laboratory ID:	09-019-03					
Diesel Range Organics	<b>4.4</b>	0.26	NWTPH-Dx	9-8-17	9-8-17	M
Lube Oil Range Organics	<b>ND</b>	1.6	NWTPH-Dx	9-8-17	9-8-17	U1
Surrogate: <i>o-Terphenyl</i>	Percent Recovery 106	Control Limits 50-150				
<b>Client ID:</b>	<b>KMW-04-8-31-17</b>					
Laboratory ID:	09-019-04					
Diesel Range Organics	<b>1.6</b>	0.27	NWTPH-Dx	9-8-17	9-8-17	M
Lube Oil Range Organics	<b>ND</b>	0.44	NWTPH-Dx	9-8-17	9-8-17	
Surrogate: <i>o-Terphenyl</i>	Percent Recovery 94	Control Limits 50-150				
<b>Client ID:</b>	<b>KMW-02R-9-1-17</b>					
Laboratory ID:	09-019-05					
Diesel Range Organics	<b>ND</b>	0.27	NWTPH-Dx	9-8-17	9-8-17	
Lube Oil Range Organics	<b>ND</b>	0.43	NWTPH-Dx	9-8-17	9-8-17	
Surrogate: <i>o-Terphenyl</i>	Percent Recovery 87	Control Limits 50-150				
<b>Client ID:</b>	<b>KMW-07-9-1-17</b>					
Laboratory ID:	09-019-06					
Diesel Range Organics	<b>ND</b>	0.28	NWTPH-Dx	9-8-17	9-8-17	
Lube Oil Range Organics	<b>ND</b>	0.45	NWTPH-Dx	9-8-17	9-8-17	
Surrogate: <i>o-Terphenyl</i>	Percent Recovery 108	Control Limits 50-150				



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Date of Report: September 14, 2017  
 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

### NWTPH-Dx

Matrix: Water  
 Units: mg/L (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-08-9-1-17</b>					
Laboratory ID:	09-019-07					
Diesel Range Organics	<b>ND</b>	0.30	NWTPH-Dx	9-8-17	9-8-17	
Lube Oil Range Organics	<b>ND</b>	0.48	NWTPH-Dx	9-8-17	9-8-17	

Surrogate:	Percent Recovery	Control Limits
<i>o-Terphenyl</i>	95	50-150

**Client ID:** KMW-03R-9-1-17  
 Laboratory ID: 09-019-08

Diesel Range Organics	<b>0.33</b>	0.27	NWTPH-Dx	9-8-17	9-8-17	M
Lube Oil Range Organics	<b>ND</b>	0.43	NWTPH-Dx	9-8-17	9-8-17	
Surrogate:	Percent Recovery	Control Limits				
<i>o-Terphenyl</i>	103	50-150				



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Date of Report: September 14, 2017  
 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

**NWTPH-Dx**  
**QUALITY CONTROL**

Matrix: Water  
 Units: mg/L (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB0908W1					
Diesel Range Organics	ND	0.25	NWTPH-Dx	9-8-17	9-8-17	
Lube Oil Range Organics	ND	0.40	NWTPH-Dx	9-8-17	9-8-17	
Surrogate: <i>o-Terphenyl</i>	Percent Recovery 85	Control Limits 50-150				

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>								
Laboratory ID:	09-019-05							
	ORIG	DUP						
Diesel Range	ND	ND	NA	NA	NA	NA	NA	NA
Lube Oil Range	ND	ND	NA	NA	NA	NA	NA	NA
Surrogate: <i>o-Terphenyl</i>				87	104	50-150		

**SPIKE BLANK**

Laboratory ID: SB0908W1

Diesel Fuel #2	0.815	1.00	NA	82	58-113	NA	NA
<i>Surrogate: <i>o-Terphenyl</i></i>							
				112	50-150		



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 Project: 14697.00008

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

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-10-8-31-17</b>					
<b>Laboratory ID:</b>	09-019-01					
Dichlorodifluoromethane	ND	4.0	EPA 8260C	9-7-17	9-7-17	
Chloromethane	ND	20	EPA 8260C	9-7-17	9-7-17	
Vinyl Chloride	ND	4.0	EPA 8260C	9-7-17	9-7-17	
Bromomethane	ND	7.2	EPA 8260C	9-7-17	9-7-17	
Chloroethane	ND	20	EPA 8260C	9-7-17	9-7-17	
Trichlorofluoromethane	ND	4.0	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloroethene	ND	4.0	EPA 8260C	9-7-17	9-7-17	
Acetone	ND	100	EPA 8260C	9-7-17	9-7-17	
Iodomethane	ND	40	EPA 8260C	9-7-17	9-7-17	
Carbon Disulfide	ND	4.0	EPA 8260C	9-7-17	9-7-17	
Methylene Chloride	ND	40	EPA 8260C	9-7-17	9-7-17	
(trans) 1,2-Dichloroethene	ND	4.0	EPA 8260C	9-7-17	9-7-17	
Methyl t-Butyl Ether	ND	4.0	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloroethane	ND	4.0	EPA 8260C	9-7-17	9-7-17	
Vinyl Acetate	ND	20	EPA 8260C	9-7-17	9-7-17	
2,2-Dichloropropane	ND	4.0	EPA 8260C	9-7-17	9-7-17	
(cis) 1,2-Dichloroethene	ND	4.0	EPA 8260C	9-7-17	9-7-17	
2-Butanone	ND	100	EPA 8260C	9-7-17	9-7-17	
Bromochloromethane	ND	4.0	EPA 8260C	9-7-17	9-7-17	
Chloroform	ND	4.0	EPA 8260C	9-7-17	9-7-17	
1,1,1-Trichloroethane	ND	4.0	EPA 8260C	9-7-17	9-7-17	
Carbon Tetrachloride	ND	4.0	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloropropene	ND	4.0	EPA 8260C	9-7-17	9-7-17	
Benzene	8.2	4.0	EPA 8260C	9-7-17	9-7-17	
1,2-Dichloroethane	ND	4.0	EPA 8260C	9-7-17	9-7-17	
Trichloroethene	ND	4.0	EPA 8260C	9-7-17	9-7-17	
1,2-Dichloropropane	ND	4.0	EPA 8260C	9-7-17	9-7-17	
Dibromomethane	ND	4.0	EPA 8260C	9-7-17	9-7-17	
Bromodichloromethane	ND	4.0	EPA 8260C	9-7-17	9-7-17	
2-Chloroethyl Vinyl Ether	ND	26	EPA 8260C	9-7-17	9-7-17	
(cis) 1,3-Dichloropropene	ND	4.0	EPA 8260C	9-7-17	9-7-17	
Methyl Isobutyl Ketone	ND	40	EPA 8260C	9-7-17	9-7-17	
Toluene	ND	20	EPA 8260C	9-7-17	9-7-17	
(trans) 1,3-Dichloropropene	ND	4.0	EPA 8260C	9-7-17	9-7-17	



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 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

**VOLATILES EPA 8260C**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-10-8-31-17</b>					
<b>Laboratory ID:</b>	09-019-01					
1,1,2-Trichloroethane	ND	4.0	EPA 8260C	9-7-17	9-7-17	
Tetrachloroethene	ND	4.0	EPA 8260C	9-7-17	9-7-17	
1,3-Dichloropropane	ND	4.0	EPA 8260C	9-7-17	9-7-17	
2-Hexanone	ND	40	EPA 8260C	9-7-17	9-7-17	
Dibromochloromethane	ND	4.0	EPA 8260C	9-7-17	9-7-17	
1,2-Dibromoethane	ND	4.0	EPA 8260C	9-7-17	9-7-17	
Chlorobenzene	ND	4.0	EPA 8260C	9-7-17	9-7-17	
1,1,1,2-Tetrachloroethane	ND	4.0	EPA 8260C	9-7-17	9-7-17	
Ethylbenzene	810	4.0	EPA 8260C	9-7-17	9-7-17	
m,p-Xylene	1100	20	EPA 8260C	9-8-17	9-8-17	
o-Xylene	22	4.0	EPA 8260C	9-7-17	9-7-17	
Styrene	ND	4.0	EPA 8260C	9-7-17	9-7-17	
Bromoform	ND	20	EPA 8260C	9-7-17	9-7-17	
Isopropylbenzene	20	4.0	EPA 8260C	9-7-17	9-7-17	
Bromobenzene	ND	4.0	EPA 8260C	9-7-17	9-7-17	
1,1,2,2-Tetrachloroethane	ND	4.0	EPA 8260C	9-7-17	9-7-17	
1,2,3-Trichloropropane	ND	4.0	EPA 8260C	9-7-17	9-7-17	
n-Propylbenzene	18	4.0	EPA 8260C	9-7-17	9-7-17	
2-Chlorotoluene	ND	4.0	EPA 8260C	9-7-17	9-7-17	
4-Chlorotoluene	ND	4.0	EPA 8260C	9-7-17	9-7-17	
1,3,5-Trimethylbenzene	23	4.0	EPA 8260C	9-7-17	9-7-17	
tert-Butylbenzene	ND	4.0	EPA 8260C	9-7-17	9-7-17	
1,2,4-Trimethylbenzene	53	4.0	EPA 8260C	9-7-17	9-7-17	
sec-Butylbenzene	ND	4.0	EPA 8260C	9-7-17	9-7-17	
1,3-Dichlorobenzene	ND	4.0	EPA 8260C	9-7-17	9-7-17	
p-Isopropyltoluene	ND	4.0	EPA 8260C	9-7-17	9-7-17	
1,4-Dichlorobenzene	ND	4.0	EPA 8260C	9-7-17	9-7-17	
1,2-Dichlorobenzene	ND	4.0	EPA 8260C	9-7-17	9-7-17	
n-Butylbenzene	ND	4.0	EPA 8260C	9-7-17	9-7-17	
1,2-Dibromo-3-chloropropane	ND	20	EPA 8260C	9-7-17	9-7-17	
1,2,4-Trichlorobenzene	ND	4.0	EPA 8260C	9-7-17	9-7-17	
Hexachlorobutadiene	ND	4.0	EPA 8260C	9-7-17	9-7-17	
Naphthalene	ND	20	EPA 8260C	9-7-17	9-7-17	
1,2,3-Trichlorobenzene	ND	4.0	EPA 8260C	9-7-17	9-7-17	
<b>Surrogate:</b>	<b>Percent Recovery</b>	<b>Control Limits</b>				
Dibromofluoromethane	99	77-129				
Toluene-d8	98	80-127				
4-Bromofluorobenzene	98	78-125				



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 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

**VOLATILES EPA 8260C**  
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Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	KMW-09-8-31-17					
Laboratory ID:	09-019-02					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Chloromethane	ND	1.0	EPA 8260C	9-7-17	9-7-17	
Vinyl Chloride	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromomethane	ND	0.36	EPA 8260C	9-7-17	9-7-17	
Chloroethane	ND	1.0	EPA 8260C	9-7-17	9-7-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Acetone	ND	5.0	EPA 8260C	9-7-17	9-7-17	
Iodomethane	ND	2.0	EPA 8260C	9-7-17	9-7-17	
Carbon Disulfide	0.21	0.20	EPA 8260C	9-7-17	9-7-17	
Methylene Chloride	ND	2.0	EPA 8260C	9-7-17	9-7-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Vinyl Acetate	ND	1.0	EPA 8260C	9-7-17	9-7-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Butanone	ND	5.0	EPA 8260C	9-7-17	9-7-17	
Bromochloromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Chloroform	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Benzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Trichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Dibromomethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromodichloromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Chloroethyl Vinyl Ether	ND	1.3	EPA 8260C	9-7-17	9-7-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	9-7-17	9-7-17	
Toluene	ND	1.0	EPA 8260C	9-7-17	9-7-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-7-17	9-7-17	



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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>KMW-09-8-31-17</b>					
<b>Laboratory ID:</b>	09-019-02					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Tetrachloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Hexanone	ND	2.0	EPA 8260C	9-7-17	9-7-17	
Dibromochloromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Chlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Ethylbenzene	0.36	0.20	EPA 8260C	9-7-17	9-7-17	
m,p-Xylene	0.69	0.40	EPA 8260C	9-7-17	9-7-17	
o-Xylene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Styrene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromoform	ND	1.0	EPA 8260C	9-7-17	9-7-17	
Isopropylbenzene	11	0.20	EPA 8260C	9-7-17	9-7-17	
Bromobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
n-Propylbenzene	12	0.20	EPA 8260C	9-7-17	9-7-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
tert-Butylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
sec-Butylbenzene	0.91	0.20	EPA 8260C	9-7-17	9-7-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
n-Butylbenzene	0.49	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	9-7-17	9-7-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Naphthalene	ND	1.0	EPA 8260C	9-7-17	9-7-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
<b>Surrogate:</b>	<b>Percent Recovery</b>	<b>Control Limits</b>				
Dibromofluoromethane	92	77-129				
Toluene-d8	95	80-127				
4-Bromofluorobenzene	99	78-125				



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 Project: 14697.00008

**VOLATILES EPA 8260C**  
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Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	KMW-06-8-31-17					
Laboratory ID:	09-019-03					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Chloromethane	ND	1.0	EPA 8260C	9-7-17	9-7-17	
Vinyl Chloride	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromomethane	ND	0.36	EPA 8260C	9-7-17	9-7-17	
Chloroethane	ND	1.0	EPA 8260C	9-7-17	9-7-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Acetone	ND	5.0	EPA 8260C	9-7-17	9-7-17	
Iodomethane	ND	2.0	EPA 8260C	9-7-17	9-7-17	
Carbon Disulfide	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Methylene Chloride	ND	2.0	EPA 8260C	9-7-17	9-7-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Vinyl Acetate	ND	1.0	EPA 8260C	9-7-17	9-7-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Butanone	ND	5.0	EPA 8260C	9-7-17	9-7-17	
Bromochloromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Chloroform	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Benzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Trichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Dibromomethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromodichloromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Chloroethyl Vinyl Ether	ND	1.3	EPA 8260C	9-7-17	9-7-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	9-7-17	9-7-17	
Toluene	ND	1.0	EPA 8260C	9-7-17	9-7-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-7-17	9-7-17	



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 Project: 14697.00008

**VOLATILES EPA 8260C**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-06-8-31-17</b>					
<b>Laboratory ID:</b>	09-019-03					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Tetrachloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Hexanone	ND	2.0	EPA 8260C	9-7-17	9-7-17	
Dibromochloromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Chlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Ethylbenzene	0.27	0.20	EPA 8260C	9-7-17	9-7-17	
m,p-Xylene	1.6	0.40	EPA 8260C	9-7-17	9-7-17	
o-Xylene	0.47	0.20	EPA 8260C	9-7-17	9-7-17	
Styrene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromoform	ND	1.0	EPA 8260C	9-7-17	9-7-17	
Isopropylbenzene	26	0.20	EPA 8260C	9-7-17	9-7-17	
Bromobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
n-Propylbenzene	37	0.20	EPA 8260C	9-7-17	9-7-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
tert-Butylbenzene	0.36	0.20	EPA 8260C	9-7-17	9-7-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
sec-Butylbenzene	3.7	0.20	EPA 8260C	9-7-17	9-7-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
n-Butylbenzene	2.3	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	9-7-17	9-7-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Naphthalene	ND	1.0	EPA 8260C	9-7-17	9-7-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
<b>Surrogate:</b>	<b>Percent Recovery</b>	<b>Control Limits</b>				
Dibromofluoromethane	100	77-129				
Toluene-d8	97	80-127				
4-Bromofluorobenzene	109	78-125				



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

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Date of Report: September 14, 2017  
 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

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

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-04-8-31-17</b>					
<b>Laboratory ID:</b>	09-019-04					
Dichlorodifluoromethane	ND	20	EPA 8260C	9-7-17	9-7-17	
Chloromethane	ND	100	EPA 8260C	9-7-17	9-7-17	
Vinyl Chloride	ND	20	EPA 8260C	9-7-17	9-7-17	
Bromomethane	ND	36	EPA 8260C	9-7-17	9-7-17	
Chloroethane	ND	100	EPA 8260C	9-7-17	9-7-17	
Trichlorofluoromethane	ND	20	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloroethene	ND	20	EPA 8260C	9-7-17	9-7-17	
Acetone	ND	500	EPA 8260C	9-7-17	9-7-17	
Iodomethane	ND	200	EPA 8260C	9-7-17	9-7-17	
Carbon Disulfide	ND	20	EPA 8260C	9-7-17	9-7-17	
Methylene Chloride	ND	200	EPA 8260C	9-7-17	9-7-17	
(trans) 1,2-Dichloroethene	ND	20	EPA 8260C	9-7-17	9-7-17	
Methyl t-Butyl Ether	ND	20	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloroethane	ND	20	EPA 8260C	9-7-17	9-7-17	
Vinyl Acetate	ND	100	EPA 8260C	9-7-17	9-7-17	
2,2-Dichloropropane	ND	20	EPA 8260C	9-7-17	9-7-17	
(cis) 1,2-Dichloroethene	ND	20	EPA 8260C	9-7-17	9-7-17	
2-Butanone	ND	500	EPA 8260C	9-7-17	9-7-17	
Bromochloromethane	ND	20	EPA 8260C	9-7-17	9-7-17	
Chloroform	ND	20	EPA 8260C	9-7-17	9-7-17	
1,1,1-Trichloroethane	ND	20	EPA 8260C	9-7-17	9-7-17	
Carbon Tetrachloride	ND	20	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloropropene	ND	20	EPA 8260C	9-7-17	9-7-17	
Benzene	ND	20	EPA 8260C	9-7-17	9-7-17	
1,2-Dichloroethane	ND	20	EPA 8260C	9-7-17	9-7-17	
Trichloroethene	ND	20	EPA 8260C	9-7-17	9-7-17	
1,2-Dichloropropane	ND	20	EPA 8260C	9-7-17	9-7-17	
Dibromomethane	ND	20	EPA 8260C	9-7-17	9-7-17	
Bromodichloromethane	ND	20	EPA 8260C	9-7-17	9-7-17	
2-Chloroethyl Vinyl Ether	ND	130	EPA 8260C	9-7-17	9-7-17	
(cis) 1,3-Dichloropropene	ND	20	EPA 8260C	9-7-17	9-7-17	
Methyl Isobutyl Ketone	ND	200	EPA 8260C	9-7-17	9-7-17	
Toluene	980	100	EPA 8260C	9-7-17	9-7-17	
(trans) 1,3-Dichloropropene	ND	20	EPA 8260C	9-7-17	9-7-17	



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 Samples Submitted: September 1, 2017  
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 Project: 14697.00008

**VOLATILES EPA 8260C**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-04-8-31-17</b>					
Laboratory ID:	09-019-04					
1,1,2-Trichloroethane	ND	20	EPA 8260C	9-7-17	9-7-17	
Tetrachloroethene	ND	20	EPA 8260C	9-7-17	9-7-17	
1,3-Dichloropropane	ND	20	EPA 8260C	9-7-17	9-7-17	
2-Hexanone	ND	200	EPA 8260C	9-7-17	9-7-17	
Dibromochloromethane	ND	20	EPA 8260C	9-7-17	9-7-17	
1,2-Dibromoethane	ND	20	EPA 8260C	9-7-17	9-7-17	
Chlorobenzene	ND	20	EPA 8260C	9-7-17	9-7-17	
1,1,1,2-Tetrachloroethane	ND	20	EPA 8260C	9-7-17	9-7-17	
Ethylbenzene	4300	20	EPA 8260C	9-7-17	9-7-17	
m,p-Xylene	7800	40	EPA 8260C	9-7-17	9-7-17	
o-Xylene	1900	20	EPA 8260C	9-7-17	9-7-17	
Styrene	ND	20	EPA 8260C	9-7-17	9-7-17	
Bromoform	ND	100	EPA 8260C	9-7-17	9-7-17	
Isopropylbenzene	63	20	EPA 8260C	9-7-17	9-7-17	
Bromobenzene	ND	20	EPA 8260C	9-7-17	9-7-17	
1,1,2,2-Tetrachloroethane	ND	20	EPA 8260C	9-7-17	9-7-17	
1,2,3-Trichloropropane	ND	20	EPA 8260C	9-7-17	9-7-17	
n-Propylbenzene	63	20	EPA 8260C	9-7-17	9-7-17	
2-Chlorotoluene	ND	20	EPA 8260C	9-7-17	9-7-17	
4-Chlorotoluene	ND	20	EPA 8260C	9-7-17	9-7-17	
1,3,5-Trimethylbenzene	30	20	EPA 8260C	9-7-17	9-7-17	
tert-Butylbenzene	ND	20	EPA 8260C	9-7-17	9-7-17	
1,2,4-Trimethylbenzene	97	20	EPA 8260C	9-7-17	9-7-17	
sec-Butylbenzene	ND	20	EPA 8260C	9-7-17	9-7-17	
1,3-Dichlorobenzene	ND	20	EPA 8260C	9-7-17	9-7-17	
p-Isopropyltoluene	ND	20	EPA 8260C	9-7-17	9-7-17	
1,4-Dichlorobenzene	ND	20	EPA 8260C	9-7-17	9-7-17	
1,2-Dichlorobenzene	ND	20	EPA 8260C	9-7-17	9-7-17	
n-Butylbenzene	ND	20	EPA 8260C	9-7-17	9-7-17	
1,2-Dibromo-3-chloropropane	ND	100	EPA 8260C	9-7-17	9-7-17	
1,2,4-Trichlorobenzene	ND	20	EPA 8260C	9-7-17	9-7-17	
Hexachlorobutadiene	ND	20	EPA 8260C	9-7-17	9-7-17	
Naphthalene	ND	100	EPA 8260C	9-7-17	9-7-17	
1,2,3-Trichlorobenzene	ND	20	EPA 8260C	9-7-17	9-7-17	
<b>Surrogate:</b>	<b>Percent Recovery</b>	<b>Control Limits</b>				
Dibromofluoromethane	97	77-129				
Toluene-d8	98	80-127				
4-Bromofluorobenzene	99	78-125				



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Date of Report: September 14, 2017  
 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

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

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-2R-9-1-17</b>					
<b>Laboratory ID:</b>	09-019-05					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Chloromethane	ND	1.0	EPA 8260C	9-7-17	9-7-17	
Vinyl Chloride	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromomethane	ND	0.36	EPA 8260C	9-7-17	9-7-17	
Chloroethane	ND	1.0	EPA 8260C	9-7-17	9-7-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Acetone	ND	5.0	EPA 8260C	9-7-17	9-7-17	
Iodomethane	ND	2.0	EPA 8260C	9-7-17	9-7-17	
Carbon Disulfide	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Methylene Chloride	ND	2.0	EPA 8260C	9-7-17	9-7-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Vinyl Acetate	ND	1.0	EPA 8260C	9-7-17	9-7-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Butanone	ND	5.0	EPA 8260C	9-7-17	9-7-17	
Bromochloromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Chloroform	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Benzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Trichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Dibromomethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromodichloromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Chloroethyl Vinyl Ether	ND	1.3	EPA 8260C	9-7-17	9-7-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	9-7-17	9-7-17	
Toluene	ND	1.0	EPA 8260C	9-7-17	9-7-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-7-17	9-7-17	



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 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

**VOLATILES EPA 8260C**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-2R-9-1-17</b>					
<b>Laboratory ID:</b>	09-019-05					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Tetrachloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Hexanone	ND	2.0	EPA 8260C	9-7-17	9-7-17	
Dibromochloromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Chlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Ethylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
m,p-Xylene	ND	0.40	EPA 8260C	9-7-17	9-7-17	
o-Xylene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Styrene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromoform	ND	1.0	EPA 8260C	9-7-17	9-7-17	
Isopropylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
n-Propylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
tert-Butylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
sec-Butylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
n-Butylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	9-7-17	9-7-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Naphthalene	ND	1.0	EPA 8260C	9-7-17	9-7-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
<b>Surrogate:</b>	<b>Percent Recovery</b>	<b>Control Limits</b>				
Dibromofluoromethane	96	77-129				
Toluene-d8	100	80-127				
4-Bromofluorobenzene	97	78-125				



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 Laboratory Reference: 1709-019  
 Project: 14697.00008

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

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-07-9-1-17</b>					
<b>Laboratory ID:</b>	09-019-06					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Chloromethane	ND	1.0	EPA 8260C	9-7-17	9-7-17	
Vinyl Chloride	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromomethane	ND	0.36	EPA 8260C	9-7-17	9-7-17	
Chloroethane	ND	1.0	EPA 8260C	9-7-17	9-7-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Acetone	ND	5.0	EPA 8260C	9-7-17	9-7-17	
Iodomethane	ND	2.0	EPA 8260C	9-7-17	9-7-17	
Carbon Disulfide	0.26	0.20	EPA 8260C	9-7-17	9-7-17	
Methylene Chloride	ND	2.0	EPA 8260C	9-7-17	9-7-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Vinyl Acetate	ND	1.0	EPA 8260C	9-7-17	9-7-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Butanone	ND	5.0	EPA 8260C	9-7-17	9-7-17	
Bromochloromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Chloroform	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Benzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Trichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Dibromomethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromodichloromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Chloroethyl Vinyl Ether	ND	1.3	EPA 8260C	9-7-17	9-7-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	9-7-17	9-7-17	
Toluene	ND	1.0	EPA 8260C	9-7-17	9-7-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-7-17	9-7-17	



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 Project: 14697.00008

**VOLATILES EPA 8260C**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-07-9-1-17</b>					
Laboratory ID:	09-019-06					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Tetrachloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Hexanone	ND	2.0	EPA 8260C	9-7-17	9-7-17	
Dibromochloromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Chlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Ethylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
m,p-Xylene	ND	0.40	EPA 8260C	9-7-17	9-7-17	
o-Xylene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Styrene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromoform	ND	1.0	EPA 8260C	9-7-17	9-7-17	
Isopropylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
n-Propylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
tert-Butylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
sec-Butylbenzene	0.29	0.20	EPA 8260C	9-7-17	9-7-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
n-Butylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	9-7-17	9-7-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Naphthalene	ND	1.0	EPA 8260C	9-7-17	9-7-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
<b>Surrogate:</b>	<b>Percent Recovery</b>	<b>Control Limits</b>				
Dibromofluoromethane	101	77-129				
Toluene-d8	103	80-127				
4-Bromofluorobenzene	103	78-125				



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Date of Report: September 14, 2017  
 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

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

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-08-9-1-17</b>					
<b>Laboratory ID:</b>	<b>09-019-07</b>					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Chloromethane	ND	1.0	EPA 8260C	9-7-17	9-7-17	
Vinyl Chloride	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromomethane	ND	0.36	EPA 8260C	9-7-17	9-7-17	
Chloroethane	ND	1.0	EPA 8260C	9-7-17	9-7-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Acetone	ND	5.0	EPA 8260C	9-7-17	9-7-17	
Iodomethane	ND	2.0	EPA 8260C	9-7-17	9-7-17	
Carbon Disulfide	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Methylene Chloride	ND	2.0	EPA 8260C	9-7-17	9-7-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Vinyl Acetate	ND	1.0	EPA 8260C	9-7-17	9-7-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Butanone	ND	5.0	EPA 8260C	9-7-17	9-7-17	
Bromochloromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Chloroform	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Benzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Trichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Dibromomethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromodichloromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Chloroethyl Vinyl Ether	ND	1.3	EPA 8260C	9-7-17	9-7-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	9-7-17	9-7-17	
Toluene	ND	1.0	EPA 8260C	9-7-17	9-7-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-7-17	9-7-17	



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 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

**VOLATILES EPA 8260C**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-08-9-1-17</b>					
Laboratory ID:	09-019-07					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Tetrachloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Hexanone	ND	2.0	EPA 8260C	9-7-17	9-7-17	
Dibromochloromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Chlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Ethylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
m,p-Xylene	ND	0.40	EPA 8260C	9-7-17	9-7-17	
o-Xylene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Styrene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromoform	ND	1.0	EPA 8260C	9-7-17	9-7-17	
Isopropylbenzene	0.79	0.20	EPA 8260C	9-7-17	9-7-17	
Bromobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
n-Propylbenzene	0.84	0.20	EPA 8260C	9-7-17	9-7-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
tert-Butylbenzene	0.24	0.20	EPA 8260C	9-7-17	9-7-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
sec-Butylbenzene	2.2	0.20	EPA 8260C	9-7-17	9-7-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
n-Butylbenzene	0.51	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	9-7-17	9-7-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Naphthalene	ND	1.0	EPA 8260C	9-7-17	9-7-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
<b>Surrogate:</b>	<b>Percent Recovery</b>	<b>Control Limits</b>				
Dibromofluoromethane	97	77-129				
Toluene-d8	102	80-127				
4-Bromofluorobenzene	100	78-125				



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 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

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

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-03R-9-1-17</b>					
<b>Laboratory ID:</b>	09-019-08					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Chloromethane	ND	1.0	EPA 8260C	9-7-17	9-7-17	
Vinyl Chloride	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromomethane	ND	0.36	EPA 8260C	9-7-17	9-7-17	
Chloroethane	ND	1.0	EPA 8260C	9-7-17	9-7-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Acetone	ND	5.0	EPA 8260C	9-7-17	9-7-17	
Iodomethane	ND	2.0	EPA 8260C	9-7-17	9-7-17	
Carbon Disulfide	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Methylene Chloride	ND	2.0	EPA 8260C	9-7-17	9-7-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Vinyl Acetate	ND	1.0	EPA 8260C	9-7-17	9-7-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Butanone	ND	5.0	EPA 8260C	9-7-17	9-7-17	
Bromochloromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Chloroform	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Benzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Trichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Dibromomethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromodichloromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Chloroethyl Vinyl Ether	ND	1.3	EPA 8260C	9-7-17	9-7-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	9-7-17	9-7-17	
Toluene	ND	1.0	EPA 8260C	9-7-17	9-7-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-7-17	9-7-17	



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 Project: 14697.00008

**VOLATILES EPA 8260C**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-03R-9-1-17</b>					
<b>Laboratory ID:</b>	09-019-08					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Tetrachloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Hexanone	ND	2.0	EPA 8260C	9-7-17	9-7-17	
Dibromochloromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Chlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Ethylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
m,p-Xylene	ND	0.40	EPA 8260C	9-7-17	9-7-17	
o-Xylene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Styrene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromoform	ND	1.0	EPA 8260C	9-7-17	9-7-17	
Isopropylbenzene	9.2	0.20	EPA 8260C	9-7-17	9-7-17	
Bromobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
n-Propylbenzene	17	0.20	EPA 8260C	9-7-17	9-7-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
tert-Butylbenzene	0.32	0.20	EPA 8260C	9-7-17	9-7-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
sec-Butylbenzene	2.2	0.20	EPA 8260C	9-7-17	9-7-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
n-Butylbenzene	0.67	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	9-7-17	9-7-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Naphthalene	ND	1.0	EPA 8260C	9-7-17	9-7-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
<b>Surrogate:</b>	<b>Percent Recovery</b>	<b>Control Limits</b>				
Dibromofluoromethane	89	77-129				
Toluene-d8	95	80-127				
4-Bromofluorobenzene	96	78-125				



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 Laboratory Reference: 1709-019  
 Project: 14697.00008

