

## **2018 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

October 8, 2019



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#### **Prepared for:**

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

#### **Prepared by:**

Wood Environment & Infrastructure Solutions, Inc. 600 University Street, Suite 600 Seattle, Washington 98101 USA T: 206-342-1760

#### October 8, 2019

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#### FORMER KELLY-MOORE MANUFACTURING FACILITY

2018 Summary of investigations and remedial actions 5400–5580 Airport Way South Seattle, Washington

October 8, 2019 Project No. 0146970060

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John Long, L.G., L.Hg. Licensed Geologist/Hydrogeologist #1354 Expiration Date: May 23, 2021



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## 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 2017 Summary of Investigations and Remedial Actions (Wood, 2018) was submitted to the Washington State Department of Ecology (Ecology) in June 2018 and provides information regarding implementation of the remedial actions discussed in the remedial investigation/feasibility study/ disproportionate cost analysis, which summarized the results of groundwater and soil sampling conducted since site activities began in 2011. Site activities conducted after June 2018 include dry and wet season groundwater sampling events, which were conducted in August 2018 and 2019, and the air sparge system began operating in May 2019. Details about the site activities conducted during this reporting period 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 in 2016.

Groundwater elevation contours for water level measurements collected in August 2018 and February 2019 are presented on Figures 2 and 3. Water level measurements collected in September 2018 and February 2019 indicate that groundwater generally flows to the west-southwest, in agreement with measurements from previous years (Wood, 2018). 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 during the 2018–2019 reporting period 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, specific 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.



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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 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-G 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 KMW-06, KMW-09, and KMW-10.

During the most recent sampling event conducted in February 2019, the concentrations of TPH-G were below the Model Toxics Control Act Method A Cleanup Level of 800 micrograms per liter ( $\mu$ g/L) (where benzene is present) except for KMW-04 and KMW-06.

TPH-D and/or TPH-O have been detected in the groundwater from all of the monitoring wells except for KMW-07 at least once since sampling began in 2011. During the most recent sampling event, the concentrations of TPH-D exceeded the Model Toxics Control Act Method A Cleanup Level of 500 µg/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. Benzene, toluene, ethylbenzene and xylenes compounds were the most frequently detected VOCs, and were predominantly detected in the central area of the site, where high concentrations of TPH-G also have been detected. These detections are most significant in the groundwater from KMW-04 and appear to have decreased in concentration over time, like the TPH-G concentrations in the groundwater from KMW-04. We expect to see VOC concentrations decrease in groundwater at KMW-04 as SVE and air sparging continue to target the western portion of the property.

Other VOC compounds detected in the groundwater are 1,2,4-trimethylbenzene and 1,3,5trimethylbenzene, 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 2018 and 2019 sampling events, which is consistent with historical results.

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## **3.3 Polycyclic aromatic hydrocarbons**

During the August 2018 and February 2019 sampling events, no PAHs were detected in the groundwater. We will continue to monitor for PAHs in groundwater during the 2019–2020 reporting period.

### 3.4 Metals

Groundwater samples were analyzed for total arsenic, chromium, copper, lead, mercury, nickel and zinc. The concentrations of metals in the groundwater samples were below cleanup levels, except for arsenic. For the 2018 and 2019 sampling events, arsenic was detected at concentrations that exceed the Ecology background level of 5.0  $\mu$ g/L at KMW-04 during the August 2018 (14.5  $\mu$ g/L) and February 2019 (17.4  $\mu$ g/L) sampling events, and at KMW-10 during the February 2019 (6.72  $\mu$ g/L) sampling event.

## 4.0 Soil vapor extraction/air sparge system operations

## 4.1 Design, installation, and operations

## 4.1.1 SVE system design & installation

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 to increase the contaminant removal rate. Both methods introduce oxygen into the subsurface, which also promotes aerobic biodegradation of residual hydrocarbons.

A series of eight horizontal SVE wells (SVE-01 through SVE-08) were installed beneath the building during redevelopment in 2015. 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. Applicable permits and construction details were included in the 2017 Summary of Investigations and Remedial Actions (Wood, 2018).

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 current configuration of the SVE and air sparge system.

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 hydrocarbons sparged from the shallow groundwater recovered by the western SVE wells. The treatment system was permitted with the Puget Sound Clean Air Agency (PSCAA) as detailed in the 2017 Summary of Investigations and Remedial Actions (Wood, 2018).

Figure 7 is a process and instrumentation diagram showing the SVE and air sparge systems and the treatment equipment. As shown, both SVE manifolds route extracted soil vapor to the CATOX treatment unit. The combined SVE and air sparging system is equipped with automatic controls and an auto-dialer that notifies Wood personnel if the CATOX system has shut down or if specific maintenance tasks are required, such as disposal of condensate water that is produced by the SVE wells. The system is equipped with a 250-gallon polyethylene tote to store condensate water produced by the SVE wells, and the tote is

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alarmed to ensure that it cannot overfill. The air sparging pump is configured to shut down immediately upon failure of the SVE system.

## 4.1.2 SVE and air sparge operations

The SVE system operated between November 2017 and October 2018. In October 2018, a vehicle crashed into the treatment compound area and the treatment system was promptly shut down until the damaged components could be inspected and repaired. The SVE system was restarted in April 2019, and operated independently until the air sparge system began operating in late May 2019. Since May 29, 2019, both the air sparge and SVE systems have been operating continuously. The SVE system is used to extract the vapors volatilized from the groundwater by the air sparge system.

## 4.2 SVE and air sparge performance evaluation

## 4.2.1 CATOX performance monitoring and regulatory compliance

Since initial startup of the SVE system in November 2017, performance monitoring vapor samples have been collected monthly from the CATOX influent vapor stream sampling port and at the effluent sample port on the emissions stack. Field forms can be found in Appendix C. In compliance with PSCAA Registration No. 29932, monthly performance monitoring samples are field-analyzed by a flame ionization detector (FID) calibrated to 100 parts per million hexane. Monthly performance monitoring samples are also submitted to a local laboratory for additional analysis of benzene. Additionally, samples are collected from each SVE well monthly and are field-analyzed by FID. Individual SVE well vapor samples were submitted for laboratory analysis of benzene during SVE startup (November 2017) and during and after air sparging startup (May 2019 and June 2019). Analytical data for individual SVE well samples are provided in Table 5. All FID results are shown in Tables 6A and 6B. All analytical results for SVE performance monitoring are provided in Appendix D.

CATOX performance is determined by its destruction removal efficiency (DRE) for removed soil vapor compounds. DRE is calculated from FID results at the CATOX influent and at the effluent emissions stack and are shown in Table 7. DREs generally exceed 97 percent and demonstrate adequate system performance. Despite generally high DREs, reported DREs fell below 97 percent during the summer of 2018. On-site investigations and communication with the CATOX manufacturer did not identify any operational issues. PSCAA issued a Notice of Violation letter on April 17, 2019, due to the decreased destruction efficiencies in mid-2018. On April 25, 2019, Wood provided a response to the Notice of Violation, including a plan for future site actions that will be taken in the event that the DRE again falls below levels required by PSCAA Registration No. 29932. This corrective action memo is included as Appendix E.

## 4.2.2 SVE and air sparge optimization and performance monitoring

Table 6a provides FID readings for SVE concentrations prior to air sparge sparge startup, and Table 6b provides FID readings for SVE concentrations with air sparging. While the air sparge system has been running continuously since May 2019, in the future we may choose to cycle the air sparge system to increase mixing in the subsurface if SVE concentrations decline.

Since 2017, an estimated 8,400 pounds of TPH (as hexane equivalent) have been removed from the subsurface by the SVE in conjunction with the air-sparge system. Table 8 summarizes the performance data, and Figure 8 shows the mass removal rate over time as estimated from site performance data and FID measurements. The mass removal rate was calculated from the measured influent concentration and system flow rate during each site visit, and is based on the total system runtime between site visits. Thus,



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the removal rate is high when influent concentration is high (such as during air sparge startup) and low when influent concentrations are low (such as at the end of the summer in 2018).

Mass removal rates are highest when SVE concentrations are high during periods of low groundwater elevations from June through October. We continue to optimize the flow from the wells to maximize the concentration of hydrocarbons at the CATOX influent. Individual well concentrations over time are shown in Figure 9. Since startup, the highest-concentration SVE wells have been SVE-03, SVE-05, SVE-10, and SVE-12. After air sparging began in May 2019, concentrations increased in all western wells except for SVE-09. We will continue to monitor SVE well concentrations and target high-concentration areas with SVE and air sparging to improve mass removal rates.

High site groundwater levels during the winter reduce the vadose zone volume and SVE influent concentrations. SVE concentrations at all wells fluctuate with the seasonal cycle, which may seasonally reduce SVE mass removal rates. Wet season operations are less efficient due to intermittent alarm conditions and periodic shutdowns associated with increased condensate production at the CATOX knockout pot. We will continue to operate the SVE and air sparge systems during the wet season and will work to optimize mass removal during dry season operations.

In addition to FID monitoring in the western parking lot during air sparging startup, indoor air samples were collected inside the northern building both prior to and after the startup of the air sparging system. Sample locations are shown on Figure 4. Samples were analyzed for VOCs using EPA Method TO-15 and analytical results are provided in Appendix F. As shown in Table 9, VOC concentrations during June 2019 (while air sparging) are comparable to and generally lower than compound concentrations during October 2018 (before air sparging). Both sets of analytical results meet permissable exposure limits for indoor air, suggesting that vapor intrusion from contaminated groundwater or volatilized contamination was not a concern on either day that indoor air sampling occurred.

## 5.0 Conclusions and recommendations

The following actions will be conducted before the end of 2019:

- Groundwater samples were collected for the dry season sampling event in August 2019.
- SVE and air sparging system inspections (including performance monitoring sampling) will occur at least monthly. On-site personnel will continue to optimize SVE and air sparging operating conditions in order to maximize mass removal rates and CATOX performance.
- Kelly-Moore and Ecology would like to continue working together to take the necessary steps to eventually obtain "No Further Action" for the site.

## 6.0 References

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 Environment & Infrastructure Solutions, Inc. (Wood), 2018, 2017 Summary of Investigations and Remedial Actions, Former Kelly-Moore Manufacturing Facility, 5400-5580 Airport Way South, Seattle, Washington, June 5.













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#### **TABLE 1: MONITORING WELL SURVEY DATA**

#### Former Kelly-Moore Manufacturing Facility, Seattle, Washington

	WCS No	orth Zone		
	NAD 8	33(91) <sup>3</sup>	Ground Surface	Top of Casing
Well Name	Northing	Easting	Elevation	Elevation
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 & Associates.
- 2. Survey completed on December 13, 2016, by Duane Hartman & Associates.
- 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>
	$21 \text{ C2}^2$	8/15/2018	9.96	11.67
KIVIVV-UZR	21.03	2/7/2019	9.17	12.46
	21 F 4 <sup>2</sup>	8/15/2018	9.93	11.61
KIVIVV-USK	21.54	2/7/2019	9.37	12.17
KN4\\4/_04	19 E6 <sup>2</sup>	8/15/2018	7.06	11.50
KIVIV-04	10.50	2/7/2019	6.60	11.96
	10 90 <sup>2</sup>	8/15/2018	8.29	11.51
	19.00	2/7/2019	7.77	12.03
KN4\\4/_07	21 62 <sup>2</sup>	8/15/2018	9.96	11.67
	21.05	2/7/2019	9.21	12.42
KVV/V-08	21 65 <sup>2</sup>	8/15/2018	10.00	11.65
KIVIV-00	21.05	2/7/2019	9.31	12.34
KV4/V/-00	10 14 <sup>3</sup>	8/15/2018	6.64	11.50
KIVIV-03	10.14	2/7/2019	6.15	11.99
KN4\4/_10	20 20 <sup>3</sup>	8/15/2018	9.01	11.38
	20.59	2/7/2019	8.65	11.74

Notes:

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

<sup>1.</sup> Elevations in feet above mean sea level.

Surveys conducted on June 30, 2016. Vertical Datum: NAVD88. Benchmark: 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

#### **TABLE 3: GROUNDWATER PARAMETERS**

Former Kelly-Moore Manufacturing Facility, Seattle, Washington

			DO	SC	ORP	Temperature	Turbidity
Monitoring Well	Date	рН	(mg/L)	(ms/cm)	(mv)	(°C)	(NTU)
	8/16/2018	5.85	0.55	0.173	274	14.5	30.8
KIVIV-UZK	2/8/2019	5.95	0.81	0.245	130.1	14.1	8.1
	8/16/2018	7.03	0.47	0.378	112	18.3	10
KIVIV-USK	2/8/2019	6.97	0.51	0.582	-87.0	16.5	20.52
	8/16/2018	6.09	0.63	0.326	99.0	18	7.9
NIVI VV-04	2/7/2019	6.22	0.62	0.341	-74.0	11.2	12.01
	8/16/2018	6.33	0.37	0.421	-39.0	19.5	18.4
KIVIVV-00	2/7/2019	6.18	0.65	0.635	-32.0	14.2	12.4
	8/16/2018	6.02	0.6	0.211	268	15.4	31
	2/8/2019	6.23	0.52	0.318	51.1	14.4	17.91
	8/16/2018	5.95	0.58	0.211	248	15.7	10.1
	2/8/2019	6.05	0.74	0.250	91.4	13.9	30.9
	8/16/2018	6.35	0.47	0.387	-24	18.4	12.2
	2/7/2019	6.47	0.57	0.400	-69	15.2	4.99
KN/W/ 10	8/16/2018	6.25	0.46	0.416	-15	16.5	21.8
	2/7/2019	6.53	0.49	0.430	-82	13.9	15.6

**Abbreviations** 

°C = degrees Celsius

DO = dissolved oxygen

mg/L = milligrams per liter

ms/cm = milli Siemens per centimeter

mv = millivolts

NTU = nephelometric turbidity unit

ORP = oxidation reduction potential

SC = specific conductivity

## TABLE 4: GROUNDWATER ANALYTICAL RESULTS<sup>1, 2</sup>

Former Kelly-Moore Manufacturing Facility, Seattle, Washington

			KMW-01						KMW-02/02R	3										KMW-03/0	3R <sup>4</sup>					
Ī															3/28/2011							11/10/2016				
Analyte	Preliminary Screening Level	3/28/2011	8/4/2011	6/7/2013	3/28/2011	8/4/2011	6/7/2013	6/30/2016	11/10/2016	9/1/2017	1/26/2018	8/16/2018	2/8/2019	3/28/2011	(D)	8/4/2011	4/4/2013	6/7/2013	3/10/2015	7/1/2016	11/10/2016	(D)	9/1/2017	1/26/2018	8/16/2018	2/8/2019
Total Metals (µg/L)																										
Arsenic	5	3.3 U	3.3 U		3.3 U	3.3 U		3.3 U	3.3 U	3.3 U	3.3 U	1.0 U	1.0 U	3.3 U	3.3 U	3.3 U				3.3 U	3.3 U	3.3 U	3.3 U	3.3 U	1.0 U	1.0 U
Chromium	100	11 U			11 U			11 U	11 U	11 U	11 U	1.0 U	1.0 U	11 U	11 U					11 U	11 U	11 U	11 U	11 U	1.0 U	1.47
Copper	640		11 U			11 U						5.0 U	5.0 U			11 U									5.0 U	5.0 U
Lead	15	1.1 U	1.1 U		1.1 U	1.1 U		1.1 U	1.1 U	1.1 U	1.1 U	1.0 U	1.0 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	1.0 U	1.0 U
Mercury	2.0							0.5 U	0.50 U	0.50 U	0.50 U	1.0 U	1.0 U							0.5 U	0.50 U	0.50 U	0.50 U	0.50 U	1.0 U	1.0 U
Nickel	100											1.82	2.15												1.46	1.67
Zinc	4800											5.0 U	5.0 U												5.0 U	5.0 U
Polycyclic Aromatic Hydrocarbons (µg/L)																										
Benzo(a)anthracene		0.0098 U	0.0095 U		0.0098 U	0.0096 U		0.0095 U	0.0094 U	0.0099 U	0.011 U	0.06 U	0.04 U	0.0096 U	0.0095 U	0.0095 U				0.018	0.0095 U	0.0095 U	0.0098 U	0.011 U	0.06 U	0.04 U
Benzo(a)pyrene	0.023	0.0098 U	0.0095 U		0.0098 U	0.0096 U		0.0095 U	0.0094 U	0.0099 U	0.011 U	0.06 U	0.04 U	0.0096 U	0.0095 U	0.0095 U				0.011	0.0095 U	0.0095 U	0.0098 U	0.011 U	0.06 U	0.04 U
Benzo(b)fluoranthene		0.0098 U	0.0095 U		0.0098 U	0.0096 U		0.0095 U	0.0094 U	0.0099 U	0.011 U	0.06 U	0.04 U	0.0096 U	0.0095 U	0.0095 U				0.011	0.0095 U	0.0095 U	0.0098 U	0.011 U	0.06 U	0.04 U
Benzo(ghi)perylene		0.0098 U	0.0095 U		0.0098 U	0.0096 U		0.0095 U	0.0094 U	0.0099 U	0.011 U	0.06 U	0.04 U	0.0096 U	0.0095 U	0.0095 U				0.014	0.0095 U	0.0095 U	0.0098 U	0.011 U	0.06 U	0.04 U
Benzo(j,k)fluoranthene		0.0098 U	0.0095 U		0.0098 U	0.0096 U		0.0095 U	0.0094 U	0.0099 U	0.011 U	0.06 U	0.04 U	0.0096 U	0.0095 U	0.0095 U				0.0095 U	0.0095 U	0.0095 U	0.0098 U	0.011 U	0.06 U	0.04 U
Chrysene		0.0098 U	0.0095 U		0.0098 U	0.0096 U		0.0095 U	0.0094 U	0.0099 U	0.011 U	0.06 U	0.04 U	0.0096 U	0.0095 U	0.0095 U				0.012	0.0095 U	0.0095 U	0.0098 U	0.011 U	0.06 U	0.04 U
Dibenz(a,h)anthracene		0.0098 U	0.0095 U		0.0098 U	0.0096 U		0.0095 U	0.0094 U	0.0099 U	0.011 U	0.06 U	0.04 U	0.0096 U	0.0095 U	0.0095 U				0.0095 U	0.0095 U	0.0095 U	0.0098 U	0.011 U	0.06 U	0.04 U
Indeno(1,2,3-cd)pyrene		0.0098 U	0.0095 U		0.0098 U	0.0096 U		0.0095 U	0.0094 U	0.0099 U	0.011 U	0.06 U	0.04 U	0.0096 U	0.0095 U	0.0095 U				0.011	0.0095 U	0.0095 U	0.0098 U	0.011 U	0.06 U	0.04 U
Total cPAHs	0.20	0.007 U	0.007 U		0.007 U	0.007 U		0.007 U	0.007 U	0.007 U	0.008 U	0.05 U	0.03 U	0.007 U	0.007 U	0.007 U				0.016	0.007 U	0.007 U	0.007 U	0.008 U	0.05 U	0.03 U
Polychlorinated Biphenyls (µg/L)																										
Aroclor 1016								0.047 U												0.047 U						
Aroclor 1221								0.047 U												0.047 U						
Aroclor 1232								0.047 U												0.047 U						
Aroclor 1242								0.047 U												0.047 U						
Aroclor 1248								0.047 U												0.047 U						
Aroclor 1254								0.047 U												0.047 U						
Aroclor 1260	0.04							0.047 U												0.047 U						
Total PCBs	0.44							0.047 U												0.047 U						
Total Petroleum Hydrocarbons (µg/L)																										
Gasoline Range Organics	500	100 U	100 U	100 U	100 U	100 U	100 U	500 U	100 U	100 U	100 U	100 U	100 U	7,700	7,000	6,100	1,800	1,100	1,100	300	130	170	270	150	290	140
Diesel Range Organics	500							260	260 U	0.27 U	260 U	50 U	60 U							660 J	310 U	280 U	330	510	350	1,700
Lube Oil	800							410 U	410 U	0.43 U	410 U	250 U	300 U							410 U	410 U	410 U	430 U	410 U	250 U	300 U
Volatile Organic Compounds (µg/L)																										
1,2,4-Trimethylbenzene	80	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	0.20 U	1.0 U	1.0 U				1.0 U	0.20 U		0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U
1,3,5-Trimethylbenzene	80	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	0.20 U	1.0 U	1.0 U				1.0 U	0.20 U		0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U
Acetone	7200	5.0 U	5.0 U	5.0 U	5.0 U	5.0 U	5.0 U	12 U	5.0 U	5.0 U	5.0 U	50 U	50 U				25 U	5.0 U		12 U	7.9	6.5	5.0 U	5.0 U	50 U	50 U
Benzene	5.0	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	0.20 U	0.35 U	0.35 U	8.1	8.4	4.0 U	1.0 U	0.41		0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.35 U	0.35 U
Ethylbenzene	700	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	0.20 U	1.0 U	1.0 U	3,100	2,700	2,400	170	27		1.4	0.33	0.34	0.20 U	0.20 U	1.0 U	1.0 U
m,p-Xylene	1600	0.40 U	0.40 U	0.40 U	0.40 U	0.40 U	0.40 U	0.40 U	0.40 U	0.40 U	0.40 U	2.0 U	2.0 U	18	18	7.1	2.0 U	0.4 U		2.1	1.0	0.97	0.40 U	0.40 U	2.0 U	2.0 U
Naphthalene	160	1.0 U	1.0 U	1.4 U	1.0 U	1.0 U	1.4 U	1.0 U	1.3 U	1.0 U	1.0 U	0.06 U	0.4 U				19	9.8		1.0 U	1.3 U	1.3 U	1.0 U	1.0 U	0.078	0.4 U
o-Xylene	1600	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	0.20 U	1.0 U	1.0 U	4.0 U	4.0 U	4.0 U	1.0 U	0.20 U		0.35	0.35	0.37	0.20 U	0.20 U	1.0 U	1.0 U
Toluene	640	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	4.0 U	4.0 U	4.0 U	5.0 U	1.0 U		1.1	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Trichloroethene	4.0	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	0.20 U	1.0 U	1.0 U				1.0 U	0.20 U		0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U
Vinyl Chloride	0.29	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	0.20 U	0.20 U	0.20 U				1.0 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

## TABLE 4: GROUNDWATER ANALYTICAL RESULTS<sup>1, 2</sup>

Former Kelly-Moore Manufacturing Facility, Seattle, Washington

						км	W-04							KMW-05				кмм	/-06					кмм	-07		
			8/4/2011				6/30/2016						1														
Analyte	3/28/2011	8/4/2011	(D)	6/7/2013	3/10/2015	6/30/2016	(D)	11/11/2016	8/31/2017	1/25/2018	8/16/2018	2/7/2019	3/28/2011	8/4/2011	6/7/2013	6/30/2016	11/11/2016	8/31/2017	1/24/2018	8/16/2018	2/7/2019	7/1/2016	11/10/2016	9/1/2017	1/26/2018	8/16/2018	2/8/2019
Total Metals (µg/L)																											
Arsenic	12	12	10			12	12	20	14	7.6	14.5	17.4	3.3 U	3.3 U		3.5	3.4	3.3 U	3.3 U	4.83	3.04	3.3 U	3.3 U	3.3 U	3.3 U	1.0 U	1.0 U
Chromium	11 U					11 U	11 U	11 U	11 U	11 U	1.0 UJ	1.58	11 U			11 U	11 U	11 U	11 U	2.48 J	2	11 U	11 U	11 U	11 U	1.0 U	1.0 U
Copper		11 U	11 U								5.0 UJ	24.7		11 U						11.9 J	19.4					5.0 U	5.0 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	1.0 U	1.51	1.1 U	1.1 U		3.7	1.1	2.7	7.3	5.61	2.04	1.1 U	1.1 U	1.1 U	1.1 U	1.0 U	1.0 U
Mercury						0.5 U	0.5 U	0.50 U	0.50 U	0.50 U	1.0 U	1.0 U				0.5 U	0.50 U	0.50 U	0.50 U	1.0 U	1.0 U	0.5 U	0.50 U	0.50 U	0.50 U	1.0 U	1.0 U
Nickel											2.43 J	1.98								1.38 J	7.79					1.0 U	1.36
Zinc											5.0 UJ	5.0 U								5.72 J	135					5.0 U	5.0 U
Polycyclic Aromatic Hydrocarbons (µg/L)																											
Benzo(a)anthracene	0.012 U	0.0095 U	0.0095 U			0.011	0.014	0.0094 U	0.0096 U	0.010 U	0.06 U	0.04 U	0.0095 U	0.0095 U		0.047	0.013	0.041	0.055	0.06 U	0.04 U	0.0095 U	0.0095 U	0.0099 U	0.011 U	0.06 U	0.04 U
Benzo(a)pyrene	0.012 U	0.0095 U	0.0095 U			0.0095 UJ	0.015 J	0.0094 U	0.0096 U	0.010 U	0.06 U	0.04 U	0.0095 U	0.0095 U		0.038	0.022	0.033	0.071	0.06 U	0.04 U	0.0095 U	0.0095 U	0.0099 U	0.011 U	0.06 U	0.04 U
Benzo(b)fluoranthene	0.012 U	0.0095 U	0.0095 U			0.0095 UJ	0.022 J	0.0094 U	0.0120	0.010 U	0.06 U	0.04 U	0.0095 U	0.0095 U		0.047	0.021	0.034	0.082	0.06 U	0.04 U	0.0095 U	0.0095 U	0.0099 U	0.011 U	0.06 U	0.04 U
Benzo(ghi)perylene	0.012 U	0.0095 U	0.0095 U			0.0095 UJ	0.023 J	0.0094 U	0.0096 U	0.010 U	0.06 U	0.04 U	0.0095 U	0.0095 U		0.041	0.024	0.025	0.056	0.06 U	0.04 U	0.0095 U	0.0095 U	0.0099 U	0.011 U	0.06 U	0.04 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	0.06 U	0.04 U	0.0095 U	0.0095 U		0.018	0.031	0.018	0.034	0.06 U	0.04 U	0.0095 U	0.0095 U	0.0099 U	0.011 U	0.06 U	0.04 U
Chrysene	0.012 U	0.0095 U	0.0095 U			0.0095 U	0.0095 U	0.0094 U	0.0100	0.010 U	0.06 U	0.04 U	0.0095 U	0.0095 U		0.035	0.028	0.032	0.062	0.06 U	0.04 U	0.0095 U	0.0095 U	0.0099 U	0.011 U	0.06 U	0.04 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	0.06 U	0.04 U	0.0095 U	0.0095 U		0.0095 U	0.0095 U	0.0095 U	0.0110	0.06 U	0.04 U	0.0095 U	0.0095 U	0.0099 U	0.011 U	0.06 U	0.04 U
Indeno(1,2,3-cd)pyrene	0.012 U	0.0095 U	0.0095 U			0.0095 UJ	0.016 J	0.0094 U	0.0096 U	0.010 U	0.06 U	0.04 U	0.0095 U	0.0095 U		0.028	0.023	0.023	0.054	0.06 U	0.04 U	0.0095 U	0.0095 U	0.0099 U	0.011 U	0.06 U	0.04 U
Total cPAHs	0.009 U	0.007 U	0.007 U			0.008	0.020 J	0.007 U	0.008	0.008 U	0.05 U	0.03 U	0.007 U	0.007 U		0.050	0.030	0.045	0.095	0.05 U	0.03 U	0.007 U	0.007 U	0.007 U	0.008 U	0.05 U	0.03 U
Polychlorinated Biphenyls (µg/L)																											
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					
Total PCBs						0.048 U	0.047 U									0.047 U						0.047 U					
Total Petroleum Hydrocarbons (µg/L)																											
Gasoline Range Organics	75,000	55,000	50,000	48,000	27,000	27,000	27,000	63,000	8,000	9,000	33,000	31,000	100 U	100 U	100 U	2,700	850	1,600	1,300	4,000	2,200	500 U	100 U	100 U	100 U	100 U	100 U
Diesel Range Organics						3,000 J	2,700 J	6,400 U	1,600	1,700	2,000	2,600				5,400 J	3,500	4,400	4,200	8,600	19,000	260 U	260 U	280 U	260 U	50 U	60 U
Lube Oil						510	870	410 U	440 U	410 U	250 U	300 U				1,500 J	1,200	1,600	600	680	790	410 U	420 U	450 U	410 U	250 U	300 U
Volatile Organic Compounds (µg/L)																											
1,2,4-Trimethylbenzene				77		44 J	65 J	160	97	110	54	67			0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U
1,3,5-Trimethylbenzene				20 U		20 U	20	52	30	50 U	17	33			0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U
Acetone				500 U		1,200 U	1,200 U	1,000 U	500 U	1300 U	50 U	500 U			5.0 U	12 U	5.0 U	5.0 U	5.0 U	50 U	50 U	12 U	6.5	5.0 U	5.0 U	50 U	50.0 U
Benzene	10	13	13	20 U		20 U	20 U	20 U	20 U	50 U	0.35 U	3.5 U	1.0 U	1.0 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.35 U	0.35 U	0.20 U	0.20 U	0.20 U	0.20 U	0.35 U	0.35 U
Ethylbenzene	5,700	3,700	3,400	3,400		3,700	4,300	5,200	4,300	4,700	2,600	2,800	1.0 U	1.0 U	0.20 U	0.38	0.25	0.27	0.20 U	1.0 U	1.0 U	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U
m,p-Xylene	12,000	8,500	7,700	6,800		7,100	7,900	12,000	7,800	12,000	6,400	6,100	1.0 U	1.0 U	0.4 U	1.4	0.92	1.6	0.42	2.0 U	2.0 U	0.4 U	0.44	0.40 U	0.40 U	2.0 U	2.0 U
Naphthalene				140 U		100 U	100 U	100 U	100 U	250 U	5.1	3.3			1.4 U	1.0 U	1.3 U	1.0 U	1.0 U	0.16	0.4 U	1.0 U	1.3 U	1.0 U	1.0 U	0.06 U	0.4 U
o-Xylene	3,400	2,100	1,900	2,200		1,700	1,700	3,600	1,900	3,600	1,500	1,300	1.0 U	1.0 U	0.20 U	0.64	0.49	0.47	0.20 U	1.0 U	1.0 U	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U
Toluene	7,400	5,800	5,500	3,800		1,400	1,300	5,300	980	5,500	610	190	1.0 U	1.0 U	10	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Trichloroethene				20 U		20 U	20 U	20 U	20 U	50 U	1.0 U	10 U			0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U
Vinyl Chloride				20 U		2.0 U	2.0 U	2.0 U	20 U	50 U	0.20 U	2.0 U			0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.2 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.2 U

## TABLE 4: GROUNDWATER ANALYTICAL RESULTS<sup>1, 2</sup>

## Former Kelly-Moore Manufacturing Facility, Seattle, Washington

			кми	V-08					KMW-09					кмм	/-10		
															1/25/2018		
Analyte	7/1/2016	11/10/2016	9/1/2017	1/26/2018	8/16/2019	2/8/2019	11/11/2016	8/31/2017	1/24/2018	8/16/2018	2/7/2019	11/11/2016	8/31/2017	1/25/2018	(D)	8/16/2018	2/7/2019
Total Metals (μg/L)																	
Arsenic	3.3 U	3.3 U	3.3 U	3.3 U	1.0 U	1.0 U	3.3 U	3.3 U	3.3 U	2.04	1.0 U	9.1	10	6.8	5.7	4.61	6.72
Chromium	11 U	11 U	11 U	11 U	1.0 U	1.0 U	11 U	11 U	11 U	1.40 J	1.12	11 U	11 U	11 U	11 U	1.35 J	2.00 J
Copper					5.0 U	5.0 U				5.62 J	5.0 U					5.0 UJ	5.0 UJ
Lead	1.1 U	1.1 U	1.1 U	1.1 U	1.0 U	1.0 U	1.1 U	1.1 U	3.0	3.1	1.0 U	1.1 U	1.1 U	1.1	1.1 U	1.0 U	1.0 U
Mercury	0.5 U	0.50 U	0.50 U	0.50 U	1.0 U	1.0 U	0.50 U	0.50 U	0.50 U	1.0 U	1.0 U	0.50 U	0.50 U	0.50 U	0.50 U	1.0 U	1.0 U
Nickel					1.11	1.99				1.25 J	1.0 U					1.0 UJ	1.0 UJ
Zinc					5.0 U	10.1				5.0 UJ	5.0 U					5.0 UJ	5.0 UJ
Polycyclic Aromatic Hydrocarbons (µg/L)																	
Benzo(a)anthracene	0.086	0.015	0.012	0.018	0.06 U	0.04 U	0.0098 U	0.0098 U	0.020	0.06 U	0.04 U	0.0098 U	0.0094 U	0.011 U	0.011 U	0.06 U	0.04 U
Benzo(a)pyrene	0.11	0.013	0.010 U	0.015	0.06 U	0.04 U	0.0098 U	0.0098 U	0.021	0.06 U	0.04 U	0.0098 U	0.0094 U	0.011 U	0.011 U	0.06 U	0.04 U
Benzo(b)fluoranthene	0.12	0.020	0.010 U	0.018	0.06 U	0.04 U	0.0098 U	0.0098 U	0.017	0.06 U	0.04 U	0.0098 U	0.0094 U	0.011 U	0.011 U	0.06 U	0.04 U
Benzo(ghi)perylene	0.1	0.0095 U	0.010 U	0.012 U	0.06 U	0.04 U	0.0098 U	0.0098 U	0.011	0.06 U	0.04 U	0.0098 U	0.0094 U	0.011 U	0.011 U	0.06 U	0.04 U
Benzo(j,k)fluoranthene	0.046	0.0095 U	0.010 U	0.012 U	0.06 U	0.04 U	0.0098 U	0.0098 U	0.015	0.06 U	0.04 U	0.0098 U	0.0094 U	0.011 U	0.011 U	0.06 U	0.04 U
Chrysene	0.09	0.042	0.012	0.028	0.06 U	0.04 U	0.0098 U	0.0098 U	0.018	0.06 U	0.04 U	0.0098 U	0.0094 U	0.011 U	0.011 U	0.06 U	0.04 U
Dibenz(a,h)anthracene	0.024	0.0095 U	0.010 U	0.012 U	0.06 U	0.04 U	0.0098 U	0.0098 U	0.011 U	0.06 U	0.04 U	0.0098 U	0.0094 U	0.011 U	0.011 U	0.06 U	0.04 U
Indeno(1,2,3-cd)pyrene	0.063	0.0095 U	0.010 U	0.012 U	0.06 U	0.04 U	0.0098 U	0.0098 U	0.0130	0.06 U	0.04 U	0.0098 U	0.0094 U	0.011 U	0.011 U	0.06 U	0.04 U
Total cPAHs	0.14	0.018	0.008	0.021	0.05 U	0.03 U	0.007 U	0.007 U	0.028	0.05 U	0.03 U	0.007 U	0.007 U	0.008 U	0.008 U	0.05 U	0.03 U
Polychlorinated Biphenyls (µg/L)																	
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																
Total PCBs	0.047 U																
Total Petroleum Hydrocarbons (µg/L)																	
Gasoline Range Organics	1,000	400	130	120	230	120	370	360	760	940	450	110	3,400	270	260	4,800	200
Diesel Range Organics	770 J	370 U	300 U	450	160	440	1,700	2,300	3,100	3,600	3,100	1,300 U	1,800	2,300	2,300	1,400	970
Lube Oil	410 U	410 U	480 U	410 U	250 U	300 U	660	810	690	360	300 U	420 U	430 U	410 U	410 U	250 U	320 U
Volatile Organic Compounds (µg/L)																	
1,2,4-Trimethylbenzene	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	0.2 U	0.20 U	0.20 U	1.0 U	1.0 U	3.7	53	2.7	5.4	38	1.0 U
1,3,5-Trimethylbenzene	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	0.2 U	0.20 U	0.20 U	1.0 U	1.0 U	0.38	23	0.9	1.7	19	1.0 U
Acetone	12 U	10	5.0 U	5.0 U	50 U	50.0 U	5.0 U	5.0 U	5.0 U	50 U	50 U	5.0 U	100 U	5.0 U	5.0 U	50 U	50 U
Benzene	0.20 U	0.20	0.20 U	0.20 U	0.35 U	0.35 U	0.2 U	0.20 U	0.20 U	0.35 U	0.35 U	0.7	8.2	0.20 U	0.20 U	1.5	0.35 U
Ethylbenzene	0.20 U	0.31	0.36	0.20 U	1.0 U	1.0 U	3.1	0.20 U	0.44	1.0 U	1.0 U	1.6	810	14	30	370	1.0 U
m,p-Xylene	0.4 U	0.76	0.69	0.40 U	2.0 U	2.0 U	0.51	0.40 U	0.40 U	2.0 U	2.0 U	11	1100	28	65	1100	6.0
Naphthalene	1.5	1.3 U	1.0 U	1.0 U	0.06 U	0.4 U	1.3 U	1.0 U	1.0 U	0.12	1.0 U	1.3 U	20 U	1.0 U	1.0 U	0.51	0.4 U
o-Xylene	0.20 U	0.34	0.20 U	0.20 U	1.0 U	1.0 U	0.2 U	0.20 U	0.20 U	1.0 U	1.0 U	0.29	22	0.42	0.58	1.0 U	1.0 U
Toluene	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	20 U	1.0	1.0	1.0 U	1.0 U
Trichloroethene	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	0.2 U	0.20 U	0.20 U	1.0 U	1.0 U	0.2 U	4.0 U	0.20 U	0.20 U	1.0 U	1.0 U
Vinyl Chloride	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.2 U	0.2 U	0.20 U	0.20 U	0.20 U	0.2 U	0.2 U	4.0 U	0.20 U	0.20 U	0.20 U	0.2 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.
- and was replaced Julie 20, 20
- 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, 3</sup>

Former Kelly-Moore Manufacturing Facility, Seattle, Washington

		Benzene	TPH as Hexane
Sample	Date <sup>4</sup>	(ppmv)	(ppmv)
	Western SVE \	Nells	
SVE-09	11/7/17	0.069	310
	5/30/19	<0.1	630
	6/4/19	<0.1	440
SVE-10	11/7/17	0.53	820 J
	5/30/19	< 0.5	3,500
	6/4/19	< 0.5	2,300
SVE-11	11/7/17	0.069	220
	5/30/19	<0.1	1,300
	6/4/19	<0.1	660
SVE-12	11/7/17	0.44	1,400 J
	5/30/19	<0.1	3,300
	6/4/19	<0.2	1,400
SVE-13	11/7/17	0.23	600 J
	5/30/19	<0.1	2,100
	6/4/19	<0.1	760
	Eastern SVE V	Vells	
SVE-02	11/7/17	< 0.03	3.4
	5/30/19	<0.1	<10
	6/4/19	<0.1	14
SVE-04	11/7/17	< 0.03	310
	5/30/19	<0.1	470
	6/4/19	<0.1	400
SVE-06	11/7/17	0.041	280
	5/30/19	<0.1	36
	6/4/19	<0.1	33
SVE-08	11/7/17	< 0.03	65
	5/30/19	<0.1	30
	6/4/19	<0.1	16
SVE-07	11/7/17	< 0.03	82
	5/30/19	<0.1	70
	6/4/19	<0.1	230
SVE-05	11/7/17	0.50	2,200 J
	5/30/19	<0.2	5,100
	6/4/19	< 0.5	3,500
SVE-03	11/7/17	1.1	1,900 J
	5/30/19	<0.2	1,900
	6/4/19	<0.1	2,400
SVE-01	11/7/17	0.14	450
	5/30/19	<0.1	10
	6/4/19	<0.1	14
Influent	11/7/17	0.18	650 J
	5/30/19	<0.1	1,100
	6/4/19	<0.1	640
Effluent	11/7/17	< 0.03	28
	5/30/19	<0.1	41
	6/4/19	< 0.1	20

#### Notes:

- Data qualifiers are as follows:
   J = the result is estimated because the
- concentration exceeded the calibration range of the instrument.
- 2. **Bold** values indicate results above the reporting limits.
- 3. Analytes that were not detected above the method detection limit are listed as less than the detection limit.
- 4. The SVE system began operating without air sparging on November 7, 2017. The SVE did not operate between October 16, 2018 and April 1, 2019. SVE system with air sparge began operating on May 29, 2019.

#### Abbreviations:

- ppmv = parts per million volume
- TPH = total petroleum hydrocarbons
- SVE = soil vapor extraction

#### TABLE 6a : FLAME IONIZATION DETECTOR READINGS, SVE OPERATIONS WITHOUT AIR SPARGING

Former Kelly-Moore Manufacturing Facility, Seattle, Washington

Well								FID	Readings <sup>1</sup> (pp	omv)							
	11/7/2017	12/13/2017	1/16/2018	2/13/2018	3/21/2018	4/24/2018	5/30/2018	6/29/2018	7/31/2018	8/31/2018	9/19/2018	10/16/2018	4/1/2019	4/3/2019	5/3/2019	5/6/2019	5/29/2019
							١	Western SVE V	Vells								
West Manifold	331	19.4	NM	NM	NM	NM	8.3	NM	NM	191	321	273.4	NM	87	2.2	NM	13.2
SVE-09	NM	NM	0	53.5	53.5	7.6	7.1	246.3	505.3	118	253	154.2	NM	162.1	1.9	NM	12.1
SVE-10	NM	54.5	55.5	258.1	258.1	14.4	70.8	794.6	1173	764	1204	828.6	NM	232	8.5	NM	54.5
SVE-11	NM	NM	6.7	0.7	0.7	6.2	7.9	30.8	116.2	90	112	166	NM	19.2	3.5	NM	4.6
SVE-12	NM	30.7	10.7	65.2	65.2	14.8	3.1	55.9	357.9	613	1005	824.4	NM	42.9	12.3	NM	47.1
SVE-13	NM	11.7	4.1	27.6	27.6	7.4	3.3	18	20	183	417	409.6	NM	3.4	1.8	NM	4.2
								Eastern SVE W	/ells								
East Manifold	540	106.5	NM	NM	NM	NM	5	NM	NM	NM	340	54.2	NM	204.4	4.9	NM	20.9
SVE-02	NM	NM	0	0.4	0	0.1	0.2	4.3	32	NM	16.5	9.7	NM	4.4	0	NM	0.2
SVE-04	NM	NM	16.6	8.3	0	0.5	0.7	94.9	137.8	NM	313	8.2	NM	54.3	0	NM	28.3
SVE-06	NM	NM	52.5	24.5	20.8	13.9	15	326.9	387.6	NM	109	8.3	NM	401.7	1	NM	11
SVE-08	NM	NM	1.9	3.5	1.2	6.3	0.4	38.4	100.6	NM	64	11.1	NM	57.7	3.2	NM	2.4
SVE-07	NM	NM	4.7	4.2	24.8	7.5	0.8	61.9	299.6	NM	98	13.5	NM	66.8	4.5	NM	7.8
SVE-05	NM	1,105	828.9	483.7	1,811	43	105	>3980	>3980	NM	NM	2,093	NM	3,222	144	NM	490
SVE-03	NM	446.8	246.7	85.9	622.8	21.8	65.8	>3980	>3980	NM	NM	2543	NM	1494	21.5	NM	147.2
SVE-01	NM	11.0	6.3	2.7	2.1	5.8	4.9	98.6	238	NM	12.7	58.4	NM	3.6	0	NM	0
Influent	358	85.6	72.3	16.4	45.2	6.8	NM	258.1	315.7	190	336	58.4	1188	165	16.1	15.4	21.6
Effluent	0	1.2	0.2	0.6	0.7	0.2	NM	17.3	28.8	5.9	17.7	11.1	9	1.1	0.1	0.1	0

#### <u>Notes:</u>

1. FID was calibrated to 100 ppmv isobutulene prior to 3/21/2018. After 3/21/2018, FID was calibrated to 100-ppmv hexane

Abbreviations:

FID = flame ionization detector NM = not measured ppmv = parts per million by volume

SVE = soil vapor extraction

#### TABLE 6b : FLAME IONIZATION DETECTOR READINGS, SVE OPERATIONS WITH AIR SPARGING

Former Kelly-Moore Manufacturing Facility, Seattle, Washington

Well		FID	Readings <sup>1,2</sup> (pp	mv)	
	5/29/2019	5/29/2019	5/30/2019	6/4/2019	7/2/2019
		Western SVE	Wells		
West Manifold	NM	NM	173	43.2	38.4
SVE-09	NM	79.2	18.1	14	15.9
SVE-10	NM	1464	216	120	69.5
SVE-11	NM	1421	99	32.6	23.8
SVE-12	NM	2101	371	82.5	52.9
SVE-13	NM	2696	235	52.1	58
		Eastern SVE	Wells		
East Manifold	NM	NM	93.3	88.9	154
SVE-02	NM	11.2	0	0	0
SVE-04	NM	9.3	18.6	14.3	30.3
SVE-06	NM	7.3	0	0	0
SVE-08	NM	6.1	1	2	0.3
SVE-07	NM	6.1	2.5	3.4	5.7
SVE-05	NM	92	393	446	457
SVE-03	NM	40.3	163.1	121.8	213
SVE-01	NM	6.3	0.6	0	4.5
Influent	2322	1058	144.3	50.4	53
Effluent	79.5	20.1	3	1.2	0

Notes:

1. FID was calibrated to 100 ppmv hexane.

2. Air sparging in the western SVE area began on May 29, 2019

Abbreviations:

FID = flame ionization detector

NM = not measured

ppmv = parts per million by volume

SVE = soil vapor extraction

#### TABLE 7: SOIL VAPOR EXTRACTION SYSTEM DESTRUCTION REMOVAL EFFICIENCIES

Former Kelly-Moore Manufacturing Facility, Seattle, Washington

	FID Reading at	FID Reading at	Destruction Removal
	Influent <sup>1</sup>	Effluent <sup>1</sup>	Efficiency <sup>2</sup>
Date	(ppm)	(ppm)	(percent)
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%
6/29/18	258.1	17.3	93.3%
7/31/18	315.7	28.8	90.9%
8/31/18	190.0	5.9	96.9%
9/19/18	336.0	17.7	94.7%
10/16/18	58.4	11.1	81.0%
4/1/19	1188.0	9.0	99.2%
4/3/19	165.0	1.1	99.3%
5/3/19	16.1	0.1	99.4%
5/6/19	15.4	0.1	99.4%
5/29/19	21.6	0.0	100.0%
5/29/19	2322.0	79.5	96.6%
5/29/19	1058.0	20.1	98.1%
5/30/19	144.3	3.0	97.9%
6/4/19	50.4	1.2	97.6%
7/2/19	53.0	0.0	100.0%

#### Notes:

- 1. After 3/21/2018, all FID measurements are calibrated to 100 ppm hexane in accordance with Puget Sound Clean Air Agency requirements.
- Destruction removal efficiency generally exceeds the Puget Sound Clean Air Agency standard of 95% removal. In mid-2018, destruction removal efficiency was reduced. Troubleshooting and investigation suggested that FID calibration drift and sampling error were the likely cause.

#### Abbreviations:

- FID = flame ionization detector
- ppm = parts per million

#### TABLE 8: SOIL VAPOR EXTRACTION SYSTEM PERFORMANCE SUMMARY

Former Kelly-Moore Manufacturing Facility, Seattle, Washington

		Destruction				
	Soil Vapor	Removal	Operating	Extraction Flow	Mass Removal Rate	Total Mass Removed
	Concentration'	Efficiency	Factor <sup>2</sup>	Rate	(Hexane equivalent)	(Hexane equivalent)
Date	(ppm as Hexane)	(%)	(% Runtime)	(SCFM) <sup>3</sup>	(lb/hr)	(lbs)
11/7/17	358.0	100%	100%	215	0.35	0
11/15/17	547.5	99%	80%	215	0.43	64
11/21/17	123.7	98%	100%	215	0.12	100
12/5/17	59.6	100%	99%	135	0.04	123
12/13/17	85.6	99%	84%	249	0.08	188
1/16/18	72.3	100%	45%	127	0.02	211
2/13/18	16.4	96%	100%	240	0.02	260
3/21/18	45.0	98%	99%	353	0.07	509
4/24/18	6.8	97%	88%	142	0.00	526
6/29/18	258.1	93%	71%	225	0.19	1441
7/31/18	315.7	91%	36%	251	0.13	2172
8/31/18	190.0	97%	100%	242	0.21	3277
9/19/18	336.0	95%	100%	83	0.13	3948
10/16/18	58.4	81%	100%	238	0.06	3949
4/1/19	1188.0	99%	100%	237	1.28	4038
4/3/19	165.0	99%	99%	236	0.18	4177
5/3/19	16.1	99%	100%	191	0.01	4189
5/6/19	15.4	99%	98%	187	0.01	4207
5/29/19	21.6	100%	93%	184	0.02	4230
5/29/19	2322.0	97%	100%	184	1.95	6910
5/29/19	1058.0	98%	100%	184	0.89	8132
5/30/19	144.3	98%	95%	161	0.10	8272
6/4/19	50.4	98%	100%	160	0.04	8328
7/2/19	53.0	100%	99%	154	0.04	8408

#### Notes:

- 1. After 3/21/2018, all FID measurements are calibrated to 100ppm Hexane in accordance with Puget Sound Clean Air Agency requirements.
- Operating factor is calculated as the percentage of operating time standard of 95% removal. In mid-2018, removal efficiency was reduced. Troubleshooting and investigation suggested that FID calibration drift and sampling error were likely the cause of this issue.
- 3. Reported flow rates on 11/7/2017, 11/15/2017, 2/13/2018, and 4/1/2019 are estimates

#### Abbreviations:

ppm = parts per million, volume

FID = flame ionization detector

NM = not measured

SCFM = standard cubic feet per minute

#### TABLE 9: AIR QUALITY MONITORING RESULTS <sup>1, 2</sup>

Former Kelly-Moore Manufacturing Facility, Seattle, Washington

				Results report	ed in microg	rams	per cubic mete	er (µg	g/m³)						
	MTCA Meth	od B Indoor	MTCA Meth	od C Indoor	KM-AS-	1-	KM-AS-1	-	KM-AS-2-		KM-AS-2	-	KM-AS-3-		KM-AS-3-
	Air Clean	up Level <sup>3</sup>	Air Clean	up Level <sup>4</sup>	10161	8	060519		101618		060519		101618		060519
Compound	Non cancer	Cancer	Non cancer	Cancer	(Above SV	E-05)	(Above SVE	-05)	(Above SVE-0	3)	(Above SVE	-03)	(Ambient)	)	(Ambient)
1,1,1-Trichloroethane	2,286	-	5,000	-	0.55	U	0.55	U	0.55	U	0.55	U	0.55	U	0.55 U
1,1,2,2-Tetrachloroethane	-	0.043	-	0.431	0.14	U	0.14	U	0.14	U	0.14	U	0.14	U	0.14 U
1,1,2-Trichloroethane	0.091	0.156	0.200	1.563	0.055	U	0.11	U	0.055	U	0.11	U	0.055	U	0.11 U
1,1-Dichloroethane	-	1.56	-	15.63	0.4	U	0.4	U	0.4	U	0.4	U	0.4	U	0.4 U
1,1-Dichloroethene	91.4	-	200.0	-	0.4	U	0.4	U	0.4	U	0.4	U	0.4	U	0.4 U
1,2,3-Trimethylbenzene	-	-	-	-	2.5	U	NA		2.5	U	NA		2.5	U	NA
1,2,4-Trichlorobenzene	0.914	-	2.000	-	0.74	U	0.74	U	0.74	U	0.74	U	0.74	U	0.74 U
1,2,4-Trimethylbenzene	3.20	-	7.00	-	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5 U
1,2-Dibromoethane	4.11	0.004	9.00	0.042	0.077	U	0.077	U	0.077	U	0.077	U	0.077	U	0.077 U
1,2-Dichlorobenzene	91.4	-	200.0	-	0.6	U	0.6	U	0.6	U	0.6	U	0.6	U	0.6 U
1,2-Dichloroethane	3.20	0.096	7.00	0.962	0.19		0.11		0.17		0.11		0.11		0.077
1,2-Dichloropropane	1.83	0.250	4.00	2.500	0.23	U	0.23	U	0.23	U	0.23	U	0.23	U	0.23 U
1,3,5-Trimethylbenzene	-	-	-	-	2.5	U	2.5	U	2.5	U	2.5	U	2.5	U	2.5 U
1,3-Butadiene	0.914	0.083	2.000	0.833	0.48		0.022	U	0.44		0.022		0.19		0.022 U
1,3-Dichlorobenzene	-	-	-	-	0.6	U	0.6	U	0.6	U	0.6	U	0.6	U	0.6 U
1,4-Dichlorobenzene	366	0.227	800	2.273	0.24	U	0.24	U	0.24	U	0.24	U	0.24	U	0.24 U
1,4-Dioxane	13.7	0.500	30.0	5.000	0.36	U	0.36	U	0.36	U	0.36	U	0.36	U	0.36 U
1-Butanol	-	-	-	-	6.1	U	NA		6.1	U	NA		6.1	U	NA
2-Butanone	2,286	-	5,000	-	15		17		15	_	17		2.9	U	2.9 U
2-Hexanone	-	-	-	-	4.1	U	4.1	U	4.1	U	4.1	U	4.1	U	4.1 U
2-Pentanone	-	-	-	-	3.5	U	NA		3.5	U	NA		3.5	U	NA
3-Hexanone	-	-	-	-	4.1	U	NA		4.1	U	NA		4.1	U	NA
3-Pentanone	-	-	-	-	3.5	U	NA		3.5	U	NA		3.5	U	NA
Acetaldehyde	4.11	1.14	9.00	11.36	87		NA		32	_	NA		9	U	NA
Acetone	14,171	-	31,000	-	47		18		41		16		17		7.5
Acetonitrile	27.4	-	60.0	-	1./	U	NA		1./	U	NA		1./	U	NA
Acrolein	-	-	-	-	2.6		1.1		1.6		0.92	U	0.92	0	0.92 U
Acrylonitrile	0.914	0.037	2.000	0.368	0.22	U	NA		0.22	U	NA		0.22	U	NA 0.22
Benzene	13.7	0.321	30.0	3.205	1.9		0.32	0	1.7		0.32	0	0.98		0.32 U
Benzyl chioride	0.457	0.051	1.000	0.510	0.052	0	0.052	0	0.052	0	0.052	0	0.052	0	0.052 0
Bromotorm	-	2.21	-	22.13	2.1	0	2.1	0	2.1	0	2.1	0	2.1	0	2.1 U
Bromomethane	2.29	-	5.00	-	1.0	0	1.0	U	1.0	0	1.0	U	1.0	0	1.6 U
Carbon disulfido	220	-	700	-	2.9		6 2		2.9		6 2		2.9	0	6.2 LL
Carbon totrachlorido	320	- 0.417	100.0	- 4 167	0.2		0.2	0	0.2	0	0.2		0.2	0	0.2 0
	43.7	0.417	700	4.107	1.0	0	0.03		1.0	0	0.03		1.7	0	2.03 0
	12 714	-	20,000	-	0.82		0.77	0	0.77		0.77		0.77		0.77 11
CFC-12	<u>45</u> 7	-	100.0	-	2 1		0.77	0	2.1	0	0.77	0	2 1	0	23
Chlorobenzene	22.9	-	50.0		0.46		0.46		0.46		0.46		0.46	11	0.46 11
Chlorodifluoromethane	22.5	-	50.000	_	1 9	0	0.40	0	1.8	0	0.40	0	1.5	0	0.40 U
Chloroethane	4 571	_	10,000	_	0.26	U	2.6	ш	0.26	ш	2.6	П	0.26	U.	26 11
Chloroform	44.8	0 109	98.0	1 087	0.35	0	0.24	Ū	0.29		0.24	0	0.14	0	0.083
Chloromethane	41.1	-	90.0	-	1.5		21	U	14		21	U	14	_	21 11
cis-1.2-Dichloroethene	-	-	-	-	0.4	U	0.4	Ŭ	0.4	U	0.4	U	0.4	U	0.4 11
cis-1.3-Dichloropropene	-	-	-	-	0.45	U	0.45	U	0.45	Ū	0.51	-	0.45	U	0.45 U
Cyclohexane	2,743	-	6,000	-	6.9	U	6.9	U	6.9	U	6.9	U	6.9	U	6.9 U
Cyclopentane	-	-	-	-	0.41		NA	-	0.29	U	NA		0.29	U	NA
Dibromochloromethane	-	0.093	-	0.926	0.085	U	0.085	U	0.085	U	0.085	U	0.085	U	0.085 U
Dichlorobromomethane	-	0.068	-	0.676	0.067	U	0.067	U	0.067	U	0.067	U	0.067	U	0.067 U
Ethanol	-	-	-	-	5,900	J	1,500	J	5,300	J	1,400	J	29	J	150 J
Ethylbenzene	457	-	1,000	-	1.3		0.43	U	1.2		0.43	U	0.61		0.43 U
F-114	_	_	_	_	0.7	Ш	0.7	U.	0.7	Ш	07	Ш	0.7	Ш	07 11

Wood Environment & Infrastructure Solutions, Inc.

