

January 12, 2022

Mr. Travis O'Rourke Municipality of Anchorage Department of Property and Facility Management 3640 East Tudor Road, Warehouse No. 1 Anchorage, AK 99507

RE: GROUNDWATER MONITORING REPORT, 5701 NORTHWOOD DRIVE, ANCHORAGE, ALASKA; ADEC FILE NO. 2100.38.536

Dear Mr. O'Rourke:

This letter report documents the July and October 2021 groundwater monitoring activities conducted at the Municipality of Anchorage (MOA) Northwood Maintenance Facility, located at 5701 Northwood Drive, Anchorage, Alaska. The site is an Alaska Department of Environmental Conservation (ADEC) contaminated site identified as "MOA Northwood Maintenance Facility Garage Bay 6". A vicinity map is included as Figure 1.

Our work was conducted in general accordance with our July 6, 2021 *Revised Work Plan for Groundwater Monitoring*, 5701 *Northwood Drive*, *Anchorage*, *Alaska*; *ADEC File No.* 2100.38.536, which was approved by Ms. Jessica Hall of the ADEC, in a letter dated July 8, 2021.

# BACKGROUND

A dry well was removed from Garage Bay 6 of the Northwood Maintenance Facility in 2001. The facility is located within the boundaries of the closed International Airport Landfill (IAL). The on-site structure and former dry well are located on the leading edge of the closed landfill. Imported fill material was placed above the landfill and the structure and other site improvements were constructed. Based on institutional knowledge of groundwater monitoring conducted at IAL, the regional groundwater flow is towards the west/southwest. Connors Lake is located about 600 feet southwest of the former dry well. According to the Alaska Department of Natural Resources Well Log Tracking System (WELTS) database, the closest downgradient drinking water well is located about 2,700 feet southwest of the site. The database also lists several potential well locations in the vicinity of the site, but it is unlikely that the majority of these wells exist due to the location of the closed landfill and wetlands.



The dry well was connected to the building's floor drain system. The dry well consisted of a 4-foot diameter concrete pipe with an open bottom. A pipe in the wall of the dry well discharged to a drainage ditch west of the facility. Following removal of the dry well, the pipe no longer discharges to the drainage ditch. Soil samples collected from the dry well excavation contained concentrations of gasoline range organics (GRO) (maximum of 376 milligrams per kilogram [mg/kg]), diesel range organics (DRO) (maximum of 7,610 mg/kg), ethylbenzene (maximum of 0.315 mg/kg), xylenes (maximum of 5.58 mg/kg), cis-1,2-dichloroethene (maximum of 0.15 mg/kg), 1,2,4-trimethylbenzene (maximum of 30.2 mg/kg), 1,3,5-trimethylbenzene (maximum of 13.7 mg/kg), 1,2-dichlorbenzene (maximum of 8.74 mg/kg) exceeding the current ADEC Method Two cleanup levels.

In December 2018, one soil boring (Boring B1) was advanced and one groundwater monitoring well (Monitoring Well B1MW) was installed adjacent to the former dry well. The approximate location of Boring B1/Well B1MW is shown on Figure 2. A soil sample collected from Boring B1 contained an estimated (J-flagged) concentration of DRO (12.8 mg/kg) less than the applicable ADEC Method Two cleanup level of 250 mg/kg. The soil sample also contained concentrations of residual range organics (RRO) (67.6 mg/kg) and 1,2-dichlorobenzene (0.0209 mg/kg) less than the ADEC Method Two cleanup levels of 10,000 mg/kg and 2.4 mg/kg, respectively. A groundwater sample collected form Well B1MW contained vinyl chloride (0. 72 micrograms per liter [ $\mu$ g/L]) at a concentration exceeding the ADEC Table C cleanup level of 0. 19  $\mu$ g/L. The groundwater sample also contained 1,230  $\mu$ g/L DRO and 1,090  $\mu$ g/L RRO which are less than the ADEC Table C cleanup levels of 1,500  $\mu$ g/L and 1,100  $\mu$ g/L, respectively. Benzene, toluene, xylenes, chloroethane, cis-1-2, dichloroethane, and dichlorodifluoromethane were also detected in the sample at concentrations less than the ADEC Table C cleanup levels.

# FIELD ACTIVITIES

The field activities consisted of collecting analytical groundwater samples and managing investigation-derived waste (IDW). SGS North America Inc. (SGS) provided analytical testing of the groundwater samples. A site plan is included as Figure 2. Copies of field notes are included in Attachment 1. The groundwater sampling activities were conducted by Ms. Schylar Healy and Mr. Zach Thon of Shannon & Wilson. Ms. Healy and Mr. Thon are Qualified Environmental Professionals in accordance with 18 Alaska Administrative Code (AAC) 75.333.



# Monitoring Well Sampling

Monitoring Well B1MW1 was sampled on July 9 and October 7, 2021. The non-disposable equipment that came into contact with groundwater was decontaminated prior to use with a non-phosphate detergent wash, a tap water rinse, and a distilled-water rinse. Prior to sampling, depth-to-water, product, and total well depth measurements were recorded using a dual phase product meter and electronic water level meter. Product was not encountered in the well.

During each event, the well was purged and sampled using a low-flow technique, with a submersible pump and disposable tubing. The submersible pump was placed within 2 feet of the surface of the groundwater column. The pump rate was adjusted with a goal of limiting the sustained water drawdown to a maximum of 0.5 foot (typical pump rate of 0.1 to 0.4 liter per minute).

During the purging process, field personnel monitored water quality parameters (temperature, specific conductivity, pH, and turbidity), purge volume, and drawdown at 3-to 5-minute intervals. The water quality instruments were calibrated prior to use, using the manufacturer's instructions. Stabilization criteria comprised of three successive readings of: temperature within 3 percent (minimum 0.2 degree Celsius), pH within 0.1-unit, specific conductivity within 3 percent, and turbidity within 10 percent or less than 10 Nephelometric Turbidity Units (NTU). Purging was considered complete when at least one well volume was removed, and the water quality parameters stabilized. A sheen was not observed on the purge water.

Analytical samples were collected by transferring water directly from the pump tubing into the laboratory supplied containers, which were the correct volumes and contain the proper preservatives, for analysis. The sample jars were filled in decreasing order of volatility. The results of the field measurements and purging data are presented in Table 1.

# Investigation Derived Waste

IDW from this project consisted of purge water and decontamination water, which was containerized in a labeled, five-gallon bucket. Shannon & Wilson coordinated with US Ecology for offsite disposal/treatment of the IDW. The signed ADEC *Transport, Treatment & Disposal Approval Form for Contaminated Media* and a copy of the waste manifest are provided in Attachment 2.



# LABORATORY ANALYSIS

During each sampling event one groundwater sample and one duplicate were submitted to the laboratory. The samples were analyzed for DRO by Alaska Method (AK) 102, RRO by AK 103, volatile organic compounds (VOCs) by Environmental Protection Agency (EPA) Method 8260D, and polynuclear aromatic hydrocarbons (PAHs) by EPA Method 8270D SIM. For quality control purposes, during each sampling event a trip blank was submitted to the laboratory and analyzed for VOCs by EPA Method 8260D. The laboratory reports and completed ADEC Laboratory Data Review Checklists (LDRCs) are provided in Attachment 3. The analytical groundwater sample results are summarized in Tables 2A and 2B. A summary of historical groundwater sample results is included in Table 3.

# DISCUSSION OF RESULTS

The analytical groundwater results were compared to the ADEC cleanup levels presented in the June 2021, 18 AAC 75 regulations. The applicable groundwater cleanup levels are established in Table C of 18 AAC 75.345.