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

Matrix: Water

Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>TRIP BLANK</b>					
<b>Laboratory ID:</b>	09-019-09					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Chloromethane	ND	1.0	EPA 8260C	9-7-17	9-7-17	
Vinyl Chloride	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromomethane	ND	0.36	EPA 8260C	9-7-17	9-7-17	
Chloroethane	ND	1.0	EPA 8260C	9-7-17	9-7-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Acetone	ND	5.0	EPA 8260C	9-7-17	9-7-17	
Iodomethane	ND	2.0	EPA 8260C	9-7-17	9-7-17	
Carbon Disulfide	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Methylene Chloride	ND	2.0	EPA 8260C	9-7-17	9-7-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Vinyl Acetate	ND	1.0	EPA 8260C	9-7-17	9-7-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Butanone	ND	5.0	EPA 8260C	9-7-17	9-7-17	
Bromochloromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Chloroform	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Benzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Trichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Dibromomethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromodichloromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Chloroethyl Vinyl Ether	ND	1.3	EPA 8260C	9-7-17	9-7-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	9-7-17	9-7-17	
Toluene	ND	1.0	EPA 8260C	9-7-17	9-7-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-7-17	9-7-17	



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

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Date of Report: September 14, 2017  
 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

**VOLATILES EPA 8260C**  
 Page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>TRIP BLANK</b>					
Laboratory ID:	09-019-09					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Tetrachloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Hexanone	ND	2.0	EPA 8260C	9-7-17	9-7-17	
Dibromochloromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Chlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Ethylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
m,p-Xylene	ND	0.40	EPA 8260C	9-7-17	9-7-17	
o-Xylene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Styrene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromoform	ND	1.0	EPA 8260C	9-7-17	9-7-17	
Isopropylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
n-Propylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
tert-Butylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
sec-Butylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
n-Butylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	9-7-17	9-7-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Naphthalene	ND	1.0	EPA 8260C	9-7-17	9-7-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
<b>Surrogate:</b>	<b>Percent Recovery</b>	<b>Control Limits</b>				
Dibromofluoromethane	98	77-129				
Toluene-d8	99	80-127				
4-Bromofluorobenzene	96	78-125				



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Date of Report: September 14, 2017  
 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

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

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0907W1					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Chloromethane	ND	1.0	EPA 8260C	9-7-17	9-7-17	
Vinyl Chloride	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromomethane	ND	0.36	EPA 8260C	9-7-17	9-7-17	
Chloroethane	ND	1.0	EPA 8260C	9-7-17	9-7-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Acetone	ND	5.0	EPA 8260C	9-7-17	9-7-17	
Iodomethane	ND	2.0	EPA 8260C	9-7-17	9-7-17	
Carbon Disulfide	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Methylene Chloride	ND	2.0	EPA 8260C	9-7-17	9-7-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Vinyl Acetate	ND	1.0	EPA 8260C	9-7-17	9-7-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Butanone	ND	5.0	EPA 8260C	9-7-17	9-7-17	
Bromochloromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Chloroform	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Benzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Trichloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Dibromomethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromodichloromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Chloroethyl Vinyl Ether	ND	1.3	EPA 8260C	9-7-17	9-7-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	9-7-17	9-7-17	
Toluene	ND	1.0	EPA 8260C	9-7-17	9-7-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-7-17	9-7-17	



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 Project: 14697.00008

**VOLATILES EPA 8260C**  
**METHOD BLANK QUALITY CONTROL**  
 Page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0907W1					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Tetrachloroethene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Hexanone	ND	2.0	EPA 8260C	9-7-17	9-7-17	
Dibromochloromethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Chlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Ethylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
m,p-Xylene	ND	0.40	EPA 8260C	9-7-17	9-7-17	
o-Xylene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Styrene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromoform	ND	1.0	EPA 8260C	9-7-17	9-7-17	
Isopropylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Bromobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	9-7-17	9-7-17	
n-Propylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
tert-Butylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
sec-Butylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
n-Butylbenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	9-7-17	9-7-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Naphthalene	ND	1.0	EPA 8260C	9-7-17	9-7-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	9-7-17	9-7-17	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	99	77-129				
Toluene-d8	102	80-127				
4-Bromofluorobenzene	97	78-125				



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 Project: 14697.00008

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

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0908W1					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	9-8-17	9-8-17	
Chloromethane	ND	1.0	EPA 8260C	9-8-17	9-8-17	
Vinyl Chloride	ND	0.20	EPA 8260C	9-8-17	9-8-17	
Bromomethane	ND	0.20	EPA 8260C	9-8-17	9-8-17	
Chloroethane	ND	1.0	EPA 8260C	9-8-17	9-8-17	
Trichlorofluoromethane	ND	0.20	EPA 8260C	9-8-17	9-8-17	
1,1-Dichloroethene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
Acetone	ND	5.0	EPA 8260C	9-8-17	9-8-17	
Iodomethane	ND	1.3	EPA 8260C	9-8-17	9-8-17	
Carbon Disulfide	ND	0.20	EPA 8260C	9-8-17	9-8-17	
Methylene Chloride	ND	2.0	EPA 8260C	9-8-17	9-8-17	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	9-8-17	9-8-17	
1,1-Dichloroethane	ND	0.20	EPA 8260C	9-8-17	9-8-17	
Vinyl Acetate	ND	1.0	EPA 8260C	9-8-17	9-8-17	
2,2-Dichloropropane	ND	0.20	EPA 8260C	9-8-17	9-8-17	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
2-Butanone	ND	5.0	EPA 8260C	9-8-17	9-8-17	
Bromochloromethane	ND	0.20	EPA 8260C	9-8-17	9-8-17	
Chloroform	ND	0.20	EPA 8260C	9-8-17	9-8-17	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	9-8-17	9-8-17	
Carbon Tetrachloride	ND	0.20	EPA 8260C	9-8-17	9-8-17	
1,1-Dichloropropene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
Benzene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
1,2-Dichloroethane	ND	0.20	EPA 8260C	9-8-17	9-8-17	
Trichloroethene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
1,2-Dichloropropane	ND	0.20	EPA 8260C	9-8-17	9-8-17	
Dibromomethane	ND	0.20	EPA 8260C	9-8-17	9-8-17	
Bromodichloromethane	ND	0.20	EPA 8260C	9-8-17	9-8-17	
2-Chloroethyl Vinyl Ether	ND	1.3	EPA 8260C	9-8-17	9-8-17	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	9-8-17	9-8-17	
Toluene	ND	1.0	EPA 8260C	9-8-17	9-8-17	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	9-8-17	9-8-17	



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 Project: 14697.00008

**VOLATILES EPA 8260C**  
**METHOD BLANK QUALITY CONTROL**  
 Page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0908W1					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	9-8-17	9-8-17	
Tetrachloroethene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
1,3-Dichloropropane	ND	0.20	EPA 8260C	9-8-17	9-8-17	
2-Hexanone	ND	2.0	EPA 8260C	9-8-17	9-8-17	
Dibromochloromethane	ND	0.20	EPA 8260C	9-8-17	9-8-17	
1,2-Dibromoethane	ND	0.20	EPA 8260C	9-8-17	9-8-17	
Chlorobenzene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-8-17	9-8-17	
Ethylbenzene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
m,p-Xylene	ND	0.40	EPA 8260C	9-8-17	9-8-17	
o-Xylene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
Styrene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
Bromoform	ND	1.0	EPA 8260C	9-8-17	9-8-17	
Isopropylbenzene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
Bromobenzene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	9-8-17	9-8-17	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	9-8-17	9-8-17	
n-Propylbenzene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
2-Chlorotoluene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
4-Chlorotoluene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
tert-Butylbenzene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
sec-Butylbenzene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
p-Isopropyltoluene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
n-Butylbenzene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	9-8-17	9-8-17	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
Hexachlorobutadiene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
Naphthalene	ND	1.0	EPA 8260C	9-8-17	9-8-17	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	9-8-17	9-8-17	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	104	77-129				
Toluene-d8	101	80-127				
4-Bromofluorobenzene	101	78-125				



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 Project: 14697.00008

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

Matrix: Water

Units: ug/L

Analyte	Result	Spike Level		Percent Recovery		Recovery Limits	RPD RPD	RPD Limit	Flags							
		Recovery	Limits													
<b>SPIKE BLANKS</b>																
Laboratory ID: SB0908W1																
		SB	SBD	SB	SBD	SB	SBD									
1,1-Dichloroethene	<b>8.95</b>	<b>8.71</b>	10.0	10.0	90	87	63-127	3	17							
Benzene	<b>9.85</b>	<b>9.67</b>	10.0	10.0	99	97	76-121	2	12							
Trichloroethene	<b>8.93</b>	<b>8.95</b>	10.0	10.0	89	90	64-120	0	15							
Toluene	<b>9.67</b>	<b>9.60</b>	10.0	10.0	97	96	82-120	1	13							
Chlorobenzene	<b>9.77</b>	<b>9.67</b>	10.0	10.0	98	97	80-120	1	14							
<i>Surrogate:</i>																
<i>Dibromofluoromethane</i>					100	104	77-129									
<i>Toluene-d8</i>					100	101	80-127									
<i>4-Bromofluorobenzene</i>					97	100	78-125									



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**VOLATILES EPA 8260C**  
**MS/MSD QUALITY CONTROL**

Matrix: Water

Units: ug/L

Analyte	Result	Spike Level	Source	Percent	Recovery	RPD	RPD	Limit	Flags
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**MATRIX SPIKES**

Laboratory ID: 09-019-05

	MS	MSD	MS	MSD	MS	MSD				
1,1-Dichloroethene	<b>8.98</b>	<b>9.66</b>	10.0	10.0	ND	90	97	65-119	7	15
Benzene	<b>9.35</b>	<b>10.3</b>	10.0	10.0	ND	94	103	75-117	10	15
Trichloroethene	<b>9.33</b>	<b>9.89</b>	10.0	10.0	ND	93	99	66-120	6	15
Toluene	<b>9.48</b>	<b>10.2</b>	10.0	10.0	ND	95	102	79-120	7	15
Chlorobenzene	<b>9.79</b>	<b>10.6</b>	10.0	10.0	ND	98	106	76-120	8	15

*Surrogate:*

*Dibromofluoromethane* 99 104 77-129

*Toluene-d8* 98 101 80-127

*4-Bromofluorobenzene* 96 98 78-125



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

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Date of Report: September 14, 2017  
 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

### PAHs EPA 8270D/SIM

Matrix: Water

Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-10-8-31-17</b>					
<b>Laboratory ID:</b>	09-019-01					
Naphthalene	<b>0.42</b>	0.094	EPA 8270D/SIM	9-5-17	9-6-17	
2-Methylnaphthalene	<b>ND</b>	0.094	EPA 8270D/SIM	9-5-17	9-6-17	
1-Methylnaphthalene	<b>ND</b>	0.094	EPA 8270D/SIM	9-5-17	9-6-17	
Acenaphthylene	<b>ND</b>	0.094	EPA 8270D/SIM	9-5-17	9-6-17	
Acenaphthene	<b>ND</b>	0.094	EPA 8270D/SIM	9-5-17	9-6-17	
Fluorene	<b>ND</b>	0.094	EPA 8270D/SIM	9-5-17	9-6-17	
Phenanthrene	<b>ND</b>	0.094	EPA 8270D/SIM	9-5-17	9-6-17	
Anthracene	<b>ND</b>	0.094	EPA 8270D/SIM	9-5-17	9-6-17	
Fluoranthene	<b>ND</b>	0.094	EPA 8270D/SIM	9-5-17	9-6-17	
Pyrene	<b>ND</b>	0.094	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[a]anthracene	<b>ND</b>	0.0094	EPA 8270D/SIM	9-5-17	9-6-17	
Chrysene	<b>ND</b>	0.0094	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[b]fluoranthene	<b>ND</b>	0.0094	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo(j,k)fluoranthene	<b>ND</b>	0.0094	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[a]pyrene	<b>ND</b>	0.0094	EPA 8270D/SIM	9-5-17	9-6-17	
Indeno(1,2,3-c,d)pyrene	<b>ND</b>	0.0094	EPA 8270D/SIM	9-5-17	9-6-17	
Dibenz[a,h]anthracene	<b>ND</b>	0.0094	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[g,h,i]perylene	<b>ND</b>	0.0094	EPA 8270D/SIM	9-5-17	9-6-17	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorobiphenyl	54	30 - 124				
Pyrene-d10	50	40 - 143				
Terphenyl-d14	75	27 - 127				



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Date of Report: September 14, 2017  
 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

### PAHs EPA 8270D/SIM

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-09-8-31-17</b>					
<b>Laboratory ID:</b>	09-019-02					
Naphthalene	ND	0.098	EPA 8270D/SIM	9-5-17	9-6-17	
2-Methylnaphthalene	ND	0.098	EPA 8270D/SIM	9-5-17	9-6-17	
1-Methylnaphthalene	0.32	0.098	EPA 8270D/SIM	9-5-17	9-6-17	
Acenaphthylene	ND	0.098	EPA 8270D/SIM	9-5-17	9-6-17	
Acenaphthene	1.0	0.098	EPA 8270D/SIM	9-5-17	9-6-17	
Fluorene	0.19	0.098	EPA 8270D/SIM	9-5-17	9-6-17	
Phenanthrene	ND	0.098	EPA 8270D/SIM	9-5-17	9-6-17	
Anthracene	0.16	0.098	EPA 8270D/SIM	9-5-17	9-6-17	
Fluoranthene	ND	0.098	EPA 8270D/SIM	9-5-17	9-6-17	
Pyrene	ND	0.098	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[a]anthracene	ND	0.0098	EPA 8270D/SIM	9-5-17	9-6-17	
Chrysene	ND	0.0098	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[b]fluoranthene	ND	0.0098	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo(j,k)fluoranthene	ND	0.0098	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[a]pyrene	ND	0.0098	EPA 8270D/SIM	9-5-17	9-6-17	
Indeno(1,2,3-c,d)pyrene	ND	0.0098	EPA 8270D/SIM	9-5-17	9-6-17	
Dibenz[a,h]anthracene	ND	0.0098	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[g,h,i]perylene	ND	0.0098	EPA 8270D/SIM	9-5-17	9-6-17	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
2-Fluorobiphenyl	33		30 - 124			
Pyrene-d10	32		40 - 143			Q
Terphenyl-d14	46		27 - 127			



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Date of Report: September 14, 2017  
 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

### PAHs EPA 8270D/SIM

Matrix: Water

Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-06-8-31-17</b>					
<b>Laboratory ID:</b>	09-019-03					
Naphthalene	<b>0.12</b>	0.095	EPA 8270D/SIM	9-5-17	9-6-17	
2-Methylnaphthalene	<b>0.79</b>	0.095	EPA 8270D/SIM	9-5-17	9-6-17	
1-Methylnaphthalene	<b>3.1</b>	0.095	EPA 8270D/SIM	9-5-17	9-6-17	
Acenaphthylene	<b>0.25</b>	0.095	EPA 8270D/SIM	9-5-17	9-6-17	
Acenaphthene	<b>1.9</b>	0.095	EPA 8270D/SIM	9-5-17	9-6-17	
Fluorene	<b>0.75</b>	0.095	EPA 8270D/SIM	9-5-17	9-6-17	
Phenanthrene	<b>0.12</b>	0.095	EPA 8270D/SIM	9-5-17	9-6-17	
Anthracene	<b>1.0</b>	0.095	EPA 8270D/SIM	9-5-17	9-6-17	
Fluoranthene	<b>0.10</b>	0.095	EPA 8270D/SIM	9-5-17	9-6-17	
Pyrene	<b>0.10</b>	0.095	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[a]anthracene	<b>0.041</b>	0.0095	EPA 8270D/SIM	9-5-17	9-6-17	
Chrysene	<b>0.032</b>	0.0095	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[b]fluoranthene	<b>0.034</b>	0.0095	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo(j,k)fluoranthene	<b>0.018</b>	0.0095	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[a]pyrene	<b>0.033</b>	0.0095	EPA 8270D/SIM	9-5-17	9-6-17	
Indeno(1,2,3-c,d)pyrene	<b>0.023</b>	0.0095	EPA 8270D/SIM	9-5-17	9-6-17	
Dibenz[a,h]anthracene	<b>ND</b>	0.0095	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[g,h,i]perylene	<b>0.025</b>	0.0095	EPA 8270D/SIM	9-5-17	9-6-17	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
2-Fluorobiphenyl	57		30 - 124			
Pyrene-d10	53		40 - 143			
Terphenyl-d14	77		27 - 127			



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Date of Report: September 14, 2017  
 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

### PAHs EPA 8270D/SIM

Matrix: Water

Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-04-8-31-17</b>					
<b>Laboratory ID:</b>	09-019-04					
Naphthalene	<b>2.9</b>	0.096	EPA 8270D/SIM	9-5-17	9-6-17	
2-Methylnaphthalene	<b>0.40</b>	0.096	EPA 8270D/SIM	9-5-17	9-6-17	
1-Methylnaphthalene	<b>0.33</b>	0.096	EPA 8270D/SIM	9-5-17	9-6-17	
Acenaphthylene	<b>ND</b>	0.096	EPA 8270D/SIM	9-5-17	9-6-17	
Acenaphthene	<b>0.35</b>	0.096	EPA 8270D/SIM	9-5-17	9-6-17	
Fluorene	<b>ND</b>	0.096	EPA 8270D/SIM	9-5-17	9-6-17	
Phenanthrene	<b>ND</b>	0.096	EPA 8270D/SIM	9-5-17	9-6-17	
Anthracene	<b>ND</b>	0.096	EPA 8270D/SIM	9-5-17	9-6-17	
Fluoranthene	<b>ND</b>	0.096	EPA 8270D/SIM	9-5-17	9-6-17	
Pyrene	<b>ND</b>	0.096	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[a]anthracene	<b>ND</b>	0.0096	EPA 8270D/SIM	9-5-17	9-6-17	
Chrysene	<b>0.010</b>	0.0096	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[b]fluoranthene	<b>0.012</b>	0.0096	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo(j,k)fluoranthene	<b>ND</b>	0.0096	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[a]pyrene	<b>ND</b>	0.0096	EPA 8270D/SIM	9-5-17	9-6-17	
Indeno(1,2,3-c,d)pyrene	<b>ND</b>	0.0096	EPA 8270D/SIM	9-5-17	9-6-17	
Dibenz[a,h]anthracene	<b>ND</b>	0.0096	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[g,h,i]perylene	<b>ND</b>	0.0096	EPA 8270D/SIM	9-5-17	9-6-17	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
2-Fluorobiphenyl	50		30 - 124			
Pyrene-d10	42		40 - 143			
Terphenyl-d14	75		27 - 127			



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 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

### PAHs EPA 8270D/SIM

Matrix: Water

Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-2R-9-1-17</b>					
<b>Laboratory ID:</b>	09-019-05					
Naphthalene	ND	0.099	EPA 8270D/SIM	9-5-17	9-6-17	
2-Methylnaphthalene	ND	0.099	EPA 8270D/SIM	9-5-17	9-6-17	
1-Methylnaphthalene	ND	0.099	EPA 8270D/SIM	9-5-17	9-6-17	
Acenaphthylene	ND	0.099	EPA 8270D/SIM	9-5-17	9-6-17	
Acenaphthene	ND	0.099	EPA 8270D/SIM	9-5-17	9-6-17	
Fluorene	ND	0.099	EPA 8270D/SIM	9-5-17	9-6-17	
Phenanthrene	ND	0.099	EPA 8270D/SIM	9-5-17	9-6-17	
Anthracene	ND	0.099	EPA 8270D/SIM	9-5-17	9-6-17	
Fluoranthene	ND	0.099	EPA 8270D/SIM	9-5-17	9-6-17	
Pyrene	ND	0.099	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[a]anthracene	ND	0.0099	EPA 8270D/SIM	9-5-17	9-6-17	
Chrysene	ND	0.0099	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[b]fluoranthene	ND	0.0099	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo(j,k)fluoranthene	ND	0.0099	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[a]pyrene	ND	0.0099	EPA 8270D/SIM	9-5-17	9-6-17	
Indeno(1,2,3-c,d)pyrene	ND	0.0099	EPA 8270D/SIM	9-5-17	9-6-17	
Dibenz[a,h]anthracene	ND	0.0099	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[g,h,i]perylene	ND	0.0099	EPA 8270D/SIM	9-5-17	9-6-17	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
2-Fluorobiphenyl	77		30 - 124			
Pyrene-d10	64		40 - 143			
Terphenyl-d14	80		27 - 127			



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 Laboratory Reference: 1709-019  
 Project: 14697.00008

### PAHs EPA 8270D/SIM

Matrix: Water

Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-07-9-1-17</b>					
<b>Laboratory ID:</b>	09-019-06					
Naphthalene	ND	0.099	EPA 8270D/SIM	9-5-17	9-6-17	
2-Methylnaphthalene	ND	0.099	EPA 8270D/SIM	9-5-17	9-6-17	
1-Methylnaphthalene	ND	0.099	EPA 8270D/SIM	9-5-17	9-6-17	
Acenaphthylene	ND	0.099	EPA 8270D/SIM	9-5-17	9-6-17	
Acenaphthene	ND	0.099	EPA 8270D/SIM	9-5-17	9-6-17	
Fluorene	ND	0.099	EPA 8270D/SIM	9-5-17	9-6-17	
Phenanthrene	ND	0.099	EPA 8270D/SIM	9-5-17	9-6-17	
Anthracene	ND	0.099	EPA 8270D/SIM	9-5-17	9-6-17	
Fluoranthene	ND	0.099	EPA 8270D/SIM	9-5-17	9-6-17	
Pyrene	ND	0.099	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[a]anthracene	ND	0.0099	EPA 8270D/SIM	9-5-17	9-6-17	
Chrysene	ND	0.0099	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[b]fluoranthene	ND	0.0099	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo(j,k)fluoranthene	ND	0.0099	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[a]pyrene	ND	0.0099	EPA 8270D/SIM	9-5-17	9-6-17	
Indeno(1,2,3-c,d)pyrene	ND	0.0099	EPA 8270D/SIM	9-5-17	9-6-17	
Dibenz[a,h]anthracene	ND	0.0099	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[g,h,i]perylene	ND	0.0099	EPA 8270D/SIM	9-5-17	9-6-17	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
2-Fluorobiphenyl	48		30 - 124			
Pyrene-d10	55		40 - 143			
Terphenyl-d14	69		27 - 127			



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 Laboratory Reference: 1709-019  
 Project: 14697.00008

### PAHs EPA 8270D/SIM

Matrix: Water

Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-08-9-1-17</b>					
<b>Laboratory ID:</b>	09-019-07					
Naphthalene	ND	0.10	EPA 8270D/SIM	9-8-17	9-8-17	
2-Methylnaphthalene	ND	0.10	EPA 8270D/SIM	9-8-17	9-8-17	
1-Methylnaphthalene	ND	0.10	EPA 8270D/SIM	9-8-17	9-8-17	
Acenaphthylene	ND	0.10	EPA 8270D/SIM	9-8-17	9-8-17	
Acenaphthene	0.18	0.10	EPA 8270D/SIM	9-8-17	9-8-17	
Fluorene	0.29	0.10	EPA 8270D/SIM	9-8-17	9-8-17	
Phenanthrene	ND	0.10	EPA 8270D/SIM	9-8-17	9-8-17	
Anthracene	ND	0.10	EPA 8270D/SIM	9-8-17	9-8-17	
Fluoranthene	0.11	0.10	EPA 8270D/SIM	9-8-17	9-8-17	
Pyrene	ND	0.10	EPA 8270D/SIM	9-8-17	9-8-17	
Benzo[a]anthracene	0.012	0.010	EPA 8270D/SIM	9-8-17	9-8-17	
Chrysene	0.012	0.010	EPA 8270D/SIM	9-8-17	9-8-17	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	9-8-17	9-8-17	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	9-8-17	9-8-17	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	9-8-17	9-8-17	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	9-8-17	9-8-17	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	9-8-17	9-8-17	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270D/SIM	9-8-17	9-8-17	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
2-Fluorobiphenyl	43		30 - 124			
Pyrene-d10	42		40 - 143			
Terphenyl-d14	47		27 - 127			



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 Project: 14697.00008

### PAHs EPA 8270D/SIM

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-03R-9-1-17</b>					
<b>Laboratory ID:</b>	09-019-08					
Naphthalene	ND	0.098	EPA 8270D/SIM	9-8-17	9-8-17	
2-Methylnaphthalene	ND	0.098	EPA 8270D/SIM	9-8-17	9-8-17	
1-Methylnaphthalene	ND	0.098	EPA 8270D/SIM	9-8-17	9-8-17	
Acenaphthylene	ND	0.098	EPA 8270D/SIM	9-8-17	9-8-17	
Acenaphthene	0.22	0.098	EPA 8270D/SIM	9-8-17	9-8-17	
Fluorene	0.29	0.098	EPA 8270D/SIM	9-8-17	9-8-17	
Phenanthrene	0.16	0.098	EPA 8270D/SIM	9-8-17	9-8-17	
Anthracene	ND	0.098	EPA 8270D/SIM	9-8-17	9-8-17	
Fluoranthene	ND	0.098	EPA 8270D/SIM	9-8-17	9-8-17	
Pyrene	ND	0.098	EPA 8270D/SIM	9-8-17	9-8-17	
Benzo[a]anthracene	ND	0.0098	EPA 8270D/SIM	9-8-17	9-8-17	
Chrysene	ND	0.0098	EPA 8270D/SIM	9-8-17	9-8-17	
Benzo[b]fluoranthene	ND	0.0098	EPA 8270D/SIM	9-8-17	9-8-17	
Benzo(j,k)fluoranthene	ND	0.0098	EPA 8270D/SIM	9-8-17	9-8-17	
Benzo[a]pyrene	ND	0.0098	EPA 8270D/SIM	9-8-17	9-8-17	
Indeno(1,2,3-c,d)pyrene	ND	0.0098	EPA 8270D/SIM	9-8-17	9-8-17	
Dibenz[a,h]anthracene	ND	0.0098	EPA 8270D/SIM	9-8-17	9-8-17	
Benzo[g,h,i]perylene	ND	0.0098	EPA 8270D/SIM	9-8-17	9-8-17	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
2-Fluorobiphenyl	76		30 - 124			
Pyrene-d10	89		40 - 143			
Terphenyl-d14	112		27 - 127			



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 Project: 14697.00008

**PAHs EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**

Matrix: Water

Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0905W1					
Naphthalene	ND	0.10	EPA 8270D/SIM	9-5-17	9-6-17	
2-Methylnaphthalene	ND	0.10	EPA 8270D/SIM	9-5-17	9-6-17	
1-Methylnaphthalene	ND	0.10	EPA 8270D/SIM	9-5-17	9-6-17	
Acenaphthylene	ND	0.10	EPA 8270D/SIM	9-5-17	9-6-17	
Acenaphthene	ND	0.10	EPA 8270D/SIM	9-5-17	9-6-17	
Fluorene	ND	0.10	EPA 8270D/SIM	9-5-17	9-6-17	
Phenanthrene	ND	0.10	EPA 8270D/SIM	9-5-17	9-6-17	
Anthracene	ND	0.10	EPA 8270D/SIM	9-5-17	9-6-17	
Fluoranthene	ND	0.10	EPA 8270D/SIM	9-5-17	9-6-17	
Pyrene	ND	0.10	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	9-5-17	9-6-17	
Chrysene	ND	0.010	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	9-5-17	9-6-17	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	9-5-17	9-6-17	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	9-5-17	9-6-17	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270D/SIM	9-5-17	9-6-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	65	30 - 124				
Pyrene-d10	64	40 - 143				
Terphenyl-d14	78	27 - 127				



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Date of Report: September 14, 2017  
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 Laboratory Reference: 1709-019  
 Project: 14697.00008

**PAHs EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**

Matrix: Water

Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0908W1					
Naphthalene	ND	0.10	EPA 8270D/SIM	9-8-17	9-8-17	
2-Methylnaphthalene	ND	0.10	EPA 8270D/SIM	9-8-17	9-8-17	
1-Methylnaphthalene	ND	0.10	EPA 8270D/SIM	9-8-17	9-8-17	
Acenaphthylene	ND	0.10	EPA 8270D/SIM	9-8-17	9-8-17	
Acenaphthene	ND	0.10	EPA 8270D/SIM	9-8-17	9-8-17	
Fluorene	ND	0.10	EPA 8270D/SIM	9-8-17	9-8-17	
Phenanthrene	ND	0.10	EPA 8270D/SIM	9-8-17	9-8-17	
Anthracene	ND	0.10	EPA 8270D/SIM	9-8-17	9-8-17	
Fluoranthene	ND	0.10	EPA 8270D/SIM	9-8-17	9-8-17	
Pyrene	ND	0.10	EPA 8270D/SIM	9-8-17	9-8-17	
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	9-8-17	9-8-17	
Chrysene	ND	0.010	EPA 8270D/SIM	9-8-17	9-8-17	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	9-8-17	9-8-17	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	9-8-17	9-8-17	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	9-8-17	9-8-17	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	9-8-17	9-8-17	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	9-8-17	9-8-17	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270D/SIM	9-8-17	9-8-17	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	76	30 - 124				
Pyrene-d10	93	40 - 143				
Terphenyl-d14	104	27 - 127				



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 Project: 14697.00008

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

Matrix: Water

Units: ug/L

Analyte	Result		Spike Level		Source	Percent	Recovery	RPD		
	Result	Recovery	Spike Level	Result	Recovery	Limits	RPD	Limit	Flags	
<b>MATRIX SPIKES</b>										
Laboratory ID:	09-019-05									
	MS	MSD	MS	MSD	MS	MSD				
Naphthalene	<b>0.272</b>	<b>0.313</b>	0.490	0.496	ND	56	63	29 - 101	14	47
Acenaphthylene	<b>0.318</b>	<b>0.341</b>	0.490	0.496	ND	65	69	20 - 117	7	50
Acenaphthene	<b>0.327</b>	<b>0.350</b>	0.490	0.496	ND	67	71	37 - 109	7	43
Fluorene	<b>0.345</b>	<b>0.372</b>	0.490	0.496	ND	70	75	47 - 108	8	34
Phenanthrene	<b>0.355</b>	<b>0.368</b>	0.490	0.496	ND	72	74	49 - 109	4	28
Anthracene	<b>0.425</b>	<b>0.463</b>	0.490	0.496	ND	87	93	34 - 140	9	32
Fluoranthene	<b>0.382</b>	<b>0.409</b>	0.490	0.496	ND	78	82	45 - 120	7	39
Pyrene	<b>0.382</b>	<b>0.405</b>	0.490	0.496	ND	78	82	42 - 133	6	39
Benzo[a]anthracene	<b>0.391</b>	<b>0.414</b>	0.490	0.496	ND	80	83	71 - 117	6	28
Chrysene	<b>0.414</b>	<b>0.445</b>	0.490	0.496	ND	84	90	53 - 110	7	25
Benzo[b]fluoranthene	<b>0.343</b>	<b>0.350</b>	0.490	0.496	ND	70	71	53 - 123	2	37
Benzo(j,k)fluoranthene	<b>0.456</b>	<b>0.498</b>	0.490	0.496	ND	93	100	52 - 119	9	41
Benzo[a]pyrene	<b>0.344</b>	<b>0.367</b>	0.490	0.496	ND	70	74	37 - 129	6	33
Indeno(1,2,3-c,d)pyrene	<b>0.365</b>	<b>0.374</b>	0.490	0.496	ND	74	75	45 - 128	2	31
Dibenz[a,h]anthracene	<b>0.378</b>	<b>0.411</b>	0.490	0.496	ND	77	83	54 - 120	8	30
Benzo[g,h,i]perylene	<b>0.386</b>	<b>0.402</b>	0.490	0.496	ND	79	81	49 - 117	4	29
<i>Surrogate:</i>										
<i>2-Fluorobiphenyl</i>						67	74	30 - 124		
<i>Pyrene-d10</i>						77	82	40 - 143		
<i>Terphenyl-d14</i>						86	91	27 - 127		