#### TABLE 9: AIR QUALITY MONITORING RESULTS <sup>1, 2</sup>

Former Kelly-Moore Manufacturing Facility, Seattle, Washington

	Results reported in micrograms per cubic meter ( $\mu g/m^3$ )														
	MTCA Method B Indoor MT		MTCA Meth	MTCA Method C Indoor		KM-AS-1-		KM-AS-1-		KM-AS-2-		KM-AS-2-		-	KM-AS-3-
	Air Cleanu	up Level <sup>3</sup>	Air Cleanu	up Level <sup>4</sup>	101618	3	060519		101618		060519	)	101618		060519
Compound	Non cancer	Cancer	Non cancer	Cancer	(Above SVI	E-05)	(Above SVE	-05)	(Above SVE-	03)	(Above SVI	-03)	(Ambient	)	(Ambient)
Hexachlorobutadiene	-	0.114	-	1.136	0.21	U	0.21	U	0.21	U	0.21	U	0.21	U	0.21 U
Hexanal	-	-	-	-	11		NA		7.9		NA		4.1	U	NA
Hexane	320	-	700	-	4.2		3.5	U	3.5	U	3.5	U	3.5	U	3.5 U
Iodomethane	-	-	-	-	0.58	U	NA		0.58	U	NA		0.58	U	NA
Isobutene	-	-	-	-	1.5		NA		1.4		NA		0.92	U	NA
Isoprene	-	-	-	-	0.51		NA		0.59		NA		0.28	U	NA
Isopropyl alcohol	-	-	-	-	13		8.6	U	11		8.6	U	8.6	U	8.6 U
m, p-Xylene	-	-	-	-	5		1.2		4.5		1.3		2		1.1
Methacrolein	-	-	-	-	2.9	U	NA		2.9	U	NA		2.9	U	NA
Methyl isobutyl ketone	1,371	-	3,000	-	4.1	U	4.1	U	4.1	U	4.1	U	4.1	U	4.1 U
Methyl t-butyl ether	1,371	9.6	3,000	96.2	1.8	U	1.8	U	1.8	U	1.8	U	1.8	U	1.8 U
Methyl vinyl ketone	-	-	-	-	2.9	U	NA		2.9	U	NA		2.9	U	NA
Methylene chloride	274	250	600	2,500	87	U	87	U	87	U	87	U	87	U	87 U
Naphthalene	1.37	0.074	3.00	0.735	0.59	J	0.14		0.55	J	0.14		0.27	J	0.079 U
o-Xylene	45.7	-	100.0	-	1.7		0.44		1.5		0.46		0.70		0.43 U
Pentanal	-	-	-	-	3.5	U	NA		3.5	U	NA		3.5	U	NA
Pentane	-	-	-	-	9		3	U	8.5		3	U	5.3		3 U
Propene	-	-	-	-	23		0.69	U	20		0.69	U	4.5		0.69 U
Styrene	457	-	1,000	-	1.3		0.85	U	1.1		0.86	U	0.85	U	0.85 U
Tetrachloroethene	18.3	9.62	40.0	96.15	0.68	U	6.8	U	0.68	U	6.8	U	0.68	U	6.8 U
Toluene	2,286	-	5,000	-	7.5		2.3		7.0		2.1		3.4		1.5
trans-1,2-Dichloroethene	27.4	-	-	-	0.4	U	0.4	U	0.4	U	0.4	U	0.4	U	0.4 U
trans-1,3-Dichloropropene	-	-	-	-	0.45	U	0.45	U	0.45	U	0.45	U	0.45	U	0.45 U
Trichloroethene	0.914	0.370	2.000	6.300	0.27	U	0.27	U	0.27	U	0.27	U	0.27	U	0.27 U
Vinyl acetate	91.4	-	200.0	-	7	U	7	U	7	U	7	U	7	U	7 U
Vinyl chloride	45.7	-	100.0	-	0.26	U	0.26	U	0.26	U	0.26	U	0.26	U	0.26 U

#### <u>Notes</u>

1. Data qualifiers are as follows

J = The result is estimated.

U = Analyte not detected above the indicated laboratory reporting limit.

2. Values in **bold** were above the outdoor ambient concentration.

 MTCA Method B Indoor Air Cleanup Levels, Draft Guidance for Evaluating Soil Vapor Intrusion in Washington State: Investigation and Remedial Action, October 2009 and updated in April 2018.

4. MTCA Method C Indoor Air Cleanup Levels, Draft Guidance for Evaluating Soil Vapor Intrusion in Washington State: Investigation and Remedial Action, October 2009 and updated in April 2018.

#### Abbreviations

- = there is no clean up level for this compound
 MTCA = Model Toxics Control Act
 NA = not analyzed



## Appendix A

WATER LEVEL MONITORING RECORD											
Project Na	me: <u>Ke</u>	ly- Moore		Projec	t and Task N	lumber: 14697					
Date: 8	16/18	Measured I	by: Luk	Le. Kerne	Instrumen	t Used: Solingt Made	1 101				
Note: For	you conve	nience, the fo	llowing abbre	eviations ma	v be used.						
P = Pum	nping	I = Inacces	ssible	D = Dedica	ated Pump						
ST = Stee	el Tape	ES = Electric	Sounder	MP = Measu	uring Point	WL = Water Level					
Well No.	Time	MP Elevation (feet)	Water Level Below MP (feet)	Water Level Elevation (feet)	Previous Water Level Below MP	Remarks	1				
KAU-OZR	1358	21.63	9.%	1.67							
KAU-03R	1353	21.54	9.93	11.61							
KMU- 04	1350	18:5 6-	7.06	11.50							
KMV- 06	1400	19.80	8.29	11.51							
KMU-07	1356	21.63	9.96	11.67		ыт. 					
KMV-08	1354	21.65	10.00	11.65							
1(mu-09	1402	18.14	6.64	11.50							
KMU-10	1404	20.39	9.01	11.38							
-											

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# MONITORING WELL/PIEZOMETER NUMBER KMW-028.

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Page 1 of

Chemical Analyses: Jee coc

Analytical Lab: Friedman Bruga

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					25.00			ame	c A
			GROU	JNDWA Low F	TER SA Iow Sa	MPLIN mpling	G LOG	whee	er eler
		MONIT	ORING WE	ELL/PIE	ZOMET	ER NU	MBER_K	MW-03R	
Project N	ame:	Kelly	Moore		1		_	5x1. 110	
Project N	umber:	14697	ē		Wea	ather Co	Date:	Overcast	
Location	Jeat	Fle WA							
Sampler:	L	Kerne			Win	d Speed	/Direction: _	MB	
				WELL	INFORM	NOITAN	1 .		
Casing D	iameter (ii	n):	2		c	Groundw	ater Elevatio	on (ft):	
Top of Ca	asing Elev	ation (ft):			0	Depth of	Well Casing	(ft):	
Initial Dep	oth to Wat	er (ft):	9.93		F	Actual Pu	irge Volume	(gal): 2.5	
weinead	Condition	1:							
;			PU	RGING	MEASU	JREME	NTS	- 2 	-
WI (ft		pH (std	SC	Tomp	OPP	DO	Turbidity		N.,
btoc)	Time	units)	(ms/cm)	(°C)	(mv)	(mg/L)	(NTUs)	Notes	
993	1210	6.89	0.412	18:3	235.5	100	7.8		1
992	1215	6.96	0.398	18.2	203	O.U.	10.7		1
9.95	1220	6.99	0.797	184	164	0.57	16.5		
9.92	1775	7.00	0.301	18.3	158	0.55	19:2	B. Shlisin me	Her Turn upside da
9.941	1230	7.02	0.382	18.2	143	0.51	ID		
9.95'	1235	7.02	0.384	18:3	133	0.50	10		2
995'	1730	7.03	0.378	18.3	122	0.49	11		
9.95'	1741	7.04	0.376	182	114	0.47	Ŋ		
9.951	1244	703	0.370	18.3	112	6.47	10		
								2	-
									_
			in the Difference						· · ·
			440						_
Sample II	D No.: K	MW- 03	R-081618	<u>A. 0.1.10</u>					
ORP/DO I	Meter Mod	lel & No.:	VET PO	MORCH IN	1				
Purge Eq	uipment U	lsed:	Geo Pum	ptf-1					
Sampling	Equipme	nt Used:	YSI Pro	PSS					
Purge Sta	art Time:		176	15		Sample (	Collection Ti	ime: 1250	
Purge Co	mpletion	Time:	174	5		Purging	Method:	SAA	101
Average I Analytica	Lab:	(mL/min)	Bang		_	Sample ( Chemica	I Analyses:	See CEC	real

0

Other Field Obse	rvations:	No	Smell,	Nev.	0	÷.	
+ Take	Dupicate	Sample	KMI	1.03R-0	081618	@ 1255	
0,		- 1 -					



Project Na	ame: _	MONIT Selly-2	ORING WE	LL/PIE	ZOMET	ER NUI	MBERK	MW-04 116/10	
Project N	umber:	4697	/		We	ather Co	nditions:	Overest	
Location:	Seattl	e, WA							
Sampler:	L. K	erner			Wir	nd Speed	/Direction:	6	
				WELL	INFORI	MATION	1		
Casing Di Top of Ca Initial Dep	ameter (i sing Elev oth to Wat	n): ration (ft): ter (ft):	7.06		C I J	Groundw Depth of V Actual Pu	ater Elevati Well Casing Irge Volume	on (ft): g (ft): ∋ (gal):Z <u>Sa(</u>	š
Wellhead	Conditio	n:	Good. Could	ld use r	rew Bul	172	-		
			PU	RGING	MEASU	JREME	NTS		
WL (ft btoc)	Time	pH (std. units)	SC (ms/cm)	Temp. (°C)	ORP (mv)	DO (mg/L)	Turbidity (NTUs)	Notes	
7.00	1350	6.23	0.370	16.7	130	1.17	7.0	J. Hinbegionin 3	Some
7.06	1355	6.16	0.353	17.0	123	0.70	9.0		
707	1400	6.11	0.326	17.2	121	0.68	9.5		
7.07	MOS	6.11	0.376	172	110	0.66	8.0		2
7.07	1408	6.09	0.325	17.5	104	0.45	7.8	Sun correspont.	
7.08	1411	6.09	0.326	17.8	61	0.6.4	7.9		
7.09	1414	6.89	0.326	18.0	99	0.65	7.9	· · · · · · · · · · · · · · · · · · ·	
		-							12
				+					
					1				
Sample ID Water Lev ORP/DO M Purge Equ	No.: el Ind. Mo leter Moc lipment L	KMW~( odel & No. lel & No.: Jsed:	24-081618 : Jel: nst YSI Pro CTCOPLO	Madel p Pss	101 	dicatal	Tusing		
Sampling	Equipme	nt Used:	TSI Pro	1755	,		ð		
Purge Sta Purge Cor	rt Time: npletion	Time: e (ml /min	1245	-		Sample ( Purging   Sample (	Collection T Method:	ime: <u>1430</u> <u>Per:stalta</u>	<u>544</u>
Analytical	Lab: F	riedan	Brugg			Chemica	Analyses:	Spe COC	17014
Other Fiel	d Observ	ations: _/	No Smell,	No she	in	1.			are a



# MONITORING WELL/PIEZOMETER NUMBER \_ KMU-06

Project Name: Kolly - Moore		
Project Number: 14697	Date: <u>B//6/18</u> Weather Conditions: Outcode	
Location: Seattle, UA		

Sampler: L.Kern

Wind Speed/Direction: \_NONE

#### WELL INFORMATION

Casing Diameter (in):	2
Top of Casing Elevation (ft):	
Initial Depth to Water (ft):	8.291
Wellhead Condition: Char	b

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

WL (ft btoc)	Time	pH (std. units)	SC (ms/cm)	Temp. (°C)	ORP (mv)	DO (mg/L)	Turbidity (NTUs)	Notes
8.29	14 39	6.30	0.502	19.4	110	0.71	23.5	Clear
8.30	1444	6.31	0.495	19.8	66	O.SL	23.5	
8,30	1449	6.32	NUTI	19.8	48	0.46	11	
8.30	1454	6.31	0.456	19.9	30	0.43	21.4	
8.30	Allood 14	\$ 6.33	0.446	19.9	12011	0.40	17.7	1
8.31'	1504	6.33	0.438	19.9	-6	0.39	17.4	
6.31'	1509	6.33	0.432	19.9	-21	0.38	19.10	
8.70'	1512	6.33	0,430	19.9	- 27	0.37	20.6	
8.30'	1515	6.33	0.427	19.9	- 32	037	18.2	
8.30'	1518	6.33	0.424	19.9	-36	0.37	18.0	
8.301	1520	6.33	0.421	19.5	- 39	0.37	18.4	
						¥1	0	
-								

Sample ID No.: <u>KMU-06-</u> Water Level Ind. Model & No.:	Colinst Model 101			
ORP/DO Meter Model & No.:	YSI Pro DSS			
Purge Equipment Used:	Geo Puno 1 uli	Dedreaded Tabin		
Sampling Equipment Used:	451 Pro DSS	0	1	
Purge Start Time:	1434	Sample Collection Time:	1520	
Purge Completion Time:	1520	Purging Method:	SAA	1
Average Purge Rate (mL/min):	200	Sample Containers Used:	LABPROVIDED 9TOT	
Analytical Lab: Fr: 1105	Brunk	Chemical Analyses: See	· (a)	
Other Field Observations: 91	Pi SC would not	stabilizeaber USminules.	5. (4) (5) (6) (6) (	



1. .

	Date: 9/16/18
Project Number: 14697	Weather Conditions:
Sampler: 2 Keyner	Wind Speed/Direction:
WELL	INFORMATION
Casing Diameter (in):	Groundwater Elevation (ft):
Top of Casing Elevation (ft):	Depth of Well Casing (ft):
	Actual Purge volume (gal):
Wellhead Condition:	
Wellhead Condition:	MEASUREMENTS

WL (ft btoc)	Time	(std. vunits)	SC (ms/cm)	Temp. (°C)	ORP (mv)	DO (mg/L)	Turbidity (NTUs)	Notes
9:96	940	6.08	0.210	15	312.5	1.06	22.4	
9.90	945	602	6.210	14.9	311.5	0.79	27.0	
9.98	9:50	600	0.214	15	308.5	0.72	27.8	
9.99	9:53	6.05	0.213	15	305	0.67	30	
9.98	9:56	6.05	0.213	15	224	0405	29	200 U.S. 10 U.S. 20
9.98	9:50	6.24	0.213	15.2	145	0.64	78	
9.98	10:02	603	0.212	15.0	7.56	6.61	30	
9.98	10:05	6.03	0.212	15	260	0.60	31	
9.98	10:08	602	0.212	15.2	714.5	0.61	30	
9.93	10:11	6.02	0.217	15.3	Udes	0.61	19	
e	10:14	6.02	0-211	15.4	167	0.6	30	
	10:17	602	0.211	15.4	268	10.60	31	
-	(0:20							1
/								36

Water Level Ind. Model & No.:	Solinst: Model 10	
ORP/DO Meter Model & No.:	VSI Pro DSA	
Purge Equipment Used:	Stoland up N	dirated Tubion.
Sampling Equipment Used:	KUSI PRO OSS	
Purge Start Time:	9:35	Sample Collection Time: GTB 10:70
Purge Completion Time:	10:12	Purging Method: 544
Average Purge Rate (mL/min)	100	Sample Containers Used: 17 9. Https
Analytical Lab: _ Con Local men	C. Baus	Chemical Analyses:Cec_OC

Project N	ame: _	Kelly-	Moore				Data: Q	111,118	
Proiect N	umber:	141.97		25	We	ather Co	Date: 0	Avecst	1
ocation	Seat	HENA			***C			JUNERS	
Sampler:	L. Ker	ne			Wir	d Speed	/Direction:	NA	
1.1				й. 					5 L
		*		WELL	INFOR	MATION	1	192	
Casing D	iameter (i	n):	2.		- (	Groundw	ater Elevati	on (ft).	
rop of Ca	sing Elev	ation (ft):	6			Depth of	Well Casing	(ft):	1
nitial Dep	oth to Wat	ter (ft):	io'			Actual Pu	rge Volume	(gal): 2591	
veiinead	Conditio	n:							
			PU	IRGING	MEAS	JREME	NTS		
	1 at 1	pН			4				
WL (ft	Time	(std.	SC (molom)	Temp.	ORP	DO	Turbidity		
DLOC)	111	( orb	(ms/cm)	(°C)	(mv)	(mg/L)	(NTUs)	Notes	1222
10.	IIIO	6.00	0.117	Dif	4.5	0.91	1200	Drainvater, Oran	e FLOC
10:01	110	5.94	0.1991	15.8	150	7.6	75		1.
0.061	1120	5.92	0,197	15.7	252	0.60	12.9	1997 - A	
9.91	1125	5.95	0,703	15.7	234	0.62	9,1		
9.91	1100	5.93	0.204	15.7	244	0.61	8-49.	3	
10,00	1131	5.94	0.208	5.7	248	0.59	9.2		
10,01	1134	2.94	0.701	15.7	149	0.50	9.5		
0.02	1157	2.95	0,211	15.7	748	0,50	10.1		
						145	N		
					-		<u></u>		
								_	
	-								
Sample ID	No · W	M11-00	ERILA A						
Vater Lev	el Ind. Mo	del & No.:	, Folinst	- Mule	1:101				
RP/DO N	leter Mod	lel & No.:	KSI Prot	Ross			12	2	
urge Equ	lipment U	sed:	Crep Ru	of fine	staltic	VI	<i>pedicated</i>	Jubing	
	- minna		1 1 00			C			

amec foster wheeler



#### MONITORING WELL/PIEZOMETER NUMBER KMU-01

Project Name:	Kelly-Moore
Project Number	14107

Location: ( Jung

Sampler:

Date: <u>8//w/10</u> Weather Conditions: <u>OVERCAS</u>

Wind Speed/Direction: NONE /NA

#### WELL INFORMATION

Casing Diameter (in): <u>7</u> Top of Casing Elevation (ft): <u>6.641</u> Initial Depth to Water (ft): <u>6.641</u> Wellhead Condition: <u>6.800</u>

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

WL (ft btoc)	Time	pH (std. units)	SC (ms/cm)	Temp. (°C)	ORP (mv)	DO (mg/L)	Turbidity (NTUs)	Notes
6.64	1540	6.32	0.447	18.1	99	0.81	35.5	
0.65	1545	6.34	0.441	18.2	61	0.58	8.0	<i>L</i>
6.65	1550	6:306	0.432	18.5	25	0.52	75	
0.65	1555	6.35	0.421	18.7	1)	0.50	7.3	
6.65	1600	6.36	0.407	18.8	-5	0.48	7.0	-
6.65	1105	6.30	0.395	19:3	-22	0.47	8.2	
0.65	1608	6.41	0.391	20.2	~28	0.47	75	
6.65	1611	6.39	0,387	20.6	-33	OUT	7.8	
10.66	1614	6.35	0.390	18.4	-27	0.47	11.0	
6.65	167	635	0.391	18.4	-23	0.47	12.0	
6.65	1620	6.35	0.387	18.4	-24	0.47	12.2	
8-65		- ve n a		33.5 8.		2.8	•	
5		2						it•
1. A.							1	5

Sample ID No .: KMU-09-1	08/018		
Water Level Ind. Model & No .:	Jolinst Model: 101		
ORP/DO Meter Model & No.:	YST Pro Das	1	
Purge Equipment Used:	Greepung HI w/ Dedicat	ed Tubin	
Sampling Equipment Used:	451 Pa 055	the start of the s	
Purge Start Time:	1535	Sample Collection Time:	1622
Purge Completion Time:	1620	Purging Method:	LAA
Average Purge Rate (mL/min):	2200	Sample Containers Used:	LAB Provides Of Tor
Analytical Lab: Friedman	Brugg	Chemical Analyses:	ca.
Other Field Observations:	Some sheen in	Purge Later	



# MONITORING WELL/PIEZOMETER NUMBER KMW-10

Project N	ame:	Kelly-1	loore					***	
							Date: 00	3/16/18	
Project N	umber:	14697			Weather Conditions: SunNF				
Location:	Location: Sest KC							01000	
Sampler:	A	Herne			Win	d Speed	Direction:	10000	
				WELL	INFORI		1		
Casing Di	iameter (iı	n):	2		c	Groundw	ater Elevati	on (ft):	
Top of Ca	sing Elev	ation (ft):	···		Γ	Depth of	Well Casing	g (ft):	
Initial Dep	oth to Wat	er (ft):	~ d		1	Actual Pu	irge Volume	e (gal): <u>2.5</u>	
weimeau	Continuition	(40)	001						
			PU	RGING	MEASU	JREME	NTS		
		pН		_			х		
WL (ft	Time	(std. units)	SC (ms/cm)	Temp.	ORP (my)	DO (mg/l)	(NTUs)	Notes	
5.00)	1650	C.74	1105	16:Z	lia	1.25	77	Notes	
	1155	6.01	6 407	17.7	125	7.92	14		
- 4	1700	6-110	19 UI4	16.10	7.5	63-3	77		
•	1705	6.27	DUIZ	16.0	48	0.53	22.8		
	1710	6.76	DUIZ	16.5	30	0.50	22.0		
	1715	6:25	0.418	16.5	10	0.47	223		
	1720110	6.25	0.415	16.5	25	0.47	21.9		
	1721	6.25	0.415	16.5	-5	0.47	71.8		
	1724	6.25	0.416	16.5	-8	0.47	21.7		
	1727	6.25	0.413	110:5	-12	0.46	22		
	1732	0.25	0.410	16.5	-15	0.46	21.8		
	A CONTRACTOR OF THE OWNER	All and the second s					2 2	19	

ample ID No.: [MU-10-08]	618		
Vater Level Ind. Model & No.:			1.00 M
DRP/DO Meter Model & No.:	ST Pro RSS	×	
Purge Equipment Used:	nes Pump		1
Sampling Equipment Used:	SI Pro RSS		
urge Start Time:	1645	Sample Collection Time:	17.30
urge Completion Time:	1730	Purging Method:	SAA
Average Purge Rate (mL/min):	200	Sample Containers Used:	LAR Por molel
Analytical Lab: Fridad Brit		Chemical Analyses: Ca	Car

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# WATER LEVEL MONITORING RECORD



Project Na	Project Name: Kelly-Moore Proje					ect and Task Number: <u>14697009</u>			
Date: 02/0	07/2019	_ Measured	by: <u>L. Kerne</u> r	•<	Instrument Used: Water Level Meter				
Note: For	Note: For you convenience, the following abbreviations may be used.								
P = Pun	nping	I = Inacce	ssible	D = Dedic	ated Pump				
ST = Stee	el Tape	ES = Electric	Sounder	MP = Measu	uring Point	WL = Water Level			
Well No.	Time	MP Elevation (feet)	Water Level Below MP (feet)	Water Level Elevation (feet)	Previous Water Level Below MP	Remarks			
KMW-02R	928	21.63'	9.17'	12.46	11.67				
KMW-03R	935	24.54'	9.37	12.17	11.61				
KMW-04	1003	18.56'	6.60	11.94	11.50				
KMW-06	938	19.80'	7.77	1203	11.51				
KMW-07	931	21.63'	9.21	12.42	11.67 .				
KMW-08	935	21.65'	9.31'	12.34	11.65				
KMW-09	940	18.14'	6.15	11.99	11.50	8			
KMW-10	943	20.39'	8.65	11.74	11.38				
		10/74			_				
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# MONITORING WELL/PIEZOMETER NUMBER- KMW-02R

**Project Name:** 

Kelly-Moore

Date: 2/8/19 Weather Conditions: INDEORS

Project Number: 14697009 Location: Seattle, WA Sampler: Lucas Kerner

Wind Speed/Direction: \_\_\_\_\_\_

#### WELL INFORMATION

Casing Diameter (in):	2"
Top of Casing Elevation (ft	): 21.63'
Initial Depth to Water (ft):	9.17
Wellhead Condition:	OK

Groundwater Elevation (ft): 12.46 Depth of Well Casing (ft): Actual Purge Volume (gal): 2 Saglic->

#### PURGING MEASUREMENTS

WL (ft btoc)	Time	pH (std. units)	SC (ms/cm)	Temp. (°C)	ORP (mv)	DO (mg/L)	Turbidity (NTUs)	Notes
9.14	923	5.99	0.257	13.9	106	1.60	104	Hoteo
9.21	926	5.96	0.750	14.0	1187	1.7.1	1070	
9.20	9219	5.96	0:240	1402	1749	111	1112	
9:21	932	5.95	0.245	14.0	17911	104	1111	
9:27	935	5.95	0.745	14.1	1785	0.90	7 24	
9.23	938	5.97	0.140	14.0	1774	0.82	6.91	
922	941	5.96	0:241	14.1	1291	0.80	781	
9:29	944	5.95	10.295	191	130.1	0.81	81	
9:25	947 24	,	123 ± 3 - 322	A	1 10/		00	
		_						
								1
	_				table -	811		
				_				N

Sample ID No .: KMW-02R- 07.0819 Water Level Ind. Model & No.: Solinst Model 101 ORP/DO Meter Model & No.: YSI-Pro Dss Purge Equipment Used: Peristaltic Pump with dedicated tubing Sampling Equipment Used: YSI Pro Dss 970 **Purge Start Time:** 945 Sample Collection Time: Purge Completion Time: 949 Purging Method: SAA Average Purge Rate (mL/min): Sample Containers Used: Lab Provided 220 numin Analytical Lab: Friedman & Bruya Inc. Chemical Analyses: See COC Other Field Observations: \_

Updated 1/31/19

Page <u>1</u> of <u>1</u>



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

**Project Name:** 

Kelly-Moore

	Date: 2/8/19
Project Number: 14697009	Weather Conditions: INDOORS MA
Location: Seattle, WA	
Sampler: Lucas Kerner	Wind Speed/Direction:

#### WELL INFORMATION

Casing Diameter (in):	2"
Top of Casing Elevation (ft	: 21.54'
Initial Depth to Water (ft):	9.32
Wellhead Condition:	MKAY

Groundwater Elevation (ft): 12-17 Depth of Well Casing (ft): Actual Purge Volume (gal): 7.991

WL (ft btoc)	Time	pH (std. units)	SC (ms/cm)	Temp. (°C)	ORP (mv)	DO (mg/L)	Turbidity (NTUs)	Notes
9.50	113	6.82	0.570	164	-49	1.43	9.80	
9.55	1116	6.91	0.617	16.5	-66	0.82	6.70	
9.56	1119	6.95	0.675	16.6	-72	0.71	7.71	
9.58	1122	8.99	0.635	16.6	- 80	0.01	11.85	
9.55	1125	7,01	0.591	16.0	-83	0.56	is.c	
9,57	1/18	703	0.581	14.6	-85	0.54	19.8	
9.57	113	6.95	0.586	16.5.	-85	0.52	20.7	
9.52	1134	6.99	0.583	16.5	86.4	0.51	70.4	10
9.57	1137	6.97	0.582	16.5	-87	0.51	70:52	
>	-							
				St	ship	/	100	
					DA		12	

Sample ID No.: KMW-03R- Ø	20819	
Water Level Ind. Model & No.:	Solinst Model 101	
ORP/DO Meter Model & No.:	YSI-Pro Dss	
Purge Equipment Used:	Peristaltic Pump with de	dicated tubing
Sampling Equipment Used:	YSI Pro Dss	
Purge Start Time: Purge Completion Time: Average Purge Rate (mL/min) Analytical Lab: Friedman & Br	1110 1137 137 175 <i>ml/nin</i>	Sample Collection Time: <u>1/40</u> Purging Method: <u>SAA</u> Sample Containers Used: <u>Lab Provided</u> Chemical Analyses: See COC
Other Field Observations:	5	



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

Project Name:

Kelly-Moore

Date: 02/07/19 Weather Conditions: 30 F SUNNY

Project Number: 14697009 Location: Seattle, WA Sampler: Lucas Kerner

Wind Speed/Direction: \_\_\_\_\_\_\_

#### WELL INFORMATION

Casing Diameter (in):	2"
Top of Casing Elevation (ft)	: 18.56'
Initial Depth to Water (ft):	6.60
Wellhead Condition:	OK

Groundwater Elevation (ft): 11.96 Depth of Well Casing (ft): Actual Purge Volume (gal): 1.5 stillers

#### PURGING MEASUREMENTS

WL (ft btoc)	Time	pH (std. units)	SC ⁄ (ms/cm)	Temp. (°C)	ORP (mv)	DO (mg/L)	Turbidity (NTUs)	Notes
6.65	1030	6.14	0.274	11.9	-10	3,15	20.50	
6.63	1032	6.23	0,785	17.1	-19	1.17	18.41	
1:64	1030	6:24	0,294	11.8	-60	0.96	18.43	
6.64	1039	6.24	0.305	11.9	-68	0,82	16.15	
6,103	1042	6.24	0.315	11.8	-74	0.75	15,91	
6.63	1045	6.21	0.31B	11.2	-72	0.74	15:30	
6.61	1046	6.22	0.334	11:2	-75	0.66.	14.64	
GUL	10.51	6:22	0.343	10.9	-76	0.641	12.98	
641	1054	6.22	0:3413	11,2	-76	0.61	12.95	
6.63	1057	6.22	0.341	11.7	-74	0.62	Rol	
	# X							· · · · · · · · · · · · · · · · · · ·
	tazille			SLU	Qu		-	6
	1106-20			Jang	Lak			
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#### Sample ID No .: KMW-04-070719

Water Level Ind. Model & No .:	Solinst Model 101						
ORP/DO Meter Model & No.:	YSI-Pro Dss	/SI-Pro Dss					
Purge Equipment Used:	Peristaltic Pump with dedic	Peristaltic Pump with dedicated tubing					
Sampling Equipment Used:	YSI Pro Dss	1 (A) A					
Purge Start Time:	10:24	Sample Collection Time:	1115				
Purge Completion Time:	10.57	Purging Method:	SAA				
Average Purge Rate (mL/min)	Hours from 150 ml/m	Sample Containers Used:	Lab Provided				
Analytical Lab: Friedman & Br	uya Inc.	Chemical Analyses: See (	200				
Other Field Observations:	Well Monument	Frozen ul vator					



### MONITORING WELL/PIEZOMETER NUMBER- KMW-06

Project Name:

Kelly-Moore

	Date: <u>PZ107/19</u>			
Project Number: 14697009	Weather Conditions: <u>472</u> SUNNY			
Location: Seattle, WA		2		
Sampler: Lucas Kerner	Wind Speed/Direction: N/A			

#### WELL'INFORMATION

Casing Diameter (in):	2"
Top of Casing Elevation (ft):	19.80'
Initial Depth to Water (ft):	7.77
Wellhead Condition:OI	it i

Groundwater Elevation (ft): <u>12,03</u> Depth of Well Casing (ft): Actual Purge Volume (gal): <u>デジム・</u>

WL (ft btoc)	Time	pH (std. units)	SC (ms/cm)	Temp. (°C)	ORP (mv)	DO (mg/L)	Turbidity (NTUs)	Notes	
7.80	1319	6.21	0, 4,23	141:3	-4.5	200	1358	12	ĺ
7.71	1322	6.20	0.640	14.2	-17.1	1.11	12:72		1
7.80	1325	6.19	0.639	14.2	-25	0.80	12.73		)
7.87	1328	6.18	0.641	14:2	-76.5	0.74	17.52	н	
7.81	1331	618	0.639	14:3	-79.2	0.67	12.45		
7.80	1334	6.18	0.635	14:2	-32	0:05	17.40		
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		$\sim$	sable	lin					
				BK					
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Sample ID No.: KMW-06-07	2711		
Water Level Ind. Model & No .:	Solinst Model 101		
ORP/DO Meter Model & No.:	YSI-Pro Dss		
Purge Equipment Used:	Peristaltic Pump with dedica	ted tubing	
Sampling Equipment Used:	YSI Pro Dss		
Purge Start Time: Purge Completion Time: Average Purge Rate (mL/min): Analytical Lab: Friedman & Br	$\frac{ 3 6}{ 334 }$	Sample Collection Time: Purging Method: Sample Containers Used:	SAA Lab Provided
Other Field Observations: _/	ISMSP Collec	for	



### MONITORING WELL/PIEZOMETER NUMBER- KMW-07

Project Name:

Location: Seattle, WA

Sampler: Lucas Kerner

Project Number: 14697009

Kelly-Moore

Date: 2/8/19 Weather Conditions: TNDOORS - NI

Wind Speed/Direction: \_\_\_\_

#### WELL INFORMATION

Casing Diameter (in):	2"
Top of Casing Elevation (ft	): 21.63'
Initial Depth to Water (ft):	9.21
Wellhead Condition:	OKAY

Groundwater Elevation (ft): 12.42 Depth of Well Casing (ft): Actual Purge Volume (gal): \_\_\_\_\_ ciallars

WL (ft btoc)	Time	pH (std. units)	SC (ms/cm)	Temp. (°C)	ORP (mv)	DO (mg/L)	Turbidity (NTUs)	Notes
9.25	1158	6:31	0.330	14.4	3.2	1.33	75.12	
9.24	1701	6.792	0.377	14.4	38	0.78	14.51	
9.25	1284	6.26	0.325	14.4	43	0.61	15.80	
9.26	1207	6.23	0.321	14.4	50	0.58	17.81	
9.75	1210	6.22	0.320	19.4	52.2	0.55	18-71	2
9.20	1213	6.73	0.319	14.4	51.9	0.52	18.66	- N
2:28	1216	6.23	0.318	14.4	51.1	0.52	17.91	
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			JUASIC	all	7			
				/				
		-						

Sample ID No .: KMW-07- 07	0819			
Water Level Ind. Model & No.:	Solinst Model 101			
ORP/DO Meter Model & No.:	YSI-Pro Dss			
Purge Equipment Used:	Peristaltic Pump with ded	icated tubing		
Sampling Equipment Used:	YSI Pro Dss			
Purge Start Time:	1155	Sample Collection Time:	1220	
Purge Completion Time:	1210	Purging Method:	SAA	
Average Purge Rate (mL/min):	175 million	Sample Containers Used:	Lab Provided	
Analytical Lab: Friedman & Br	uya Inc.	Chemical Analyses: See	COC	100
Other Field Observations:		S		1227



### MONITORING WELL/PIEZOMETER NUMBER- KMW-08

**Project Name:** 

Kelly-Moore

Date:	210	119	
Weather Conditions	: <u>'</u> †	NDEDKS/NIA	

Project Number: 14697009 Location: Seattle, WA Sampler: Lucas Kerner

Wind Speed/Direction:

### WELL INFORMATION

Casing Diameter (in):	2"
Top of Casing Elevation (ft)	21.65'
Initial Depth to Water (ft):	9.31
Wellhead Condition: /9/	AY

Groundwater Elevation (ft): 12.34 Depth of Well Casing (ft): Actual Purge Volume (gal): SSIL

WL (ft btoc)	Time	pH (std. units)	SC (ms/cm)	Temp. (°C)	ORP (mv)	DO (mg/L)	Turbidity (NTUs)	Notes
9.38	1070	6.13	0.200	13.7	78	28	100	
9:3	1023	6.05	0.241	13.8	09	1.34	5.2	
9,39	1076	6.04	0:74	13.0	97	1.00	44	·
9.40	1029	6.04	0.240	13.8	110	0.78	350	
9.40	1032	6.05	0-249	13.9	96	0.79	31.7	_
9,40	1035	6.05	0.297	13.9	95:3	0.75	31.2	
9.40	1033	6.05	0.250	13.9	91.4	0:74	3.0.9	
~	111 2° 50							
			SLIL					
			2thily	0	16			
				-22				
14								
12 2 12								

Nater Level Ind. Model & No.: Solinst Model 101	
Purge Equipment Used: Peristaltic Pump with dedi Sampling Equipment Used: YSI Pro Dss	icated tubing
Purge Start Time: <u>1017</u> Purge Completion Time: <u>1038</u> Average Purge Rate (mL/min): <u>700 pt //min</u> Analytical Lab: <u>Friedman &amp; Bruya Inc.</u>	Sample Collection Time: Purging Method: Sample Containers Used: Lab Provided Chemical Analyses: See COC
Other Field Observations:	



#### MONITORING WELL/PIEZOMETER NUMBER- KMW-09

**Project Name:** 

Kelly-Moore

	Date: <u>107/07/19</u>			
Project Number: <u>14697009</u>	Weather Conditions: 390 Clouds			
Location: Seattle, WA	đ			
Sampler: Lucas Kerner	Wind Speed/Direction: NW 4MPH			

#### WELL INFORMATION

Casing Diameter (in):	2"
Top of Casing Elevation (ft):	18.14'
Initial Depth to Water (ft):	615
Wellhead Condition:	GLOOD

Groundwater Elevation (ft): <u>11,99'</u> Depth of Well Casing (ft): Actual Purge Volume (gal): <u>1,59'</u>

#### PURGING MEASUREMENTS

WL (ft btoc)	Time	pH (std. units)	SC (ms/cm)	Temp. (°C)	ORP (mv)	DO (mg/L)	Turbidity (NTUs)	Notes
615	1433	6.58	0.404	19:2	-57	4.50	6.80	
6.17	M36	648	0,401	15.Z	-60	1.20	5:75	
6.19	1439	6.48	0.400	15:2	-65	0.79	5.40	
6.18	144	6.47	0.400	15.2	·ida	0.63	5.09	
6.19	1445	647	0.400	15.2	-1.0	063	5.54	t
6.19	1548	6.47	0:391	15:2	-la	OSB	4.95	
6.19	1551	6.47	0.400	15.2	-69	0.57	4.99	
~								
			SL	bi				
				gion	11	1		
		14			21		5	

Sampling Equipment Used: 1	SI Pro Lice		
Purge Start Time: Purge Completion Time: Average Purge Rate (mL/min): Analytical Lab: <u>Friedman &amp; Bru</u>	1430 1551 125 ya Inc.	Sample Collection Time: Purging Method: SA Sample Containers Used: La Chemical Analyses: See COC	515 AA ab Provided C
Other Field Observations:		а,	k.

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#### MONITORING WELL/PIEZOMETER NUMBER- KMW-10

Project Name: Kelly-Moore

	Date:				
Project Number: 14697009	Weather Conditions: 300 SUNNY				
Location: Seattle, WA					
Sampler: Lucas Kerner	Wind Speed/Direction: N/A				

#### WELL INFORMATION

Casing Diameter (in):	2"
Top of Casing Elevation (ft	): 20.39'
Initial Depth to Water (ft):	8.65
Wellhead Condition:	OKAY

Groundwater Elevation (ft): <u>11.74'</u> Depth of Well Casing (ft): Actual Purge Volume (gal): <u>-1.5 Sar Zgal</u>

2:0010

#### PURGING MEASUREMENTS

WL (ft btoc)	Time	pH (std. units)	SC (ms/cm)	Temp. (°C)	ORP (mv)	DO (mg/L)	Turbidity (NTUs)	Notes
	1203	6.54	0.423	13.1	-44	2.8	61	
8.65	1206	6.56	0.430	13.1	-59.41	1.13	472	
8.H	1209	6.55	0.432	13.4	-68	0.84	28	
8.68	17.12	6.54	0.430	13.6	-698	0.74	41	
8.68	1215	654	0.429	13.7	-71.0	0.70	37	
8.68	1218	6.54	0.433	13.8	-732	0.64	27.5	
0.60	1221	6.54	0.432	12.5	-74	0.65	29.9	
8.69	1229	6.54	0.429	14:2	-77	0.54	23.0	
8.71	1227	6.54	0.427	14.3	-79	0.51	16	
6:71	1230	6.54	0.429	14.0	-30.3	0.51	16.1	
8.71	1733	6.54	0.479	13.9	-81	0.50	15.6	
869	1136	6.53	0.1130	13.9	-82	0.49	15.6	
	173						5	p.
	1242							
	1245							

Sample ID No .: KMW-10- 02071	19
Water Level Ind. Model & No .: Sol	linst Model 101
ORP/DO Meter Model & No.: YSI	I-Pro Dss
Purge Equipment Used: Per	ristaltic Pump with dedicated tubing
Sampling Equipment Used: YSI	I Pro Dss
Purge Start Time:	2.00 Sample Collection Time: 1245
Purge Completion Time:	17.36 Purging Method: SAA
Average Purge Rate (mL/min):	Sample Containers Used: Lab Provided
Analytical Lab: Friedman & Bruya	Inc. Chemical Analyses: See COC
Other Field Observations:	up Talken = Kmw-10-9-0207@19 @1250

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# **Appendix B**

#### ENVIRONMENTAL CHEMISTS

James E. Bruya, Ph.D. Yelena Aravkina, M.S. Michael Erdahl, B.S. Arina Podnozova, B.S. Eric Young, B.S. 3012 16th Avenue West Seattle, WA 98119-2029 (206) 285-8282 fbi@isomedia.com www.friedmanandbruya.com

February 18, 2019

Crystal Thimsen, Project Manager Wood Environment & Infrastructure Solutions, Inc. One Union Square 600 University Street, Suite 600 Seattle, WA 98101

Dear Ms Thimsen:

Included are the results from the testing of material submitted on February 8, 2019 from the Kelly Moore 14697009, F&BI 902134 project. There are 45 pages included in this report. Any samples that may remain are currently scheduled for disposal in 30 days, or as directed by the Chain of Custody document. If you would like us to return your samples or arrange for long term storage at our offices, please contact us as soon as possible.

We appreciate this opportunity to be of service to you and hope you will call if you should have any questions.

Sincerely,

FRIEDMAN & BRUYA, INC.

Michael Erdahl Project Manager

Enclosures WEI0218R.DOC

#### ENVIRONMENTAL CHEMISTS

#### CASE NARRATIVE

This case narrative encompasses samples received on February 8, 2019 by Friedman & Bruya, Inc. from the Wood Environment & Infrastructure Solutions Kelly Moore 14697009, F&BI 902134 project. Samples were logged in under the laboratory ID's listed below.

<u>Laboratory ID</u>	Wood Environment & Infrastructure Solutions
902134 -01	KMW-02R-020819
902134 -02	KMW-03R-020819
902134 -03	KMW-04-020719
902134 -04	KMW-06-020719
902134 -05	KMW-08-020819
902134 -06	KMW-09-020719
902134 -07	KMW-10-020719
902134 -08	KMW-10-9-020719
902134 -09	KMW-7-020819
902134 -10	Trip Blanks

A 6020B internal standard failed the acceptance criteria for samples KMW-10-020719 and KMW-10-9-020719. The samples were diluted and reanalyzed with acceptable results. Both data sets were reported.

The 8270D matrix spike and matrix spike duplicate failed the relative percent difference for several compounds. The analytes were not detected therefore the data were acceptable.

1,1-Dichloroethane in the 8260C laboratory control sample exceeded the acceptance criteria. The analyte was not detected in the sample, therefore the data were acceptable.

All other quality control requirements were acceptable.

#### ENVIRONMENTAL CHEMISTS

Date of Report: 02/18/19 Date Received: 02/08/19 Project: Kelly Moore 14697009, F&BI 902134 Date Extracted: 02/12/19 Date Analyzed: 02/12/19

#### RESULTS FROM THE ANALYSIS OF WATER SAMPLES FOR TOTAL PETROLEUM HYDROCARBONS AS GASOLINE USING METHOD NWTPH-Gx

Results Reported as ug/L (ppb)

		Surrogate
Sample ID Laboratory ID	Gasoline Range	( <u>% Recovery</u> ) (Limit 51-134)
KMW-02R-020819 902134-01	<100	92
KMW-03R-020819 902134-02	140	97
KMW-04-020719 902134-03 1/10	31,000	108
KMW-06-020719 902134-04	2,200	ip
KMW-08-020819 902134-05	120	96
KMW-09-020719 902134-06	450	107
KMW-10-020719 902134-07	200	109
KMW-10-9-020719 902134-08	210	109
KMW-7-020819 902134-09	<100	94
Method Blank <sup>09-321 MB</sup>	<100	103

#### ENVIRONMENTAL CHEMISTS

Date of Report: 02/18/19 Date Received: 02/08/19 Project: Kelly Moore 14697009, F&BI 902134 Date Extracted: 02/12/19 Date Analyzed: 02/12/19

#### RESULTS FROM THE ANALYSIS OF WATER SAMPLES FOR TOTAL PETROLEUM HYDROCARBONS AS DIESEL AND MOTOR OIL USING METHOD NWTPH-Dx

Results Reported as ug/L (ppb)

Sample ID Laboratory ID	Diesel Range (C10-C25)	Motor Oil Range (C25-C36)	Surrogate <u>(% Recovery)</u> (Limit 47-140)
KMW-02R-020819 902134-01 1/1.2	<60	<300	95
KMW-03R-020819 902134-02 1/1.2	1,700 x	<300	110
KMW-04-020719 902134-03 1/1.2	2,600 x	<300	64
KMW-06-020719 902134-04 1/1.2	19,000 x	790 x	63
KMW-08-020819 902134-05 1/1.2	440 x	<300	102
KMW-09-020719 902134-06 1/1.2	3,100 x	<300	64
KMW-10-020719 902134-07 1/1.3	970 x	<320	75
KMW-10-9-020719 902134-08 1/1.2	1,300 x	<300	105
KMW-7-020819 902134-09 1/1.2	<60	<300	106
Method Blank	<60	<300	100

### ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-02R-0 02/08/19 02/12/19 02/13/19 Water ug/L (ppb)	020819	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Infrastructure Solutions Kelly Moore 14697009 902134-01 1/2 021311.D GCMS6 VM
Surrogates: Anthracene-d10 Benzo(a)anthracene	-d12	% Recovery: 99 112	Lower Limit: 31 25	Upper Limit: 160 165
Compounds:		Concentration ug/L (ppb)		
Naphthalene		< 0.4		
Acenaphthylene		< 0.04		
Acenaphthene		< 0.04		
Fluorene		< 0.04		
Phenanthrene		< 0.04		
Anthracene		< 0.04		
Fluoranthene		< 0.04		
Pyrene		< 0.04		
Benz(a)anthracene		< 0.04		
Chrysene		< 0.04		
Benzo(a)pyrene		< 0.04		
Benzo(b)fluoranther	ne	< 0.04		
Benzo(k)fluoranthei	ne	< 0.04		
Indeno(1,2,3-cd)pyro	ene	< 0.04		
Dibenz(a,h)anthrac	ene	< 0.04		
Benzo(g,h,i)perylene	9	< 0.04		

### ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-03R-0 02/08/19 02/12/19 02/13/19 Water ug/L (ppb)	020819	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Infrastructure Solutions Kelly Moore 14697009 902134-02 1/2 021312.D GCMS6 VM
Surrogates: Anthracene-d10 Benzo(a)anthracene	-d12	% Recovery: 97 111	Lower Limit: 31 25	Upper Limit: 160 165
Compounds:		Concentration ug/L (ppb)		
Naphthalene		< 0.4		
Acenaphthylene		1.7		
Acenaphthene		0.14		
Fluorene		0.099		
Phenanthrene		0.041		
Anthracene		< 0.04		
Fluoranthene		< 0.04		
Pyrene		< 0.04		
Benz(a)anthracene		< 0.04		
Chrysene		< 0.04		
Benzo(a)pyrene		< 0.04		
Benzo(b)fluoranther	ne	< 0.04		
Benzo(k)fluoranther	ne	< 0.04		
Indeno(1,2,3-cd)pyre	ene	< 0.04		
Dibenz(a,h)anthrace	ene	< 0.04		
Benzo(g,h,i)perylene	è.	< 0.04		

### ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-04-02 02/08/19 02/12/19 02/13/19 Water ug/L (ppb)	0719	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Infrastructure Solutions Kelly Moore 14697009 902134-03 1/2 021313.D GCMS6 VM
Surrogates: Anthracene-d10 Benzo(a)anthracene	-d12	% Recovery: 80 80	Lower Limit: 31 25	Upper Limit: 160 165
Compounds:		Concentration ug/L (ppb)		
Naphthalene		3.3		
Acenaphthylene		< 0.04		
Acenaphthene		0.069		
Fluorene		< 0.04		
Phenanthrene		< 0.04		
Anthracene		< 0.04		
Fluoranthene		< 0.04		
Pyrene		< 0.04		
Benz(a)anthracene		< 0.04		
Chrysene		< 0.04		
Benzo(a)pyrene		< 0.04		
Benzo(b)fluoranther	ie	< 0.04		
Benzo(k)fluoranther	ne	< 0.04		
Indeno(1,2,3-cd)pyre	ene	< 0.04		
Dibenz(a,h)anthrace	ene	< 0.04		
Benzo(g,h,i)perylene	9	< 0.04		

### ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-06-02 02/08/19 02/12/19 02/13/19 Water ug/L (ppb)	0719	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Infrastructure Solutions Kelly Moore 14697009 902134-04 1/2 021314.D GCMS6 VM
Surrogates: Anthracene-d10 Benzo(a)anthracene	-d12	% Recovery: 83 87	Lower Limit: 31 25	Upper Limit: 160 165
Compounds:		Concentration		
compounds.		dg/L (ppb)		
Naphthalene		< 0.4		
Acenaphthylene		28 ve		
Acenaphthene		0.71		
Fluorene		0.36		
Phenanthrene		< 0.04		
Anthracene		< 0.04		
Fluoranthene		0.11		
Pyrene		0.15		
Benz(a)anthracene		< 0.04		
Chrysene		< 0.04		
Benzo(a)pyrene		< 0.04		
Benzo(b)fluoranther	ne	< 0.04		
Benzo(k)fluoranther	ne	< 0.04		
Indeno(1,2,3-cd)pyre	ene	< 0.04		
Dibenz(a,h)anthrace	ene	< 0.04		
Benzo(g,h,i)perylene	<u>)</u>	< 0.04		
-				

### ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-06-02 02/08/19 02/12/19 02/14/19 Water ug/L (ppb)	0719	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Infrastructure Solutions Kelly Moore 14697009 902134-04 1/100 021411.D GCMS6 VM
Surrogates: Anthracene-d10 Benzo(a)anthracene	-d12	% Recovery: 80 d 74 d	Lower Limit: 31 25	Upper Limit: 160 165
Compounds:		Concentration ug/L (ppb)		
Naphthalene		<20		
Acenaphthylene		38		
Acenaphthene		<2		
Fluorene		<2		
Phenanthrene		<2		
Anthracene		<2		
Fluoranthene		<2		
Pyrene		<2		
Benz(a)anthracene		<2		
Chrysene		<2		
Benzo(a)pyrene		<2		
Benzo(b)fluoranther	ne	<2		
Benzo(k)fluoranthei	ne	<2		
Indeno(1,2,3-cd)pyro	ene	<2		
Dibenz(a,h)anthrac	ene	<2		
Benzo(g,h,i)perylen	9	<2		

### ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-08-02 02/08/19 02/12/19 02/13/19 Water ug/L (ppb)	0819	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Infrastructure Solutions Kelly Moore 14697009 902134-05 1/2 021317.D GCMS6 VM
Surrogates: Anthracene-d10 Benzo(a)anthracene	-d12	% Recovery: 95 109	Lower Limit: 31 25	Upper Limit: 160 165
Compounds:		Concentration		
compoundo		a8.2 (hha)		
Naphthalene		< 0.4		
Acenaphthylene		< 0.04		
Acenaphthene		0.15		
Fluorene		0.26		
Phenanthrene		< 0.04		
Anthracene		0.095		
Fluoranthene		0.14		
Pyrene		0.20		
Benz(a)anthracene		< 0.04		
Chrysene		< 0.04		
Benzo(a)pyrene		< 0.04		
Benzo(b)fluoranthen	e	< 0.04		
Benzo(k)fluoranther	ne	< 0.04		
Indeno(1,2,3-cd)pyre	ene	< 0.04		
Dibenz(a,h)anthrace	ene	< 0.04		
Benzo(g,h,i)perylene	e e e e e e e e e e e e e e e e e e e	< 0.04		

### ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-09-02 02/08/19 02/12/19 02/13/19 Water ug/L (ppb)	0719	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Infrastructure Solutions Kelly Moore 14697009 902134-06 1/2 021318.D GCMS6 VM
Surrogates: Anthracene-d10 Benzo(a)anthracene	-d12	% Recovery: 83 93	Lower Limit: 31 25	Upper Limit: 160 165
Compounds:		Concentration ug/L (ppb)		
Naphthalene		<0.4		
Acenaphthylene		4.5		
Acenaphthene		2.4		
Fluorene		0.34		
Phenanthrene		< 0.04		
Anthracene		< 0.04		
Fluoranthene		< 0.04		
Pyrene		< 0.04		
Benz(a)anthracene		< 0.04		
Chrysene		< 0.04		
Benzo(a)pyrene		< 0.04		
Benzo(b)fluoranther	ne	< 0.04		
Benzo(k)fluoranther	ne	< 0.04		
Indeno(1,2,3-cd)pyre	ene	< 0.04		
Dibenz(a,h)anthrac	ene	< 0.04		
Benzo(g,h,i)perylene	<del>j</del>	< 0.04		

### ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-10-02 02/08/19 02/12/19 02/13/19 Water ug/L (ppb)	0719	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Infrastructure Solutions Kelly Moore 14697009 902134-07 1/2 021319.D GCMS6 VM
Omts.	ug/L (ppb)		Lower	Unner
Surrogates: Anthracene-d10 Benzo(a)anthracene	-d12	% Recovery: 78 90	Limit: 31 25	Limit: 160 165
Compounds:		Concentration ug/L (ppb)		
Naphthalene		< 0.4		
Acenaphthylene		< 0.04		
Acenaphthene		0.041		
Fluorene		< 0.04		
Phenanthrene		< 0.04		
Anthracene		< 0.04		
Fluoranthene		< 0.04		
Pyrene		< 0.04		
Benz(a)anthracene		< 0.04		
Chrysene		< 0.04		
Benzo(a)pyrene		< 0.04		
Benzo(b)fluoranther	ne	< 0.04		
Benzo(k)fluoranther	ne	< 0.04		
Indeno(1,2,3-cd)pyre	ene	< 0.04		
Dibenz(a,h)anthrac	ene	< 0.04		
Benzo(g,h,i)perylene	9	< 0.04		

### ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-10-9-0 02/08/19 02/12/19 02/13/19 Water ug/L (ppb)	020719	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Infrastructure Solutions Kelly Moore 14697009 902134-08 1/2 021320.D GCMS6 VM
Surrogates: Anthracene-d10 Benzo(a)anthracene	-d12	% Recovery: 89 98	Lower Limit: 31 25	Upper Limit: 160 165
Compounds:		Concentration ug/L (ppb)		
Naphthalene		< 0.4		
Acenaphthylene		< 0.04		
Acenaphthene		0.049		
Fluorene		< 0.04		
Phenanthrene		< 0.04		
Anthracene		< 0.04		
Fluoranthene		< 0.04		
Pyrene		< 0.04		
Benz(a)anthracene		< 0.04		
Chrysene		< 0.04		
Benzo(a)pyrene		< 0.04		
Benzo(b)fluoranther	ne	< 0.04		
Benzo(k)fluoranther	ne	< 0.04		
Indeno(1,2,3-cd)pyre	ene	< 0.04		
Dibenz(a,h)anthrace	ene	< 0.04		
Benzo(g,h,i)perylene	è	< 0.04		

### ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-7-020 02/08/19 02/12/19 02/13/19 Water ug/L (ppb)	819	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Infrastructure Solutions Kelly Moore 14697009 902134-09 1/2 021321.D GCMS6 VM
Surrogates: Anthracene-d10 Benzo(a)anthracene-	·d12	% Recovery: 96 106	Lower Limit: 31 25	Upper Limit: 160 165
Compounds:		Concentration ug/L (ppb)		
Naphthalene		< 0.4		
Acenaphthylene		< 0.04		
Acenaphthene		< 0.04		
Fluorene		< 0.04		
Phenanthrene		< 0.04		
Anthracene		< 0.04		
Fluoranthene		< 0.04		
Pyrene		< 0.04		
Benz(a)anthracene		< 0.04		
Chrysene		< 0.04		
Benzo(a)pyrene		< 0.04		
Benzo(b)fluoranthen	e	< 0.04		
Benzo(k)fluoranthen	ie	< 0.04		
Indeno(1,2,3-cd)pyre	ne	< 0.04		
Dibenz(a,h)anthrace	ene	< 0.04		
Benzo(g,h,i)perylene	•	< 0.04		

### ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	Method Blan Not Applicab 02/12/19 02/13/19 Water ug/L (ppb)	k le	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Envir Kelly Moore 09-359 mb 021306.D GCMS6 VM	ronment & Infrastructure Solutions e 14697009
Surrogates: Anthracene-d10 Banza(a)anthracene	d12	% Recovery: 96	Lower Limit: 31	VIVI	Upper Limit: 160
Compounds:	U12	Concentration ug/L (ppb)	23		105
Naphthalene		<0.2			
Acenaphthylene		< 0.02			
Acenaphthene		< 0.02			
Fluorene		< 0.02			
Phenanthrene		< 0.02			
Anthracene		< 0.02			
Fluoranthene		< 0.02			
Pyrene		< 0.02			
Benz(a)anthracene		< 0.02			
Chrysene		< 0.02			
Benzo(a)pyrene		< 0.02			
Benzo(b)fluoranthen	e	< 0.02			
Benzo(k)fluoranthen	e	< 0.02			
Indeno(1,2,3-cd)pyre	ne	< 0.02			
Dibenz(a,h)anthrace	ene	< 0.02			
Benzo(g,h,i)perylene		< 0.02			

### ENVIRONMENTAL CHEMISTS

# Analysis For Total Metals By EPA Method 6020B

Client ID: Date Received: Date Extracted: Date Analyzed: Matrix:	KMW-02R-020819 02/08/19 02/12/19 02/12/19 Water	Client: Project: Lab ID: Data File: Instrument:	Wood Environment & Infrastructure Solutions Kelly Moore 14697009 902134-01 902134-01.096 ICPMS2
Units:	ug/L (ppb)	Operator:	SP
Analyte:	Concentration ug/L (ppb)		
Arsenic	<1		
Chromium	<1		
Copper	<5		
Lead	<1		
Mercury	<1		
Nickel	2.15		
Zinc	<5		

### ENVIRONMENTAL CHEMISTS

# Analysis For Total Metals By EPA Method 6020B

Client ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-03R-020819 02/08/19 02/12/19 02/12/19 Water ug/L (ppb)	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Infrastructure Solutions Kelly Moore 14697009 902134-02 902134-02.103 ICPMS2 SP
Analyte:	Concentration ug/L (ppb)		
Arsenic	<1		
Chromium	1.47		
Copper	<5		
Lead	<1		
Mercury	<1		
Nickel	1.67		
Zinc	<5		

### ENVIRONMENTAL CHEMISTS

# Analysis For Total Metals By EPA Method 6020B

Client ID:	KMW-04-020719	Client:	Wood Environment & Infrastructure Solutions
Date Received:	02/08/19	Project:	Kelly Moore 14697009
Date Extracted:	02/12/19	Lab ID:	902134-03
Date Analyzed:	02/12/19	Data File:	902134-03.104
Matrix:	Water	Instrument:	ICPMS2
Units:	ug/L (ppb)	Operator:	SP
	Concentration		
Analyte:	ug/L (ppb)		
Arsenic	17.4		
Chromium	1.58		
Copper	24.7		
Lead	1.51		
Mercury	<1		
Nickel	1.98		
Zinc	<5		
## ENVIRONMENTAL CHEMISTS

Client ID: Date Received: Date Extracted: Date Analyzed: Matrix:	KMW-06-020719 02/08/19 02/12/19 02/12/19 Water	Client: Project: Lab ID: Data File: Instrument:	Wood Environment & Infrastructure Solutions Kelly Moore 14697009 902134-04 902134-04.105 ICPMS2
Units:	ug/L (ppb)	Operator:	SP
Analyte:	Concentration ug/L (ppb)		
Arsenic	3.04		
Chromium	2.00		
Copper	19.4		
Lead	2.04		
Mercury	<1		
Nickel	7.79		
Zinc	135		

### ENVIRONMENTAL CHEMISTS

Client ID: Date Received: Date Extracted: Date Analyzed: Matrix:	KMW-08-020819 02/08/19 02/12/19 02/12/19 Water	Client: Project: Lab ID: Data File: Instrument:	Wood Environment & Infrastructure Solutions Kelly Moore 14697009 902134-05 902134-05.109 ICPMS2
Units:	ug/L (ppb)	Operator:	SP
Analyte:	Concentration ug/L (ppb)	-	
Arsenic	<1		
Chromium	<1		
Copper	<5		
Lead	<1		
Mercury	<1		
Nickel	1.99		
Zinc	10.1		

## ENVIRONMENTAL CHEMISTS

Client ID: Date Received: Date Extracted: Date Analyzed: Matrix:	KMW-09-020719 02/08/19 02/12/19 02/12/19 Water	Client: Project: Lab ID: Data File: Instrument:	Wood Environment & Infrastructure Solutions Kelly Moore 14697009 902134-06 902134-06.110 ICPMS2
Units:	ug/L (ppb)	Operator:	SP
Analyte:	Concentration ug/L (ppb)		
Arsenic	<1		
Chromium	1.12		
Copper	<5		
Lead	<1		
Mercury	<1		
Nickel	<1		
Zinc	<5		

## ENVIRONMENTAL CHEMISTS

Client ID:	KMW-10-020719	Client:	Wood Environment & Infrastructure Solutions
Date Received:	02/08/19	Project:	Kelly Moore 14697009
Date Extracted:	02/12/19	Lab ID:	902134-07
Date Analyzed:	02/12/19	Data File:	902134-07.111
Matrix:	Water	Instrument:	ICPMS2
Units:	ug/L (ppb)	Operator:	SP
	Concentration		
Analyte:	ug/L (ppb)		
Arsenic	6.72		
Chromium	2.00 J		
Copper	<5 J		
Lead	<1		
Mercury	<1		
Nickel	<1 J		
Zinc	<5 J		

## ENVIRONMENTAL CHEMISTS

Client ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-10-020719 02/08/19 02/12/19 02/13/19 Water ug/L (ppb)	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Infrastructure Solutions Kelly Moore 14697009 902134-07 x2 902134-07 x2.030 ICPMS2 SP
Analyte:	Concentration ug/L (ppb)		
Arsenic	6.55		
Chromium	2.33		
Copper	<10		
Lead	<2		
Mercury	<2		
Nickel	<2		
Zinc	<10		

## ENVIRONMENTAL CHEMISTS

Client ID: Date Received: Date Extracted:	KMW-10-9-020719 02/08/19 02/12/19	Client: Project: Lab ID:	Wood Environment & Infrastructure Solutions Kelly Moore 14697009 902134-08
Date Analyzed:	02/12/19	Data File:	902134-08.112
Matrix:	water	Instrument:	ICPMIS2
Units:	ug/L (ppb)	Operator:	SP
Analyte:	Concentration ug/L (ppb)		
Arsenic	6.52		
Chromium	2.00 J		
Copper	<5 J		
Lead	<1		
Mercury	<1		
Nickel	<1 J		
Zinc	<5 J		

## ENVIRONMENTAL CHEMISTS

Client ID: Date Received: Date Extracted: Date Analyzed:	KMW-10-9-020719 02/08/19 02/12/19 02/13/19	Client: Project: Lab ID: Data File:	Wood Environment & Infrastructure Solutions Kelly Moore 14697009 902134-08 x2 902134-08 x2.031
Inite	waler	Operatory	
Units:	ug/L (ppb)	Operator:	SP
Analyte:	Concentration ug/L (ppb)		
Arsenic	6.38		
Chromium	2.17		
Copper	<10		
Lead	<2		
Mercury	<2		
Nickel	<2		
Zinc	<10		

## ENVIRONMENTAL CHEMISTS

Client ID: Date Received: Date Extracted:	KMW-7-020819 02/08/19 02/12/19	Client: Project: Lab ID:	Wood Environment & Infrastructure Solutions Kelly Moore 14697009 902134-09
Date Analyzed:	02/12/19	Data File:	902134-09.113
Matrix:	Water	Instrument:	ICPMS2
Units:	ug/L (ppb)	Operator:	SP
Analyte:	Concentration ug/L (ppb)		
Arsenic	<1		
Chromium	<1		
Copper	<5		
Lead	<1		
Mercury	<1		
Nickel	1.36		
Zinc	<5		

## ENVIRONMENTAL CHEMISTS

Client ID:	Method Blank	Client:	Wood Environment & Infrastructure Solutions
Date Received:	NA	Project:	Kelly Moore 14697009
Date Extracted:	02/12/19	Lab ID:	I9-95 mb
Date Analyzed:	02/12/19	Data File:	I9-95 mb.094
Matrix:	Water	Instrument:	ICPMS2
Units:	ug/L (ppb)	Operator:	SP
	Concentration		
Analyte:	ug/L (ppb)		
Arsenic	<1		
Chromium	<1		
Copper	<5		
Lead	<1		
Mercury	<1		
Nickel	<1		
Zinc	<5		

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-02R-0 02/08/19 02/14/15 02/14/19 Water ug/L (ppb)	020819	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Kelly Moore 14697009 902134-01 021421.D GCMS4 MS	& Infrastructure Solutions 9
			Lower	Upper	
Surrogates:		% Recovery:	Limit:	Limit:	
1,2-Dichloroethane-d	4	100	57	121	
Toluene-d8		100	63	127	
4-Bromofluorobenzer	ne	100	60	133	
		Concentration			Concentration
Compounds:		ug/L (ppb)	Compour	nds:	ug/L (ppb)
Dichlorodifluorometl	nane	<1	1,3-Dichl	oropropane	<1
Chloromethane		<10	Tetrachl	oroethene	<1
Vinyl chloride		< 0.2	Dibromo	chloromethane	<1
Bromomethane		<1	1,2-Dibro	omoethane (EDB)	<1
Chloroethane		<1	Chlorobe	nzene	<1
Trichlorofluorometh	ane	<1	Ethvlber	izene	<1
Acetone		<50	1,1,1,2-T	etrachloroethane	<1
1,1-Dichloroethene		<1	m,p-Xyle	ene	<2
Hexane		<1	o-Xylene		<1
Methylene chloride		<5	Stvrene		<1
Methyl t-butyl ether	(MTBE)	<1	Isopropy	lbenzene	<1
trans-1.2-Dichloroet	hene	<1	Bromofor	rm	<1
1.1-Dichloroethane		<1	n-Propyl	benzene	<1
2.2-Dichloropropane		<1	Bromobe	nzene	<1
cis-1.2-Dichloroether	ne	<1	1.3.5-Tri	methylbenzene	<1
Chloroform		<1	1.1.2.2-T	etrachloroethane	<1
2-Butanone (MEK)		<10	1.2.3-Tri	chloropropane	<1
1.2-Dichloroethane (	EDC)	<1	2-Chloro	toluene	<1
1.1.1-Trichloroethan	e	<1	4-Chloro	toluene	<1
1.1-Dichloropropene		<1	tert-Buty	lbenzene	<1
Carbon tetrachloride		<1	1.2.4-Tri	methylbenzene	<1
Benzene		< 0.35	sec-Buty	lbenzene	<1
Trichloroethene		<1	p-Isopror	ovltoluene	<1
1.2-Dichloropropane		<1	1.3-Dich	lorobenzene	<1
Bromodichlorometha	nne	<1	1 4-Dich	lorobenzene	<1
Dibromomethane		<1	1 2-Dich	lorobenzene	<1
4-Methyl-2-pentanor	e	<10	1,2-Dibro	omo-3-chloropropane	<10
cis-1.3-Dichloroprope	ene	<1	1,2.4-Tri	chlorobenzene	<1
Toluene		<1	Hexachle	probutadiene	<1
trans-1.3-Dichloropr	opene	<1	Nanhtha	lene	<1
1.1.2-Trichloroethan	e	<1	1.2.3-Tri	chlorobenzene	<1
2-Hexanone		<10	, ,		

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-03R-0 02/08/19 02/14/15 02/14/19 Water ug/L (ppb)	020819	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Kelly Moore 14697009 902134-02 021422.D GCMS4 MS	Infrastructure Solutions
			Lower	Upper	
Surrogates:		% Recovery:	Limit:	Limit:	
1,2-Dichloroethane-o	14	98	57	121	
Toluene-d8		99	63	127	
4-Bromofluorobenze	ne	99	60	133	
		Concentration			Concentration
Compounds:		ug/L (ppb)	Compour	nds:	ug/L (ppb)
Dichlorodifluoromet	hane	<1	1.3-Dich	oropropane	<1
Chloromethane		<10	Tetrachl	oroethene	<1
Vinvl chloride		< 0.2	Dibromo	chloromethane	<1
Bromomethane		<1	1.2-Dibro	omoethane (EDB)	<1
Chloroethane		<1	Chlorobe	nzene	<1
Trichlorofluorometh	ane	<1	Ethvlber	izene	<1
Acetone		<50	1,1,1,2-T	etrachloroethane	<1
1,1-Dichloroethene		<1	m,p-Xyle	ene	<2
Hexane		<1	o-Xylene		<1
Methylene chloride		<5	Styrene		<1
Methyl t-butyl ether	(MTBE)	<1	Isopropy	lbenzene	<1
trans-1,2-Dichloroet	hene	<1	Bromofo	rm	<1
1,1-Dichloroethane		<1	n-Propyl	benzene	1.5
2,2-Dichloropropane		<1	Bromobe	nzene	<1
cis-1,2-Dichloroether	ne	<1	1,3,5-Tri	methylbenzene	<1
Chloroform		<1	1,1,2,2-T	'etrachloroethane	<1
2-Butanone (MEK)		<10	1,2,3-Tri	chloropropane	<1
1,2-Dichloroethane (	EDC)	<1	2-Chloro	toluene	<1
1,1,1-Trichloroethar	ne	<1	4-Chloro	toluene	<1
1,1-Dichloropropene		<1	tert-Buty	lbenzene	<1
Carbon tetrachloride	<u>e</u>	<1	1,2,4-Tri	methylbenzene	<1
Benzene		< 0.35	sec-Buty	lbenzene	<1
Trichloroethene		<1	p-Isoprop	yltoluene	<1
1,2-Dichloropropane		<1	1,3-Dich	lorobenzene	<1
Bromodichlorometha	ane	<1	1,4-Dich	lorobenzene	<1
Dibromomethane		<1	1,2-Dichl	lorobenzene	<1
4-Methyl-2-pentanor	ne	<10	1,2-Dibro	omo-3-chloropropane	<10
cis-1,3-Dichloroprop	ene	<1	1,2,4-Tri	chlorobenzene	<1
Toluene		<1	Hexachle	orobutadiene	<1
trans-1,3-Dichloropr	opene	<1	Naphtha	lene	<1
1,1,2-Trichloroethar	ne	<1	1,2,3-Tri	chlorobenzene	<1
2-Hexanone		<10			

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-04-02 02/08/19 02/14/15 02/14/19 Water ug/L (ppb)	0719	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Kelly Moore 14697009 902134-03 1/10 021431.D GCMS4 MS	Infrastructure Solutions
			Lower	Upper	
Surrogates:		% Recovery:	Limit:	Limit:	
1,2-Dichloroethane-c	14	103	57	121	
Toluene-d8		94	63	127	
4-Bromofluorobenzer	ne	90	60	133	
		Concentration			Concentration
Compounds:		ug/L (ppb)	Compour	nds:	ug/L (ppb)
Dichlorodifluoromet	hane	<10	1.3-Dichl	oropropane	<10
Chloromethane		<100	Tetrachl	proethene	<10
Vinvl chloride		<2	Dibromo	chloromethane	<10
Bromomethane		<10	1.2-Dibro	omoethane (EDB)	<10
Chloroethane		<10	Chlorobe	nzene	<10
Trichlorofluorometh	ane	<10	Ethylben	zene	2.500 ve
Acetone		<500	1,1,1,2-T	etrachloroethane	<10
1,1-Dichloroethene		<10	m,p-Xyle	ene	5,400 ve
Hexane		<10	o-Xylene		1,200
Methylene chloride		<50	Styrene		<10
Methyl t-butyl ether	(MTBE)	<10	Isopropy	lbenzene	26
trans-1,2-Dichloroet	hene	<10	Bromofor	rm	<10
1,1-Dichloroethane		<10	n-Propyl	benzene	23
2,2-Dichloropropane		<10	Bromobe	nzene	<10
cis-1,2-Dichloroether	ne	<10	1,3,5-Tri	methylbenzene	33
Chloroform		<10	1,1,2,2-T	etrachloroethane	<10
2-Butanone (MEK)		<100	1,2,3-Tri	chloropropane	<10
1,2-Dichloroethane (	EDC)	<10	2-Chloro	toluene	<10
1,1,1-Trichloroethan	ie	<10	4-Chlorot	toluene	<10
1,1-Dichloropropene		<10	tert-Buty	lbenzene	<10
Carbon tetrachloride	<u>)</u>	<10	1,2,4-Tri	methylbenzene	67
Benzene		<3.5	sec-Buty	lbenzene	<10
Trichloroethene		<10	p-Isoprop	yltoluene	<10
1,2-Dichloropropane		<10	1,3-Dichl	orobenzene	<10
Bromodichlorometha	ane	<10	1,4-Dichl	orobenzene	<10
Dibromomethane		<10	1,2-Dichl	orobenzene	<10
4-Methyl-2-pentanor	ne	<100	1,2-Dibro	omo-3-chloropropane	<100
cis-1,3-Dichloroprope	ene	<10	1,2,4-Tri	chlorobenzene	<10
Toluene		190	Hexachle	orobutadiene	<10
trans-1,3-Dichloropr	opene	<10	Naphtha	lene	<10
1,1,2-Trichloroethan	ie	<10	1,2,3-Tri	chlorobenzene	<10
2-Hexanone		<100			

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-04-02 02/08/19 02/14/15 02/14/19 Water ug/L (ppb)	0719	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Kelly Moore 14697009 902134-03 1/100 021430.D GCMS4 MS	: Infrastructure Solutions )
			Lower	Upper	
Surrogates:	• -	% Recovery:	Limit:	Limit:	
1,2-Dichloroethane-c	14	99	57	121	
Toluene-d8		100	63	127	
4-Bromofluorobenzei	ne	102	60	133	
		Concentration			Concentration
Compounds:		ug/L (ppb)	Compour	nds:	ug/L (ppb)
Dichlorodifluoromet	hane	<100	1,3-Dichl	oropropane	<100
Chloromethane		<1,000	Tetrachle	oroethene	<100
Vinyl chloride		<20	Dibromo	chloromethane	<100
Bromomethane		<100	1,2-Dibro	omoethane (EDB)	<100
Chloroethane		<100	Chlorobe	nzene	<100
Trichlorofluorometh	ane	<100	Ethylben	zene	2,800
Acetone		<5,000	1,1,1,2-T	etrachloroethane	<100
1,1-Dichloroethene		<100	m,p-Xyle	ene	6,100
Hexane		<100	o-Xylene		1,300
Methylene chloride		<500	Styrene		<100
Methyl t-butyl ether	(MTBE)	<100	Isopropylbenzene		<100
trans-1,2-Dichloroet	hene	<100	Bromoform		<100
1,1-Dichloroethane		<100	n-Propylbenzene		<100
2,2-Dichloropropane		<100	Bromobe	nzene	<100
cis-1,2-Dichloroether	ne	<100	1,3,5-Tri	methylbenzene	<100
Chloroform		<100	1,1,2,2-T	etrachloroethane	<100
2-Butanone (MEK)		<1,000	1,2,3-Tri	chloropropane	<100
1,2-Dichloroethane (	EDC)	<100	2-Chloro	toluene	<100
1,1,1-Trichloroethan	ne	<100	4-Chloro	toluene	<100
1,1-Dichloropropene		<100	tert-Buty	lbenzene	<100
Carbon tetrachloride	9	<100	1,2,4-Tri	methylbenzene	<100
Benzene		<35	sec-Buty	lbenzene	<100
Trichloroethene		<100	p-Isoprop	oyltoluene	<100
1,2-Dichloropropane		<100	1,3-Dichl	orobenzene	<100
Bromodichlorometha	ane	<100	1,4-Dichl	orobenzene	<100
Dibromomethane		<100	1,2-Dichl	orobenzene	<100
4-Methyl-2-pentanor	ne	<1,000	1,2-Dibro	omo-3-chloropropane	<1,000
cis-1,3-Dichloroprop	ene	<100	1,2,4-Tri	chlorobenzene	<100
Toluene		190	Hexachle	probutadiene	<100
trans-1,3-Dichloropr	opene	<100	Naphtha	lene	<100
1,1,2-Trichloroethan	ne	<100	1,2,3-Tri	chlorobenzene	<100
2-Hexanone		<1,000			

## ENVIRONMENTAL CHEMISTS

Client Sample ID:KMW-06-020719Client:Wood Environment & InfrastructDate Received:02/08/19Project:Kelly Moore 14697009Date Extracted:02/14/15Lab ID:902134-04Date Analyzed:02/14/19Data File:021428.DMatrix:WaterInstrument:GCMS4Units:ug/L (ppb)Operator:MS	ture Solutions
Lower Upper	
Surrogates: % Recovery: Limit: Limit:	
1,2-Dichloroethane-d4 99 57 121	
Toluene-d8 99 63 127	
4-Bromofluorobenzene 106 60 133	
Concentration Concent	ration
Compounds: ug/L (ppb) Compounds: ug/L (j	ppb)
Dichlorodifluoromethane <1 1.3-Dichloropropane <1	
Chloromethane <10 Tetrachloroethene <1	
Vinyl chloride <0.2 Dibromochloromethane <1	
Bromomethane <1 1.2-Dibromoethane (EDB) <1	
Chloroethane <1 Chlorobenzene <1	
Trichlorofluoromethane <1 Ethylbenzene <1	
Acetone <50 1.1.1.2-Tetrachloroethane <1	
1.1-Dichloroethene <1 m.p-Xylene <2	
Hexane <1 o-Xvlene <1	
Methylene chloride <5 Styrene <1	
Methyl t-butyl ether (MTBE) <1 Isopropylbenzene 8.5	3
trans-1.2-Dichloroethene <1 Bromoform <1	
1.1-Dichloroethane <1 n-Propylbenzene 12	
2.2-Dichloropropane <1 Bromobenzene <1	
cis-1,2-Dichloroethene <1 1.3.5-Trimethylbenzene <1	
Chloroform <1 1122-Tetrachloroethane <1	
2-Butanone (MEK) <10 1.2.3-Trichloropropane <1	
1 2-Dichloroethane (FDC) <1 2-Chlorotoluene <1	
1 1 1-Trichloroethane <1 4-Chlorotoluene <1	
1 1-Dichloropropene <1 tert-Butylbenzene <1	
Carbon tetrachloride <1 1 2 4-Trimethylbenzene <1	
Benzene <0.35 sec-Butylbenzene 1.5	2
Trichloroethene <1 p-Isopropyltoluene <1	~
1 2-Dichloropropane <1 1 3-Dichlorobenzene <1	
Bromodichloromethane <1 14-Dichlorobenzene <1	
Dibromomethane <1 1 2-Dichlorobenzene <1	
4-Methyl-2-pentanone <10 1 2-Dibromo-3-chloropropane <10	
cis-1 3-Dichloropropene <1 1 2 4-Trichlorobenzene <1	
Toluene <1 Hexachlorobutadiene <1	
trans-1 3-Dichloropropene <1 Nanhthalene <1	
1 1 2-Trichloroethane <1 1 2 3-Trichlorobenzene <1	
2-Hexanone <10	

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-08-02 02/08/19 02/14/15 02/14/19 Water ug/L (ppb)	0819	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Kelly Moore 14697009 902134-05 021423.D GCMS4 MS	Infrastructure Solutions
			Lower	Upper	
Surrogates:		% Recovery:	Limit:	Limit:	
1,2-Dichloroethane-o	14	107	57	121	
Toluene-d8		98	63	127	
4-Bromofluorobenze	ne	95	60	133	
		Concentration			Concentration
Compounds:		ug/L (ppb)	Compour	nds:	ug/L (ppb)
Dichlorodifluoromet	hane	<1	1.3-Dich	oropropane	<1
Chloromethane		<10	Tetrachl	oroethene	<1
Vinvl chloride		< 0.2	Dibromo	chloromethane	<1
Bromomethane		<1	1.2-Dibro	omoethane (EDB)	<1
Chloroethane		<1	Chlorobe	nzene	<1
Trichlorofluorometh	ane	<1	Ethylber	izene	<1
Acetone		<50	1.1.1.2-T	etrachloroethane	<1
1,1-Dichloroethene		<1	m,p-Xyle	ene	<2
Hexane		<1	o-Xylene		<1
Methylene chloride		<5	Styrene		<1
Methyl t-butyl ether	(MTBE)	<1	Isopropy	lbenzene	<1
trans-1,2-Dichloroet	hene	<1	Bromofo	rm	<1
1,1-Dichloroethane		<1	n-Propyl	benzene	<1
2,2-Dichloropropane		<1	Bromobe	nzene	<1
cis-1,2-Dichloroethe	ne	<1	1,3,5-Tri	methylbenzene	<1
Chloroform		<1	1,1,2,2-T	'etrachloroethane	<1
2-Butanone (MEK)		<10	1,2,3-Tri	chloropropane	<1
1,2-Dichloroethane (	EDC)	<1	2-Chloro	toluene	<1
1,1,1-Trichloroethar	ne	<1	4-Chloro	toluene	<1
1,1-Dichloropropene		<1	tert-Buty	lbenzene	<1
Carbon tetrachloride	<u>e</u>	<1	1,2,4-Tri	methylbenzene	<1
Benzene		< 0.35	sec-Buty	lbenzene	<1
Trichloroethene		<1	p-Isoprop	yltoluene	<1
1,2-Dichloropropane		<1	1,3-Dich	lorobenzene	<1
Bromodichlorometha	ane	<1	1,4-Dich	lorobenzene	<1
Dibromomethane		<1	1,2-Dich	lorobenzene	<1
4-Methyl-2-pentanor	ne	<10	1,2-Dibro	omo-3-chloropropane	<10
cis-1,3-Dichloroprop	ene	<1	1,2,4-Tri	chlorobenzene	<1
Toluene		<1	Hexachle	orobutadiene	<1
trans-1,3-Dichloropr	opene	<1	Naphtha	lene	<1
1,1,2-Trichloroethar	ne	<1	1,2,3-Tri	chlorobenzene	<1
2-Hexanone		<10			

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-09-02 02/08/19 02/14/15 02/14/19 Water ug/L (ppb)	0719	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Kelly Moore 14697009 902134-06 021424.D GCMS4 MS	Infrastructure Solutions
			Lower	Upper	
Surrogates:		% Recovery:	Limit:	Limit:	
1,2-Dichloroethane-o	14	101	57	121	
Toluene-d8		107	63	127	
4-Bromofluorobenzer	ne	100	60	133	
		Concentration			Concentration
Compounds:		ug/L (ppb)	Compour	nds:	ug/L (ppb)
Dichlorodifluoromet	hane	<1	1.3-Dichl	oropropape	<1
Chloromethane	nane	<10	Tetrachl	proethene	<1
Vinvl chloride		<0.2	Dibromo	chloromethane	<1
Bromomethane		<1	1 2-Dibro	omoethane (EDB)	<1
Chloroethane		<1	Chlorobe	nzene	<1
Trichlorofluorometh	ane	<1	Ethylber	zene	<1
Acetone	une	<50	1.1.1.2-T	etrachloroethane	<1
1.1-Dichloroethene		<1	m.p-Xvle	ne	<2
Hexane		<1	o-Xvlene		<1
Methylene chloride		<5	Styrene		<1
Methyl t-butyl ether	(MTBE)	<1	Isopropy	lbenzene	1.6
trans-1.2-Dichloroet	hene	<1	Bromofor	m	<1
1,1-Dichloroethane		<1	n-Propyl	benzene	1.6
2,2-Dichloropropane		<1	Bromobe	nzene	<1
cis-1,2-Dichloroether	ne	<1	1,3,5-Tri	methylbenzene	<1
Chloroform		<1	1,1,2,2-T	etrachloroethane	<1
2-Butanone (MEK)		<10	1,2,3-Tri	chloropropane	<1
1,2-Dichloroethane (	EDC)	<1	2-Chloro	toluene	<1
1,1,1-Trichloroethan	ie	<1	4-Chloro	toluene	<1
1,1-Dichloropropene		<1	tert-Buty	lbenzene	<1
Carbon tetrachloride	)	<1	1,2,4-Tri	methylbenzene	<1
Benzene		< 0.35	sec-Buty	lbenzene	<1
Trichloroethene		<1	p-Isoprop	yltoluene	<1
1,2-Dichloropropane		<1	1,3-Dichl	orobenzene	<1
Bromodichlorometha	ane	<1	1,4-Dichl	orobenzene	<1
Dibromomethane		<1	1,2-Dichl	orobenzene	<1
4-Methyl-2-pentanor	ne	<10	1,2-Dibro	omo-3-chloropropane	<10
cis-1,3-Dichloroprop	ene	<1	1,2,4-Tri	chlorobenzene	<1
Toluene		<1	Hexachle	orobutadiene	<1
trans-1,3-Dichloropr	opene	<1	Naphtha	lene	<1
1,1,2-Trichloroethan	ie	<1	1,2,3-Tri	chlorobenzene	<1
2-Hexanone		<10			

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-10-02 02/08/19 02/14/15 02/14/19 Water ug/L (ppb)	0719	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Kelly Moore 14697009 902134-07 021425.D GCMS4 MS	2 Infrastructure Solutions )
			Lower	Upper	
Surrogates:		% Recovery:	Limit:	Limit:	
1,2-Dichloroethane-o	14	99	57	121	
Toluene-d8		106	63	127	
4-Bromofluorobenze	ne	105	60	133	
		Concentration			Concentration
Compounds:		ug/L (ppb)	Compour	nds:	ug/L (ppb)
Dichlorodifluoromet	hane	<1	1.3-Dichl	oropropane	<1
Chloromethane		<10	Tetrachl	oroethene	<1
Vinvl chloride		< 0.2	Dibromo	chloromethane	<1
Bromomethane		<1	1.2-Dibro	omoethane (EDB)	<1
Chloroethane		<1	Chlorobe	nzene	<1
Trichlorofluorometh	ane	<1	Ethylber	izene	<1
Acetone		<50	1.1.1.2-T	'etrachloroethane	<1
1.1-Dichloroethene		<1	m.p-Xvle	ene	6.0
Hexane		<1	o-Xvlene		<1
Methylene chloride		<5	Styrene		<1
Methyl t-butyl ether	(MTBE)	<1	Isopropy	lbenzene	10
trans-1.2-Dichloroet	hene	<1	Bromofor	rm	<1
1.1-Dichloroethane		<1	n-Propyl	benzene	8.5
2.2-Dichloropropane		<1	Bromobe	nzene	<1
cis-1.2-Dichloroethe	ne	<1	1.3.5-Tri	methylbenzene	<1
Chloroform		<1	1,1,2,2-T	etrachloroethane	<1
2-Butanone (MEK)		<10	1.2.3-Tri	chloropropane	<1
1.2-Dichloroethane (	EDC)	<1	2-Chloro	toluene	<1
1.1.1-Trichloroethar	22 0) 10	<1	4-Chlorot	toluene	<1
1.1-Dichloropropene		<1	tert-Buty	lbenzene	<1
Carbon tetrachloride	ć	<1	1.2.4-Tri	methylbenzene	<1
Benzene	-	< 0.35	sec-Buty	lbenzene	<1
Trichloroethene		<1	p-Isopror	ovltoluene	<1
1.2-Dichloropropane		<1	1.3-Dichl	lorobenzene	<1
Bromodichlorometha	ane	<1	1.4-Dichl	orobenzene	<1
Dibromomethane		<1	1,2-Dichl	lorobenzene	<1
4-Methyl-2-pentanor	ne	<10	1,2 Dibro	omo-3-chloropropane	<10
cis-1 3-Dichloroprop	ene	<1	1,2 2.151	chlorobenzene	<1
Toluene		<1	Hexachle	probutadiene	<1
trans-1.3-Dichloropr	opene	<1	Nanhtha	lene	<1
1.1.2-Trichloroethar		<1	1 2 3-Tri	chlorobenzene	<1
2-Hexanone	-	<10	1,2,0 111		

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-10-9-0 02/08/19 02/14/15 02/14/19 Water ug/L (ppb)	020719	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Kelly Moore 14697009 902134-08 021426.D GCMS4 MS	Infrastructure Solutions
			Lower	Upper	
Surrogates:		% Recovery:	Limit:	Limit:	
1,2-Dichloroethane-c	14	99	57	121	
Toluene-d8		104	63	127	
4-Bromofluorobenzer	ne	99	60	133	
		Concentration			Concentration
Compounds:		ug/L (ppb)	Compour	nds:	ug/L (ppb)
Dichlorodifluoromet	hane	<1	1.3-Dichl	oropropane	<1
Chloromethane		<10	Tetrachl	proethene	<1
Vinvl chloride		< 0.2	Dibromo	chloromethane	<1
Bromomethane		<1	1.2-Dibro	omoethane (EDB)	<1
Chloroethane		<1	Chlorobe	nzene	<1
Trichlorofluorometh	ane	<1	Ethylben	izene	<1
Acetone		<50	1,1,1,2-T	etrachloroethane	<1
1,1-Dichloroethene		<1	m,p-Xyle	ene	6.1
Hexane		<1	o-Xylene		<1
Methylene chloride		<5	Styrene		<1
Methyl t-butyl ether	(MTBE)	<1	Isopropy	lbenzene	9.1
trans-1,2-Dichloroet	hene	<1	Bromofor	rm	<1
1,1-Dichloroethane		<1	n-Propyl	benzene	7.5
2,2-Dichloropropane		<1	Bromobe	nzene	<1
cis-1,2-Dichloroether	ne	<1	1,3,5-Tri	methylbenzene	<1
Chloroform		<1	1,1,2,2-T	etrachloroethane	<1
2-Butanone (MEK)		<10	1,2,3-Tri	chloropropane	<1
1,2-Dichloroethane (	EDC)	<1	2-Chloro	toluene	<1
1,1,1-Trichloroethan	ie	<1	4-Chloro	toluene	<1
1,1-Dichloropropene		<1	tert-Buty	lbenzene	<1
Carbon tetrachloride	<u>)</u>	<1	1,2,4-Tri	methylbenzene	<1
Benzene		< 0.35	sec-Buty	lbenzene	<1
Trichloroethene		<1	p-Isoprop	oyltoluene	<1
1,2-Dichloropropane		<1	1,3-Dichl	orobenzene	<1
Bromodichlorometha	ane	<1	1,4-Dichl	orobenzene	<1
Dibromomethane		<1	1,2-Dichl	orobenzene	<1
4-Methyl-2-pentanor	ne	<10	1,2-Dibro	omo-3-chloropropane	<10
cis-1,3-Dichloroprop	ene	<1	1,2,4-Tri	chlorobenzene	<1
Toluene		<1	Hexachle	orobutadiene	<1
trans-1,3-Dichloropr	opene	<1	Naphtha	lene	<1
1,1,2-Trichloroethan	ie	<1	1,2,3-Tri	chlorobenzene	<1
2-Hexanone		<10			

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-7-020 02/08/19 02/14/15 02/14/19 Water ug/L (ppb)	819	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Kelly Moore 14697009 902134-09 021427.D GCMS4 MS	Infrastructure Solutions
			Lower	Upper	
Surrogates:		% Recovery:	Limit:	Limit:	
1,2-Dichloroethane-o	14	100	57	121	
Toluene-d8		94	63	127	
4-Bromofluorobenze	ne	99	60	133	
		Concentration			Concentration
Compounds:		ug/L (ppb)	Compour	nds:	ug/L (ppb)
Dichlorodifluoromet	hane	<1	1.3-Dich	loropropane	<1
Chloromethane		<10	Tetrachl	oroethene	<1
Vinvl chloride		< 0.2	Dibromo	chloromethane	<1
Bromomethane		<1	1,2-Dibro	omoethane (EDB)	<1
Chloroethane		<1	Chlorobe	nzene	<1
Trichlorofluorometh	ane	<1	Ethylber	izene	<1
Acetone		<50	1,1,1,2-T	etrachloroethane	<1
1,1-Dichloroethene		<1	m,p-Xyle	ene	<2
Hexane		<1	o-Xylene		<1
Methylene chloride		<5	Styrene		<1
Methyl t-butyl ether	· (MTBE)	<1	Isopropy	lbenzene	<1
trans-1,2-Dichloroet	hene	<1	Bromofo	rm	<1
1,1-Dichloroethane		<1	n-Propyl	benzene	<1
2,2-Dichloropropane	1	<1	Bromobe	nzene	<1
cis-1,2-Dichloroethe	ne	<1	1,3,5-Tri	methylbenzene	<1
Chloroform		<1	1,1,2,2-T	etrachloroethane	<1
2-Butanone (MEK)		<10	1,2,3-Tri	chloropropane	<1
1,2-Dichloroethane (	EDC)	<1	2-Chloro	toluene	<1
1,1,1-Trichloroethar	ne	<1	4-Chloro	toluene	<1
1,1-Dichloropropene		<1	tert-Buty	lbenzene	<1
Carbon tetrachloride	e	<1	1,2,4-Tri	methylbenzene	<1
Benzene		< 0.35	sec-Buty	lbenzene	<1
Trichloroethene		<1	p-Isoprop	oyltoluene	<1
1,2-Dichloropropane	:	<1	1,3-Dichl	lorobenzene	<1
Bromodichlorometha	ane	<1	1,4-Dich	lorobenzene	<1
Dibromomethane		<1	1,2-Dichl	lorobenzene	<1
4-Methyl-2-pentanor	ne	<10	1,2-Dibro	omo-3-chloropropane	<10
cis-1,3-Dichloroprop	ene	<1	1,2,4-Tri	chlorobenzene	<1
Toluene		<1	Hexachle	orobutadiene	<1
trans-1,3-Dichloropr	ropene	<1	Naphtha	lene	<1
1,1,2-Trichloroethar	ne	<1	1,2,3-Tri	chlorobenzene	<1
2-Hexanone		<10			

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	Trip Blanks 02/08/19 02/14/15 02/14/19 Water ug/L (ppb)		Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Kelly Moore 14697009 902134-10 021429.D GCMS4 MS	Infrastructure Solutions
			Lower	Upper	
Surrogates:		% Recovery:	Limit:	Limit:	
1,2-Dichloroethane-o	14	99	57	121	
Toluene-d8		103	63	127	
4-Bromofluorobenze	ne	100	60	133	
		Concentration			Concentration
Compounds:		ug/L (ppb)	Compour	nds:	ug/L (ppb)
Dichlorodifluoromet	hane	<1	1.3-Dichl	oropropane	<1
Chloromethane		<10	Tetrachl	oroethene	<1
Vinyl chloride		< 0.2	Dibromo	chloromethane	<1
Bromomethane		<1	1,2-Dibro	omoethane (EDB)	<1
Chloroethane		<1	Chlorobe	nzene	<1
Trichlorofluorometh	ane	<1	Ethylber	izene	<1
Acetone		<50	1,1,1,2-T	etrachloroethane	<1
1,1-Dichloroethene		<1	m,p-Xyle	ene	<2
Hexane		<1	o-Xylene		<1
Methylene chloride		<5	Styrene		<1
Methyl t-butyl ether	· (MTBE)	<1	Isopropylbenzene		<1
trans-1,2-Dichloroet	hene	<1	Bromoform		<1
1,1-Dichloroethane		<1	n-Propylbenzene		<1
2,2-Dichloropropane		<1	Bromobe	nzene	<1
cis-1,2-Dichloroethe	ne	<1	1,3,5-Tri	methylbenzene	<1
Chloroform		<1	1,1,2,2-T	etrachloroethane	<1
2-Butanone (MEK)		<10	1,2,3-Tri	chloropropane	<1
1,2-Dichloroethane (	EDC)	<1	2-Chloro	toluene	<1
1,1,1-Trichloroethar	ne	<1	4-Chloro	toluene	<1
1,1-Dichloropropene		<1	tert-Buty	lbenzene	<1
Carbon tetrachloride	<del>j</del>	<1	1,2,4-Tri	methylbenzene	<1
Benzene		< 0.35	sec-Buty	lbenzene	<1
Trichloroethene		<1	p-Isoprop	oyltoluene	<1
1,2-Dichloropropane		<1	1,3-Dichl	lorobenzene	<1
Bromodichlorometha	ane	<1	1,4-Dich	lorobenzene	<1
Dibromomethane		<1	1,2-Dichl	lorobenzene	<1
4-Methyl-2-pentanor	ne	<10	1,2-Dibro	omo-3-chloropropane	<10
cis-1,3-Dichloroprop	ene	<1	1,2,4-Tri	chlorobenzene	<1
Toluene		<1	Hexachle	probutadiene	<1
trans-1,3-Dichloropr	opene	<1	Naphtha	lene	<1
1,1,2-Trichloroethar	ne	<1	1,2,3-Tri	chlorobenzene	<1
2-Hexanone		<10			

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	Method Blan Not Applicat 02/14/19 02/14/19 Water ug/L (ppb)	k le	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Kelly Moore 1469700 09-0281 mb 021413.D GCMS4 MS	& Infrastructure Solutions 19
			Lower	Upper	
Surrogates:		% Recovery:	Limit:	Limit:	
1,2-Dichloroethane-d	4	105	57	121	
Toluene-d8		102	63	127	
4-Bromofluorobenzer	ne	103	60	133	
		Concentration			Concentration
Compounds:		ug/L (ppb)	Compour	nds:	ug/L (ppb)
Dichlorodifluorometh	nane	<1	1,3-Dichl	oropropane	<1
Chloromethane		<10	Tetrachl	oroethene	<1
Vinvl chloride		<0.2	Dibromo	chloromethane	<1
Bromomethane		<1	1.2-Dibro	omoethane (EDB)	<1
Chloroethane		<1	Chlorobe	nzene	<1
Trichlorofluorometha	ane	<1	Ethvlber	zene	<1
Acetone		<50	1.1.1.2-T	etrachloroethane	<1
1.1-Dichloroethene		<1	m.p-Xvle	ene	<2
Hexane		<1	o-Xvlene		<1
Methylene chloride		<5	Styrene		<1
Methyl t-butyl ether	(MTBE)	<1	Isopropy	lbenzene	<1
trans-1.2-Dichloroeth	nene	<1	Bromofor	rm	<1
1.1-Dichloroethane		<1	n-Propyl	benzene	<1
2.2-Dichloropropane		<1	Bromobe	nzene	<1
cis-1.2-Dichloroether	ne	<1	1.3.5-Tri	methylbenzene	<1
Chloroform		<1	1.1.2.2-T	etrachloroethane	<1
2-Butanone (MEK)		<10	1.2.3-Tri	chloropropane	<1
1.2-Dichloroethane (I	EDC)	<1	2-Chloro	toluene	<1
1.1.1-Trichloroethan	e	<1	4-Chloro	toluene	<1
1.1-Dichloropropene		<1	tert-Buty	lbenzene	<1
Carbon tetrachloride		<1	1.2.4-Tri	methylbenzene	<1
Benzene		< 0.35	sec-Buty	lbenzene	<1
Trichloroethene		<1	p-Isopror	ovltoluene	<1
1.2-Dichloropropane		<1	1.3-Dich	orobenzene	<1
Bromodichlorometha	ine	<1	1.4-Dich	orobenzene	<1
Dibromomethane		<1	1.2-Dich	orobenzene	<1
4-Methyl-2-pentanon	е	<10	1.2-Dibro	omo-3-chloropropane	<10
cis-1.3-Dichloroprope	ene	<1	1.2.4-Tri	chlorobenzene	<1
Toluene	-	<1	Hexachle	probutadiene	<1
trans-1.3-Dichloropro	opene	<1	Naphtha	lene	<1
1,1,2-Trichloroethan	e	<1	1.2.3-Tri	chlorobenzene	<1
2-Hexanone		<10			

#### ENVIRONMENTAL CHEMISTS

Date of Report: 02/18/19 Date Received: 02/08/19 Project: Kelly Moore 14697009, F&BI 902134

#### QUALITY ASSURANCE RESULTS FOR THE ANALYSIS OF WATER SAMPLES FOR TPH AS GASOLINE USING METHOD NWTPH-Gx

Laboratory (	Code: 902134-04 (Ma	trix Spike)					
-		_		Percent	Percent		
			Sample	Recovery	Recovery	Acceptance	RPD
Analyte	Reporting Units	Spike Level	Result	MS	MSD	Criteria	(Limit 20)
Gasoline	ug/L (ppb)	1,000	2,200	140 b	154 b	53-117	10
Laboratory (	Code: Laboratory Co	ntrol Sample					
			Percent				
	Reportir	ng Spike	Recovery	Acceptance	e		
Analyte	Units	Level	LCS	Criteria			
Gasoline	ug/L (pp	b) 1,000	103	69-134			

#### ENVIRONMENTAL CHEMISTS

Date of Report: 02/18/19 Date Received: 02/08/19 Project: Kelly Moore 14697009, F&BI 902134

#### QUALITY ASSURANCE RESULTS FOR THE ANALYSIS OF WATER SAMPLES FOR TOTAL PETROLEUM HYDROCARBONS AS DIESEL EXTENDED USING METHOD NWTPH-Dx

Laboratory Code: 90	2134-04 (Matrix	Spike)					
·		-		Percent	Percent		
	Reporting	Spike	Sample	Recovery	Recovery	Acceptance	e RPD
Analyte	Units	Level	Result	MS	MSD	Criteria	(Limit 20)
Diesel Extended	ug/L (ppb)	3,000	19,000	140 b	0 b	64-141	200 b
Laboratory Code: La	aboratory Contro	l Sample					
			Percent	Percent			
	Reporting	Spike	Recovery	/ Recovery	/ Accept	ance F	RPD
Analyte	Units	Level	LCS	LCSD	Crite	ria (Lin	nit 20)
Diesel Extended	ug/L (ppb)	3,000	105	108	61-1	33	3

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#### ENVIRONMENTAL CHEMISTS

Date of Report: 02/18/19 Date Received: 02/08/19 Project: Kelly Moore 14697009, F&BI 902134

#### QUALITY ASSURANCE RESULTS FOR THE ANALYSIS OF WATER SAMPLES FOR PAHS BY EPA METHOD 8270D SIM

Laboratory Code: 902134-04 1/2 (Matrix Spike)

			Sample	Percent	Percent		
	Reporting	Spike	Result	Recovery	Recovery	Acceptance	RPD
Analyte	Units	Level	(Wet wt)	MS	MSD	Criteria	(Limit 20)
Naphthalene	ug/L (ppb)	1	< 0.4	84	80	10-172	5
Acenaphthylene	ug/L (ppb)	1	28	26 b	2 b	38-137	171 b
Acenaphthene	ug/L (ppb)	1	0.71	49 b	53 b	20-150	8 b
Fluorene	ug/L (ppb)	1	0.36	49 b	50 b	10-181	2 b
Phenanthrene	ug/L (ppb)	1	< 0.04	67	67	58-109	0
Anthracene	ug/L (ppb)	1	< 0.04	70	71	47-114	1
Fluoranthene	ug/L (ppb)	1	0.11	71	77	10-171	8
Pyrene	ug/L (ppb)	1	0.15	65	74	63-107	13
Benz(a)anthracene	ug/L (ppb)	1	< 0.04	53 vo	69	60-93	26 vo
Chrysene	ug/L (ppb)	1	< 0.04	48 vo	62	60-102	25 vo
Benzo(b)fluoranthene	ug/L (ppb)	1	< 0.04	45 vo	65	62-91	36 vo
Benzo(k)fluoranthene	ug/L (ppb)	1	< 0.04	38 vo	67	51-98	55 vo
Benzo(a)pyrene	ug/L (ppb)	1	< 0.04	42 vo	64	60-86	42 vo
Indeno(1,2,3-cd)pyrene	ug/L (ppb)	1	< 0.04	34	60	10-98	55 vo
Dibenz(a,h)anthracene	ug/L (ppb)	1	< 0.04	31	58	10-97	61 vo
Benzo(g,h,i)perylene	ug/L (ppb)	1	< 0.04	31	54	10-102	54 vo

Laboratory Code: Laboratory Control Sample

, i i i i i i i i i i i i i i i i i i i			Percent	Percent		
	Reporting	Spike	Recovery LCS	Recovery	Acceptance	RPD
Analyte	Units	Level	-	LCSD	Criteria	(Limit 20)
Naphthalene	ug/L (ppb)	1	88	92	67-116	4
Acenaphthylene	ug/L (ppb)	1	100	103	65-119	3
Acenaphthene	ug/L (ppb)	1	98	102	66-118	4
Fluorene	ug/L (ppb)	1	99	107	64-125	8
Phenanthrene	ug/L (ppb)	1	88	89	67-120	1
Anthracene	ug/L (ppb)	1	91	96	65-122	5
Fluoranthene	ug/L (ppb)	1	88	97	65-127	10
Pyrene	ug/L (ppb)	1	96	92	62-130	4
Benz(a)anthracene	ug/L (ppb)	1	96	100	60-118	4
Chrysene	ug/L (ppb)	1	91	94	66-125	3
Benzo(b)fluoranthene	ug/L (ppb)	1	103	101	55-135	2
Benzo(k)fluoranthene	ug/L (ppb)	1	97	105	62-125	8
Benzo(a)pyrene	ug/L (ppb)	1	96	100	58-127	4
Indeno(1,2,3-cd)pyrene	ug/L (ppb)	1	104	105	36-142	1
Dibenz(a,h)anthracene	ug/L (ppb)	1	87	93	37-133	7
Benzo(g,h,i)perylene	ug/L (ppb)	1	90	97	34-135	7

#### ENVIRONMENTAL CHEMISTS

Date of Report: 02/18/19 Date Received: 02/08/19 Project: Kelly Moore 14697009, F&BI 902134

#### QUALITY ASSURANCE RESULTS FOR THE ANALYSIS OF WATER SAMPLES FOR TOTAL METALS USING EPA METHOD 6020B

Laboratory Code: 902134-04 (Matrix Spike)

Laboratory couct	Reporting	Sniko	Sample	Percent	Percent	Accentance	RPD
Analyte	Units	Level	Result	MS	MSD	Criteria	(Limit 20)
Arsenic	ug/L (ppb)	10	3.04	109	110	75-125	1
Chromium	ug/L (ppb)	20	2.00	88	89	75-125	1
Copper	ug/L (ppb)	20	19.4	86	89	75-125	3
Lead	ug/L (ppb)	10	2.04	101	102	75-125	1
Mercury	ug/L (ppb)	5	<1	114	117	75-125	3
Nickel	ug/L (ppb)	20	7.79	84	86	75-125	2
Zinc	ug/L (ppb)	50	135	96	101	75-125	5

Laboratory Code: Laboratory Control Sample

		Percent	
Reporting	Spike	Recovery	Acceptance
Units	Level	LCS	Criteria
ug/L (ppb)	10	100	80-120
ug/L (ppb)	20	101	80-120
ug/L (ppb)	20	103	80-120
ug/L (ppb)	10	102	80-120
ug/L (ppb)	5	116	80-120
ug/L (ppb)	20	102	80-120
ug/L (ppb)	50	105	80-120
	Reporting Units ug/L (ppb) ug/L (ppb) ug/L (ppb) ug/L (ppb) ug/L (ppb) ug/L (ppb) ug/L (ppb)	Reporting Units Spike Level   ug/L (ppb) 10   ug/L (ppb) 20   ug/L (ppb) 20   ug/L (ppb) 10   ug/L (ppb) 5   ug/L (ppb) 20   ug/L (ppb) 5   ug/L (ppb) 20   ug/L (ppb) 5   ug/L (ppb) 50	Reporting Spike Recovery   Units Level LCS   ug/L (ppb) 10 100   ug/L (ppb) 20 101   ug/L (ppb) 20 103   ug/L (ppb) 10 102   ug/L (ppb) 5 116   ug/L (ppb) 20 102   ug/L (ppb) 50 105

#### ENVIRONMENTAL CHEMISTS

Date of Report: 02/18/19 Date Received: 02/08/19 Project: Kelly Moore 14697009, F&BI 902134

#### QUALITY ASSURANCE RESULTS FOR THE ANALYSIS OF WATER SAMPLES FOR VOLATILES BY EPA METHOD 8260C

Laboratory Code: 902134-04 (Matrix Spike)

U C	•			Percent	Percent		
	Reporting	Spike	Sample	Recovery	Recovery	Acceptance	RPD
Analyte	Units	Level	Result	MS	MSD	Criteria	(Limit 20)
Dichlorodifluoromethane	ug/L (ppb)	50	<1	102	102	10-172	0
Chloromethane	ug/L (ppb)	50	<10	98	99	25-166	1
Vinyl chloride Bromomethane	ug/L (ppb)	50 50	<0.2	101	101	36-166	0
Chloroethane	ug/L (ppb)	50	<1	90	95 97	47-169	1
Trichlorofluoromethane	ug/L (ppb)	50	<1	105	101	44-165	4
Acetone	ug/L (ppb)	250	<50	83	83	10-182	0
1,1-Dichloroethene	ug/L (ppb)	50	<1	102	99	60-136	3
Hexane	ug/L (ppb)	50	<1	99	93	52-150	6
Metnylene chloride Methyl t hutyl other (MTRE)	ug/L (ppb)	50 50	<5	106	109	67-132 74 197	3
trans-1 2-Dichloroethene	ug/L (ppb)	50	<1	97	98	72-129	1
1.1-Dichloroethane	ug/L (ppb)	50	<1	100	94	70-128	6
2,2-Dichloropropane	ug/L (ppb)	50	<1	100	95	36-154	5
cis-1,2-Dichloroethene	ug/L (ppb)	50	<1	97	93	71-127	4
Chloroform	ug/L (ppb)	50	<1	96	89	65-132	8
2-Butanone (MEK) 1.2 Disblaresthans (EDC)	ug/L (ppb)	250	<10	89	84	10-129	6
1.1.1-Trichloroethane	ug/L (ppb)	50	<1	98	99 89	60-146	3 10
1.1-Dichloropropene	ug/L (ppb)	50	<1	98	95	69-133	3
Carbon tetrachloride	ug/L (ppb)	50	<1	100	96	56-152	4
Benzene	ug/L (ppb)	50	< 0.35	98	95	76-125	3
Trichloroethene	ug/L (ppb)	50	<1	92	94	66-135	2
1,2-Dichloropropane	ug/L (ppb)	50	<1	88	95	78-125	8
Dibromomethane	ug/L (ppb)	50 50	<1	91 87	99 94	66-141	8
4-Methyl-2-pentanone	ug/L (ppb)	250	<10	90	99	10-185	10
cis-1,3-Dichloropropene	ug/L (ppb)	50	<1	90	96	72-132	6
Toluene	ug/L (ppb)	50	<1	91	98	76-122	7
trans-1,3-Dichloropropene	ug/L (ppb)	50	<1	88	101	76-130	14
1,1,2-Trichloroethane	ug/L (ppb)	50	<1	97	101	68-131	4
2-Hexanone 1.3 Dichleropropane	ug/L (ppb)	250	<10	102	102	10-185	0
Tetrachloroethene	ug/L (ppb)	50	<1	100	90	10-226	1
Dibromochloromethane	ug/L (ppb)	50	<1	107	106	70-139	1
1,2-Dibromoethane (EDB)	ug/L (ppb)	50	<1	101	101	69-134	0
Chlorobenzene	ug/L (ppb)	50	<1	95	95	77-122	0
Ethylbenzene	ug/L (ppb)	50	<1	99	95	69-135	4
1,1,1,2-1 etrachloroethane	ug/L (ppb)	50	<1	98	101	73-137	3
n,p-Aylene	ug/L (ppb)	50	<2	90	90	60-140	0
Styrene	ug/L (ppb)	50	<1	98	02 97	71-133	1
Isopropylbenzene	ug/L (ppb)	50	8.3	93	93	65-142	0
Bromoform	ug/L (ppb)	50	<1	107	106	65-142	1
n-Propylbenzene	ug/L (ppb)	50	12	108 b	103 b	58-144	5 b
Bromobenzene 1.3.5 Trimothylbonzono	ug/L (ppb)	50 50	<1	107	105	75-124	2
1,3,3-Trimethyldenzene 1,1,2,2-Tetrachloroethane	ug/L (ppb)	50	<1	103	98 119	51-154	1
1.2.3-Trichloropropane	ug/L (ppb)	50	<1	108	104	53-150	4
2-Chlorotoluene	ug/L (ppb)	50	<1	106	99	66-127	7
4-Chlorotoluene	ug/L (ppb)	50	<1	105	99	65-130	6
tert-Butylbenzene	ug/L (ppb)	50	<1	98	95	65-137	3
1,2,4-Trimethylbenzene	ug/L (ppb)	50	<1	100	95	59-146	5
sec-Bulyidenzene	ug/L (ppb)	50 50	1.2	98	97	64-140 65-141	2
1.3-Dichlorobenzene	ug/L (ppb)	50	<1	97	98	72-123	1
1,4-Dichlorobenzene	ug/L (ppb)	50	<1	95	95	69-126	Ō
1,2-Dichlorobenzene	ug/L (ppb)	50	<1	96	94	69-128	2
1,2-Dibromo-3-chloropropane	ug/L (ppb)	50	<10	116	124	32-164	7
1,2,4-Trichlorobenzene	ug/L (ppb)	50	<1	106	105	66-136	1
Hexachiorobutadiene	ug/L (ppb)	50 50	<1	104	103	60-143	1
1.2.3-Trichlorobenzene	ug/L (ppb)	50	<1	108	108	69-148	0
, ,	-0 - (FF-)						-

#### ENVIRONMENTAL CHEMISTS

Date of Report: 02/18/19 Date Received: 02/08/19 Project: Kelly Moore 14697009, F&BI 902134

#### QUALITY ASSURANCE RESULTS FOR THE ANALYSIS OF WATER SAMPLES FOR VOLATILES BY EPA METHOD 8260C

Laboratory Code: Laboratory Control Sample

c c	•		Percent	
	Reporting	Spike	Recovery	Acceptance
Analyte	Units	Level	LCS	Criteria
Dichlorodifluoromethane	ug/L (ppb)	50	121	25-158
Chloromethane Vincel ablantide	ug/L (ppb)	50	112	45-156
Bromomethane	ug/L (ppb)	50 50	110	55-143
Chloroethane	ug/L (ppb)	50	116	58-146
Trichlorofluoromethane	ug/L (ppb)	250	119	50-150
Acetone	ug/L (ppb)	250	93	53-131
1,1-Dichloroethene	ug/L (ppb)	50	116	67-136
Hexane Methodana ablanida	ug/L (ppb)	50	116	57-137
Methylene chloride Methyl t-butyl ether (MTBE)	ug/L (ppb)	50 50	128	59-148 64-147
trans-1 2-Dichloroethene	ug/L (ppb)	50	122	68-128
1.1-Dichloroethane	ug/L (ppb)	50	122 vo	79-121
2,2-Dichloropropane	ug/L (ppb)	50	129	55-143
cis-1,2-Dichloroethene	ug/L (ppb)	50	119	80-123
Chloroform	ug/L (ppb)	50	114	80-121
2-Butanone (MEK)	ug/L (ppb)	250	97	57-149
1,2-Dichloroethane (EDC)	ug/L (ppb)	50	101	73-132
1 1-Dichloropropene	ug/L (ppb)	50	106	77-129
Carbon tetrachloride	ug/L (ppb)	50	110	75-158
Benzene	ug/L (ppb)	50	101	69-134
Trichloroethene	ug/L (ppb)	50	96	80-120
1,2-Dichloropropane	ug/L (ppb)	50	98	77-123
Bromodichloromethane	ug/L (ppb)	50	96	81-133
A Mothyl 2 pontanono	ug/L (ppb)	50 250	93	82-125
cis-1 3-Dichloropropene	ug/L (ppb)	50	93	82-132
Toluene	ug/L (ppb)	50	100	72-122
trans-1,3-Dichloropropene	ug/L (ppb)	50	88	80-136
1,1,2-Trichloroethane	ug/L (ppb)	50	96	75-124
2-Hexanone	ug/L (ppb)	250	79	60-136
1,3-Dichloropropane	ug/L (ppb)	50	87	76-126
1 etrachioroethene Dibromochloromothano	ug/L (ppb)	50 50	103	76-121 84 122
1.2-Dibromoethane (FDB)	ug/L (ppb)	50	88	82-125
Chlorobenzene	ug/L (ppb)	50	94	83-114
Ethylbenzene	ug/L (ppb)	50	103	77-124
1,1,1,2-Tetrachloroethane	ug/L (ppb)	50	114	84-127
m,p-Xylene	ug/L (ppb)	100	101	83-125
o-Xylene	ug/L (ppb)	50	105	81-121
Styrene	ug/L (ppb)	50	100	84-119
Bromoform	ug/L (ppb)	50	104	74-136
n-Propylbenzene	ug/L (ppb)	50	99	74-126
Bromobenzene	ug/L (ppb)	50	96	80-121
1,3,5-Trimethylbenzene	ug/L (ppb)	50	103	78-123
1,1,2,2-Tetrachloroethane	ug/L (ppb)	50	98	66-126
1,2,3-Trichloropropane	ug/L (ppb)	50	92	67-124
2-Chlorotoluene	ug/L (ppb)	50	101	77-127
4-Chiofololuene tert-Butylbenzene	ug/L (ppb)	50	95 104	70-120 80-123
1.2.4-Trimethylbenzene	ug/L (ppb)	50	104	79-122
sec-Butylbenzene	ug/L (ppb)	50	109	80-125
p-Isopropyltoluene	ug/L (ppb)	50	106	81-123
1,3-Dichlorobenzene	ug/L (ppb)	50	98	85-116
1,4-Dichlorobenzene	ug/L (ppb)	50	95	84-121
1,2-Dicnioropenzene	ug/L (ppb)	50	98	85-116
1,2-Didiono-3-Chloropropane 1 2 4-Trichlorobenzene	ug/L (ppb)	50 50	110	07-141 79-130
Hexachlorobutadiene	ug/L (ppb)	50	110	53-141
Naphthalene	ug/L (ppb)	50	112	64-133
1,2,3-Trichlorobenzene	ug/L (ppb)	50	111	65-136

#### ENVIRONMENTAL CHEMISTS

### **Data Qualifiers & Definitions**

a - The analyte was detected at a level less than five times the reporting limit. The RPD results may not provide reliable information on the variability of the analysis.