# July 2021 Samples

Vinyl chloride (0.381  $\mu$ g/L]) was measured in Sample B1MW at a concentration exceeding the ADEC Table C cleanup level of 0.19  $\mu$ g/L. DRO, RRO, benzene, and naphthalene were measured at concentrations less than the ADEC cleanup levels. The remaining tested analytes were not detected in the duplicate groundwater sample set.

## October 2021 Samples

Vinyl chloride (0.481  $\mu$ g/L) was measured in Sample B1MW at a concentration exceeding the ADEC Table C cleanup level. DRO, RRO, benzene, and cis-1,2-dichloroethene were measured at concentrations less than the ADEC cleanup levels. The remaining tested analytes were not detected in the duplicate groundwater sample set.

# **Quality Assurance Samples**

The project laboratory implements on-going quality assurance/quality control procedures to evaluate conformance to ADEC data quality objectives (DQOs). Internal laboratory controls to assess data quality for this project include surrogates, method blanks, matrix spike/matrix spike duplicates (MS/MSD), and laboratory control sample/laboratory control sample duplicates (LCS/LCSD) to assess precision, accuracy, and matrix bias. If a DQO was not



met, the project laboratory provides a brief narrative concerning the problem in the case narrative of their laboratory reports (See Attachment 3).

The method blank associated with project Sample B1MW (July 2021 sampling event) contained estimated concentrations of fluorene (0.0171 J  $\mu g/L$ ) and phenanthrene (0.0243 J  $\mu g/L$ ) less than the LOQ. Samples are flagged "B" in Table 2A when the reported sample concentration is within 10 times the reported method blank concentration. Fluorene and phenanthrene were non-detect in Sample B1MW, therefore, the sample concentrations are reported as non-detect and flagging is not required. The method blank associated with project samples B1MW and B11MW (July sampling event) contained an estimated concentration of DRO (0.337 J  $\mu g/L$ ) less than the LOQ. The concentrations of DRO detected in Samples B1MW (859  $\mu g/L$ ) and B11MW (1,020  $\mu g/L$ ) are within five times the reported method blank concentration, therefore, the sample concentrations are reported and flagged "B".

The method blank associated with Sample B1MW (October sampling event) contained estimated concentrations of 2-methylnaphthalene (0.0163 J  $\mu$ g/L) and phenanthrene (0.0214 J  $\mu$ g/L) less than the LOQ. The concentrations of 2-methylnaphthalene and phenanthrene detected in Sample B1MW and the method blank are reported at levels less than the LOQ, therefore, the sample concentrations are reported as non-detect at the LOQ and flagged "B".

The relative percent difference (RPD) between the project sample and associated duplicate results is a measure of precision affected by matrix heterogeneity, sampling technique, and laboratory analyses. The ADEC recommends an RPD of less than 50 percent for duplicate soil samples and 30 percent for duplicate groundwater samples. All of the RPDs are within the ADEC recommended DQO of 30 percent for groundwater for both sampling events.

Laboratory-prepared trip blank samples accompanied the project sample bottles from the laboratory to the site during sampling activities and back again to SGS. The trip blank from each sampling event did not contain detectable concentrations of VOCs.

Shannon & Wilson conducted a limited data assessment to review the laboratory's compliance with precision, accuracy, sensitivity, and completeness to the data quality objectives. Shannon & Wilson reviewed the SGS data deliverables and completed the ADEC's Laboratory Data Review Checklist for each data package, which is included in Attachment 3. No non-conformances that would adversely affect the data quality or usability of the data were noted, with the exceptions discussed above.



# CONCLUSIONS/RECOMMENDATIONS

The groundwater samples collected during the July and October 2021 sampling events contained vinyl chloride concentrations exceeding the ADEC cleanup levels. Vinyl chloride concentrations have exceeded ADEC cleanup levels in all three sampling events, with the highest concentration detected in the December 2018 sampling event. DRO and RRO were detected in the groundwater samples collected during each sampling event at concentrations less than ADEC cleanup levels. We recommend continuing biannual groundwater sampling of Monitoring Well B1MW, to further evaluate contaminant concentration trends.

# CLOSURE/LIMITATIONS

This report was prepared for the exclusive use of our client and their representatives in the study of this site. The findings we have presented in this report are based on the limited sampling and analyses that we conducted. They should not be construed as a definite conclusion regarding the site's groundwater quality. As a result, the sampling and analyses performed is the basis for our professional judgment as to the environmental characteristics of this site, and in no way guarantees that an agency or its staff will reach the same conclusions as Shannon & Wilson, Inc. The data presented in this report should be considered representative of the time of our site assessment. Changes in site conditions can occur over time, due to natural forces or human activity. In addition, changes in government codes, regulations, or laws may occur. Because of such changes beyond our control, our observations and interpretations for this site may need to be revised.

Shannon & Wilson has prepared the information in Attachment 4, "Important Information About Your Geotechnical/Environmental Report," to assist you and others in understanding the use and limitations of our reports.



We appreciate this opportunity to be of service. Please call Dan P. McMahon or the undersigned at 907-561-2120 with questions or comments concerning the contents of this report.

Sincerely,

SHANNON & WILSON

Chris Pepe

**Environmental Scientist** 

Enc. Tables 1, 2A, 2B, and 3; Figures 1 and 2; and Attachments 1, 2, 3, and 4

TABLE 1
GROUNDWATER MONITORING WELL SAMPLE LOG

	Monitor	ing Well		
	B1MW	B1MW		
Water Level Measuremment Data				
Date Water Level Measured	7/9/2021	10/7/2021		
Time Water Level Measured	11:28	11:39		
Measured Depth to Water (ft below TOC)	19.21	21.21		
Sampling Data				
Date Sampled	7/9/2021	10/7/2021		
Time Sampled	12:30	12:30		
Measured Depth to Water (ft below TOC)	19.21	21.21		
Total Depth of Well (ft below TOC)	27.60	27.60		
Water Column in Well (ft)	8.39	6.39		
Gallons per Foot	0.16	0.16		
Water Column Volume (gallons)	1.34	1.02		
Total Volume Pumped (gallons)	1.4	2.0		
Sampling Method	Submersible Pump	Submersible Pump		
Diameter of Well Casing	2-inch	2-inch		
Water Quality Data				
Temperature (°C)	9.20	10.29		
pH (Standard Units)	6.04	6.07		
Specific Conductivity (μS/cm)	1,060	689		
Turbidity, (NTU)	0.0	8.3		
Remarks	Duplicate Sample B11MW	Duplicate Sample B11MW		

## Notes:

Water quality parameters were measured with Horiba/YSI 556 and turbidimeter water quality instruments.

TOC = Top of casing

ft = Feet

SP = Submersible Pump

°C = Degrees Celsius

 $\mu S/cm = Microsiemens per Centimeter$ 

NTU = Nephelometric Turbidity Units

# TABLE 2A JULY 2021 GROUNDWATER SAMPLE ANALYTICAL RESULTS

			Well ID, Sample ID number, and Water Depth Feet BTOC (See Table 1)		
			Monito	ring Well	Trip Blank
Parameter Tested	Method*	Cleanup Level**	B1MW 19.21	B11MW~ 19.21	WTB -
Diesel Range Organics (DRO) - μg/L	AK 102	1,500	859 B	1,020 B	-
Residual Range Organics (RRO) - μg/L	AK 103	1,100	901	1,070	-
Volatile Organic Compounds (VOCs)					
Benzene - μg/L	EPA 8260D	4.6	0.660	-	< 0.200
Toluene - μg/L	EPA 8260D	1,100	< 0.500	-	< 0.500
Ethylbenzene - μg/L	EPA 8260D	15	< 0.500	-	< 0.500
Xylenes (total) - μg/L	EPA 8260D	190	<1.50	-	<1.50
Vinyl chloride - μg/L	EPA 8260D	0.19	0.381	-	< 0.0750
Other VOCs - µg/L	EPA 8260D	various	ND	-	ND
Polynuclear Aromatic Hydrocarbons (PAHs)					-
Naphthalene - μg/L	EPA 8270D-SIM	1.7	0.0439 J	-	-
Other PAHs - µg/L	EPA 8270D SIM	various	ND	-	-

#### Notes:

\* = See Attachment 3 for compounds tested, methods, and laboratory reporting limits
 \*\* = Groundwater cleanup levels are listed in Table C, 18 AAC 75.345 (June 2021)
 - Sample ID number preceded by "101528-" on the chain of custody form.