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 Project: 14697.00008

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

Matrix: Water

Units: ug/L

Analyte	Result		Spike Level		Percent Recovery		Recovery Limits	RPD	RPD Limit	Flags				
<b>SPIKE BLANKS</b>														
Laboratory ID:	SB0905W1													
	SB	SBD	SB	SBD	SB	SBD								
Naphthalene	<b>0.267</b>	<b>0.273</b>	0.500	0.500	53	55	29 - 101	2	47					
Acenaphthylene	<b>0.286</b>	<b>0.293</b>	0.500	0.500	57	59	20 - 117	2	50					
Acenaphthene	<b>0.313</b>	<b>0.297</b>	0.500	0.500	63	59	37 - 109	5	43					
Fluorene	<b>0.310</b>	<b>0.322</b>	0.500	0.500	62	64	47 - 108	4	34					
Phenanthrene	<b>0.324</b>	<b>0.325</b>	0.500	0.500	65	65	49 - 109	0	28					
Anthracene	<b>0.345</b>	<b>0.344</b>	0.500	0.500	69	69	34 - 140	0	32					
Fluoranthene	<b>0.345</b>	<b>0.344</b>	0.500	0.500	69	69	45 - 120	0	39					
Pyrene	<b>0.396</b>	<b>0.356</b>	0.500	0.500	79	71	42 - 133	11	39					
Benzo[a]anthracene	<b>0.373</b>	<b>0.374</b>	0.500	0.500	75	75	71 - 117	0	28					
Chrysene	<b>0.323</b>	<b>0.335</b>	0.500	0.500	65	67	53 - 110	4	25					
Benzo[b]fluoranthene	<b>0.341</b>	<b>0.310</b>	0.500	0.500	68	62	53 - 123	10	37					
Benzo(j,k)fluoranthene	<b>0.339</b>	<b>0.320</b>	0.500	0.500	68	64	52 - 119	6	41					
Benzo[a]pyrene	<b>0.335</b>	<b>0.341</b>	0.500	0.500	67	68	37 - 129	2	33					
Indeno(1,2,3-c,d)pyrene	<b>0.346</b>	<b>0.305</b>	0.500	0.500	69	61	45 - 128	13	31					
Dibenz[a,h]anthracene	<b>0.345</b>	<b>0.308</b>	0.500	0.500	69	62	54 - 120	11	30					
Benzo[g,h,i]perylene	<b>0.338</b>	<b>0.296</b>	0.500	0.500	68	59	49 - 117	13	29					
<i>Surrogate:</i>														
<i>2-Fluorobiphenyl</i>					63	60	30 - 124							
<i>Pyrene-d10</i>					65	65	40 - 143							
<i>Terphenyl-d14</i>					84	77	27 - 127							



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 Laboratory Reference: 1709-019  
 Project: 14697.00008

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

Matrix: Water

Units: ug/L

Analyte	Result		Spike Level		Percent Recovery		Recovery Limits	RPD	RPD Limit	Flags				
<b>SPIKE BLANKS</b>														
Laboratory ID:	SB0908W1													
	SB	SBD	SB	SBD	SB	SBD								
Naphthalene	<b>0.368</b>	<b>0.366</b>	0.500	0.500	74	73	29 - 101	1	47					
Acenaphthylene	<b>0.434</b>	<b>0.433</b>	0.500	0.500	87	87	20 - 117	0	50					
Acenaphthene	<b>0.411</b>	<b>0.416</b>	0.500	0.500	82	83	37 - 109	1	43					
Fluorene	<b>0.455</b>	<b>0.445</b>	0.500	0.500	91	89	47 - 108	2	34					
Phenanthrene	<b>0.430</b>	<b>0.422</b>	0.500	0.500	86	84	49 - 109	2	28					
Anthracene	<b>0.454</b>	<b>0.514</b>	0.500	0.500	91	103	34 - 140	12	32					
Fluoranthene	<b>0.492</b>	<b>0.488</b>	0.500	0.500	98	98	45 - 120	1	39					
Pyrene	<b>0.495</b>	<b>0.483</b>	0.500	0.500	99	97	42 - 133	2	39					
Benzo[a]anthracene	<b>0.503</b>	<b>0.520</b>	0.500	0.500	101	104	71 - 117	3	28					
Chrysene	<b>0.467</b>	<b>0.486</b>	0.500	0.500	93	97	53 - 110	4	25					
Benzo[b]fluoranthene	<b>0.503</b>	<b>0.521</b>	0.500	0.500	101	104	53 - 123	4	37					
Benzo(j,k)fluoranthene	<b>0.471</b>	<b>0.503</b>	0.500	0.500	94	101	52 - 119	7	41					
Benzo[a]pyrene	<b>0.466</b>	<b>0.478</b>	0.500	0.500	93	96	37 - 129	3	33					
Indeno(1,2,3-c,d)pyrene	<b>0.487</b>	<b>0.491</b>	0.500	0.500	97	98	45 - 128	1	31					
Dibenz[a,h]anthracene	<b>0.506</b>	<b>0.520</b>	0.500	0.500	101	104	54 - 120	3	30					
Benzo[g,h,i]perylene	<b>0.501</b>	<b>0.518</b>	0.500	0.500	100	104	49 - 117	3	29					
<i>Surrogate:</i>														
<i>2-Fluorobiphenyl</i>					91	94	30 - 124							
<i>Pyrene-d10</i>					97	96	40 - 143							
<i>Terphenyl-d14</i>					110	110	27 - 127							



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Date of Report: September 14, 2017  
 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

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

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	EPA Method	Date Prepared	Date Analyzed	Flags
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Lab ID: 09-019-01  
**Client ID:** KMW-10-8-31-17

Arsenic	<b>10</b>	3.3	200.8	9-12-17	9-12-17
Barium	<b>ND</b>	28	200.8	9-12-17	9-12-17
Cadmium	<b>ND</b>	4.4	200.8	9-12-17	9-12-17
Chromium	<b>ND</b>	11	200.8	9-12-17	9-12-17
Lead	<b>ND</b>	1.1	200.8	9-12-17	9-12-17
Mercury	<b>ND</b>	0.50	7470A	9-7-17	9-7-17
Selenium	<b>ND</b>	5.6	200.8	9-12-17	9-12-17
Silver	<b>ND</b>	11	200.8	9-12-17	9-12-17

Lab ID: 09-019-02  
**Client ID:** KMW-09-8-31-17

Arsenic	<b>ND</b>	3.3	200.8	9-12-17	9-12-17
Barium	<b>ND</b>	28	200.8	9-12-17	9-12-17
Cadmium	<b>ND</b>	4.4	200.8	9-12-17	9-12-17
Chromium	<b>ND</b>	11	200.8	9-12-17	9-12-17
Lead	<b>ND</b>	1.1	200.8	9-12-17	9-12-17
Mercury	<b>ND</b>	0.50	7470A	9-7-17	9-7-17
Selenium	<b>ND</b>	5.6	200.8	9-12-17	9-12-17
Silver	<b>ND</b>	11	200.8	9-12-17	9-12-17



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 Project: 14697.00008

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

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	EPA Method	Date Prepared	Date Analyzed	Flags
Lab ID:	09-019-03					
<b>Client ID:</b>	<b>KMW-06-8-31-17</b>					
Arsenic	<b>ND</b>	3.3	200.8	9-12-17	9-12-17	
Barium	<b>ND</b>	28	200.8	9-12-17	9-12-17	
Cadmium	<b>ND</b>	4.4	200.8	9-12-17	9-12-17	
Chromium	<b>ND</b>	11	200.8	9-12-17	9-12-17	
Lead	<b>2.7</b>	1.1	200.8	9-12-17	9-12-17	
Mercury	<b>ND</b>	0.50	7470A	9-7-17	9-7-17	
Selenium	<b>ND</b>	5.6	200.8	9-12-17	9-12-17	
Silver	<b>ND</b>	11	200.8	9-12-17	9-12-17	

Lab ID: 09-019-04  
**Client ID:** KMW-04-8-31-17

Arsenic	<b>14</b>	3.3	200.8	9-12-17	9-12-17
Barium	<b>ND</b>	28	200.8	9-12-17	9-12-17
Cadmium	<b>ND</b>	4.4	200.8	9-12-17	9-12-17
Chromium	<b>ND</b>	11	200.8	9-12-17	9-12-17
Lead	<b>ND</b>	1.1	200.8	9-12-17	9-12-17
Mercury	<b>ND</b>	0.50	7470A	9-7-17	9-7-17
Selenium	<b>ND</b>	5.6	200.8	9-12-17	9-12-17
Silver	<b>ND</b>	11	200.8	9-12-17	9-12-17



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

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	EPA Method	Prepared	Date	Analyzed	Date	Flags
Lab ID:	09-019-05							
<b>Client ID:</b>	<b>KMW-02R-9-1-17</b>							
Arsenic	ND	3.3	200.8	9-12-17	9-12-17			
Barium	ND	28	200.8	9-12-17	9-12-17			
Cadmium	ND	4.4	200.8	9-12-17	9-12-17			
Chromium	ND	11	200.8	9-12-17	9-12-17			
Lead	ND	1.1	200.8	9-12-17	9-12-17			
Mercury	ND	0.50	7470A	9-7-17	9-7-17			
Selenium	ND	5.6	200.8	9-12-17	9-12-17			
Silver	ND	11	200.8	9-12-17	9-12-17			

Lab ID: 09-019-06  
**Client ID:** KMW-07-9-1-17

Arsenic	ND	3.3	200.8	9-12-17	9-12-17			
Barium	ND	28	200.8	9-12-17	9-12-17			
Cadmium	ND	4.4	200.8	9-12-17	9-12-17			
Chromium	ND	11	200.8	9-12-17	9-12-17			
Lead	ND	1.1	200.8	9-12-17	9-12-17			
Mercury	ND	0.50	7470A	9-7-17	9-7-17			
Selenium	ND	5.6	200.8	9-12-17	9-12-17			
Silver	ND	11	200.8	9-12-17	9-12-17			



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 Laboratory Reference: 1709-019  
 Project: 14697.00008

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

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	EPA Method	Date Prepared	Date Analyzed	Flags
Lab ID:	09-019-07					
<b>Client ID:</b>	<b>KMW-08-9-1-17</b>					
Arsenic	ND	3.3	200.8	9-12-17	9-12-17	
Barium	ND	28	200.8	9-12-17	9-12-17	
Cadmium	ND	4.4	200.8	9-12-17	9-12-17	
Chromium	ND	11	200.8	9-12-17	9-12-17	
Lead	ND	1.1	200.8	9-12-17	9-12-17	
Mercury	ND	0.50	7470A	9-7-17	9-7-17	
Selenium	ND	5.6	200.8	9-12-17	9-12-17	
Silver	ND	11	200.8	9-12-17	9-12-17	

Lab ID: 09-019-08  
**Client ID:** KMW-03R-9-1-17

Arsenic	ND	3.3	200.8	9-12-17	9-12-17	
Barium	ND	28	200.8	9-12-17	9-12-17	
Cadmium	ND	4.4	200.8	9-12-17	9-12-17	
Chromium	ND	11	200.8	9-12-17	9-12-17	
Lead	ND	1.1	200.8	9-12-17	9-12-17	
Mercury	ND	0.50	7470A	9-7-17	9-7-17	
Selenium	ND	5.6	200.8	9-12-17	9-12-17	
Silver	ND	11	200.8	9-12-17	9-12-17	



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 Project: 14697.00008

**TOTAL METALS**  
**EPA 200.8/7470A**  
**METHOD BLANK QUALITY CONTROL**

Date Extracted: 9-7&12-17  
 Date Analyzed: 9-7&12-17  
 Matrix: Water  
 Units: ug/L (ppb)  
 Lab ID: MB0907W1&MB0912WM1

Analyte	Method	Result	PQL
Arsenic	200.8	<b>ND</b>	3.3
Barium	200.8	<b>ND</b>	28
Cadmium	200.8	<b>ND</b>	4.4
Chromium	200.8	<b>ND</b>	11
Lead	200.8	<b>ND</b>	1.1
Mercury	7470A	<b>ND</b>	0.50
Selenium	200.8	<b>ND</b>	5.6
Silver	200.8	<b>ND</b>	11



Date of Report: September 14, 2017  
 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

**TOTAL METALS  
EPA 200.8/7470A  
DUPLICATE QUALITY CONTROL**

Date Extracted: 9-7&12-17  
 Date Analyzed: 9-7&12-17

Matrix: Water  
 Units: ug/L (ppb)

Lab ID: 09-019-05

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Arsenic	ND	ND	NA	3.3	
Barium	ND	ND	NA	28	
Cadmium	ND	ND	NA	4.4	
Chromium	ND	ND	NA	11	
Lead	ND	ND	NA	1.1	
Mercury	ND	ND	NA	0.50	
Selenium	ND	ND	NA	5.6	
Silver	ND	ND	NA	11	



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

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Date of Report: September 14, 2017  
 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

**TOTAL METALS**  
**EPA 200.8/7470A**  
**MS/MSD QUALITY CONTROL**

Date Extracted: 9-7&12-17  
 Date Analyzed: 9-7&12-17

Matrix: Water  
 Units: ug/L (ppb)

Lab ID: 09-019-05

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Arsenic	222	<b>242</b>	109	<b>245</b>	110	1	
Barium	222	<b>225</b>	101	<b>227</b>	102	1	
Cadmium	222	<b>242</b>	109	<b>244</b>	110	1	
Chromium	222	<b>201</b>	90	<b>207</b>	93	3	
Lead	222	<b>225</b>	101	<b>229</b>	103	2	
Mercury	12.5	<b>12.9</b>	103	<b>12.8</b>	102	1	
Selenium	222	<b>255</b>	115	<b>251</b>	113	2	
Silver	222	<b>234</b>	106	<b>243</b>	109	4	



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 Laboratory Reference: 1709-019  
 Project: 14697.00008

**TOTAL METALS**  
**EPA 200.8/7470A**  
**SPIKE BLANK QUALITY CONTROL**

Date Extracted: 9-7&12-17  
 Date Analyzed: 9-7&12-17

Matrix: Water  
 Units: ug/L (ppb)

Lab ID: SB0907W1&SB0912WM1

Analyte	Method	Spike Level	Result	Percent Recovery
Arsenic	200.8	222	<b>226</b>	102
Barium	200.8	222	<b>214</b>	96
Cadmium	200.8	222	<b>228</b>	103
Chromium	200.8	222	<b>195</b>	88
Lead	200.8	222	<b>221</b>	100
Mercury	7470A	12.5	<b>13.1</b>	105
Selenium	200.8	222	<b>238</b>	107
Silver	200.8	222	<b>226</b>	102



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Date of Report: September 14, 2017  
 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

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

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	EPA Method	Date Prepared	Date Analyzed	Flags
Lab ID:	09-019-01					
<b>Client ID:</b>	<b>KMW-10-8-31-17</b>					
Arsenic	ND	3.0	200.8	9-1-17	9-12-17	
Barium	ND	25	200.8	9-1-17	9-12-17	
Cadmium	ND	4.0	200.8	9-1-17	9-12-17	
Chromium	ND	10	200.8	9-1-17	9-12-17	
Lead	ND	1.0	200.8	9-1-17	9-12-17	
Mercury	ND	0.50	7470A	9-1-17	9-7-17	
Selenium	ND	5.0	200.8	9-1-17	9-12-17	
Silver	ND	10	200.8	9-1-17	9-12-17	
Lab ID:	09-019-02					
<b>Client ID:</b>	<b>KMW-09-8-31-17</b>					
Arsenic	ND	3.0	200.8	9-1-17	9-12-17	
Barium	ND	25	200.8	9-1-17	9-12-17	
Cadmium	ND	4.0	200.8	9-1-17	9-12-17	
Chromium	ND	10	200.8	9-1-17	9-12-17	
Lead	ND	1.0	200.8	9-1-17	9-12-17	
Mercury	ND	0.50	7470A	9-1-17	9-7-17	
Selenium	ND	5.0	200.8	9-1-17	9-12-17	
Silver	ND	10	200.8	9-1-17	9-12-17	



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Date of Report: September 14, 2017  
 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

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

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	EPA Method	Prepared	Date	Date Analyzed	Flags
Lab ID:	09-019-03						
<b>Client ID:</b>	<b>KMW-06-8-31-17</b>						
Arsenic	<b>ND</b>	3.0	200.8	9-1-17	9-12-17		
Barium	<b>ND</b>	25	200.8	9-1-17	9-12-17		
Cadmium	<b>ND</b>	4.0	200.8	9-1-17	9-12-17		
Chromium	<b>ND</b>	10	200.8	9-1-17	9-12-17		
Lead	<b>1.1</b>	1.0	200.8	9-1-17	9-12-17		
Mercury	<b>ND</b>	0.50	7470A	9-1-17	9-7-17		
Selenium	<b>ND</b>	5.0	200.8	9-1-17	9-12-17		
Silver	<b>ND</b>	10	200.8	9-1-17	9-12-17		

Lab ID: 09-019-04  
**Client ID:** KMW-04-8-31-17

Arsenic	<b>5.1</b>	3.0	200.8	9-1-17	9-12-17		
Barium	<b>ND</b>	25	200.8	9-1-17	9-12-17		
Cadmium	<b>ND</b>	4.0	200.8	9-1-17	9-12-17		
Chromium	<b>ND</b>	10	200.8	9-1-17	9-12-17		
Lead	<b>ND</b>	1.0	200.8	9-1-17	9-12-17		
Mercury	<b>ND</b>	0.50	7470A	9-1-17	9-7-17		
Selenium	<b>ND</b>	5.0	200.8	9-1-17	9-12-17		
Silver	<b>ND</b>	10	200.8	9-1-17	9-12-17		



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Date of Report: September 14, 2017  
 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

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

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	EPA Method	Prepared	Date	Date Analyzed	Flags
Lab ID:	09-019-05						
<b>Client ID:</b>	<b>KMW-02R-9-1-17</b>						
Arsenic	<b>ND</b>	3.0	200.8	9-1-17	9-12-17		
Barium	<b>ND</b>	25	200.8	9-1-17	9-12-17		
Cadmium	<b>ND</b>	4.0	200.8	9-1-17	9-12-17		
Chromium	<b>ND</b>	10	200.8	9-1-17	9-12-17		
Lead	<b>ND</b>	1.0	200.8	9-12-17	9-12-17		
Mercury	<b>ND</b>	0.50	7470A	9-1-17	9-7-17		
Selenium	<b>ND</b>	5.0	200.8	9-1-17	9-12-17		
Silver	<b>ND</b>	10	200.8	9-1-17	9-12-17		

Lab ID: 09-019-06  
**Client ID:** KMW-07-9-1-17

Arsenic	<b>ND</b>	3.0	200.8	9-1-17	9-12-17		
Barium	<b>ND</b>	25	200.8	9-1-17	9-12-17		
Cadmium	<b>ND</b>	4.0	200.8	9-1-17	9-12-17		
Chromium	<b>ND</b>	10	200.8	9-1-17	9-12-17		
Lead	<b>ND</b>	1.0	200.8	9-12-17	9-12-17		
Mercury	<b>ND</b>	0.50	7470A	9-1-17	9-7-17		
Selenium	<b>ND</b>	5.0	200.8	9-1-17	9-12-17		
Silver	<b>ND</b>	10	200.8	9-1-17	9-12-17		



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Date of Report: September 14, 2017  
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 Laboratory Reference: 1709-019  
 Project: 14697.00008

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

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	EPA Method	Prepared	Date	Date Analyzed	Flags
Lab ID:	09-019-07						
<b>Client ID:</b>	<b>KMW-08-9-1-17</b>						
Arsenic	ND	3.0	200.8	9-1-17	9-12-17		
Barium	ND	25	200.8	9-1-17	9-12-17		
Cadmium	ND	4.0	200.8	9-1-17	9-12-17		
Chromium	ND	10	200.8	9-1-17	9-12-17		
Lead	ND	1.0	200.8	9-1-17	9-12-17		
Mercury	ND	0.50	7470A	9-1-17	9-7-17		
Selenium	ND	5.0	200.8	9-1-17	9-12-17		
Silver	ND	10	200.8	9-1-17	9-12-17		

Lab ID: 09-019-08  
**Client ID:** KMW-03R-9-1-17

Arsenic	ND	3.0	200.8	9-1-17	9-12-17	
Barium	ND	25	200.8	9-1-17	9-12-17	
Cadmium	ND	4.0	200.8	9-1-17	9-12-17	
Chromium	ND	10	200.8	9-1-17	9-12-17	
Lead	ND	1.0	200.8	9-12-17	9-12-17	
Mercury	ND	0.50	7470A	9-1-17	9-7-17	
Selenium	ND	5.0	200.8	9-1-17	9-12-17	
Silver	ND	10	200.8	9-1-17	9-12-17	



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 Samples Submitted: September 1, 2017  
 Laboratory Reference: 1709-019  
 Project: 14697.00008

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

Date Filtered: 9-1-17  
 Date Analyzed: 9-7&12-17

Matrix: Water  
 Units: ug/L (ppb)

Lab ID: MB0901F1

Analyte	Method	Result	PQL
Arsenic	200.8	<b>ND</b>	3.0
Barium	200.8	<b>ND</b>	25
Cadmium	200.8	<b>ND</b>	4.0
Chromium	200.8	<b>ND</b>	10
Lead	200.8	<b>ND</b>	1.0
Mercury	7470A	<b>ND</b>	0.50
Selenium	200.8	<b>ND</b>	5.0
Silver	200.8	<b>ND</b>	10



Date of Report: September 14, 2017  
Samples Submitted: September 1, 2017  
Laboratory Reference: 1709-019  
Project: 14697.00008

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

Date Filtered: 9-12-17  
Date Analyzed: 9-12-17

Matrix: Water  
Units: ug/L (ppb)

Lab ID: MB0912F1

Analyte	Method	Result	PQL
Lead	200.8	ND	1.0



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 Project: 14697.00008

**DISSOLVED METALS**  
**EPA 200.8/7470A**  
**DUPPLICATE QUALITY CONTROL**

Date Filtered: 9-1&12-17  
 Date Analyzed: 9-7&12-17

Matrix: Water  
 Units: ug/L (ppb)

Lab ID: 09-019-05

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Arsenic	<b>ND</b>	<b>ND</b>	NA	3.0	
Barium	<b>ND</b>	<b>ND</b>	NA	25	
Cadmium	<b>ND</b>	<b>ND</b>	NA	4.0	
Chromium	<b>ND</b>	<b>ND</b>	NA	10	
Lead	<b>ND</b>	<b>ND</b>	NA	1.0	
Mercury	<b>ND</b>	<b>ND</b>	NA	0.50	
Selenium	<b>ND</b>	<b>ND</b>	NA	5.0	
Silver	<b>ND</b>	<b>ND</b>	NA	10	




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 Laboratory Reference: 1709-019  
 Project: 14697.00008

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

Date Filtered: 9-1&12-17  
 Date Analyzed: 9-7&12-17

Matrix: Water  
 Units: ug/L (ppb)

Lab ID: 09-019-05

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Arsenic	200	<b>214</b>	107	<b>213</b>	107	0	
Barium	200	<b>193</b>	97	<b>194</b>	97	1	
Cadmium	200	<b>212</b>	106	<b>215</b>	107	1	
Chromium	200	<b>177</b>	89	<b>180</b>	90	1	
Lead	200	<b>194</b>	97	<b>194</b>	97	0	
Mercury	12.5	<b>12.9</b>	103	<b>13.1</b>	105	1	
Selenium	200	<b>214</b>	107	<b>220</b>	110	3	
Silver	200	<b>172</b>	86	<b>170</b>	85	1	



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**DISSOLVED METALS**  
**EPA 200.8/7470A**  
**SPIKE BLANK QUALITY CONTROL**

Date Filtered: 9-1-17  
 Date Analyzed: 9-7&12-17

Matrix: Water  
 Units: ug/L (ppb)

Lab ID: SB0901F1

Analyte	Method	Spike Level	Result	Percent Recovery
Arsenic	200.8	200	<b>208</b>	104
Barium	200.8	200	<b>189</b>	95
Cadmium	200.8	200	<b>212</b>	106
Chromium	200.8	200	<b>178</b>	89
Lead	200.8	200	<b>195</b>	97
Mercury	7470A	12.5	<b>12.8</b>	102
Selenium	200.8	200	<b>214</b>	107
Silver	200.8	200	<b>170</b>	85



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Date of Report: September 14, 2017  
Samples Submitted: September 1, 2017  
Laboratory Reference: 1709-019  
Project: 14697.00008

**DISSOLVED LEAD  
EPA 200.8  
SPIKE BLANK QUALITY CONTROL**

Date Filtered: 9-12-17  
Date Analyzed: 9-12-17

Matrix: Water  
Units: ug/L (ppb)

Lab ID: SB0912F1

Analyte	Method	Spike Level	Result	Percent Recovery
Lead	200.8	200	195	98



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### 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



## Chain of Custody

 Page 1 of 1

Turnaround Request (in working days)				Laboratory Number:
(Check One)				09-019

 Company **AFLAC FW**  
 Project Number: **14697.00008**

 Project Name: **Kelly-Moore**  
 Project Manager: **Cristy Thomsen**

 Sampled by: **Brian Haerlie**
 Standard (7 Days)  
 NTPH analysis (5 Days)

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

Lab ID	Sample Identification	Date Sampled	Time Sampled	Matrix	Number of Containers
1	KMW-10-8-31-17	8/31/17	1200	W	11
2	KMW-09-8-31-17	8/31/17	1330	W	11
3	KMW-06-8-31-17	8/31/17	1510	W	11
4	KMW-04-8-31-17	8/31/17	1650	W	11
5	KMW-02R-9-1-17	9/1/17	0935	W	13
6	KMW-07-9-1-17	9/1/17	1150	W	11
7	KMW-08-9-1-17	9/1/17	1340	W	11
8	KMW-03R-9-1-17	9/1/17	1440	W	11
9	TRIP BLANK	-	-	W	11

NWTPH-HCID	X
NWTPH-Gx/BTEX	
NWTPH-Gx	X
NWTPH-Dx ( <input type="checkbox"/> Acid / SG Clean-up)	
Volatiles 8260C	X
Halogenated Volatiles 8260C	X
EDB EPA 8011 (Waters Only)	
Semivolatiles 8270D/SIM (with low-level PAHs)	
PAHs 8270D/SIM (low-level)	
PCBs 8082A	
Organochlorine Pesticides 8081B	
Organophosphorus Pesticides 8270D/SIM	
Chlorinated Acid Herbicides 8151A	
Total RCRA Metals	
Total MTCA Metals	
TCLP Metals	
HEM (oil and grease) 1664A	
Total Metals - RRA 8	X
Dissolved Metals - RRA 8	X
% Moisture	

Signature	Company	Date	Time	Comments/Special Instructions
Relinquished	AFLAC FW	9/1/17	1505	lab will filter dissolved metals,
Received	Speddy	9/1/17	1505	dissolved metals unrepresented
Relinquished	OSCE	9/1/17	1620	KMW-02R was used
Received				
Relinquished				
Received				
Reviewed/Dates				

 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

February 2, 2018

Kathleen Goodman  
AMEC Foster Wheeler  
Environment and Infrastructure, Inc.  
One Union Square  
600 University Street, Suite 600  
Seattle, WA 98101

Re: Analytical Data for Project 0146970060  
Laboratory Reference No. 1801-261

Dear Kathleen:

Enclosed are the analytical results and associated quality control data for samples submitted on January 25, 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" followed by a cursive surname.

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

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Date of Report: February 2, 2018  
Samples Submitted: January 25, 2018  
Laboratory Reference: 1801-261  
Project: 0146970060

#### Case Narrative

Samples were collected on January 24, 2018 and received by the laboratory on January 25, 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.