 ${\bf b}$  - The analyte was spiked at a level that was less than five times that present in the sample. Matrix spike recoveries may not be meaningful.

ca - The calibration results for the analyte were outside of acceptance criteria. The value reported is an estimate.

c - The presence of the analyte may be due to carryover from previous sample injections.

cf - The sample was centrifuged prior to analysis.

d - The sample was diluted. Detection limits were raised and surrogate recoveries may not be meaningful.

dv - Insufficient sample volume was available to achieve normal reporting limits.

f - The sample was laboratory filtered prior to analysis.

fb - The analyte was detected in the method blank.

fc - The analyte is a common laboratory and field contaminant.

 $hr\ \text{-}\ The\ sample\ and\ duplicate\ were\ reextracted\ and\ reanalyzed.\ RPD\ results\ were\ still\ outside\ of\ control\ limits.\ Variability\ is\ attributed\ to\ sample\ inhomogeneity.$ 

hs - Headspace was present in the container used for analysis.

ht – The analysis was performed outside the method or client-specified holding time requirement.

ip - Recovery fell outside of control limits due to sample matrix effects.

j - The analyte concentration is reported below the lowest calibration standard. The value reported is an estimate.

 ${\rm J}$  - The internal standard associated with the analyte is out of control limits. The reported concentration is an estimate.

jl - The laboratory control sample(s) percent recovery and/or RPD were out of control limits. The reported concentration should be considered an estimate.

js - The surrogate associated with the analyte is out of control limits. The reported concentration should be considered an estimate.

lc - The presence of the analyte is likely due to laboratory contamination.

L - The reported concentration was generated from a library search.

nm - The analyte was not detected in one or more of the duplicate analyses. Therefore, calculation of the RPD is not applicable.

pc - The sample was received with incorrect preservation or in a container not approved by the method. The value reported should be considered an estimate.

ve - The analyte response exceeded the valid instrument calibration range. The value reported is an estimate.

vo - The value reported fell outside the control limits established for this analyte.

x - The sample chromatographic pattern does not resemble the fuel standard used for quantitation.

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Sample ID	Lab ID	Date Sampled	Time Sampled	Sampl Type	e #of Jars	TPH-HCID	TPH-Diesel	TPH-Casoline	BTEX by 8021B	VOCs by 8260C	SVOCs by 8270D	PAHs 8270D SIN	Total Metals			No	otes
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KMU-08-020819	05 .	2/8/19	1040		9									64	MAG		
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#### ENVIRONMENTAL CHEMISTS

James E. Bruya, Ph.D. Yelena Aravkina, M.S. Michael Erdahl, B.S. Arina Podnozova, B.S. Eric Young, B.S. 3012 16th Avenue West Seattle, WA 98119-2029 (206) 285-8282 fbi@isomedia.com www.friedmanandbruya.com

August 27, 2018

Crystal Thimsen, Project Manager Wood Environment & Infrastructure Solutions, Inc. One Union Square 600 University Street, Suite 600 Seattle, WA 98101

Dear Ms Thimsen:

Included are the results from the testing of material submitted on August 17, 2018 from the Kelly-Moore, F&BI 808402 project. There are 46 pages included in this report. Any samples that may remain are currently scheduled for disposal in 30 days. If you would like us to return your samples or arrange for long term storage at our offices, please contact us as soon as possible.

We appreciate this opportunity to be of service to you and hope you will call if you should have any questions.

Sincerely,

FRIEDMAN & BRUYA, INC.

Gale

Michael Erdahl Project Manager

Enclosures WEI0827R.DOC

#### ENVIRONMENTAL CHEMISTS

#### CASE NARRATIVE

This case narrative encompasses samples received on August 17, 2018 by Friedman & Bruya, Inc. from the Wood Environment & Infrastructure Solutions Kelly-Moore, F&BI 808402 project. Samples were logged in under the laboratory ID's listed below.

Laboratory ID	Wood Environment & Infrastructure Solutions
808402 -01	KMW-02R-081618
808402 -02	KMW-03R-081618
808402 -03	KMW-04-081618
808402 -04	KMW-06-081618
808402 -05	KMW-07-081618
808402 -06	KMW-08-081618
808402 -07	KMW-09-081618
808402 -08	KMW-10-081618
808402 -09	KMW-03R-9-081618

A 6020A internal standard failed the acceptance criteria for sample KMW-04-081618, KMW-06-081618, KMW-09-081618, and KMW-10-081618 due to matrix interferences. The data were flagged accordingly. The sample was diluted and reanalyzed.

The 8260C calibration standard failed the acceptance criteria for 2-butanone and 2hexanone. The data were flagged accordingly.

Phenanthrene was detected in the 8270D SIM method blank at a level within 10 times the concentration detected in several samples. The data were flagged accordingly.

All other quality control requirements were acceptable.

#### ENVIRONMENTAL CHEMISTS

Date of Report: 08/27/18 Date Received: 08/17/18 Project: Kelly-Moore, F&BI 808402 Date Extracted: 08/20/18 Date Analyzed: 08/20/18 and 08/21/18

#### RESULTS FROM THE ANALYSIS OF WATER SAMPLES FOR TOTAL PETROLEUM HYDROCARBONS AS GASOLINE USING METHOD NWTPH-Gx

Results Reported as ug/L (ppb)

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<u>Sample ID</u> Laboratory ID	Gasoline Range	Surrogate ( <u>% Recovery)</u> (Limit 51-134)
KMW-02R-081618 808402-01	<100	101
KMW-03R-081618 808402-02	290	114
KMW-04-081618 808402-03 1/10	33,000	123
KMW-06-081618 808402-04	4,000	ip
KMW-07-081618 808402-05	<100	96
KMW-08-081618 808402-06	230	109
KMW-09-081618 808402-07	940	114
KMW-10-081618 808402-08 1/10	4,800	108
KMW-03R-9-081618 808402-09	380	126
Method Blank 08-1759 MB	<100	96

#### ENVIRONMENTAL CHEMISTS

Date of Report: 08/27/18 Date Received: 08/17/18 Project: Kelly-Moore, F&BI 808402 Date Extracted: 08/17/18 Date Analyzed: 08/17/18

#### RESULTS FROM THE ANALYSIS OF WATER SAMPLES FOR TOTAL PETROLEUM HYDROCARBONS AS DIESEL AND MOTOR OIL USING METHOD NWTPH-Dx

Results Reported as ug/L (ppb)

<u>Sample ID</u> Laboratory ID	Diesel Range (C10-C25)	<u>Motor Oil Range</u> (C <sub>25</sub> -C <sub>36</sub> )	Surrogate <u>(% Recovery)</u> (Limit 41-152)
KMW-02R-081618 808402-01	<50	<250	70
KMW-03R-081618 808402-02	350 x	<250	73
KMW-04-081618 808402-03	2,000 x	<250	69
KMW-06-081618 808402-04	8,600	680 x	71
KMW-07-081618 808402-05	<50	<250	62
KMW-08-081618 808402-06	160 x	<250	62
KMW-09-081618 808402-07	3,600	360 x	67
KMW-10-081618 808402-08	1,400 x	<250	73
KMW-03R-9-081618 808402-09	450 x	<250	70
Method Blank	<50	<250	55

08-1856 MB

## ENVIRONMENTAL CHEMISTS

Client ID: Date Received: Date Extracted: Date Analyzed: Matrix:	KMW-02R-081618 08/17/18 08/20/18 08/20/18 Water	Client: Project: Lab ID: Data File: Instrument:	Wood Environment & Infrastructure Solutions Kelly-Moore, F&BI 808402 808402-01 808402-01.049 ICPMS2
Units:	ug/L (ppb)	Operator:	SP
Analyte:	Concentration ug/L (ppb)		
Arsenic	<1		
Chromium	<1		
Copper	<5		
Lead	<1		
Mercury	<1		
Nickel	1.82		
Zinc	<5		

## ENVIRONMENTAL CHEMISTS

Client ID:	KMW-03R-081618	Client:	Wood Environment & Infrastructure Solutions
Date Received:	08/17/18	Project:	Kelly-Moore, F&BI 808402
Date Extracted:	08/20/18	Lab ID:	808402-02
Date Analyzed:	08/20/18	Data File:	808402-02.050
Matrix:	Water	Instrument:	ICPMS2
Units:	ug/L (ppb)	Operator:	SP
	Concentration		
Analyte:	ug/L (ppb)		
Arsenic	<1		
Chromium	<1		
Copper	<5		
Lead	<1		
Mercury	<1		
Nickel	1.46		
Zinc	<5		

## ENVIRONMENTAL CHEMISTS

Client ID: Date Received: Date Extracted:	KMW-04-081618 08/17/18 08/20/18	Client: Project: Lab ID:	Wood Environment & Infrastructure Solutions Kelly-Moore, F&BI 808402 808402-03
Date Analyzed:	08/20/18 Water	Data File:	000402-03.001
Matrix:	water	instrument:	ICPMISZ
Units:	ug/L (ppb)	Operator:	SP
Analyte:	Concentration ug/L (ppb)		
Arsenic	14.5		
Chromium	<1 J		
Copper	<5 J		
Lead	<1		
Mercury	<1		
Nickel	2.43 J		
Zinc	<5 J		
## ENVIRONMENTAL CHEMISTS

Client ID: Date Received: Date Extracted: Date Analyzed:	KMW-04-081618 08/17/18 08/20/18 08/20/18 Watan	Client: Project: Lab ID: Data File:	Wood Environment & Infrastructure Solutions Kelly-Moore, F&BI 808402 808402-03 x10 808402-03 x10.095 ICDMS2
Unite:	water	Operator:	
Units:	ug/L (ppb)	Operator:	SP
Analyte:	Concentration ug/L (ppb)		
Arsenic	14.5		
Chromium	<10		
Copper	<50		
Lead	<10		
Mercury	<10		
Nickel	<10		
Zinc	<50		

## ENVIRONMENTAL CHEMISTS

Client ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-06-081618 08/17/18 08/20/18 08/20/18 Water ug/L (ppb)	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Infrastructure Solutions Kelly-Moore, F&BI 808402 808402-04 808402-04.052 ICPMS2 SP
Analyte:	Concentration ug/L (ppb)		
Arsenic	4.83		
Chromium	2.48 J		
Copper	11.9 J		
Lead	5.61		
Mercury	<1		
Nickel	1.38 J		
Zinc	5.72 J		

## ENVIRONMENTAL CHEMISTS

Client ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-06-081618 08/17/18 08/20/18 08/20/18 Water ug/L (ppb)	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Infrastructure Solutions Kelly-Moore, F&BI 808402 808402-04 x10 808402-04 x10.107 ICPMS2 SP
Analyte:	Concentration ug/L (ppb)		
Arsenic	<10		
Chromium	<10		
Copper	<50		
Lead	<10		
Mercury	<10		
Nickel	<10		
Zinc	<50		

## ENVIRONMENTAL CHEMISTS

Client ID: Date Received: Date Extracted: Date Analyzed:	KMW-07-081618 08/17/18 08/20/18 08/20/18	Client: Project: Lab ID: Data File:	Wood Environment & Infrastructure Solutions Kelly-Moore, F&BI 808402 808402-05 808402-05.054
Matrix:	Water	Instrument:	ICPMS2
Units:	ug/L (ppb)	Operator:	SP
Analyte:	Concentration ug/L (ppb)		
Arsenic	<1		
Chromium	<1		
Copper	<5		
Lead	<1		
Mercury	<1		
Nickel	<1		
Zinc	<5		

## ENVIRONMENTAL CHEMISTS

Client ID: Date Received: Date Extracted: Date Analyzed: Matrix:	KMW-08-081618 08/17/18 08/20/18 08/20/18 Water	Client: Project: Lab ID: Data File: Instrument:	Wood Environment & Infrastructure Solutions Kelly-Moore, F&BI 808402 808402-06 808402-06.086 ICPMS2
Units:	ug/L (ppb)	Operator:	SP
Analyte:	Concentration ug/L (ppb)		
Arsenic	<1		
Chromium	<1		
Copper	<5		
Lead	<1		
Mercury	<1		
Nickel	1.11		
Zinc	<5		

## ENVIRONMENTAL CHEMISTS

Client ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-09-081618 08/17/18 08/20/18 08/20/18 Water ug/L (ppb)	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Infrastructure Solutions Kelly-Moore, F&BI 808402 808402-07 808402-07.087 ICPMS2 SP
Analyte:	Concentration ug/L (ppb)		
Arsenic	2.04		
Chromium	1.40 J		
Copper	5.62 J		
Lead	3.10		
Mercury	<1		
Nickel	1.25 J		
Zinc	<5 J		

## ENVIRONMENTAL CHEMISTS

Client ID: Date Received: Date Extracted: Date Analyzed: Matrix:	KMW-09-081618 08/17/18 08/20/18 08/21/18 Water	Client: Project: Lab ID: Data File: Instrument:	Wood Environment & Infrastructure Solutions Kelly-Moore, F&BI 808402 808402-07 x10 808402-07 x10.030 ICPMS2
Units:	ug/L (ppb)	Operator:	SP
Analyte:	Concentration ug/L (ppb)		
Arsenic	<10		
Chromium	<10		
Copper	<50		
Lead	<10		
Mercury	<10		
Nickel	<10		
Zinc	<50		

## ENVIRONMENTAL CHEMISTS

Client ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-10-081618 08/17/18 08/20/18 08/20/18 Water ug/L (ppb)	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Infrastructure Solutions Kelly-Moore, F&BI 808402 808402-08 808402-08.088 ICPMS2 SP
Analyte:	Concentration ug/L (ppb)		
Arsenic	4.61		
Chromium	1.35 J		
Copper	<5 J		
Lead	<1		
Mercury	<1		
Nickel	<1 J		
Zinc	<5 J		

## ENVIRONMENTAL CHEMISTS

Client ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-10-081618 08/17/18 08/20/18 08/22/18 Water ug/L (ppb)	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Infrastructure Solutions Kelly-Moore, F&BI 808402 808402-08 x10 808402-08 x10.043 ICPMS2 SP
Analyte:	Concentration ug/L (ppb)		
Arsenic	<10		
Chromium	<10		
Copper	<50		
Lead	<10		
Mercury	<10		
Nickel	<10		
Zinc	<50		

## ENVIRONMENTAL CHEMISTS

Client ID: Date Received: Date Extracted: Date Analyzed: Matrix:	KMW-03R-9-081618 08/17/18 08/20/18 08/20/18 Water	Client: Project: Lab ID: Data File: Instrument:	Wood Environment & Infrastructure Solutions Kelly-Moore, F&BI 808402 808402-09 808402-09.089 ICPMS2
Units:	ug/L (ppb)	Operator:	SP
Analyte:	Concentration ug/L (ppb)		
Arsenic	<1		
Chromium	<1		
Copper	<5		
Lead	<1		
Mercury	<1		
Nickel	1.82		
Zinc	<5		

## ENVIRONMENTAL CHEMISTS

Client ID:	Method Blank	Client:	Wood Environment & Infrastructure Solutions
Date Received:	NA	Project:	Kelly-Moore, F&BI 808402
Date Extracted:	08/20/18	Lab ID:	I8-535 mb
Date Analyzed:	08/20/18	Data File:	I8-535 mb.053
Matrix:	Water	Instrument:	ICPMS2
Units:	ug/L (ppb)	Operator:	SP
	Concentration		
Analyte:	ug/L (ppb)		
Arsenic	<1		
Chromium	<1		
Copper	<5		
Lead	<1		
Mercury	<1		
Nickel	<1		
Zinc	<5		

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-02R-0 08/17/18 08/17/18 08/17/18 Water ug/L (ppb)	)81618	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Kelly-Moore, F&BI 8 808402-01 081730.D GCMS4 JS	& Infrastructure Solutions 08402
			Lower	Upper	
Surrogates:		% Recovery:	Limit:	Limit:	
1,2-Dichloroethane-o	14	101	57	121	
Toluene-d8		97	63	127	
4-Bromofluorobenzer	ne	95	60	133	
		Concentration			Concentration
Compounds:		ug/L (ppb)	Compour	nds:	ug/L (ppb)
Dichlorodifluoromet	hane	<1	1.3-Dichl	oropropane	<1
Chloromethane		<10	Tetrachl	proethene	<1
Vinvl chloride		< 0.2	Dibromo	chloromethane	<1
Bromomethane		<1	1.2-Dibro	omoethane (EDB)	<1
Chloroethane		<1	Chlorobe	nzene	<1
Trichlorofluorometh	ane	<1	Ethvlben	zene	<1
Acetone		<50	1,1,1,2-T	etrachloroethane	<1
1,1-Dichloroethene		<1	m,p-Xyle	ene	<2
Hexane		<1	o-Xylene		<1
Methylene chloride		<5	Styrene		<1
Methyl t-butyl ether (MTBE)		<1	Isopropylbenzene		<1
trans-1,2-Dichloroethene		<1	Bromoform		<1
1.1-Dichloroethane		<1	n-Propylbenzene		<1
2,2-Dichloropropane		<1	Bromobenzene		<1
cis-1,2-Dichloroether	ne	<1	1,3,5-Trimethylbenzene		<1
Chloroform		<1	1,1,2,2-Tetrachloroethane		<1
2-Butanone (MEK)		<10 ca	1,2,3-Tri	chloropropane	<1
1,2-Dichloroethane (	EDC)	<1	2-Chloro	toluene	<1
1,1,1-Trichloroethan	ie	<1	4-Chloro	toluene	<1
1,1-Dichloropropene		<1	tert-Buty	lbenzene	<1
Carbon tetrachloride	)	<1	1,2,4-Tri	methylbenzene	<1
Benzene		< 0.35	sec-Buty	lbenzene	<1
Trichloroethene		<1	p-Isoprop	yltoluene	<1
1,2-Dichloropropane		<1	1,3-Dichl	orobenzene	<1
Bromodichlorometha	ane	<1	1,4-Dichl	orobenzene	<1
Dibromomethane		<1	1,2-Dichl	orobenzene	<1
4-Methyl-2-pentanor	ne	<10	1,2-Dibro	omo-3-chloropropane	<10
cis-1,3-Dichloroprop	ene	<1	1,2,4-Trichlorobenzene		<1
Toluene		<1	Hexachle	orobutadiene	<1
trans-1,3-Dichloropr	opene	<1	Naphtha	lene	<1
1,1,2-Trichloroethan	ie	<1	1,2,3-Tri	chlorobenzene	<1
2-Hexanone		<10 ca			

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-03R-0 08/17/18 08/17/18 08/17/18 Water ug/L (ppb)	)81618	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Kelly-Moore, F&BI 8 808402-02 081731.D GCMS4 JS	& Infrastructure Solutions 08402
			Lower	Upper	
Surrogates:		% Recovery:	Limit:	Limit:	
1,2-Dichloroethane-c	14	100	57	121	
Toluene-d8		98	63	127	
4-Bromofluorobenzer	ne	97	60	133	
		Concentration			Concentration
Compounds:		ug/L (ppb)	Compour	nds:	ug/L (ppb)
Dichlorodifluoromet	hane	<1	1 3-Dichl	oropropape	<1
Chloromethane	lianc	<10	Tetrachl	oroethene	<1
Vinvl chloride		<0.2	Dibromo	chloromethane	<1
Bromomethane		<0.2	1 2-Dibro	omoethane (FDB)	<1
Chloroethane		<1	Chlorobe	nzene	<1
Trichlorofluorometh	ane	<1	Ethylber	izene	<1
Acetone	une	<50	1 1 1 2-T	etrachloroethane	<1
1 1-Dichloroethene		<1	m p-Xvle	ne	<2
Hexane		<1	o-Xvlene		<1
Methylene chloride		<5	Styrene		<1
Methyl t-butyl ether	(MTBE)	<1	Isopropy	lbenzene	3.6
trans-1.2-Dichloroet	hene	<1	Bromofor	rm	<1
1.1-Dichloroethane		<1	n-Propyl	benzene	7.4
2.2-Dichloropropane		<1	Bromobe	nzene	<1
cis-1.2-Dichloroether	ne	<1	1.3.5-Tri	methylbenzene	<1
Chloroform		<1	1,0,0 11 1.1.2.2-T	etrachloroethane	<1
2-Butanone (MEK)		<10 ca	1.2.3-Tri	chloropropane	<1
1.2-Dichloroethane (	EDC)	<1	2-Chloro	toluene	<1
1.1.1-Trichloroethan	)e	<1	4-Chloro	toluene	<1
1.1-Dichloropropene		<1	tert-Buty	lbenzene	<1
Carbon tetrachloride	<u>)</u>	<1	1.2.4-Tri	methylbenzene	<1
Benzene		< 0.35	sec-Buty	lbenzene	1.8
Trichloroethene		<1	p-Isoprop	ovltoluene	<1
1,2-Dichloropropane		<1	1,3-Dichl	orobenzene	<1
Bromodichlorometha	ane	<1	1,4-Dichl	orobenzene	<1
Dibromomethane		<1	1,2-Dichl	orobenzene	<1
4-Methyl-2-pentanor	ne	<10	1,2-Dibro	omo-3-chloropropane	<10
cis-1,3-Dichloroprope	ene	<1	1,2,4-Tri	chlorobenzene	<1
Toluene		<1	Hexachlo	orobutadiene	<1
trans-1,3-Dichloropr	opene	<1	Naphtha	lene	<1
1,1,2-Trichloroethan	ie	<1	1,2,3-Tri	chlorobenzene	<1
2-Hexanone		<10 ca			

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-04-08 08/17/18 08/17/18 08/17/18 Water ug/L (ppb)	1618	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Kelly-Moore, F&BI 8 808402-03 081732.D GCMS4 JS	& Infrastructure Solutions 08402
			Lower	Upper	
Surrogates:		% Recovery:	Limit:	Limit:	
1,2-Dichloroethane-c	14	101	57	121	
Toluene-d8		95	63	127	
4-Bromofluorobenzer	ne	84	60	133	
		Concentration			Concentration
Compounds:		ug/L (ppb)	Compour	nds:	ug/L (ppb)
Dichlorodifluoromet	nane	<1	1.3-Dichl	oropropane	<1
Chloromethane	lane	<10	Tetrachl	oroethene	<1
Vinvl chloride		<02	Dibromo	chloromethane	<1
Bromomethane		<1	1 2-Dibro	omoethane (FDB)	<1
Chloroethane		<1	Chlorobe	nzene	<1
Trichlorofluorometh	ane	<1	Ethylber	izene	860 ve
Acetone	une	<50	1 1 1 2-T	etrachloroethane	<1
1 1-Dichloroethene		<1	m p-Xvle	ne	2.800 ve
Hexane		43	o-Xvlene		1 100 ve
Methylene chloride		<5	Styrene		<1
Methyl t-butyl ether	(MTBE)	<1	Isopropy	lbenzene	21
trans-1.2-Dichloroet	hene	<1	Bromofor	rm	<1
1.1-Dichloroethane		<1	n-Propyl	benzene	19
2.2-Dichloropropane		<1	Bromobe	nzene	<1
cis-1.2-Dichloroether	ne	<1	1.3.5-Tri	methylbenzene	17
Chloroform		<1	1,0,0 11 1.1.2.2-T	etrachloroethane	<1
2-Butanone (MEK)		<10 ca	1.2.3-Tri	chloropropane	<1
1.2-Dichloroethane (	EDC)	<1	2-Chloro	toluene	<1
1.1.1-Trichloroethan	)e	<1	4-Chloro	toluene	<1
1.1-Dichloropropene		<1	tert-Buty	lbenzene	<1
Carbon tetrachloride	<u>)</u>	<1	1.2.4-Tri	methylbenzene	54
Benzene		< 0.35	sec-Buty	lbenzene	1.4
Trichloroethene		<1	p-Isoprop	oyltoluene	<1
1,2-Dichloropropane		<1	1,3-Dichl	orobenzene	<1
Bromodichlorometha	ane	<1	1,4-Dichl	orobenzene	<1
Dibromomethane		<1	1,2-Dichl	orobenzene	<1
4-Methyl-2-pentanor	ne	<10	1,2-Dibro	omo-3-chloropropane	<10
cis-1,3-Dichloroprop	ene	<1	1,2,4-Tri	chlorobenzene	<1
Toluene		600 ve	Hexachlo	orobutadiene	<1
trans-1,3-Dichloropr	opene	<1	Naphtha	lene	5.1
1,1,2-Trichloroethan	ie	<1	1,2,3-Tri	chlorobenzene	<1
2-Hexanone		<10 ca			

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-04-08 08/17/18 08/17/18 08/20/18 Water ug/L (ppb)	1618	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Kelly-Moore, F&BI 8 808402-03 1/100 082017.D GCMS4 JS	& Infrastructure Solutions 08402
			Lower	Upper	
Surrogates:		% Recovery:	Limit:	Limit:	
1,2-Dichloroethane-c	14	99	57	121	
Toluene-d8		100	63	127	
4-Bromofluorobenzer	ne	97	60	133	
		Concentration			Concentration
Compounds:		ug/L (ppb)	Compour	nds:	ug/L (ppb)
Dichlorodifluoromet	hane	<100	1,3-Dichl	oropropane	<100
Chloromethane		<1,000	Tetrachl	proethene	<100
Vinyl chloride		<20	Dibromo	chloromethane	<100
Bromomethane		<100	1,2-Dibro	omoethane (EDB)	<100
Chloroethane		<100	Chlorobe	nzene	<100
Trichlorofluorometh	ane	<100	Ethylben	zene	2,600
Acetone		<5,000	1,1,1,2-T	etrachloroethane	<100
1,1-Dichloroethene		<100	m,p-Xyle	ne	6,400
Hexane		<100	o-Xylene		1,500
Methylene chloride		<500	Styrene		<100
Methyl t-butyl ether	(MTBE)	<100	Isopropy	lbenzene	<100
trans-1,2-Dichloroet	hene	<100	Bromofor	rm	<100
1,1-Dichloroethane		<100	n-Propyl	benzene	<100
2,2-Dichloropropane		<100	Bromobe	nzene	<100
cis-1,2-Dichloroether	ne	<100	1,3,5-Tri	methylbenzene	<100
Chloroform		<100	1,1,2,2-T	etrachloroethane	<100
2-Butanone (MEK)		<1,000	1,2,3-Tri	chloropropane	<100
1,2-Dichloroethane (	EDC)	<100	2-Chloro	toluene	<100
1,1,1-Trichloroethan	ie	<100	4-Chloro	toluene	<100
1,1-Dichloropropene		<100	tert-Buty	lbenzene	<100
Carbon tetrachloride	ġ.	<100	1,2,4-Tri	methylbenzene	<100
Benzene		<35	sec-Buty	lbenzene	<100
Trichloroethene		<100	p-Isoprop	yltoluene	<100
1,2-Dichloropropane		<100	1,3-Dichl	orobenzene	<100
Bromodichlorometha	ane	<100	1,4-Dichl	orobenzene	<100
Dibromomethane		<100	1,2-Dichl	orobenzene	<100
4-Methyl-2-pentanor	ne	<1,000	1,2-Dibro	omo-3-chloropropane	<1,000
cis-1,3-Dichloroprope	ene	<100	1,2,4-Tri	chlorobenzene	<100
Toluene		610	Hexachle	orobutadiene	<100
trans-1,3-Dichloropr	opene	<100	Naphtha	lene	<100
1,1,2-Trichloroethan	ie	<100	1,2,3-Tri	chlorobenzene	<100
2-Hexanone		<1,000			

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-06-08 08/17/18 08/17/18 08/20/18 Water ug/L (ppb)	1618	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Kelly-Moore, F&BI & 808402-04 082009.D GCMS4 JS	& Infrastructure Solutions 08402
			Lower	Upper	
Surrogates:		% Recovery:	Limit:	Limit:	
1,2-Dichloroethane-c	14	100	57	121	
Toluene-d8		97	63	127	
4-Bromofluorobenzer	ne	106	60	133	
		Concentration			Concentration
Compounds:		ug/L (ppb)	Compour	nds:	ug/L (ppb)
Dichlorodifluoromet	hane	<1	1 3-Dichl	oropropane	<1
Chloromethane	liane	<10	Tetrachl	proethene	<1
Vinvl chloride		<0.2	Dibromo	chloromethane	<1
Bromomethane		<1	1 2-Dibro	omoethane (EDB)	<1
Chloroethane		<1	Chlorobe	nzene	<1
Trichlorofluorometh	ane	<1	Ethylber	izene	<1
Acetone	une	<50	1.1.1.2-T	etrachloroethane	<1
1.1-Dichloroethene		<1	m.p-Xvle	ne	<2
Hexane		<1	o-Xvlene		<1
Methylene chloride		<5	Styrene		<1
Methyl t-butyl ether	(MTBE)	<1	Isopropy	lbenzene	18
trans-1.2-Dichloroet	hene	<1	Bromofor	rm	<1
1,1-Dichloroethane		<1	n-Propyl	benzene	25
2,2-Dichloropropane		<1	Bromobe	nzene	<1
cis-1,2-Dichloroether	ne	<1	1,3,5-Tri	methylbenzene	<1
Chloroform		<1	1,1,2,2-T	etrachloroethane	<1
2-Butanone (MEK)		<10	1,2,3-Tri	chloropropane	<1
1,2-Dichloroethane (	EDC)	<1	2-Chloro	toluene	<1
1,1,1-Trichloroethan	ie	<1	4-Chloro	toluene	<1
1,1-Dichloropropene		<1	tert-Buty	lbenzene	<1
Carbon tetrachloride	)	<1	1,2,4-Tri	methylbenzene	<1
Benzene		< 0.35	sec-Buty	lbenzene	2.6
Trichloroethene		<1	p-Isoprop	yltoluene	<1
1,2-Dichloropropane		<1	1,3-Dichl	orobenzene	<1
Bromodichlorometha	ane	<1	1,4-Dichl	orobenzene	<1
Dibromomethane		<1	1,2-Dichl	orobenzene	<1
4-Methyl-2-pentanor	ne	<10	1,2-Dibro	omo-3-chloropropane	<10
cis-1,3-Dichloroprop	ene	<1	1,2,4-Tri	chlorobenzene	<1
Toluene		<1	Hexachle	orobutadiene	<1
trans-1,3-Dichloropr	opene	<1	Naphtha	lene	<1
1,1,2-Trichloroethan	ie	<1	1,2,3-Tri	chlorobenzene	<1
2-Hexanone		<10			

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-07-08 08/17/18 08/17/18 08/17/18 Water ug/L (ppb)	1618	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Kelly-Moore, F&BI 8 808402-05 081734.D GCMS4 JS	& Infrastructure Solutions 08402
			Lower	Upper	
Surrogates:		% Recovery:	Limit:	Limit:	
1,2-Dichloroethane-c	14	101	57	121	
Toluene-d8		99	63	127	
4-Bromofluorobenzer	ne	96	60	133	
		Concentration			Concentration
Compounds:		ug/L (ppb)	Compour	nds:	ug/L (ppb)
Dichlorodifluoromet	hane	<1	1.3-Dichl	oropropane	<1
Chloromethane	liune	<10	Tetrachl	proethene	<1
Vinvl chloride		< 0.2	Dibromo	chloromethane	<1
Bromomethane		<1	1.2-Dibro	omoethane (EDB)	<1
Chloroethane		<1	Chlorobe	nzene	<1
Trichlorofluorometh	ane	<1	Ethylber	izene	<1
Acetone		<50	1.1.1.2-T	etrachloroethane	<1
1.1-Dichloroethene		<1	m.p-Xvle	ne	<2
Hexane		<1	o-Xvlene		<1
Methylene chloride		<5	Styrene		<1
Methyl t-butyl ether	(MTBE)	<1	Isopropy	lbenzene	<1
trans-1.2-Dichloroet	hene	<1	Bromofor	rm	<1
1,1-Dichloroethane		<1	n-Propyl	benzene	<1
2,2-Dichloropropane		<1	Bromobe	nzene	<1
cis-1,2-Dichloroether	ne	<1	1,3,5-Tri	methylbenzene	<1
Chloroform		<1	1,1,2,2-T	etrachloroethane	<1
2-Butanone (MEK)		<10 ca	1,2,3-Tri	chloropropane	<1
1,2-Dichloroethane (	EDC)	<1	2-Chloro	toluene	<1
1,1,1-Trichloroethan	ie	<1	4-Chloro	toluene	<1
1,1-Dichloropropene		<1	tert-Buty	lbenzene	<1
Carbon tetrachloride	è.	<1	1,2,4-Tri	methylbenzene	<1
Benzene		< 0.35	sec-Buty	lbenzene	<1
Trichloroethene		<1	p-Isoprop	yltoluene	<1
1,2-Dichloropropane		<1	1,3-Dichl	orobenzene	<1
Bromodichlorometha	ane	<1	1,4-Dichl	orobenzene	<1
Dibromomethane		<1	1,2-Dichl	orobenzene	<1
4-Methyl-2-pentanor	ne	<10	1,2-Dibro	omo-3-chloropropane	<10
cis-1,3-Dichloroprop	ene	<1	1,2,4-Tri	chlorobenzene	<1
Toluene		<1	Hexachle	orobutadiene	<1
trans-1,3-Dichloropr	opene	<1	Naphtha	lene	<1
1,1,2-Trichloroethan	ie	<1	1,2,3-Tri	chlorobenzene	<1
2-Hexanone		<10 ca			

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-08-08 08/17/18 08/17/18 08/17/18 Water ug/L (ppb)	1618	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Kelly-Moore, F&BI & 808402-06 081735.D GCMS4 JS	& Infrastructure Solutions 08402
			Lower	Upper	
Surrogates:		% Recovery:	Limit:	Limit:	
1,2-Dichloroethane-o	14	101	57	121	
Toluene-d8		98	63	127	
4-Bromofluorobenze	ne	96	60	133	
		Concentration			Concentration
Compounds:		ug/L (ppb)	Compour	nds:	ug/L (ppb)
Dichlorodifluoromet	hane	<1	1,3-Dichl	loropropane	<1
Chloromethane		<10	Tetrachl	oroethene	<1
Vinyl chloride		< 0.2	Dibromo	chloromethane	<1
Bromomethane		<1	1,2-Dibro	omoethane (EDB)	<1
Chloroethane		<1	Chlorobe	enzene	<1
Trichlorofluorometh	ane	<1	Ethylber	nzene	<1
Acetone		<50	1,1,1,2-T	'etrachloroethane	<1
1,1-Dichloroethene		<1	m,p-Xyle	ene	<2
Hexane		<1	o-Xylene		<1
Methylene chloride		<5	Styrene		<1
Methyl t-butyl ether	· (MTBE)	<1	Isopropy	lbenzene	<1
trans-1,2-Dichloroet	hene	<1	Bromofo	rm	<1
1,1-Dichloroethane		<1	n-Propyl	benzene	<1
2,2-Dichloropropane		<1	Bromobe	enzene	<1
cis-1,2-Dichloroethe	ne	<1	1,3,5-Tri	methylbenzene	<1
Chloroform		<1	1,1,2,2-T	etrachloroethane	<1
2-Butanone (MEK)		<10 ca	1,2,3-Tri	chloropropane	<1
1,2-Dichloroethane (	EDC)	<1	2-Chloro	toluene	<1
1,1,1-Trichloroethar	ne	<1	4-Chloro	toluene	<1
1,1-Dichloropropene		<1	tert-Buty	lbenzene	<1
Carbon tetrachloride	9	<1	1,2,4-Tri	methylbenzene	<1
Benzene		< 0.35	sec-Buty	lbenzene	<1
Trichloroethene		<1	p-Isoprop	oyltoluene	<1
1,2-Dichloropropane		<1	1,3-Dichl	lorobenzene	<1
Bromodichlorometha	ane	<1	1,4-Dichl	lorobenzene	<1
Dibromomethane		<1	1,2-Dichl	lorobenzene	<1
4-Methyl-2-pentanor	ne	<10	1,2-Dibro	omo-3-chloropropane	<10
cis-1,3-Dichloroprop	ene	<1	1,2,4-Tri	chlorobenzene	<1
Toluene		<1	Hexachle	orobutadiene	<1
trans-1,3-Dichloropr	opene	<1	Naphtha	lene	<1
1,1,2-Trichloroethar	ne	<1	1,2,3-Tri	chlorobenzene	<1
2-Hexanone		<10 ca			

## ENVIRONMENTAL CHEMISTS

LowerUpper1.2-Dichloroethane d499571211.2-Dichloroethane d497631274-Bromofluorobenzene9660133ComeentrationCompounds:ug/L (ppb)Compounds:ug/L (ppb)Ug/L (ppb)Dichlorodifluoromethane<11,3-Dichloropropane<1Chloromethane<11,2-Dibmoochloromethane<1Vinyl chloride<0.2Dibmoochloromethane<1Bromomethane<11,2-Dibmoochloromethane<1Trichlorofluoromethane<11,1-Dichloropropane<1Chloroethane<11,2-Dibromochloromethane<1Trichlorofluoromethane<11,1-Zietrachloroethane<1Acctone<501,1,1.2-Tetrachloroethane<11,1-Dichloroethene<1m,p.Xylene<2Hexane<1oXylene<11,1-Dichloroethene<1Bromoform<11,2-Dichloroethene<1Bromoform<11,1-Dichloroethene<1Bromoform<11,1-Dichloroethene<11,3.5-Trimethylbenzene<11,1-Dichloroethene<11,1,2,2-Tetrachloroethane<11,2-Dichloroethene<11,1,2,2-Tetrachloroethane<11,1-Dichloroethene<11,1,2,2-Tetrachloroethane<11,1-Dichloroethene<11,2,3-Trichloropropane<11,1-Dichloroethene<11,2,2-Tetrachloroethane<11,1-Dichloroethene<1	Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-09-08 08/17/18 08/17/18 08/17/18 Water ug/L (ppb)	1618	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Kelly-Moore, F&BI & 808402-07 081736.D GCMS4 JS	& Infrastructure Solutions 08402
Surrogates:% Recovery:Limit:Limit:1.2-Dichloroethane-d499571211 oluene-d897631274-Bromofluorobenzene9660133ConcentrationConcentrationCompounds:ug/L (ppb)Compounds:ug/L (ppb)Dichlorodifluoromethane<1				Lower	Upper	
$1.2 \text{-Dichoroethane-d4} 99 57 121$ Toluene-d8 97 63 127 Toluene-d8 97 63 127 A-Bromofluorobenzene 96 60 133 $\frac{\text{Concentration}}{\text{Compounds:}} \frac{\text{Concentration}}{\text{ug/L} (ppb)} Compounds: ug/L (ppb)$ Dichlorodifluoromethane <1 1.3-Dichloropropane <1 Chloromethane <1 1.3-Dichloropropane <1 Chloromethane <1 1.2-Dibromochloromethane <1 Chloromethane <1 1.2-Dibromochloromethane <1 Chloromethane <1 1.2-Dibromochloromethane <1 Chloroethane <1 1.2-Dibromochloromethane <1 Chloroethane <1 Chlorobenzene <1 Chloroform <1 Chloroethane <1 Propylbenzene <1 Chloroform <1 Chloroethane <1 Romoform <1 Chloroform <1 Chloro	Surrogates:		% Recovery:	Limit:	Limit:	
Toluened897631274-Bromofluorobenzene9660133ConcentrationCompounds:ug/L (ppb)Dichlorodifluoromethane<1	1,2-Dichloroethane-o	14	99	57	121	
4-Bromofluorobenzene9660133Concentration Qompounds:Concentration ug/L (ppb)Compounds:ug/L (ppb)Dichlorodifluoromethane<1	Toluene-d8		97	63	127	
ConcentrationConcentrationCompounds:ug/L (ppb)Dichlorodifluoromethane<1	4-Bromofluorobenze	ne	96	60	133	
Compounds:ug/L (ppb)Compounds:ug/L (ppb)Dichlorodifluoromethane<1			Concentration			Concentration
$ \begin{array}{c c c c c c c c c c c c c c c c c c c $	Compounds:		ug/L (ppb)	Compour	nds:	ug/L (ppb)
Chloromethane<10Tetrachloroethene<1Vinyl chloride<0.2	Dichlorodifluoromet	hane	<1	1.3-Dichl	oropropane	<1
Vinyl chloride $0.2$ Dibromochloromethane $-1$ Bromomethane $1$ $1,2$ -Dibromochlane (EDB) $-1$ Chloroethane $-1$ Chlorobenzene $-1$ Chloroethane $-1$ Chlorobenzene $-1$ Acetone $-50$ $1,1,1,2$ -Tetrachloroethane $-1$ $1,1$ -Dichloroethene $-1$ $m,p,Xylene$ $-2$ Hexane $-1$ $o,Xylene$ $-1$ Methylene chloride $-5$ Styrene $-1$ Methyl t-butyl ether (MTBE) $-1$ Isopropylbenzene $11$ $1,1$ -Dichloroethene $-1$ $n$ -Propylbenzene $11$ $1,1$ -Dichloroethene $-1$ Bromoform $-1$ $2,2$ -Dichloroethene $-1$ $n$ -Propylbenzene $11$ $2,2$ -Dichloroethene $-1$ $n$ -Propylbenzene $-1$ $1,1$ -Dichloroethane $-1$ $n$ -Propylbenzene $-1$ $2,2$ -Dichloroethene $-1$ $n$ -Propylbenzene $-1$ $2,2$ -Dichloroethene $-1$ $1,2,3$ -Trichloroethane $-1$ $2,2$ -Dichloroethane $-1$ $1,2,3$ -Trichloroethane $-1$ $2,2$ -Dichloroethane $-1$ $1,2,3$ -Trichloroethane $-1$ $1,1,1$ -Trichloroethane $-1$ $1,2,3$ -Trichloroethane $-1$ $1,1,1$ -Trichloroethane $-1$ $1,2,3$ -Trichloroethane $-1$ $1,1,1$ -Trichloroethane $-1$ $1,2,4$ -Trimethylbenzene $-1$ $1,1,1$ -Trichloroethane $-1$ $1,2,4$ -Trimethylbenzene $-1$ $1,1,1$ -Dichloropenzene $-1$ $1,2,2$ -Trimethyl	Chloromethane		<10	Tetrachl	oroethene	<1
Bromomethane<11,2-Dibromothane (EDB)<1Chloroethane<1	Vinvl chloride		< 0.2	Dibromo	chloromethane	<1
Chlorothane<1Chlorobenzene<1Chlorothane<1	Bromomethane		<1	1.2-Dibro	omoethane (EDB)	<1
$\begin{array}{c c c c c c c c c c c c c c c c c c c $	Chloroethane		<1	Chlorobe	nzene	<1
Acetone<501.1.1.2-Tetrachloroethane<11.1-Dichloroethene<1	Trichlorofluorometh	ane	<1	Ethylber	izene	<1
1.1-Dichloroethene<1m.p.Xylene<2Hexane<1	Acetone		<50	1.1.1.2-T	'etrachloroethane	<1
Hexane1or Aylene1Methylene chloride<5	1.1-Dichloroethene		<1	m.p-Xvle	ne	<2
Methylene chloride  Methylene chloride Styrene<1	Hexane		<1	o-Xvlene		<1
Mathyl t-butyl ether (MTBE)Isopropylbenzene11trans-1,2-Dichloroethene<1	Methylene chloride		<5	Styrene		<1
Intans-1,2-Dichloroethene<1Bromoform<11,1-Dichloroethane<1	Methyl t-butyl ether	(MTBE)	<1	Isopropy	lbenzene	11
International formationInternational formationInternational formation1.1-Dichloroethane<1	trans-1 2-Dichloroet	hene	<1	Bromofor	rm	<1
2,2-Dichloropropane<1Bromobenzene<12,2-Dichloropropane<1	1.1-Dichloroethane		<1	n-Propyl	benzene	11
In Driving PopulationInternational ConstructionInternational ConstructionCharlen Charlen ConstructionInternational ConstructionInternational ConstructionInternational ConstructionChloroformInternational ConstructionInternational ConstructionInternational ConstructionInternational Construction2-Butanone (MEK)International ConstructionInternational ConstructionInternational ConstructionInternational Construction1,2-Dichloroethane (EDC)International ConstructionInternational ConstructionInternational ConstructionInternational Construction1,1-DichloroptopeneInternational ConstructionInternational ConstructionInternational ConstructionInternational Construction1,1-DichloroptopeneInternational ConstructionInternational ConstructionInternational ConstructionInternational Construction1,1-DichloroptopeneInternational ConstructionInternational ConstructionInternational ConstructionInternational Construction1,1-DichloroptopeneInternational ConstructionInternational Constru	2 2-Dichloropropane		<1	Bromobe	nzene	<1
Chloroform<11,1,2,2-Tetrachloroethane<12-Butanone (MEK)<10 ca	cis-1 2-Dichloroethe	ne	<1	1 3 5-Tri	methylbenzene	<1
2-Butanone (MEK)<10 ca1,2,3-Trichloropropane<12-Butanone (MEK)<10 ca	Chloroform		<1	1,0,0 111 1 1 2 2-T	etrachloroethane	<1
1,2-Dichloroethane (EDC)<12-Chlorotoluene<11,1,1-Trichloroethane<1	2-Butanone (MEK)		<10 ca	1, 1, 2, 2 1 1 2 3-Tri	chloropropane	<1
1,1,1-Trichloroethane<14-Chlorotoluche<11,1,1-Trichloroethane<1	1 2-Dichloroethane (	FDC)	<10 cu	2-Chloro	toluene	<1
1,1,1 Trichlorobrunc<11 tert Butylbenzene<11,1-Dichloropropene<1	1 1 1-Trichloroethar		<1	2 Chlorot 4-Chlorot	toluene	<1
The Diction oproprior<1Carbon tetrachloride<1(1)Carbon tetrachloride<1	1 1-Dichloropropene		<1	tert-Buty	Ilhenzene	<1
Benzene<0.35sec-Butylbenzene<1Trichloroethene<1	Carbon tetrachloride	2	<1	1 9 4-Tri	methylbenzene	<1
Definition<1p-Isopropyltoluene<1Trichloroethene<1	Renzene	<i>,</i>	<0.35	sec-Buty	lhenzene	<1
Internoroctifience<1p isopropyronation<11,2-Dichloropropane<1	Trichloroethene		<0.55	n-Isopror	vltoluone	<1
Fight Diction oproprint<1Fight Diction operation<1Bromodichloromethane<1	1 2-Dichloropropage		<1	1 3-Dichl	lorobenzene	<1
Dibromodultifier<11,4 Dichlorobenzene<1Dibromomethane<1	Bromodichlorometh:	ano	<1	1,0-Dichi 1 <i>1</i> -Dichi	lorobenzene	<1
A-Methyl-2-pentanone<101,2-Dibromo-3-chloropropane<10cis-1,3-Dichloropropene<1	Dibromomethane	ane	<1	1,4 Dichi 1 2-Dichi	lorobenzene	<1
A Methyl 2 perturbation<101,2 Distributes entorspropriate<10cis-1,3-Dichloropropene<1	A-Methyl-2-pentanor	10	<10	1,2-Dichi 1 2-Dibro	mo-3-chloropropape	<10
Toluene<1Hexachlorobutadiene<1trans-1,3-Dichloropropene<1	cis-1 3-Dichloroprop	ano	<10	1,2-Dibio 1 9 <i>1</i> -Tri	chlorobonzono	<1
Trans-1,3-Dichloropropene<1Naphthalene<11,1,2-Trichloroethane<1	Toluene		<1	Hovachle	rohutadiene	<1
1,1,2-Trichloroethane<11,2,3-Trichlorobenzene<12 Hovapopo<10 ca	trans-1 3-Dichloropr	onene	<1	Nanhtha	lene	<1
$\frac{1}{2 \text{ Hovenono}} = \frac{10 \text{ co}}{2 \text{ Hovenono}}$	1 1 2-Trichloroothar	opene	<1	1 9 2-Tri	chlorobenzene	<1
	2-Hexanone		<10 ca	1,~,0-111		<b>~1</b>