- Not applicable or sample not tested for this analyte

~ = Duplicate of Sample B1MW

901 = Analyte detected

| Column | C

 $\mu$ g/L = micrograms per liter BTOC = Below top of casing

B = Compound detected in method blank at an estimated concentration and may potentially affect the sample result.

ND = Analyte not detected

# TABLE 2B OCTOBER 2021 GROUNDWATER SAMPLE ANALYTICAL RESULTS

			Well ID, Sample ID number, and Water Depth Feet BTOC (See Table 1)		
			Monitor	Trip Blank	
		Cleanup	B1MW	WTB	
Parameter Tested	Method*	Level**	21.21	21.21	-
Diesel Range Organics (DRO) - μg/L	AK 102	1,500	701	609	-
Residual Range Organics (RRO) - μg/L	AK 103	1,100	752	760	-
Volatile Organic Compounds (VOCs)					
Benzene - μg/L	EPA 8260D	4.6	0.921	-	< 0.200
Toluene - μg/L	EPA 8260D	1,100	< 0.500	-	< 0.500
Ethylbenzene - μg/L	EPA 8260D	15	< 0.500	-	< 0.500
Xylenes (total) - μg/L	EPA 8260D	190	<1.50	-	<1.50
cis-1,2-Dichloroethene - μg/L	EPA 8260D	4.7	0.317 J	-	< 0.500
Vinyl chloride - μg/L	EPA 8260D	0.19	0.418	-	< 0.0750
Other VOCs - µg/L	EPA 8260D	various	ND	-	ND
Polynuclear Aromatic Hydrocarbons (PAHs)					-
2-Methylnaphthalene - μg/L	EPA 8270D-SIM	36	<0.0481 B	-	-
Phenanthrene - μg/L	EPA 8270D-SIM	170	<0.0481 B	-	-
Other PAHs - µg/L	EPA 8270D SIM	various	ND	-	-

#### Notes:

- \* = See Attachment 3 for compounds tested, methods, and laboratory reporting limits
- \*\* = Groundwater cleanup levels are listed in Table C, 18 AAC 75.345 (June 2021)
- ^ = Sample ID number preceded by "101528-" on the chain of custody form.
- Not applicable or sample not tested for this analyte
- ~ = Duplicate of Sample B1MW
- 752 = Analyte detected
- **0.418** = Reported concentration exceeds the ADEC cleanup level
- < 0.500 = Analyte not detected; laboratory limit of detection of 0.500 mg/L
- $\mu g/L$  = micrograms per liter
- BTOC = Below top of casing
  - B = Compound detected in method blank at an estimated concentration and may potentially affect the sample result.
- ND = Analyte not detected

TABLE 3
HISTORICAL WATER ANALYTICAL RESULTS

		Donth to	Parameter Tested and ADEC Cleanup Level in μg/L*			
Monitoring Well	Date	Depth to Groundwater (feet BTOC)	DRO 1,500	RRO 2,200	Vinyl Chloride 0.19	
B1MW	12/28/2018	22.28	1,230	1,090	0.720	
B1MW~	7/9/2021	19.21	1,020 B	1,070	0.381	
B1MW~	10/7/2021	21.21	701	760	0.418	

Notes:

\* = Groundwater cleanup levels are listed in Table C, 18 AAC 75.345 (June 2021)

BTOC = Below top of casing

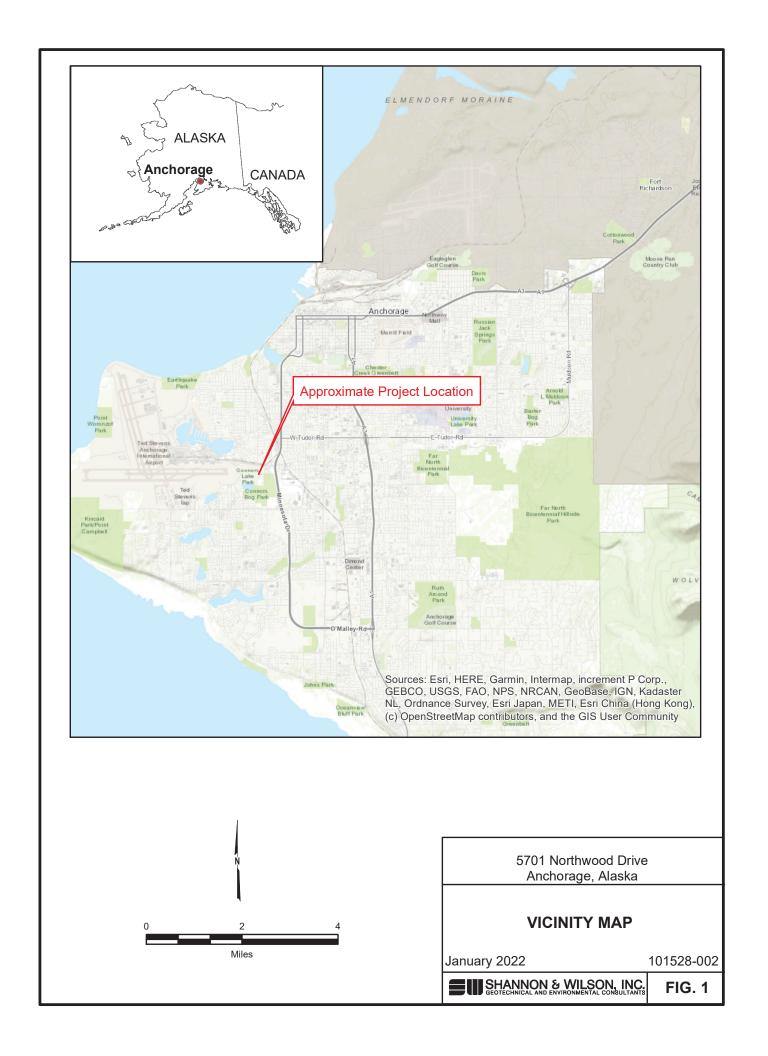
~ = Listed value based on highest concentration in duplicate sample set.

 $\mu g/L$  = micrograms per liter

701 = Analyte detected at a concentration less than the applicable ADEC cleanup level.

**0.418** = Reported concentration exceeds the ADEC cleanup level

B = Analyte concentration potentially affected by compound detected in method blank.

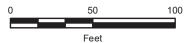






B1/B1MW

Approximate location of Monitoring Well B1MW



5701 Northwood Drive Anchorage, Alaska

# **SITE PLAN**

January 2022

101528-002

SHANNON & WILSON, INC.