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Date of Report: February 2, 2018  
 Samples Submitted: January 25, 2018  
 Laboratory Reference: 1801-261  
 Project: 0146970060

### NWTPH-Gx

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-06-012418</b>					
Laboratory ID:	01-261-01					
Gasoline	<b>1300</b>	100	NWTPH-Gx	1-26-18	1-26-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	81	66-114				
<b>Client ID:</b>	<b>KMW-09-012418</b>					
Laboratory ID:	01-261-02					
Gasoline	<b>760</b>	100	NWTPH-Gx	1-26-18	1-26-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	78	66-114				



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 Laboratory Reference: 1801-261  
 Project: 0146970060

**NWTPH-Gx**  
**QUALITY CONTROL**

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB0126W1					
Gasoline	ND	100	NWTPH-Gx	1-26-18	1-26-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	94	66-114				
Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD RPD Limit Flags
<b>DUPPLICATE</b>						
Laboratory ID:	01-271-01					
	ORIG	DUP				
Gasoline	265	256	NA	NA	NA	3 30
Surrogate:						
Fluorobenzene				89	81	66-114



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 Laboratory Reference: 1801-261  
 Project: 0146970060

### NWTPH-Dx

Matrix: Water  
 Units: mg/L (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-06-012418</b>					
Laboratory ID:	01-261-01					
Diesel Range Organics	<b>4.2</b>	0.26	NWTPH-Dx	1-29-18	1-29-18	
Lube Oil Range Organics	<b>0.60</b>	0.41	NWTPH-Dx	1-29-18	1-29-18	

Surrogate:	Percent Recovery	Control Limits
<i>o-Terphenyl</i>	84	50-150

**Client ID:** KMW-09-012418  
**Laboratory ID:** 01-261-02

Diesel Range Organics	<b>3.1</b>	0.26	NWTPH-Dx	1-29-18	1-29-18
Lube Oil Range Organics	<b>0.69</b>	0.41	NWTPH-Dx	1-29-18	1-29-18
Surrogate:	Percent Recovery	Control Limits			

*o-Terphenyl* 87 50-150



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 Project: 0146970060

**NWTPH-Dx**  
**QUALITY CONTROL**

Matrix: Water  
 Units: mg/L (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB0129W1					
Diesel Range Organics	ND	0.25	NWTPH-Dx	1-29-18	1-29-18	
Lube Oil Range Organics	ND	0.40	NWTPH-Dx	1-29-18	1-29-18	
Surrogate: <i>o-Terphenyl</i>	Percent Recovery 93	Control Limits 50-150				

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>								
Laboratory ID:	01-271-01							
	ORIG	DUP						
Diesel Range Organics	2.30	2.22	NA	NA	NA	NA	4	NA
Lube Oil Range	ND	ND	NA	NA	NA	NA	NA	NA
Surrogate: <i>o-Terphenyl</i>				83	80	50-150		

**SPIKE BLANK**

Laboratory ID: SB0129W1

Diesel Fuel #2	0.784	1.00	NA	78	58-113	NA	NA
<i>Surrogate: <i>o-Terphenyl</i></i>							
				90	50-150		



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Date of Report: February 2, 2018  
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 Laboratory Reference: 1801-261  
 Project: 0146970060

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

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-06-012418</b>					
<b>Laboratory ID:</b>	01-261-01					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloromethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Vinyl Chloride	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromomethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloroethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Acetone	ND	5.0	EPA 8260C	1-29-18	1-29-18	
Iodomethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Carbon Disulfide	1.1	0.20	EPA 8260C	1-29-18	1-29-18	
Methylene Chloride	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Vinyl Acetate	ND	1.0	EPA 8260C	1-29-18	1-29-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Butanone	ND	5.0	EPA 8260C	1-29-18	1-29-18	
Bromochloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloroform	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Benzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Trichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Dibromomethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromodichloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	1-29-18	1-29-18	
Toluene	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	



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Date of Report: February 2, 2018  
 Samples Submitted: January 25, 2018  
 Laboratory Reference: 1801-261  
 Project: 0146970060

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

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-06-012418</b>					
<b>Laboratory ID:</b>	01-261-01					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Tetrachloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Hexanone	ND	2.0	EPA 8260C	1-29-18	1-29-18	
Dibromochloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Ethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
m,p-Xylene	0.42	0.40	EPA 8260C	1-29-18	1-29-18	
o-Xylene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Styrene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromoform	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Isopropylbenzene	8.1	0.20	EPA 8260C	1-29-18	1-29-18	
Bromobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
n-Propylbenzene	9.5	0.20	EPA 8260C	1-29-18	1-29-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
sec-Butylbenzene	1.2	0.20	EPA 8260C	1-29-18	1-29-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
n-Butylbenzene	1.0	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Hexachlorobutadiene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Naphthalene	ND	1.0	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
<b>Surrogate:</b>	<b>Percent Recovery</b>	<b>Control Limits</b>				
Dibromofluoromethane	99	75-127				
Toluene-d8	102	80-127				
4-Bromofluorobenzene	114	78-125				



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Date of Report: February 2, 2018  
 Samples Submitted: January 25, 2018  
 Laboratory Reference: 1801-261  
 Project: 0146970060

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

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-09-012418</b>					
<b>Laboratory ID:</b>	01-261-02					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloromethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Vinyl Chloride	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromomethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloroethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Acetone	ND	5.0	EPA 8260C	1-29-18	1-29-18	
Iodomethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Carbon Disulfide	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methylene Chloride	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Vinyl Acetate	ND	1.0	EPA 8260C	1-29-18	1-29-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Butanone	ND	5.0	EPA 8260C	1-29-18	1-29-18	
Bromochloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloroform	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Benzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Trichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Dibromomethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromodichloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	1-29-18	1-29-18	
Toluene	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	



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 Project: 0146970060

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

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-09-012418</b>					
<b>Laboratory ID:</b>	01-261-02					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Tetrachloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Hexanone	ND	2.0	EPA 8260C	1-29-18	1-29-18	
Dibromochloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Ethylbenzene	0.44	0.20	EPA 8260C	1-29-18	1-29-18	
m,p-Xylene	ND	0.40	EPA 8260C	1-29-18	1-29-18	
o-Xylene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Styrene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromoform	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Isopropylbenzene	7.1	0.20	EPA 8260C	1-29-18	1-29-18	
Bromobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
n-Propylbenzene	7.4	0.20	EPA 8260C	1-29-18	1-29-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
sec-Butylbenzene	0.70	0.20	EPA 8260C	1-29-18	1-29-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
n-Butylbenzene	0.52	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Hexachlorobutadiene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Naphthalene	ND	1.0	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
<b>Surrogate:</b>	<b>Percent Recovery</b>	<b>Control Limits</b>				
Dibromofluoromethane	103	75-127				
Toluene-d8	102	80-127				
4-Bromofluorobenzene	104	78-125				



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 Project: 0146970060

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

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0129W1					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloromethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Vinyl Chloride	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromomethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloroethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Acetone	ND	5.0	EPA 8260C	1-29-18	1-29-18	
Iodomethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Carbon Disulfide	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methylene Chloride	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Vinyl Acetate	ND	1.0	EPA 8260C	1-29-18	1-29-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Butanone	ND	5.0	EPA 8260C	1-29-18	1-29-18	
Bromochloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloroform	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Benzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Trichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Dibromomethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromodichloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	1-29-18	1-29-18	
Toluene	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	



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

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Date of Report: February 2, 2018  
 Samples Submitted: January 25, 2018  
 Laboratory Reference: 1801-261  
 Project: 0146970060

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

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0129W1					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Tetrachloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Hexanone	ND	2.0	EPA 8260C	1-29-18	1-29-18	
Dibromochloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Ethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
m,p-Xylene	ND	0.40	EPA 8260C	1-29-18	1-29-18	
o-Xylene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Styrene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromoform	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Isopropylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
n-Propylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
n-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Hexachlorobutadiene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Naphthalene	ND	1.0	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	102	75-127				
Toluene-d8	107	80-127				
4-Bromofluorobenzene	96	78-125				



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Date of Report: February 2, 2018  
 Samples Submitted: January 25, 2018  
 Laboratory Reference: 1801-261  
 Project: 0146970060

**VOLATILES by EPA 8260C**  
**MS/MSD QUALITY CONTROL**

Matrix: Water

Units: ug/L

Analyte	Result	Spike Level	Source	Percent	Recovery	RPD	RPD	Limit	Flags
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**MATRIX SPIKES**

Laboratory ID: 01-271-01

	MS	MSD	MS	MSD	MS	MSD			
1,1-Dichloroethene	<b>8.70</b>	<b>9.68</b>	10.0	10.0	ND	87	97	58-121	11
Benzene	<b>9.39</b>	<b>10.4</b>	10.0	10.0	ND	94	104	67-122	10
Trichloroethene	<b>8.37</b>	<b>9.52</b>	10.0	10.0	ND	84	95	60-120	13
Toluene	<b>9.42</b>	<b>10.5</b>	10.0	10.0	ND	94	105	66-131	11
Chlorobenzene	<b>8.95</b>	<b>9.76</b>	10.0	10.0	ND	90	98	73-120	9

*Surrogate:*

*Dibromofluoromethane* 94 97 75-127

*Toluene-d8* 95 100 80-127

*4-Bromofluorobenzene* 94 99 78-125



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 Laboratory Reference: 1801-261  
 Project: 0146970060

**VOLATILES by EPA 8260C**  
**SPIKE BLANK QUALITY CONTROL**

Matrix: Water  
 Units: ug/L

Analyte	Result	Spike Level	Percent Recovery	Recovery Limits	Flags
<b>SPIKE BLANK</b>					
Laboratory ID: SB0129W1					
1,1-Dichloroethene	<b>9.35</b>	10.0	94	63-126	
Benzene	<b>9.55</b>	10.0	96	78-122	
Trichloroethene	<b>9.18</b>	10.0	92	63-120	
Toluene	<b>10.0</b>	10.0	100	79-124	
Chlorobenzene	<b>9.60</b>	10.0	96	78-120	
<i>Surrogate:</i>					
<i>Dibromofluoromethane</i>			97	75-127	
<i>Toluene-d8</i>			102	80-127	
<i>4-Bromofluorobenzene</i>			94	78-125	



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 Laboratory Reference: 1801-261  
 Project: 0146970060

### PAHs EPA 8270D/SIM

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-06-012418</b>					
<b>Laboratory ID:</b>	01-261-01					
Naphthalene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
2-Methylnaphthalene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
1-Methylnaphthalene	1.3	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Acenaphthylene	0.13	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Acenaphthene	2.5	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Fluorene	0.56	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Phenanthrene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Anthracene	0.41	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Fluoranthene	0.12	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Pyrene	0.22	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[a]anthracene	0.055	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Chrysene	0.062	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[b]fluoranthene	0.082	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo(j,k)fluoranthene	0.034	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[a]pyrene	0.071	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Indeno(1,2,3-c,d)pyrene	0.054	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Dibenz[a,h]anthracene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[g,h,i]perylene	0.056	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
2-Fluorobiphenyl	96		25 - 107			
Pyrene-d10	70		28 - 103			
Terphenyl-d14	112		36 - 129			



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Date of Report: February 2, 2018  
 Samples Submitted: January 25, 2018  
 Laboratory Reference: 1801-261  
 Project: 0146970060

### PAHs EPA 8270D/SIM

Matrix: Water

Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-09-012418</b>					
<b>Laboratory ID:</b>	01-261-02					
Naphthalene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
2-Methylnaphthalene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
1-Methylnaphthalene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Acenaphthylene	0.32	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Acenaphthene	2.4	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Fluorene	0.22	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Phenanthrene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Anthracene	0.62	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Fluoranthene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Pyrene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[a]anthracene	0.020	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Chrysene	0.018	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[b]fluoranthene	0.017	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo(j,k)fluoranthene	0.015	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[a]pyrene	0.021	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Indeno(1,2,3-c,d)pyrene	0.013	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Dibenz[a,h]anthracene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[g,h,i]perylene	0.011	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
2-Fluorobiphenyl	76		25 - 107			
Pyrene-d10	70		28 - 103			
Terphenyl-d14	111		36 - 129			



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 Laboratory Reference: 1801-261  
 Project: 0146970060

**PAHs EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**

Matrix: Water

Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0130W1					
Naphthalene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
2-Methylnaphthalene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
1-Methylnaphthalene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Acenaphthylene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Acenaphthene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Fluorene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Phenanthrene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Anthracene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Fluoranthene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Pyrene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Chrysene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	56	25 - 107				
Pyrene-d10	75	28 - 103				
Terphenyl-d14	84	36 - 129				



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 Project: 0146970060

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

Matrix: Water

Units: ug/L

Analyte	Result	Spike Level	Source	Percent	Recovery	RPD	RPD Limit	Flags				
			Result	Recovery	Limits							
<b>MATRIX SPIKES</b>												
Laboratory ID:		01-271-01										
	MS	MSD	MS	MSD	MS	MSD						
Naphthalene	<b>0.421</b>	<b>0.469</b>	0.588	0.569	ND	72	82	27 - 106				
Acenaphthylene	<b>0.518</b>	<b>0.526</b>	0.588	0.569	ND	88	92	20 - 117				
Acenaphthene	<b>0.648</b>	<b>0.582</b>	0.588	0.569	ND	110	102	30 - 114				
Fluorene	<b>0.552</b>	<b>0.529</b>	0.588	0.569	ND	94	93	36 - 116				
Phenanthrene	<b>0.494</b>	<b>0.501</b>	0.588	0.569	ND	84	88	31 - 122				
Anthracene	<b>0.589</b>	<b>0.586</b>	0.588	0.569	ND	100	103	33 - 144				
Fluoranthene	<b>0.498</b>	<b>0.505</b>	0.588	0.569	ND	85	89	44 - 120				
Pyrene	<b>0.506</b>	<b>0.517</b>	0.588	0.569	ND	86	91	40 - 130				
Benzo[a]anthracene	<b>0.571</b>	<b>0.595</b>	0.588	0.569	ND	97	105	47 - 131				
Chrysene	<b>0.556</b>	<b>0.547</b>	0.588	0.569	ND	95	96	48 - 120				
Benzo[b]fluoranthene	<b>0.560</b>	<b>0.563</b>	0.588	0.569	ND	95	99	42 - 128				
Benzo(j,k)fluoranthene	<b>0.537</b>	<b>0.555</b>	0.588	0.569	ND	91	98	46 - 121				
Benzo[a]pyrene	<b>0.556</b>	<b>0.570</b>	0.588	0.569	ND	95	100	34 - 121				
Indeno(1,2,3-c,d)pyrene	<b>0.551</b>	<b>0.573</b>	0.588	0.569	ND	94	101	39 - 128				
Dibenz[a,h]anthracene	<b>0.557</b>	<b>0.572</b>	0.588	0.569	ND	95	101	39 - 125				
Benzo[g,h,i]perylene	<b>0.522</b>	<b>0.535</b>	0.588	0.569	ND	89	94	41 - 122				
<i>Surrogate:</i>												
<i>2-Fluorobiphenyl</i>					53	58	25 - 107					
<i>Pyrene-d10</i>					71	74	28 - 103					
<i>Terphenyl-d14</i>					98	101	36 - 129					



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 Laboratory Reference: 1801-261  
 Project: 0146970060

**PAHs EPA 8270D/SIM**  
**SPIKE BLANK QUALITY CONTROL**

Matrix: Water  
 Units: ug/L

Analyte	Result	Spike Level	Percent Recovery	Recovery Limits	Flags
<b>SPIKE BLANK</b>					
Laboratory ID:	SB0130W1				
	SB	SB	SB		
Naphthalene	<b>0.360</b>	0.500	72	27 - 106	
Acenaphthylene	<b>0.387</b>	0.500	77	20 - 117	
Acenaphthene	<b>0.461</b>	0.500	92	30 - 114	
Fluorene	<b>0.454</b>	0.500	91	36 - 116	
Phenanthrene	<b>0.458</b>	0.500	92	31 - 122	
Anthracene	<b>0.480</b>	0.500	96	33 - 144	
Fluoranthene	<b>0.488</b>	0.500	98	44 - 120	
Pyrene	<b>0.566</b>	0.500	113	40 - 130	
Benzo[a]anthracene	<b>0.503</b>	0.500	101	47 - 131	
Chrysene	<b>0.498</b>	0.500	100	48 - 120	
Benzo[b]fluoranthene	<b>0.491</b>	0.500	98	42 - 128	
Benzo(j,k)fluoranthene	<b>0.489</b>	0.500	98	46 - 121	
Benzo[a]pyrene	<b>0.502</b>	0.500	100	34 - 121	
Indeno(1,2,3-c,d)pyrene	<b>0.522</b>	0.500	104	39 - 128	
Dibenz[a,h]anthracene	<b>0.528</b>	0.500	106	39 - 125	
Benzo[g,h,i]perylene	<b>0.518</b>	0.500	104	41 - 122	
<i>Surrogate:</i>					
<i>2-Fluorobiphenyl</i>			83	25 - 107	
<i>Pyrene-d10</i>			85	28 - 103	
<i>Terphenyl-d14</i>			103	36 - 129	



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 Project: 0146970060

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

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	EPA Method	Date Prepared	Date Analyzed	Flags
Lab ID:	01-261-01					
<b>Client ID:</b>	<b>KMW-06-012418</b>					
Arsenic	<b>ND</b>	3.3	200.8	1-31-18	1-31-18	
Barium	<b>ND</b>	28	200.8	1-31-18	1-31-18	
Cadmium	<b>ND</b>	4.4	200.8	1-31-18	1-31-18	
Chromium	<b>ND</b>	11	200.8	1-31-18	1-31-18	
Lead	<b>7.3</b>	1.1	200.8	1-31-18	1-31-18	
Mercury	<b>ND</b>	0.50	7470A	1-30-18	1-30-18	
Selenium	<b>ND</b>	5.6	200.8	1-31-18	1-31-18	
Silver	<b>ND</b>	11	200.8	1-31-18	1-31-18	

Lab ID: 01-261-02  
**Client ID:** KMW-09-012418

Arsenic	<b>ND</b>	3.3	200.8	1-31-18	1-31-18
Barium	<b>ND</b>	28	200.8	1-31-18	1-31-18
Cadmium	<b>ND</b>	4.4	200.8	1-31-18	1-31-18
Chromium	<b>ND</b>	11	200.8	1-31-18	1-31-18
Lead	<b>3.0</b>	1.1	200.8	1-31-18	1-31-18
Mercury	<b>ND</b>	0.50	7470A	1-30-18	1-30-18
Selenium	<b>ND</b>	5.6	200.8	1-31-18	1-31-18
Silver	<b>ND</b>	11	200.8	1-31-18	1-31-18



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

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Date of Report: February 2, 2018  
 Samples Submitted: January 25, 2018  
 Laboratory Reference: 1801-261  
 Project: 0146970060

**TOTAL METALS**  
**EPA 200.8/7470A**  
**METHOD BLANK QUALITY CONTROL**

Date Extracted: 1-30&31-18  
 Date Analyzed: 1-30&31-18  
 Matrix: Water  
 Units: ug/L (ppb)  
 Lab ID: MB0130W1&MB0131WM1

Analyte	Method	Result	PQL
Arsenic	200.8	<b>ND</b>	3.3
Barium	200.8	<b>ND</b>	28
Cadmium	200.8	<b>ND</b>	4.4
Chromium	200.8	<b>ND</b>	11
Lead	200.8	<b>ND</b>	1.1
Mercury	7470A	<b>ND</b>	0.50
Selenium	200.8	<b>ND</b>	5.6
Silver	200.8	<b>ND</b>	11



Date of Report: February 2, 2018  
 Samples Submitted: January 25, 2018  
 Laboratory Reference: 1801-261  
 Project: 0146970060

**TOTAL METALS  
EPA 200.8/7470A  
DUPLICATE QUALITY CONTROL**

Date Extracted: 1-30&31-18  
 Date Analyzed: 1-30&31-18

Matrix: Water  
 Units: ug/L (ppb)

Lab ID: 01-271-01

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Arsenic	<b>6.78</b>	<b>6.33</b>	7	3.3	
Barium	<b>ND</b>	<b>ND</b>	NA	28	
Cadmium	<b>ND</b>	<b>ND</b>	NA	4.4	
Chromium	<b>ND</b>	<b>ND</b>	NA	11	
Lead	<b>1.12</b>	<b>ND</b>	NA	1.1	
Mercury	<b>ND</b>	<b>ND</b>	NA	0.50	
Selenium	<b>ND</b>	<b>ND</b>	NA	5.6	
Silver	<b>ND</b>	<b>ND</b>	NA	11	



Date of Report: February 2, 2018  
 Samples Submitted: January 25, 2018  
 Laboratory Reference: 1801-261  
 Project: 0146970060

**TOTAL METALS**  
**EPA 200.8/7470A**  
**MS/MSD QUALITY CONTROL**

Date Extracted: 1-30&31-18  
 Date Analyzed: 1-30&31-18

Matrix: Water  
 Units: ug/L (ppb)

Lab ID: 01-271-01

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Arsenic	222	<b>227</b>	99	<b>233</b>	102	3	
Barium	222	<b>224</b>	101	<b>222</b>	100	1	
Cadmium	222	<b>228</b>	103	<b>231</b>	104	1	
Chromium	222	<b>211</b>	95	<b>216</b>	97	2	
Lead	222	<b>207</b>	93	<b>207</b>	93	0	
Mercury	12.5	<b>12.3</b>	98	<b>12.4</b>	99	1	
Selenium	222	<b>231</b>	104	<b>232</b>	104	0	
Silver	222	<b>196</b>	88	<b>202</b>	91	3	



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 Samples Submitted: January 25, 2018  
 Laboratory Reference: 1801-261  
 Project: 0146970060

**TOTAL METALS**  
**EPA 200.8/7470A**  
**SPIKE BLANK QUALITY CONTROL**

Date Extracted: 1-30&31-18  
 Date Analyzed: 1-30&31-18  
 Matrix: Water  
 Units: ug/L (ppb)  
 Lab ID: SB0130W1&SB0131WM1

Analyte	Method	Spike Level	SB Result	Percent Recovery
Arsenic	200.8	222	<b>228</b>	103
Barium	200.8	222	<b>212</b>	96
Cadmium	200.8	222	<b>224</b>	101
Chromium	200.8	222	<b>210</b>	95
Lead	200.8	222	<b>217</b>	98
Mercury	7470A	12.5	<b>12.6</b>	101
Selenium	200.8	222	<b>240</b>	108
Silver	200.8	222	<b>207</b>	93





### 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





## Chain of Custody

Analytical Laboratory Testing Services  
14648 NE 95th Street • Redmond WA 98052

14040 NE 35th Street • Redmond, WA 98052  
Phone: (425) 883-3881 • [www.onsite-env.com](http://www.onsite-env.com)

Company:

**HMC Foster Wheeler**

Project Name: 014697

**Kelly Moore**  
Project Manager

Project Manager  
Kathleen Goodman

Sampled by:  
**K. Black**



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

February 2, 2018

Kathleen Goodman  
AMEC Foster Wheeler  
Environment and Infrastructure, Inc.  
One Union Square  
600 University Street, Suite 600  
Seattle, WA 98101

Re: Analytical Data for Project 0146970060  
Laboratory Reference No. 1801-271

Dear Kathleen:

Enclosed are the analytical results and associated quality control data for samples submitted on January 25, 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" followed by a cursive surname.

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

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and is intended only for the use of the individual or company to whom it is addressed.

Date of Report: February 2, 2018  
Samples Submitted: January 25, 2018  
Laboratory Reference: 1801-271  
Project: 0146970060

#### Case Narrative

Samples were collected on January 25, 2018 and received by the laboratory on January 25, 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.



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Date of Report: February 2, 2018  
 Samples Submitted: January 25, 2018  
 Laboratory Reference: 1801-271  
 Project: 0146970060

### NWTPH-Gx

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-10-012518</b>					
Laboratory ID:	01-271-01					
Gasoline	<b>270</b>	100	NWTPH-Gx	1-26-18	1-26-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	89	66-114				
<b>Client ID:</b>	<b>KMW-109-012518</b>					
Laboratory ID:	01-271-02					
Gasoline	<b>260</b>	100	NWTPH-Gx	1-26-18	1-26-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	82	66-114				
<b>Client ID:</b>	<b>Trip Blank</b>					
Laboratory ID:	01-271-03					
Gasoline	<b>ND</b>	100	NWTPH-Gx	1-26-18	1-26-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	93	66-114				
<b>Client ID:</b>	<b>KMW-04-012518</b>					
Laboratory ID:	01-271-04					
Gasoline	<b>9000</b>	100	NWTPH-Gx	1-26-18	1-26-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	79	66-114				



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 Laboratory Reference: 1801-271  
 Project: 0146970060

**NWTPH-Gx**  
**QUALITY CONTROL**

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB0126W1					
Gasoline	ND	100	NWTPH-Gx	1-26-18	1-26-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	94	66-114				
Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD RPD Limit Flags
<b>DUPLICATE</b>						
Laboratory ID:	01-271-01					
	ORIG	DUP				
Gasoline	265	256	NA	NA	NA	3 30
Surrogate:						
Fluorobenzene				89 81	66-114	



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Date of Report: February 2, 2018  
 Samples Submitted: January 25, 2018  
 Laboratory Reference: 1801-271  
 Project: 0146970060

### NWTPH-Dx

Matrix: Water  
 Units: mg/L (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-10-012518</b>					
<b>Laboratory ID:</b>	01-271-01					
Diesel Range Organics	<b>2.3</b>	0.26	NWTPH-Dx	1-29-18	1-29-18	
Lube Oil Range Organics	<b>ND</b>	0.41	NWTPH-Dx	1-29-18	1-29-18	

<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>
<i>o-Terphenyl</i>	83	50-150

**Client ID:** **KMW-109-012518**  
**Laboratory ID:** 01-271-02

Diesel Range Organics	<b>2.3</b>	0.26	NWTPH-Dx	1-29-18	1-29-18
Lube Oil Range Organics	<b>ND</b>	0.41	NWTPH-Dx	1-29-18	1-29-18
<i>Surrogate:</i>	<i>Percent Recovery</i>				
<i>o-Terphenyl</i>	85		<i>Control Limits</i>		

**Client ID:** **KMW-04-012518**  
**Laboratory ID:** 01-271-04

Diesel Range Organics	<b>1.7</b>	0.26	NWTPH-Dx	1-29-18	1-29-18
Lube Oil Range Organics	<b>ND</b>	0.41	NWTPH-Dx	1-29-18	1-29-18
<i>Surrogate:</i>	<i>Percent Recovery</i>				
<i>o-Terphenyl</i>	71		<i>Control Limits</i>		



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Date of Report: February 2, 2018  
 Samples Submitted: January 25, 2018  
 Laboratory Reference: 1801-271  
 Project: 0146970060

**NWTPH-Dx**  
**QUALITY CONTROL**

Matrix: Water  
 Units: mg/L (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB0129W1					
Diesel Range Organics	ND	0.25	NWTPH-Dx	1-29-18	1-29-18	
Lube Oil Range Organics	ND	0.40	NWTPH-Dx	1-29-18	1-29-18	
Surrogate: <i>o-Terphenyl</i>	Percent Recovery 93	Control Limits 50-150				

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>								
Laboratory ID:	01-271-01							
	ORIG	DUP						
Diesel Range Organics	2.30	2.22	NA	NA	NA	NA	4	NA
Lube Oil Range	ND	ND	NA	NA	NA	NA	NA	NA
Surrogate: <i>o-Terphenyl</i>				83	80	50-150		

**SPIKE BLANK**

Laboratory ID: SB0129W1

Diesel Fuel #2	0.784	1.00	NA	78	58-113	NA	NA
<i>Surrogate: <i>o-Terphenyl</i></i>							
				90	50-150		



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Date of Report: February 2, 2018  
 Samples Submitted: January 25, 2018  
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 Project: 0146970060

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

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-10-012518</b>					
<b>Laboratory ID:</b>	01-271-01					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloromethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Vinyl Chloride	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromomethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloroethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Acetone	ND	5.0	EPA 8260C	1-29-18	1-29-18	
Iodomethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Carbon Disulfide	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methylene Chloride	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Vinyl Acetate	ND	1.0	EPA 8260C	1-29-18	1-29-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Butanone	ND	5.0	EPA 8260C	1-29-18	1-29-18	
Bromochloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloroform	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Benzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Trichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Dibromomethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromodichloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	1-29-18	1-29-18	
Toluene	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	



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Date of Report: February 2, 2018  
 Samples Submitted: January 25, 2018  
 Laboratory Reference: 1801-271  
 Project: 0146970060

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

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-10-012518</b>					
<b>Laboratory ID:</b>	01-271-01					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Tetrachloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Hexanone	ND	2.0	EPA 8260C	1-29-18	1-29-18	
Dibromochloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Ethylbenzene	14	0.20	EPA 8260C	1-29-18	1-29-18	
m,p-Xylene	28	0.40	EPA 8260C	1-29-18	1-29-18	
o-Xylene	0.42	0.20	EPA 8260C	1-29-18	1-29-18	
Styrene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromoform	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Isopropylbenzene	3.8	0.20	EPA 8260C	1-29-18	1-29-18	
Bromobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
n-Propylbenzene	3.6	0.20	EPA 8260C	1-29-18	1-29-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3,5-Trimethylbenzene	0.90	0.20	EPA 8260C	1-29-18	1-29-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trimethylbenzene	2.7	0.20	EPA 8260C	1-29-18	1-29-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
n-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Hexachlorobutadiene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Naphthalene	ND	1.0	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
<b>Surrogate:</b>	<b>Percent Recovery</b>	<b>Control Limits</b>				
Dibromofluoromethane	93	75-127				
Toluene-d8	97	80-127				
4-Bromofluorobenzene	95	78-125				



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Date of Report: February 2, 2018  
 Samples Submitted: January 25, 2018  
 Laboratory Reference: 1801-271  
 Project: 0146970060

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

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-109-012518</b>					
<b>Laboratory ID:</b>	01-271-02					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloromethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Vinyl Chloride	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromomethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloroethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Acetone	ND	5.0	EPA 8260C	1-29-18	1-29-18	
Iodomethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Carbon Disulfide	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methylene Chloride	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Vinyl Acetate	ND	1.0	EPA 8260C	1-29-18	1-29-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Butanone	ND	5.0	EPA 8260C	1-29-18	1-29-18	
Bromochloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloroform	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Benzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Trichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Dibromomethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromodichloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	1-29-18	1-29-18	
Toluene	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	



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

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Date of Report: February 2, 2018  
 Samples Submitted: January 25, 2018  
 Laboratory Reference: 1801-271  
 Project: 0146970060

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

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-109-012518</b>					
<b>Laboratory ID:</b>	01-271-02					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Tetrachloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Hexanone	ND	2.0	EPA 8260C	1-29-18	1-29-18	
Dibromochloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Ethylbenzene	30	0.20	EPA 8260C	1-29-18	1-29-18	
m,p-Xylene	65	0.40	EPA 8260C	1-29-18	1-29-18	
o-Xylene	0.58	0.20	EPA 8260C	1-29-18	1-29-18	
Styrene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromoform	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Isopropylbenzene	6.5	0.20	EPA 8260C	1-29-18	1-29-18	
Bromobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
n-Propylbenzene	5.6	0.20	EPA 8260C	1-29-18	1-29-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3,5-Trimethylbenzene	1.7	0.20	EPA 8260C	1-29-18	1-29-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trimethylbenzene	5.4	0.20	EPA 8260C	1-29-18	1-29-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
n-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Hexachlorobutadiene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Naphthalene	ND	1.0	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
<b>Surrogate:</b>	<b>Percent Recovery</b>	<b>Control Limits</b>				
Dibromofluoromethane	101	75-127				
Toluene-d8	102	80-127				
4-Bromofluorobenzene	106	78-125				



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

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Date of Report: February 2, 2018  
 Samples Submitted: January 25, 2018  
 Laboratory Reference: 1801-271  
 Project: 0146970060

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

Matrix: Water

Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>Trip Blank</b>					
<b>Laboratory ID:</b>	01-271-03					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloromethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Vinyl Chloride	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromomethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloroethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Acetone	ND	5.0	EPA 8260C	1-29-18	1-29-18	
Iodomethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Carbon Disulfide	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methylene Chloride	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Vinyl Acetate	ND	1.0	EPA 8260C	1-29-18	1-29-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Butanone	ND	5.0	EPA 8260C	1-29-18	1-29-18	
Bromochloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloroform	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Benzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Trichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Dibromomethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromodichloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	1-29-18	1-29-18	
Toluene	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	



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

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Date of Report: February 2, 2018  
 Samples Submitted: January 25, 2018  
 Laboratory Reference: 1801-271  
 Project: 0146970060

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

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>Trip Blank</b>					
Laboratory ID:	01-271-03					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Tetrachloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Hexanone	ND	2.0	EPA 8260C	1-29-18	1-29-18	
Dibromochloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Ethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
m,p-Xylene	ND	0.40	EPA 8260C	1-29-18	1-29-18	
o-Xylene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Styrene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromoform	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Isopropylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
n-Propylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
n-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Hexachlorobutadiene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Naphthalene	ND	1.0	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
<b>Surrogate:</b>	<b>Percent Recovery</b>	<b>Control Limits</b>				
Dibromofluoromethane	100	75-127				
Toluene-d8	105	80-127				
4-Bromofluorobenzene	96	78-125				



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Date of Report: February 2, 2018  
 Samples Submitted: January 25, 2018  
 Laboratory Reference: 1801-271  
 Project: 0146970060