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-10-08 08/17/18 08/17/18 08/17/18 Water ug/L (ppb)	1618	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Kelly-Moore, F&BI 8 808402-08 081737.D GCMS4 JS	& Infrastructure Solutions 08402
			Lower	Upper	
Surrogates:		% Recovery:	Limit:	Limit:	
1,2-Dichloroethane-c	14	100	57	121	
Toluene-d8		98	63	127	
4-Bromofluorobenzer	ne	93	60	133	
		Concentration			Concentration
Compounds:		ug/L (ppb)	Compour	nds:	ug/L (ppb)
Dichlorodifluoromet	hane	<1	1.3-Dichl	oropropane	<1
Chloromethane		<10	Tetrachl	proethene	<1
Vinvl chloride		< 0.2	Dibromo	chloromethane	<1
Bromomethane		<1	1.2-Dibro	omoethane (EDB)	<1
Chloroethane		<1	Chlorobe	nzene	<1
Trichlorofluorometh	ane	<1	Ethylber	izene	320 ve
Acetone		<50	1.1.1.2-T	etrachloroethane	<1
1.1-Dichloroethene		<1	m.p-Xvle	ne	930 ve
Hexane		<1	o-Xvlene		<1
Methylene chloride		<5	Styrene		<1
Methyl t-butyl ether	(MTBE)	<1	Isopropy	lbenzene	14
trans-1 2-Dichloroet	hene	<1	Bromofor	rm	<1
1 1-Dichloroethane		<1	n-Propyl	henzene	13
2 2-Dichloropropane		<1	Bromobe	nzene	<1
cis-1 2-Dichloroethe	ne	<1	1 3 5-Tri	methylbenzene	18
Chloroform		<1	1,0,0 111 1 1 2 2-T	etrachloroethane	<1
2-Butanone (MEK)		<10 ca	1, 1, 2, 2 1 1 2 3-Tri	chloropropane	<1
1 2-Dichloroethane (	FDC)	<10 cu	2-Chlorot	toluene	<1
1 1 1-Trichloroethan		<1	2 Chlorot 4-Chlorot	toluene	<1
1 1-Dichloropropene		<1	tert-Buty	Ibenzene	<1
Carbon tetrachloride	<b>`</b>	<1	1 9 4-Tri	methylbenzene	36
Renzene	<i>,</i>	15	sec-Butv	lhenzene	<1
Trichloroethene		<1.0 <1	n-Isopror	vltoluene	<1
1 2-Dichloronronane		<1	1 3-Dichl	orobenzene	<1
Bromodichlorometh:	ane	<1	1,0 Dichi 1 4-Dichi	orobenzene	<1
Dibromomethane		<1	1,4 Dichi 1 2-Dichi	orobenzene	<1
4-Methyl-2-pentanor	1e	<10	1,2 Diem 1 2-Dibro	mo-3-chloropropane	<10
cis-1 3-Dichloroprop	ne	<1	1,2 Dibit 1 2 4-Tri	chlorobenzene	<1
Toluene		<1	Hevachle	robutadiene	<1
trans-1 3-Dichloropr	onene	<1	Nanhtha	lene	<1
1 1 2-Trichloroethar	opene Ip	<1	1 9 3-Tri	chlorobenzene	<1
2-Hexanone		<10 ca	1,~,0 111		· •

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-10-08 08/17/18 08/17/18 08/20/18 Water ug/L (ppb)	1618	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Kelly-Moore, F&BI & 808402-08 1/10 082016.D GCMS4 JS	& Infrastructure Solutions 08402
			Lower	Upper	
Surrogates:		% Recovery:	Limit:	Limit:	
1,2-Dichloroethane-c	14	101	57	121	
Toluene-d8		100	63	127	
4-Bromofluorobenzer	ne	97	60	133	
		Concentration			Concentration
Compounds:		ug/L (ppb)	Compour	nds:	ug/L (ppb)
Diablana diffusanamati	h o m o	.10			.10
Chlanamathana	nane	<10	I,3-DICNI Totro obl	oropropane	<10
Vinue ablanida		<100	Dihnama	oroethene	<10
Vinyi chioride		<2	DIDFOIIIO	chioromethane	<10
Chloroothono		<10	1,2-DIDIC Chloroho	niloeuliane (EDD)	<10
Trichlorofluoromoth	0.00	<10	Ethylbor	nzene	<10
Acotopo	alle	<10		atrachlaraathana	570 <10
1 1 Dichloroothono		< 10	$1, 1, 1, 2^{-1}$		<10
Lovono		<10	ni, p-Ayle	ine	1,100 <10
Mathulana ahlarida		<10	0-Aylene		<10
Methyletie chioride		<0	Iconropu	bonzono	<10
trong 1 2 Dishlenget	(WIIDE)	<10	Bromofor	idenzene	10
1 1 Disklangstham	nene	<10	DI UIIIUIUI	'III	<10
1,1-Dichloroethane		<10	n-Propyn Deservatio	benzene	14
2,2-Dichloropropane		<10	Bromode	nzene	<10
Cls-1,2-Dichloroethei	ne	<10	1,3,5-111	metnyibenzene	19
Chloroform		<10	1,1,2,2-1	etrachioroethane	<10
2-Butanone (MEK)		<100	1,2,3-1ri	chioropropane	<10
1,2-Dichloroethane (	EDC)	<10	2-Chlorot	toluene	<10
1,1,1-1richloroethan	ie	<10	4-Chloro	toluene	<10
I, I-Dichloropropene		<10	tert-Buty	lbenzene	<10
Carbon tetrachloride	<u>)</u>	<10	1,2,4-1ri	methylbenzene	38
Benzene		<3.5	sec-Buty	lbenzene	<10
Trichloroethene		<10	p-Isoprop	oyltoluene	<10
1,2-Dichloropropane		<10	1,3-Dichl	orobenzene	<10
Bromodichlorometha	ane	<10	1,4-Dichl	orobenzene	<10
Dibromomethane		<10	1,2-Dichl	orobenzene	<10
4-Methyl-2-pentanor	ne	<100	1,2-Dibro	omo-3-chloropropane	<100
cis-1,3-Dichloroprop	ene	<10	1,2,4-Tri	chlorobenzene	<10
Toluene		<10	Hexachle	orobutadiene	<10
trans-1,3-Dichloropr	opene	<10	Naphtha	lene	<10
1,1,2-Trichloroethan	ie	<10	1,2,3-Tri	chlorobenzene	<10
2-Hexanone		<100			

## ENVIRONMENTAL CHEMISTS

LowerUpperSurrogates:% Recovery:Limit:Limit:1,2-Dichloroethane-d410257121Tolucne-d897631274-Bronofluorobenzene9760133Compounds:ug/L (ppb)Compounds:ug/L (ppb)Dichloroethane<11,3-Dichloropropane<1Chloromethane<10Tetrachloroethane<1Chloromethane<11,2-Dibromochloromethane<1Srichloroethane<11,2-Dibromochloromethane<1Chlorobenzene<11,2-Dibromochloromethane<1Chlorobenzene<11,2-Dibromochloromethane<1Chlorobenzene<11,1-Dichloroethane<1Chlorobenzene<11,1-Dichloroethane<1Acetone<501,1,1.2-Tetrachloroethane<1Rexane<1 $0$ ,Xylene<1Hexane<1Bromoform<11,1-Dichloroethane<1Bromoform<11,2-Dichloroethane<1Bromoform<11,1-Dichloroethane<1Bromoform<11,2-Dichloroethane<1Bromoform<11,1-Dichloroethane<1Bromoform<11,1-Dichloroethane<1Bromoform<11,1-Dichloroethane<1Bromoform<11,1-Dichloroethane<1Bromoform<11,1-Dichloroethane<1Bromoform<11,1-Dichloroethane<1Bromoform<	Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-03R-9 08/17/18 08/17/18 08/18/18 Water ug/L (ppb)	0-081618	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Kelly-Moore, F&BI & 808402-09 081738.D GCMS4 JS	& Infrastructure Solutions 08402
				Lower	Upper	
$1,2:Dichloroethane-d4 102 57 121 \\ Tohuene-d8 97 63 127 \\ -4:Bromofluorobenzene 97 60 133 \\ \hline Concentration Compounds: ug/L (ppb) Compounds: ug/L (ppb) \\ Dichlorodifluoromethane <1 1,3:Dichloropropane <1 \\ Chloromethane <1 1,3:Dichloropropane <1 \\ Chloromethane <1 1,2:Dibromoethane <1 \\ Chloromethane <1 1,2:Dibromoethane <1 \\ Chlorobenzene <1 \\ Chloro$	Surrogates:		% Recovery:	Limit:	Limit:	
Toluened897631274-Bromofluorobenzene9760133Compounds:ConcentrationCompounds:ug/L (ppb)Dichlorodifluoromethane<1	1,2-Dichloroethane-c	14	102	57	121	
4-Bromofluorobenzene9760133Concentration Qorpounds:Concentration ug/L (ppb)Compounds:ug/L (ppb)Dichlorodifluoromethane<1	Toluene-d8		97	63	127	
ConcentrationConcentrationConcentrationCompounds:ug/L (ppb)Compounds:ug/L (ppb)Dichlorodifluoromethane<1	4-Bromofluorobenzer	ne	97	60	133	
Compounds:ug/L (ppb)Compounds:ug/L (ppb)Dichlorodifluoromethane<1			Concentration			Concentration
$ \begin{array}{c c c c c c c c c c c c c c c c c c c $	Compounds:		ug/L (ppb)	Compour	nds:	ug/L (ppb)
Chloromethane<10Tetrachloroethene<1Vinyl chloride<0.2	Dichlorodifluoromet	hane	<1	1.3-Dichl	oropropane	<1
Vinyl chloride $0.2$ Dibromochloromethane $1$ Bromomethane $1$ $1,2$ -Dibromochloromethane (EDB) $<1$ Chloroethane $<1$ Chlorobenzene $<1$ Trichlorofluoromethane $<1$ Ethylbenzene $<1$ Acctone $<50$ $1,1,1,2$ -Tetrachloroethane $<1$ $1,1$ -Dichloroethene $<1$ $m,p.Xylene$ $<2$ Hexane $<1$ $o.Xylene$ $<1$ Methylen chloride $<5$ Styrene $<1$ Methyl -butyl ether (MTBE) $<1$ Isopropylbenzene $<1$ $1,1$ -Dichloroethene $<1$ $n$ -Propylbenzene $<1$ $1,1$ -Dichloroethene $<1$ Bromoform $<1$ $1,1$ -Dichloropapane $<1$ Bromobenzene $<1$ $cis-1,2$ -Dichloroethene $<1$ $1,3,5$ -Trimethylbenzene $<1$ $2,2$ -Dichloropropane $<1$ $1,2,3$ -Trichloropropane $<1$ $2,2$ -Dichloroethene $<1$ $1,2,3$ -Trichloropropane $<1$ $1,1,1$ -Trichloroethane $<1$ $1,2,3$ -Trichloropropane $<1$ $1,1,1$ -Trichloroethane $<1$ $1,2,4$ -Trimethylbenzene $<1$ $1,1,1$ -Trichloroethane $<1$ $1,2,4$ -Trimethylbenzene $<1$ $1,2$ -Dichloropopane $<1$ $1,3$ -Dichlorobenzene $<1$ $1,1$ -Dichlorophenzene $<1$ $1,2,4$ -Trimethylbenzene $<1$ $1,2$ -Dichlorophenzene $<1$ $1,2,2$ -Trimethylbenzene $<1$ $1,2$ -Dichlorophenzene $<1$ $1,2,2$ -Trimethylbenzene $<1$ $1,1,1$ -Trichloroethane	Chloromethane		<10	Tetrachle	proethene	<1
Bromomethane (II 1,2-Dibromothane (EDB) (I Chloroethane (ID Chlorobenzene (ID Chlorobenzene (ID Chloroethane (ID Chlorobenzene (ID Chlorobenzene (ID Chloroethane (ID Chloroet	Vinvl chloride		<0.2	Dibromo	chloromethane	<1
Chlorothane(1)Chlorothane (EE)(1)Trichlorofluoromethane(1)Chlorobenzene(1)Trichlorofluoromethane(1)Ethylbenzene(1)Acctone(50) $1, 1, 1, 2$ -Tetrachloroethane(1) $1, 1$ -Dichloroethene(1)m, p-Xylene(2)Hexane(1)m, p-Xylene(1)Methyl t-butyl ether (MTBE)(1)Isopropylbenzene(1) $1, 1$ -Dichloroethene(1)Bromoform(1) $1, 1$ -Dichloroethene(1)Bromoform(1) $1, 1$ -Dichloroethene(1)Bromobenzene(1) $2, 2$ -Dichloroethene(1) $1, 3, 5$ -Trimethylbenzene(1) $2, 2$ -Dichloroethene(1) $1, 2, 2$ -Tetrachloroethane(1) $2, 2$ -Dichloroethene(1) $1, 2, 2$ -Tetrachloroethane(1) $2, 2$ -Dichloroethane(1) $1, 2, 2$ -Tetrachloroethane(1) $2, 2$ -Dichloroethane(1) $1, 2, 2, 7$ -Tetrachloroethane(1) $2, 2$ -Dichloroethane(1) $1, 2, 2, 7$ -Tetrachloroethane(1) $1, 2$ -Dichloroethane(1) $2, 3, 7$ -Triholoropropane(1) $1, 2, 1, 1, 1$ -Trichloroethane(2)(1)(2) $1, 1, 1, 1$ -Trichloroethane(1) $1, 2, 4, 7$ -Trimethylbenzene(1) $1, 1, 1, 1, 1$ -Trichloroethane(1) $1, 2, 4, 7, 7, 7, 7, 7, 7, 7, 7, 7, 7, 7, 7, 7,$	Bromomethane		<1	1 2-Dibro	omoethane (EDB)	<1
$\begin{array}{c c c c c c c c c c c c c c c c c c c $	Chloroethane		<1	Chlorobe	nzene	<1
Acetone<501.1.2.Tetrachloroethane<11,1-Dichloroethane<1	Trichlorofluorometh	ane	<1	Ethylber	izene	<1
1.1-Dichloroethene<1n, P.Xylene<2Hexane<1	Acetone		<50	1.1.1.2-T	etrachloroethane	<1
Hexane </td <td>1 1-Dichloroethene</td> <td></td> <td>&lt;1</td> <td>m p-Xvle</td> <td>ne</td> <td>&lt;2</td>	1 1-Dichloroethene		<1	m p-Xvle	ne	<2
Name1111Methylene chloride<5	Hexane		<1	o-Xvlene		<1
InterviewIsoBoyneIsoMethyl t-butyl ether (MTBE)<1	Methylene chloride		<5	Styrene		<1
Internation of the problem of the second s	Methyl t-butyl ether	(MTBE)	<1	Isopropy	lbenzene	4 1
InterformationImage: ConstructionImage: ConstructionImage: Construction1,1-Dichloroethane<1	trans-1 2-Dichloroet	hene	<1	Bromofor	rm	<1
1) Formulation111000 model2,2-Dichloropropane<1	1 1-Dichloroethane		<1	n-Propyl	benzene	8.5
1.1. Dramoproprime1Dramoproprime1cis-1,2-Dichloroethene11,3,5-Trimethylbenzene1Chloroform11,1,2,2-Tetrachloroethane12-Butanone (MEK)10 ca1,2,3-Trichloropropane11,2-Dichloroethane (EDC)12-Chlorotoluene11,1.Trichloroethane14-Chlorotoluene11,1.Trichloropropene1tert-Butylbenzene11,1.Dichloropropene11,2,4-Trimethylbenzene1Carbon tetrachloride11,2,4-Trimethylbenzene1Benzene<0.35	2 2-Dichloropropage		<1	Bromobe	nzene	<1
Chloroform<11,0,0 Trimetry bench<1Chloroform<1	cis-1 2-Dichloroethe	ne	<1	1 3 5-Tri	methylbenzene	<1
2-Butanone (MEK)<10 ca1,2,3-Trichloropropane<12-Butanone (MEK)<10 ca	Chloroform		<1	1,0,0 111 1 1 2 2-T	etrachloroethane	<1
1,2-Dichloroethane (EDC)<12-Chlorotoluene<11,1-Trichloroethane (EDC)<1	2-Butanone (MFK)		<10 ca	1, 1, 2, 2, 7 1 2 3-Tri	chloropropane	<1
1,1. Dictinition (LDC)(1)2 Chilorobutant(1)1,1.1-Trichloroethane<1	1 2-Dichloroethane (	FDC)	<10 cu	2-Chloro	toluene	<1
1,1,1 Theorematic<11 construct<11,1-Dichloropropene<1	1 1 1-Trichloroethan		<1	2 Chlorot 4-Chlorot	toluene	<1
The bick of property<1Carbon tetrachloride<11,2,4-Trimethylbenzene<1Benzene<0.35	1 1-Dichloronronene		<1	tert-Buty	Ibenzene	<1
Benzene<0.35sec-Butylbenzene1.9Trichloroethene<1	Carbon tetrachloride	<b>`</b>	<1	1 9 4-Tri	methylbenzene	<1
DefinitionstateDefinition1.0Trichloroethene<1	Renzene	<i>,</i>	<0.35	sec-Buty	lhenzene	19
11.2-Dichloropropane<11,3-Dichlorobenzene<11,2-Dichloropropane<1	Trichloroethene		<0.00	n-Isonror	vltoluene	<1
T,2-Dichlorophyphile<11,3-Dichlorobenzene<1Bromodichloromethane<1	1 2-Dichloropropana		<1	1 3-Dich	orobenzene	<1
Dibromodicinition<11,4 Dichlorobenzene<1Dibromomethane<1	Bromodichlorometh:	ano	<1	1,5-Dichi 1 <i>4</i> -Dichi	orobenzene	<1
4-Methyl-2-pentanone<1	Dibromomethane		<1	1,4 Dich	orobenzene	<1
A Methyl 2 pentatione<101,2 Diblomo 5 chloropropriopane<10cis-1,3-Dichloropropene<1	A-Methyl-2-pentanor		<10	1,2-Dichi 1 2-Dibro	ma-3-chloropropapa	<10
Toluene<1Hexachlorobutadiene<1Toluene<1	cis-1 3-Dichloroprop	ne	<10	1.2-DIDIC 1.9 1_Tri	chlorohenzene	<1
trans-1,3-Dichloropropene<1Naphthalene<11,1,2-Trichloroethane<1	Toluene		<1	Hovachle	nohutadiene	<1
1,1,2-Trichloroethane<1Napitulalelle<11,1,2,3-Trichlorobenzene<1	trans_1 3_Dichloropr	onene	~1	Nanhtha	lana	~1
	1 1 2-Trichlaroothan	opene	<1	1 9 2-Tri	chlorohenzene	<1
2-Hexanone <10 ca	2-Hexanone		<10 ca	1,~,0-111		<b>~1</b>

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	Method Blar Not Applicat 08/17/18 08/17/18 Water ug/L (ppb)	ık ble	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Kelly-Moore, F&BI 8 08-1788 mb 081710.D GCMS4 JS	& Infrastructure Solutions 08402
			Lower	Upper	
Surrogates:		% Recovery:	Limit:	Limit:	
1,2-Dichloroethane-o	14	101	57	121	
Toluene-d8		98	63	127	
4-Bromofluorobenzer	ne	96	60	133	
		Concentration			Concentration
Compounds:		ug/L (ppb)	Compour	nds:	ug/L (ppb)
Dichlorodifluoromet	hane	<1	1,3-Dichl	loropropane	<1
Chloromethane		<10	Tetrachl	oroethene	<1
Vinyl chloride		< 0.2	Dibromo	chloromethane	<1
Bromomethane		<1	1,2-Dibro	omoethane (EDB)	<1
Chloroethane		<1	Chlorobe	nzene	<1
Trichlorofluorometh	ane	<1	Ethylber	izene	<1
Acetone		<50	1,1,1,2-T	etrachloroethane	<1
1,1-Dichloroethene		<1	m,p-Xyle	ene	<2
Hexane		<1	o-Xylene		<1
Methylene chloride		<5	Styrene		<1
Methyl t-butyl ether	(MTBE)	<1	Isopropy	lbenzene	<1
trans-1,2-Dichloroet	hene	<1	Bromofo	rm	<1
1,1-Dichloroethane		<1	n-Propyl	benzene	<1
2,2-Dichloropropane		<1	Bromobe	nzene	<1
cis-1,2-Dichloroether	ne	<1	1,3,5-Tri	methylbenzene	<1
Chloroform		<1	1,1,2,2-T	etrachloroethane	<1
2-Butanone (MEK)		<10	1,2,3-Tri	chloropropane	<1
1,2-Dichloroethane (	EDC)	<1	2-Chloro	toluene	<1
1,1,1-Trichloroethan	ie	<1	4-Chloro	toluene	<1
1,1-Dichloropropene		<1	tert-Buty	lbenzene	<1
Carbon tetrachloride	è	<1	1,2,4-Tri	methylbenzene	<1
Benzene		< 0.35	sec-Buty	lbenzene	<1
Trichloroethene		<1	p-Isoprop	yltoluene	<1
1,2-Dichloropropane		<1	1,3-Dich	lorobenzene	<1
Bromodichlorometha	ane	<1	1,4-Dichl	lorobenzene	<1
Dibromomethane		<1	1,2-Dichl	lorobenzene	<1
4-Methyl-2-pentanor	ne	<10	1,2-Dibro	omo-3-chloropropane	<10
cis-1,3-Dichloroprop	ene	<1	1,2,4-Tri	chlorobenzene	<1
Toluene		<1	Hexachle	orobutadiene	<1
trans-1,3-Dichloropr	opene	<1	Naphtha	lene	<1
1,1,2-Trichloroethan	ie	<1	1,2,3-Tri	chlorobenzene	<1
2-Hexanone		<10			

#### ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-02R-0 08/17/18 08/17/18 08/17/18 Water ug/L (ppb)	81618	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Infrastructure Solutions Kelly-Moore, F&BI 808402 808402-01 1/2 081709.D GCMS6 VM
Surrogates: Anthracene-d10 Benzo(a)anthracene	-d12	% Recovery: 109 117	Lower Limit: 31 25	Upper Limit: 160 165
Compounds:		ug/L (ppb)		
Naphthalene		< 0.06		
Acenaphthylene		< 0.06		
Acenaphthene		< 0.06		
Fluorene		< 0.06		
Phenanthrene		< 0.06		
Anthracene		< 0.06		
Fluoranthene		< 0.06		
Pyrene		< 0.06		
Benz(a)anthracene		< 0.06		
Chrysene		< 0.06		
Benzo(a)pyrene		< 0.06		
Benzo(b)fluoranthen	ie	< 0.06		
Benzo(k)fluoranther	ne	< 0.06		
Indeno(1,2,3-cd)pyre	ene	< 0.06		
Dibenz(a,h)anthrace	ene	< 0.06		
Benzo(g,h,i)perylene	<u>,</u>	< 0.06		

#### ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-03R-0 08/17/18 08/17/18 08/17/18 Water ug/L (ppb)	81618	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Infrastructure Solutions Kelly-Moore, F&BI 808402 808402-02 1/2 081710.D GCMS6 VM
Surrogates: Anthracene-d10 Benzo(a)anthracene	-d12	% Recovery: 107 115	Lower Limit: 31 25	Upper Limit: 160 165
Compounds:		Concentration ug/L (ppb)		
Naphthalene		0.078		
Acenaphthylene		< 0.06		
Acenaphthene		0.27		
Fluorene		0.18		
Phenanthrene		0.074 fb		
Anthracene		< 0.06		
Fluoranthene		< 0.06		
Pyrene		< 0.06		
Benz(a)anthracene		< 0.06		
Chrysene		< 0.06		
Benzo(a)pyrene		< 0.06		
Benzo(b)fluoranther	ne	< 0.06		
Benzo(k)fluoranther	ne	< 0.06		
Indeno(1,2,3-cd)pyre	ene	< 0.06		
Dibenz(a,h)anthrac	ene	< 0.06		
Benzo(g,h,i)perylene	<del>j</del>	< 0.06		

#### ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-04-081 08/17/18 08/17/18 08/17/18 Water ug/L (ppb)	618	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Infrastructure Solutions Kelly-Moore, F&BI 808402 808402-03 1/2 081711.D GCMS6 VM
Surrogates: Anthracene-d10 Benzo(a)anthracene-d	d12	% Recovery: 106 108	Lower Limit: 31 25	Upper Limit: 160 165
Compounds:		Concentration ug/L (ppb)		
Naphthalene		3.9		
Acenaphthylene		< 0.06		
Acenaphthene		0.082		
Fluorene		< 0.06		
Phenanthrene		0.071 fb		
Anthracene		< 0.06		
Fluoranthene		< 0.06		
Pyrene		< 0.06		
Benz(a)anthracene		< 0.06		
Chrysene		< 0.06		
Benzo(a)pyrene		< 0.06		
Benzo(b)fluoranthene	9	< 0.06		
Benzo(k)fluoranthene	e	< 0.06		
Indeno(1,2,3-cd)pyrei	ne	< 0.06		
Dibenz(a,h)anthrace	ne	< 0.06		
Benzo(g,h,i)perylene		< 0.06		

#### ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-06-08 08/17/18 08/17/18 08/17/18 Water ug/L (ppb)	1618	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Infrastructure Solutions Kelly-Moore, F&BI 808402 808402-04 1/2 081712.D GCMS6 VM
Surrogates: Anthracene-d10 Benzo(a)anthracene	-d12	% Recovery: 102 105	Lower Limit: 31 25	Upper Limit: 160 165
Compounds:		Concentration ug/L (ppb)		
Naphthalene		0.16		
Acenaphthylene		8.2		
Acenaphthene		2.2		
Fluorene		0.79		
Phenanthrene		0.20		
Anthracene		0.12		
Fluoranthene		0.18		
Pyrene		0.21		
Benz(a)anthracene		< 0.06		
Chrysene		< 0.06		
Benzo(a)pyrene		< 0.06		
Benzo(b)fluoranther	ne	< 0.06		
Benzo(k)fluoranther	ne	< 0.06		
Indeno(1,2,3-cd)pyre	ene	< 0.06		
Dibenz(a,h)anthrace	ene	< 0.06		
Benzo(g,h,i)perylene	e	< 0.06		

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-07-08 08/17/18 08/17/18 08/17/18 Water ug/L (ppb)	1618	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Infrastructure Solutions Kelly-Moore, F&BI 808402 808402-05 1/2 081713.D GCMS6 VM
Surrogates: Anthracene-d10 Benzo(a)anthracene	-d12	% Recovery: 109 108	Lower Limit: 31 25	Upper Limit: 160 165
Compounds:		Concentration ug/L (ppb)		
Nonhtholono		-0 0C		
		< 0.00		
Acenaphthylene		<0.06		
Acenaphthene		< 0.06		
Fluorene		< 0.06		
Phenanthrene		< 0.06		
Anthracene		< 0.06		
Fluoranthene		< 0.06		
Pyrene		< 0.06		
Benz(a)anthracene		< 0.06		
Chrysene		< 0.06		
Benzo(a)pyrene		< 0.06		
Benzo(b)fluoranthen	e	< 0.06		
Benzo(k)fluoranther	ne	< 0.06		
Indeno(1,2,3-cd)pyre	ene	< 0.06		
Dibenz(a,h)anthrace	ene	< 0.06		
Benzo(g,h,i)perylene	<u>)</u>	< 0.06		

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-08-08 08/17/18 08/17/18 08/17/18 Water ug/L (ppb)	1618	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Infrastructure Solutions Kelly-Moore, F&BI 808402 808402-06 1/2 081716.D GCMS6 VM
Surrogates: Anthracene-d10 Benzo(a)anthracene	-d12	% Recovery: 106 122	Lower Limit: 31 25	Upper Limit: 160 165
Compounds:		Concentration ug/L (ppb)		
Naphthalene		< 0.06		
Acenaphthylene		< 0.06		
Acenaphthene		0.11		
Fluorene		0.34		
Phenanthrene		< 0.06		
Anthracene		0.16		
Fluoranthene		0.24		
Pyrene		0.23		
Benz(a)anthracene		< 0.06		
Chrysene		< 0.06		
Benzo(a)pyrene		< 0.06		
Benzo(b)fluoranther	ne	< 0.06		
Benzo(k)fluoranther	ne	< 0.06		
Indeno(1,2,3-cd)pyre	ene	< 0.06		
Dibenz(a,h)anthrac	ene	< 0.06		
Benzo(g,h,i)perylene	<u>)</u>	< 0.06		

#### ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-09-081 08/17/18 08/17/18 08/17/18 Water ug/L (ppb)	1618	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Infrastructure Solutions Kelly-Moore, F&BI 808402 808402-07 1/2 081717.D GCMS6 VM
Surrogates: Anthracene-d10 Benzo(a)anthracene-	d12	% Recovery: 104 109	Lower Limit: 31 25	Upper Limit: 160 165
		Concentration		
Compounds:		ug/L (ppb)		
Naphthalene		0.12		
Acenaphthylene		< 0.06		
Acenaphthene		3.8		
Fluorene		0.69		
Phenanthrene		0.062 fb		
Anthracene		< 0.06		
Fluoranthene		< 0.06		
Pyrene		< 0.06		
Benz(a)anthracene		< 0.06		
Chrysene		< 0.06		
Benzo(a)pyrene		< 0.06		
Benzo(b)fluoranthene	e	< 0.06		
Benzo(k)fluoranthen	e	< 0.06		
Indeno(1,2,3-cd)pyrei	ne	< 0.06		
Dibenz(a,h)anthrace	ne	< 0.06		
Benzo(g,h,i)perylene		< 0.06		

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-10-08 08/17/18 08/17/18 08/17/18 Water ug/L (ppb)	1618	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Infrastructure Solutions Kelly-Moore, F&BI 808402 808402-08 1/2 081718.D GCMS6 VM
Surrogates: Anthracene-d10 Benzo(a)anthracene	-d12	% Recovery: 108 116	Lower Limit: 31 25	Upper Limit: 160 165
Compounds:		Concentration ug/L (ppb)		
Naphthalene		0.51		
Acenaphthylene		0.069		
Acenaphthene		0.061		
Fluorene		< 0.06		
Phenanthrene		< 0.06		
Anthracene		< 0.06		
Fluoranthene		< 0.06		
Pyrene		< 0.06		
Benz(a)anthracene		< 0.06		
Chrysene		< 0.06		
Benzo(a)pyrene		< 0.06		
Benzo(b)fluoranther	ne	< 0.06		
Benzo(k)fluoranther	ne	< 0.06		
Indeno(1,2,3-cd)pyre	ene	< 0.06		
Dibenz(a,h)anthrace	ene	< 0.06		
Benzo(g,h,i)perylene	<u>)</u>	< 0.06		

#### ENVIRONMENTAL CHEMISTS

Surrogates: Anthracene-d10 Benzo(a)anthracene-d12Kecovery: 108Lower Limit: Limit: Limit: Limit: Limit: Limit: Dimit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Limit: Lim	Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	KMW-03R-9 08/17/18 08/17/18 08/17/18 Water ug/L (ppb)	-081618	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Infrastructure Solutions Kelly-Moore, F&BI 808402 808402-09 1/2 081719.D GCMS6 VM
ConcentrationCompounds:ug/L (ppb)Naphthalene<0.06	Surrogates: Anthracene-d10 Benzo(a)anthracene-	d12	% Recovery: 108 116	Lower Limit: 31 25	Upper Limit: 160 165
Naphthalene         <0.06	Compounds:		Concentration ug/L (ppb)		
Acenaphthylene         <0.06	Naphthalene		< 0.06		
Acenaphthene         0.27           Fluorene         0.17           Phenanthrene         0.066 fb           Anthracene         <0.06	Acenaphthylene		< 0.06		
Fluorene       0.17         Phenanthrene       0.066 fb         Anthracene       <0.06	Acenaphthene		0.27		
Phenanthrene         0.066 fb           Anthracene         <0.06	Fluorene		0.17		
Anthracene       <0.06	Phenanthrene		0.066 fb		
Fluoranthene       <0.06	Anthracene		< 0.06		
Pyrene         <0.06	Fluoranthene		< 0.06		
Benz(a) anthracene<0.06Chrysene<0.06	Pyrene		< 0.06		
Chrysene         <0.06	Benz(a)anthracene		< 0.06		
Benzo(a)pyrene<0.06Benzo(b)fluoranthene<0.06	Chrysene		< 0.06		
Benzo(b)fluoranthene<0.06Benzo(k)fluoranthene<0.06	Benzo(a)pyrene		< 0.06		
Benzo(k)fluoranthene<0.06Indeno(1,2,3-cd)pyrene<0.06	Benzo(b)fluoranthen	e	< 0.06		
Indeno(1,2,3-cd)pyrene<0.06Dibenz(a,h)anthracene<0.06	Benzo(k)fluoranthen	e	< 0.06		
Dibenz(a,h)anthracene <0.06 Benzo(g,h,i)perylene <0.06	Indeno(1,2,3-cd)pyre	ne	< 0.06		
Benzo(g,h,i)perylene <0.06	Dibenz(a,h)anthrace	ene	< 0.06		
	Benzo(g,h,i)perylene		< 0.06		

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Extracted: Date Analyzed: Matrix: Units:	Method Blan Not Applicabl 08/17/18 08/17/18 Water ug/L (ppb)	k le	Client: Project: Lab ID: Data File: Instrument: Operator:	Wood Environment & Infrastructure Solutions Kelly-Moore, F&BI 808402 08-1857 mb 081706.D GCMS6 VM
Surrogates: Anthracene-d10 Benzo(a)anthracene-	d12	% Recovery: 105 113	Lower Limit: 31 25	Upper Limit: 160 165
Compounds:		Concentration ug/L (ppb)		
Naphthalene		< 0.03		
Acenaphthylene		< 0.03		
Acenaphthene		< 0.03		
Fluorene		< 0.03		
Phenanthrene		< 0.03		
Anthracene		< 0.03		
Fluoranthene		< 0.03		
Pyrene		< 0.03		
Benz(a)anthracene		< 0.03		
Chrysene		< 0.03		
Benzo(a)pyrene		< 0.03		
Benzo(b)fluoranthen	е	< 0.03		
Benzo(k)fluoranthen	e	< 0.03		
Indeno(1,2,3-cd)pyre	ne	< 0.03		
Dibenz(a,h)anthrace	ne	< 0.03		
Benzo(g,h,i)perylene		< 0.03		

#### ENVIRONMENTAL CHEMISTS

Date of Report: 08/27/18 Date Received: 08/17/18 Project: Kelly-Moore, F&BI 808402

#### QUALITY ASSURANCE RESULTS FOR THE ANALYSIS OF WATER SAMPLES FOR TPH AS GASOLINE USING METHOD NWTPH-Gx

Laboratory Code: 8	08402-05 (Mat	rix Spike	e)				
-		-		Percent	Percent		
	Reporting	Spike	Sample	Recovery	y Recovery	Acceptance	RPD
Analyte	Units	Level	Result	MS	MSD	Criteria	(Limit 20)
Gasoline	ug/L (ppb)	1,000	<100	97	100	53-117	3
Laboratory Code: L	aboratory Cor	trol Sam	ple				
			Per	rcent			
	Reporting Spike Recovery Acceptar						
Analyte	Units	Le	vel L	CS	Criteria		
Gasoline	ug/L (ppł	o) 1,0	00	96	69-134		

#### ENVIRONMENTAL CHEMISTS

Date of Report: 08/27/18 Date Received: 08/17/18 Project: Kelly-Moore, F&BI 808402

#### QUALITY ASSURANCE RESULTS FOR THE ANALYSIS OF WATER SAMPLES FOR TOTAL PETROLEUM HYDROCARBONS AS DIESEL EXTENDED USING METHOD NWTPH-Dx

Laboratory Code: 80	08402-05 (Matrix	Spike)					
Ū				Percent	Percent		
	Reporting	Spike	Sample	Recovery	Recovery	Acceptance	RPD
Analyte	Units	Level	Result	MS	MSD	Criteria	(Limit 20)
Diesel Extended	ug/L (ppb)	2,500	<50	91	84	50-150	8
Laboratory Code: Laboratory	aboratory Contro	l Sample					
			Percent				
	Reporting	Spike	Recovery	y Accept	ance		
Analyte	Units	Level	LCS	Crite	ria		
Diesel Extended	ug/L (ppb)	2,500	84	63-14	42		

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#### ENVIRONMENTAL CHEMISTS

Date of Report: 08/27/18 Date Received: 08/17/18 Project: Kelly-Moore, F&BI 808402

#### QUALITY ASSURANCE RESULTS FOR THE ANALYSIS OF WATER SAMPLES FOR TOTAL METALS USING EPA METHOD 6020B

Laboratory Code: 808402-05 (Matrix Spike)

		F	)	Percent	Percent		
	Reporting	Spike	Sample	Recovery	Recovery	Acceptance	RPD
Analyte	Units	Level	Result	MS	MSD	Criteria	(Limit 20)
Arsenic	ug/L (ppb)	10	<1	109	102	75-125	7
Chromium	ug/L (ppb)	20	<1	95	94	75-125	1
Copper	ug/L (ppb)	20	<5	93	89	75-125	4
Lead	ug/L (ppb)	10	<1	93	90	75-125	3
Mercury	ug/L (ppb)	5	<1	92	91	75-125	1
Nickel	ug/L (ppb)	20	<1	94	91	75-125	3
Zinc	ug/L (ppb)	50	<5	97	90	75-125	7

Laboratory Code: Laboratory Control Sample

		Percent	
Reporting	Spike	Recovery	Acceptance
Units	Level	LCS	Criteria
ug/L (ppb)	10	97	80-120
ug/L (ppb)	20	107	80-120
ug/L (ppb)	20	106	80-120
ug/L (ppb)	10	101	80-120
ug/L (ppb)	5	98	80-120
ug/L (ppb)	20	107	80-120
ug/L (ppb)	50	92	80-120
	Reporting Units ug/L (ppb) ug/L (ppb) ug/L (ppb) ug/L (ppb) ug/L (ppb) ug/L (ppb) ug/L (ppb)	Reporting Units         Spike Level           ug/L (ppb)         10           ug/L (ppb)         20           ug/L (ppb)         20           ug/L (ppb)         10           ug/L (ppb)         5           ug/L (ppb)         20           ug/L (ppb)         5           ug/L (ppb)         20           ug/L (ppb)         5           ug/L (ppb)         50	Reporting         Spike         Recovery           Units         Level         LCS           ug/L (ppb)         10         97           ug/L (ppb)         20         107           ug/L (ppb)         20         106           ug/L (ppb)         10         97           ug/L (ppb)         20         106           ug/L (ppb)         5         98           ug/L (ppb)         20         107           ug/L (ppb)         50         92
#### ENVIRONMENTAL CHEMISTS

Date of Report: 08/27/18 Date Received: 08/17/18 Project: Kelly-Moore, F&BI 808402

#### QUALITY ASSURANCE RESULTS FOR THE ANALYSIS OF WATER SAMPLES FOR VOLATILES BY EPA METHOD 8260C

Laboratory Code: 808402-05 (Matrix Spike)

U U	•			Percent	Percent		
	Reporting	Spike	Sample	Recovery	Recovery	Acceptance	RPD
Analyte	Units	Level	Result	MS	MSD	Criteria	(Limit 20)
Dichlorodifluoromethane	ug/L (ppb)	50	<1	100	101	10-172	1
Chloromethan e	ug/L (ppb)	50	<10	90	92	25-166	2
Vinyl chloride	ug/L (ppb)	50	<0.2	95	97	36-166	2
Bromomethane	ug/L (ppb)	50	<1	105	106	47-169	1
Trichlorofluoromethane	ug/L (ppb)	50	<1	102	103	40-100	1
Acetone	ug/L (ppb)	250	<50	90	89	10-182	1
1,1-Dichloroethene	ug/L (ppb)	50	<1	110	109	60-136	1
Hexane	ug/L (ppb)	50	<1	98	100	52-150	2
Methylene chloride	ug/L (ppb)	50	<5	106	104	67-132	2
Methyl t-butyl ether (MTBE)	ug/L (ppb)	50	<1	106	104	74-127	2
1 1-Dichloroethane	ug/L (ppb)	50	<1	100	105	70-128	1
2.2-Dichloropropane	ug/L (ppb)	50	<1	130	121	36-154	7
cis-1,2-Dichloroethene	ug/L (ppb)	50	<1	103	101	71-127	2
Chloroform	ug/L (ppb)	50	<1	101	100	65-132	1
2-Butanone (MEK)	ug/L (ppb)	250	<10	89	92	10-129	3
1,2-Dichloroethane (EDC)	ug/L (ppb)	50	<1	95	96	69-133	1
1,1,1-111CHI010etHalle	ug/L (ppb)	50	<1	109	108	69-133	1
Carbon tetrachloride	ug/L (ppb)	50	<1	100	100	56-152	1
Benzene	ug/L (ppb)	50	< 0.35	98	97	76-125	1
Trichloroethene	ug/L (ppb)	50	<1	97	97	66-135	0
1,2-Dichloropropane	ug/L (ppb)	50	<1	95	97	78-125	2
Bromodichloromethane	ug/L (ppb)	50	<1	99	99	61-150	0
A-Methyl-2-pentanone	ug/L (ppb)	250	<1	97	97 101	10-141	0
cis-1 3-Dichloropropene	ug/L (ppb)	50	<10	96	99	72-132	3
Toluene	ug/L (ppb)	50	<1	97	96	76-122	1
trans-1,3-Dichloropropene	ug/L (ppb)	50	<1	96	99	76-130	3
1,1,2-Trichloroethane	ug/L (ppb)	50	<1	93	97	68-131	4
2-Hexanone	ug/L (ppb)	250	<10	82	93	10-185	13
1,3-Dichloropropane	ug/L (ppb)	50 50	<1	94 104	97	/1-128	3
Dibromochloromethane	ug/L (ppb)	50	<1	98	103	70-139	2
1,2-Dibromoethane (EDB)	ug/L (ppb)	50	<1	94	98	69-134	$\tilde{4}$
Chlorobenzene	ug/L (ppb)	50	<1	96	97	77-122	1
Ethylbenzene	ug/L (ppb)	50	<1	96	97	69-135	1
1,1,1,2-Tetrachloroethane	ug/L (ppb)	50	<1	105	103	73-137	2
m,p-Xylene	ug/L (ppb)	100	<2	98	99	69-135	1
Styrene	ug/L (ppb)	50	<1	98	101	71-133	3
Isopropylbenzene	ug/L (ppb)	50	<1	101	99	65-142	2
Bromoform	ug/L (ppb)	50	<1	95	98	65-142	3
n-Propylbenzene	ug/L (ppb)	50	<1	99	99	58-144	0
Bromobenzene	ug/L (ppb)	50	<1	97	99	75-124	2
1,3,5-1 rimethylbenzene	ug/L (ppb)	50	<1	102	100	66-137	2
1,1,2,2- Tetracilloropenane	ug/L (ppb)	50	<1	93 88	93	53-150	6
2-Chlorotoluene	ug/L (ppb)	50	<1	99	98	66-127	1
4-Chlorotoluene	ug/L (ppb)	50	<1	96	98	65-130	2
tert-Butylbenzene	ug/L (ppb)	50	<1	102	96	65-137	6
1,2,4-Trimethylbenzene	ug/L (ppb)	50	<1	102	99	59-146	3
sec-Butylbenzene	ug/L (ppb)	50	<1	102	99	64-140	3
1.3-Dichlorobenzene	ug/L (ppb)	50 50	<1	101	99 07	00-141 72-123	2
1.4-Dichlorobenzene	ug/L (ppb)	50	<1	95	96	69-126	1
1,2-Dichlorobenzene	ug/L (ppb)	50	<1	98	97	69-128	1
1,2-Dibromo-3-chloropropane	ug/L (ppb)	50	<10	96	93	32-164	3
1,2,4-Trichlorobenzene	ug/L (ppb)	50	<1	120	114	66-136	5
Hexachlorobutadiene	ug/L (ppb)	50	<1	114	109	60-143	4
Naphthalene	ug/L (ppb)	50	<1	110	105	44-164	5
1,2,3-1110110F0Denzene	ug/L (ppb)	50	<1	119	114	09-148	4

#### ENVIRONMENTAL CHEMISTS

Date of Report: 08/27/18 Date Received: 08/17/18 Project: Kelly-Moore, F&BI 808402

#### QUALITY ASSURANCE RESULTS FOR THE ANALYSIS OF WATER SAMPLES FOR VOLATILES BY EPA METHOD 8260C

Laboratory Code: Laboratory Control Sample

c c	•		Percent	Percent		
	Reporting	Spike	Recovery	Recovery	Acceptance	RPD
Analyte	Units	Level	LCS	LCSD	Criteria	(Limit 20)
Dichlorodifluoromethane	ug/L (ppb)	50	104	107	25-158	3
Chloromethane	ug/L (ppb)	50	92	96	45-156	4
Vinyi chioride Bromomothano	ug/L (ppb)	50	98 107	101	50-154 55-143	3
Chloroethane	ug/L (ppb)	50	95	97	53-145 58-146	4
Trichlorofluoromethane	ug/L (ppb)	250	104	104	50-150	õ
Acetone	ug/L (ppb)	250	93	89	53-131	4
1,1-Dichloroethene	ug/L (ppb)	50	109	107	67-136	2
Hexane	ug/L (ppb)	50	98	96	57-137	2
Methylene chloride	ug/L (ppb)	50	105	103	39-148	2
metnyl t-butyl etner (MIBE)	ug/L (ppb)	50	105	105	68 128	0
1 1-Dichloroethane	ug/L (ppb)	50	105	98	79-120	2
2.2-Dichloropropane	ug/L (ppb)	50	123	126	55-143	2
cis-1,2-Dichloroethene	ug/L (ppb)	50	101	100	80-123	1
Chloroform	ug/L (ppb)	50	100	98	80-121	2
2-Butanone (MEK)	ug/L (ppb)	250	96	94	57-149	2
1,2-Dichloroethane (EDC)	ug/L (ppb)	50	97	95	73-132	2
1,1,1-Trichloroethane	ug/L (ppb)	50	109	109	83-130	0
I,I-Dichloropropene	ug/L (ppb)	50	100	98	77-129	2
Carbon tetrachioride	ug/L (ppb)	50	106	107	70-108	1
Trichloroethene	ug/L (ppb)	50	98	96	80-120	2
1.2-Dichloropropane	ug/L (ppb)	50	98	96	77-123	2
Bromodichloromethane	ug/L (ppb)	50	100	98	81-133	2
Dibromomethane	ug/L (ppb)	50	98	97	82-125	1
4-Methyl-2-pentanone	ug/L (ppb)	250	103	101	65-138	2
cis-1,3-Dichloropropene	ug/L (ppb)	50	100	97	82-132	3
Toluene	ug/L (ppb)	50	97	95	72-122	2
trans-1,3-Dichloropropene	ug/L (ppb)	50	98	95	80-136	3
2-Hevenone	ug/L (ppb)	250	97	94	70-124 60-136	3
1.3-Dichloropropane	ug/L (ppb)	50	98	95	76-126	3
Tetrachloroethene	ug/L (ppb)	50	104	102	76-121	2
Dibromochloromethane	ug/L (ppb)	50	100	97	84-133	3
1,2-Dibromoethane (EDB)	ug/L (ppb)	50	98	96	82-125	2
Chlorobenzene	ug/L (ppb)	50	97	95	83-114	2
Ethylbenzene	ug/L (ppb)	50	100	97	77-124	3
1,1,1,2-1 etrachioroethane	ug/L (ppb)	50	102	101	84-127	1
n,p-Aylene	ug/L (ppb)	50	101	99 100	81-121	ے 1
Styrene	ug/L (ppb)	50	100	100	84-119	0
Isopropylbenzene	ug/L (ppb)	50	100	98	85-117	2
Bromoform	ug/L (ppb)	50	98	97	74-136	1
n-Propylbenzene	ug/L (ppb)	50	100	96	74-126	4
Bromobenzene	ug/L (ppb)	50	99	97	80-121	2
1,3,5-Trimethylbenzene	ug/L (ppb)	50	100	98	78-123	2
1,1,2,2-1 etrachioroethane	ug/L (ppb)	50	96	94	66-126	2
2-Chlorotoluene	ug/L (ppb)	50	93	91	07-124	23
4-Chlorotoluene	ug/L (ppb)	50	98	95	78-128	3
tert-Butylbenzene	ug/L (ppb)	50	98	96	80-123	2
1,2,4-Trimethylbenzene	ug/L (ppb)	50	100	98	79-122	2
sec-Butylbenzene	ug/L (ppb)	50	100	97	80-125	3
p-Isopropyltoluene	ug/L (ppb)	50	100	97	81-123	3
1,3-Dichlorobenzene	ug/L (ppb)	50	97	95	85-116	2
1,4-Dichlorobenzene	ug/L (ppb)	50	96	94	84-121	2
1,2-Dichiolopenzene 1 2-Dibromo-3-chloropropane	ug/L (ppb)	50	97	90 95	80-110 57-141	2
1.2.4-Trichlorobenzene	ug/L (ppb)	50	115	115	72-130	0
Hexachlorobutadiene	ug/L (ppb)	50	109	110	53-141	1
Naphthalene	ug/L (ppb)	50	107	108	64-133	1
1,2,3-Trichlorobenzene	ug/L (ppb)	50	114	116	65-136	2

#### ENVIRONMENTAL CHEMISTS

Date of Report: 08/27/18 Date Received: 08/17/18 Project: Kelly-Moore, F&BI 808402

#### QUALITY ASSURANCE RESULTS FOR THE ANALYSIS OF WATER SAMPLES FOR PAHS BY EPA METHOD 8270D SIM

Laboratory Code: 808402-05 1/2 (Matrix Spike)

		(pine)	~ .	-	-		
			Sample	Percent	Percent		
	Reporting	Spike	Result	Recovery	Recovery	Acceptance	RPD
Analyte	Units	Level	(Wet wt)	MS	MSD	Criteria	(Limit 20)
Naphthalene	ug/L (ppb)	1	< 0.06	94	92	10-172	2
Acenaphthylene	ug/L (ppb)	1	< 0.06	100	99	38-137	1
Acenaphthene	ug/L (ppb)	1	< 0.06	95	94	20-150	1
Fluorene	ug/L (ppb)	1	< 0.06	99	99	10-181	0
Phenanthrene	ug/L (ppb)	1	< 0.06	96	95	58-109	1
Anthracene	ug/L (ppb)	1	< 0.06	99	101	47-114	2
Fluoranthene	ug/L (ppb)	1	< 0.06	99	98	10-171	1
Pyrene	ug/L (ppb)	1	< 0.06	96	96	63-107	0
Benz(a)anthracene	ug/L (ppb)	1	< 0.06	92	92	60-93	0
Chrysene	ug/L (ppb)	1	< 0.06	92	90	60-102	2
Benzo(b)fluoranthene	ug/L (ppb)	1	< 0.06	68	66	62-91	3
Benzo(k)fluoranthene	ug/L (ppb)	1	< 0.06	63	61	51-98	3
Benzo(a)pyrene	ug/L (ppb)	1	< 0.06	64	61	60-86	5
Indeno(1,2,3-cd)pyrene	ug/L (ppb)	1	< 0.06	36	31	10-98	15
Dibenz(a,h)anthracene	ug/L (ppb)	1	< 0.06	38	32	10-97	17
Benzo(g,h,i)perylene	ug/L (ppb)	1	< 0.06	37	31	10-102	18

Laboratory Code: Laboratory Control Sample

<i>y</i>	ľ		Percent	Percent		
	Reporting	Spike	Recovery LCS	Recovery	Acceptance	RPD
Analyte	Units	Level	-	LCSD	Criteria	(Limit 20)
Naphthalene	ug/L (ppb)	1	90	94	67-116	4
Acenaphthylene	ug/L (ppb)	1	90	100	65-119	11
Acenaphthene	ug/L (ppb)	1	90	96	66-118	6
Fluorene	ug/L (ppb)	1	90	98	64-125	9
Phenanthrene	ug/L (ppb)	1	93	97	67-120	4
Anthracene	ug/L (ppb)	1	94	99	65-122	5
Fluoranthene	ug/L (ppb)	1	93	97	65-127	4
Pyrene	ug/L (ppb)	1	97	95	62-130	2
Benz(a)anthracene	ug/L (ppb)	1	95	98	60-118	3
Chrysene	ug/L (ppb)	1	96	100	66-125	4
Benzo(b)fluoranthene	ug/L (ppb)	1	101	98	55-135	3
Benzo(k)fluoranthene	ug/L (ppb)	1	92	98	62-125	6
Benzo(a)pyrene	ug/L (ppb)	1	93	97	58-127	4
Indeno(1,2,3-cd)pyrene	ug/L (ppb)	1	99	104	36-142	5
Dibenz(a,h)anthracene	ug/L (ppb)	1	93	104	37-133	11
Benzo(g,h,i)perylene	ug/L (ppb)	1	94	100	34-135	6

ENVIRONMENTAL CHEMISTS

#### **Data Qualifiers & Definitions**

a - The analyte was detected at a level less than five times the reporting limit. The RPD results may not provide reliable information on the variability of the analysis.

b - The analyte was spiked at a level that was less than five times that present in the sample. Matrix spike recoveries may not be meaningful.

ca - The calibration results for the analyte were outside of acceptance criteria. The value reported is an estimate.

c - The presence of the analyte may be due to carryover from previous sample injections.

cf - The sample was centrifuged prior to analysis.

d - The sample was diluted. Detection limits were raised and surrogate recoveries may not be meaningful.

dv - Insufficient sample volume was available to achieve normal reporting limits.

f - The sample was laboratory filtered prior to analysis.

fb - The analyte was detected in the method blank.

fc - The compound is a common laboratory and field contaminant.

 $hr\ \text{-}\ The\ sample\ and\ duplicate\ were\ reextracted\ and\ reanalyzed.\ RPD\ results\ were\ still\ outside\ of\ control\ limits.\ Variability\ is\ attributed\ to\ sample\ inhomogeneity.$ 

hs - Headspace was present in the container used for analysis.

ht – The analysis was performed outside the method or client-specified holding time requirement.

 ${\rm ip}$  - Recovery fell outside of control limits. Compounds in the sample matrix interfered with the quantitation of the analyte.

j - The analyte concentration is reported below the lowest calibration standard. The value reported is an estimate.

 ${\rm J}$  - The internal standard associated with the analyte is out of control limits. The reported concentration is an estimate.

jl - The laboratory control sample(s) percent recovery and/or RPD were out of control limits. The reported concentration should be considered an estimate.

js - The surrogate associated with the analyte is out of control limits. The reported concentration should be considered an estimate.

lc - The presence of the analyte is likely due to laboratory contamination.

L - The reported concentration was generated from a library search.

nm - The analyte was not detected in one or more of the duplicate analyses. Therefore, calculation of the RPD is not applicable.

pc - The sample was received with incorrect preservation or in a container not approved by the method. The value reported should be considered an estimate.

ve - The analyte response exceeded the valid instrument calibration range. The value reported is an estimate.

vo - The value reported fell outside the control limits established for this analyte.

x - The sample chromatographic pattern does not resemble the fuel standard used for quantitation.