GEOTECHNICAL AND ENVIRONMENTAL CONSULTANTS

FIG. 2

# SHANNON & WILSON, INC.

# **ATTACHMENT 1**

FIELD NOTES

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	Shannon & Wilson, Inc.	LOW-FLOW WA	ALER SAWIPLI	NG LUG		
	Job No: 181518	Location: 5701	bookAtrol	Weather: 55	Cloudy	
	Well No.: Brand			m: 0 1,	1. 1275	
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	Total Depth of Well Below MP:	27.40	Product Thickn	ess, if noted:	1.0	
	Depth-to-Water (DTW) Below MI		(Total Double of	CYVall Dalaw MD	DTW/Dalary	· ·
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(	Gallons in Well:	1.3767	(water Commin	III Well x Gallons	s per 100t)	·
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	Max. Drawdown (generally 0.3 ft):		Pump Rate:			
	Well Purged Dry:	Yes 🗆 No 📮		l Purged Dry Log	)	
Time:		TW Drawdown	Temp: Sp. Co.			ORP: Turb:
I IIIIC.	The second secon	SMP): (ft):	(°C) (uS/c)			(mV) (NTU)
1145	< 0.1 0.1 19	.21 0	9.07 150		5.66	0.0
1150	0.25 0.2 19	.21 0	8 79 1090		5.79	2.0
1155	0.5 0.2 17	.21	9:05 1090		555	0.0
1200		5.24 3	9.13 108	<u> </u>	5.81	0.0
1205			3.17 1030		5.89	0.0
1210	1.0 02 15	^	9.16 100		4.02	0.0
1210					2	
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		SING VOLUMES (GAI AR SPACE VOLUME (				•



# LOW-FLOW WATER SAMPLING LOG

Continued from previous page

Job No: 101828 Location: BIMW Site: 5701 Northwood Dane Well No.: 13 mm Date: 7/9/14

Time: 1215 1220 1225 1223 SAM	1.35	Pump Rate (L/min): <u>0.1</u> 0.1  3.30	DTW (ft BMP): 19.22 19.21 19.21		Temp: (°C) 9.16 9.16 5.17	Sp. Cond (uS/cm) 10 60 10 10 10 10 10 10 10 10 10 10 10 10 10	(mg/L)	pH: (S.U.) (6.09 (6.04 (6.34)	ORP: (mV)	Turb: (NTU) 0.0 0.0
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	Interval (minutes)	Pump Rate (mL/min):	Drawdown (ft):	Temp: (°C)	Sp. Cond.: (uS/cm)	DO (mg/L)	pH: (S.U.)	ORP: (mV)	Turb: (NTU)	
ADEC May 2010)	3 to 5	100 to 150	<0.0328	±3% or ±0.2	±3%	±10%	±0.1	±10	±10%	
EPA Jan. 2010)	5	50	<0.3	±3%	±3%	±10%	±0.1	±10	±10% or <5 N	NTU

EPA guidance requires all parameters to stabilize for 3 consecutive readings before sampling. If not stable within 2 hours, collect sample.

ADEC guidance requires 3 parameters (4 if using temperature) to stabilize for 3 consecutive readings before sampling.



Sampling Personnel:

# LOW-FLOW WATER SAMPLING LOG

Shannon & Wilson, Inc. Location: 5701 Northwood Job No: <u>101</u>528 Weather: Well No.: BIMW Time Completed: 13:15 Time Started: 11:35 Date: 10-7-21 Develop End Time: (24 hour break) Develop Date: INITIAL GROUNDWATER LEVEL DATA Date of Depth Measurement: \0-7-2\ 11:39 Time of Depth Measurement: Measuring Point (MP): Top of PVC Casing / Top of Steel Protective Casing / Other: Diameter of Casing: Well Screen Interval: 27.60 Product Thickness, if noted: Total Depth of Well Below MP: 21.21 Depth-to-Water (DTW) Below MP: (Total Depth of Well Below MP - DTW Below MP) Water Column in Well: 6.39 0.16 Gallons per foot: Gallons in Well: 1.02 (Water Column in Well x Gallons per foot) **PURGING DATA** Time Completed: 13:05 11:50 Time Started: Date Purged: 10-7-21 Three Well Volumes: 3.07 (Gallons in Well x 3) Depth of Pump (generally 2 ft from bottom): ~24 Gallons Purged: 2.0 Pump Rate: Max. Drawdown (generally 0.3 ft): 0.03 (If yes, use Well Purged Dry Log) Yes 🗆 No 🖭 Well Purged Dry: ORP: Temp: DO! pH: Turb: Gallons: Pump Rate DTW Drawdown Sp. Cond.: Time: (uS/cm) (mg/L) (S.U.) (mV) (NTU) (L/min): (ft BMP): (ft): (°C) (,99 5.36 21.24 8.79 11:55 0.3 0,2 0.03 10.52 9.24 696 5.56 10.50 0.2 21.24 12:00 0.03 9.78 10.28 692 12:05 0.9 0.2 21.24 0.03 5. BO. 5.92 9.11 690 21.24 0.03 10.02 12:10 0.2 21.24 5.99 8-77 1.5 10.14 690 12:15 0.2 0.08 8.12 21.24 10.22 689 6.04 12:20 0.2 0.03 10.29 689 6.07 8.25 0.03 21.24 12:25 2.0 0.2 SAMPLING DATA Odor: \_ Suffer Color: Clear 101528-BIMW Time / Date: \_\_\_\2:30 Sample Designation: 10-7-21 Time / Date: 13:00 101528 - 811MW 10-7-21 QC Sample Designation: Time / Date: QA Sample Designation: Evacuation Method: Submersible Pump / Other: Double Whole Sampling Method: Submersible Pump / Other: Double whate Water Quality Instruments Used/Manufacturer/Model Number YSI 556 + Micro TPW Calibration Info (Time, Ranges, etc) \\'\5 Full. Bring another for next Remarks: IDW bucket stored on site is Sampling Event

WELL CASING VOLUMES (GAL/FT): 1" = 0.04 2" = 0.16 4" = 0.65 ANNULAR SPACE VOLUME (GAL/FT): 4" casing and 2" well = 0.23

# SHANNON & WILSON, INC.

# **ATTACHMENT 2**

DISPOSAL DOCUMENTS



# ALASKA DEPARTMENT OF ENVIRONMENTAL CONSERVATION DIVISION OF SPILL PREVENTION AND RESPONSE

**Contaminated Sites and Prevention Preparedness and Response Programs** 

# Contaminated Media Transport and Treatment or Disposal Approval Form

DEC HAZARD/SPILL ID #	NAME OF CON	TAMIN	ATED SITE OR SPILI					
		NAME OF CONTAMINATED SITE OR SPILL						
25942		MOA I	Northwood Maintenai	nce Facility Garage Bay 6				
CONTAMINATED SITE OR S	SPILL LOCATION	N – ADI	DRESS OR OTHER AP	PROPRIATE DESCRIPTION				
	5701 North	nwood	Drive, Anchorage, Al	< 99502				
CURRENT PHYSICAL LOCATION OF MEDIA SOURCE OF THE CONTAMINATION (DAY TANK, WASH BAY, FIRE TRAINING PIT, LUST, ETC.)								
5701 Northwood Drive, Anchorage, AK 99502			Former Dry Well					
CONTAMINANTS OF CONC	ERN	ESTI	MATED VOLUME	DATE(S) GENERATED				
DRO, RRO, VOCs	, PAHs	1, 5-g	allon bucket of water	July and October 2021				
POST TREATMENT ANALY	SIS REQUIRED (S	such as (	GRO, DRO, RRO, VOCs,	metals, PFAS, and/or Chlorinated Solvents)				
			NA					
COMMENTS OR OTHER IM	PORTANT INFO	RMATI	ON					
Purge water generated while sampling Monitoring Well B1MW. Water will be processed in a wastewater treatment unit at the NRC Alaska facility (2020 Viking Drive, Anchorage, AK 99501)								

TREATMENT FACILITY, LANDFILL, AND/OR FINAL DESTINATION OF MEDIA	PHYSICAL ADDRESS/PHONE NUMBER
US Ecology	2020 Viking Drive, Anchorage, Alaska 99501 / (907) 258-1558
RESPONSIBLE PARTY	ADDRESS/PHONE NUMBER
Municipality of Anchorage/ Travis O'Rourke	3640 East Tudor Road, Warehouse No. 1 / 907-350-2784
WASTE MANAGEMENT CO. / ORGANIZER	ADDRESS/PHONE NUMBER
US Ecology	2020 Viking Drive, Anchorage, Alaska 99501 / (907) 258-1558

<sup>\*</sup>Note, disposal of polluted soil in a landfill requires prior approval from the landfill operator and ADEC Solid Waste Program.