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

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-04-012518</b>					
<b>Laboratory ID:</b>	01-271-04					
Dichlorodifluoromethane	ND	50	EPA 8260C	1-29-18	1-29-18	
Chloromethane	ND	250	EPA 8260C	1-29-18	1-29-18	
Vinyl Chloride	ND	50	EPA 8260C	1-29-18	1-29-18	
Bromomethane	ND	50	EPA 8260C	1-29-18	1-29-18	
Chloroethane	ND	250	EPA 8260C	1-29-18	1-29-18	
Trichlorofluoromethane	ND	50	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethene	ND	50	EPA 8260C	1-29-18	1-29-18	
Acetone	ND	1300	EPA 8260C	1-29-18	1-29-18	
Iodomethane	ND	250	EPA 8260C	1-29-18	1-29-18	
Carbon Disulfide	ND	50	EPA 8260C	1-29-18	1-29-18	
Methylene Chloride	ND	250	EPA 8260C	1-29-18	1-29-18	
(trans) 1,2-Dichloroethene	ND	50	EPA 8260C	1-29-18	1-29-18	
Methyl t-Butyl Ether	ND	50	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethane	ND	50	EPA 8260C	1-29-18	1-29-18	
Vinyl Acetate	ND	250	EPA 8260C	1-29-18	1-29-18	
2,2-Dichloropropane	ND	50	EPA 8260C	1-29-18	1-29-18	
(cis) 1,2-Dichloroethene	ND	50	EPA 8260C	1-29-18	1-29-18	
2-Butanone	ND	1300	EPA 8260C	1-29-18	1-29-18	
Bromochloromethane	ND	50	EPA 8260C	1-29-18	1-29-18	
Chloroform	ND	50	EPA 8260C	1-29-18	1-29-18	
1,1,1-Trichloroethane	ND	50	EPA 8260C	1-29-18	1-29-18	
Carbon Tetrachloride	ND	50	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloropropene	ND	50	EPA 8260C	1-29-18	1-29-18	
Benzene	ND	50	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloroethane	ND	50	EPA 8260C	1-29-18	1-29-18	
Trichloroethene	ND	50	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloropropane	ND	50	EPA 8260C	1-29-18	1-29-18	
Dibromomethane	ND	50	EPA 8260C	1-29-18	1-29-18	
Bromodichloromethane	ND	50	EPA 8260C	1-29-18	1-29-18	
2-Chloroethyl Vinyl Ether	ND	250	EPA 8260C	1-29-18	1-29-18	
(cis) 1,3-Dichloropropene	ND	50	EPA 8260C	1-29-18	1-29-18	
Methyl Isobutyl Ketone	ND	500	EPA 8260C	1-29-18	1-29-18	
Toluene	5500	250	EPA 8260C	1-29-18	1-29-18	
(trans) 1,3-Dichloropropene	ND	50	EPA 8260C	1-29-18	1-29-18	



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Date of Report: February 2, 2018  
 Samples Submitted: January 25, 2018  
 Laboratory Reference: 1801-271  
 Project: 0146970060

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

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-04-012518</b>					
Laboratory ID:	01-271-04					
1,1,2-Trichloroethane	ND	50	EPA 8260C	1-29-18	1-29-18	
Tetrachloroethene	ND	50	EPA 8260C	1-29-18	1-29-18	
1,3-Dichloropropane	ND	50	EPA 8260C	1-29-18	1-29-18	
2-Hexanone	ND	500	EPA 8260C	1-29-18	1-29-18	
Dibromochloromethane	ND	50	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromoethane	ND	50	EPA 8260C	1-29-18	1-29-18	
Chlorobenzene	ND	50	EPA 8260C	1-29-18	1-29-18	
1,1,1,2-Tetrachloroethane	ND	50	EPA 8260C	1-29-18	1-29-18	
Ethylbenzene	4700	50	EPA 8260C	1-29-18	1-29-18	
m,p-Xylene	12000	100	EPA 8260C	1-29-18	1-29-18	
o-Xylene	3600	50	EPA 8260C	1-29-18	1-29-18	
Styrene	ND	50	EPA 8260C	1-29-18	1-29-18	
Bromoform	ND	250	EPA 8260C	1-29-18	1-29-18	
Isopropylbenzene	ND	50	EPA 8260C	1-29-18	1-29-18	
Bromobenzene	ND	50	EPA 8260C	1-29-18	1-29-18	
1,1,2,2-Tetrachloroethane	ND	50	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichloropropane	ND	50	EPA 8260C	1-29-18	1-29-18	
n-Propylbenzene	ND	50	EPA 8260C	1-29-18	1-29-18	
2-Chlorotoluene	ND	50	EPA 8260C	1-29-18	1-29-18	
4-Chlorotoluene	ND	50	EPA 8260C	1-29-18	1-29-18	
1,3,5-Trimethylbenzene	ND	50	EPA 8260C	1-29-18	1-29-18	
tert-Butylbenzene	ND	50	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trimethylbenzene	110	50	EPA 8260C	1-29-18	1-29-18	
sec-Butylbenzene	ND	50	EPA 8260C	1-29-18	1-29-18	
1,3-Dichlorobenzene	ND	50	EPA 8260C	1-29-18	1-29-18	
p-Isopropyltoluene	ND	50	EPA 8260C	1-29-18	1-29-18	
1,4-Dichlorobenzene	ND	50	EPA 8260C	1-29-18	1-29-18	
1,2-Dichlorobenzene	ND	50	EPA 8260C	1-29-18	1-29-18	
n-Butylbenzene	ND	50	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromo-3-chloropropane	ND	250	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trichlorobenzene	ND	50	EPA 8260C	1-29-18	1-29-18	
Hexachlorobutadiene	ND	50	EPA 8260C	1-29-18	1-29-18	
Naphthalene	ND	250	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichlorobenzene	ND	50	EPA 8260C	1-29-18	1-29-18	
<b>Surrogate:</b>	<b>Percent Recovery</b>	<b>Control Limits</b>				
Dibromofluoromethane	97	75-127				
Toluene-d8	100	80-127				
4-Bromofluorobenzene	101	78-125				



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

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Date of Report: February 2, 2018  
 Samples Submitted: January 25, 2018  
 Laboratory Reference: 1801-271  
 Project: 0146970060

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

Matrix: Water

Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0129W1					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloromethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Vinyl Chloride	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromomethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloroethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Acetone	ND	5.0	EPA 8260C	1-29-18	1-29-18	
Iodomethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Carbon Disulfide	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methylene Chloride	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Vinyl Acetate	ND	1.0	EPA 8260C	1-29-18	1-29-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Butanone	ND	5.0	EPA 8260C	1-29-18	1-29-18	
Bromochloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloroform	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Benzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Trichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Dibromomethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromodichloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	1-29-18	1-29-18	
Toluene	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	



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Date of Report: February 2, 2018  
 Samples Submitted: January 25, 2018  
 Laboratory Reference: 1801-271  
 Project: 0146970060

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

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0129W1					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Tetrachloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Hexanone	ND	2.0	EPA 8260C	1-29-18	1-29-18	
Dibromochloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Ethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
m,p-Xylene	ND	0.40	EPA 8260C	1-29-18	1-29-18	
o-Xylene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Styrene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromoform	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Isopropylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
n-Propylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
n-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Hexachlorobutadiene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Naphthalene	ND	1.0	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	102	75-127				
Toluene-d8	107	80-127				
4-Bromofluorobenzene	96	78-125				



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Date of Report: February 2, 2018  
 Samples Submitted: January 25, 2018  
 Laboratory Reference: 1801-271  
 Project: 0146970060

**VOLATILES by EPA 8260C**  
**MS/MSD QUALITY CONTROL**

Matrix: Water

Units: ug/L

Analyte	Result	Spike Level	Source	Percent	Recovery	Limits	RPD	RPD Limit	Flags						
			Result	Recovery	Recovery										
<b>MATRIX SPIKES</b>															
Laboratory ID: 01-271-01															
	MS	MSD	MS	MSD	MS	MSD									
1,1-Dichloroethene	<b>8.70</b>	<b>9.68</b>	10.0	10.0	ND	87	97	58-121	11						
Benzene	<b>9.39</b>	<b>10.4</b>	10.0	10.0	ND	94	104	67-122	10						
Trichloroethene	<b>8.37</b>	<b>9.52</b>	10.0	10.0	ND	84	95	60-120	13						
Toluene	<b>9.42</b>	<b>10.5</b>	10.0	10.0	ND	94	105	66-131	11						
Chlorobenzene	<b>8.95</b>	<b>9.76</b>	10.0	10.0	ND	90	98	73-120	9						
<i>Surrogate:</i>															
<i>Dibromofluoromethane</i>						94	97	75-127							
<i>Toluene-d8</i>						95	100	80-127							
<i>4-Bromofluorobenzene</i>						94	99	78-125							



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 Laboratory Reference: 1801-271  
 Project: 0146970060

**VOLATILES by EPA 8260C**  
**SPIKE BLANK QUALITY CONTROL**

Matrix: Water  
 Units: ug/L

Analyte	Result	Spike Level	Percent Recovery	Recovery Limits	Flags
<b>SPIKE BLANK</b>					
Laboratory ID:	SB0129W1				
1,1-Dichloroethene	<b>9.35</b>	10.0	94	63-126	
Benzene	<b>9.55</b>	10.0	96	78-122	
Trichloroethene	<b>9.18</b>	10.0	92	63-120	
Toluene	<b>10.0</b>	10.0	100	79-124	
Chlorobenzene	<b>9.60</b>	10.0	96	78-120	
<i>Surrogate:</i>					
<i>Dibromofluoromethane</i>			97	75-127	
<i>Toluene-d8</i>			102	80-127	
<i>4-Bromofluorobenzene</i>			94	78-125	



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 Laboratory Reference: 1801-271  
 Project: 0146970060

### PAHs EPA 8270D/SIM

Matrix: Water

Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-10-012518</b>					
<b>Laboratory ID:</b>	01-271-01					
Naphthalene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
2-Methylnaphthalene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
1-Methylnaphthalene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Acenaphthylene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Acenaphthene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Fluorene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Phenanthrene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Anthracene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Fluoranthene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Pyrene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[a]anthracene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Chrysene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[b]fluoranthene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo(j,k)fluoranthene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[a]pyrene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Indeno(1,2,3-c,d)pyrene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Dibenz[a,h]anthracene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[g,h,i]perylene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
2-Fluorobiphenyl	58		25 - 107			
Pyrene-d10	71		28 - 103			
Terphenyl-d14	96		36 - 129			



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Date of Report: February 2, 2018  
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 Laboratory Reference: 1801-271  
 Project: 0146970060

### PAHs EPA 8270D/SIM

Matrix: Water

Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-109-012518</b>					
<b>Laboratory ID:</b>	01-271-02					
Naphthalene	<b>0.15</b>	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
2-Methylnaphthalene	<b>ND</b>	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
1-Methylnaphthalene	<b>ND</b>	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Acenaphthylene	<b>ND</b>	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Acenaphthene	<b>ND</b>	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Fluorene	<b>ND</b>	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Phenanthrene	<b>ND</b>	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Anthracene	<b>0.13</b>	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Fluoranthene	<b>ND</b>	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Pyrene	<b>ND</b>	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[a]anthracene	<b>ND</b>	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Chrysene	<b>ND</b>	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[b]fluoranthene	<b>ND</b>	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo(j,k)fluoranthene	<b>ND</b>	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[a]pyrene	<b>ND</b>	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Indeno(1,2,3-c,d)pyrene	<b>ND</b>	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Dibenz[a,h]anthracene	<b>ND</b>	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[g,h,i]perylene	<b>ND</b>	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
2-Fluorobiphenyl	60		25 - 107			
Pyrene-d10	75		28 - 103			
Terphenyl-d14	106		36 - 129			



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Date of Report: February 2, 2018  
 Samples Submitted: January 25, 2018  
 Laboratory Reference: 1801-271  
 Project: 0146970060

### PAHs EPA 8270D/SIM

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-04-012518</b>					
<b>Laboratory ID:</b>	01-271-04					
Naphthalene	3.7	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
2-Methylnaphthalene	0.46	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
1-Methylnaphthalene	0.36	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Acenaphthylene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Acenaphthene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Fluorene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Phenanthrene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Anthracene	0.30	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Fluoranthene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Pyrene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Chrysene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
2-Fluorobiphenyl	54		25 - 107			
Pyrene-d10	65		28 - 103			
Terphenyl-d14	95		36 - 129			



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 Laboratory Reference: 1801-271  
 Project: 0146970060

**PAHs EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**

Matrix: Water

Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0130W1					
Naphthalene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
2-Methylnaphthalene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
1-Methylnaphthalene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Acenaphthylene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Acenaphthene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Fluorene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Phenanthrene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Anthracene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Fluoranthene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Pyrene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Chrysene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	56	25 - 107				
Pyrene-d10	75	28 - 103				
Terphenyl-d14	84	36 - 129				



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 Project: 0146970060

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

Matrix: Water

Units: ug/L

Analyte	Result	Spike Level	Source	Percent	Recovery	RPD	RPD Limit	Flags				
			Result	Recovery	Limits							
<b>MATRIX SPIKES</b>												
Laboratory ID:		01-271-01										
	MS	MSD	MS	MSD	MS	MSD						
Naphthalene	<b>0.421</b>	<b>0.469</b>	0.588	0.569	ND	72	82	27 - 106				
Acenaphthylene	<b>0.518</b>	<b>0.526</b>	0.588	0.569	ND	88	92	20 - 117				
Acenaphthene	<b>0.648</b>	<b>0.582</b>	0.588	0.569	ND	110	102	30 - 114				
Fluorene	<b>0.552</b>	<b>0.529</b>	0.588	0.569	ND	94	93	36 - 116				
Phenanthrene	<b>0.494</b>	<b>0.501</b>	0.588	0.569	ND	84	88	31 - 122				
Anthracene	<b>0.589</b>	<b>0.586</b>	0.588	0.569	ND	100	103	33 - 144				
Fluoranthene	<b>0.498</b>	<b>0.505</b>	0.588	0.569	ND	85	89	44 - 120				
Pyrene	<b>0.506</b>	<b>0.517</b>	0.588	0.569	ND	86	91	40 - 130				
Benzo[a]anthracene	<b>0.571</b>	<b>0.595</b>	0.588	0.569	ND	97	105	47 - 131				
Chrysene	<b>0.556</b>	<b>0.547</b>	0.588	0.569	ND	95	96	48 - 120				
Benzo[b]fluoranthene	<b>0.560</b>	<b>0.563</b>	0.588	0.569	ND	95	99	42 - 128				
Benzo(j,k)fluoranthene	<b>0.537</b>	<b>0.555</b>	0.588	0.569	ND	91	98	46 - 121				
Benzo[a]pyrene	<b>0.556</b>	<b>0.570</b>	0.588	0.569	ND	95	100	34 - 121				
Indeno(1,2,3-c,d)pyrene	<b>0.551</b>	<b>0.573</b>	0.588	0.569	ND	94	101	39 - 128				
Dibenz[a,h]anthracene	<b>0.557</b>	<b>0.572</b>	0.588	0.569	ND	95	101	39 - 125				
Benzo[g,h,i]perylene	<b>0.522</b>	<b>0.535</b>	0.588	0.569	ND	89	94	41 - 122				
<i>Surrogate:</i>												
<i>2-Fluorobiphenyl</i>					53	58	25 - 107					
<i>Pyrene-d10</i>					71	74	28 - 103					
<i>Terphenyl-d14</i>					98	101	36 - 129					



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

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Date of Report: February 2, 2018  
 Samples Submitted: January 25, 2018  
 Laboratory Reference: 1801-271  
 Project: 0146970060

**PAHs EPA 8270D/SIM**  
**SPIKE BLANK QUALITY CONTROL**

Matrix: Water  
 Units: ug/L

Analyte	Result	Spike Level	Percent Recovery	Recovery Limits	Flags
<b>SPIKE BLANK</b>					
Laboratory ID:	SB0130W1				
	SB	SB	SB		
Naphthalene	<b>0.360</b>	0.500	72	27 - 106	
Acenaphthylene	<b>0.387</b>	0.500	77	20 - 117	
Acenaphthene	<b>0.461</b>	0.500	92	30 - 114	
Fluorene	<b>0.454</b>	0.500	91	36 - 116	
Phenanthrene	<b>0.458</b>	0.500	92	31 - 122	
Anthracene	<b>0.480</b>	0.500	96	33 - 144	
Fluoranthene	<b>0.488</b>	0.500	98	44 - 120	
Pyrene	<b>0.566</b>	0.500	113	40 - 130	
Benzo[a]anthracene	<b>0.503</b>	0.500	101	47 - 131	
Chrysene	<b>0.498</b>	0.500	100	48 - 120	
Benzo[b]fluoranthene	<b>0.491</b>	0.500	98	42 - 128	
Benzo(j,k)fluoranthene	<b>0.489</b>	0.500	98	46 - 121	
Benzo[a]pyrene	<b>0.502</b>	0.500	100	34 - 121	
Indeno(1,2,3-c,d)pyrene	<b>0.522</b>	0.500	104	39 - 128	
Dibenz[a,h]anthracene	<b>0.528</b>	0.500	106	39 - 125	
Benzo[g,h,i]perylene	<b>0.518</b>	0.500	104	41 - 122	
<i>Surrogate:</i>					
<i>2-Fluorobiphenyl</i>			83	25 - 107	
<i>Pyrene-d10</i>			85	28 - 103	
<i>Terphenyl-d14</i>			103	36 - 129	



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 Laboratory Reference: 1801-271  
 Project: 0146970060

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

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	EPA Method	Date Prepared	Date Analyzed	Flags
Lab ID:	01-261-01					
<b>Client ID:</b>	<b>KMW-06-012418</b>					
Arsenic	<b>ND</b>	3.3	200.8	1-31-18	1-31-18	
Barium	<b>ND</b>	28	200.8	1-31-18	1-31-18	
Cadmium	<b>ND</b>	4.4	200.8	1-31-18	1-31-18	
Chromium	<b>ND</b>	11	200.8	1-31-18	1-31-18	
Lead	<b>7.3</b>	1.1	200.8	1-31-18	1-31-18	
Mercury	<b>ND</b>	0.50	7470A	1-30-18	1-30-18	
Selenium	<b>ND</b>	5.6	200.8	1-31-18	1-31-18	
Silver	<b>ND</b>	11	200.8	1-31-18	1-31-18	

Lab ID: 01-261-02  
**Client ID:** KMW-09-012418

Arsenic	<b>ND</b>	3.3	200.8	1-31-18	1-31-18
Barium	<b>ND</b>	28	200.8	1-31-18	1-31-18
Cadmium	<b>ND</b>	4.4	200.8	1-31-18	1-31-18
Chromium	<b>ND</b>	11	200.8	1-31-18	1-31-18
Lead	<b>3.0</b>	1.1	200.8	1-31-18	1-31-18
Mercury	<b>ND</b>	0.50	7470A	1-30-18	1-30-18
Selenium	<b>ND</b>	5.6	200.8	1-31-18	1-31-18
Silver	<b>ND</b>	11	200.8	1-31-18	1-31-18



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 Laboratory Reference: 1801-271  
 Project: 0146970060

**TOTAL METALS**  
**EPA 200.8/7470A**  
**METHOD BLANK QUALITY CONTROL**

Date Extracted: 1-30&31-18  
 Date Analyzed: 1-30&31-18  
 Matrix: Water  
 Units: ug/L (ppb)  
 Lab ID: MB0130W1&MB0131WM1

Analyte	Method	Result	PQL
Arsenic	200.8	<b>ND</b>	3.3
Barium	200.8	<b>ND</b>	28
Cadmium	200.8	<b>ND</b>	4.4
Chromium	200.8	<b>ND</b>	11
Lead	200.8	<b>ND</b>	1.1
Mercury	7470A	<b>ND</b>	0.50
Selenium	200.8	<b>ND</b>	5.6
Silver	200.8	<b>ND</b>	11



Date of Report: February 2, 2018  
 Samples Submitted: January 25, 2018  
 Laboratory Reference: 1801-271  
 Project: 0146970060

**TOTAL METALS  
EPA 200.8/7470A  
DUPLICATE QUALITY CONTROL**

Date Extracted: 1-30&31-18  
 Date Analyzed: 1-30&31-18

Matrix: Water  
 Units: ug/L (ppb)

Lab ID: 01-271-01

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Arsenic	<b>6.78</b>	<b>6.33</b>	7	3.3	
Barium	<b>ND</b>	<b>ND</b>	NA	28	
Cadmium	<b>ND</b>	<b>ND</b>	NA	4.4	
Chromium	<b>ND</b>	<b>ND</b>	NA	11	
Lead	<b>1.12</b>	<b>ND</b>	NA	1.1	
Mercury	<b>ND</b>	<b>ND</b>	NA	0.50	
Selenium	<b>ND</b>	<b>ND</b>	NA	5.6	
Silver	<b>ND</b>	<b>ND</b>	NA	11	



Date of Report: February 2, 2018  
 Samples Submitted: January 25, 2018  
 Laboratory Reference: 1801-271  
 Project: 0146970060

**TOTAL METALS**  
**EPA 200.8/7470A**  
**MS/MSD QUALITY CONTROL**

Date Extracted: 1-30&31-18  
 Date Analyzed: 1-30&31-18

Matrix: Water  
 Units: ug/L (ppb)

Lab ID: 01-271-01

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Arsenic	222	<b>227</b>	99	<b>233</b>	102	3	
Barium	222	<b>224</b>	101	<b>222</b>	100	1	
Cadmium	222	<b>228</b>	103	<b>231</b>	104	1	
Chromium	222	<b>211</b>	95	<b>216</b>	97	2	
Lead	222	<b>207</b>	93	<b>207</b>	93	0	
Mercury	12.5	<b>12.3</b>	98	<b>12.4</b>	99	1	
Selenium	222	<b>231</b>	104	<b>232</b>	104	0	
Silver	222	<b>196</b>	88	<b>202</b>	91	3	



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 Laboratory Reference: 1801-271  
 Project: 0146970060

**TOTAL METALS**  
**EPA 200.8/7470A**  
**SPIKE BLANK QUALITY CONTROL**

Date Extracted: 1-30&31-18  
 Date Analyzed: 1-30&31-18  
 Matrix: Water  
 Units: ug/L (ppb)  
 Lab ID: SB0130W1&SB0131WM1

Analyte	Method	Spike Level	SB Result	Percent Recovery
Arsenic	200.8	222	<b>228</b>	103
Barium	200.8	222	<b>212</b>	96
Cadmium	200.8	222	<b>224</b>	101
Chromium	200.8	222	<b>210</b>	95
Lead	200.8	222	<b>217</b>	98
Mercury	7470A	12.5	<b>12.6</b>	101
Selenium	200.8	222	<b>240</b>	108
Silver	200.8	222	<b>207</b>	93





### Data Qualifiers and Abbreviations

- A - Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
- B - The analyte indicated was also found in the blank sample.
- C - The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
- E - The value reported exceeds the quantitation range and is an estimate.
- F - Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
- H - The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
- I - Compound recovery is outside of the control limits.
- J - The value reported was below the practical quantitation limit. The value is an estimate.
- K - Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
- L - The RPD is outside of the control limits.
- M - Hydrocarbons in the gasoline range are impacting the diesel range result.
- M1 - Hydrocarbons in the gasoline range (toluene-naphthalene) are present in the sample.
- N - Hydrocarbons in the lube oil range are impacting the diesel range result.
- N1 - Hydrocarbons in diesel range are impacting lube oil range results.
- O - Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
- P - The RPD of the detected concentrations between the two columns is greater than 40.
- Q - Surrogate recovery is outside of the control limits.
- S - Surrogate recovery data is not available due to the necessary dilution of the sample.
- T - The sample chromatogram is not similar to a typical \_\_\_\_\_.
- U - The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- U1 - The practical quantitation limit is elevated due to interferences present in the sample.
- V - Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
- W - Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
- X - Sample extract treated with a mercury cleanup procedure.
- X1 - Sample extract treated with a sulfuric acid/silica gel cleanup procedure.
- Y - The calibration verification for this analyte exceeded the 20% drift specified in method 8260C, and therefore the reported result should be considered an estimate. The overall performance of the calibration verification standard met the acceptance criteria of the method.
- Z -
- ND - Not Detected at PQL
- PQL - Practical Quantitation Limit
- RPD - Relative Percent Difference





**OnSite  
Environmental Inc.**

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

## Chain of Custody

Page 1 of 1

Turnaround Request (in working days)				Laboratory Number: 01-271		
(Check One)						
<input type="checkbox"/> Same Day <input type="checkbox"/> 1 Day <input checked="" type="checkbox"/> 2 Days <input type="checkbox"/> 3 Days <input checked="" type="checkbox"/> Standard (7 Days) (TPH analysis 5 Days)						
<input type="checkbox"/> (other) _____						
Lab ID	Sample Identification	Date Sampled	Time Sampled	Matrix	Number of Containers	
1	KMW-10-012518 **	1/25/18	0930	W	19	NWTPH-HCID
2	KMW-104-012518	1/25/18	1030	W	10	NWTPH-Gx/BTEX
3	trip Blank	-	-	W	3	NWTPH-Gx
4	KMW-04-012518	1/25/18	1215	W	10	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 80
						Total MTCA Metals 80
						TCLP Metals
						HEM (oil and grease) 1664A
						% Moisture
Signature	Company	Date	Time	Comments/Special Instructions		
Relinquished Received	Anne Feser Wheeler's Analytical Services Inc.	1/25/18	1340	** Extra volume for NO/MOD		
Relinquished Received	Ames FW Office of Crustal Geology (William Young) Anne Feser Wheeler's Analytical Services Inc.	1/25/18 1/20/18	1511 1521			
Relinquished Received	John Rose	1/25/18	1602			
Reviewed/Date				Data Package: Standard <input type="checkbox"/> Level III <input type="checkbox"/> Level IV <input type="checkbox"/>		
Reviewed/Date				Chromatograms with final report <input type="checkbox"/> Electronic Data Deliverables (EDDs) <input type="checkbox"/>		



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

February 2, 2018

Kathleen Goodman  
AMEC Foster Wheeler  
Environment and Infrastructure, Inc.  
One Union Square  
600 University Street, Suite 600  
Seattle, WA 98101

Re: Analytical Data for Project 0146970060  
Laboratory Reference No. 1801-282

Dear Kathleen:

Enclosed are the analytical results and associated quality control data for samples submitted on January 26, 2018.

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

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

Sincerely,

A handwritten signature in black ink, appearing to read "DB" followed by a cursive surname.

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,  
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Date of Report: February 2, 2018  
Samples Submitted: January 26, 2018  
Laboratory Reference: 1801-282  
Project: 0146970060

### Case Narrative

Samples were collected on January 26, 2018 and received by the laboratory on January 26, 2018. They were maintained at the laboratory at a temperature of 2°C to 6°C.

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

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



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 Project: 0146970060

### NWTPH-Gx

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-03R-012618</b>					
Laboratory ID:	01-282-01					
Gasoline	<b>150</b>	100	NWTPH-Gx	1-29-18	1-29-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	78	66-114				
<b>Client ID:</b>	<b>KMW-08-012618</b>					
Laboratory ID:	01-282-02					
Gasoline	<b>120</b>	100	NWTPH-Gx	1-29-18	1-29-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	72	66-114				
<b>Client ID:</b>	<b>KMW-07-012618</b>					
Laboratory ID:	01-282-03					
Gasoline	<b>ND</b>	100	NWTPH-Gx	1-29-18	1-29-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	92	66-114				
<b>Client ID:</b>	<b>KMW-02R-012618</b>					
Laboratory ID:	01-282-04					
Gasoline	<b>ND</b>	100	NWTPH-Gx	1-29-18	1-29-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	77	66-114				
<b>Client ID:</b>	<b>Trip Blanks</b>					
Laboratory ID:	01-282-05					
Gasoline	<b>ND</b>	100	NWTPH-Gx	1-29-18	1-29-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	77	66-114				



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 Project: 0146970060

**NWTPH-Gx**  
**QUALITY CONTROL**

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB0129W2					
Gasoline	ND	100	NWTPH-Gx	1-29-18	1-29-18	
Surrogate:	Percent Recovery	Control Limits				
Fluorobenzene	92	66-114				
Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD RPD Limit Flags
<b>DUPPLICATE</b>						
Laboratory ID:	01-282-03					
	ORIG	DUP				
Gasoline	ND	ND	NA	NA	NA	NA 30
Surrogate:						
Fluorobenzene				92	92	66-114



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 Laboratory Reference: 1801-282  
 Project: 0146970060

### NWTPH-Dx

Matrix: Water  
 Units: mg/L (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-03R-012618</b>					
Laboratory ID:	01-282-01					
Diesel Range Organics	<b>0.51</b>	0.26	NWTPH-Dx	1-29-18	1-29-18	
Lube Oil Range Organics	<b>ND</b>	0.41	NWTPH-Dx	1-29-18	1-29-18	
Surrogate: <i>o-Terphenyl</i>	Percent Recovery 80	Control Limits 50-150				
<b>Client ID:</b>	<b>KMW-08-012618</b>					
Laboratory ID:	01-282-02					
Diesel Range Organics	<b>0.45</b>	0.25	NWTPH-Dx	1-29-18	1-29-18	
Lube Oil Range Organics	<b>ND</b>	0.41	NWTPH-Dx	1-29-18	1-29-18	
Surrogate: <i>o-Terphenyl</i>	Percent Recovery 98	Control Limits 50-150				
<b>Client ID:</b>	<b>KMW-07-012618</b>					
Laboratory ID:	01-282-03					
Diesel Range Organics	<b>ND</b>	0.26	NWTPH-Dx	1-29-18	1-29-18	
Lube Oil Range Organics	<b>ND</b>	0.41	NWTPH-Dx	1-29-18	1-29-18	
Surrogate: <i>o-Terphenyl</i>	Percent Recovery 94	Control Limits 50-150				
<b>Client ID:</b>	<b>KMW-02R-012618</b>					
Laboratory ID:	01-282-04					
Diesel Range Organics	<b>ND</b>	0.26	NWTPH-Dx	1-29-18	1-29-18	
Lube Oil Range Organics	<b>ND</b>	0.41	NWTPH-Dx	1-29-18	1-29-18	
Surrogate: <i>o-Terphenyl</i>	Percent Recovery 95	Control Limits 50-150				



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 Project: 0146970060

**NWTPH-Dx**  
**QUALITY CONTROL**

Matrix: Water  
 Units: mg/L (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB0129W1					
Diesel Range Organics	ND	0.25	NWTPH-Dx	1-29-18	1-29-18	
Lube Oil Range Organics	ND	0.40	NWTPH-Dx	1-29-18	1-29-18	
Surrogate: <i>o-Terphenyl</i>	Percent Recovery 93	Control Limits 50-150				

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>								
Laboratory ID:	01-271-01							
	ORIG	DUP						
Diesel Range Organics	2.30	2.22	NA	NA	NA	NA	4	NA
Lube Oil Range	ND	ND	NA	NA	NA	NA	NA	NA
Surrogate: <i>o-Terphenyl</i>				83	80	50-150		

**SPIKE BLANK**

Laboratory ID: SB0129W1

Diesel Fuel #2	0.784	1.00	NA	78	58-113	NA	NA
<i>Surrogate: <i>o-Terphenyl</i></i>							
				90	50-150		



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 Project: 0146970060

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

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-03R-012618</b>					
<b>Laboratory ID:</b>	01-282-01					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloromethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Vinyl Chloride	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromomethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloroethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Acetone	ND	5.0	EPA 8260C	1-29-18	1-29-18	
Iodomethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Carbon Disulfide	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methylene Chloride	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Vinyl Acetate	ND	1.0	EPA 8260C	1-29-18	1-29-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Butanone	ND	5.0	EPA 8260C	1-29-18	1-29-18	
Bromochloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloroform	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Benzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Trichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Dibromomethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromodichloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	1-29-18	1-29-18	
Toluene	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	



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

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Date of Report: February 2, 2018  
 Samples Submitted: January 26, 2018  
 Laboratory Reference: 1801-282  
 Project: 0146970060

**VOLATILES EPA 8260C**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-03R-012618</b>					
<b>Laboratory ID:</b>	01-282-01					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Tetrachloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Hexanone	ND	2.0	EPA 8260C	1-29-18	1-29-18	
Dibromochloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Ethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
m,p-Xylene	ND	0.40	EPA 8260C	1-29-18	1-29-18	
o-Xylene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Styrene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromoform	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Isopropylbenzene	7.2	0.20	EPA 8260C	1-29-18	1-29-18	
Bromobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
n-Propylbenzene	12	0.20	EPA 8260C	1-29-18	1-29-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
tert-Butylbenzene	0.22	0.20	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
sec-Butylbenzene	1.7	0.20	EPA 8260C	1-29-18	1-29-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
n-Butylbenzene	0.50	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Hexachlorobutadiene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Naphthalene	ND	1.0	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
<b>Surrogate:</b>	<b>Percent Recovery</b>	<b>Control Limits</b>				
Dibromofluoromethane	100	75-127				
Toluene-d8	100	80-127				
4-Bromofluorobenzene	100	78-125				