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Seattle, WA 98119-2029	Relinquished by:		<b>K</b> 3															
Ph. (206) 285-8282	Received by:				· · · ·													



# Appendix C

#### Visual/Audio Inspection

Item	Inspected (Y/N)	Condition (Cracks, leaks, non-operational gauges, etc.)
Above Ground Piping	Y	No issurs
Control Pump (Regenerative Blower)	Y	(On) Off)
Entrainment Pump (Transfer Pump)	X	(Aujo / Hand / Off)
Pressure Gauges/Flow Meters	Y	All from themas
Knockout Tank (record level)	7 30	% full
Knockout Water Tote (record level)	Y 60	% full
Dilution Valve Status	Y	(1041)
Recirculation Valve Status	¥	Orm

#### **CATOX Screen Readings**

Item	Reading	Operating Range
Catox In (T <sub>1</sub> )	663 °F	>650°F
Catox Out (T <sub>2</sub> )	637 °F	600 – 650 °F
Heat Ex (T <sub>3</sub> )	35% °F	300 – 400 °F
Flow	225 SCFM	<300 SCFM
LEL	13 %	5-15 %

#### System Gauge Readings

1

ltem	Reading
FE – 1	2.9 "WC
PI – 1	O "WC (vacuum)
TI – 1	59"F °F
FE-2	1,2 "WC

#### **FID Measurements**

Location	Time	FID Reading (ppm)	Valve Position (record notch)	Vacuum ("WC )	Differential Pressure ("WC)
Western Manifold	-	-			
SVE – 13	1846	18.0	3	415	0.05
SVE - 12	0857	55.9	6	4.6	016
SVE – 11	0857	30.8	3	3.0	0.02
SVE – 10	MOL	794.6	6	Ø.0	500
SVE - 09	Q911	246.3	. 6	3.8	0.11
Eastern manifold	-		The second second second		
SVE - 01	0917	98,6	3	2:2	0.14
SVE - 03	0420	>3980	6	3.4	0.05
SVE - 05	allou	73988	6	26	0.10
SVE - 07	0425	61.9	6	17.5	2.97
SVE - 08	0922	38.4	2	25	020
SVE - 06	0026	326.9	3	28	0.18
SVE - 04	QUIPD	<u>44.9</u>	3	2.9	0.09
SVE - 02	0946	4.3	3	3.8	0.08
SVE Influent	PASS	2551	Carlos and a state of		Substanting the second
SVE Effluent	1006	17.3		and the set	and the second

Influent Sample ID: <u>1070629(8</u> Influent Sample Time: <u>0959</u>

Effluent Sample ID: <u>EFF067916</u> Effluent Sample Time: <u>1010</u>

Field Representative (Print and Sign): William Young

\_ Date of Visit: 6/ willing me

#### Visual/Audio Inspection

ltem	Inspected (Y/N)	Condition (Cracks, leaks, non-operational gauges, etc.)
Above Ground Piping	X	No TEGNA
Control Pump (Regenerative Blower)	Y	(Oh / Off)
Entrainment Pump (Transfer Pump)	X	(Auto / Hand / Off)
Pressure Gauges/Flow Meters	r	Q11 funi?
Knockout Tank (record level)	-7/10/13	% full
Knockout Water Tote (record level)	7/10	% full
Dilution Valve Status	Diorn Gisably	
Recirculation Valve Status	Stin Within	

# AFTER MODIFICATIONS > \* recine 1 difution Gir-CATOX Screen Readings

Item	Readi	ing	Operating Range
Catox In (T <sub>1</sub> )	552	٥F	>650°F
Catox Out (T <sub>2</sub> )	670	۰F	600 – 650 °F
Heat Ex (T <sub>3</sub> )	345	٥F	300 - 400 °F
Flow	235 SC	FM	<300 SCFM
LEL	17	%	5-15 %

Item	Reading
FE – 1	"WC
PI – 1	"WC (vacuum)
TI – 1	٥F
FE-2	"WC

#### FID Measurements

Location	Time	FID Reading (ppm)	Valve Position (record notch)	Vacuum (''WC )	Differential Pressure ("WC)
Western Manifold					
SVE - 13					
SVE - 12					
SVE - 11					
SVE - 10					
SVE - 09					
Eastern manifold				The States	
SVE - 01		т. Т.			
SVE - 03					
SVE - 05					
SVE - 07					
SVE - 08		-			
SVE - 06					
SVE - 04					~
SVE - 02	1				
SVE Influent					
SVE Effluent				and the second second	/

Influent Sample ID: Influent Sample Time:

Effluent Sample ID: Effluent Sample Time:

\_\_\_\_ Date of Visit: 7/12/16

Field Representative (Print and Sign): <u>William Young</u>

#### Visual/Audio Inspection

	Inspected	Condition (Cracks, leaks, non-operational
Item	(Y/N)	gauges, etc.)
Above Ground Piping	X	Nº 1341-7
Control Pump (Regenerative Blower)	V	((On/ Off)
Entrainment Pump (Transfer Pump)	Ý	(Auto / Hand / Off)
Pressure Gauges/Flow Meters	Y	1-aling
Knockout Tank (record level)	Q	% full
Knockout Water Tote (record level)	70	% full
Dilution Valve Status	X	oxin
Recirculation Valve Status	X	APON

#### **CATOX Screen Readings**

Item	Reading		Operating Range	
Catox In (T1)	636	°F	>650°F	
Catox Out (T <sub>2</sub> )	634	٩F	600 – 650 °F	
Heat Ex (T <sub>3</sub> )	352	°F	300 – 400 °F	
Flow	251 SC	FM	<300 SCFM	
LEL	14	%	5-15 %	

#### System Gauge Readings

ltem	Reading		
FE – 1	1.4 "WC		
PI – 1	vacuum)		
TI – 1	62 °F		
FE-2	1-3 "WC		

#### **FID Measurements**

Location	Time	FID Reading (ppm)	Valve Position (record notch)	Vacuum (''WC )	Differential Pressure (''WC )
Western Manifold		<i>~</i>			
SVE – 13	1000	29.9	3	3.2	0.17
SVE – 12	1005	357.9	6	1.5	007
SVE - 11	jano	116.2	3	ic	8-10
SVE – 10	inis	11 72	6	0.1	0.01
SVE – 09	1021	505.3	6	15	0.11
Eastern manifold	$\sim$	)			
SVE - 01	ing	238	ζ	8.6	9.07
SVE - 03	10.9	73950	6	0	9.00
SVE – 05	1035	>3960	6	0	0.02
SVE - 07	1039	71116	6	21	0.05
SVE – 08	lour	Ian.s	3	0	0.07
SVE - 06	1055	367.6	3	0.5	0.00
SVE - 04	1058	137.8	3	0	0.06
SVE - 02	11.00	32.0	2	3.0	9.05
SVE Influent	1110	315.7			
SVE Effluent	120	25.5		Destroyed and	

Influent Sample ID: <u>1NF 073118</u> Influent Sample Time: <u>1111</u>

Effluent Sample ID: <u>EFF07311k</u> Effluent Sample Time: <u>1120</u>

Field Representative (Print and Sign): William Young Date of Visit: \_\_\_\_\_

#### Visual/Audio Inspection

ltem	Inspected (Y/N)	Condition (Cracks, leaks, non-operational gauges, etc.)
Above Ground Piping	N'	appa
Control Pump (Regenerative Blower)	1	(On) Off)
Entrainment Pump (Transfer Pump)	1	(Auto) / Hand / Off)
Pressure Gauges/Flow Meters	y	Sping advage Same full so Call (2)
Knockout Tank (record level)	y	% full 10/2 (0.200 and 1000
Knockout Water Tote (record level)		% full (109 - 109
Dilution Valve Status	И	men
Recirculation Valve Status	V	open ino reevaluation

#### **CATOX Screen Readings**

ltem	Reading	Operating Range
Catox In (T <sub>1</sub> )	680 °F	>650°F
Catox Out (T <sub>2</sub> )	(133 °F	600 - 650 °F
Heat Ex (T <sub>3</sub> )	356 °F	300 - 400 °F
Flow	242SCFM	<300 SCFM
LEL	5 %	5-15 %

System Gaug	e Readings	1. Andres -
ltem	Reading	1 Allato
FE – 1	WC	Dig K /kalse / set
PI – 1	-1.9 "WC (vacuum)	gange=1.8"
TI – 1	(da °F	manometer= 3.4" WC
FE-2	14 "WC	

#### **FID Measurements**

FID Measurement	t <u>s</u>		)	readings t	nom Her	Pagange (discrepance
Location	Time	FID Reading (ppm)	Valve Position (record notch)	Vacuum (''WC )	Differential Pressure ("WC)	
Western Manifold	1220	191			and the second second second	
SVE – 13	1218	183	3	0.8	0.01	
SVE – 12	1216	613	6	1.10	0.03	-
SVE – 11	1214	90	3	0,5	0.03	9.0
SVE – 10	1212	764	()	1.10	0.02	1
SVE – 09	1210	118	10	1.6	0.04	nd
Eastern manifold		_		110	0/01	0.7
SVE - 01				The third of the Article		
SVE - 03	not	IN DOLGIA FED	hup to look it	in the second second	0	
SVE - 05	and	La horitona	hand blood	eg weeke	ina	
SVE - 07	_ and	1 acong vry	A A A A A A A A A A A A A A A A A A A	per chi		
SVE - 08			A IOD	- 94		
SVE - 06		AV	0			
SVE - 04						
SVE - 02	/					
SVE Influent	2.12	1917	P. State State State State State	and the state of the state of the		
SVE Effluent	221	6.9				

Effluent Sample Time: 1220 Influent Sample Time: 35110 Field Representative (Print and Sign): Sey Davidge Date of Visit: 8/31/2018

#### Visual/Audio Inspection

Kelly Moore 09-19-18 5400 AZEPORT WAY SD. SEATTLE, WA 98108

VIOLAND MICHAELE	Inspected	Condition (Cracks, leaks, non-operational
Item	(Y/N)	gauges, etc.)
Above Ground Piping	yes	
Control Pump (Regenerative Blower)	ines	(On / Off) ON
Entrainment Pump (Transfer Pump)	yes_	(Auto / Hand / Off) Auto
Pressure Gauges/Flow Meters	yes_	PI-1is "of Hg-Voc, Not "H2D Voc.
Knockout Tank (record level)	للركال	% full 5% on 3 - 3 - 3
Knockout Water Tote (record level)	110 22	% full 90% Jull @ 110 get - 250 gal Mox,
Dilution Valve Status	·yes	open approx 50% - mehanged
Recirculation Valve Status	400	Open ? 1. ununour - uncharged

#### **CATOX Screen Readings**

lterre	Peoding	Operating
	neading	
Catox In (T1)	651°F	>650%
Catox Out (T <sub>2</sub> )	645 °F	600 – 650 °F
Heat Ex (T <sub>3</sub> )	353 °F	300 - 400 °F
Gt Flowerm	\$3 SCFM	<300 SCFM
LEL	13 *	5-15 %

#### System Gauge Readings

Item	Reading	
FE – 1	3.0 "WC	
PI-1 1.0"#a	(vacuum) 📌	""5
TI-1	ie38. ⁰F	
FE-2	1.5 "WC	]

#### FID Measurements

≁

		FID B	ading	Valve Position	Vacuum	Differential Pressure
Location	Time	FIP (PI	m) <del>303</del> -	(record notch)	("WC)	("WC)
Western Manifold	1100 400	321	740			
SVE - 13	1127/10	417	246	100 to open death	12.2"uc	0.00000
SVE - 12	1123 tha	1005	629	100"/ . apan - 6	2.3 WC	0,03 "WC_
SVE-11	112040	112	<b>8</b> 9	50% open -3	15,9"00	0,00"WC
SVE - 10	11 (4) 445	1204	589	100% apon - 6	Ooil "we	0.00"wc_
SVE - 09	1106#125	253	164	100% open - le	0.0 WC	0-05 "టడ
Eastern manifold	1226 1125	340	169			
SVE - 01	1229 1185	12.7	12.1	50% open - 3	0.0 "we	0.48"WC_
SVE - 03	1235 HRS	FLAME	1140	60% epers - 4	0.0 402	0.00 WC
SVE-05	1241 NR5	FIRGE	1061	60% open - 4	1.0 Hwc_	0.00 " WC
SVE - 07	1248 #0	98	93	75%000 - 4.5	26.5°wc	0,00° we
SVE - 08	1250 125	64	SI	75% open -4,5	0.05 WC	0,00" WC
SVE - 06	1258425	109	106.5	60% cars - 3	1.00 'wc	0.00 " WC
SVE - 04	1301 HES	313	226	50/ 0 AM - 3	0.08 "wc.	0.00° W C
SVE - 02	1304 HRS	16.5	10.2	50% open -3	4.00" wc	0.00" 00
SVE Influent	1334	336	187		建高度。	
SVE Effluent	1330	17.7	16.2			

Influent Sample ID: <u>INF.091918</u> Influent Sample Time: <u>1410 HR5</u> Effluent Sample Time: <u>1355</u> George Hacand Field Representative (Print and Sign): <u>George Hacand</u> SVE-03- FID-FLAME OUT WARNING 2^7ZMES -NO FID READING SVE-05 FID - FLAME OUT WARNING 207ZMES -NO FID READING

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A. T. O. T.	The second		SAMPL	ERS (signo	uture)									ר [	)	Page			<u>_</u>	<b>-</b> ,
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Company WOOD ENV	nonmentel		PROJE	CTNAME							P	)#		11		H		<u> </u>		
An Contrainer	Ť., < <del>†</del> . 5.,	To LAY	Kee	Lu Moore	2_										Rush (	charg	es autho	orized t	oy:	
Address (AL) Manuela	tag an i am		REMAR	IKS	<u> </u>	<u>.</u>				IN	voi	CE	TO	┥╞		SAM	PLE DI	SPOSA	T.	-
City, State, ZIP Seat	te WA 9811	01	-											1	2Disp	ose a	ifter 30 (	days		
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Sample ID <sub>ý</sub> :	Lab ID	Date Sampled	Time Sampled	Sample Type	# of Jars	TPH-HCID	TPH-Diesel	TPH-Gasoline	BTEX by 80211	VOCs by 82600	SVOCs by 8270	PAHs 8270D SI	BENZENE					Notes	\$	
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3012 16 <sup>th</sup> Avenue West	Received by.	T	- <u>*</u>	Mar	1	7		به مر بر	ζ_			Ē	151	7	-	-	9/19	1181	530	۱ ۱
Seattle, WA 98119-2029	Relinquished by:	* <u> </u>		- Mai	<i></i>	~	ng j	16.1.	22	•	<b></b>	•			-		1			1
Ph. (206) 285-8282	Received by:				. <u> </u>															
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Visual/Audio	Inspection
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ltem	Inspected (Y/N)	Condition (Cracks, leaks, non-operational gauges, etc.)
Above Ground Piping	Y	NO ISSNEZ
Control Pump (Regenerative Blower)	Y	(On) / Off)
Entrainment Pump (Transfer Pump)	Ý	(Auto / Hand / Off)
Pressure Gauges/Flow Meters	γ	All working
Knockout Tank (record level)	Y	% full
Knockout Water Tote (record level)	Ý	% full < \$'%
Dilution Valve Status	Y	CH? T
Recirculation Valve Status	Y	Open

#### CATOX Screen Readings

Readi	ing	Operating Range
660	٥F	>650°F
653	۰F	600 – 650 °F
348	٩F	300 - 400 °F
734 SC	FM	<300 SCFM
1-	%	5-15 %
	Readi 660 663 348 736 SC	Reading           660         °F           663         °F           348         °F           348         °F           348         °F           348         °F           348         °F           348         °F

#### System Gauge Readings

Item	Reading
FE – 1	27 "WC
PI – 1	O "WC (vacuum)
TI – 1	50 °F
FE-2	1.3 "WC

#### **FID Measurements**

Location	Time	FID Reading (ppm)	Valve Position (record notch)	Vacuum ("WC )	Differential Pressure ("WC )
Western Manifold	1925	273,4			
SVE – 13	1020	409.5	3	20.1	9.6
SVE – 12	1016	824.4.	6	4.9	0.5
SVE – 11	1012	166.0	3	22.5	9.G
SVE - 10	1004	826.5	5	2.1	WER4 0.2
SVE - 09	0.955	154.2	<b>Ø</b> 6	1.6.61.9	were 0.3
Eastern manifold	1030	54.2			
SVE - 01	1034	58.4	- 3	0.5	0.1
SVE - 03	1038	2542	36	0.5	0.6
SVE - 05	1042	2093	3-6	0.5	0.3
SVE - 07	1045	13.5	3	. ,	0.7
SVE - 08	10219	17.1	3	1.0	Da
SVE - 06	1063	8.3	3	2.5	0.8
SVE - 04	1067	8.2	3	1.0	0.5
SVE - 02	102	9.7	Э	3.0	Q.VX
SVE Influent	1107	58.4			
SVE Effluent	UIN	11-1			

Influent Sample ID: $\frac{1}{115}$  $\frac{1}{125}$ Effluent Sample ID: $\frac{1}{125}$ Influent Sample Time: $\frac{1}{125}$  $\frac{1}{125}$ 

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Field Representative (Print and Sign): William Young Date of Visit: 10-16-16

#### Melly Moore 4/112019 SVE System Monthly Inspection Log

Visual/Audio Inspection SVE - CATOX RESTMETED TODAY @ 1100-HES

ltem	Inspected (Y/N)	Condition (Cracks, leaks, non-operational gauges, etc.)
Above Ground Piping	13	it- Vacuum Gauges on Manifold are broken.
Control Pump (Regenerative Blower)	- y	(On/Off) on in Anto
Entrainment Pump (Transfer Pump)	y y	(Auto/Hand/Off) operational us Auto.
Pressure Gauges/Flow Meters		Vacuum Courses need replaced at least -4.
Knockout Tank (record level)	4	% full w.
Knockout Water Tote (record level)		% full (0 %,
Dilution Valve Status	yes	100% CLOSED
Recirculation Valve Status	njen	50'1, open

#### CATOX Screen Readings

ltem	Reading	Operating Range
Catox In (T1)	650 °F	>650ºF
Catox Out (T2)	633 °F	600 – 650 °F
Heat Ex (T <sub>3</sub> )	342 °F	300 – 400 °F
Flow	SCFM	<300 SCFM
LEL	8 *	5-15 %

#### System Gauge Readings

item	Reading
FE - 1	2.25 "WC
PI-1	Uto "WC (vacuum)
TI – 1	52 °F
FE-2	0, 4 WC

# PROVED AIR SPARCE ALARM CONTROLS LOW PRESSURE CUTOUT & 0.0 PSI, HIGH PRESSURE CUTOUT & 23,0 PSI

#### FID Measurements

HIGH TEMP CUT OUT C 154 F GETTING) NOT PROVOD, Differential Pressure Valve Position Vacuum **FID Reading** ("WC) ("WC) (record notch) (ppm) Time Location 建汽车部 NIA NA Western Manifold NIA NIA SVE - 13 NIA SVE - 12 SVE - 11 SVE - 10 SVE - 09 Eastern manifold Щ. N/A-N ) A-SVE - 01 NIA SVE - 03 SVE - 05 SVE - 07 SVE - 08 SVE - 06 SVE - 04 SVE - 02 1205 1138 PPM SVE Influent -1100-C1200 - 9.0 PPM SVE Effluent - 561 -0 Effluent Sample ID: Nonse Traces NONE TAXEN Influent Sample ID: \_ Effluent Sample Time: \_

Field Representative (Print and Sign):]

Influent Sample Time:

\_\_\_\_ Date of Visit: 04-01-2019

### SVE System Monthly Inspection Log 04-03 - 2019

lion	Inspected (Y/N)	Condition (Cracks, leaks, non-operational gauges, etc.)
	<u> </u>	A strange Course of It Anserous Beoren
Above Ground Piping	<u> </u>	4- VACIUM DAUGES ON VE MANLFOOD, DECKON
Control Pump (Regenerative Blower)	yes	(On / Off)
Entrainment Pump (Transfer Pump)	485	(Auto / Hand / Off) ANTO
Pressure Gauges/Flow Meters	YE5	
Knockout Tank (record level)	<u>, 4€5</u>	% full -0
Knockout Water Tote (record level)	455	% full 10'1
Dilution Valve Status	455	100%. CLOSED
Recirculation Valve Status	455	100%. CLOSED

CATOX Screen Readings @iz.3	SG- His
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Item	Reading	Operating Range
Catox In (T1)	664 °F	>650°F
Catox Out (T <sub>2</sub> )	630 °F	600 – 650 °F
Heat Ex (T <sub>3</sub> )	337 °F	300 – 400 °F
Flow	6,61 SCFM	<300 SCFM
	4 %	5-15 %
HOURME	TER 69	ا د

System Gauge Readings @ 1236 423

OTOLOTT GRANGE		
Item	Reading	
FE – 1	4.99 WC	
PI – 1	1,25 "₩ <del>C</del> (vac <u>uum)</u> ≤ <b>4</b>	чŚ
Tl – 1	51, <u>°F</u>	
FE-2	1.0 "WC	

FID Measurements VALUE RECORDED FROM SEROLING LED DISPLAY - # 4 Spot.

Location	Time	FID Reading (ppm)	Valve Position (record notch)	Vacuum ("WC)	Differential Pressure (''WC)
Western Manifold	1258.11	\$7.0PPM			
SVE - 13	j256-Ha	3.4 PPM	50% - man - 4	14.5 Baover	0.023"
SVE - 12	1254-144	4219 PPM_	100°/100 -7-	5,5	0.011
SVE - 11	1253-14	19.2.PPM	Schopen-4	21"- Broken	0.034
SVE - 10	1251-44	232,0 PPM	100% open - 7	5.75"	<u>0-003"</u>
SVE 09	1250-Hu	162.1 PPM	100 / 10 per - 7	1.0 4	0.040
Eastern manifold	1304-13	204.4 PPM			
SVE - 01	1305-113	3.6 PPM	50% open -4	2,5"	0.036
SVE - 03	1307-144	1494. 191	50% open -4	3.5"	4,380"
SVE - 05	1309-110	3222, 884	leDil. open - 5	2.25	0.001
SVE - 07	1311-145	66.8 PPM	75%, open - 6	730" Broken	0,001
SVE - 08	1313-HAZ.	57.7 ppm	75% open - le	4,5	0.002"
SVE - 06	1315-140	401.7 PPM	60% spin - 5	4.01	0.002"
SVE - 04	1316.443	54.3 PPM_	SO% open - H	4,5"	0.003
SVE 02	1318-143	4.4 ppm	Soll open	5,25"	0.032"
SVE Influent	1151	165 PPM			
SVE Effluent	1139	1.1 PPM			

Influent Sample ID: <u><i>INP</i></u>	-040319 1157 HES	2-TUXOLAR- - BAGS -	E
·····		C . 14	

Effluent Sample ID: <u>EPF - 040319</u> 2 - TEDLAR · Effluent Sample Time: <u>1143 - HeS</u> BAGS

\_\_\_\_ Date of Visit: <u>04.03-2019</u> Field Representative (Print and Sign): George Hagen

Visual/Audio inspection (	LA INDORED	Jon division (Ornalia Jacka non-operational
Item	Vinspected (Y/N)	gauges, etc.)
Above Ground Piping	حعرب	
Control Pump (Regenerative Blower)	yes	(On / Off) 0 N
Entrainment Pump (Transfer Pump)	yes	(Auto / Hand / Off)
Pressure Gauges/Flow Meters	yes	SEVERAL VACUUM · MANIFOLD GAUGES BROKED.
Knockout Tank (record level)	wes	% full 40 /,
Knockout Water Tote (record level)	yes	% full 10'/.
Dilution Valve Status	yes	100% Closed
Recirculation Valve Status	wes	100%. Closed

#### CATOX Screen Readings

		Operating
ltem	Reading	Range
Catox In (T1)	<i>`</i> 455`°₽	>650ºF
Catox Out (T <sub>2</sub> )	660 °F	600 – 650 °F
Heat Ex (T <sub>3</sub> )	350 F	300 – 400 °F
Flow	191 SCFM	<300 SCFM
	3 *	5-15 %

#### System Gauge Readings

ltem	Reading	
FE – 1	? "WC	
PI – 1	1 "We <sup>c</sup> ?	d q
	(vacuum)	
TI-1	53 °F	
FE-2	1 "WC	

#### How Mater 788.62 Hrs

#### FID Measurements

	Time	FID Reading	Valve Position (record notch)	Vacuum ("WC)	Differential Pressure ("WC)
Western Manifold	1233 His	2.2 PPM			
SVE - 13	1244	1-8 PPM	4	<u>9.5°</u> wc	o. o. o.
SVE - 12	12.2	12.3 PPM	<b>1</b>	6" WC	0.02" WC
SVE - 11	1240	3.5P1M	4	Arinen - 23"	0.03400
SVE - 10	1238	8.5PPM	·	4.5" wc_	0,02" 200
SVE - 09	1236	1.9PPM_	7. 100%, 072	4 "wc	0.05 to D
Eastern manifold	1308-141	4.9 PPM			
SVE - 01	1310	0.0PPM		2.50° UC	0.05 " WC
SVE - 03	1312	21,5 PPM_	5	3.25 "42	0.01 0.0
SVE - 05	1314	144.2 PPM	5	4.50"24	0.009 "WC
SVE - 07	1316	4.5 444	6	730" Beoken	0,009 "WC
SVE-08	1318	3,2 PPM	6	4.50 "We	0.006 WC
SVE - 06	1320	I.O PPM	6	4.50 "WC_	0.01 WC
SVE - 04	1322	0.0 PPM	<u> </u>	3,50"WC	0.01"00
SVE - 02	1324	0.0 PPM	5	5.50"WE	0.05 "22
SVE Influent	1030	16,1 PPM			
SVE Effluent	1025	\$9.1 PPM			

Influent Sample ID: INF - 050319 Influent Sample Time: 1030 - HRS

Effluent Sample ID: EFF -050319 Effluent Sample Time: 1025 - HES

Field Representative (Print and Sign): George Hagan Date of Visit: 5/3/2019

Visual/Audio Inspection Kerve	MOORE	, SEATTLE, 5400 AIRPORT WAY YOUTH.
	Inspected	Condition (Cracks, leaks, non-operational
Item	(Y/N)	gauges, etc.)
Above Ground Piping	yes	
Control Pump (Regenerative Blower)	yes	(On) Off)
Entrainment Pump (Transfer Pump)	yes	(Auto / Hand / Off)
Pressure Gauges/Flow Meters	yes	C Manifold's figuras Broken
Knockout Tank (record level)	5823	% full 5 % Full
Knockout Water Tote (record level)	ues	% full 10 /. Fall
Dilution Valve Status	iges	100% closed
Recirculation Valve Status	yes	100% Closed

#### CATOX Screen Readings

Item	Reading	Operating Range
Catox In (T1)	640 °F	>650°F
Catox Out (T <sub>2</sub> )	606 °F	600 – 650 °F
Heat Ex (T <sub>3</sub> )	343 °F	300 – 400 °F
Flow	187 SCFM	<300 SCFM
LEL	2 %	5-15 %

#### System Gauge Readings

Hour Meter 860.02 Hom.

Item	Reading
FE – 1	"WC
PI – 1	2 (vacuum) H
TI – 1	59 °F
FE-2	O, q "WC

Maisture K.O. Tonce Runphead + wige Strainer, Calibrated the LEL Sensor, Replaced FID Measurements Voc's Via FZA. Recorded System Date, Monitored op's, all OK House Keeping done around area.

Location	cation Time (ppm) Valve Position (record notch)		Position d notch)	Vacuum (''WC )		Pressure (''WC )				
Western Manifold	N	A	N	A						
SVE - 13				/	N	1A	N	IA	N	14
SVE - 12						/				
SVE - 11										
SVE - 10										
SVE - 09										
Eastern manifold					小小小				and a second	
SVE - 01		1								
SVE - 03										
SVE - 05										
SVE - 07										
SVE - 08										
SVE - 06										
SVE - 04										
SVE - 02		1		7	2			at		4-
SVE Influent	130	-3	15	.4PPM						
SVE Effluent	13:	20	0	· 1 PPM					1	

 Influent Sample ID:
 No
 Sample

 Influent Sample Time:
 Influent Sample Time:
 Influent Sample Time:

 Field Representative (Print and Sign):
 George Hacked
 Date of Visit:

 Server Hacked
 George Hacked

# SVE System Monthly Inspection Log Kelly more 05-28-19 - Alorn response Visual/Audio Inspection (130 the . System of Confe annual Restanted & monitored.

Item	Inspected (Y/N)	Condition (Cracks, leaks, non-operational gauges, etc.)
Above Ground Piping	ises	2 3 TO 4 Voc Courges Bod.
Control Pump (Regenerative Blower)	yes	(Op// Off)
Entrainment Pump (Transfer Pump)	yes	(Auto) / Hand / Off)
Pressure Gauges/Flow Meters	yes	
Knockout Tank (record level)	yes	% full _3 '/.
Knockout Water Tote (record level)	yes	% full 10/,
Dilution Valve Status	yes	100%. Closed
Recirculation Valve Status	yes	100%. Closed

#### CATOX Screen Readings

Reading	Operating Range
613 °F	>650°F
585 °F	600 – 650 °F
341 ºF	300 – 400 °F
120 SCFM	<300 SCFM
1 %	5-15 %
	Reading ゆ13 °F 585 °F ふ41 °F 180 SCFM / <sup>%</sup>

#### System Gauge Readings

ltem	Reading	
FE – 1	4.30 WC	
PI – 1	(vacuum)	"they
TI – 1	€0° °F	7
FE-2	0,50 "WC	

How meter 437 1347.26

#### **FID Measurements**

Location	Time	FID Reading (ppm)	Valve Position (record notch)	Vacuum (''WC )	Differential Pressure (''WC)
Western Manifold	1	1			
SVE – 13			No manifold	dates too	cell :
SVE - 12			System Re	stent V	monitor,
SVE - 11			Ceitt		
SVE - 10		$\backslash$ /			
SVE - 09		$\backslash$			
Eastern manifold			the second states and		
SVE - 01		$\square$			
SVE – 03					
SVE - 05					
SVE - 07					
SVE - 08	/	$\left( \right)$			
SVE - 06					
SVE - 04					
SVE - 02					
SVE Influent					
SVE Effluent	1				
Influent Sample ID: _ Influent Sample Time	NJA 		Effluent Sample Effluent Sample	ID: <u>م إ</u> كم Time:	
Field Representative	(Print and S	Sign): Bunge Alag	Date o	f Visit: <u>७५-२</u>	8-19

ALE SPARCE START UP SHAKE DOWN. SVE System Monthly inspection LOG 05-29-2019 - Kelly Moore Solo Auport Way S, Visual/Audio Inspection Shate - Baytal & Lundray - WOOD. MARCY - Kelly Moore,

Item	Inspected (Y/N)	Condition (Cracks, leaks, non-operational gauges, etc.)
Above Ground Piping	yes	
Control Pump (Regenerative Blower)	yes_	((On) Off)
Entrainment Pump (Transfer Pump)	yes	(Auto) / Hand / Off)
Pressure Gauges/Flow Meters	نيعه	5- Gauges Replaced Today
Knockout Tank (record level)	yes	% full 3'1. Fuel
Knockout Water Tote (record level)	res	% full 10% Full
Dilution Valve Status	ues	100% llosed
Recirculation Valve Status	yes	100% closed

#### CATOX Screen Readings

		Operating
Item	Reading	Range
Catox In (T1)	617 °F	>650°F
Catox Out (T2)	580 °F	600 – 650 °F
Heat Ex (T <sub>3</sub> )	342 °F	300 – 400 °F
Flow	184 SCFM	<300 SCFM
LEL	1 %	5-15 %

#### System Gauge Readings

Item	Reading
FE – 1	4.6 "WC
PI-1	(vaouum)
TI – 1	<u>58 °</u> ₽
FE-2	0,4 "WC

#### FID Measurements

Location	Time	FID Reading (ppm)	Valve Position (record notch)	Vacuum (''WC )	Differential Pressure ("WC)	1536 PD readings
Western Manifold	0943	13.2 Pp1				see Lucisse
SVE - 13	0956	4,2 PPM	4	3ªwc	0.015	Log GOOK
SVE - 12	0954	47,1 ppm	7	S"WC	0.00	Nales
SVE - 11	0953	4,6		31000	0.025	4 /
SVE - 10	0951	54,5000	F	<u>5.5°ພc</u>	0,002	4 }
SVE - 09	0945	12,1990	<u> 구</u>	S"WC	3.03	
Eastern manifold	1052	20,9 PPM				
SVE - 01	1053	0,0 PPM		3ª wc	0,013	6.5 6
SVE - 03	1055	147.2PPM	5	3.75 WC	0.016	40,500
SVE - 05	1057	490 PPM	5	4tuc	0.006	42.0pp
SVE - 07	1101	2. 8 PPM		5 ° 0 C	0.048	6. ppm
SVE - 08	1102-	2.4 PPM_		4" WC_	0,005	6.1 ppm
SVE - 06	1004	11.0 PPM	5	3.60"WC	0,004	Tispr.
SVE - 04	1105	28,30 PPM	+4	4.50 04	0.005	9,5 ppm
SVE - 02	1100	0.2 994	4	4"WC	0.021	_ 11,2 ppv
SVE Influent	0940	21.6 PPM				
SVE Effluent	0938	0.0PP.4				
Influent Sample ID:			Effluent Sample Effluent Sample	ID: Time:	A  29~19	

Field Representative (Print and Sign): <u>Comparent date of the</u> Air Spange Chouser Started C1120 Hrs.

	Air	Sponge	Stanted @	1120 Hus				
	Read	ing - many	19.4 <sup>H~s</sup>			کنی	Tem	
Time	HSI 1000 9 PSP	- Asz	A53	ASY	ASS	Blower Temp 1 per	CCOLLER CCOLLER DE + Temp	
11 32	5.5 J SEFM PSI	5.0 1.0 SCFM PSZ	5   1.5	5./1.5	5.5 /1.0	175° D	79'8/215	
1325	6.25/2.25 Sefmi PSZ	4,75/1.0 SCFM / P5#	4,75 1,5 SCEM PSI	5 1. SCFM PSL	5 1.5 SEFA PSF	165 F O TEMP PSS	84'F 1.5 Tenp (psz	
1400 Hy	ZNE	REASOD AT	r Fear To	10 SCFM @	EACH W	REPOINT		LEL @1
415	REDUCE	D ANFLO	5 To 850	PM I AS	4 + 45 5	1		
		LINDSA	5 Tanzasio Ne	rees freein This	s Point	FORWARD		
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#### SVE System Monthly Inspection Log 5-30 19 Kelly More

Item	Inspected (Y/N)	Condition (Cracks, leaks, non-operational gauges, etc.)
Above Ground Piping	yes	yes
Control Pump (Regenerative Blower)	unes.	(On / Off) on
Entrainment Pump (Transfer Pump)	yey	(Auto / Hand / Off) Auto
Pressure Gauges/Flow Meters	ues	
Knockout Tank (record level)	ier	% full 5 %
Knockout Water Tote (record level)	yes	% full 10'/,
Dilution Valve Status	uses	100% Closed
Recirculation Valve Status	yes	100% Closed

#### **CATOX Screen Readings**

Item	Reading	Operating Range
Catox In (T1)	632 °F	>650°F
Catox Out (T2)	626 °F	600 – 650 °F
Heat Ex (T <sub>3</sub> )	362 ºF	300 – 400 °F
Flow	161 SCFM	<300 SCFM
LEL	3 %	5-15 %

#### System Gauge Readings

Item	Reading
FE – 1	4.3" "WC
PI – 1	パキ、S <sup>ペ</sup> "WC (vacuum)
TI – 1	59′°F
FE-2	0,4" "WC

# Air Sponge stort 5.29-19, 157 Sompling event Post Stortup,

HR METER 1393+16M @ 1134. HAS FID Measurements

Location	Time	FID Reading (ppm)	Valve Position (record notch)	Vacuum (''WC )	Differential Pressure (''WC )	Som pile
Western Manifold	0906	173, PPM			Colored Series	- and
SVE - 13	0902	235, PPM	7	5.0	0,074"wc	0904
SVE - 12	0858	371, PPM	7	8,5"	0.009 "wc	0900
SVE - 11	0854	99, PPM	7	5,60"	ONIZWC	0835
SVE - 10	0848	216, PPM	オ	7.25"	0,000 02	0 850
SVE - 09	0842	18,1 PPM	4	4,251	01022 WC	0844
Eastern manifold	0943	9.3 . PPM				
SVE - 01	0945	0.6 Ppm	2	0,50 00	-0,004"WC	0948
SVE - 03	0950	163,1 PPM	2	4,75"wc	0.013 WC	0952
SVE - 05	0954	393, PPM	2	4"wc	0.000"00	0956
SVE - 07	0958	2.5 PPM	2	715"WC	0,003"wc	1000
SVE - 08	1002	1.0	2	8"WC	-0.004 WC	1004
SVE - 06	1006	0.0 PPM	2	24 w C	- 0,003 We	1008
SVE - 04	1010	18.6 PPM	2	215"WC	- 0,002 WC	1012
SVE - 02	1014	0-0-PM	2	S"wc	- 0.002"wc	1016
SVE Influent	0834	144,3 PPM				
SVE Effluent	0820	3.0 ppm			Salah Manakan	

Influent Sample ID: <u>053019 - TN</u>F Influent Sample Time: <u>0835 - Has</u>

Effluent Sample ID: <u>053019-EFF</u> Effluent Sample Time: <u>0330 - Hiz</u>5

Field Representative (Print and Sign): George Hagan Date of Visit: 05-30-19

AIR SPARGE DATA ON BACK

POST COULING TEMP - BI'F PSI - 3,5 PSI

WELL	FLOW SCFM	PRESSURE PSI
ASPI	8,0	3,25
ASPZ	10,5	3,00
ASP3	10,5	3.00
4928	9,5	3,00
ASP5	9,5	2,75

#### SVE System Monthly Inspection Log. Kelly Moore. Date: 06-04-19

#### Visual/Audio Inspection. Located at; 5400 Airport Way South Seattle, WA

ltem	Inspected (Y/N)	Condition (Cracks, leaks, non-operational gauges, etc.)
Above Ground Piping	izes	
Control Pump (Regenerative Blower)	ies	(On / Off) ON
Entrainment Pump (Transfer Pump)	ues	(Auto / Hand / Off) Auto
Pressure Gauges/Flow Meters	ney	
Knockout Tank (record level)	yes	% full 5 /,
Knockout Water Tote (record level)	حعم	% full io '/,
Dilution Valve Status	yer	
Recirculation Valve Status	izes	

#### **CATOX Screen Readings**

Item	Units	Reading	Operating Range
Catox In (T1)	٩F	613	>650
Catox Out (T <sub>2</sub> )	٥F	610	600 - 650
Heat Ex (T <sub>3</sub> )	٩F	353	300 - 400
Flow	SCFM	154	<300
LEL	%	1	5-15
	24		10 m 1 2 3

0850 BLAPTEST FID 99.4 PAM

## FID Measurements

#### System Gauge Readings

ltem	Units	Reading
FE – 1	"WC	3.95"WC
PI – 1	"WC (vacuum)	14.5°wc
TI – 1	°F	40'F
FE-2	''WC	0.4" we

Location	Time	FID Reading (ppm)	Valve Position (record notch)	Vacuum ("WC )	Differential Pressure ("WC)	Somple
Western Manifold	0930	43-2 PPM			0.0221	( Und
SVE – 13	0951	52,1 PPM	7	5,5 "we	SH 5.652 DC	0953
SVE - 12	0946	82.5 PPM	7	5.75 WC	0.013 "WC	0949
SVE – 11	0942	32.6 PPM	7	Gilese	0.190 02	0944
SVE - 10	0938	120, PP 17	7	6"00	0.000"wc	0940
SVE - 09	0932	14,0PPM	4	1,75"WC	0.011 "We	0934
Eastern manifold	1049	88,9 PPM	Marine Marine			1
SVE - 01	1058	OLOPPW	2_	0.50"WC	0.004400	1102
>> SVE - 03 ·	1103	121, 8 PPM	GH 2 TO 3	4.50 We	0.012 "we	1105
🔹 SVE – 05 -	1107	446.0 PPM	GH 2 TO 4	4,50 me	01083"00	1109
SVE - 07	1112	3.4 PPM	2	9.00" we	0,007"00	1114 1405
SVE – 08 –	1115	2.0 PPM	2	9.00"we	0,000"WC	1117
SVE - 06 -	1119	D, O PPM	2	2,25"WC	0,002 4WC	1121
SVE - 04	1124	14,3 PPM	2	2,25"we	0,000 "wc	lale
SVE - 02	1128	0, OPEN	2	4.00 We	0.000 6	1129
SVE Influent	10901	50,4 PPM	and the second states			
SVE Effluent	0853	1.2 PPM				

Influent Sample ID: <u>INF-060419</u> Influent Sample Time: <u>0903</u>

Effluent Sample ID: <u>EFF - 060419</u> Effluent Sample Time: <u>0856</u>

Field Representative (Print and Sign): GEORGE HAGAN Date of Visit: 6-4-19

SUE-05 Value opened to notch of SUE-03 value opened to notch 3

Page 10f 2

#### AS System Monthly Inspection Log Visual/Audio Inspection

Item	Inspected (Y/N)	Condition (Cracks, leaks, non-operational gauges, etc.)
Above Ground Piping	yes	
Regenerative Blower	yes	(Auto / Hand / Off) Auto
Heat Exchanger	yes	(Auto / Hand / Off) Allo
Pressure Gauges/Flow Meters	yes	
Vent Valve Status	yes	open 80'/,

#### System Gauge Readings

Be	fore Hea	at Exchange	r	After	Heat Ex	changer		
ltem	Units	Reading	Operating Range	Item	Units	Reading	Operating Range	GH
PI – 3	psi	5.25 PSZ	0 TO 30	PI – 4	psi	4,25 FSZ	0-530	GH
TI – 3	٥F	163 F	150 - 200	TI – 4	٩F	79'F	150 - 200	

#### **Air Flow Monitoring**

Location	Time	Valve Position (record appx angle)	Pressure (psi)	Air Flow (SCFM )
AS – 1	1006 Hus	open 20%	3.0	8.0
AS – 2	1007 Hy	open 25%	2.75	10,50
AS – 3	1008 445	open 25%	3.0	10.50
AS – 4	10091445	open 20%	3.0	9,50
AS – 5	1010 Hrs	open 20%	2.75	9,50

Additional Notes.

See Log entry with today's date.

Field Representative (Print and Sign): George Hacan Date of Visit: 6-4-19

Page 20f2

#### SVE System Monthly Inspection Log. Kelly Moore. Date: 06-05-19 9

Item	Inspected (Y/N)	Condition (Cracks, leaks, non-operational U gauges, etc.)
Above Ground Piping	lyes	
Control Pump (Regenerative Blower)	yes	(On) Off)
Entrainment Pump (Transfer Pump)	Les	(Auto) Hand / Off)
Pressure Gauges/Flow Meters	ies	
Knockout Tank (record level)	nes	% full 3'. Ful
Knockout Water Tote (record level)	ner	% full 10 % Fell
Dilution Valve Status	Les	
Recirculation Valve Status	wet	

#### CATOX Screen Readings

Item	Units	Reading	Operating Range
Catox In (T1)	٩F	613	>650
Catox Out (T <sub>2</sub> )	٩F	623	600 - 650
Heat Ex (T <sub>3</sub> )	٥F	361	300 - 400
Flow	SCFM	160	<300
LEL	%	3	5-15

#### System Gauge Readings

Item	Units	Reading
FE – 1	"WC	3.71
PI – 1	"WC (vacuum)	14 " H2D
TI – 1	٥F	61'F
FE-2	"WC	0.03"420

FID Measurements LEL CALEBRATED TODAY,

Location	Time	FID Reading (ppm)	Valve Position (record notch)	Vacuum ("WC)	Differential Pressure (''WC )
Western Manifold			The second second	and the second second	
SVE – 13	No Ma	pelold dates	today.		
SVE – 12		NO FID	date targe		
SVE – 11			and weight		
SVE - 10					
SVE - 09					
Eastern manifold					
SVE - 01					
SVE - 03					
SVE - 05					
SVE - 07					
SVE - 08					
SVE - 06					
SVE - 04					
SVE - 02					
SVE Influent			The second second		AND STREET
SVE Effluent					

Influent Sample ID: N A	Effluent Sample ID: N//
Influent Sample Time:	Effluent Sample Time:
Gun	are those in 1
Field Representative (Print and Sign):	Date of Visit: 06/05 / 2019

#### AS System Monthly Inspection Log Visual/Audio Inspection

Condition (Cracks, leaks, non-operationa Inspected Item (Y/N) gauges, etc.) Above Ground Piping yes **Regenerative Blower** (Auto / Hand / Off) yes Heat Exchanger (Auto / Hand / Off) Pressure Gauges/Flow Meters ues openves Vent Valve Status open go/ HR. METER. HEAT EXCHANCER

the moter AIR Sparce Pump 189.7/10

Reading

5.0

165 F

System Gauge Readings Before Heat Exchanger

Units

psi

٩F

After Heat Exchanger

Operating Range	ltem	Units	Reading	Operating Range
0	PI – 4	psi		0 - 5
150 – 200	TI – 4	٩F	84'F	150 – 200

#### Air Flow Monitoring

Item

PI - 3

TI - 3

Location	Time	Valve Position (record appx angle)	Pressure (psi)	Air Flow (SCFM )
AS – 1	1465	20'/, open	3.0	8,25
AS – 2	1416	25% open	2.5	10.75
AS – 3	1417	25% open	3,0	10,50
AS – 4	1418	25% open	3,0	9.75
AS – 5	1419	zo'l. open	2.5	9,75

Additional Notes.

0 eclod 8 Co umma samples Ware d build ing exterior one milke Monthly LEL Calibration. Change Con The ch econder per un

Date of Visit: 06-05-2019

Kelly Mone 06)05/19

190.24



# Appendix D

#### ENVIRONMENTAL CHEMISTS

James E. Bruya, Ph.D. Yelena Aravkina, M.S. Michael Erdahl, B.S. Arina Podnozova, B.S. Eric Young, B.S. 3012 16th Avenue West Seattle, WA 98119-2029 (206) 285-8282 fbi@isomedia.com www.friedmanandbruya.com

May 2, 2018

Crystal Thimsen, Project Manager Wood Environment & Infrastructure Solutions, Inc. One Union Square 600 University Street, Suite 600 Seattle, WA 98101

Dear Ms Thimsen:

Included are the results from the testing of material submitted on April 25, 2018 from the Kelley Moore 0146970060.00009, F&BI 804419 project. There are 4 pages included in this report.

We appreciate this opportunity to be of service to you and hope you will call if you should have any questions.

Sincerely,

FRIEDMAN & BRUYA, INC.

Michael Erdahl Project Manager

Enclosures AFW0502R.DOC

#### ENVIRONMENTAL CHEMISTS

#### CASE NARRATIVE

This case narrative encompasses samples received on April 25, 2018 by Friedman & Bruya, Inc. from the Wood Environment & Infrastructure Solutions Kelley Moore 0146970060.00009, F&BI 804419 project. Samples were logged in under the laboratory ID's listed below.

<u>Laboratory ID</u>	Wood Environment & Infrastructure Solutions
804419 -01	WF-042418
804419 -02	EFF-042418

All quality control requirements were acceptable.

#### ENVIRONMENTAL CHEMISTS

Date of Report: 05/02/18 Date Received: 04/25/18 Project: Kelley Moore 0146970060.00009, F&BI 804419 Date Extracted: 04/26/18 Date Analyzed: 04/26/18

#### RESULTS FROM THE ANALYSIS OF AIR SAMPLES FOR BENZENE AND TPH AS GASOLINE USING MODIFIED METHODS 8021B AND NWTPH-Gx

Sample ID Laboratory ID	<u>Benzene</u>	Gasoline <u>Range</u>	Surrogate ( <u>% Recovery)</u> (Limit 50-150)
WF-042418 804419-01	<0.1	80	82
EFF-042418 804419-02	<0.1	15	82
Method Blank 08-870 MB	<0.1	<10	82

Results Reported as mg/m<sup>3</sup>

#### ENVIRONMENTAL CHEMISTS

Date of Report: 05/02/18 Date Received: 04/25/18 Project: Kelley Moore 0146970060.00009, F&BI 804419

#### QUALITY ASSURANCE RESULTS FOR THE ANALYSIS OF AIR SAMPLES FOR BENZENE AND TPH AS GASOLINE USING MODIFIED EPA METHOD 8021B AND NWTPH-Gx

Laboratory Code:	804419-02 (Duplica	ate)		
	Reporting	Sample	Duplicate	RPD
Analyte	Units	Result	Result	(Limit 20)
Benzene	mg/m³	<0.1	<0.1	nm
Gasoline	mg/m <sup>3</sup>	15	17	12
	_			

Laboratory Code: Laboratory Control Sample

			Percent	
	Reporting	Spike	Recovery	Acceptance
Analyte	Units	Level	LCS	Criteria
Benzene	mg/m <sup>3</sup>	5.0	86	70-130
Gasoline	mg/m <sup>3</sup>	100	137	86-144

ENVIRONMENTAL CHEMISTS

#### **Data Qualifiers & Definitions**

a - The analyte was detected at a level less than five times the reporting limit. The RPD results may not provide reliable information on the variability of the analysis.

 ${\bf b}$  - The analyte was spiked at a level that was less than five times that present in the sample. Matrix spike recoveries may not be meaningful.

ca - The calibration results for the analyte were outside of acceptance criteria. The value reported is an estimate.

c - The presence of the analyte may be due to carryover from previous sample injections.

cf - The sample was centrifuged prior to analysis.

d - The sample was diluted. Detection limits were raised and surrogate recoveries may not be meaningful.

dv - Insufficient sample volume was available to achieve normal reporting limits.

f - The sample was laboratory filtered prior to analysis.

fb - The analyte was detected in the method blank.

fc - The compound is a common laboratory and field contaminant.

hr - The sample and duplicate were reextracted and reanalyzed. RPD results were still outside of control limits. Variability is attributed to sample inhomogeneity.

hs - Headspace was present in the container used for analysis.

ht – The analysis was performed outside the method or client-specified holding time requirement.

 ${\rm ip}$  - Recovery fell outside of control limits. Compounds in the sample matrix interfered with the quantitation of the analyte.

j - The analyte concentration is reported below the lowest calibration standard. The value reported is an estimate.

 ${\rm J}$  - The internal standard associated with the analyte is out of control limits. The reported concentration is an estimate.

jl - The laboratory control sample(s) percent recovery and/or RPD were out of control limits. The reported concentration should be considered an estimate.

js - The surrogate associated with the analyte is out of control limits. The reported concentration should be considered an estimate.

lc - The presence of the analyte is likely due to laboratory contamination.

L - The reported concentration was generated from a library search.

nm - The analyte was not detected in one or more of the duplicate analyses. Therefore, calculation of the RPD is not applicable.

pc - The sample was received with incorrect preservation or in a container not approved by the method. The value reported should be considered an estimate.

ve - The analyte response exceeded the valid instrument calibration range. The value reported is an estimate.

vo - The value reported fell outside the control limits established for this analyte.

x - The sample chromatographic pattern does not resemble the fuel standard used for quantitation.

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Limits of Liability: ARI will perform all requested services in accordance with appropriate methodology following ARI Standard Operating Procedures and the ARI Quality Assurance Program. This program meets standards for the industry. The total liability of ARI, its officers, agents, employees, or successors, arising out of or in connection with the requested services, shall not exceed the Invoiced amount for said services. The acceptance by the client of a proposal for services by ARI release ARI from any liability in excess thereof, not withstanding any provision to the contrary in any contract, purchase order or co-signed agreement between ARI and the Client.

Sample Retention Policy: All samples submitted to ARI will be appropriately discarded no sconer than 90 days after receipt or 60 days after submission of hardcopy data, whichever is longer, unless alternate retention schedules have been established by work-order or contract.

#### ENVIRONMENTAL CHEMISTS

James E. Bruya, Ph.D. Yelena Aravkina, M.S. Michael Erdahl, B.S. Arina Podnozova, B.S. Eric Young, B.S. 3012 16th Avenue West Seattle, WA 98119-2029 (206) 285-8282 fbi@isomedia.com www.friedmanandbruya.com

June 6, 2018

Crystal Thimsen, Project Manager Wood Environment & Infrastructure Solutions, Inc. One Union Square 600 University Street, Suite 600 Seattle, WA 98101

Dear Ms Thimsen:

Included are the results from the testing of material submitted on May 30, 2018 from the 14697, F&BI 805514 project. There are 4 pages included in this report. Any samples that may remain are currently scheduled for disposal in 30 days. If you would like us to return your samples or arrange for long term storage at our offices, please contact us as soon as possible.

We appreciate this opportunity to be of service to you and hope you will call if you should have any questions.

Sincerely,

FRIEDMAN & BRUYA, INC.

Cale

Michael Erdahl Project Manager

Enclosures

WEI0606R.DOC

#### ENVIRONMENTAL CHEMISTS

#### CASE NARRATIVE

This case narrative encompasses samples received on May 30, 2018 by Friedman & Bruya, Inc. from the Wood Environment & Infrastructure Solutions 14697, F&BI 805514 project. Samples were logged in under the laboratory ID's listed below.

Laboratory ID	Wood Environment & Infrastructure Solution
805514 -01	INF-053018
805514 -02	EFF-053018

All quality control requirements were acceptable.

#### ENVIRONMENTAL CHEMISTS

Date of Report: 06/06/18 Date Received: 05/30/18 Project: 14697, F&BI 805514 Date Extracted: 06/01/18 Date Analyzed: 06/01/18

#### RESULTS FROM THE ANALYSIS OF AIR SAMPLES FOR BENZENE AND TPH AS GASOLINE USING MODIFIED METHODS 8021B AND NWTPH-Gx

Sample ID Laboratory ID	<u>Benzene</u>	Gasoline <u>Range</u>	Surrogate ( <u>% Recovery)</u> (Limit 50-150)
INF-053018 805514-01	<0.1	170	103
EFF-053018 805514-02	<0.1	160	102
Method Blank 08-1136 MB2	<0.1	<10	89

Results Reported as mg/m<sup>3</sup>
### ENVIRONMENTAL CHEMISTS

Date of Report: 06/06/18 Date Received: 05/30/18 Project: 14697, F&BI 805514

### QUALITY ASSURANCE RESULTS FOR THE ANALYSIS OF AIR SAMPLES FOR BENZENE AND TPH AS GASOLINE USING MODIFIED EPA METHOD 8021B AND NWTPH-Gx

Laboratory Code:	805501-01 (Duplica	ate)		
	Reporting	Sample	Duplicate	RPD
Analyte	Units	Result	Result	(Limit 20)
Benzene	mg/m³	<0.1	<0.1	nm
Gasoline	mg/m <sup>3</sup>	<10	<10	nm

Laboratory Code: Laboratory Control Sample

			Percent	
	Reporting	Spike	Recovery	Acceptance
Analyte	Units	Level	LCS	Criteria
Benzene	mg/m <sup>3</sup>	5.0	95	70-130
Gasoline	mg/m <sup>3</sup>	100	109	70-130

ENVIRONMENTAL CHEMISTS

# **Data Qualifiers & Definitions**

a - The analyte was detected at a level less than five times the reporting limit. The RPD results may not provide reliable information on the variability of the analysis.

 ${\bf b}$  - The analyte was spiked at a level that was less than five times that present in the sample. Matrix spike recoveries may not be meaningful.

ca - The calibration results for the analyte were outside of acceptance criteria. The value reported is an estimate.

c - The presence of the analyte may be due to carryover from previous sample injections.

cf - The sample was centrifuged prior to analysis.

d - The sample was diluted. Detection limits were raised and surrogate recoveries may not be meaningful.

dv - Insufficient sample volume was available to achieve normal reporting limits.

f - The sample was laboratory filtered prior to analysis.

fb - The analyte was detected in the method blank.

fc - The compound is a common laboratory and field contaminant.

hr - The sample and duplicate were reextracted and reanalyzed. RPD results were still outside of control limits. Variability is attributed to sample inhomogeneity.

hs - Headspace was present in the container used for analysis.

ht – The analysis was performed outside the method or client-specified holding time requirement.

ip - Recovery fell outside of control limits. Compounds in the sample matrix interfered with the quantitation of the analyte.

j - The analyte concentration is reported below the lowest calibration standard. The value reported is an estimate.

 ${\rm J}$  - The internal standard associated with the analyte is out of control limits. The reported concentration is an estimate.

jl - The laboratory control sample(s) percent recovery and/or RPD were out of control limits. The reported concentration should be considered an estimate.

js - The surrogate associated with the analyte is out of control limits. The reported concentration should be considered an estimate.

lc - The presence of the analyte is likely due to laboratory contamination.

L - The reported concentration was generated from a library search.

nm - The analyte was not detected in one or more of the duplicate analyses. Therefore, calculation of the RPD is not applicable.

pc - The sample was received with incorrect preservation or in a container not approved by the method. The value reported should be considered an estimate.

ve - The analyte response exceeded the valid instrument calibration range. The value reported is an estimate.

vo - The value reported fell outside the control limits established for this analyte.

x - The sample chromatographic pattern does not resemble the fuel standard used for quantitation.

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Limits of Liability: ARI will perform all requested services in accordance with appropriate methodology following ARI Standard Operating Procedures and the ARI Quality Assurance Program. This program meets standards for the industry. The total liability of ARI, its officers, agents, employees, or successors, arising out of or in connection with the requested services, shall not exceed the Invoiced amount for said services. The acceptance by the client of a proposal for services by ARI release ARI from any liability in excess thereof, not withstanding any provision to the contrary in any contract, purchase order or co-signed agreement between ARI and the Client.