# Alec Rizzo

Name of the Person Requesting Approval (printed)

Alec Rizzo

Signature

Digitally signed by Alec Rizzo
Date: 2021.11.18 09:38:53 -09'00'

Title/Association
11/18/21

907-561-2120

Date

Phone Number

------DEC USE ONLY------

Based on the information provided, ADEC approves transport of the above mentioned material. The Responsible Party or their consultant must submit to the DEC Project Manager a copy of weight receipts of the loads transported and a post treatment analytical report, if disposed of at an approved treatment facility. The contaminated soil shall be transported as a covered load in compliance with 18 AAC 60.015.

# Jessica Hall

DEC Project Manager Name (printed)

Jessica Hall

Digitally signed by Jessica Hall Date: 2021.11.18 10:56:19 -09'00'

Environmental Project Specialist III

Environmental Staff/ Shannon & Wilson

Project Manager Title

11/18/2021

907-269-7553

Date Phone Number

Signature

172449-MA

SENCY	CALL 800	-899-46/2 """		
NOI	N-HAZ/	ARDOUS	WASTE	<b>MANIFEST</b>

Plea	se print or type (Form designed for use on elite (1	2 pitch) typewriter)					
	NON-HAZARDOUS	1. Generator's US EPA ID No	).		Manifest		2. Page 1
	WASTE MANIFEST	VSQG			Document No.	172449A	of 1
	3. Generator's Name and Mailing Address						
	MUNICIPALITY OF ANCHORAGE	- STREET MAINT	MUNICIPALITY OF AN	CHOR	AGE - STE	REET N	
1	5701 NORTHWOOD DRIVE		5701 NORTHWOOD D	RIVE			
	ANCHORAGE, AK 99517 4. Generator's Phone (		ANCHORAGE, AK 995	517			
	5. Transporter 1 Company Name	6.	US EPA ID Number		A. State Trans	nortorio ID	
	CLIENT DELIVERED	I.	OO E! A ID Number				
	7. Transporter 2 Company Name		LIO EDA ID Mussia		B. Transporter		
400	7. Transporter 2 Company Name	8. I	US EPA ID Number		C. State Trans		
	0.00				D. Transporter		
	Designated Facility Name and Site Address	10.	US EPA ID Number		E. State Facilit	y's ID	
	US ECOLOGY ALASKA LLC						
	2020 VIKING DRIVE ANCHORAGE, AK 99501		AKR000004184		F. Facility's Ph	one 907-258-1558	
			ANK000004104			907-230-1330	
	11. WASTE DESCRIPTION			Co	ntainers	13. Total	14. Unit
F				No.	Type	Quantity	Wt./Vol.
	Hly						
	MATERIAL NOT REGULATED	BY D.O.T.			DE	42	
				1	DF	10.	P
G	b.						
E							
Й							
G M Z M Z H O K	C.						
A							
J							
9	d.						
7							
	G. Additional Descriptions for Materials Listed Above				H. Handling Co.	des for Wastes Listed Above	
				Da	_	ass for tractor Elotod / Boyo	
1	EA0301 PETROLEUM CONTAN	MINATED WATER (	DF05)	+13	9604		
		*		- 1	D391	2	
	15. Special Handling Instructions and Additional Inform	nation			0010	:U1	
	-						
	Shipper's Certification: This is to	,			9	,	
	packaged, marked and labeled, a		ndition for transportatior	n acco	rding to the	e applicable regula	itions
-	of the Department of Transportati	on					
	THE PERSON NAMED AND POST ASSESSMENT		N ANNU ANNU ANNU A	THE A	Salar Assess	ATTENDED AND ADDRESS	- AND AND AND A
-	139 ASSE ASSE ASSE ASSE ASSE					AND AND AND	
	16. GENERATOR'S CERTIFICATION: I hereby certify in proper condition for transport. The materials des	that the contents of this shipme cribed on this manifest are not s	ent are fully and accurately described a subject to federal hazardous waste req	ınd are in a ulations.	Ill respects		
			T				Date
	Printed/Typed Name		Signature			Month	Day Year
	1110		oun				18 2
T	17. Transporter 1 Acknowledgement of Receipt of Mat	erials					Date
A	Printed/Typed Name		Signature	_		Month	
S L	HEL KIZZO		apr				18 21
TRANSPORTER	18. Transporter 2 Acknowledgement of Receipt of Mat	erials					Date
Ξ.	Printed/Typed Name		Signature			Month	Day Year
Ŕ							
F	19. Discrepancy Indication Space						
A							
C							
!	20. Facility Owner or Operator: Certification of receipt of	of the waste materials covered b	y this manifest, except as noted in item	n 19.			
			. ,				Date
÷ŀ	Printed/Typed Name		Signature			Month	Day Year
γt	Day Girard		Marie (NV)		)	1)	110 101
- 1	I MINI UIII WI U						17 1

NON-HAZARDOUS WASTE

C	E	N	E	P	٨	T	0	D	

**MUNICIPALITY OF ANCHORAGE - STREET MAINT** 

**5701 NORTHWOOD DRIVE** ANCHORAGE, AK 99517

**DISPOSAL FACILITY:** 

**US ECOLOGY ALASKA LLC** 

**2020 VIKING DRIVE** 

ANCHORAGE, AK 99501

**EPA ID NUMBER:** 

**VSQG** 

MANIFEST/DOCUMENT #:

172449A

DATE OF DISPOSAL/RECYCLE: NOV-19-2021

LINE

**WASTE DESCRIPTION** 

CONTAINERS

UOM

1

PETROLEUM CONTAMINATED WATER (DF05)

I certify, on behalf of the above listed treatment facility, that to the best of my knowledge, the above described waste was managed in compliance with all applicable laws, regulations, permits and licenses on the date listed above.

PREPARED BY:

SIGNATURE:

# **ATTACHMENT 3**

# RESULTS OF ANALYTICAL TESTING BY SGS NORTH AMERICA INC. OF ANCHORAGE, ALASKA AND

ADEC LABORATORY DATA REVIEW CHECKLISTS



## **Laboratory Report of Analysis**

To: Shannon & Wilson, Inc.

5430 Fairbanks Street, Suite 3 Anchorage, AK 99518 (907)433-3223

Report Number: 1214128

Client Project: 101528-002 5701 NorthwoodDrive

Dear Dan McMahon,

Enclosed are the results of the analytical services performed under the referenced project for the received samples and associated QC as applicable. The samples are certified to meet the requirements of the National Environmental Laboratory Accreditation Conference Standards. Copies of this report and supporting data will be retained in our files for a period of ten years in the event they are required for future reference. All results are intended to be used in their entirety and SGS is not responsible for use of less than the complete report. Any samples submitted to our laboratory will be retained for a maximum of fourteen (14) days from the date of this report unless other archiving requirements were included in the quote.