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

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Date of Report: February 2, 2018  
 Samples Submitted: January 26, 2018  
 Laboratory Reference: 1801-282  
 Project: 0146970060

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

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-08-012618</b>					
<b>Laboratory ID:</b>	01-282-02					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloromethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Vinyl Chloride	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromomethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloroethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Acetone	ND	5.0	EPA 8260C	1-29-18	1-29-18	
Iodomethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Carbon Disulfide	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methylene Chloride	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Vinyl Acetate	ND	1.0	EPA 8260C	1-29-18	1-29-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Butanone	ND	5.0	EPA 8260C	1-29-18	1-29-18	
Bromochloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloroform	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Benzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Trichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Dibromomethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromodichloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	1-29-18	1-29-18	
Toluene	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	



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Date of Report: February 2, 2018  
 Samples Submitted: January 26, 2018  
 Laboratory Reference: 1801-282  
 Project: 0146970060

**VOLATILES EPA 8260C**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-08-012618</b>					
<b>Laboratory ID:</b>	01-282-02					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Tetrachloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Hexanone	ND	2.0	EPA 8260C	1-29-18	1-29-18	
Dibromochloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Ethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
m,p-Xylene	ND	0.40	EPA 8260C	1-29-18	1-29-18	
o-Xylene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Styrene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromoform	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Isopropylbenzene	0.95	0.20	EPA 8260C	1-29-18	1-29-18	
Bromobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
n-Propylbenzene	0.85	0.20	EPA 8260C	1-29-18	1-29-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
tert-Butylbenzene	0.29	0.20	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
sec-Butylbenzene	2.3	0.20	EPA 8260C	1-29-18	1-29-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
n-Butylbenzene	0.39	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Hexachlorobutadiene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Naphthalene	ND	1.0	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
<b>Surrogate:</b>	<b>Percent Recovery</b>	<b>Control Limits</b>				
Dibromofluoromethane	100	75-127				
Toluene-d8	101	80-127				
4-Bromofluorobenzene	101	78-125				



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Date of Report: February 2, 2018  
 Samples Submitted: January 26, 2018  
 Laboratory Reference: 1801-282  
 Project: 0146970060

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

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-07-012618</b>					
<b>Laboratory ID:</b>	01-282-03					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloromethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Vinyl Chloride	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromomethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloroethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Acetone	ND	5.0	EPA 8260C	1-29-18	1-29-18	
Iodomethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Carbon Disulfide	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methylene Chloride	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Vinyl Acetate	ND	1.0	EPA 8260C	1-29-18	1-29-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Butanone	ND	5.0	EPA 8260C	1-29-18	1-29-18	
Bromochloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloroform	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Benzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Trichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Dibromomethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromodichloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	1-29-18	1-29-18	
Toluene	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	



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Date of Report: February 2, 2018  
 Samples Submitted: January 26, 2018  
 Laboratory Reference: 1801-282  
 Project: 0146970060

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

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-07-012618</b>					
<b>Laboratory ID:</b>	01-282-03					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Tetrachloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Hexanone	ND	2.0	EPA 8260C	1-29-18	1-29-18	
Dibromochloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Ethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
m,p-Xylene	ND	0.40	EPA 8260C	1-29-18	1-29-18	
o-Xylene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Styrene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromoform	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Isopropylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
n-Propylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
n-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Hexachlorobutadiene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Naphthalene	ND	1.0	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
<b>Surrogate:</b>	<b>Percent Recovery</b>	<b>Control Limits</b>				
Dibromofluoromethane	101	75-127				
Toluene-d8	103	80-127				
4-Bromofluorobenzene	98	78-125				



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Date of Report: February 2, 2018  
 Samples Submitted: January 26, 2018  
 Laboratory Reference: 1801-282  
 Project: 0146970060

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

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-02R-012618</b>					
<b>Laboratory ID:</b>	01-282-04					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloromethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Vinyl Chloride	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromomethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloroethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Acetone	ND	5.0	EPA 8260C	1-29-18	1-29-18	
Iodomethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Carbon Disulfide	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methylene Chloride	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Vinyl Acetate	ND	1.0	EPA 8260C	1-29-18	1-29-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Butanone	ND	5.0	EPA 8260C	1-29-18	1-29-18	
Bromochloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloroform	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Benzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Trichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Dibromomethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromodichloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	1-29-18	1-29-18	
Toluene	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	



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

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Date of Report: February 2, 2018  
 Samples Submitted: January 26, 2018  
 Laboratory Reference: 1801-282  
 Project: 0146970060

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

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-02R-012618</b>					
<b>Laboratory ID:</b>	01-282-04					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Tetrachloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Hexanone	ND	2.0	EPA 8260C	1-29-18	1-29-18	
Dibromochloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Ethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
m,p-Xylene	ND	0.40	EPA 8260C	1-29-18	1-29-18	
o-Xylene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Styrene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromoform	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Isopropylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
n-Propylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
n-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Hexachlorobutadiene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Naphthalene	ND	1.0	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
<b>Surrogate:</b>	<b>Percent Recovery</b>	<b>Control Limits</b>				
Dibromofluoromethane	102	75-127				
Toluene-d8	103	80-127				
4-Bromofluorobenzene	100	78-125				



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

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Date of Report: February 2, 2018  
 Samples Submitted: January 26, 2018  
 Laboratory Reference: 1801-282  
 Project: 0146970060

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

Matrix: Water

Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>Trip Blanks</b>					
<b>Laboratory ID:</b>	01-282-05					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloromethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Vinyl Chloride	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromomethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloroethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Acetone	ND	5.0	EPA 8260C	1-29-18	1-29-18	
Iodomethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Carbon Disulfide	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methylene Chloride	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Vinyl Acetate	ND	1.0	EPA 8260C	1-29-18	1-29-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Butanone	ND	5.0	EPA 8260C	1-29-18	1-29-18	
Bromochloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloroform	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Benzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Trichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Dibromomethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromodichloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	1-29-18	1-29-18	
Toluene	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	



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Date of Report: February 2, 2018  
 Samples Submitted: January 26, 2018  
 Laboratory Reference: 1801-282  
 Project: 0146970060

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

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>Trip Blanks</b>					
Laboratory ID:	01-282-05					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Tetrachloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Hexanone	ND	2.0	EPA 8260C	1-29-18	1-29-18	
Dibromochloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Ethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
m,p-Xylene	ND	0.40	EPA 8260C	1-29-18	1-29-18	
o-Xylene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Styrene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromoform	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Isopropylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
n-Propylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
n-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Hexachlorobutadiene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Naphthalene	ND	1.0	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
<b>Surrogate:</b>	<b>Percent Recovery</b>	<b>Control Limits</b>				
Dibromofluoromethane	100	75-127				
Toluene-d8	104	80-127				
4-Bromofluorobenzene	95	78-125				



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Date of Report: February 2, 2018  
 Samples Submitted: January 26, 2018  
 Laboratory Reference: 1801-282  
 Project: 0146970060

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

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0129W1					
Dichlorodifluoromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloromethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Vinyl Chloride	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromomethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloroethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Trichlorofluoromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Acetone	ND	5.0	EPA 8260C	1-29-18	1-29-18	
Iodomethane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Carbon Disulfide	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methylene Chloride	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methyl t-Butyl Ether	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Vinyl Acetate	ND	1.0	EPA 8260C	1-29-18	1-29-18	
2,2-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Butanone	ND	5.0	EPA 8260C	1-29-18	1-29-18	
Bromochloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chloroform	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,1-Trichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Carbon Tetrachloride	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Benzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Trichloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Dibromomethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromodichloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260C	1-29-18	1-29-18	
Toluene	ND	1.0	EPA 8260C	1-29-18	1-29-18	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260C	1-29-18	1-29-18	



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Date of Report: February 2, 2018  
 Samples Submitted: January 26, 2018  
 Laboratory Reference: 1801-282  
 Project: 0146970060

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

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0129W1					
1,1,2-Trichloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Tetrachloroethene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3-Dichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Hexanone	ND	2.0	EPA 8260C	1-29-18	1-29-18	
Dibromochloromethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromoethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Chlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Ethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
m,p-Xylene	ND	0.40	EPA 8260C	1-29-18	1-29-18	
o-Xylene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Styrene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromoform	ND	1.0	EPA 8260C	1-29-18	1-29-18	
Isopropylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Bromobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichloropropane	ND	0.20	EPA 8260C	1-29-18	1-29-18	
n-Propylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
2-Chlorotoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
4-Chlorotoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
tert-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
sec-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,3-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
p-Isopropyltoluene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,4-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
n-Butylbenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260C	1-29-18	1-29-18	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Hexachlorobutadiene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Naphthalene	ND	1.0	EPA 8260C	1-29-18	1-29-18	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260C	1-29-18	1-29-18	
Surrogate:	Percent Recovery	Control Limits				
Dibromofluoromethane	102	75-127				
Toluene-d8	107	80-127				
4-Bromofluorobenzene	96	78-125				



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: February 2, 2018  
 Samples Submitted: January 26, 2018  
 Laboratory Reference: 1801-282  
 Project: 0146970060

**VOLATILES by EPA 8260C**  
**MS/MSD QUALITY CONTROL**

Matrix: Water

Units: ug/L

Analyte	Result	Spike Level	Source	Percent	Recovery	RPD	RPD	Limit	Flags
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**MATRIX SPIKES**

Laboratory ID: 01-271-01

	MS	MSD	MS	MSD	MS	MSD			
1,1-Dichloroethene	<b>8.70</b>	<b>9.68</b>	10.0	10.0	ND	87	97	58-121	11
Benzene	<b>9.39</b>	<b>10.4</b>	10.0	10.0	ND	94	104	67-122	10
Trichloroethene	<b>8.37</b>	<b>9.52</b>	10.0	10.0	ND	84	95	60-120	13
Toluene	<b>9.42</b>	<b>10.5</b>	10.0	10.0	ND	94	105	66-131	11
Chlorobenzene	<b>8.95</b>	<b>9.76</b>	10.0	10.0	ND	90	98	73-120	9

*Surrogate:*

*Dibromofluoromethane* 94 97 75-127

*Toluene-d8* 95 100 80-127

*4-Bromofluorobenzene* 94 99 78-125



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 Samples Submitted: January 26, 2018  
 Laboratory Reference: 1801-282  
 Project: 0146970060

**VOLATILES by EPA 8260C**  
**SPIKE BLANK QUALITY CONTROL**

Matrix: Water  
 Units: ug/L

Analyte	Result	Spike Level	Percent Recovery	Recovery Limits	Flags
<b>SPIKE BLANK</b>					
Laboratory ID: SB0129W1					
1,1-Dichloroethene	<b>9.35</b>	10.0	94	63-126	
Benzene	<b>9.55</b>	10.0	96	78-122	
Trichloroethene	<b>9.18</b>	10.0	92	63-120	
Toluene	<b>10.0</b>	10.0	100	79-124	
Chlorobenzene	<b>9.60</b>	10.0	96	78-120	
<i>Surrogate:</i>					
<i>Dibromofluoromethane</i>			97	75-127	
<i>Toluene-d8</i>			102	80-127	
<i>4-Bromofluorobenzene</i>			94	78-125	



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Date of Report: February 2, 2018  
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 Laboratory Reference: 1801-282  
 Project: 0146970060

### PAHs EPA 8270D/SIM

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-03R-012618</b>					
<b>Laboratory ID:</b>	01-282-01					
Naphthalene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
2-Methylnaphthalene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
1-Methylnaphthalene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Acenaphthylene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Acenaphthene	0.22	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Fluorene	0.18	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Phenanthrene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Anthracene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Fluoranthene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Pyrene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[a]anthracene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Chrysene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[b]fluoranthene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo(j,k)fluoranthene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[a]pyrene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Indeno(1,2,3-c,d)pyrene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Dibenz[a,h]anthracene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[g,h,i]perylene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
2-Fluorobiphenyl	57		25 - 107			
Pyrene-d10	83		28 - 103			
Terphenyl-d14	96		36 - 129			



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Date of Report: February 2, 2018  
 Samples Submitted: January 26, 2018  
 Laboratory Reference: 1801-282  
 Project: 0146970060

### PAHs EPA 8270D/SIM

Matrix: Water

Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-08-012618</b>					
<b>Laboratory ID:</b>	01-282-02					
Naphthalene	ND	0.12	EPA 8270D/SIM	1-30-18	1-30-18	
2-Methylnaphthalene	ND	0.12	EPA 8270D/SIM	1-30-18	1-30-18	
1-Methylnaphthalene	ND	0.12	EPA 8270D/SIM	1-30-18	1-30-18	
Acenaphthylene	ND	0.12	EPA 8270D/SIM	1-30-18	1-30-18	
Acenaphthene	0.29	0.12	EPA 8270D/SIM	1-30-18	1-30-18	
Fluorene	0.33	0.12	EPA 8270D/SIM	1-30-18	1-30-18	
Phenanthrene	ND	0.12	EPA 8270D/SIM	1-30-18	1-30-18	
Anthracene	0.13	0.12	EPA 8270D/SIM	1-30-18	1-30-18	
Fluoranthene	0.20	0.12	EPA 8270D/SIM	1-30-18	1-30-18	
Pyrene	0.22	0.12	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[a]anthracene	0.018	0.012	EPA 8270D/SIM	1-30-18	1-30-18	
Chrysene	0.028	0.012	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[b]fluoranthene	0.018	0.012	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo(j,k)fluoranthene	ND	0.012	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[a]pyrene	0.015	0.012	EPA 8270D/SIM	1-30-18	1-30-18	
Indeno(1,2,3-c,d)pyrene	ND	0.012	EPA 8270D/SIM	1-30-18	1-30-18	
Dibenz[a,h]anthracene	ND	0.012	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[g,h,i]perylene	ND	0.012	EPA 8270D/SIM	1-30-18	1-30-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
2-Fluorobiphenyl	51		25 - 107			
Pyrene-d10	73		28 - 103			
Terphenyl-d14	84		36 - 129			



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Date of Report: February 2, 2018  
 Samples Submitted: January 26, 2018  
 Laboratory Reference: 1801-282  
 Project: 0146970060

### PAHs EPA 8270D/SIM

Matrix: Water

Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-07-012618</b>					
<b>Laboratory ID:</b>	01-282-03					
Naphthalene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
2-Methylnaphthalene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
1-Methylnaphthalene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Acenaphthylene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Acenaphthene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Fluorene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Phenanthrene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Anthracene	0.11	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Fluoranthene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Pyrene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[a]anthracene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Chrysene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[b]fluoranthene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo(j,k)fluoranthene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[a]pyrene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Indeno(1,2,3-c,d)pyrene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Dibenz[a,h]anthracene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[g,h,i]perylene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
2-Fluorobiphenyl	66		25 - 107			
Pyrene-d10	83		28 - 103			
Terphenyl-d14	93		36 - 129			



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 Samples Submitted: January 26, 2018  
 Laboratory Reference: 1801-282  
 Project: 0146970060

### PAHs EPA 8270D/SIM

Matrix: Water

Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>KMW-02R-012618</b>					
<b>Laboratory ID:</b>	01-282-04					
Naphthalene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
2-Methylnaphthalene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
1-Methylnaphthalene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Acenaphthylene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Acenaphthene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Fluorene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Phenanthrene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Anthracene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Fluoranthene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Pyrene	ND	0.11	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[a]anthracene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Chrysene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[b]fluoranthene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo(j,k)fluoranthene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[a]pyrene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Indeno(1,2,3-c,d)pyrene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Dibenz[a,h]anthracene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[g,h,i]perylene	ND	0.011	EPA 8270D/SIM	1-30-18	1-30-18	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
2-Fluorobiphenyl	64		25 - 107			
Pyrene-d10	76		28 - 103			
Terphenyl-d14	87		36 - 129			



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 Laboratory Reference: 1801-282  
 Project: 0146970060

**PAHs EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**

Matrix: Water

Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0130W1					
Naphthalene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
2-Methylnaphthalene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
1-Methylnaphthalene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Acenaphthylene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Acenaphthene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Fluorene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Phenanthrene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Anthracene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Fluoranthene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Pyrene	ND	0.10	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[a]anthracene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Chrysene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[b]fluoranthene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[a]pyrene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270D/SIM	1-30-18	1-30-18	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	56	25 - 107				
Pyrene-d10	75	28 - 103				
Terphenyl-d14	84	36 - 129				



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 Project: 0146970060

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

Matrix: Water

Units: ug/L

Analyte	Result	Spike Level	Source	Percent	Recovery	RPD	RPD Limit	Flags				
			Result	Recovery	Limits							
<b>MATRIX SPIKES</b>												
Laboratory ID:	01-271-01											
	MS	MSD	MS	MSD	MS	MSD						
Naphthalene	<b>0.421</b>	<b>0.469</b>	0.588	0.569	ND	72	82	27 - 106				
Acenaphthylene	<b>0.518</b>	<b>0.526</b>	0.588	0.569	ND	88	92	20 - 117				
Acenaphthene	<b>0.648</b>	<b>0.582</b>	0.588	0.569	ND	110	102	30 - 114				
Fluorene	<b>0.552</b>	<b>0.529</b>	0.588	0.569	ND	94	93	36 - 116				
Phenanthrene	<b>0.494</b>	<b>0.501</b>	0.588	0.569	ND	84	88	31 - 122				
Anthracene	<b>0.589</b>	<b>0.586</b>	0.588	0.569	ND	100	103	33 - 144				
Fluoranthene	<b>0.498</b>	<b>0.505</b>	0.588	0.569	ND	85	89	44 - 120				
Pyrene	<b>0.506</b>	<b>0.517</b>	0.588	0.569	ND	86	91	40 - 130				
Benzo[a]anthracene	<b>0.571</b>	<b>0.595</b>	0.588	0.569	ND	97	105	47 - 131				
Chrysene	<b>0.556</b>	<b>0.547</b>	0.588	0.569	ND	95	96	48 - 120				
Benzo[b]fluoranthene	<b>0.560</b>	<b>0.563</b>	0.588	0.569	ND	95	99	42 - 128				
Benzo(j,k)fluoranthene	<b>0.537</b>	<b>0.555</b>	0.588	0.569	ND	91	98	46 - 121				
Benzo[a]pyrene	<b>0.556</b>	<b>0.570</b>	0.588	0.569	ND	95	100	34 - 121				
Indeno(1,2,3-c,d)pyrene	<b>0.551</b>	<b>0.573</b>	0.588	0.569	ND	94	101	39 - 128				
Dibenz[a,h]anthracene	<b>0.557</b>	<b>0.572</b>	0.588	0.569	ND	95	101	39 - 125				
Benzo[g,h,i]perylene	<b>0.522</b>	<b>0.535</b>	0.588	0.569	ND	89	94	41 - 122				
Surrogate:												
2-Fluorobiphenyl						53	58	25 - 107				
Pyrene-d10						71	74	28 - 103				
Terphenyl-d14						98	101	36 - 129				



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 Project: 0146970060

**PAHs EPA 8270D/SIM**  
**SPIKE BLANK QUALITY CONTROL**

Matrix: Water  
 Units: ug/L

Analyte	Result	Spike Level	Percent Recovery	Recovery Limits	Flags
<b>SPIKE BLANK</b>					
Laboratory ID:	SB0130W1				
	SB	SB	SB		
Naphthalene	<b>0.360</b>	0.500	72	27 - 106	
Acenaphthylene	<b>0.387</b>	0.500	77	20 - 117	
Acenaphthene	<b>0.461</b>	0.500	92	30 - 114	
Fluorene	<b>0.454</b>	0.500	91	36 - 116	
Phenanthrene	<b>0.458</b>	0.500	92	31 - 122	
Anthracene	<b>0.480</b>	0.500	96	33 - 144	
Fluoranthene	<b>0.488</b>	0.500	98	44 - 120	
Pyrene	<b>0.566</b>	0.500	113	40 - 130	
Benzo[a]anthracene	<b>0.503</b>	0.500	101	47 - 131	
Chrysene	<b>0.498</b>	0.500	100	48 - 120	
Benzo[b]fluoranthene	<b>0.491</b>	0.500	98	42 - 128	
Benzo(j,k)fluoranthene	<b>0.489</b>	0.500	98	46 - 121	
Benzo[a]pyrene	<b>0.502</b>	0.500	100	34 - 121	
Indeno(1,2,3-c,d)pyrene	<b>0.522</b>	0.500	104	39 - 128	
Dibenz[a,h]anthracene	<b>0.528</b>	0.500	106	39 - 125	
Benzo[g,h,i]perylene	<b>0.518</b>	0.500	104	41 - 122	
<i>Surrogate:</i>					
<i>2-Fluorobiphenyl</i>			83	25 - 107	
<i>Pyrene-d10</i>			85	28 - 103	
<i>Terphenyl-d14</i>			103	36 - 129	



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 Project: 0146970060

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

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	EPA Method	Date Prepared	Date Analyzed	Flags
Lab ID:	01-282-01					
<b>Client ID:</b>	<b>KMW-03R-012618</b>					
Arsenic	<b>ND</b>	3.3	200.8	1-31-18	1-31-18	
Barium	<b>ND</b>	28	200.8	1-31-18	1-31-18	
Cadmium	<b>ND</b>	4.4	200.8	1-31-18	1-31-18	
Chromium	<b>ND</b>	11	200.8	1-31-18	1-31-18	
Lead	<b>ND</b>	1.1	200.8	1-31-18	1-31-18	
Mercury	<b>ND</b>	0.50	7470A	1-30-18	1-30-18	
Selenium	<b>ND</b>	5.6	200.8	1-31-18	1-31-18	
Silver	<b>ND</b>	11	200.8	1-31-18	1-31-18	

Lab ID: 01-282-02  
**Client ID:** KMW-08-012618

Arsenic	<b>ND</b>	3.3	200.8	1-31-18	1-31-18
Barium	<b>45</b>	28	200.8	1-31-18	1-31-18
Cadmium	<b>ND</b>	4.4	200.8	1-31-18	1-31-18
Chromium	<b>ND</b>	11	200.8	1-31-18	1-31-18
Lead	<b>ND</b>	1.1	200.8	1-31-18	1-31-18
Mercury	<b>ND</b>	0.50	7470A	1-30-18	1-30-18
Selenium	<b>ND</b>	5.6	200.8	1-31-18	1-31-18
Silver	<b>ND</b>	11	200.8	1-31-18	1-31-18



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: February 2, 2018  
 Samples Submitted: January 26, 2018  
 Laboratory Reference: 1801-282  
 Project: 0146970060

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

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	EPA Method	Prepared	Date	Analyzed	Date	Flags
Lab ID:	01-282-03							
<b>Client ID:</b>	<b>KMW-07-012618</b>							
Arsenic	ND	3.3	200.8	1-31-18	1-31-18			
Barium	ND	28	200.8	1-31-18	1-31-18			
Cadmium	ND	4.4	200.8	1-31-18	1-31-18			
Chromium	ND	11	200.8	1-31-18	1-31-18			
Lead	ND	1.1	200.8	1-31-18	1-31-18			
Mercury	ND	0.50	7470A	1-30-18	1-30-18			
Selenium	ND	5.6	200.8	1-31-18	1-31-18			
Silver	ND	11	200.8	1-31-18	1-31-18			

Lab ID: 01-282-04  
**Client ID:** KMW-02R-012618

Arsenic	ND	3.3	200.8	1-31-18	1-31-18			
Barium	ND	28	200.8	1-31-18	1-31-18			
Cadmium	ND	4.4	200.8	1-31-18	1-31-18			
Chromium	ND	11	200.8	1-31-18	1-31-18			
Lead	ND	1.1	200.8	1-31-18	1-31-18			
Mercury	ND	0.50	7470A	1-30-18	1-30-18			
Selenium	ND	5.6	200.8	1-31-18	1-31-18			
Silver	ND	11	200.8	1-31-18	1-31-18			



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This report pertains to the samples analyzed in accordance with the chain of custody,  
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Date of Report: February 2, 2018  
 Samples Submitted: January 26, 2018  
 Laboratory Reference: 1801-282  
 Project: 0146970060

**TOTAL METALS**  
**EPA 200.8/7470A**  
**METHOD BLANK QUALITY CONTROL**

Date Extracted: 1-30&31-18  
 Date Analyzed: 1-30&31-18  
 Matrix: Water  
 Units: ug/L (ppb)  
 Lab ID: MB0130W1&MB0131WM1

Analyte	Method	Result	PQL
Arsenic	200.8	<b>ND</b>	3.3
Barium	200.8	<b>ND</b>	28
Cadmium	200.8	<b>ND</b>	4.4
Chromium	200.8	<b>ND</b>	11
Lead	200.8	<b>ND</b>	1.1
Mercury	7470A	<b>ND</b>	0.50
Selenium	200.8	<b>ND</b>	5.6
Silver	200.8	<b>ND</b>	11



Date of Report: February 2, 2018  
 Samples Submitted: January 26, 2018  
 Laboratory Reference: 1801-282  
 Project: 0146970060

**TOTAL METALS  
EPA 200.8/7470A  
DUPLICATE QUALITY CONTROL**

Date Extracted: 1-30&31-18  
 Date Analyzed: 1-30&31-18

Matrix: Water  
 Units: ug/L (ppb)

Lab ID: 01-271-01

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Arsenic	<b>6.78</b>	<b>6.33</b>	7	3.3	
Barium	<b>ND</b>	<b>ND</b>	NA	28	
Cadmium	<b>ND</b>	<b>ND</b>	NA	4.4	
Chromium	<b>ND</b>	<b>ND</b>	NA	11	
Lead	<b>1.12</b>	<b>ND</b>	NA	1.1	
Mercury	<b>ND</b>	<b>ND</b>	NA	0.50	
Selenium	<b>ND</b>	<b>ND</b>	NA	5.6	
Silver	<b>ND</b>	<b>ND</b>	NA	11	



Date of Report: February 2, 2018  
 Samples Submitted: January 26, 2018  
 Laboratory Reference: 1801-282  
 Project: 0146970060

**TOTAL METALS**  
**EPA 200.8/7470A**  
**MS/MSD QUALITY CONTROL**

Date Extracted: 1-30&31-18  
 Date Analyzed: 1-30&31-18

Matrix: Water  
 Units: ug/L (ppb)

Lab ID: 01-271-01

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Arsenic	222	<b>227</b>	99	<b>233</b>	102	3	
Barium	222	<b>224</b>	101	<b>222</b>	100	1	
Cadmium	222	<b>228</b>	103	<b>231</b>	104	1	
Chromium	222	<b>211</b>	95	<b>216</b>	97	2	
Lead	222	<b>207</b>	93	<b>207</b>	93	0	
Mercury	12.5	<b>12.3</b>	98	<b>12.4</b>	99	1	
Selenium	222	<b>231</b>	104	<b>232</b>	104	0	
Silver	222	<b>196</b>	88	<b>202</b>	91	3	



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,  
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Date of Report: February 2, 2018  
 Samples Submitted: January 26, 2018  
 Laboratory Reference: 1801-282  
 Project: 0146970060

**TOTAL METALS**  
**EPA 200.8/7470A**  
**SPIKE BLANK QUALITY CONTROL**

Date Extracted: 1-30&31-18  
 Date Analyzed: 1-30&31-18  
 Matrix: Water  
 Units: ug/L (ppb)  
 Lab ID: SB0130W1&SB0131WM1

Analyte	Method	Spike Level	SB Result	Percent Recovery
Arsenic	200.8	222	<b>228</b>	103
Barium	200.8	222	<b>212</b>	96
Cadmium	200.8	222	<b>224</b>	101
Chromium	200.8	222	<b>210</b>	95
Lead	200.8	222	<b>217</b>	98
Mercury	7470A	12.5	<b>12.6</b>	101
Selenium	200.8	222	<b>240</b>	108
Silver	200.8	222	<b>207</b>	93





### 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



## Chain of Custody

 Page 1 of 1

 Turnaround Request  
(in working days)  
(Check One)

 Company: **Ameec Foster Wheeler**  
 Project Number: **0146917**

 Project Name: **Kelli Moore**  
 Project Manager: **Kathleen Goodman**

 Sampled by: **K. Black**
 Standard (7 Days)  
(TPH analysis 5 Days)  
  
 \_\_\_\_\_  
(other)

 Laboratory Number: **01 - 282**

Lab ID	Sample Identification	Date Sampled	Time Sampled	Matrix	Number of Containers
1	KMW-03R-012618	1/24/18	0945	W	10
2	KMW-08-012618	1/26/18	1100	W	10
3	KMW-07-012618	1/26/18	1200	W	10
4	KMW-02R-012618	1/26/18	1255	W	10
5	Trip Blanks	—	—	W	4

Comments/Special Instructions	Date	Time	Comments/Special Instructions	Date	Time	Comments/Special Instructions
Relinquished	<i>Michele Morello</i>	Ameec Foster Wheeler	1/26/18	1442		
Received			Office Sealed w/ packing tape	1/26/18	1442	
Relinquished	<i>Engineer Cao</i>	Ameec Foster Wheeler	1/26/18	15.02		
Received	<i>John Gorme</i>	ALPHA	1/26/18	1502		
Relinquished	<i>John Gorme</i>	ALPHA	1/26/18	16:18		
Received	<i>John Gorme</i>	ALPHA	1/26/18	16:18		
Reviewed/Date						

 Data Package: Standard  Level III  Level IV 

 Chromatograms with final report  Electronic Data Deliverables (EDDs)

**wood.**

## **Appendix C**

Please print, sign and return to the Department of Ecology

## RESOURCE PROTECTION WELL REPORT

(SUBMIT ONE WELL REPORT PER WELL INSTALLED) SVE-01

CURRENT Notice of Intent No. RE11442

Construction/Decommission ("x" in box)

Construction

Decommission

ORIGINAL INSTALLATION Notice of Intent Number:

RE11442

Consulting Firm AMEC Foster Wheeler

Unique Ecology Well IDTag No. BCB316

WELL CONSTRUCTION CERTIFICATION: I constructed and/or accept responsibility for construction of this well, and its compliance with all Washington well construction standards. Materials used and the information reported above are true to my best knowledge and belief.

Driller  Engineer  Trainee

Name (Print Last, First Name) Hsieh, Pat

Driller/Engineer/Trainee Signature 

Driller or Trainee License No. 45361

Type of Well ("x in box)

Resource Protection

Geotech Soil Boring

Property Owner Robert Stetson - Kelley-Moore Paint Company Inc.

Site Address 5400 Airport Way South

City Seattle County King

Location SW 1/4-1/4 NE 1/4 Sec 20 Twp 24N R 4E

EWM  or WWM

Lat/Long (s, t, r) Lat Deg \_\_\_\_ Min \_\_\_\_ Sec \_\_\_\_  
still REQUIRED) Long Deg \_\_\_\_ Min \_\_\_\_ Sec \_\_\_\_

Tax Parcel No. 3868400270

Cased or Uncased Diameter 4" Static Level NA

Work/Decommission Start Date 6/24/15

Work/Decommission Completed Date 6/25/15

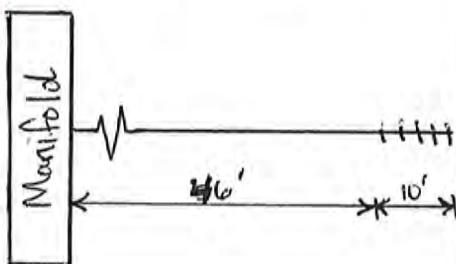
If trainee, licensed driller's Signature and License Number:

### Construction Design

### Well Data

### Formation Description

Horizontal Soil Vapor extraction well.