Sample Retention Policy: All samples submitted to ARI will be appropriately discarded no sconer than 90 days after receipt or 60 days after submission of hardcopy data, whichever is longer, unless alternate retention schedules have been established by work-order or contract.

#### ENVIRONMENTAL CHEMISTS

James E. Bruya, Ph.D. Yelena Aravkina, M.S. Michael Erdahl, B.S. Arina Podnozova, B.S. Eric Young, B.S. 3012 16th Avenue West Seattle, WA 98119-2029 (206) 285-8282 fbi@isomedia.com www.friedmanandbruya.com

July 10, 2018

Crystal Thimsen, Project Manager Wood Environment & Infrastructure Solutions, Inc. One Union Square 600 University Street, Suite 600 Seattle, WA 98101

Dear Ms Thimsen:

Included are the results from the testing of material submitted on June 29, 2018 from the Kelly Moore 14697, F&BI 806572 project. There are 4 pages included in this report. Any samples that may remain are currently scheduled for disposal in 30 days. If you would like us to return your samples or arrange for long term storage at our offices, please contact us as soon as possible.

We appreciate this opportunity to be of service to you and hope you will call if you should have any questions.

Sincerely,

FRIEDMAN & BRUYA, INC.

Cale

Michael Erdahl Project Manager

Enclosures WEI0710R.DOC

## ENVIRONMENTAL CHEMISTS

## CASE NARRATIVE

This case narrative encompasses samples received on June 29, 2018 by Friedman & Bruya, Inc. from the Wood Environment & Infrastructure Solutions Kelly Moore 14697, F&BI 806572 project. Samples were logged in under the laboratory ID's listed below.

Laboratory ID	Wood Environment & Infrastructure Solution
806572 -01	INF062918
806572 -02	EFF062918

All quality control requirements were acceptable.

#### ENVIRONMENTAL CHEMISTS

Date of Report: 07/10/18 Date Received: 06/29/18 Project: Kelly Moore 14697, F&BI 806572 Date Extracted: 07/02/18 Date Analyzed: 07/02/18

### RESULTS FROM THE ANALYSIS OF AIR SAMPLES FOR BENZENE, TOLUENE, ETHYLBENZENE, XYLENES AND TPH AS GASOLINE USING MODIFIED METHODS 8021B AND NWTPH-Gx

Results Reported as mg/m<sup>3</sup>

<u>Sample ID</u> Laboratory ID	<u>Benzene</u>	Gasoline <u>Range</u>	Surrogate ( <u>% Recovery</u> ) (Limit 50-150)
INF 062918 806572-01	<0.1	530	107
EFF062918 806572-02	<0.1	520	103
Method Blank 08-1389 MB	<0.1	<10	75

#### ENVIRONMENTAL CHEMISTS

Date of Report: 07/10/18 Date Received: 06/29/18 Project: Kelly Moore 14697, F&BI 806572

### QUALITY ASSURANCE RESULTS FOR THE ANALYSIS OF AIR SAMPLES FOR BENZENE, TOLUENE, ETHYLBENZENE, XYLENES, AND TPH AS GASOLINE USING MODIFIED EPA METHOD 8021B AND NWTPH-Gx

Laboratory Code:	806572-02 (Duplica	ate)		
-	Reporting	Sample	Duplicate	RPD
Analyte	Units	Result	Result	(Limit 20)
Benzene	mg/m <sup>3</sup>	< 0.1	<0.1	nm
Gasoline	mg/m <sup>3</sup>	520	530	2

Laboratory Code: Laboratory Control Sample

			Percent	
	Reporting	Spike	Recovery	Acceptance
Analyte	Units	Level	LCS	Criteria
Benzene	mg/m³	5.0	81	70-130
Gasoline	mg/m <sup>3</sup>	100	137	86-144

ENVIRONMENTAL CHEMISTS

# **Data Qualifiers & Definitions**

a - The analyte was detected at a level less than five times the reporting limit. The RPD results may not provide reliable information on the variability of the analysis.

 ${\bf b}$  - The analyte was spiked at a level that was less than five times that present in the sample. Matrix spike recoveries may not be meaningful.

ca - The calibration results for the analyte were outside of acceptance criteria. The value reported is an estimate.

c - The presence of the analyte may be due to carryover from previous sample injections.

cf - The sample was centrifuged prior to analysis.

d - The sample was diluted. Detection limits were raised and surrogate recoveries may not be meaningful.

dv - Insufficient sample volume was available to achieve normal reporting limits.

f - The sample was laboratory filtered prior to analysis.

fb - The analyte was detected in the method blank.

fc - The compound is a common laboratory and field contaminant.

 $hr\ \text{-}\ The\ sample\ and\ duplicate\ were\ reextracted\ and\ reanalyzed.\ RPD\ results\ were\ still\ outside\ of\ control\ limits.\ Variability\ is\ attributed\ to\ sample\ inhomogeneity.$ 

hs - Headspace was present in the container used for analysis.

ht – The analysis was performed outside the method or client-specified holding time requirement.

 ${\rm ip}$  - Recovery fell outside of control limits. Compounds in the sample matrix interfered with the quantitation of the analyte.

j - The analyte concentration is reported below the lowest calibration standard. The value reported is an estimate.

 ${\rm J}$  - The internal standard associated with the analyte is out of control limits. The reported concentration is an estimate.

jl - The laboratory control sample(s) percent recovery and/or RPD were out of control limits. The reported concentration should be considered an estimate.

js - The surrogate associated with the analyte is out of control limits. The reported concentration should be considered an estimate.

lc - The presence of the analyte is likely due to laboratory contamination.

L - The reported concentration was generated from a library search.

nm - The analyte was not detected in one or more of the duplicate analyses. Therefore, calculation of the RPD is not applicable.

pc - The sample was received with incorrect preservation or in a container not approved by the method. The value reported should be considered an estimate.

ve - The analyte response exceeded the valid instrument calibration range. The value reported is an estimate.

vo - The value reported fell outside the control limits established for this analyte.

x - The sample chromatographic pattern does not resemble the fuel standard used for quantitation.

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#### ENVIRONMENTAL CHEMISTS

James E. Bruya, Ph.D. Yelena Aravkina, M.S. Michael Erdahl, B.S. Arina Podnozova, B.S. Eric Young, B.S. 3012 16th Avenue West Seattle, WA 98119-2029 (206) 285-8282 fbi@isomedia.com www.friedmanandbruya.com

August 6, 2018

Crystal Thimsen, Project Manager Wood Environment & Infrastructure Solutions, Inc. One Union Square 600 University Street, Suite 600 Seattle, WA 98101

Dear Ms Thimsen:

Included are the results from the testing of material submitted on July 31, 2018 from the Kelly Moore, F&BI 807615 project. There are 4 pages included in this report. Any samples that may remain are currently scheduled for disposal in 30 days. If you would like us to return your samples or arrange for long term storage at our offices, please contact us as soon as possible.

We appreciate this opportunity to be of service to you and hope you will call if you should have any questions.

Sincerely,

FRIEDMAN & BRUYA, INC.

Cale

Michael Erdahl Project Manager

Enclosures WEI0806R.DOC

## ENVIRONMENTAL CHEMISTS

## CASE NARRATIVE

This case narrative encompasses samples received on July 31, 2018 by Friedman & Bruya, Inc. from the Wood Environment & Infrastructure Solutions Kelly Moore, F&BI 807615 project. Samples were logged in under the laboratory ID's listed below.

Laboratory ID	Wood Environment & Infrastructure Solutions
807615 -01	EFF073118
807615 -02	INF073118

All quality control requirements were acceptable.

### ENVIRONMENTAL CHEMISTS

Date of Report: 08/06/18 Date Received: 07/31/18 Project: Kelly Moore, F&BI 807615 Date Extracted: 08/01/18 Date Analyzed: 08/01/18

### RESULTS FROM THE ANALYSIS OF AIR SAMPLES FOR BENZENE AND TPH AS GASOLINE USING MODIFIED METHODS 8021B AND NWTPH-Gx

Sample ID Laboratory ID	<u>Benzene</u>	Gasoline <u>Range</u>	Surrogate ( <u>% Recovery)</u> (Limit 50-150)
EFF073118 807615-01	<0.1	450	123
INF073118 807615-02	<0.1	730	139
Method Blank 08-1653 MB	<0.1	<10	90

Results Reported as mg/m<sup>3</sup>

#### ENVIRONMENTAL CHEMISTS

Date of Report: 08/06/18 Date Received: 07/31/18 Project: Kelly Moore, F&BI 807615

### QUALITY ASSURANCE RESULTS FOR THE ANALYSIS OF AIR SAMPLES FOR BENZENE AND TPH AS GASOLINE USING MODIFIED EPA METHOD 8021B AND NWTPH-Gx

Laboratory Code:	807615-01 (Duplica	ate)		
-	Reporting	Sample	Duplicate	RPD
Analyte	Units	Result	Result	(Limit 20)
Benzene	mg/m <sup>3</sup>	<0.1	< 0.1	nm
Gasoline	mg/m <sup>3</sup>	450	460	3
Laboratory Code:	Laboratory Contro	l Sample		

-	-		Percent	
	Reporting	Spike	Recovery	Acceptance
Analyte	Units	Level	LCS	Criteria
Benzene	mg/m³	5.0	96	70-130
Gasoline	mg/m <sup>3</sup>	100	125	86-144

ENVIRONMENTAL CHEMISTS

# **Data Qualifiers & Definitions**

a - The analyte was detected at a level less than five times the reporting limit. The RPD results may not provide reliable information on the variability of the analysis.

 ${\bf b}$  - The analyte was spiked at a level that was less than five times that present in the sample. Matrix spike recoveries may not be meaningful.

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cf - The sample was centrifuged prior to analysis.

d - The sample was diluted. Detection limits were raised and surrogate recoveries may not be meaningful.

dv - Insufficient sample volume was available to achieve normal reporting limits.

f - The sample was laboratory filtered prior to analysis.

fb - The analyte was detected in the method blank.

fc - The compound is a common laboratory and field contaminant.

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vo - The value reported fell outside the control limits established for this analyte.

x - The sample chromatographic pattern does not resemble the fuel standard used for quantitation.

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#### ENVIRONMENTAL CHEMISTS

James E. Bruya, Ph.D. Yelena Aravkina, M.S. Michael Erdahl, B.S. Arina Podnozova, B.S. Eric Young, B.S. 3012 16th Avenue West Seattle, WA 98119-2029 (206) 285-8282 fbi@isomedia.com www.friedmanandbruya.com

September 7, 2018

Crystal Thimsen, Project Manager Wood Environment & Infrastructure Solutions, Inc. One Union Square 600 University Street, Suite 600 Seattle, WA 98101

Dear Ms Thimsen:

Included are the results from the testing of material submitted on August 31, 2018 from the Kelly-Moore, F&BI 808717 project. There are 4 pages included in this report.

We appreciate this opportunity to be of service to you and hope you will call if you should have any questions.

Sincerely,

FRIEDMAN & BRUYA, INC.

Michael Erdahl Project Manager

Enclosures WEI0907R.DOC

## ENVIRONMENTAL CHEMISTS

## CASE NARRATIVE

This case narrative encompasses samples received on August 31, 2018 by Friedman & Bruya, Inc. from the Wood Environment & Infrastructure Solutions Kelly-Moore, F&BI 808717 project. Samples were logged in under the laboratory ID's listed below.

Laboratory ID	Wood Environment & Infrastructure Solutions
808717 -01	EFF-083118
808717 -02	INF-083118

All quality control requirements were acceptable.

### ENVIRONMENTAL CHEMISTS

Date of Report: 09/07/18 Date Received: 08/31/18 Project: Kelly-Moore, F&BI 808717 Date Extracted: 08/31/18 Date Analyzed: 08/31/18

### RESULTS FROM THE ANALYSIS OF AIR SAMPLES FOR BENZENE AND TPH AS GASOLINE USING MODIFIED METHODS 8021B AND NWTPH-Gx

Results Reported as mg/m<sup>3</sup>

<u>Sample ID</u> Laboratory ID	<u>Benzene</u>	Gasoline <u>Range</u>	Surrogate ( <u>% Recovery)</u> (Limit 50-150)
EFF-083118 808717-01	<0.1	23	89
INF-083118 808717-02	<0.1	550	131
Method Blank 08-1941 MB	<0.1	<10	87

### ENVIRONMENTAL CHEMISTS

Date of Report: 09/07/18 Date Received: 08/31/18 Project: Kelly-Moore, F&BI 808717

### QUALITY ASSURANCE RESULTS FOR THE ANALYSIS OF AIR SAMPLES FOR BENZENE AND TPH AS GASOLINE USING MODIFIED EPA METHOD 8021B AND NWTPH-Gx

Laboratory Code:	808717-01 (Duplica	ate)		
	Reporting	Sample	Duplicate	RPD
Analyte	Units	Result	Result	(Limit 20)
Benzene	mg/m³	< 0.1	< 0.1	nm
Gasoline	mg/m <sup>3</sup>	23	46	67 a
	-			

Laboratory Code: Laboratory Control Sample

		Percent	
Reporting	Spike	Recovery	Acceptance
Units	Level	LCS	Criteria
mg/m <sup>3</sup>	5.0	90	70-130
mg/m <sup>3</sup>	100	128	86-144
	Reporting Units mg/m <sup>3</sup> mg/m <sup>3</sup>	Reporting UnitsSpike Levelmg/m³5.0 mg/m³	Reporting UnitsSpike LevelRecovery LCSmg/m³5.090mg/m³100128

#### ENVIRONMENTAL CHEMISTS

# **Data Qualifiers & Definitions**

a - The analyte was detected at a level less than five times the reporting limit. The RPD results may not provide reliable information on the variability of the analysis.

 ${\bf b}$  - The analyte was spiked at a level that was less than five times that present in the sample. Matrix spike recoveries may not be meaningful.

ca - The calibration results for the analyte were outside of acceptance criteria. The value reported is an estimate.

c - The presence of the analyte may be due to carryover from previous sample injections.

cf - The sample was centrifuged prior to analysis.

d - The sample was diluted. Detection limits were raised and surrogate recoveries may not be meaningful.

dv - Insufficient sample volume was available to achieve normal reporting limits.

f - The sample was laboratory filtered prior to analysis.

fb - The analyte was detected in the method blank.

fc - The compound is a common laboratory and field contaminant.

 $hr\ \text{-}\ The\ sample\ and\ duplicate\ were\ reextracted\ and\ reanalyzed.\ RPD\ results\ were\ still\ outside\ of\ control\ limits.\ Variability\ is\ attributed\ to\ sample\ inhomogeneity.$ 

hs - Headspace was present in the container used for analysis.

ht – The analysis was performed outside the method or client-specified holding time requirement.

 $\ensuremath{\text{ip}}$  - Recovery fell outside of control limits. Compounds in the sample matrix interfered with the quantitation of the analyte.

j - The analyte concentration is reported below the lowest calibration standard. The value reported is an estimate.

 ${\rm J}$  - The internal standard associated with the analyte is out of control limits. The reported concentration is an estimate.

jl - The laboratory control sample(s) percent recovery and/or RPD were out of control limits. The reported concentration should be considered an estimate.

js - The surrogate associated with the analyte is out of control limits. The reported concentration should be considered an estimate.

lc - The presence of the analyte is likely due to laboratory contamination.

L - The reported concentration was generated from a library search.

nm - The analyte was not detected in one or more of the duplicate analyses. Therefore, calculation of the RPD is not applicable.

pc - The sample was received with incorrect preservation or in a container not approved by the method. The value reported should be considered an estimate.

ve - The analyte response exceeded the valid instrument calibration range. The value reported is an estimate.

vo - The value reported fell outside the control limits established for this analyte.

x - The sample chromatographic pattern does not resemble the fuel standard used for quantitation.

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Samples received at \_\_\_\_°C

#### ENVIRONMENTAL CHEMISTS

James E. Bruya, Ph.D. Yelena Aravkina, M.S. Michael Erdahl, B.S. Arina Podnozova, B.S. Eric Young, B.S. 3012 16th Avenue West Seattle, WA 98119-2029 (206) 285-8282 fbi@isomedia.com www.friedmanandbruya.com

September 25, 2018

Crystal Thimsen, Project Manager Wood Environment & Infrastructure Solutions, Inc. One Union Square 600 University Street, Suite 600 Seattle, WA 98101

Dear Ms Thimsen:

Included are the results from the testing of material submitted on September 19, 2018 from the Kelly Moore, F&BI 809326 project. There are 4 pages included in this report. Any samples that may remain are currently scheduled for disposal in 30 days. If you would like us to return your samples or arrange for long term storage at our offices, please contact us as soon as possible.

We appreciate this opportunity to be of service to you and hope you will call if you should have any questions.

Sincerely,

FRIEDMAN & BRUYA, INC.

Colo

Michael Erdahl Project Manager

Enclosures WEI0925R.DOC

## ENVIRONMENTAL CHEMISTS

## CASE NARRATIVE

This case narrative encompasses samples received on September 19, 2018 by Friedman & Bruya, Inc. from the Wood Environment & Infrastructure Solutions Kelly Moore, F&BI 809326 project. Samples were logged in under the laboratory ID's listed below.

Laboratory ID	Wood Environment & Infrastructure Solutions
809326 -01	EFF-091918
809326 -02	INF-091918

All quality control requirements were acceptable.

### ENVIRONMENTAL CHEMISTS

Date of Report: 09/25/18 Date Received: 09/19/18 Project: Kelly Moore, F&BI 809326 Date Extracted: 09/20/18 Date Analyzed: 09/20/18

### RESULTS FROM THE ANALYSIS OF AIR SAMPLES FOR BENZENE AND TPH AS GASOLINE USING MODIFIED METHODS 8021B AND NWTPH-Gx

<u>Sample ID</u> Laboratory ID	<u>Benzene</u>	Gasoline <u>Range</u>	Surrogate ( <u>% Recovery)</u> (Limit 50-150)
EFF-091918 809326-01	<0.1	59	82
INF-091918 809326-02	<0.1	470	109
Method Blank 08-1970 MB	<0.1	<10	79

Results Reported as mg/m<sup>3</sup>

### ENVIRONMENTAL CHEMISTS

Date of Report: 09/25/18 Date Received: 09/19/18 Project: Kelly Moore, F&BI 809326

### QUALITY ASSURANCE RESULTS FOR THE ANALYSIS OF AIR SAMPLES FOR BENZENE AND TPH AS GASOLINE USING MODIFIED EPA METHOD 8021B AND NWTPH-Gx

Laboratory Code:	809326-01 (Dupli	cate)		
-	Reporting	Sample	Duplicate	RPD
Analyte	Units	Result	Result	(Limit 20)
Benzene	mg/m <sup>3</sup>	< 0.1	< 0.1	nm
Gasoline	mg/m <sup>3</sup>	59	56	5
Laboratory Code:	Laboratory Contr	ol Sample		

			Percent	
	Reporting	Spike	Recovery	Acceptance
Analyte	Units	Level	LCS	Criteria
Benzene	mg/m <sup>3</sup>	5.0	83	70-130
Gasoline	mg/m <sup>3</sup>	100	141	86-144

#### ENVIRONMENTAL CHEMISTS

## **Data Qualifiers & Definitions**

a - The analyte was detected at a level less than five times the reporting limit. The RPD results may not provide reliable information on the variability of the analysis.

 ${\bf b}$  - The analyte was spiked at a level that was less than five times that present in the sample. Matrix spike recoveries may not be meaningful.

ca - The calibration results for the analyte were outside of acceptance criteria. The value reported is an estimate.

c - The presence of the analyte may be due to carryover from previous sample injections.

cf - The sample was centrifuged prior to analysis.

 ${\rm d}$  - The sample was diluted. Detection limits were raised and surrogate recoveries may not be meaningful.

dv - Insufficient sample volume was available to achieve normal reporting limits.

f - The sample was laboratory filtered prior to analysis.

fb - The analyte was detected in the method blank.

fc - The compound is a common laboratory and field contaminant.

hr - The sample and duplicate were reextracted and reanalyzed. RPD results were still outside of control limits. Variability is attributed to sample inhomogeneity.

hs - Headspace was present in the container used for analysis.

ht – The analysis was performed outside the method or client-specified holding time requirement.

ip - Recovery fell outside of control limits. Compounds in the sample matrix interfered with the quantitation of the analyte.

j - The analyte concentration is reported below the lowest calibration standard. The value reported is an estimate.

J - The internal standard associated with the analyte is out of control limits. The reported concentration is an estimate.

jl - The laboratory control sample(s) percent recovery and/or RPD were out of control limits. The reported concentration should be considered an estimate.

js - The surrogate associated with the analyte is out of control limits. The reported concentration should be considered an estimate.

lc - The presence of the analyte is likely due to laboratory contamination.

L - The reported concentration was generated from a library search.

nm - The analyte was not detected in one or more of the duplicate analyses. Therefore, calculation of the RPD is not applicable.

pc - The sample was received with incorrect preservation or in a container not approved by the method. The value reported should be considered an estimate.

ve - The analyte response exceeded the valid instrument calibration range. The value reported is an estimate.

vo - The value reported fell outside the control limits established for this analyte.

x - The sample chromatographic pattern does not resemble the fuel standard used for quantitation.

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#### ENVIRONMENTAL CHEMISTS

James E. Bruya, Ph.D. Yelena Aravkina, M.S. Michael Erdahl, B.S. Arina Podnozova, B.S. Eric Young, B.S. 3012 16th Avenue West Seattle, WA 98119-2029 (206) 285-8282 fbi@isomedia.com www.friedmanandbruya.com

October 23, 2018

Crystal Thimsen, Project Manager Wood Environment & Infrastructure Solutions, Inc. One Union Square 600 University Street, Suite 600 Seattle, WA 98101

Dear Ms Thimsen:

Included are the results from the testing of material submitted on October 16, 2018 from the Kelly Moore 14697, F&BI 810312 project. There are 4 pages included in this report. Any samples that may remain are currently scheduled for disposal in 30 days. If you would like us to return your samples or arrange for long term storage at our offices, please contact us as soon as possible.

We appreciate this opportunity to be of service to you and hope you will call if you should have any questions.

Sincerely,

FRIEDMAN & BRUYA, INC.

Gal

Michael Erdahl Project Manager

Enclosures WEI1023R.DOC

## ENVIRONMENTAL CHEMISTS

## CASE NARRATIVE

This case narrative encompasses samples received on October 16, 2018 by Friedman & Bruya, Inc. from the Wood Environment & Infrastructure Solutions Kelly Moore 14697, F&BI 810312 project. Samples were logged in under the laboratory ID's listed below.

Laboratory ID	Wood Environment & Infrastructure Solution
810312 -01	INF-101618
810312 -02	EFF-101618

All quality control requirements were acceptable.

### ENVIRONMENTAL CHEMISTS

Date of Report: 10/23/18 Date Received: 10/16/18 Project: Kelly Moore 14697, F&BI 810312 Date Extracted: 10/19/18 Date Analyzed: 10/19/18

### RESULTS FROM THE ANALYSIS OF AIR SAMPLES FOR BENZENE AND TPH AS GASOLINE USING MODIFIED METHODS 8021B AND NWTPH-Gx

Sample ID Laboratory ID	<u>Benzene</u>	Gasoline <u>Range</u>	Surrogate ( <u>% Recovery)</u> (Limit 50-150)
INF-101618 810312-01	<0.1	1,900	149
EFF-101618 810312-02	<0.1	460	101
Method Blank 08-2356 MB	<0.1	<10	72

Results Reported as mg/m<sup>3</sup>

#### ENVIRONMENTAL CHEMISTS

Date of Report: 10/23/18 Date Received: 10/16/18 Project: Kelly Moore 14697, F&BI 810312

### QUALITY ASSURANCE RESULTS FOR THE ANALYSIS OF AIR SAMPLES FOR BENZENE AND TPH AS GASOLINE USING MODIFIED EPA METHOD 8021B AND NWTPH-Gx

Laboratory Code:	810312-01 (Duplica	ate)		
	Reporting	Sample	Duplicate	RPD
Analyte	Units	Result	Result	(Limit 20)
Benzene	mg/m <sup>3</sup>	< 0.1	<0.1	nm
Gasoline	mg/m <sup>3</sup>	1,900	2,100 ve	10
	0			
Laboratory Code:	Laboratory Contro	l Sample		

			Percent	
	Reporting	Spike	Recovery	Acceptance
Analyte	Units	Level	LCS	Criteria
Benzene	mg/m³	5.0	75	70-130
Gasoline	mg/m <sup>3</sup>	100	139	86-144

ENVIRONMENTAL CHEMISTS

# **Data Qualifiers & Definitions**

a - The analyte was detected at a level less than five times the reporting limit. The RPD results may not provide reliable information on the variability of the analysis.

 ${\bf b}$  - The analyte was spiked at a level that was less than five times that present in the sample. Matrix spike recoveries may not be meaningful.

ca - The calibration results for the analyte were outside of acceptance criteria. The value reported is an estimate.

c - The presence of the analyte may be due to carryover from previous sample injections.

cf - The sample was centrifuged prior to analysis.

d - The sample was diluted. Detection limits were raised and surrogate recoveries may not be meaningful.

dv - Insufficient sample volume was available to achieve normal reporting limits.

f - The sample was laboratory filtered prior to analysis.

fb - The analyte was detected in the method blank.

fc - The compound is a common laboratory and field contaminant.

 $hr\ \text{-}\ The\ sample\ and\ duplicate\ were\ reextracted\ and\ reanalyzed.\ RPD\ results\ were\ still\ outside\ of\ control\ limits.\ Variability\ is\ attributed\ to\ sample\ inhomogeneity.$ 

hs - Headspace was present in the container used for analysis.

ht - The analysis was performed outside the method or client-specified holding time requirement.

 ${\rm ip}$  - Recovery fell outside of control limits. Compounds in the sample matrix interfered with the quantitation of the analyte.

j - The analyte concentration is reported below the lowest calibration standard. The value reported is an estimate.

 ${\rm J}$  - The internal standard associated with the analyte is out of control limits. The reported concentration is an estimate.

jl - The laboratory control sample(s) percent recovery and/or RPD were out of control limits. The reported concentration should be considered an estimate.

js - The surrogate associated with the analyte is out of control limits. The reported concentration should be considered an estimate.

lc - The presence of the analyte is likely due to laboratory contamination.

L - The reported concentration was generated from a library search.

nm - The analyte was not detected in one or more of the duplicate analyses. Therefore, calculation of the RPD is not applicable.

pc - The sample was received with incorrect preservation or in a container not approved by the method. The value reported should be considered an estimate.

ve - The analyte response exceeded the valid instrument calibration range. The value reported is an estimate.

vo - The value reported fell outside the control limits established for this analyte.

x - The sample chromatographic pattern does not resemble the fuel standard used for quantitation.

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#### ENVIRONMENTAL CHEMISTS

James E. Bruya, Ph.D. Yelena Aravkina, M.S. Michael Erdahl, B.S. Arina Podnozova, B.S. Eric Young, B.S. 3012 16th Avenue West Seattle, WA 98119-2029 (206) 285-8282 fbi@isomedia.com www.friedmanandbruya.com

April 9, 2019

Crystal Thimsen, Project Manager Wood Environment & Infrastructure Solutions, Inc. One Union Square 600 University Street, Suite 600 Seattle, WA 98101

Dear Ms Thimsen:

Included are the results from the testing of material submitted on April 4, 2019 from the Kelly Moore, F&BI 904101 project. There are 4 pages included in this report.

We appreciate this opportunity to be of service to you and hope you will call if you should have any questions.

Sincerely,

FRIEDMAN & BRUYA, INC.

-le

Michael Erdahl Project Manager

Enclosures WEI0409R.DOC

## ENVIRONMENTAL CHEMISTS

## CASE NARRATIVE

This case narrative encompasses samples received on April 4, 2019 by Friedman & Bruya, Inc. from the Wood Environment & Infrastructure Solutions Kelly Moore, F&BI 904101 project. Samples were logged in under the laboratory ID's listed below.

Laboratory ID	Wood Environment & Infrastructure Solutions
904101 -01	EFF-040319
904101 -02	INF-040319

All quality control requirements were acceptable.

## ENVIRONMENTAL CHEMISTS

Date of Report: 04/09/19 Date Received: 04/04/19 Project: Kelly Moore, F&BI 904101 Date Extracted: 04/05/19 Date Analyzed: 04/05/19

#### RESULTS FROM THE ANALYSIS OF AIR SAMPLES FOR BENZENE, TOLUENE, ETHYLBENZENE, XYLENES AND TPH AS GASOLINE USING MODIFIED METHODS 8021B AND NWTPH-Gx

Results	Reported	as	mg/m <sup>3</sup>	
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Sample ID Laboratory ID	<u>Benzene</u>	<u>Toluene</u>	Ethyl <u>Benzene</u>	Total <u>Xylenes</u>	Gasoline <u>Range</u>	Surrogate ( <u>% Recovery</u> ) (Limit 50-150)
EFF-040319 904101-01	<0.1	<0.1	<0.1	< 0.3	10	105
INF-040319 904101-02	<0.1	0.59	5.6	15	670	ip
Method Blank 09-529 MB	<0.1	<0.1	<0.1	<0.3	<10	98
#### ENVIRONMENTAL CHEMISTS

Date of Report: 04/09/19 Date Received: 04/04/19 Project: Kelly Moore, F&BI 904101

#### QUALITY ASSURANCE RESULTS FOR THE ANALYSIS OF AIR SAMPLES FOR BENZENE, TOLUENE, ETHYLBENZENE, XYLENES, AND TPH AS GASOLINE USING MODIFIED EPA METHOD 8021B AND NWTPH-Gx

Laboratory Code: 904101-01 (Duplicate)								
	Reporting	Sample	Duplicate	RPD				
Analyte	Units	Result	Result	(Limit 20)				
Benzene	mg/m <sup>3</sup>	< 0.1	< 0.1	nm				
Toluene	mg/m <sup>3</sup>	< 0.1	< 0.1	nm				
Ethylbenzene	mg/m <sup>3</sup>	< 0.1	< 0.1	nm				
Xylenes	mg/m <sup>3</sup>	< 0.3	< 0.3	nm				
Gasoline	mg/m <sup>3</sup>	10	<10	nm				

Laboratory Code: Laboratory Control Sample

			Percent	
	Reporting	Spike	Recovery	Acceptance
Analyte	Units	Level	LCS	Criteria
Benzene	mg/m <sup>3</sup>	5.0	92	70-130
Toluene	mg/m <sup>3</sup>	5.0	96	70-130
Ethylbenzene	mg/m <sup>3</sup>	5.0	112	70-130
Xylenes	mg/m <sup>3</sup>	15	109	70-130
Gasoline	mg/m <sup>3</sup>	100	95	86-144

#### ENVIRONMENTAL CHEMISTS

## **Data Qualifiers & Definitions**

a - The analyte was detected at a level less than five times the reporting limit. The RPD results may not provide reliable information on the variability of the analysis.

 ${\bf b}$  - The analyte was spiked at a level that was less than five times that present in the sample. Matrix spike recoveries may not be meaningful.

ca - The calibration results for the analyte were outside of acceptance criteria. The value reported is an estimate.

c - The presence of the analyte may be due to carryover from previous sample injections.

cf - The sample was centrifuged prior to analysis.

 ${\rm d}$  - The sample was diluted. Detection limits were raised and surrogate recoveries may not be meaningful.

dv - Insufficient sample volume was available to achieve normal reporting limits.

f - The sample was laboratory filtered prior to analysis.

fb - The analyte was detected in the method blank.

fc - The analyte is a common laboratory and field contaminant.

hr - The sample and duplicate were reextracted and reanalyzed. RPD results were still outside of control limits. Variability is attributed to sample inhomogeneity.

hs - Headspace was present in the container used for analysis.

ht – The analysis was performed outside the method or client-specified holding time requirement.

ip - Recovery fell outside of control limits due to sample matrix effects.

j - The analyte concentration is reported below the lowest calibration standard. The value reported is an estimate.

 ${\rm J}$  - The internal standard associated with the analyte is out of control limits. The reported concentration is an estimate.

jl - The laboratory control sample(s) percent recovery and/or RPD were out of control limits. The reported concentration should be considered an estimate.

js - The surrogate associated with the analyte is out of control limits. The reported concentration should be considered an estimate.

lc - The presence of the analyte is likely due to laboratory contamination.

L - The reported concentration was generated from a library search.

nm - The analyte was not detected in one or more of the duplicate analyses. Therefore, calculation of the RPD is not applicable.

pc - The sample was received with incorrect preservation or in a container not approved by the method. The value reported should be considered an estimate.

ve - The analyte response exceeded the valid instrument calibration range. The value reported is an estimate.

vo - The value reported fell outside the control limits established for this analyte.

x - The sample chromatographic pattern does not resemble the fuel standard used for quantitation.

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#### ENVIRONMENTAL CHEMISTS

James E. Bruya, Ph.D. Yelena Aravkina, M.S. Michael Erdahl, B.S. Arina Podnozova, B.S. Eric Young, B.S. 3012 16th Avenue West Seattle, WA 98119-2029 (206) 285-8282 fbi@isomedia.com www.friedmanandbruya.com

May 9, 2019

Crystal Thimsen, Project Manager Wood Environment & Infrastructure Solutions, Inc. One Union Square 600 University Street, Suite 600 Seattle, WA 98101

Dear Ms Thimsen:

Included are the results from the testing of material submitted on May 6, 2019 from the Kelly Moore, F&BI 905100 project. There are 4 pages included in this report. Any samples that may remain are currently scheduled for disposal in 30 days, or as directed by the Chain of Custody document. If you would like us to return your samples or arrange for long term storage at our offices, please contact us as soon as possible.

We appreciate this opportunity to be of service to you and hope you will call if you should have any questions.

Sincerely,

FRIEDMAN & BRUYA, INC.

f Colo

Michael Erdahl Project Manager

Enclosures WEI0509R.DOC

#### ENVIRONMENTAL CHEMISTS

#### CASE NARRATIVE

This case narrative encompasses samples received on May 6, 2019 by Friedman & Bruya, Inc. from the Wood Environment & Infrastructure Solutions Kelly Moore, F&BI 905100 project. Samples were logged in under the laboratory ID's listed below.

<u>Laboratory ID</u>	Wood Environment & Infrastructure Solutions
905100 -01	EFF-050319
905100 -02	INF-050319

All quality control requirements were acceptable.

#### ENVIRONMENTAL CHEMISTS

Date of Report: 05/09/19 Date Received: 05/06/19 Project: Kelly Moore, F&BI 905100 Date Extracted: 05/06/19 Date Analyzed: 05/06/19

# RESULTS FROM THE ANALYSIS OF AIR SAMPLES FOR BENZENE USING MODIFIED METHODS 8021B

Results Reported as mg/m<sup>3</sup>

<u>Sample ID</u> Laboratory ID	<u>Benzene</u>	Surrogate ( <u>% Recovery)</u> (Limit 50-150)
EFF-050319 905100-01	<0.1	114
INF-050319 905100-02	< 0.1	127
Method Blank <sup>09-833 MB</sup>	<0.1	113

#### ENVIRONMENTAL CHEMISTS

Date of Report: 05/09/19 Date Received: 05/06/19 Project: Kelly Moore, F&BI 905100

#### QUALITY ASSURANCE RESULTS FOR THE ANALYSIS OF AIR SAMPLES FOR BENZENE USING MODIFIED EPA METHOD 8021B

Laboratory Code: 9	05100-01 (Dupl	icate)			
	Reporting	Samp	le Du	plicate	$\operatorname{RPD}$
Analyte	Units	Resu	lt R	esult	(Limit 20)
Benzene	mg/m <sup>3</sup>	< 0.1	. <	< 0.1	nm
Laboratory Code: L	aboratory Cont	rol Sampl	e Percent		
	Reporting	Spike	Recovery	Acceptance	
Analyte	Units	Level	LCS	Criteria	_
Benzene	mg/m <sup>3</sup>	5.0	86	70-130	_

#### ENVIRONMENTAL CHEMISTS

# **Data Qualifiers & Definitions**

a - The analyte was detected at a level less than five times the reporting limit. The RPD results may not provide reliable information on the variability of the analysis.

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ca - The calibration results for the analyte were outside of acceptance criteria. The value reported is an estimate.

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cf - The sample was centrifuged prior to analysis.

d - The sample was diluted. Detection limits were raised and surrogate recoveries may not be meaningful.

dv - Insufficient sample volume was available to achieve normal reporting limits.

f - The sample was laboratory filtered prior to analysis.

fb - The analyte was detected in the method blank.

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ve - The analyte response exceeded the valid instrument calibration range. The value reported is an estimate.

vo - The value reported fell outside the control limits established for this analyte.

x - The sample chromatographic pattern does not resemble the fuel standard used for quantitation.

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#### ENVIRONMENTAL CHEMISTS

James E. Bruya, Ph.D. Yelena Aravkina, M.S. Michael Erdahl, B.S. Arina Podnozova, B.S. Eric Young, B.S. 3012 16th Avenue West Seattle, WA 98119-2029 (206) 285-8282 fbi@isomedia.com www.friedmanandbruya.com

June 10, 2019

Crystal Thimsen, Project Manager Wood Environment & Infrastructure Solutions, Inc. One Union Square 600 University Street, Suite 600 Seattle, WA 98101

Dear Ms Thimsen:

Included are the results from the testing of material submitted on June 4, 2019 from the Kelly Moore, F&BI 906031 project. There are 5 pages included in this report. Any samples that may remain are currently scheduled for disposal in 30 days, or as directed by the Chain of Custody document. If you would like us to return your samples or arrange for long term storage at our offices, please contact us as soon as possible.

We appreciate this opportunity to be of service to you and hope you will call if you should have any questions.

Sincerely,

FRIEDMAN & BRUYA, INC.

f Colo

Michael Erdahl Project Manager

Enclosures WEI0610R.DOC

#### ENVIRONMENTAL CHEMISTS

## CASE NARRATIVE

This case narrative encompasses samples received on June 4, 2019 by Friedman & Bruya, Inc. from the Wood Environment & Infrastructure Solutions Kelly Moore, F&BI 906031 project. Samples were logged in under the laboratory ID's listed below.

<u>Laboratory ID</u>	Wood Environment & Infrastructure Solutions
906031 -01	EFF-060419
906031 -02	INF-060419
906031 -03	SVE09-060419
906031 -04	SVE10-060419
906031 -05	SVE11-060419
906031 -06	SVE12-060419
906031 -07	SVE13-060419
906031 -08	SVE01-060419
906031 -09	SVE03-060419
906031 -10	SVE05-060419
906031 -11	SVE07-060419
906031 -12	SVE08-060419
906031 -13	SVE06-060419
906031 -14	SVE04-060419
906031 -15	SVE02-060419

All quality control requirements were acceptable.

#### ENVIRONMENTAL CHEMISTS

Date of Report: 06/10/19 Date Received: 06/04/19 Project: Kelly Moore, F&BI 906031 Date Extracted: 06/04/19 Date Analyzed: 06/05/19

# **RESULTS FROM THE ANALYSIS OF AIR SAMPLES** FOR BENZENE AND TPH AS GASOLINE **USING MODIFIED METHODS 8021B AND NWTPH-Gx**

<u>Sample ID</u> Laboratory ID	<u>Benzene</u>	Gasoline <u>Range</u>	Surrogate ( <u>% Recovery</u> ) (Limit 50-150)
EFF-060419 906031-01	<0.1	20	86
INF-060419 906031-02	<0.1	640	113
SVE09-060419 906031-03	<0.1	440	105
SVE10-060419 906031-04 1/5	< 0.5	2,300	101
${\mathop{\rm SVE11-060419}}_{{}_{906031-05}}$	<0.1	660	100
SVE12-060419 906031-06 1/2	< 0.2	1,400	98
SVE13-060419 906031-07	<0.1	760	91
SVE01-060419 906031-08	<0.1	14	85
SVE03-060419 906031-09	< 0.1	2,400	ip
SVE05-060419 906031-10 1/5	< 0.5	3,500	130

Results Reported as mg/m<sup>3</sup>

#### ENVIRONMENTAL CHEMISTS

Date of Report: 06/10/19 Date Received: 06/04/19 Project: Kelly Moore, F&BI 906031 Date Extracted: 06/04/19 Date Analyzed: 06/05/19

#### RESULTS FROM THE ANALYSIS OF AIR SAMPLES FOR BENZENE AND TPH AS GASOLINE USING MODIFIED METHODS 8021B AND NWTPH-Gx Results Reported as mg/m<sup>3</sup>

	_	Gasoline	Surrogate
<u>Sample ID</u> Laboratory ID	<u>Benzene</u>	<u>Range</u>	( <u>% Recovery</u> ) (Limit 50-150)
SVE07-060419 906031-11	< 0.1	230	91
SVE08-060419 906031-12	< 0.1	16	85
SVE06-060419 906031-13	<0.1	33	89
SVE04-060419 906031-14	<0.1	400	103
SVE02-060419 906031-15	<0.1	14	79
Method Blank <sup>09-1280 MB</sup>	< 0.1	<10	76

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#### ENVIRONMENTAL CHEMISTS

Date of Report: 06/10/19 Date Received: 06/04/19 Project: Kelly Moore, F&BI 906031

#### QUALITY ASSURANCE RESULTS FOR THE ANALYSIS OF AIR SAMPLES FOR BENZENE AND TPH AS GASOLINE USING MODIFIED EPA METHOD 8021B AND NWTPH-Gx

Laboratory Code: 906031-01 (Duplicate)								
	Reporting	Sample	Duplicate	RPD				
Analyte	Units	Result	Result	(Limit 20)				
Benzene	mg/m <sup>3</sup>	< 0.1	< 0.1	nm				
Gasoline	mg/m <sup>3</sup>	20	19	5				
Laboratory Code: Laboratory Control Sample								

			Percent	
	Reporting	Spike	Recovery	Acceptance
Analyte	Units	Level	LCS	Criteria
Benzene	mg/m <sup>3</sup>	5.0	91	70-130
Gasoline	mg/m <sup>3</sup>	100	131	86-144

#### ENVIRONMENTAL CHEMISTS

# **Data Qualifiers & Definitions**

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ca - The calibration results for the analyte were outside of acceptance criteria. The value reported is an estimate.

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cf - The sample was centrifuged prior to analysis.

d - The sample was diluted. Detection limits were raised and surrogate recoveries may not be meaningful.

dv - Insufficient sample volume was available to achieve normal reporting limits.

f - The sample was laboratory filtered prior to analysis.

fb - The analyte was detected in the method blank.

fc - The analyte is a common laboratory and field contaminant.

hr - The sample and duplicate were reextracted and reanalyzed. RPD results were still outside of control limits. Variability is attributed to sample inhomogeneity.

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vo - The value reported fell outside the control limits established for this analyte.

x - The sample chromatographic pattern does not resemble the fuel standard used for quantitation.

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Sample ID	Lab ID	Date Sampled	Time Sampled	Sample Type	# of Jars	TPH-HCID	TPH-Diesel	TPH-Gasoline	BTEX by 8021B	VOCs by 8260C	SVOCa by 8270E	PAHs 8270D SIN	BENZEWE				No	tes		
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#### ENVIRONMENTAL CHEMISTS

James E. Bruya, Ph.D. Yelena Aravkina, M.S. Michael Erdahl, B.S. Arina Podnozova, B.S. Eric Young, B.S. 3012 16th Avenue West Seattle, WA 98119-2029 (206) 285-8282 fbi@isomedia.com www.friedmanandbruya.com

July 10, 2019

Crystal Thimsen, Project Manager Wood Environment & Infrastructure Solutions, Inc. One Union Square 600 University Street, Suite 600 Seattle, WA 98101

Dear Ms Thimsen:

Included are the results from the testing of material submitted on July 5, 2019 from the Kelly Moore, F&BI 907102 project. There are 4 pages included in this report. Any samples that may remain are currently scheduled for disposal in 30 days, or as directed by the Chain of Custody document. If you would like us to return your samples or arrange for long term storage at our offices, please contact us as soon as possible.

We appreciate this opportunity to be of service to you and hope you will call if you should have any questions.

Sincerely,

FRIEDMAN & BRUYA, INC.

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Michael Erdahl Project Manager

Enclosures WEI0710R.DOC

#### ENVIRONMENTAL CHEMISTS

#### CASE NARRATIVE

This case narrative encompasses samples received on July 5, 2019 by Friedman & Bruya, Inc. from the Wood Environment & Infrastructure Solutions Kelly Moore, F&BI 907102 project. Samples were logged in under the laboratory ID's listed below.

Laboratory ID	Wood Environment & Infrastructure Solutions
907102 -01	070219-EFF
907102 -02	070219-INF

All quality control requirements were acceptable.

#### ENVIRONMENTAL CHEMISTS

Date of Report: 07/10/19 Date Received: 07/05/19 Project: Kelly Moore, F&BI 907102 Date Extracted: 07/05/19 Date Analyzed: 07/05/19

# RESULTS FROM THE ANALYSIS OF AIR SAMPLES FOR BENZENE USING MODIFIED METHODS 8021B

Results Reported as mg/m<sup>3</sup>

<u>Sample ID</u> Laboratory ID	<u>Benzene</u>	Surrogate ( <u>% Recovery)</u> (Limit 50-150)
070219-EFF 907102-01	< 0.1	88
070219-INF 907102-02	<0.1	93
Method Blank	< 0.1	92

#### ENVIRONMENTAL CHEMISTS

Date of Report: 07/10/19 Date Received: 07/05/19 Project: Kelly Moore, F&BI 907102

#### QUALITY ASSURANCE RESULTS FOR THE ANALYSIS OF AIR SAMPLES FOR BENZENE USING MODIFIED EPA METHOD 8021B

Laboratory Code: Laboratory Control Sample

			Percent	Percent		
	Reporting	Spike	Recovery	Recovery	Acceptance	$\operatorname{RPD}$
Analyte	Units	Level	LCS	LCSD	Criteria	(Limit 20)
Benzene	mg/m <sup>3</sup>	5.0	97	97	70-130	0

#### ENVIRONMENTAL CHEMISTS

# **Data Qualifiers & Definitions**

a - The analyte was detected at a level less than five times the reporting limit. The RPD results may not provide reliable information on the variability of the analysis.

b - The analyte was spiked at a level that was less than five times that present in the sample. Matrix spike recoveries may not be meaningful.

ca - The calibration results for the analyte were outside of acceptance criteria. The value reported is an estimate.

c - The presence of the analyte may be due to carryover from previous sample injections.

cf - The sample was centrifuged prior to analysis.

d - The sample was diluted. Detection limits were raised and surrogate recoveries may not be meaningful.

dv - Insufficient sample volume was available to achieve normal reporting limits.

f - The sample was laboratory filtered prior to analysis.

fb - The analyte was detected in the method blank.

fc - The analyte is a common laboratory and field contaminant.

hr - The sample and duplicate were reextracted and reanalyzed. RPD results were still outside of control limits. Variability is attributed to sample inhomogeneity.

hs - Headspace was present in the container used for analysis.

ht – The analysis was performed outside the method or client-specified holding time requirement.

ip - Recovery fell outside of control limits due to sample matrix effects.

j - The analyte concentration is reported below the lowest calibration standard. The value reported is an estimate.

 ${\rm J}$  - The internal standard associated with the analyte is out of control limits. The reported concentration is an estimate.

jl - The laboratory control sample(s) percent recovery and/or RPD were out of control limits. The reported concentration should be considered an estimate.

js - The surrogate associated with the analyte is out of control limits. The reported concentration should be considered an estimate.

lc - The presence of the analyte is likely due to laboratory contamination.

L - The reported concentration was generated from a library search.

nm - The analyte was not detected in one or more of the duplicate analyses. Therefore, calculation of the RPD is not applicable.

pc - The sample was received with incorrect preservation or in a container not approved by the method. The value reported should be considered an estimate.

ve - The analyte response exceeded the valid instrument calibration range. The value reported is an estimate.

vo - The value reported fell outside the control limits established for this analyte.

x - The sample chromatographic pattern does not resemble the fuel standard used for quantitation.

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Sample ID:		Lab ID	Date Sampled	Time Sampled	Sample Type	# of Jars	TPH-HCID	TPH-Diesel	TPH-Gasoline	BTEX by 8021B	VOCs by 8260C	SVOCe by 8270D	PAHs 8270D SIM	BENZENE				N	otes	
070219-EFF		OL	07-02-19	0942-HES	TEDLAR	1								X				FID-0	.Oppm	
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Seattle, WA 98119-2029	Relinq	uished by: $\int dh n L$	or8 d	RF 1	ن ول	m/	on	g-			T	11 11					8 7/5/1	9 13:00		
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Samples received at 22 τ



# Appendix E



Wood Environment & Infrastructure Solutions, Inc. 600 University Street, Suite 600 Seattle, Washington 98101 USA T: 206-342-1760 www.woodplc.com

April 25, 2019

Project 014697

Ivan Rivera Inspector Puget Sound Clean Air Agency 1904 3rd Avenue, Suite 105 Seattle, WA 98101

Subject: Corrective Actions Memo Former Kelly-Moore Manufacturing Facility 5400–5800 Airport Way South Seattle, Washington

#### Dear Ivan,

Wood Environment & Infrastructure Solutions, Inc. (Wood) has prepared this corrective action memo on behalf of Kelly-Moore Paint Company, Inc. (Kelly-Moore), in response to the Notice of Violation that was issued on April 17, 2019 due to destruction removal efficiencies (DREs) that did not meet the limits in Section 4 of operating permit number 11291.

In addition to the corrective actions detailed below, we contacted the manufacturer of the catalytic oxidation (Catox) unit, Drewelow Remediation Equipment, to confirm that the catalyst was new when installed at the site and that the correct catalyst was selected for site's compounds of concern. The catalyst manufacturer confirmed the correct catalyst was selected. The manufacturer also stated that the life of the catalyst should be up to 30 years, and that we should be able to achieve the permit-required DREs using that Catox unit.

Wood is implementing the following corrective actions in response to the exceedances:

- 1. Monthly inspections will include monthly calibration of the lower explosive limit meter. The manufacturer has provided an up-to-date procedure for this unit.
- 2. Total petroleum hydrocarbon (TPH) readings will be conducted using a handheld flame ionization detector (FID):
  - a. Sampling will continue to be conducted monthly during routine monthly inspections.
  - b. New tubing will be used each month for collecting the samples.
  - c. Samples will be collected first from the effluent port and second from the influent port.
  - d. Influent and effluent samples will be collected in a Tedlar bag, and TPH concentrations will be measured and recorded with an FID calibrated to hexane per permit condition #7. The influent and effluent samples will be collected in the same Tedlar bag used for the FID screening and will be sent to the laboratory for benzene analysis by U.S. Environmental Protection Agency Method 8021.



Ivan Rivera Washington State Department of Ecology April 25, 2019 Page 2 of 2

3. The DREs will be calculated on site using TPH concentrations measured by the FID, and will be communicated to the project manager before leaving the site. The spreadsheet used to calculate the DREs will be programmed to flag any exceedances in the operating permit conditions.

In case of future exceedances of the DRE, the following corrective actions will be implemented:

- 1. Ensure the FID is calibrated correctly and working properly.
- 2. Clean the catalyst using the following procedure:
  - a. Raise "Catox-In" set-point from 650 to 950 degrees Fahrenheit (<sup>0</sup>F) on Clean Air only adjusting the recirculation valve as needed.
  - b. Run at 950°F for at least 60 minutes.
  - c. Reduce the set-point from 950 to  $650^{\circ}$ F.
  - d. Open process valve.
  - e. Close dilution air based on "Catox-Out" temperature.
  - f. Do not exceed 950°F Catox-Out temperature.
- 3. Run Catox-In set-point at 700°F instead of 650°F. Increasing the temperature should improve the DRE.
- 4. Reduce flow rate.

If these steps are completed, and the DRE does not change, Wood will shut off the Catox unit and allow it to cool. After the unit is cool and it is safe, Wood will inspect the catalyst to check for any uneven distribution of the catalyst and will inspect the condition of the refractory material inside the catalyst chamber. Assuming the refractory material is intact, and if the catalyst is not evenly distributed, we will re-distribute it evenly and restart the unit.

If the DRE condition is still not resolved after redistribution of the catalyst, we will contact the manufacturer and obtain a new catalyst, which will be installed before the Catox unit is restarted.

We feel that our discussions directly with the Catox manufacturer has provided us with a solid understanding of the approach we will take in the future to operate the Catox unit correctly to avoid future exceedances of the DRE permit conditions and to identify the corrective actions we outlined above. We hope that these corrective actions address your concerns and reassure you that there will be no more violations of the DRE permit conditions.

Sincerely,

#### Wood Environment & Infrastructure Solutions, Inc.

Crystal Thimsen Senior Scientist Direct Tel: (206) 838-8469 E-mail: crystal.thimsen@woodplc.com

John Long, LHg.

Senior Associate Hydrogeologist Direct Tel: (206) 342-1979 E-mail: John.Long3@woodplc.com

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# **Appendix F**

#### ENVIRONMENTAL CHEMISTS

James E. Bruya, Ph.D. Yelena Aravkina, M.S. Michael Erdahl, B.S. Arina Podnozova, B.S. Eric Young, B.S. 3012 16th Avenue West Seattle, WA 98119-2029 (206) 285-8282 fbi@isomedia.com www.friedmanandbruya.com

October 26, 2018

Crystal Thimsen, Project Manager Wood Environment & Infrastructure Solutions, Inc. One Union Square 600 University Street, Suite 600 Seattle, WA 98101

Dear Ms Thimsen:

Included are the results from the testing of material submitted on October 16, 2018 from the Kelly Moore 14697, F&BI 810322 project. There are 8 pages included in this report.

We appreciate this opportunity to be of service to you and hope you will call if you should have any questions.

Sincerely,

FRIEDMAN & BRUYA, INC.

Michael Erdahl Project Manager

Enclosures WEI1026R.DOC

#### ENVIRONMENTAL CHEMISTS

#### CASE NARRATIVE

This case narrative encompasses samples received on October 16, 2018 by Friedman & Bruya, Inc. from the Wood Environment & Infrastructure Solutions Kelly Moore 14697, F&BI 810322 project. Samples were logged in under the laboratory ID's listed below.

Laboratory ID	Wood Environment & Infrastructure Solutions
810322 -01	KM-AS-1-101618
810322 -02	KM-AS-2-101618
810322 -03	KM-AS-3-101618

Naphthalene was detected in the TO-15 method blank at a level greater than one tenth the concentration detected in the samples. The data were flagged accordingly.

Ethanol exceeded the calibration range in samples KM-AS-1-101618 and KM-AS-2-101618 and carried over into sample KM-AS-3-101618. The data were flagged accordingly.

All other quality control requirements were acceptable.