If there are any questions about the report or services performed during this project, please call Justin at (907) 562-2343. We will be happy to answer any questions or concerns which you may have.

Thank you for using SGS North America Inc. for your analytical services. We look forward to working with you again on any additional analytical needs.

Sincerely,

SGS North America Inc.

Justin Nelson

2021.07.29 14:22:22 -08'00'

Justin Nelson Project Manager Justin.Nelson@sgs.com Date

Print Date: 07/29/2021 10:21:01AM Results via Engage



#### **Case Narrative**

SGS Client: **Shannon & Wilson, Inc.**SGS Project: **1214128** 

Project Name/Site: 101528-002 5701 NorthwoodDrive

Project Contact: Dan McMahon

Refer to sample receipt form for information on sample condition.

# MB for HBN 1822789 [XXX/45210] (1624808) MB

AK102 - DRO is detect in the MB greater than one-half the LOQ, but less than the LOQ.

#### 1214101019MSD (1622577) MSD

8270D SIM - PAH MS/MSD RPD for 2-methylnaphthalene does not meet QC criteria. This analyte was not detected above the LOQ in the parent sample.

\*QC comments may be associated with the field samples found in this report. When applicable, comments will be applied to associated field samples.

Print Date: 07/29/2021 10:21:03AM



#### **Laboratory Qualifiers**

Enclosed are the analytical results associated with the above work order. The results apply to the samples as received. All results are intended to be used in their entirety and SGS is not responsible for use of less than the complete report. This document is issued by the Company under its General Conditions of Service accessible at <a href="http://www.sgs.com/en/Terms-and-Conditions.aspx">http://www.sgs.com/en/Terms-and-Conditions.aspx</a>. Attention is drawn to the limitation of liability, indenmification and jurisdiction issues defined therein.

Any holder of this document is advised that information contained hereon reflects the Company's findings at the time of its intervention only and within the limits of Client's instructions, if any. The Company's sole responsibility is to its Client and this document does not exonerate parties to a transaction from exercising all their rights and obligations under the transaction documents. Any unauthorized alteration, forgery or falsification of the context or appearance of this document is unlawful and offenders may be prosecuted to the fullest extent of the law.

SGS maintains a formal Quality Assurance/Quality Control (QA/QC) program. A copy of our Quality Assurance Plan (QAP), which outlines this program, is available at your request. The laboratory certification numbers are AK00971 DW Chemistry (Provisionally Certified as of 05/27/2021 for Mercury by EPA200.8, Nitrate as N by SM 4500NO3-F and VOCs by EPA 524.2) & Microbiology & 17-021 (CS) for ADEC and 2944.01 for DOD ELAP/ISO17025 (RCRA methods: 1020B, 1311, 3010A, 3050B, 3520C, 3550C, 5030B, 5035A, 6020B, 7470A, 7471B, 8015C, 8021B, 8082A, 8260D, 8270D, 8270D-SIM, 9040C, 9045D, 9056A, 9060A, AK101 and AK102/103). SGS is only certified for the analytes listed on our Drinking Water Certification (DW methods: 200.8, 2130B, 2320B, 2510B, 300.0, 4500-CN-C,E, 4500-H-B, 4500-NO3-F, 4500-P-E and 524.2) and only those analytes will be reported to the State of Alaska for compliance. Except as specifically noted, all statements and data in this report are in conformance to the provisions set forth by the

The following descriptors or qualifiers may be found in your report:

\* The analyte has exceeded allowable regulatory or control limits.

! Surrogate out of control limits.

B Indicates the analyte is found in a blank associated with the sample.

CCV/CVA/CVB Continuing Calibration Verification
CCCV/CVC/CVCA/CVCB Closing Continuing Calibration Verification

CL Control Limit

DF Analytical Dilution Factor

DL Detection Limit (i.e., maximum method detection limit)
E The analyte result is above the calibrated range.

GT Greater Than
IB Instrument Blank

ICV Initial Calibration Verification

J The quantitation is an estimation.

LCS(D) Laboratory Control Spike (Duplicate)

LLQC/LLIQC Low Level Quantitation Check

LOD Limit of Detection (i.e., 1/2 of the LOQ)

LOQ Limit of Quantitation (i.e., reporting or practical quantitation limit)

LT Less Than MB Method Blank

MS(D) Matrix Spike (Duplicate)

ND Indicates the analyte is not detected.

RPD Relative Percent Difference
TNTC Too Numerous To Count

U Indicates the analyte was analyzed for but not detected.

Note: Sample summaries which include a result for "Total Solids" have already been adjusted for moisture content.

All DRO/RRO analyses are integrated per SOP.

Print Date: 07/29/2021 10:21:05AM

SGS North America Inc. | 200 West Potter Drive, Anchorage, AK 99518 | t 907.562.2343 f 907.561.5301 www.us.sgs.com



# **Sample Summary**

Client Sample ID	Lab Sample ID	<u>Collected</u>	Received	<u>Matrix</u>
101528-B1MW	1214128001	07/09/2021	07/09/2021	Water (Surface, Eff., Ground)
101528-B11MW	1214128002	07/09/2021	07/09/2021	Water (Surface, Eff., Ground)
101528-WTB	1214128003	07/09/2021	07/09/2021	Water (Surface, Eff., Ground)

MethodMethod Description8270D SIM LV (PAH)8270 PAH SIM GC/MS LVAK102DRO/RRO Low Volume WaterAK103DRO/RRO Low Volume WaterSW8260DVolatile Organic Compounds (W) FULL

Print Date: 07/29/2021 10:21:07AM



# **Detectable Results Summary**

Client Sample ID: 101528-B1MW			
Lab Sample ID: 1214128001	<u>Parameter</u>	Result	<u>Units</u>
Polynuclear Aromatics GC/MS	Naphthalene	0.0439J	ug/L
Semivolatile Organic Fuels	Diesel Range Organics	0.859	mg/L
	Residual Range Organics	0.901	mg/L
Volatile GC/MS	Benzene	0.660	ug/L
	Vinyl chloride	0.381	ug/L
Client Sample ID: 101528-B11MW			
Lab Sample ID: 1214128002	<u>Parameter</u>	Result	<u>Units</u>
Semivolatile Organic Fuels	Diesel Range Organics	1.02	mg/L
	Residual Range Organics	1.07	mg/L

Print Date: 07/29/2021 10:21:08AM

5 of 36



Client Sample ID: 101528-B1MW

Client Project ID: 101528-002 5701 NorthwoodDrive

Lab Sample ID: 1214128001 Lab Project ID: 1214128 Collection Date: 07/09/21 12:30 Received Date: 07/09/21 14:46 Matrix: Water (Surface, Eff., Ground)

Solids (%): Location:

# Results by Polynuclear Aromatics GC/MS

						<u>Allowable</u>	
<u>Parameter</u>	Result Qual	LOQ/CL	<u>DL</u>	<u>Units</u>	<u>DF</u>	<u>Limits</u>	Date Analyzed
1-Methylnaphthalene	0.0236 U	0.0472	0.0142	ug/L	1		07/22/21 02:50
2-Methylnaphthalene	0.0236 U	0.0472	0.0142	ug/L	1		07/22/21 02:50
Acenaphthene	0.0236 U	0.0472	0.0142	ug/L	1		07/22/21 02:50
Acenaphthylene	0.0236 U	0.0472	0.0142	ug/L	1		07/22/21 02:50
Anthracene	0.0236 U	0.0472	0.0142	ug/L	1		07/22/21 02:50
Benzo(a)Anthracene	0.0236 U	0.0472	0.0142	ug/L	1		07/22/21 02:50
Benzo[a]pyrene	0.00945 U	0.0189	0.00585	ug/L	1		07/22/21 02:50
Benzo[b]Fluoranthene	0.0236 U	0.0472	0.0142	ug/L	1		07/22/21 02:50
Benzo[g,h,i]perylene	0.0236 U	0.0472	0.0142	ug/L	1		07/22/21 02:50
Benzo[k]fluoranthene	0.0236 U	0.0472	0.0142	ug/L	1		07/22/21 02:50
Chrysene	0.0236 U	0.0472	0.0142	ug/L	1		07/22/21 02:50
Dibenzo[a,h]anthracene	0.00945 U	0.0189	0.00585	ug/L	1		07/22/21 02:50
Fluoranthene	0.0236 U	0.0472	0.0142	ug/L	1		07/22/21 02:50
Fluorene	0.0236 U	0.0472	0.0142	ug/L	1		07/22/21 02:50
Indeno[1,2,3-c,d] pyrene	0.0236 U	0.0472	0.0142	ug/L	1		07/22/21 02:50
Naphthalene	0.0439 J	0.0943	0.0292	ug/L	1		07/22/21 02:50
Phenanthrene	0.0236 U	0.0472	0.0142	ug/L	1		07/22/21 02:50
Pyrene	0.0236 U	0.0472	0.0142	ug/L	1		07/22/21 02:50
Surrogates							
2-Methylnaphthalene-d10 (surr)	71.7	42-86		%	1		07/22/21 02:50
Fluoranthene-d10 (surr)	90.6	50-97		%	1		07/22/21 02:50

#### **Batch Information**

Analytical Batch: XMS12765

Analytical Method: 8270D SIM LV (PAH)

Analyst: LAW

Analytical Date/Time: 07/22/21 02:50 Container ID: 1214128001-C Prep Batch: XXX45148
Prep Method: SW3535A
Prep Date/Time: 07/13/21 07:51
Prep Initial Wt./Vol.: 265 mL
Prep Extract Vol: 1 mL

Print Date: 07/29/2021 10:21:10AM



Client Sample ID: 101528-B1MW

Client Project ID: 101528-002 5701 NorthwoodDrive

Lab Sample ID: 1214128001 Lab Project ID: 1214128 Collection Date: 07/09/21 12:30 Received Date: 07/09/21 14:46 Matrix: Water (Surface, Eff., Ground)

Solids (%): Location:

# Results by Semivolatile Organic Fuels

						<u>Allowable</u>	
<u>Parameter</u>	Result Qual	LOQ/CL	<u>DL</u>	<u>Units</u>	<u>DF</u>	<u>Limits</u>	Date Analyzed
Diesel Range Organics	0.859	0.577	0.173	mg/L	1		07/22/21 18:04
Surrogates							
5a Androstane (surr)	96	50-150		%	1		07/22/21 18:04

#### **Batch Information**

Analytical Batch: XFC16015 Analytical Method: AK102

Analyst: IVM

Analytical Date/Time: 07/22/21 18:04 Container ID: 1214128001-A Prep Batch: XXX45210
Prep Method: SW3520C
Prep Date/Time: 07/21/21 18:36
Prep Initial Wt./Vol.: 260 mL
Prep Extract Vol: 1 mL

						<u>Allowable</u>	
<u>Parameter</u>	Result Qual	LOQ/CL	<u>DL</u>	<u>Units</u>	<u>DF</u>	<u>Limits</u>	Date Analyzed
Residual Range Organics	0.901	0.481	0.144	mg/L	1		07/22/21 18:04
Surrogates							
n-Triacontane-d62 (surr)	99.2	50-150		%	1		07/22/21 18:04

#### **Batch Information**

Analytical Batch: XFC16015 Analytical Method: AK103

Analyst: IVM

Analytical Date/Time: 07/22/21 18:04 Container ID: 1214128001-A Prep Batch: XXX45210
Prep Method: SW3520C
Prep Date/Time: 07/21/21 18:36
Prep Initial Wt./Vol.: 260 mL
Prep Extract Vol: 1 mL

Print Date: 07/29/2021 10:21:10AM



Client Sample ID: 101528-B1MW

Client Project ID: 101528-002 5701 NorthwoodDrive

Lab Sample ID: 1214128001 Lab Project ID: 1214128 Collection Date: 07/09/21 12:30 Received Date: 07/09/21 14:46 Matrix: Water (Surface, Eff., Ground)

Solids (%): Location:

# Results by Volatile GC/MS

5 .	D 110 1	1.00/01			D.F.	Allowable	5.4.4
Parameter	<u>Result Qual</u> 0.250 U	<u>LOQ/CL</u> 0.500	<u>DL</u> 0.150	<u>Units</u>	<u>DF</u> 1	<u>Limits</u>	<u>Date Analyzed</u> 07/21/21 19:04
1,1,1,2-Tetrachloroethane				ug/L			
1,1,1-Trichloroethane	0.500 U	1.00	0.310	ug/L	1		07/21/21 19:04
1,1,2,2-Tetrachloroethane	0.250 U	0.500	0.150	ug/L	1		07/21/21 19:04
1,1,2-Trichloroethane	0.200 U	0.400	0.120	ug/L	1		07/21/21 19:04
1,1-Dichloroethane	0.500 U	1.00	0.310	ug/L	1		07/21/21 19:04
1,1-Dichloroethene	0.500 U	1.00	0.310	ug/L	1		07/21/21 19:04
1,1-Dichloropropene	0.500 U	1.00	0.310	ug/L	1		07/21/21 19:04
1,2,3-Trichlorobenzene	0.500 U	1.00	0.310	ug/L	1		07/21/21 19:04
1,2,3-Trichloropropane	0.500 U	1.00	0.310	ug/L	1		07/21/21 19:04
1,2,4-Trichlorobenzene	0.500 U	1.00	0.310	ug/L	1		07/21/21 19:04
1,2,4-Trimethylbenzene	0.500 U	1.00	0.310	ug/L	1		07/21/21 19:04
1,2-Dibromo-3-chloropropane	5.00 U	10.0	3.10	ug/L	1		07/21/21 19:04
1,2-Dibromoethane	0.0375 U	0.0750	0.0180	ug/L	1		07/21/21 19:04
1,2-Dichlorobenzene	0.500 U	1.00	0.310	ug/L	1		07/21/21 19:04
1,2-Dichloroethane	0.250 U	0.500	0.150	ug/L	1		07/21/21 19:04
1,2-Dichloropropane	0.500 U	1.00	0.310	ug/L	1		07/21/21 19:04
1,3,5-Trimethylbenzene	0.500 U	1.00	0.310	ug/L	1		07/21/21 19:04
1,3-Dichlorobenzene	0.500 U	1.00	0.310	ug/L	1		07/21/21 19:04
1,3-Dichloropropane	0.250 U	0.500	0.150	ug/L	1		07/21/21 19:04
1,4-Dichlorobenzene	0.250 U	0.500	0.150	ug/L	1		07/21/21 19:04
2,2-Dichloropropane	0.500 U	1.00	0.310	ug/L	1		07/21/21 19:04
2-Butanone (MEK)	5.00 U	10.0	3.10	ug/L	1		07/21/21 19:04
2-Chlorotoluene	0.500 U	1.00	0.310	ug/L	1		07/21/21 19:04
2-Hexanone	5.00 U	10.0	3.10	ug/L	1		07/21/21 19:04
4-Chlorotoluene	0.500 U	1.00	0.310	ug/L	1		07/21/21 19:04
4-Isopropyltoluene	0.500 U	1.00	0.310	ug/L	1		07/21/21 19:04
4-Methyl-2-pentanone (MIBK)	5.00 U	10.0	3.10	ug/L	1		07/21/21 19:04
Benzene	0.660	0.400	0.120	ug/L	1		07/21/21 19:04
Bromobenzene	0.500 U	1.00	0.310	ug/L	1		07/21/21 19:04
Bromochloromethane	0.500 U	1.00	0.310	ug/L	1		07/21/21 19:04
Bromodichloromethane	0.250 U	0.500	0.150	ug/L	1		07/21/21 19:04
Bromoform	0.500 U	1.00	0.310	ug/L	1		07/21/21 19:04
Bromomethane	2.50 U	5.00	2.00	ug/L	1		07/21/21 19:04
Carbon disulfide	5.00 U	10.0	3.10	ug/L	1		07/21/21 19:04
Carbon tetrachloride	0.500 U	1.00	0.310	ug/L	1		07/21/21 19:04
Chlorobenzene	0.250 U	0.500	0.150	ug/L	1		07/21/21 19:04
Chloroethane	0.500 U	1.00	0.310	ug/L	1		07/21/21 19:04
Officiocations	0.000 0	1.00	0.010	ug/L	'		07721721 13.04