Name: SVE-01

Length: 10'-0"

Type: 0.020" slot  
Screen sch. 40 pvc

Well screened within  
vadose zone at  
7' below ground surface

Black silty sand,  
medium sands.

Please print, sign and return to the Department of Ecology

## RESOURCE PROTECTION WELL REPORT

(SUBMIT ONE WELL REPORT PER WELL INSTALLED) SVE-02

CURRENT Notice of Intent No. RE11442

Construction/Decommission ("x" in box)

- Construction
- Decommission

ORIGINAL INSTALLATION Notice of Intent Number:

RE11442

Consulting Firm AMEC Foster Wheeler

Unique Ecology Well IDTag No. BCB317

WELL CONSTRUCTION CERTIFICATION: I constructed and/or accept responsibility for construction of this well, and its compliance with all Washington well construction standards. Materials used and the information reported above are true to my best knowledge and belief.

Driller  Engineer  Trainee

Name (Print Last, First Name) Hsieh, Pat

Driller/Engineer/Trainee Signature 

Driller or Trainee License No. 45361

Type of Well ("x in box)

- Resource Protection
- Geotech Soil Boring

Property Owner Robert Stetson - Kelley-Moore Paint Company Inc.

Site Address 5400 Airport Way South

City Seattle County King

Location SW 1/4-1/4 NE 1/4 Sec 20 Twp 24N R 4E

EWM  or WWM

Lat/Long (s, t, r) Lat Deg \_\_\_\_ Min \_\_\_\_ Sec \_\_\_\_  
still REQUIRED Long Deg \_\_\_\_ Min \_\_\_\_ Sec \_\_\_\_

Tax Parcel No. 3868400270

Cased or Uncased Diameter 4" Static Level NA

Work/Decommission Start Date 6/24/15

Work/Decommission Completed Date 6/25/15

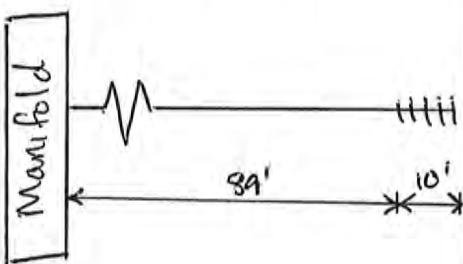
If trainee, licensed driller's Signature and License Number:

### Construction Design

### Well Data

### Formation Description

Horizontal Soil Vapor extraction well



Name: SVE-02

Length: 10'-0"

Type: 0.020" slot screen  
sch. 40 pvc

Well screened within  
Vadose zone at 7'  
below ground surface

Brown gravelly sands

SCALE: 1" = N/A PAGE 2 OF 10

Please print, sign and return to the Department of Ecology

## RESOURCE PROTECTION WELL REPORT

(SUBMIT ONE WELL REPORT PER WELL INSTALLED) SVE-03

CURRENT Notice of Intent No. RE11442

Construction/Decommission ("x" in box)

Construction

Decommission

ORIGINAL INSTALLATION Notice of Intent Number:

RE11442

Consulting Firm AMEC Foster Wheeler

Unique Ecology Well IDTag No. BCB3i8

WELL CONSTRUCTION CERTIFICATION: I constructed and/or accept responsibility for construction of this well, and its compliance with all Washington well construction standards. Materials used and the information reported above are true to my best knowledge and belief.

Driller  Engineer  Trainee

Name (Print Last, First Name) Hsieh, Pat

Driller/Engineer/Trainee Signature 

Driller or Trainee License No. 45361

Type of Well ("x in box)

Resource Protection

Geotech Soil Boring

Property Owner Robert Stetson - Kelley-Moore Paint Company Inc.

Site Address 5400 Airport Way South

City Seattle County King

Location SW 1/4-1/4 NE 1/4 Sec 20 Twp 24N R 4E

EWM  or WWM

Lat/Long (s, t, r) Lat Deg \_\_\_\_\_ Min \_\_\_\_\_ Sec \_\_\_\_\_  
still REQUIRED Long Deg \_\_\_\_\_ Min \_\_\_\_\_ Sec \_\_\_\_\_

Tax Parcel No. 3868400270

Cased or Uncased Diameter 4" Static Level NA

Work/Decommission Start Date 6/23/15

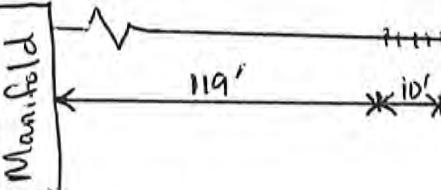
Work/Decommission Completed Date 6/25/15

If trainee, licensed driller's Signature and License Number:

### Construction Design

### Well Data

### Formation Description

Horizontal <del>Soil</del> Soil Vapor extraction well 	Name: SVE-03 Length: 10' - 0" Type: 0.020" slot screen sch. 40 PVC Well screened within vadose zone at 7' below ground surface	Black silty, medium sands.
---	---	----------------------------

SCALE: 1"= N/A PAGE 3 OF 10

Please print, sign and return to the Department of Ecology

## RESOURCE PROTECTION WELL REPORT

(SUBMIT ONE WELL REPORT PER WELL INSTALLED) SVE-04

CURRENT Notice of Intent No. RE11442

Construction/Decommission ("x" in box)

Construction

Decommission

ORIGINAL INSTALLATION Notice of Intent Number:

RE11442

Consulting Firm AMEC Foster Wheeler

Unique Ecology Well IDTag No. BC8322

WELL CONSTRUCTION CERTIFICATION: I constructed and/or accept responsibility for construction of this well, and its compliance with all Washington well construction standards. Materials used and the information reported above are true to my best knowledge and belief.

Driller  Engineer  Trainee

Name (Print Last, First Name) Hsieh, Pat

Driller/Engineer/Trainee Signature 

Driller or Trainee License No. 45341

Type of Well ("x in box)

Resource Protection

Geotech Soil Boring

Property Owner Robert Stetson - Kelley-Moore Paint Company Inc.

Site Address 5400 Airport Way South

City Seattle County King

Location SW 1/4-1/4 NE 1/4 Sec 20 Twp 24N R 4E

EWM  or WWM

Lat/Long (s, t, r still REQUIRED) Lat Deg \_\_\_\_ Min \_\_\_\_ Sec \_\_\_\_ Long Deg \_\_\_\_ Min \_\_\_\_ Sec \_\_\_\_

Tax Parcel No. 3868400270

Cased or Uncased Diameter 4" Static Level NA

Work/Decommission Start Date 6/23/15

Work/Decommission Completed Date 6/25/15

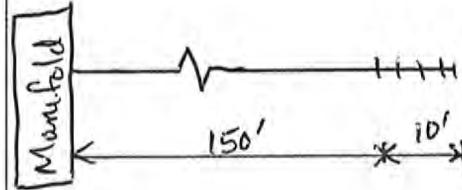
If trainee, licensed driller's Signature and License Number:

### Construction Design

### Well Data

### Formation Description

Horizontal soil vapor extraction well



Name: SVE-04

Length: 10'-0"

Type: 0.020" slot screen  
sch. 40 PVC

Well screened within  
vadose zone at 7'  
below ground surface

Black silty medium  
sands

Please print, sign and return to the Department of Ecology

## RESOURCE PROTECTION WELL REPORT

(SUBMIT ONE WELL REPORT PER WELL INSTALLED) SVE-05

CURRENT Notice of Intent No. RE11442

Construction/Decommission ("x" in box)

Construction

Decommission

ORIGINAL INSTALLATION Notice of Intent Number:

RE11442

Consulting Firm AMEC Foster Wheeler

Unique Ecology Well IDTag No. BC8323

WELL CONSTRUCTION CERTIFICATION: I constructed and/or accept responsibility for construction of this well, and its compliance with all Washington well construction standards. Materials used and the information reported above are true to my best knowledge and belief.

Driller  Engineer  Trainee

Name (Print Last, First Name) Hsieh, Pat

Driller/Engineer /Trainee Signature 

Driller or Trainee License No. 45561

Type of Well ("x in box)

Resource Protection

Geotech Soil Boring

Property Owner Robert Stetson - Kelley-Moore Paint Company Inc.

Site Address 5400 Airport Way South

City Seattle County King

Location SW 1/4-1/4 NE 1/4 Sec 20 Twp 24N R 4E

EWM  or WWM

Lat/Long (s, t, r) Lat Deg \_\_\_\_ Min \_\_\_\_ Sec \_\_\_\_  
still REQUIRED Long Deg \_\_\_\_ Min \_\_\_\_ Sec \_\_\_\_

Tax Parcel No. 3868400270

Cased or Uncased Diameter 4" Static Level N/A

Work/Decommission Start Date 6/22/15

Work/Decommission Completed Date 6/25/15

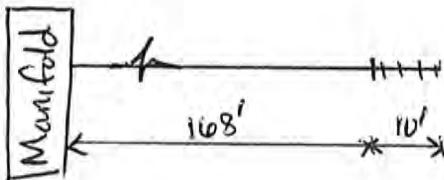
If trainee, licensed driller's Signature and License Number:

### Construction Design

### Well Data

### Formation Description

Horizontal soil vapor extraction well



Name: SVE-05

Length: 10'-0"

Type: 0.020" slot screen  
Sch. 40 PVC

Well screened within  
vadose zone at 7'  
below ground surface.

Black silty medium  
sands.

Please print, sign and return to the Department of Ecology

## RESOURCE PROTECTION WELL REPORT

(SUBMIT ONE WELL REPORT PER WELL INSTALLED) SVE-06

Construction/Decommission ("x" in box)

- Construction
- Decommission

ORIGINAL INSTALLATION Notice of Intent Number:

RE11442

Consulting Firm AMEC Foster Wheeler

Unique Ecology Well IDTag No. BCB325

WELL CONSTRUCTION CERTIFICATION: I constructed and/or accept responsibility for construction of this well, and its compliance with all Washington well construction standards. Materials used and the information reported above are true to my best knowledge and belief.

Driller  Engineer  Trainee

Name (Print Last, First Name) Hsieh, Pat

Driller/Engineer /Trainee Signature 

Driller or Trainee License No. 45361

Type of Well ("x in box)

- Resource Protection
- Geotech Soil Boring

Property Owner Robert Stetson - Kelley-Moore Paint Company Inc.

Site Address 5400 Airport Way South

City Seattle County King

Location SW 1/4-1/4 NE 1/4 Sec 20 Twn 24N R 4E

EWM  or WWM

Lat/Long (s, t, r) Lat Deg \_\_\_\_ Min \_\_\_\_ Sec \_\_\_\_\_  
still REQUIRED Long Deg \_\_\_\_ Min \_\_\_\_ Sec \_\_\_\_\_

Tax Parcel No. 3868400270

Cased or Uncased Diameter 4" Static Level NA

Work/Decommission Start Date 6/19/15

Work/Decommission Completed Date 6/25/15

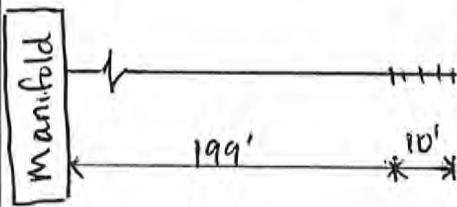
If trainee, licensed driller's Signature and License Number:

### Construction Design

### Well Data

### Formation Description

Horizontal Soil Vapor extraction well



Name: SVE-06

Length: 10'-0"

Type: 0.020" Slot Screen  
Sch 40 PVC

Well screened in vadose zone at 7' below ground surface

Black silty medium sands.

SCALE: 1" = N/A PAGE 16 OF 10



Please print, sign and return to the Department of Ecology

## RESOURCE PROTECTION WELL REPORT

CURRENT Notice of Intent No. RE11442

(SUBMIT ONE WELL REPORT PER WELL INSTALLED) SVE-08

Construction/Decommission ("x" in box)

- Construction  
 Decommission

ORIGINAL INSTALLATION Notice of Intent Number:

RE11442

Consulting Firm AMEC Foster Wheeler

Unique Ecology Well IDTag No. BCB 327

WELL CONSTRUCTION CERTIFICATION: I constructed and/or accept responsibility for construction of this well, and its compliance with all Washington well construction standards. Materials used and the information reported above are true to my best knowledge and belief.

Driller  Engineer  Trainee

Name (Print Last, First Name) Hsieh, Paul

Driller/Engineer/Trainee Signature 

Driller or Trainee License No. 45561

Type of Well ("x in box")

- Resource Protection  
 Geotech Soil Boring

Property Owner Robert Stetson - Kelley-Moore Paint Company Inc.

Site Address 5400 Airport Way South

City Seattle County King

Location SW 1/4-1/4 NE 1/4 Sec 20 Twn 24N R 4E

EWM  or WWM

Lat/Long (s, t, r) still REQUIRED Lat Deg \_\_\_\_ Min \_\_\_\_ Sec \_\_\_\_  
Long Deg \_\_\_\_ Min \_\_\_\_ Sec \_\_\_\_

Tax Parcel No. 3868400270

Cased or Uncased Diameter 4" Static Level NA

Work/Decommission Start Date 4/18/15

Work/Decommission Completed Date 6/25/15

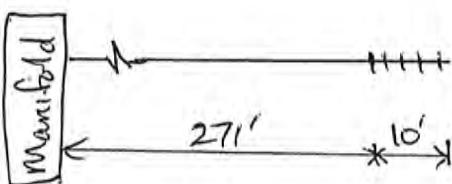
If trainee, licensed driller's Signature and License Number:

### Construction Design

### Well Data

### Formation Description

Horizontal Soil Vapor Extraction well



Name: SVE-08  
Length: 10'-0"  
Type: 0.020" slot screen  
Sch 40 PVC

Well screened within vadose zone at 7' below ground surface.

Black silty medium sands

SCALE: 1" = N/A PAGE 8 OF 10

Please print, sign and return to the Department of Ecology

## RESOURCE PROTECTION WELL REPORT

(SUBMIT ONE WELL REPORT PER WELL INSTALLED) SVE-09

CURRENT Notice of Intent No. RE11442

Construction/Decommission ("x" in box)

- Construction  
 Decommission

ORIGINAL INSTALLATION Notice of Intent Number:

RE11442

Consulting Firm AMEC Foster Wheeler

Unique Ecology Well IDTag No. BC8328

WELL CONSTRUCTION CERTIFICATION: I constructed and/or accept responsibility for construction of this well, and its compliance with all Washington well construction standards. Materials used and the information reported above are true to my best knowledge and belief.

- Driller  Engineer  Trainee

Name (Print Last, First Name) Hsieh, Pat

Driller/Engineer/Trainee Signature 

Driller or Trainee License No. 45361

Type of Well ("x in box")

- Resource Protection  
 Geotech Soil Boring

Property Owner Robert Stetson - Kelley-Moore Paint Company Inc.

Site Address 5400 Airport Way South

City Seattle County King

Location SW 1/4-1/4 NE 1/4 Sec 20 Twp 24N R 4E

EWM  or WWM

Lat/Long (s, t, r still REQUIRED) Lat Deg \_\_\_\_ Min \_\_\_\_ Sec \_\_\_\_ Long Deg \_\_\_\_ Min \_\_\_\_ Sec \_\_\_\_

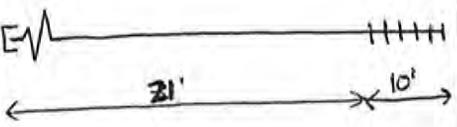
Tax Parcel No. 3868400270

Cased or Uncased Diameter 4" Static Level NA

Work/Decommission Start Date 10/15/2015

Work/Decommission Completed Date 10/15/2015

If trainee, licensed driller's Signature and License Number:

Construction Design	Well Data	Formation Description
Horizontal soil vapor extraction well.    Name: SVE-09 Screen length: 10'-0" Type: 0.020" slotted screen, Sch. 40 PVC  Well is screened within vadose zone at 4' below ground surface.  Well head is capped and buried; it will be connected to treatment manifold during system commissioning		Black silty sand, medium sands.

SCALE: 1"= NA PAGE 9 OF 10

Please print, sign and return to the Department of Ecology

## RESOURCE PROTECTION WELL REPORT

CURRENT Notice of Intent No. RE11442

(SUBMIT ONE WELL REPORT PER WELL INSTALLED) SVE-10

Construction/Decommission ("x" in box)

- Construction
- Decommission

ORIGINAL INSTALLATION Notice of Intent Number:

RE11442

Consulting Firm AMEC Foster Wheeler

Unique Ecology Well IDTag No. BCB 332

WELL CONSTRUCTION CERTIFICATION: I constructed and/or accept responsibility for construction of this well, and its compliance with all Washington well construction standards. Materials used and the information reported above are true to my best knowledge and belief.

Driller  Engineer  Trainee

Name (Print Last, First Name) Hsieh, Pat

Driller/Engineer/Trainee Signature 

Driller or Trainee License No. 45361

Type of Well ("x in box")

- Resource Protection
- Geotech Soil Boring

Property Owner Robert Stetson - Kelley-Moore Paint Company Inc.

Site Address 5400 Airport Way South

City Seattle County King

Location SW 1/4-1/4 NE 1/4 Sec 20 Twn 24N R 4E

EWM  or WWM

Lat/Long (s, t, r) still REQUIRED Lat Deg \_\_\_\_ Min \_\_\_\_ Sec \_\_\_\_ Long Deg \_\_\_\_ Min \_\_\_\_ Sec \_\_\_\_

Tax Parcel No. 3868400270

Cased or Uncased Diameter 4" Static Level NA

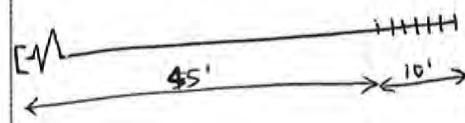
Work/Decommission Start Date 10/15/2015

Work/Decommission Completed Date 10/19/2015

If trainee, licensed driller's Signature and License Number:

### Construction Design

Horizontal soil vapor extraction well.



### Well Data

Name: SVE-10  
Screen length: 10'-0"  
Type: 0.020" slotted screen  
Sch. 40 PVC

well is screened within vadose zone at 4' below ground surface.

Well head is capped and buried; it will be connected to treatment manifold during system commissioning.

### Formation Description

Black silty sand, medium sands.

Please print, sign and return to the Department of Ecology

## RESOURCE PROTECTION WELL REPORT

(SUBMIT ONE WELL REPORT PER WELL INSTALLED) SVE-11

CURRENT Notice of Intent No. RE11868

Construction/Decommission ("x" in box)

- Construction
- Decommission

ORIGINAL INSTALLATION Notice of Intent Number:

RE11868

Consulting Firm AMEC Foster Wheeler

Unique Ecology Well IDTag No. BCB333

WELL CONSTRUCTION CERTIFICATION: I constructed and/or accept responsibility for construction of this well, and its compliance with all Washington well construction standards. Materials used and the information reported above are true to my best knowledge and belief.

Driller  Engineer  Trainee

Name (Print Last, First Name) Hsieh, Pat

Driller/Engineer/Trainee Signature 

Driller or Trainee License No. 45361

Type of Well ("x in box")

- Resource Protection
- Geotech Soil Boring

Property Owner Robert Stetson - Kelley-Moore Paint Company Inc.

Site Address 5400 Airport Way South

City Seattle County King

Location SW 1/4-1/4 NE 1/4 Sec 20 Twp 24N R 4E

EWM  or WWM

Lat/Long (s, t, r) Lat Deg \_\_\_\_ Min \_\_\_\_ Sec \_\_\_\_  
still REQUIRED) Long Deg \_\_\_\_ Min \_\_\_\_ Sec \_\_\_\_

Tax Parcel No. 3868400270

Cased or Uncased Diameter 4" Static Level \_\_\_\_\_

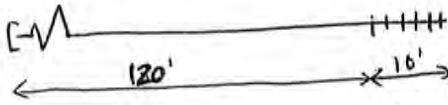
Work/Decommission Start Date 10/15/2015

Work/Decommission Completed Date 10/21/2015

If trainee, licensed driller's Signature and License Number:

### Construction Design

Horizontal soil vapor extraction well.



### Well Data

Name: SVE-11  
Screen length: 10'-0"  
Type: 0.020" slotted screen,  
Sch. 40 PVC

Well is screened within vadose zone at 4' below ground surface.

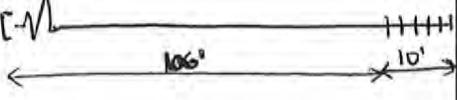
Well head is capped and buried; it will be connected to treatment manifold during system commissioning.

### Formation Description

Black silty sand, medium sands.

Please print, sign and return to the Department of Ecology

## RESOURCE PROTECTION WELL REPORT

(SUBMIT ONE WELL REPORT PER WELL INSTALLED) SVE-12		CURRENT Notice of Intent No. RE11868
Construction/Decommission ("x" in box)		
<input checked="" type="checkbox"/> Construction <input type="checkbox"/> Decommission		
ORIGINAL INSTALLATION Notice of Intent Number: RE11868		
Consulting Firm AMEC Foster Wheeler		
Unique Ecology Well IDTag No. BJK722		
WELL CONSTRUCTION CERTIFICATION: I constructed and/or accept responsibility for construction of this well, and its compliance with all Washington well construction standards. Materials used and the information reported above are true to my best knowledge and belief.		
<input type="checkbox"/> Driller <input checked="" type="checkbox"/> Engineer <input type="checkbox"/> Trainee Name (Print Last, First Name) Hsieh, Pat Driller/Engineer/Trainee Signature  Driller or Trainee License No. 45361		
If trainee, licensed driller's Signature and License Number:		
Construction Design	Well Data	Formation Description
Horizontal well for soil vapor extraction. 	Name: SVE-12 Screen length: 10'-0" Type: 0.020" slotted screen, Sch. 40 PVC  Well is screened within the vadose zone at 4' below ground surface.  Well head is capped and buried; it will be connected to the treatment manifold during system commissioning.	Black silty medium sands.
SCALE: 1" = NA PAGE 2 OF 2		

PROJECT: Former Kelly-Moore Paint Company - Air Sparge Well Installation Seattle, Washington				<b>Log of Well No. AS-1</b>				
BORING LOCATION: S. parking lot, S. south end of trench by system area fence				GROUND SURFACE ELEVATION AND DATUM: Ground surface				
DRILLING CONTRACTOR: Cascade Drilling, Inc.				DATE STARTED: 11/4/16		DATE FINISHED: 11/4/16		
DRILLING METHOD: Hollow-stem auger				TOTAL DEPTH (ft.): 19.0		SCREEN INTERVAL (ft.): 13.5-18.5		
DRILLING EQUIPMENT: CME 75				DEPTH TO FIRST WATER:	COMPL.	CASING:	2" Schedule 40 PVC	
SAMPLING METHOD: Split spoon sampler				LOGGED BY: K. Black				
HAMMER WEIGHT: 300 lb		DROP: 30"		RESPONSIBLE PROFESSIONAL: J. Long		REG. NO. L.Hg. 1354		
DEPTH (feet)	SAMPLE No.	SAMPLE No.	OVM Reading	DESCRIPTION NAME (USCS): color, moist, % by wt., plast. density, structure, cementation, react. w/HCl, geo. inter.		WELL CONSTRUCTION DETAILS AND/OR DRILLING REMARKS		
				Surface Elevation: Not surveyed				
0				ASPHALTIC CONCRETE		Traffic Rated Well Box		
1				SILTY SAND (SM): very dark grayish brown (2.5Y 3/2), moist, 75% medium to coarse sand, 20% low plasticity fines, 5% fine gravel		Concrete to Ground Surface		
2								
3								
4	2	1	1	44.4				
5				POORLY-GRADED SAND (SP): very dark gray (2.5Y 3/1), moist, 95% fine to coarse sand, 5% low plasticity fines		Bentonite Grout		
6								
7	0	0	1	5346	hydrocarbon odor, sheen observed		2" Sch. 40 Casing	
8				wet, light sheen				
9								
10						20/40 Filter Sand		
11								
12								
13						2/12 Filter Sand		
14	3	4	5	20.9			2" Sch. 40 PVC 0.010 Screen	
15								
OAKWELL (REV. 8/2011)								
<b>amec</b>				Project No. 14697		Page 1 of 2		

## Log of Well No. AS-1 (cont'd)

DEPTH (feet)	SAMPLES			DESCRIPTION NAME (USCS): color, moist, % by wt., plast. density, structure, cementation, react. w/HCl, geo. inter.	WELL CONSTRUCTION DETAILS AND/OR DRILLING REMARKS
	Sample No.	Sample	Blows/ Foot		
15		2 5 5		POORLY-GRADED SAND (SP): Continued	
16			4 6 7	SILTY SAND (SM): very dark gray (2.5Y 3/1), wet, 85% fine to medium sand, 15% low plasticity fines	
17					
18				POORLY-GRADED SAND (SP): very dark gray (2.5Y 3/1), wet, 95% fine to medium sand, 5% low plasticity fines, trace coarse gravel	
19				Ecology Well ID BJV 327	
20					
21					
22					
23					
24					
25					
26					
27					
28					
29					
30					
31					
32					
33					

OAKWELLV (REV. 8/2011)

PROJECT: Former Kelly-Moore Paint Company - Air Sparge Well Installation Seattle, Washington				<b>Log of Well No. AS-2</b>			
BORING LOCATION: S. parking lot, 20ft west of silo				GROUND SURFACE ELEVATION AND DATUM: Ground surface			
DRILLING CONTRACTOR: Cascade Drilling, Inc.				DATE STARTED: 11/3/16		DATE FINISHED: 11/3/16	
DRILLING METHOD: Hollow-stem auger				TOTAL DEPTH (ft.): 19.0		SCREEN INTERVAL (ft.): 13.5-18.5	
DRILLING EQUIPMENT: CME 75				DEPTH TO FIRST WATER:	COMPL.	CASING:	2" Schedule 40 PVC
SAMPLING METHOD: Split spoon sampler				LOGGED BY: K. Black			
HAMMER WEIGHT: 300 lb		DROP: 30"		RESPONSIBLE PROFESSIONAL: J. Long		REG. NO. L.Hg. 1354	
DEPTH (feet)	SAMPLES Sample No.	SAMPLES Sample No.	Blows/ Foot	OVM Reading	DESCRIPTION NAME (USCS): color, moist, % by wt., plast. density, structure, cementation, react. w/HCl, geo. inter.	WELL CONSTRUCTION DETAILS AND/OR DRILLING REMARKS	
					Surface Elevation: Not surveyed		
0					ASPHALTIC CONCRETE		
1					SILTY SAND with GRAVEL (SM): light brownish gray (2.5Y 6/2), moist, 65% fine to coarse sand, %20 low plasticity fines, 15% fine to coarse gravel		
2							
3					POORLY-GRADED SAND with SILT (SP-SM): very dark gray (2.5Y 3/1), moist, 90% fine to coarse sand, 10% low plasticity fines		
4							
5							
6							
7							
8							
9					wet, trace fine to coarse gravels		
10							
11					woody debris		
12							
13							
14							
15							
OAKWELLV (REV. 8/2011)							
<b>amec</b>				Project No. 14697		Page 1 of 2	

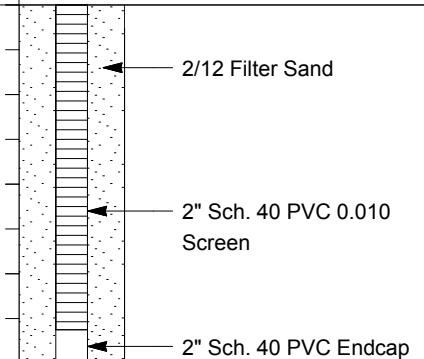
## Log of Well No. AS-2 (cont'd)

DEPTH (feet)	SAMPLES			DESCRIPTION NAME (USCS): color, moist, % by wt., plast. density, structure, cementation, react. w/HCl, geo. inter.	WELL CONSTRUCTION DETAILS AND/OR DRILLING REMARKS
	Sample No.	Sample	Blows/ Foot		
15				POORLY-GRADED SAND with SILT (SP-SM): Continued woody debris	
16					
17				SILTY SAND (SM): dark gray (2.5Y 4/1), wet, 70% fine sand, 30% low plasticity fines	
18					
19				Ecology Well ID BJV 328	
20					
21					
22					
23					
24					
25					
26					
27					
28					
29					
30					
31					
32					
33					

OAKWELLV (REV. 8/2011)

PROJECT: Former Kelly-Moore Paint Company - Air Sparge Well Installation Seattle, Washington				<b>Log of Well No. AS-3</b>			
BORING LOCATION: Mid parking lot, NW of silo				GROUND SURFACE ELEVATION AND DATUM: Ground surface			
DRILLING CONTRACTOR: Cascade Drilling, Inc.				DATE STARTED: 11/3/16		DATE FINISHED: 11/3/16	
DRILLING METHOD: Hollow-stem auger				TOTAL DEPTH (ft.): 19.0		SCREEN INTERVAL (ft.): 13.5-18.5	
DRILLING EQUIPMENT: CME 75				DEPTH TO FIRST WATER:	COMPL.	CASING:	2" Schedule 40 PVC
SAMPLING METHOD: Split spoon sampler				LOGGED BY: K. Black			
HAMMER WEIGHT: 300 lb		DROP: 30"		RESPONSIBLE PROFESSIONAL: J. Long		REG. NO. L.Hg. 1354	
DEPTH (feet)	SAMPLES Sample No.	SAMPLES Sample No.	Blows/ Foot	OVM Reading	DESCRIPTION NAME (USCS): color, moist, % by wt., plast. density, structure, cementation, react. w/HCl, geo. inter.	WELL CONSTRUCTION DETAILS AND/OR DRILLING REMARKS	
					Surface Elevation: Not surveyed		
0					ASPHALTIC CONCRETE		
1					SILTY SAND with GRAVEL (SM): light brownish gray (2.5Y 6/2), moist, 50% fine to coarse sand, 30% fine to coarse gravel, 20% low plasticity fines, trace cobles		
2							
3							
4							
5							
6							
7							
8							
9							
10							
11							
12							
13							
14							
15							
OAKWELL (REV. 8/2011)							
<b>amec</b>				Project No. 14697		Page 1 of 2	

## Log of Well No. AS-3 (cont'd)

DEPTH (feet)	SAMPLES			DESCRIPTION NAME (USCS): color, moist, % by wt., plast. density, structure, cementation, react. w/HCl, geo. inter.	WELL CONSTRUCTION DETAILS AND/OR DRILLING REMARKS
	Sample No.	Sample	Blows/ Foot		
15			5 5 6	2.6  POORLY-GRADED SAND (SP): Continued	
16					
17			5 5 5	1.8  SILTY SAND (SM): very dark gray (2.5Y 3/1), wet, 85% fine to medium sand, 15% low plasticity fines	
18					
19				Ecology Well ID BJV 325	
20					
21					
22					
23					
24					
25					
26					
27					
28					
29					
30					
31					
32					
33					