Client Sample ID:KM-ADate Received:10/16Date Collected:10/16Date Analyzed:10/19Matrix:AirUnits:ug/matrix:	.S-1-101618 /18 /18 /18 3	Client: Project Lab ID Data F Instrui Operat	Wood Environment :: Kelly Moore 14697, :: 810322-01 ile: 101908.D ment: GCMS7 or: MS	t & Infrastruct F&BI 810322	ure Solutio	ons
Surrogates:	% Recovery:	Lower Limit:	Upper Limit:			
4-Bromofluorobenzene	100	70	130			
	Co	ncentration				
	Co	ncentration				
Compounds:	ug/m3	ppbv	Compounds:	ug/m3	ppbv	
Chlorodifluoromethane	1.9	0.53	1-Butanol	<6.1	<2	
Propene	23	13	Carbon tetrachloride	< 0.63	<0.1	
Dichlorodifluoromethane	3.1	0.62	Benzene	1.9	0.60	
Chloromethane	1.5	0.71	Cyclohexane	<6.9	<2	
F-114	< 0.7	< 0.1	2-Pentanone	<3.5	<1	
Isobutene	1.5	0.65	3-Pentanone	<3.5	<1	
Acetaldehyde	87	48	Pentanal	<3.5	<1	
Vinyl chloride	< 0.26	< 0.1	1,2-Dichloropropane	< 0.23	< 0.05	
1,3-Butadiene	0.48	0.22	1,4-Dioxane	< 0.36	< 0.1	
Bromomethane	<1.6	< 0.4	Bromodichloromethane	< 0.067	< 0.01	
Chloroethane	< 0.26	< 0.1	Trichloroethene	< 0.27	< 0.05	
Ethanol	5,900 ve	3,100 ve	cis-1,3-Dichloropropene	< 0.45	< 0.1	
Acetonitrile	<1.7	<1	4-Methyl-2-pentanone	<4.1	<1	
Acrolein	2.6	1.2	trans-1,3-Dichloropropene	< 0.45	< 0.1	
Acrylonitrile	< 0.22	< 0.1	Toluene	7.5	2.0	
Pentane	9.0	3.1	1,1,2-Trichloroethane	< 0.055	< 0.01	
Trichlorofluoromethane	1.9	0.34	3-Hexanone	<4.1	<1	
Acetone	47	20	2-Hexanone	<4.1	<1	
2-Propanol	13	5.2	Hexanal	11	2.6	
Isoprene	0.51	0.18	Tetrachloroethene	< 0.68	< 0.1	
Iodomethane	< 0.58	< 0.1	Dibromochloromethane	< 0.085	< 0.01	
1.1-Dichloroethene	< 0.4	< 0.1	1.2-Dibromoethane (EDB)	< 0.077	< 0.01	
Methacrolein	<2.9	<1	Chlorobenzene	< 0.46	< 0.1	
trans-1,2-Dichloroethene	< 0.4	< 0.1	Ethylbenzene	1.3	0.31	
Cvclopentane	0.41	0.14	1.1.2.2-Tetrachloroethane	< 0.14	< 0.02	
Methyl vinyl ketone	<2.9	<1	m.p-Xvlene	5.0	1.1	
Butanal	<2.9	<1	o-Xvlene	1.7	0.38	
Methylene chloride	<87	<25	Styrene	1.3	0.30	
CFC-113	0.82	0.11	Bromoform	<2.1	< 0.2	
Carbon disulfide	<6.2	<2	Benzvl chloride	< 0.052	< 0.01	
Methyl t-butyl ether (MTI	3E) <1.8	< 0.5	1.3.5-Trimethylbenzene	<2.5	< 0.5	
Vinvl acetate	<7	<2	1.2.4-Trimethylbenzene	<2.5	< 0.5	
1.1-Dichloroethane	< 0.4	< 0.1	1.3-Dichlorobenzene	< 0.6	< 0.1	
cis-1.2-Dichloroethene	< 0.4	< 0.1	1.4-Dichlorobenzene	< 0.24	< 0.04	
Hexane	4.2	1.2	1.2.3-Trimethylbenzene	<2.5	< 0.5	
Chloroform	0.35	0.072	1.2-Dichlorobenzene	<0.6	< 0.1	
2-Butanone (MEK)	15	5 0	1.2.4-Trichlorobenzene	<0.74	< 0.1	
1.2-Dichloroethane (FDC)	0 19	0.046	Naphthalene	0 59 fb	0.11 fb	
1.1.1-Trichloroethane	< 0.55	< 0.1	Hexachlorobutadiene	< 0.21	< 0.02	
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Client Sample ID:KDate Received:1Date Collected:1Date Analyzed:1Matrix:AUnits:u	M-AS-2-101618 0/16/18 0/16/18 0/19/18 .ir g/m3	Client: Projec Lab II Data F Instru Opera	Wood Environment t: Kelly Moore 14697, D: 810322-02 File: 101909.D ment: GCMS7 tor: MS	& Infrastruct F&BI 810322	ure Solution	S
Surrogates:	% Recovery:	Lower Limit: 70	Upper Limit: 130			
+ Di omonuoi obchizent	100		100			
		oncentration				
Commence des	U(	Incentration	Commune day			
Compounds:	ug/m3	ррву	Compounds:	ug/m3	ррру	
Chlorodifluoromethan	ne 1.8	0.49	1-Butanol	< 6.1	<2	
Propene	20	11	Carbon tetrachloride	< 0.63	<0.1	
Dichlorodifluorometha	ane 3.1	0.63	Benzene	1.7	0.52	
Chloromethane	1.4	0.70	Cyclohexane	<6.9	<2	
F-114	<0.7	< 0.1	2-Pentanone	<3.5	<1	
Isobutene	1.4	0.60	3-Pentanone	<3.5	<1	
Acetaldehyde	32	18	Pentanal	<3.5	<1	
Vinyl chloride	< 0.26	<0.1	1,2-Dichloropropane	< 0.23	< 0.05	
1,3-Butadiene	0.44	0.20	1,4-Dioxane	< 0.36	<0.1	
Bromomethane	<1.6	< 0.4	Bromodichloromethane	< 0.067	< 0.01	
Chloroethane	< 0.26	<0.1	Trichloroethene	< 0.27	< 0.05	
Ethanol	5,300 ve	2,800 ve	cis-1,3-Dichloropropene	< 0.45	<0.1	
Acetonitrile	<1.7	<1	4-Methyl-2-pentanone	<4.1	<1	
Acrolein	1.6	0.69	trans-1,3-Dichloropropene	< 0.45	<0.1	
Acrylonitrile	< 0.22	<0.1	Toluene	7.0	1.9	
Pentane	8.5	2.9	1,1,2-Trichloroethane	< 0.055	< 0.01	
Trichlorofluorometha	ne 1.8	0.32	3-Hexanone	<4.1	<1	
Acetone	41	17	2-Hexanone	<4.1	<1	
2-Propanol	11	4.3	Hexanal	7.9	1.9	
Isoprene	0.59	0.21	Tetrachloroethene	< 0.68	< 0.1	
Iodomethane	< 0.58	<0.1	Dibromochloromethane	< 0.085	< 0.01	
1,1-Dichloroethene	< 0.4	<0.1	1,2-Dibromoethane (EDB)	< 0.077	< 0.01	
Methacrolein	<2.9	<1	Chlorobenzene	< 0.46	< 0.1	
trans-1,2-Dichloroeth	ene <0.4	< 0.1	Ethylbenzene	1.2	0.28	
Cyclopentane	< 0.29	< 0.1	1,1,2,2-Tetrachloroethane	< 0.14	< 0.02	
Methyl vinyl ketone	<2.9	<1	m,p-Xylene	4.5	1.0	
Butanal	<2.9	<1	o-Xylene	1.5	0.34	
Methylene chloride	<87	<25	Styrene	1.1	0.25	
CFC-113	< 0.77	< 0.1	Bromoform	<2.1	< 0.2	
Carbon disulfide	<6.2	<2	Benzyl chloride	< 0.052	< 0.01	
Methyl t-butyl ether (	(MTBE) <1.8	< 0.5	1,3,5-Trimethylbenzene	<2.5	< 0.5	
Vinyl acetate	<7	<2	1,2,4-Trimethylbenzene	<2.5	< 0.5	
1,1-Dichloroethane	< 0.4	< 0.1	1,3-Dichlorobenzene	< 0.6	< 0.1	
cis-1,2-Dichloroethene	e <0.4	<0.1	1,4-Dichlorobenzene	< 0.24	< 0.04	
Hexane	<3.5	<1	1,2,3-Trimethylbenzene	<2.5	< 0.5	
Chloroform	0.29	0.060	1,2-Dichlorobenzene	<0.6	< 0.1	
2-Butanone (MEK)	15	5.0	1,2,4-Trichlorobenzene	< 0.74	< 0.1	
1,2-Dichloroethane (E	DC) 0.17	0.041	Naphthalene	0.55 fb	0.10 fb	
1,1,1-Trichloroethane	< 0.55	< 0.1	Hexachlorobutadiene	< 0.21	< 0.02	

Client Sample ID:KM-Date Received:10/1Date Collected:10/1Date Analyzed:10/1Matrix:AirUnits:ug/n	AS-3-101618 6/18 6/18 9/18 n3	Client Projec Lab II Data I Instru Opera	:: Wood Environment ct: Kelly Moore 14697, D: 810322-03 File: 101910.D iment: GCMS7 itor: MS	: & Infrastruc F&BI 810322	ture Solutions
	%	Lower	Upper		
Surrogates:	Recovery:	Limit:	Limit:		
4-Bromofluorobenzene	101	70	130		
	Cor	ncentration			
		ncentration		./ 0	1
Compounds:	ug/m3	ppbv	Compounds:	ug/m3	ppbv
Chlorodifluoromethane	1.5	0.42	1-Butanol	<6.1	<2
Propene	4.5	2.6	Carbon tetrachloride	< 0.63	<0.1
Dichlorodifluoromethane	e 3.1	0.62	Benzene	0.98	0.31
Chloromethane	1.4	0.66	Cyclohexane	<6.9	<2
F-114	< 0.7	< 0.1	2-Pentanone	<3.5	<1
Isobutene	< 0.92	< 0.4	3-Pentanone	<3.5	<1
Acetaldehyde	<9	<5	Pentanal	<3.5	<1
Vinyl chloride	< 0.26	<0.1	1,2-Dichloropropane	< 0.23	< 0.05
1,3-Butadiene	0.19	0.086	1,4-Dioxane	< 0.36	< 0.1
Bromomethane	<1.6	< 0.4	Bromodichloromethane	< 0.067	< 0.01
Chloroethane	< 0.26	< 0.1	Trichloroethene	< 0.27	< 0.05
Ethanol	29 с	15 c	cis-1,3-Dichloropropene	< 0.45	< 0.1
Acetonitrile	<1.7	<1	4-Methyl-2-pentanone	<4.1	<1
Acrolein	< 0.92	< 0.4	trans-1.3-Dichloropropene	< 0.45	< 0.1
Acrylonitrile	< 0.22	< 0.1	Toluene	3.4	0.91
Pentane	5.3	1.8	1.1.2-Trichloroethane	< 0.055	< 0.01
Trichlorofluoromethane	17	0.30	3-Hexanone	< 4 1	<1
Acetone	17	7.0	2-Hexanone	< 4 1	<1
2-Propanol	< 8.6	< 3.5	Hexanal	< 4 1	<1
Isoprene	<0.8	<0.0	Tetrachloroethene	< 0.68	<01
Iodomethane	<0.20	<0.1	Dibromochloromethane	<0.00	<0.1
1 1-Dichloroothono	<0.00	<0.1	1 2 Dibromoothano (FDB)	<0.003	<0.01
Mothacroloin	<0.4	<0.1	Chlorobonzono	<0.077	<0.01
trans 1.2 Dichloroothone	$\sim 2.5$	<0.1	Ethylbonzono	<0.40	<0.1
Cyclopoptano	<pre>- &lt;0.4</pre>	<0.1	1 1 2 2 Totrachloroothano	0.01 <0.14	-0.14 -0.09
Mothyl vinyl kotono	<0.23	<0.1	m n Yylono	<0.14 2 0	<0.02 0.47
Rutopol	<2.9	<1	a Yulono	2.0	0.47
Dutana Mathylana ahlarida	<2.9	<1	0-Aylelle Sturono	0.70 -0.95	0.10
	<07	<2.5	Bromoform	< 0.03	< 0.2
CFC-113 Corbon digulfida	< 0.77	< 0.1	Dromul ablanda	<2.1	< 0.2
Varbul t hutul athon (M	<0.2 FDE) 1.0	< 2 - 0 E	1 2 5 Trimethallorgene	< 0.052	< 0.01
Methyl t-butyl ether (MI	IDE) <1.8	<0.5	1,3,5-1 rimethyldenzene	<2.5	< 0.5
vinyl acetate	</td <td>&lt;2</td> <td>1,2,4-1rimetnyidenzene</td> <td>&lt;2.5</td> <td>&lt; 0.5</td>	<2	1,2,4-1rimetnyidenzene	<2.5	< 0.5
1,1-Dicnioroethane	< 0.4	<0.1	1,3-Dichlorobenzene	< 0.6	<0.1
cis-1,2-Dichloroethene	< 0.4	<0.1	1,4-Dichlorobenzene	< 0.24	< 0.04
Hexane	<3.5	1>	1,2,3-1rimethylbenzene	<2.5	<0.5
Chlorotorm	0.14	0.028	1,2-Dichlorobenzene	< 0.6	< 0.1
2-Butanone (MEK)	<2.9	<1	1,2,4-Trichlorobenzene	< 0.74	<0.1
1,2-Dichloroethane (EDC	.) 0.11	0.028	Naphthalene	0.27 fb	0.052 fb
1,1,1-Trichloroethane	< 0.55	< 0.1	Hexachlorobutadiene	< 0.21	< 0.02

Client Sample ID: Date Received: Date Collected: Date Analyzed: Matrix: Units:	Method Bl Not Applie Not Applie 10/19/18 Air ug/m3	lank cable cable	Client Projec Lab II Data F Instru Opera	: t: ): File: ment: tor:	Wood Environment Kelly Moore 14697, 08-2400 mb 101907.D GCMS7 MS	& Infrastructu F&BI 810322	ire Solutions
Surrogates: 4-Bromofluorobenze	Re	% covery: 97	Lower Limit: 70	Upper Limit: 130			
		-					
~ .		Co	ncentration	~		Cone	centration
Compounds:		ug/m3	ppbv	Compo	unds:	ug/m3	ppbv
Chlorodifluorometh	ane	< 0.35	< 0.1	1-Buta	nol	<6.1	<2
Propene		< 0.69	< 0.4	Carbor	ı tetrachloride	< 0.63	< 0.1
Dichlorodifluoromet	thane	< 0.49	< 0.1	Benzer	ne	< 0.32	< 0.1
Chloromethane		< 0.21	< 0.1	Cycloh	exane	<6.9	<2
F-114		< 0.7	< 0.1	2-Pent	anone	<3.5	<1
Isobutene		< 0.92	< 0.4	3-Pent	anone	<3.5	<1
Acetaldehyde		<9	<5	Pentar	al	<3.5	<1
Vinvl chloride		< 0.26	< 0.1	1.2-Dic	hloropropane	< 0.23	< 0.05
1.3-Butadiene		< 0.022	< 0.01	1.4-Dic	oxane	< 0.36	< 0.1
Bromomethane		<1.6	< 0.4	Bromo	dichloromethane	< 0.067	< 0.01
Chloroethane		< 0.26	< 0.1	Trichle	proethene	< 0.27	< 0.05
Ethanol		<7.5	<4	cis-1.3	Dichloropropene	< 0.45	< 0.1
Acetonitrile		<1.7	<1	4-Meth	vl-2-pentanone	<4.1	<1
Acrolein		< 0.92	< 0.4	trans-1	3-Dichloropropene	<0.45	< 0.1
Acrylonitrile		<0.22	< 0.1	Toluen	ρ	< 0.38	< 0.1
Pentane		<3	<1	1 1 2-T	richloroethane	<0.05	< 0.01
Trichlorofluorometh	hane	<0.56	<01	3-Heva	none	<4 1	<0.01
Acetone	liane	<0.00	< 2	2-Heya	none	<4.1	<1
2-Propanol		<1.0 <8.6	~3 5	Heven	al	<4.1	<1
Isopropo		<0.0	<0.5	Tetrac	hloroothono	29.1 20.68	<01
Iodomothano		<0.20	<0.1	Dibron	achloromothano	<0.00	<0.1
1 1 Dichloroothono		<0.00	<0.1	1 2 Dibi	romoothang (FDB)	<0.005	< 0.01
Mothacroloin		<0.4 <2.0	<0.1	T,2-Dit Chloro	honzono	<0.077	<0.01
trans 1.2 Dichloroo	thono	<0.1	<0.1	Ethylb	onzono	<0.40	<0.1
Cyclopoptapo	ulelle	<0.4	<0.1	1 1 9 9	Totrachloroothano	< 0.43	<0.1
Mothyl vinyl koton	0	<0.29	<0.1	1, 1, 2, 2 m n V	long	<0.14	<0.02
Butanal	e	<2.9 <2.0	<1	nı, p-Ay	hene	<0.87	<0.2
Mathylana ahlarida		~2.3	<1	0-Aylei Sturon	ie o	< 0.45	<0.1
CEC 112		<077	<2.5	Bromo	form	< 0.05	<0.2
Carbon disulfido		<0.11	<0.1	Boord	oblorido	<0.052	<0.2
Mothyl t butyl otho	n (MTDE)	< 0.2	<0.5	1 2 5 T	rimethylhenzone	< 0.052	<0.01
Vinyl costate	I (MIIDE)	<1.0	<0.5	1,3,3-1	rimethylbenzene	<2.5	<0.5
1 1 Dichloroothono		<0.4	<0.1	1,2,4-1	hlorohonzono	<2.5	<0.5
is 1.2 Dichleroothe	200	<0.4	<0.1	1,3-Dic	hlorobonzono	<0.0	<0.1
Hovano		<0.4 2 5	<v.1< td=""><td>1,4-DIC 192T</td><td>rimothylbonzono</td><td>&lt;0.24 ン9 だ</td><td>&lt;0.04</td></v.1<>	1,4-DIC 192T	rimothylbonzono	<0.24 ン9 だ	<0.04
Chloroform		<0.0 _0.010	<1 <0.01	1,2,3-1 1 9 סיר	hlorobonzono	<2.3 20 G	<0.5
2 Putanana (MEV)		<0.049 ∠0.0	<0.01 -1	1,2-DIC	niorobenzene	< U. U - O 74	< 0.1
2-Dutanone (IVIEA)		<2.9 20.04	<1 <0.01	1,2,4-1 Nonh+1	ncilioropenzene	<0.74	<0.1
1 1 1 Trichlorootho	(LDC)	<0.04 <0.55	<0.01	Havas	alorobutadiana	<0.1 -0.91	<0.02 <0.02
1,1,1-11 ICHIOFOetha	ne	<0.00	<0.1	пехас	norobutaulelle	<0.21	<0.02

#### ENVIRONMENTAL CHEMISTS

Date of Report: 10/26/18 Date Received: 10/16/18 Project: Kelly Moore 14697, F&BI 810322

#### QUALITY ASSURANCE RESULTS FOR THE ANALYSIS OF AIR SAMPLES FOR VOLATILES BY METHOD TO-15

Laboratory Code: Laboratory Control Sample

ů ů	-		Percent	
	Reporting	Spike	Recovery	Acceptance
Analyte	Units	Level	LCS	Criteria
Chlorodifluoromethane	ppbv	5	113	70-130
Propene	ppbv	5	98	70-130
Dichlorodifluoromethane	ppbv	5	99	70-130
Chloromethane	ppbv	5	109	70-130
F-114	ppbv	5	108	70-130
Isobutene	ppbv	5	113	70-130
Acetaldehyde	ppbv	5	126	70-130
Vinyl chloride	ppbv	5	108	70-130
1,3-Butadiene	ppbv	5	120	70-130
Bromomethane	ppbv	5	111	70-130
Chloroethane	ppbv	5	107	70-130
Ethanol	ppbv	5	115	70-130
Acetonitrile	ppbv	5	115	70-130
Acrolein	ppbv	5	116	70-130
Acrylonitrile	ppbv	5	109	70-130
Pentane	ppbv	5	114	70-130
Trichlorofluoromethane	ppbv	5	103	70-130
Acetone	ppbv	5	104	70-130
2-Propanol	ppbv	5	113	70-130
Isoprene	ppbv	5	108	70-130
Iodomethane	ppbv	5	95	70-130
1,1-Dichloroethene	ppbv	5	98	70-130
Methacrolein	ppbv	5	106	70-130
trans-1,2-Dichloroethene	ppbv	5	99	70-130
Cyclopentane	ppbv	5	118	70-130
Methyl vinyl ketone	ppbv	5	117	70-130
Butanal	ppbv	5	99	70-130
Methylene chloride	ppbv	5	93	70-130
CFC-113	ppbv	5	100	70-130
Carbon disulfide	ppbv	5	98	70-130
Methyl t-butyl ether (MTBE)	ppbv	5	103	70-130
Vinyl acetate	ppbv	5	109	70-130
1,1-Dichloroethane	ppbv	5	107	70-130
cis-1,2-Dichloroethene	ppbv	5	95	70-130
Hexane	ppbv	5	113	70-130
Chloroform	ppbv	5	105	70-130
2-Butanone (MEK)	ppbv	5	104	70-130
1,2-Dichloroethane (EDC)	ppbv	5	106	70-130
1,1,1-Trichloroethane	ppbv	5	105	70-130

#### ENVIRONMENTAL CHEMISTS

Date of Report: 10/26/18 Date Received: 10/16/18 Project: Kelly Moore 14697, F&BI 810322

#### QUALITY ASSURANCE RESULTS FOR THE ANALYSIS OF AIR SAMPLES FOR VOLATILES BY METHOD TO-15

Laboratory Code: Laboratory Control Sample (continued)

Laboratory Couc. Laboratory Contr	of Sample (co	Jitiliucu)	Percent	
	Reporting	Snike	Recovery	Accentance
Analyte	Units	I evel	LCS	Criteria
1-Butanol	nnby	5	101	70-130
Carbon tetrachloride	ppbv	5	101	70-130
Benzene	ppbv	5	105	70-130
Cyclobexane	ppbv	5	103	70-130
2.Pentanone	ppbv	5	120	70 130
3-Pentanone	ppbv	5	120	70-130
Pontanal	ppbv	5	108	70-130
1 2-Dichloropropane	ppbv	5	105	70-130
1,2-Dictitor opropane	ppbv	5	105	70-130
Bromodichloromothano	ppbv	5	107	70-130
Trichloroothono	ppbv	5	08	70-130
cis 1.3 Dichloropropono	ppbv	5	90 02	70-130
4 Mothyl 2 pontanono	ppbv	5	92 104	70-130
4-methyl-2-pentanone	ppbv	5	104	70-130
Taluana	ppbv	5	100	70-130
1 1 2 Trichloresthere	ppbv	5	90 102	70-130
2 Hovenone	ppov	5	103	70-130
	ppov	5	100	70-130
	ppov	5	110	70-130
Hexanal Tatra shlara athan a	ppov	5 F		70-130
l'etrachioroethene	рроу	5 ~	96	70-130
	pppv	5	114	70-130
1,2-Dibromoetnane (EDB)	ppbv	5	109	70-130
Chlorobenzene	ppbv	5	99	70-130
Ethylbenzene	ppbv	5	88	70-130
1,1,2,2,-Tetrachloroethane	ppbv	5	104	70-130
m,p-Xylene	ppbv	10	95	70-130
o-Xylene	ppbv	5	103	70-130
Styrene	ppbv	5	91	70-130
Bromoform	ppbv	5	97	70-130
Benzyl chloride	ppbv	5	112	70-130
1,3,5-Trimethylbenzene	ppbv	5	88	70-130
1,2,4-Trimethylbenzene	ppbv	5	88	70-130
1,3-Dichlorobenzene	ppbv	5	94	70-130
1,4-Dichlorobenzene	ppbv	5	103	70-130
1,2,3-Trimethylbenzene	ppbv	5	95	70-130
1,2-Dichlorobenzene	ppbv	5	99	70-130
1,2,4-Trichlorobenzene	ppbv	5	84	70-130
Naphthalene	ppbv	5	89	70-130
Hexachloro-1,3-butadiene	ppbv	5	91	70-130

#### ENVIRONMENTAL CHEMISTS

# **Data Qualifiers & Definitions**

a - The analyte was detected at a level less than five times the reporting limit. The RPD results may not provide reliable information on the variability of the analysis.

 ${\bf b}$  - The analyte was spiked at a level that was less than five times that present in the sample. Matrix spike recoveries may not be meaningful.

ca - The calibration results for the analyte were outside of acceptance criteria. The value reported is an estimate.

c - The presence of the analyte may be due to carryover from previous sample injections.

cf - The sample was centrifuged prior to analysis.

d - The sample was diluted. Detection limits were raised and surrogate recoveries may not be meaningful.

dv - Insufficient sample volume was available to achieve normal reporting limits.

f - The sample was laboratory filtered prior to analysis.

fb - The analyte was detected in the method blank.

fc - The compound is a common laboratory and field contaminant.

hr - The sample and duplicate were reextracted and reanalyzed. RPD results were still outside of control limits. Variability is attributed to sample inhomogeneity.

hs - Headspace was present in the container used for analysis.

ht – The analysis was performed outside the method or client-specified holding time requirement.

ip - Recovery fell outside of control limits. Compounds in the sample matrix interfered with the quantitation of the analyte.

j - The analyte concentration is reported below the lowest calibration standard. The value reported is an estimate.

 ${\rm J}$  - The internal standard associated with the analyte is out of control limits. The reported concentration is an estimate.

jl - The laboratory control sample(s) percent recovery and/or RPD were out of control limits. The reported concentration should be considered an estimate.

js - The surrogate associated with the analyte is out of control limits. The reported concentration should be considered an estimate.

lc - The presence of the analyte is likely due to laboratory contamination.

L - The reported concentration was generated from a library search.

nm - The analyte was not detected in one or more of the duplicate analyses. Therefore, calculation of the RPD is not applicable.

pc - The sample was received with incorrect preservation or in a container not approved by the method. The value reported should be considered an estimate.

ve - The analyte response exceeded the valid instrument calibration range. The value reported is an estimate.

vo - The value reported fell outside the control limits established for this analyte.

x - The sample chromatographic pattern does not resemble the fuel standard used for quantitation.
810322				SAMP	LE CE	IAIN (	)F CU	STOI	ΟY		ь.đ	E	intulat
Report To (1954) Thin	in.			SAMPLERS (signature)									
Company Work				PROJ	PROJECT NAME					-	2 PC	)#	TURNAROUND TIME
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City, State, ZIP Sottie	No	No 1610 REPORTING LIEVEL				IN	VOI	CE T	TO SAMPLE DISPOSAL				
Phone 781-724-1014 Email CryGht Throw Manual F				indo Sub	⊂ Indoor Air ⊂ Deep Soil Gas ⊖ Sub Slab/Soil Gas ⊃ SVE/Grab					· · · ·		· .	C Other
	1	1	1	1	]		1	1	ANA	LYSI	S RE	QUE	ESTED
Sample Name	Lab ID	Canister ID	Flow Contr. ID	Date	Field Initial Press.	Field Initial	Field Final Press.	Field Final	TO-15 Full Scan	TO-15 BTEXN	TO-15 eVOCs		
KM-45-1-10616	01	21492	07051	10/16	76.0	11me 14(4	7 <i>0</i>	17ime					Notes
KM-13-2-101616	02	18570	07545	19/16	29.5	02/12	30	1628	$\frac{1}{2}$				
KM-A5-3-191618	03	18563	(17652	1916	26.1	2410	6.8	1710	$\hat{\boldsymbol{\lambda}}$				
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<b>Addama,</b>												** <b>3</b> ********	Samples received at _20 °c

Friedman & Bruya, Inc.	SIGNATURE	PRINT NAME	COMPANY		
3013 16th Avenue West	Relinquished by:		<u>COMPANY</u>	DATE	TIME
Seattle, WA 98119-2029	Received by:	William Young	41pm	10/16	1741
Ph. (206) 285-8282	Relinquished by:	- Der Shiman	EBI	10/16/14	(7:41
Fax (206) 283-5044	Received by:	است می ورد اور		6 6 6	
FORMS\COC\COUTO-16.DOC					

#### ENVIRONMENTAL CHEMISTS

James E. Bruya, Ph.D. Yelena Aravkina, M.S. Michael Erdahl, B.S. Arina Podnozova, B.S. Eric Young, B.S. 3012 16th Avenue West Seattle, WA 98119-2029 (206) 285-8282 fbi@isomedia.com www.friedmanandbruya.com

June 18, 2019

Crystal Thimsen, Project Manager Wood Environment & Infrastructure Solutions, Inc. One Union Square 600 University Street, Suite 600 Seattle, WA 98101

Dear Ms Thimsen:

Included are the results from the testing of material submitted on June 7, 2019 from the Kelly Moore, F&BI 906142 project. There are 8 pages included in this report.

We appreciate this opportunity to be of service to you and hope you will call if you should have any questions.

Sincerely,

FRIEDMAN & BRUYA, INC.

Glu

Michael Erdahl Project Manager

Enclosures WEI0618R.DOC

#### ENVIRONMENTAL CHEMISTS

#### CASE NARRATIVE

This case narrative encompasses samples received on June 7, 2019 by Friedman & Bruya, Inc. from the Wood Environment & Infrastructure Solutions Kelly Moore, F&BI 906142 project. Samples were logged in under the laboratory ID's listed below.

<u>Laboratory ID</u>	Wood Environment & Infrastructure Solutions
906142 -01	AS1-060519
906142 -02	AS2-060519
906142 -03	AS3-060519

The TO-15 calibration standard failed the acceptance criteria for bromomethane. In addition, the ethanol concentration for all samples exceeded the calibration range. The data were flagged accordingly.

All other quality control requirements were acceptable.

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Collected: Date Analyzed: Matrix: Units:	AS1-060519 06/07/19 06/05/30 06/10/19 Air ug/m3	Clier Proj Lab Data Instr Oper	nt: ect: ID: a File: rument: rator:	Wood Environment & Kelly Moore, F&BI 9 906142-01 061020.D GCMS7 MS	& Infrastructu 06142	re
	%	Lower	Upper			
Surrogates:	Recovery:	Limit:	Limit:			
4-Bromofluorobenze	ene 88	70	130			
	Concen	tration			Concen	tration
Compounds:	ug/m3	ppbv	Compo	unds:	ug/m3	ppbv
Propene	< 0.69	< 0.4	1,2-Dic	hloropropane	< 0.23	< 0.05
Dichlorodifluorome	thane 2.2	0.44	1,4-Dio	xane	< 0.36	< 0.1
Chloromethane	<2.1	<1	2,2,4-T	rimethylpentane	<4.7	<1
F-114	< 0.7	< 0.1	Methyl	methacrylate	<4.1	<1
Vinyl chloride	< 0.26	< 0.1	Heptar	ie	<4.1	<1
1,3-Butadiene	< 0.022	< 0.01	Bromo	lichloromethane	< 0.067	< 0.01
Butane	<2.4	<1	Trichlo	roethene	< 0.27	< 0.05
Bromomethane	<1.6 ca	<0.4 ca	cis-1,3-	Dichloropropene	< 0.45	< 0.1
Chloroethane	<2.6	<1	4-Meth	yl-2-pentanone	<4.1	<1
Vinyl bromide	< 0.44	< 0.1	trans-1	,3-Dichloropropene	< 0.45	< 0.1
Ethanol	1,500 ve	770 ve	Toluen	е	2.3	0.62
Acrolein	1.1	0.49	1,1,2-T	richloroethane	< 0.11	< 0.02
Pentane	<3	<1	2-Hexa	none	<4.1	<1
Trichlorofluorometh	nane <2.2	< 0.4	Tetrack	nloroethene	<6.8	<1
Acetone	18	7.7	Dibrom	ochloromethane	< 0.085	< 0.01
2-Propanol	<8.6	<3.5	1,2 <b>-</b> Dib	romoethane (EDB)	< 0.077	< 0.01
1,1-Dichloroethene	< 0.4	< 0.1	Chlorol	oenzene	< 0.46	< 0.1
trans-1,2-Dichloroe	thene <0.4	< 0.1	Ethylbo	enzene	< 0.43	< 0.1
Methylene chloride	<87	<25	1,1,2,2-	Tetrachloroethane	< 0.14	< 0.02
t-Butyl alcohol (TB	A) <12	<4	Nonane	Э	<5.2	<1
3-Chloropropene	<1.3	< 0.4	Isoprop	ylbenzene	<2.5	< 0.5
CFC-113	< 0.77	< 0.1	2-Chloi	rotoluene	<5.2	<1
Carbon disulfide	< 6.2	<2	Propyll	oenzene	<2.5	< 0.5
Methyl t-butyl ethe	r (MTBE) <1.8	< 0.5	4-Ethy	toluene	<2.5	< 0.5
Vinyl acetate	<7	<2	m,p-Xy	lene	1.2	0.27
1,1-Dichloroethane	< 0.4	< 0.1	o-Xyler	ie	0.44	0.10
cis-1,2-Dichloroethe	ene <0.4	< 0.1	Styrene	9	< 0.85	< 0.2
Hexane	<3.5	<1	Bromof	orm	<2.1	< 0.2
Chloroform	0.24	0.049	Benzyl	chloride	< 0.052	< 0.01
Ethyl acetate	14	3.8	1,3,5-T	rimethylbenzene	<2.5	< 0.5
Tetrahydrofuran	2.4	0.80	1,2,4-T	rimethylbenzene	<2.5	< 0.5
2-Butanone (MEK)	17	5.8	1,3-Dic	hlorobenzene	< 0.6	< 0.1
1,2-Dichloroethane	(EDC) 0.11	0.026	1,4-Dic	hlorobenzene	< 0.24	< 0.04
1,1,1-Trichloroethan	ne <0.55	< 0.1	1,2-Dic	hlorobenzene	< 0.6	< 0.1
Carbon tetrachlorid	le <0.63	< 0.1	1,2,4-T	richlorobenzene	< 0.74	< 0.1
Benzene	< 0.32	< 0.1	Naphth	nalene	0.14	0.026
Cyclohexane	< 6.9	<2	Hexach	lorobutadiene	< 0.21	< 0.02

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Collected: Date Analyzed: Matrix: Units:	AS2-060519 06/07/19 06/05/30 06/11/19 Air ug/m3	Clie Proj Lab Dat: Inst Ope	nt: lect: ID: a File: rument: rator:	Wood Environment of Kelly Moore, F&BI 9 906142-02 061021.D GCMS7 MS	& Infrastructu 906142	re
	%	Lower	Upper			
Surrogates:	Recovery:	Limit:	Limit:			
4-Bromofluorobenze	ene 100	70	130			
	Concer	ntration			Concen	tration
Compounds:	ug/m3	ppbv	Compo	unds:	ug/m3	ppbv
Propene	<0.69	< 0.4	1,2-Dic	hloropropane	< 0.23	< 0.05
Dichlorodifluorome	thane 2.3	0.46	1,4-Dio	xane	< 0.36	< 0.1
Chloromethane	<2.1	<1	2,2,4-T	rimethylpentane	<4.7	<1
F-114	< 0.7	< 0.1	Methyl	methacrylate	<4.1	<1
Vinyl chloride	< 0.26	< 0.1	Heptar	ie	<4.1	<1
1,3-Butadiene	< 0.022	< 0.01	Bromo	lichloromethane	< 0.067	< 0.01
Butane	<2.4	<1	Trichlo	roethene	< 0.27	< 0.05
Bromomethane	<1.6 ca	<0.4 ca	cis-1,3-	Dichloropropene	0.51	0.11
Chloroethane	<2.6	<1	4-Meth	yl-2-pentanone	<4.1	<1
Vinyl bromide	< 0.44	< 0.1	trans-1	,3-Dichloropropene	< 0.45	< 0.1
Ethanol	1,400 ve	750 ve	Toluen	е	2.1	0.57
Acrolein	< 0.92	< 0.4	1, 1, 2-T	richloroethane	< 0.11	< 0.02
Pentane	<3	<1	2-Hexa	none	<4.1	<1
Trichlorofluorometh	nane <2.2	< 0.4	Tetrach	nloroethene	<6.8	<1
Acetone	16	6.8	Dibrom	ochloromethane	< 0.085	< 0.01
2-Propanol	<8.6	<3.5	1,2-Dib	romoethane (EDB)	< 0.077	< 0.01
1,1-Dichloroethene	< 0.4	< 0.1	Chlorol	penzene	< 0.46	< 0.1
trans-1,2-Dichloroe	thene <0.4	< 0.1	Ethylbo	enzene	< 0.43	< 0.1
Methylene chloride	<87	<25	1, 1, 2, 2.	Tetrachloroethane	< 0.14	< 0.02
t-Butyl alcohol (TB	A) <12	<4	Nonane	Э	< 5.2	<1
3-Chloropropene	<1.3	< 0.4	Isoprop	ylbenzene	<2.5	< 0.5
CFC-113	< 0.77	< 0.1	2-Chlor	otoluene	< 5.2	<1
Carbon disulfide	< 6.2	<2	Propyll	oenzene	<2.5	< 0.5
Methyl t-butyl ethe	r (MTBE) <1.8	< 0.5	4-Ethy	toluene	<2.5	< 0.5
Vinyl acetate	<7	<2	m,p-Xy	lene	1.3	0.29
1,1-Dichloroethane	< 0.4	< 0.1	o-Xyler	ie	0.46	0.11
cis-1,2-Dichloroethe	ene <0.4	< 0.1	Styren	Э	< 0.85	< 0.2
Hexane	<3.5	<1	Bromof	form	<2.1	< 0.2
Chloroform	0.24	0.049	Benzyl	chloride	< 0.052	< 0.01
Ethyl acetate	13	3.5	1,3,5-T	rimethylbenzene	<2.5	< 0.5
Tetrahydrofuran	2.2	0.75	1,2,4-T	rimethylbenzene	<2.5	< 0.5
2-Butanone (MEK)	17	5.8	1,3-Dic	hlorobenzene	< 0.6	< 0.1
1,2-Dichloroethane	(EDC) 0.11	0.026	1,4-Dic	hlorobenzene	< 0.24	< 0.04
1,1,1-Trichloroethan	ne <0.55	< 0.1	1,2-Dic	hlorobenzene	< 0.6	< 0.1
Carbon tetrachlorid	le <0.63	< 0.1	1,2,4-T	richlorobenzene	< 0.74	< 0.1
Benzene	< 0.32	< 0.1	Naphtł	nalene	0.14	0.027
Cyclohexane	< 6.9	<2	Hexach	lorobutadiene	< 0.21	< 0.02

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Collected: Date Analyzed: Matrix: Units:	AS3-0605 06/07/19 06/05/30 06/11/19 Air ug/m3	19	Clien Projec Lab I Data Instru Opera	t: ct: D: File: ument: ator:	Wood Environment Kelly Moore, F&BI 9 906142-03 061022.D GCMS7 MS	& Infrastructu 906142	re
		%	Lower	Upper			
Surrogates:	Re	covery:	Limit:	Limit:			
4-Bromofluorobenzer	ne	96	70	130			
		Concen	tration			Concen	tration
Compounds:		ug/m3	ppbv	Compo	ounds:	ug/m3	ppbv
Propene		< 0.69	< 0.4	1,2-Dic	chloropropane	< 0.23	< 0.05
Dichlorodifluorometh	nane	2.3	0.46	1,4-Dio	oxane	< 0.36	< 0.1
Chloromethane		<2.1	<1	2,2,4-T	rimethylpentane	<4.7	<1
F-114		< 0.7	< 0.1	Methy	l methacrylate	<4.1	<1
Vinyl chloride		< 0.26	< 0.1	Heptai	ne	<4.1	<1
1,3-Butadiene		< 0.022	< 0.01	Bromo	dichloromethane	< 0.067	< 0.01
Butane		<2.4	<1	Trichlo	oroethene	< 0.27	< 0.05
Bromomethane		<1.6 ca	<0.4 ca	cis-1,3-	Dichloropropene	< 0.45	< 0.1
Chloroethane		<2.6	<1	4-Meth	yl-2-pentanone	<4.1	<1
Vinyl bromide		< 0.44	< 0.1	trans-1	1,3-Dichloropropene	< 0.45	< 0.1
Ethanol		150 ve	82 ve	Toluen	e	1.5	0.41
Acrolein		< 0.92	< 0.4	1, 1, 2-T	richloroethane	< 0.11	< 0.02
Pentane		<3	<1	2-Hexa	anone	<4.1	<1
Trichlorofluorometha	ane	<2.2	< 0.4	Tetrac	hloroethene	< 6.8	<1
Acetone		7.5	3.2	Dibron	nochloromethane	< 0.085	< 0.01
2-Propanol		<8.6	<3.5	1,2-Dib	promoethane (EDB)	< 0.077	< 0.01
1,1-Dichloroethene		< 0.4	< 0.1	Chloro	benzene	< 0.46	< 0.1
trans-1,2-Dichloroeth	nene	< 0.4	< 0.1	Ethylb	enzene	< 0.43	< 0.1
Methylene chloride		<87	<25	1.1.2.2	-Tetrachloroethane	< 0.14	< 0.02
t-Butyl alcohol (TBA	)	<12	<4	Nonan	e	<5.2	<1
3-Chloropropene	,	<1.3	< 0.4	Isoproi	ovlbenzene	<2.5	< 0.5
CFC-113		< 0.77	< 0.1	2-Chlo	rotoluene	<5.2	<1
Carbon disulfide		< 6.2	<2	Propvl	benzene	<2.5	< 0.5
Methyl t-butyl ether	(MTBE)	<1.8	< 0.5	4-Ethv	ltoluene	<2.5	< 0.5
Vinvl acetate		<7	<2	m.p-Xy	vlene	1.1	0.24
1.1-Dichloroethane		< 0.4	< 0.1	o-Xvlei	ne	< 0.43	< 0.1
cis-1.2-Dichloroether	ne	< 0.4	< 0.1	Styren	e	< 0.85	< 0.2
Hexane	-	<3.5	<1	Bromo	form	<2.1	< 0.2
Chloroform		0.083	$0.017^{-1}$	Benzvl	chloride	< 0.052	< 0.01
Ethyl acetate		<7.2	<2	1.3.5-T	rimethylbenzene	<2.5	< 0.5
Tetrahydrofuran		< 0.29	< 0.1	1.2.4-T	rimethylbenzene	<2.5	< 0.5
2-Butanone (MEK)		<2.9	<1	1.3-Dic	chlorobenzene	< 0.6	< 0.1
1,2-Dichloroethane (	EDC)	0.077	0.019	1.4-Dic	chlorobenzene	< 0.24	< 0.04
1.1.1-Trichloroethan	e	< 0.55	< 0.1	1.2-Die	chlorobenzene	<0.6	< 0.1
Carbon tetrachloride	-	<0.63	<0.1	1.2.4-T	richlorobenzene	< 0.74	<0.1
Benzene		< 0.32	< 0.1	Napht	halene	< 0.079	< 0.015
Cyclohexane		<6.9	<2	Hexacl	nlorobutadiene	< 0.21	< 0.02

## ENVIRONMENTAL CHEMISTS

Client Sample ID: Date Received: Date Collected: Date Analyzed: Matrix: Units:	Method E Not Appli Not Appli 06/10/19 Air	Blank icable icable	Clien Proje Lab I Data Instru	t: ct: D: File: ument:	Wood Environment & Kelly Moore, F&BI 9 09-1324 mb 061013.D GCMS7 MS	ire	
Onits.	ug/III0		Opera	at01.	WIG		
		%	Lower	Upper			
Surrogates:	R	ecovery:	Limit:	Limit:			
4-Bromofluorobenzo	ene	89	70	130			
		Concen	tration			Concer	tration
Compounds:		ug/m3	ppbv	Compo	ounds:	ug/m3	ppbv
Propene		< 0.69	< 0.4	1,2-Die	chloropropane	< 0.23	< 0.05
Dichlorodifluorome	thane	< 0.49	< 0.1	1,4-Die	oxane	< 0.36	< 0.1
Chloromethane		<2.1	<1	2,2,4-T	rimethylpentane	<4.7	<1
F-114		< 0.7	< 0.1	Methy	l methacrylate	<4.1	<1
Vinyl chloride		< 0.26	< 0.1	Heptar	ne	<4.1	<1
1,3-Butadiene		< 0.022	< 0.01	Bromo	dichloromethane	< 0.067	< 0.01
Butane		<2.4	<1	Trichle	oroethene	< 0.27	< 0.05
Bromomethane		<1.6 ca	<0.4 ca	cis-1,3	-Dichloropropene	< 0.45	< 0.1
Chloroethane <		<2.6	<1	4-Meth	nyl-2-pentanone	<4.1	<1
Vinyl bromide <		< 0.44	< 0.1	trans-1	1,3-Dichloropropene	< 0.45	< 0.1
Ethanol			<4	Toluen	le	< 0.38	< 0.1
Acrolein		< 0.92	< 0.4	1,1,2 <b>-</b> T	richloroethane	< 0.11	< 0.02
Pentane		<3	<1	2-Hexa	anone	<4.1	<1
Trichlorofluorometh	nane	<2.2	< 0.4	Tetrac	hloroethene	< 6.8	<1
Acetone		<4.8	<2	Dibron	nochloromethane	< 0.085	< 0.01
2-Propanol		<8.6	<3.5	1,2-Dil	promoethane (EDB)	< 0.077	< 0.01
1,1-Dichloroethene		< 0.4	< 0.1	Chloro	benzene	< 0.46	< 0.1
trans-1,2-Dichloroe	thene	< 0.4	< 0.1	Ethylb	enzene	< 0.43	< 0.1
Methylene chloride		<87	<25	1,1,2,2	-Tetrachloroethane	< 0.14	< 0.02
t-Butyl alcohol (TB.	A)	<12	<4	Nonan	e	< 5.2	<1
3-Chloropropene	,	<1.3	< 0.4	Isopro	pylbenzene	<2.5	< 0.5
CFC-113		< 0.77	< 0.1	2-Chlo	rotoluene	< 5.2	<1
Carbon disulfide		< 6.2	<2	Propyl	benzene	<2.5	< 0.5
Methyl t-butyl ethe	r (MTBE)	<1.8	< 0.5	4-Ethy	ltoluene	<2.5	< 0.5
Vinyl acetate	````	<7	<2	m,p-Xy	vlene	< 0.87	< 0.2
1,1-Dichloroethane		< 0.4	< 0.1	o-Xylei	ne	< 0.43	< 0.1
cis-1,2-Dichloroethe	ene	< 0.4	< 0.1	Styren	e	< 0.85	< 0.2
Hexane		<3.5	<1	Bromo	form	<2.1	< 0.2
Chloroform		< 0.049	< 0.01	Benzyl	chloride	< 0.052	< 0.01
Ethyl acetate		<7.2	<2	1,3,5-T	rimethylbenzene	<2.5	< 0.5
Tetrahydrofuran		< 0.29	< 0.1	1,2,4-T	rimethylbenzene	<2.5	< 0.5
2-Butanone (MEK)		<2.9	<1	1,3-Dio	chlorobenzene	< 0.6	< 0.1
1,2-Dichloroethane	(EDC)	< 0.04	< 0.01	1,4-Dio	chlorobenzene	< 0.24	< 0.04
1,1,1-Trichloroetha	ne	< 0.55	< 0.1	1,2-Dio	chlorobenzene	< 0.6	< 0.1
Carbon tetrachlorid	le	< 0.63	< 0.1	1,2,4-T	richlorobenzene	< 0.74	< 0.1
Benzene		< 0.32	< 0.1	Napht	halene	< 0.079	< 0.015
Cyclohexane		<6.9	<2	Hexacl	hlorobutadiene	< 0.21	< 0.02

#### ENVIRONMENTAL CHEMISTS

Date of Report: 06/18/19 Date Received: 06/07/19 Project: Kelly Moore, F&BI 906142

#### QUALITY ASSURANCE RESULTS FOR THE ANALYSIS OF AIR SAMPLES FOR VOLATILES BY METHOD TO-15

Laboratory Code: Laboratory Control Sample

	_		Percent					
	Reporting	Spike	Recovery	Acceptance				
Analyte	Units	Level	LCS	Criteria				
Propene	ppbv	5	98	70-130				
Dichlorodifluoromethane	ppbv	<b>5</b>	99	70-130				
Chloromethane	ppbv	<b>5</b>	96	70-130				
F-114	ppbv	<b>5</b>	96	70-130				
Vinyl chloride	ppbv	<b>5</b>	97	70-130				
1,3-Butadiene	ppbv	5	95	70-130				
Butane	ppbv	<b>5</b>	98	70-130				
Bromomethane	ppbv	<b>5</b>	69 vo	70-130				
Chloroethane	ppbv	<b>5</b>	99	70-130				
Ethanol	ppbv	<b>5</b>	89	70-130				
Acrolein	ppbv	<b>5</b>	87	70-130				
Pentane	ppbv	<b>5</b>	94	70-130				
Trichlorofluoromethane	ppbv	<b>5</b>	102	70-130				
Acetone	ppbv	<b>5</b>	89	70-130				
2-Propanol	ppbv	<b>5</b>	96	70-130				
1,1-Dichloroethene	ppbv	<b>5</b>	99	70-130				
trans-1,2-Dichloroethene	ppbv	<b>5</b>	97	70-130				
Methylene chloride	$\operatorname{ppbv}$	<b>5</b>	78	70-130				
t-Butyl alcohol (TBA)	$\operatorname{ppbv}$	<b>5</b>	96	70-130				
3-Chloropropene	$\operatorname{ppbv}$	<b>5</b>	93	70-130				
CFC-113	$\operatorname{ppbv}$	<b>5</b>	96	70-130				
Carbon disulfide	$\operatorname{ppbv}$	<b>5</b>	92	70-130				
Methyl t-butyl ether (MTBE)	$\operatorname{ppbv}$	<b>5</b>	84	70 - 130				
Vinyl acetate	$\operatorname{ppbv}$	<b>5</b>	97	70 - 130				
1,1-Dichloroethane	$\operatorname{ppbv}$	<b>5</b>	96	70 - 130				
cis-1,2-Dichloroethene	$\operatorname{ppbv}$	<b>5</b>	98	70-130				
Hexane	$\operatorname{ppbv}$	<b>5</b>	95	70-130				
Chloroform	$\operatorname{ppbv}$	<b>5</b>	96	70-130				
Ethyl acetate	$\operatorname{ppbv}$	<b>5</b>	77	70-130				
Tetrahydrofuran	$\operatorname{ppbv}$	<b>5</b>	87	70-130				
2-Butanone (MEK)	$\operatorname{ppbv}$	<b>5</b>	99	70-130				
1,2-Dichloroethane (EDC)	$\operatorname{ppbv}$	<b>5</b>	99	70-130				
1,1,1-Trichloroethane	$\operatorname{ppbv}$	<b>5</b>	96	70-130				
Carbon tetrachloride	$\operatorname{ppbv}$	<b>5</b>	96	70-130				
Benzene	$\operatorname{ppbv}$	<b>5</b>	89	70-130				
Cyclohexane	$\operatorname{ppbv}$	<b>5</b>	90	70-130				
1,2-Dichloropropane	$\operatorname{ppbv}$	<b>5</b>	98	70-130				
1,4-Dioxane	$\operatorname{ppbv}$	<b>5</b>	113	70-130				
2,2,4-Trimethylpentane	$\operatorname{ppbv}$	<b>5</b>	100	70-130				
Methyl methacrylate	ppbv	<b>5</b>	100	70-130				

#### ENVIRONMENTAL CHEMISTS

Date of Report: 06/18/19 Date Received: 06/07/19 Project: Kelly Moore, F&BI 906142

#### QUALITY ASSURANCE RESULTS FOR THE ANALYSIS OF AIR SAMPLES FOR VOLATILES BY METHOD TO-15

Laboratory Code: Laboratory Control Sample

	<b>I</b>	Percent						
	Reporting	Spike	Recovery	Acceptance				
Analyte	Units	Level	LCS	Criteria				
Heptane	ppbv	5	98	70-130				
Bromodichloromethane	ppbv	<b>5</b>	103	70-130				
Trichloroethene	$\operatorname{ppbv}$	<b>5</b>	97	70-130				
cis-1,3-Dichloropropene	$\operatorname{ppbv}$	<b>5</b>	99	70-130				
4-Methyl-2-pentanone	$\operatorname{ppbv}$	<b>5</b>	107	70-130				
trans-1,3-Dichloropropene	$\operatorname{ppbv}$	<b>5</b>	101	70-130				
Toluene	$\operatorname{ppbv}$	<b>5</b>	90	70-130				
1,1,2-Trichloroethane	$\operatorname{ppbv}$	<b>5</b>	100	70-130				
2-Hexanone	$\operatorname{ppbv}$	<b>5</b>	104	70-130				
Tetrachloroethene	$\operatorname{ppbv}$	<b>5</b>	95	70-130				
Dibromochloromethane	$\operatorname{ppbv}$	<b>5</b>	101	70-130				
1,2-Dibromoethane (EDB)	$\operatorname{ppbv}$	<b>5</b>	99	70-130				
Chlorobenzene	$\operatorname{ppbv}$	<b>5</b>	100	70-130				
Ethylbenzene	ppbv	<b>5</b>	94	70-130				
1,1,2,2-Tetrachloroethane	ppbv	<b>5</b>	89	70-130				
Nonane	ppbv	<b>5</b>	76	70-130				
Isopropylbenzene	$\operatorname{ppbv}$	<b>5</b>	96	70-130				
2-Chlorotoluene	$\operatorname{ppbv}$	<b>5</b>	92	70-130				
Propylbenzene	$\operatorname{ppbv}$	<b>5</b>	89	70-130				
4-Ethyltoluene	$\operatorname{ppbv}$	<b>5</b>	89	70-130				
m,p-Xylene	$\operatorname{ppbv}$	10	93	70-130				
o-Xylene	$\operatorname{ppbv}$	<b>5</b>	95	70-130				
Styrene	$\operatorname{ppbv}$	<b>5</b>	94	70-130				
Bromoform	$\operatorname{ppbv}$	<b>5</b>	91	70-130				
Benzyl chloride	$\operatorname{ppbv}$	<b>5</b>	102	70-130				
1,3,5-Trimethylbenzene	$\operatorname{ppbv}$	<b>5</b>	90	70-130				
1,2,4-Trimethylbenzene	$\operatorname{ppbv}$	<b>5</b>	81	70-130				
1,3-Dichlorobenzene	$\operatorname{ppbv}$	<b>5</b>	90	70-130				
1,4-Dichlorobenzene	$\operatorname{ppbv}$	<b>5</b>	92	70-130				
1,2-Dichlorobenzene	$\operatorname{ppbv}$	<b>5</b>	87	70-130				
1,2,4-Trichlorobenzene	$\operatorname{ppbv}$	<b>5</b>	92	70-130				
Naphthalene	ppbv	<b>5</b>	76	70-130				
Hexachlorobutadiene	$\operatorname{ppbv}$	<b>5</b>	86	70-130				

#### ENVIRONMENTAL CHEMISTS

### **Data Qualifiers & Definitions**

a - The analyte was detected at a level less than five times the reporting limit. The RPD results may not provide reliable information on the variability of the analysis.

b - The analyte was spiked at a level that was less than five times that present in the sample. Matrix spike recoveries may not be meaningful.

ca - The calibration results for the analyte were outside of acceptance criteria. The value reported is an estimate.

c - The presence of the analyte may be due to carryover from previous sample injections.

cf - The sample was centrifuged prior to analysis.

d - The sample was diluted. Detection limits were raised and surrogate recoveries may not be meaningful.

dv - Insufficient sample volume was available to achieve normal reporting limits.

f - The sample was laboratory filtered prior to analysis.

fb - The analyte was detected in the method blank.

fc - The analyte is a common laboratory and field contaminant.

hr - The sample and duplicate were reextracted and reanalyzed. RPD results were still outside of control limits. Variability is attributed to sample inhomogeneity.

hs - Headspace was present in the container used for analysis.

ht – The analysis was performed outside the method or client-specified holding time requirement.

ip - Recovery fell outside of control limits due to sample matrix effects.

j - The analyte concentration is reported below the lowest calibration standard. The value reported is an estimate.

 ${\rm J}$  - The internal standard associated with the analyte is out of control limits. The reported concentration is an estimate.

jl - The laboratory control sample(s) percent recovery and/or RPD were out of control limits. The reported concentration should be considered an estimate.

js - The surrogate associated with the analyte is out of control limits. The reported concentration should be considered an estimate.

lc - The presence of the analyte is likely due to laboratory contamination.

L - The reported concentration was generated from a library search.

nm - The analyte was not detected in one or more of the duplicate analyses. Therefore, calculation of the RPD is not applicable.

pc - The sample was received with incorrect preservation or in a container not approved by the method. The value reported should be considered an estimate.

ve - The analyte response exceeded the valid instrument calibration range. The value reported is an estimate.

vo - The value reported fell outside the control limits established for this analyte.

x - The sample chromatographic pattern does not resemble the fuel standard used for quantitation.

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