PROJECT: Former Kelly-Moore Paint Company - Air Sparge Well Installation Seattle, Washington				<b>Log of Well No. AS-4</b>			
BORING LOCATION: Mid parking lot, between AS-3 and AS-5				GROUND SURFACE ELEVATION AND DATUM: Ground surface			
DRILLING CONTRACTOR: Cascade Drilling, Inc.				DATE STARTED: 11/3/16		DATE FINISHED: 11/3/16	
DRILLING METHOD: Hollow-stem auger				TOTAL DEPTH (ft.): 19.0		SCREEN INTERVAL (ft.): 13.5-18.5	
DRILLING EQUIPMENT: CME 75				DEPTH TO FIRST WATER:	COMPL.	CASING:	2" Schedule 40 PVC
SAMPLING METHOD: Split spoon sampler				LOGGED BY: K. Black			
HAMMER WEIGHT: 300 lb		DROP: 30"		RESPONSIBLE PROFESSIONAL: J. Long		REG. NO. L.Hg. 1354	
DEPTH (feet)	SAMPLES Sample No.	SAMPLES Sample No.	Blows/ Foot	OVM Reading	DESCRIPTION NAME (USCS): color, moist, % by wt., plast. density, structure, cementation, react. w/HCl, geo. inter.	WELL CONSTRUCTION DETAILS AND/OR DRILLING REMARKS	
					Surface Elevation: Not surveyed		
0					ASPHALTIC CONCRETE		
1					SILTY SAND (SM): grayish brown (2.5Y 5/2), moist, 75% fine to coarse sand, 20% low plasticity fines, 5% fine to coarse gravel, some woody debris		
2							
3							
4							
5							
6							
7							
8							
9							
10							
11							
12							
13							
14							
15							
OAKWELL (REV. 8/2011)							
<b>amec</b>				Project No. 14697		Page 1 of 2	

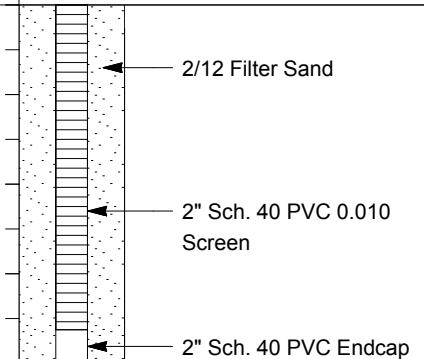
## Log of Well No. AS-4 (cont'd)

DEPTH (feet)	SAMPLES			DESCRIPTION NAME (USCS): color, moist, % by wt., plast. density, structure, cementation, react. w/HCl, geo. inter.	WELL CONSTRUCTION DETAILS AND/OR DRILLING REMARKS
	Sample No.	Sample	Blows/ Foot		
15				SILTY SAND (SM): Continued	
16				POORLY-GRADED SAND (SP): very dark gray (2.5Y 3/1), wet, 95% fine to medium sand, 5% low plasticity fines	
17				SILTY SAND (SM): very dark gray (2.5Y 3/1), wet, 85% fine to medium sand, 15% low to medium plasticity fines	
18					
19				Ecology Well ID BJV 326	
20					
21					
22					
23					
24					
25					
26					
27					
28					
29					
30					
31					
32					
33					

OAKWELLV (REV. 8/2011)

PROJECT: Former Kelly-Moore Paint Company - Air Sparge Well Installation Seattle, Washington				<b>Log of Well No. AS-5</b>			
BORING LOCATION: Mid parking lot, north end of trash system				GROUND SURFACE ELEVATION AND DATUM: Ground surface			
DRILLING CONTRACTOR: Cascade Drilling, Inc.				DATE STARTED: 11/4/16		DATE FINISHED: 11/4/16	
DRILLING METHOD: Hollow-stem auger				TOTAL DEPTH (ft.): 19.0		SCREEN INTERVAL (ft.): 13.5-18.5	
DRILLING EQUIPMENT: CME 75				DEPTH TO FIRST WATER:	COMPL.	CASING:	2" Schedule 40 PVC
SAMPLING METHOD: Split spoon sampler				LOGGED BY: K. Black			
HAMMER WEIGHT: 300 lb		DROP: 30"		RESPONSIBLE PROFESSIONAL: J. Long		REG. NO. L.Hg. 1354	
DEPTH (feet)	SAMPLE No.	SAMPLE No.	OVM Reading	DESCRIPTION NAME (USCS): color, moist, % by wt., plast. density, structure, cementation, react. w/HCl, geo. inter.		WELL CONSTRUCTION DETAILS AND/OR DRILLING REMARKS	
				Surface Elevation: Not surveyed			
0				ASPHALTIC CONCRETE		Traffic Rated Well Box	
1				SILTY SAND with GRAVEL (SM): dark olive brown (2.5Y 3/3), moist, 60% fine to medium sand, 25% low plasticity fines, 15% fine to coarse gravel		Concrete to Ground Surface	
2							
3							
4							
5	5	5	6	POORLY-GRADED SAND (SP): very dark gray (2.5Y 3/1), moist, 95% fine to medium sand, 5% low plasticity fines		Bentonite Grout	
6							
7	2	2	3	medium to coarse sand, hydrocarbon odor sheen observed		2" Sch. 40 Casing	
8							
9	3	4	4	wet		20/40 Filter Sand	
10							
11							
12							
13							
14	2	3	3	sheen observed		2/12 Filter Sand	
15						2" Sch. 40 PVC 0.010 Screen	

## Log of Well No. AS-5 (cont'd)

DEPTH (feet)	SAMPLES			DESCRIPTION NAME (USCS): color, moist, % by wt., plast. density, structure, cementation, react. w/HCl, geo. inter.	WELL CONSTRUCTION DETAILS AND/OR DRILLING REMARKS
	Sample No.	Sample	Blows/ Foot		
15		2 2 3	251.3	POORLY-GRADED SAND (SP): Continued	
16					
17		3 3 3	0.5	fine gravel lens	
18					
19				Ecology Well ID BJV 328	
20					
21					
22					
23					
24					
25					
26					
27					
28					
29					
30					
31					
32					
33					

**wood.**

## **Appendix D**



STATE OF WASHINGTON  
DEPARTMENT OF ECOLOGY

PO Box 47600 • Olympia, WA 98504-7600 • 360-407-6000

711 for Washington Relay Service • Persons with a speech disability can call 877-833-6341

January 18, 2017

Mary Logue  
Kelly Moore Paint Company  
301 West Hurst Blvd.  
Hurst, TX 76053

RE: Registration with the Underground Injection Control (UIC) Program, Kelly-Moore Paint Co., 5400  
Airport Way S., Seattle, WA 98108

Dear Ms. Logue:

This letter is to acknowledge receipt of your registration form received November 22, 2016 to register the above-mentioned site with the UIC program. The UIC wells are rule authorized and do not need a permit to operate. The site is registered as UIC site number 33439.

The project includes:

- Injecting ambient air at five locations at depths between 13 to 18 feet below ground surface. Air pressure will not exceed 100 psi. The air sparging should volatize the gasoline out of the water column and a vapor extraction system will be used to collect the volatized contaminants.
- Down gradient groundwater monitoring will ensure the contaminants are not migrating offsite due to the addition of air to the system.

Air sparging is intended to improve groundwater quality. Characterize, manage, and monitor the site to minimize risk and prevent unforeseen degradation of groundwater quality. Mobilized metals or other substances, or hazardous bi-products, are not allowed to migrate beyond the site property boundary.

The two UIC Program requirements for rule authorization are, the UIC wells must be registered and the discharge from the well must meet the nonendangerment standard, of WAC 173-218-080.

Please call me at (360) 407-6143 if you have any questions. Additional information can also be found at our website <http://www.ecy.wa.gov/programs/wq/grndwtr/uic/index.html>.

Sincerely,

*Mary Shaleen Hansen*

Mary Shaleen-Hansen  
Water Quality Program

Cc: Koorus Tahghighi, Amec Foster Wheeler



**wood.**

## **Appendix E**

# Puget Sound Clean Air Agency

Notice of  
Construction No.

11291

HEREBY ISSUES AN ORDER OF APPROVAL  
TO CONSTRUCT, INSTALL, OR ESTABLISH

Registration No. 29932

Date FEB 22 2017

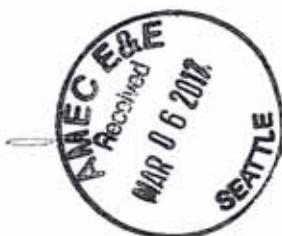
Soil and groundwater remediation site of petroleum contaminated site. This project includes an air sparge (AS) and soil vapor extraction (SVE) system equipped with either an electric catalytic oxidizer or granular activated carbon system that will be used to treat the contaminated vapors before being released to the atmosphere.

APPLICANT

Amec Foster Wheeler  
600 University St, Ste 600  
Seattle, WA 98101

OWNER

Kelly Moore Paint Company  
5400 Airport Way S  
Seattle, WA 98108



INSTALLATION ADDRESS

Kelly Moore Paint Company, 5400 Airport Way S, Seattle, WA 98108

THIS ORDER IS ISSUED SUBJECT TO THE FOLLOWING RESTRICTIONS AND CONDITIONS

1. Approval is hereby granted as provided in Article 6 of Regulation I of the Puget Sound Clean Air Agency to the applicant to install or establish the equipment, device or process described hereon at the INSTALLATION ADDRESS in accordance with the plans and specifications on file in the Engineering Division of the Puget Sound Clean Air Agency.
2. This approval does not relieve the applicant or owner of any requirement of any other governmental agency.

**Soil Remediation**

3. The owner and/or operator shall vent all vapors from the soil vapor extraction (SVE) system to a catalytic oxidizer or activated carbon control system prior to discharge to the atmosphere. When using an activated carbon control system, the system shall consist of at least two carbon canisters arranged in series. The SVE system shall not be used to process halogenated compounds or Resource Conservation Recovery Act (RCRA) hazardous waste.
4. The control efficiency of the SVE system abatement device shall meet the following requirements, as applicable:
  - a.  $\geq 97\%$  if inlet TPH  $\geq 200$  ppmv, measured as hexane or its equivalent; or
  - b.  $\geq 90\%$  if inlet TPH  $< 200$  ppmv, measured as hexane or its equivalent; or
  - c.  $\leq 10$  ppmv at the outlet of the control device, measured as hexane or its equivalent.
5. The owner or operator shall meet the following operating requirements:
  - a. The maximum influent flow rate to the abatement systems shall not exceed 300 standard cubic feet per minute (scfm).
  - b. The owner and/or operator shall only use electric power to operate the catalytic oxidizer.
  - c. The SVE system shall be operated with a minimum catalytic oxidizer temperature of at least 650

# Order of Approval for NC No. 11291

FEB 2 2017

- degrees Fahrenheit.
- d. The SVE system shall be equipped with an alarm that shuts the SVE system off when the actual temperature readings fall below the minimum operating temperature in condition 5.c.
  - 6. The catalytic oxidizer shall be equipped with continuous temperature measuring and recording instrumentation to demonstrate compliance with the minimum inlet temperature requirements of Condition No. 5.c.
  - 7. To demonstrate compliance with Condition No. 2 of this order, the owner or operator shall measure the inlet and exhaust gas streams by use of a hand held instrument capable of detecting concentrations at the levels expected, EPA Reference Method 8260B, EPA Method 8021, EPA Method TO-15 or other equivalent method approved by the agency at least once per month after initial start-up as follows:
    - a. Analyze inlet gas stream to determine the flow rate and the concentration of TPH and Benzene present.
    - b. Analyze exhaust gas to determine the flow rate, and the concentration of TPH and Benzene present.
    - c. Calculate the control efficiency based on the inlet and exhaust concentrations.
  - The SVE system shall not contain a valve or any other device which will either dilute or restrict the flow of the soil gases unless the position of the device can be measured and controlled. If a device is installed, its position must be measured and recorded any time a test sample is taken which will be used to calculate either the mass flow rate of VOCs into the atmosphere or the destruction efficiency of the control device.
  - 8. Breakthrough: During operation of the activated carbon vessels, the owner or operator shall contemporaneously monitor the gas stream by use of a hand held instrument capable of detecting concentrations at the levels expected to prevent breakthrough at least once per week at the following locations:
    - a. At the lead carbon vessel inlet;
    - b. At the inlet to the last carbon vessel in series (outlet of lead carbon vessel);
    - c. Outlet of the last carbon vessel prior to venting to the atmosphere.The owner/operator of this source may propose for Agency approval, based on actual measurements taken at the site during operation of the source, that the monitoring schedule be changed based on the decline in organic emissions and/or the demonstrated breakthrough rate of the carbon vessels.
  - 9. The owner or operator shall immediately change out the first carbon bed with unspent carbon upon breakthrough defined as the detection at its outlet of the higher of the following:
    - a. 10 % of the inlet stream concentration to the carbon vessel;
    - b. 10 ppmv measured as hexane or its equivalent.
  - 10. Spent carbon removed from the SVE system shall be stored in closed containers prior to removal from the site.
  - 11. The owner or operator may operate the soil vapor extraction system without any controls when inlet sampling data from two or more consecutive months shows the following:
    - a. Pre-control TPH emission rate is equal to or less than 2.74 lbs/day.
    - b. Pre-control Benzene emission rate is equal to or less than 0.018 lbs/day.
  - 12. The owner or operator shall maintain records of the following information:
    - a. Hours and time of operation of the extraction system and control devices.
    - b. The results of analysis or monitoring performed as required by condition 5.
    - c. The control efficiency calculation results.
    - d. A summary of the temperature readings data on a monthly basis.
    - e. When operating the activated carbon vessels, the date change out occurred and the number of carbon vessel(s) changed.

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13. The owner or operator shall report any non-compliance with any condition of this order to the Agency no later than 30 days in which it is first discovered. The owner or operator shall detail the corrective action taken and include the data showing the exceedance as well as the time of occurrence in the submittal.
14. Records required to be maintained by this Order of Approval shall be kept for at least two years from the date of generation, and made available to Puget Sound Clean Air Agency personnel upon request.

## APPEAL RIGHTS

Pursuant to Puget Sound Clean Air Agency's Regulation I, Section 3.17 and RCW 43.21B.310, this Order may be appealed to the Pollution Control Hearings Board (PCHB). To appeal to the PCHB, a written notice of appeal must be filed with the PCHB and a copy served upon Puget Sound Clean Air Agency within 30 days of the date the applicant receives this Order.



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Ralph Munoz  
Reviewing Engineer



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Carole Cenci  
Compliance Manager

# Notice of Completion

## WARNING:

Regulation I, Section 6.09, requires that the owner or applicant notify the Agency of the completion of the work covered by the application and when its operation will begin. This form is provided for your convenience to assist you in complying with this part of the Regulation.

### APPLICANT or OWNER SECTION

Mail to: Puget Sound Clean Air Agency  
Compliance Division  
1904 3rd Ave, Ste 105  
Seattle, WA 98101-3317

Notice of Construction No. 11291

Registration No. 29932

The project described below was completed on October 31, 2017.

Senior Associate    206-342-1779    11/17/17

Signature of Owner and/or Applicant

Title

Phone

Date

### FOR AGENCY USE ONLY

#### Project Description

**Soil and groundwater remediation site of petroleum contaminated site. This project includes an air sparge (AS) and soil vapor extraction (SVE) system equipped with either an electric catalytic oxidizer or granular activated carbon system that will be used to treat the contaminated vapors before being released to the atmosphere.**

#### Applicant

**Amec Foster Wheeler**  
600 University St, Ste 600  
Seattle, WA 98101

#### Owner

**Kelly Moore Paint Company**  
5400 Airport Way S  
Seattle, WA 98108

#### Location

**Kelly Moore Paint Company, 5400 Airport Way S, Seattle, WA 98108**

Inspector Check     Engineer \_\_\_\_\_ and Inspector check.

Follow up \_\_\_\_\_ (Estimated completion date plus 7)

Date Inspected \_\_\_\_\_

Inspector \_\_\_\_\_

Remarks \_\_\_\_\_

# Notice of Completion for NC No. 11291

## CONDITIONS

1. Approval is hereby granted as provided in Article 6 of Regulation I of the Puget Sound Clean Air Agency to the applicant to install or establish the equipment, device or process described hereon at the INSTALLATION ADDRESS in accordance with the plans and specifications on file in the Engineering Division of the Puget Sound Clean Air Agency.
2. This approval does not relieve the applicant or owner of any requirement of any other governmental agency.

### Soil Remediation

3. The owner and/or operator shall vent all vapors from the soil vapor extraction (SVE) system to a catalytic oxidizer or activated carbon control system prior to discharge to the atmosphere. When using an activated carbon control system, the system shall consist of at least two carbon canisters arranged in series. The SVE system shall not be used to process halogenated compounds or Resource Conservation Recovery Act (RCRA) hazardous waste.
4. The control efficiency of the SVE system abatement device shall meet the following requirements, as applicable:
  - a.  $\geq 97\%$  if inlet TPH  $\geq 200$  ppmv, measured as hexane or its equivalent; or
  - b.  $\geq 90\%$  if inlet TPH  $< 200$  ppmv, measured as hexane or its equivalent; or
  - c.  $\leq 10$  ppmv at the outlet of the control device, measured as hexane or its equivalent.
5. The owner or operator shall meet the following operating requirements:
  - a. The maximum influent flow rate to the abatement systems shall not exceed 300 standard cubic feet per minute (scfm).
  - b. The owner and/or operator shall only use electric power to operate the catalytic oxidizer.
  - c. The SVE system shall be operated with a minimum catalytic oxidizer temperature of at least 650 degrees Fahrenheit.
  - d. The SVE system shall be equipped with an alarm that shuts the SVE system off when the actual temperature readings fall below the minimum operating temperature in condition 5.c.
6. The catalytic oxidizer shall be equipped with continuous temperature measuring and recording instrumentation to demonstrate compliance with the minimum inlet temperature requirements of Condition No. 5.c.
7. To demonstrate compliance with Condition No. 2 of this order, the owner or operator shall measure the inlet and exhaust gas streams by use of a hand held instrument capable of detecting concentrations at the levels expected, EPA Reference Method 8260B, EPA Method 8021, EPA Method TO-15 or other equivalent method approved by the agency at least once per month after initial start-up as follows:
  - a. Analyze inlet gas stream to determine the flow rate and the concentration of TPH and Benzene present.
  - b. Analyze exhaust gas to determine the flow rate, and the concentration of TPH and Benzene present.
  - c. Calculate the control efficiency based on the inlet and exhaust concentrations.The SVE system shall not contain a valve or any other device which will either dilute or restrict the flow of the soil gases unless the position of the device can be measured and controlled. If a device is installed, its position must be measured and recorded any time a test sample is taken which will be used to calculate either the mass flow rate of VOCs into the atmosphere or the destruction efficiency of the control device.

8. Breakthrough: During operation of the activated carbon vessels, the owner or operator shall contemporaneously monitor the gas stream by use of a hand held instrument capable of detecting concentrations at the levels expected to prevent breakthrough at least once per week at the following locations:
  - a. At the lead carbon vessel inlet;
  - b. At the inlet to the last carbon vessel in series (outlet of lead carbon vessel);
  - c. Outlet of the last carbon vessel prior to venting to the atmosphere.

The owner/operator of this source may propose for Agency approval, based on actual measurements taken at the site during operation of the source, that the monitoring schedule be changed based on the decline in organic emissions and/or the demonstrated breakthrough rate of the carbon vessels.

## Notice of Completion for NC No. 11291

9. The owner or operator shall immediately change out the first carbon bed with unspent carbon upon breakthrough defined as the detection at its outlet of the higher of the following:
  - a. 10 % of the inlet stream concentration to the carbon vessel;
  - b. 10 ppmv measured as hexane or its equivalent.
10. Spent carbon removed from the SVE system shall be stored in closed containers prior to removal from the site.
11. The owner or operator may operate the soil vapor extraction system without any controls when inlet sampling data from two or more consecutive months shows the following:
  - a. Pre-control TPH emission rate is equal to or less than 2.74 lbs/day.
  - b. Pre-control Benzene emission rate is equal to or less than 0.018 lbs/day.
12. The owner or operator shall maintain records of the following information:
  - a. Hours and time of operation of the extraction system and control devices.
  - b. The results of analysis or monitoring performed as required by condition 5.
  - c. The control efficiency calculation results.
  - d. A summary of the temperature readings data on an monthly basis.
  - e. When operating the activated carbon vessels, the date change out occurred and the number of carbon vessel(s) changed.
13. The owner or operator shall report any non-compliance with any condition of this order to the Agency no later than 30 days in which it is first discovered. The owner or operator shall detail the corrective action taken and include the data showing the exceedance as well as the time of occurrence in the submittal.
14. Records required to be maintained by this Order of Approval shall be kept for at least two years from the date of generation, and made available to Puget Sound Clean Air Agency personnel upon request.

**wood.**

## **Appendix F**

TABLE F-1: SELECTION OF PRELIMINARY SOIL SCREENING LEVELS

Former Kelly-Moore Manufacturing Facility  
Seattle, Washington

All values are in milligrams per kilogram (mg/kg)

Compound	MTCA Cleanup Levels Protective of Human Health				MTCA Simplified TEE Industrial or Commercial Land Use	Puget Sound Regional Background Level	Final Applicable Practical Quantitation Limit	Preliminary Soil Screening Level				
	Method A - Unrestricted Land Use	Method B		Protection of Groundwater								
		Noncancer	Cancer									
<b>Metals</b>												
Arsenic	20	24	0.667	2.92	20.0	7.3	2.0	7.3				
Cadmium	2.0	80	--	0.69	36.0	0.77	0.20	0.77				
Copper	--	3,200	--	284	550	36	0.20	284				
Lead	250	--	--	3000	220	24	2.0	220				
Mercury	2.0	24	--	2.09	9.0	0.07	0.02	2.0				
Zinc	--	24,000	--	5970	570	85	1.0	570				
<b>Total Petroleum Hydrocarbons<sup>2</sup></b>												
Diesel Range Organics	2,000	--	--	--	15,000	--	5.0	2,000				
Lube Oil	2,000	--	-	--	--	--	10	2,000				
Gasoline Range Organics <sup>3</sup>	30/100	--	--	--	12,000	--	4.0	30/100				
<b>Volatile Organic Compounds</b>												
Benzene	0.03	320	18.2	0.027	--	--	0.001	0.027				
Ethylbenzene	6.00	8000	--	6.05	--	--	0.001	6.0				
Naphthalene	5.00	1600	--	4.45	--	--	0.001	4.45				
m,p-Xylene	--	--	--	13.1	--	--	0.001	13.1				
o-Xylene	--	16000	--	14.4	--	--	0.001	14.4				
Total xylenes	9.00	16000	--	--	--	--	0.002	9.0				
Toluene	7.00	6400	--	4.52	--	--	0.005	4.52				
Trichloroethene	0.03	40	12.0	0.025	--	--	0.001	0.025				
<b>Semivolatile Organic Compounds</b>												
Benzo[a]anthracene	--	--	1.37	0.858	--	--	0.0017	0.858				
Benzo[a]pyrene	0.10	24.00	0.188	2.326	300	--	0.0017	0.10				
Benzo[b]fluoranthene	--	--	1.37	2.95	--	--	0.0017	1.37				
Benzo[k]fluoranthene	--	--	13.7	29.5	--	--	0.0017	13.7				
Chrysene	--	--	137	95.5	--	--	0.0017	95.5				
Dibenz[a,h]anthracene	--	--	0.137	0.429	--	--	0.0017	0.137				
Dibenzofuran	--	80.0	--	--	--	--	0.0017	80.0				
Indeno[1,2,3-cd]pyrene	--	--	1.37	8.319	--	--	0.0017	1.37				
Total cPAHs	0.10	24.0	0.19	--	--	--	0.00	0.10				

Abbreviations

-- = not applicable

mg/kg = milligrams per kilogram

MTCA = Model Toxics Control Act

TEE = terrestrial ecological evaluation

TPH = total petroleum hydrocarbons

**TABLE F-2: SELECTION OF PRELIMINARY GROUNDWATER SCREENING LEVELS**

Former Kelly-Moore Manufacturing Facility  
Seattle, Washington

All values are in micrgrams per liter ( $\mu\text{g/L}$ )

<b>Constituents</b>	<b>Groundwater Method A</b>	<b>Groundwater Method B Non cancer</b>	<b>Groundwater Method B Cancer</b>	<b>GW (MCL - 40 CFR 141, Level)</b>	<b>GW (MCL - 40 CFR 141, NC Goal)</b>	<b>GW (MCL - Board of Health, 246-290)</b>	<b>Minimum Screening Level</b>	<b>PQLs</b>	<b>Natural Background</b>	<b>Preliminary Screening Level</b>
<b>Metals</b>										
Arsenic	5.0	4.8	0.06	10	--	10	0.06	0.5	5.0	5.0
Chromium	50	24000	--	100	100	100	50	0.2	--	50
Copper	--	640	--	1300	1300	1300	640	0.1	--	640
Lead	15	--	--	15	--	15	15	0.02	--	15.0
Mercury	2.0	2.4	--	2.0	2.0	2.0	2.0	0.1	--	2.0
Nickel	--	320	--	--	--	100	100	0.2	--	100
Zinc	--	4800	--	--	--	--	4,800	0.5	--	4,800
<b>Total Petroleum Hydrocarbons<sup>1</sup></b>										
Diesel Range Organics	500	--	--	--	--	--	500	250	--	500
Lube Oil	500	--	--	--	--	--	500	400	--	500
Gasoline Range Organics	800/1000	--	--	--	--	--	0.00	100	--	800
<b>Volatile Organic Compounds</b>										
1,2,4-Trimethylbenzene	--	80	--	--	--	--	80	0.20	--	80
1,3,5-Trimethylbenzene	--	80	--	--	--	--	80	0.20	--	80
Acetone	--	7,200	--	--	--	--	7,200	5.0	--	7,200
Benzene	5	32	0.8	5.0	--	5.0	0.8	0.20	--	0.80
Ethylbenzene	700	800	--	700	700	700	700	0.20	--	700
Naphthalene	160	160	--	--	--	--	160	1.0	--	160
m,p-Xylene	--	1,600	--	--	--	--	1,600	0.40	--	1,600
o-Xylene	--	1,600	--	--	--	--	1,600	0.20	--	1,600
Toluene	1000	640	--	1,000	1,000	1,000	640	1.0	--	640
Trichloroethene	5	4.0	0.5	5.0	--	5.0	0.54	0.20	--	0.54
Vinyl chloride	0.2	24.0	0.03	2.0	--	2.0	0.03	0.20	--	0.03
<b>Semivolatile Organic Compounds</b>										
Benzo[a]anthracene	--	--	--	--	--	--	0.00	0.01	--	--
Benzo[a]pyrene	0.10	4.8	0.023	0.20	--	0.20	0.02	0.01	--	0.02
Benzo[b]fluoranthene	--	--	--	--	--	--	0.00	0.01	--	--
Benzo[g,h,i]perylene	--	--	--	--	--	--	0.00	0.01	--	--
Benzo[k]fluoranthene	--	--	--	--	--	--	0.00	0.01	--	--
Chrysene	--	--	--	--	--	--	0.00	0.01	--	--
Dibenz[a,h]anthracene	--	--	--	--	--	--	0.00	0.01	--	--
Indeno[1,2,3-cd]pyrene	--	--	--	--	--	--	0.00	0.01	--	--
Total cPAH TEQ	0.10	4.80	0.02	0.20	--	0.20	0.02	0.01	--	0.20
<b>Polychlorinated Biphenyls</b>										
Total PCBs	--	--	0.04	0.50	--	--	0.04	0.01	--	0.04

**Abbreviations**

-- = not applicable

$\mu\text{g/L}$  = micrograms per liter

CFR = Code of Federal Regulations

cPAH = carcinogenic polycyclic aromatic hydrocarbon

PCBs = polychlorinated biphenyls

PQL = laboratory practical quantitation limit

TEQ = toxicity equivalence factor

**TABLE F-3: SELECTION OF PRELIMINARY SUB-SLAB SOIL VAPOR SCREENING LEVELS**

Former Kelly-Moore Manufacturing Facility

Seattle, Washington

All values are in micrgrams per liter ( $\mu\text{g}/\text{m}^3$ )

<b>Constituents</b>	<b>Sub-Slab Soil Gas Screening Level Method B Noncancer</b>	<b>Sub-Slab Soil Gas Screening Level Method B Cancer</b>	<b>Deep Soil Gas Screening Level Method B Noncancer</b>	<b>Deep Soil Gas Screening Level Method B Cancer</b>
<b>Metals</b>				
Arsenic	--	--	--	--
Cadmium	--	--	--	--
Chromium	--	--	--	--
Copper	--	--	--	--
Lead	--	--	--	--
Mercury	4.6	--	13.7	--
Zinc	--	--	--	--
<b>Total Petroleum Hydrocarbons<sup>1</sup></b>				
Diesel Range Organics	--	--	--	--
Lube Oil	--	--	--	--
Gasoline Range Organics	--	--	--	--
<b>Volatile Organic Compounds</b>				
Benzene	457	11	1,371	32.1
Ethylbenzene	15,238	--	45,714	--
Dibenzofuran	--	--	--	--
m,p-Xylene	1,524	--	4,571	--
o-Xylene	1,524	--	4,571	--
Toluene	76,190	--	228,571	--
Trichloroethene	30.5	12.3	91.4	37.0
<b>Semivolatile Organic Compounds</b>				
Benzo[a]anthracene	--	--	--	--
Benzo[a]pyrene	--	--	--	--
Benzo[b]fluoranthene	--	--	--	--
Benzo[g,h,i]perylene	--	--	--	--
Benzo[k]fluoranthene	--	--	--	--
Chrysene	--	--	--	--
Dibenz[a,h]anthracene	--	--	--	--
Indeno[1,2,3-cd]pyrene	--	--	--	--
Total cPAH TEQ	--	--	--	--

Abbreviations

-- = not applicable

 $\mu\text{g}/\text{m}^3$  = micrograms per cubic meter

cPAH = carcinogenic polycyclic aromatic hydrocarbon

TEQ = toxicity equivalence factor

**TABLE F-4: SELECTION OF PRELIMINARY GROUNDWATER VAPOR SCREENING LEVELS**

Former Kelly-Moore Manufacturing Facility

Seattle, Washington

All values are in micrgrams per liter ( $\mu\text{g/L}$ )

<b>Constituents</b>	<b>Groundwater Method B Non-cancer</b>	<b>Groundwater Method B Cancer</b>
<b>Metals</b>		
Arsenic	--	--
Chromium	--	--
Copper	--	--
Lead	--	--
Mercury	0.89	--
Nickel	--	--
Zinc	--	--
<b>Total Petroleum Hydrocarbons<sup>1</sup></b>		
Gasoline Range Organics	--	--
<b>Volatile Organic Compounds</b>		
1,2,4-Trimethylbenzene	28	--
1,3,5-Trimethylbenzene	--	--
Acetone	--	--
Benzene	103	2.4
Ethylbenzene	2,783	--
Naphthalene	167	8.9
m,p-Xylene	310	--
o-Xylene	440	--
Toluene	15,584	--
Trichloroethene	3.8	1.6
Vinyl chloride	56.7	0.35
<b>Semivolatile Organic Compounds</b>		
Benzo[a]anthracene	--	--
Benzo[a]pyrene	--	--
Benzo[b]fluoranthene	--	--
Benzo[g,h,i]perylene	--	--
Benzo[k]fluoranthene	--	--
Chrysene	--	--
Dibenz[a,h]anthracene	--	--
Indeno[1,2,3-cd]pyrene	--	--
Total cPAH TEQ	--	--
<b>Polychlorinated Biphenyls</b>		
Total PCBs	--	--

Abbreviations

-- = not applicable

 $\mu\text{g/L}$  = micrograms per liter

PCBs = polychlorinated biphenyls