Print Date: 07/29/2021 10:21:10AM



Client Sample ID: 101528-B1MW

Client Project ID: 101528-002 5701 NorthwoodDrive

Lab Sample ID: 1214128001 Lab Project ID: 1214128 Collection Date: 07/09/21 12:30 Received Date: 07/09/21 14:46 Matrix: Water (Surface, Eff., Ground)

Solids (%): Location:

# Results by Volatile GC/MS

Parameter	Result Qual	LOQ/CL	DL	Units	DF	<u>Allowable</u> <u>Limits</u> <u>Date A</u>	nalvze
<u>-arameter</u> Chloroform	0.500 U	1.00	<u>DL</u> 0.310	ug/L	1	<u>Limits</u> <u>Date Ai</u> 07/21/2	-
Chloromethane	0.500 U	1.00	0.310	ug/L ug/L	1	07/21/2	
cis-1,2-Dichloroethene	0.500 U	1.00	0.310	ug/L ug/L	1	07/21/2	
cis-1,3-Dichloropropene	0.250 U	0.500	0.310	ug/L ug/L	1	07/21/2	
Dibromochloromethane	0.250 U	0.500	0.150	ug/L ug/L	1	07/21/2	
Dibromocnioromethane	0.500 U	1.00	0.130	J	1	07/21/2	
Disconformethane  Dischlorodifluoromethane	0.500 U	1.00	0.310	ug/L	1	07/21/2	
				ug/L			
Ethylbenzene	0.500 U	1.00	0.310	ug/L	1	07/21/2	
Freon-113	5.00 U	10.0	3.10	ug/L	1	07/21/2	
lexachlorobutadiene	0.500 U	1.00	0.310	ug/L	1	07/21/2	
sopropylbenzene (Cumene)	0.500 U	1.00	0.310	ug/L	1	07/21/2	
Methylene chloride	5.00 U	10.0	3.10	ug/L	1	07/21/2	
Methyl-t-butyl ether	5.00 U	10.0	3.10	ug/L	1	07/21/2	
Naphthalene	0.500 U	1.00	0.310	ug/L	1	07/21/2	
n-Butylbenzene	0.500 U	1.00	0.310	ug/L	1	07/21/2	
n-Propylbenzene	0.500 U	1.00	0.310	ug/L	1	07/21/2	
o-Xylene	0.500 U	1.00	0.310	ug/L	1	07/21/2	:1 19:
P & M -Xylene	1.00 U	2.00	0.620	ug/L	1	07/21/2	:1 19:
sec-Butylbenzene	0.500 U	1.00	0.310	ug/L	1	07/21/2	11 19:
Styrene	0.500 U	1.00	0.310	ug/L	1	07/21/2	:1 19:
ert-Butylbenzene	0.500 U	1.00	0.310	ug/L	1	07/21/2	:1 19:
Tetrachloroethene	0.500 U	1.00	0.310	ug/L	1	07/21/2	:1 19:
Гoluene	0.500 U	1.00	0.310	ug/L	1	07/21/2	1 19:
rans-1,2-Dichloroethene	0.500 U	1.00	0.310	ug/L	1	07/21/2	11 19:
rans-1,3-Dichloropropene	0.500 U	1.00	0.310	ug/L	1	07/21/2	<u>:</u> 1 19:
Trichloroethene	0.500 U	1.00	0.310	ug/L	1	07/21/2	1 19:
Trichlorofluoromethane	0.500 U	1.00	0.310	ug/L	1	07/21/2	1 19:
/inyl acetate	5.00 U	10.0	3.10	ug/L	1	07/21/2	<u>:</u> 1 19:
/inyl chloride	0.381	0.150	0.0500	ug/L	1	07/21/2	<u>:</u> 1 19:
Kylenes (total)	1.50 U	3.00	1.00	ug/L	1	07/21/2	<u>:</u> 1 19:
ırrogates							
1,2-Dichloroethane-D4 (surr)	103	81-118		%	1	07/21/2	<u>:</u> 1 19:
1-Bromofluorobenzene (surr)	102	85-114		%	1	07/21/2	<u>:</u> 1 19:
Foluene-d8 (surr)	99.2	89-112		%	1	07/21/2	<u>:</u> 1 19:

Print Date: 07/29/2021 10:21:10AM



Client Sample ID: 101528-B1MW

Client Project ID: 101528-002 5701 NorthwoodDrive

Lab Sample ID: 1214128001 Lab Project ID: 1214128 Collection Date: 07/09/21 12:30 Received Date: 07/09/21 14:46 Matrix: Water (Surface, Eff., Ground)

Solids (%): Location:

# Results by Volatile GC/MS

# **Batch Information**

Analytical Batch: VMS20951 Analytical Method: SW8260D

Analyst: JMG

Analytical Date/Time: 07/21/21 19:04 Container ID: 1214128001-E Prep Batch: VXX37469
Prep Method: SW5030B
Prep Date/Time: 07/21/21 06:00
Prep Initial Wt./Vol.: 5 mL
Prep Extract Vol: 5 mL

Print Date: 07/29/2021 10:21:10AM J flagging is activated



Client Sample ID: 101528-B11MW

Client Project ID: 101528-002 5701 NorthwoodDrive

Lab Sample ID: 1214128002 Lab Project ID: 1214128 Collection Date: 07/09/21 14:30 Received Date: 07/09/21 14:46 Matrix: Water (Surface, Eff., Ground)

Solids (%): Location:

# Results by Semivolatile Organic Fuels

						<u>Allowable</u>	
<u>Parameter</u>	Result Qual	LOQ/CL	<u>DL</u>	<u>Units</u>	<u>DF</u>	<u>Limits</u>	Date Analyzed
Diesel Range Organics	1.02	0.566	0.170	mg/L	1		07/22/21 18:14
Surrogates							
5a Androstane (surr)	93.5	50-150		%	1		07/22/21 18:14

#### **Batch Information**

Analytical Batch: XFC16015 Analytical Method: AK102

Analyst: IVM

Analytical Date/Time: 07/22/21 18:14 Container ID: 1214128002-A Prep Batch: XXX45210
Prep Method: SW3520C
Prep Date/Time: 07/21/21 18:36
Prep Initial Wt./Vol.: 265 mL
Prep Extract Vol: 1 mL

						<u>Allowable</u>	
<u>Parameter</u>	Result Qual	LOQ/CL	<u>DL</u>	<u>Units</u>	<u>DF</u>	<u>Limits</u>	Date Analyzed
Residual Range Organics	1.07	0.472	0.142	mg/L	1		07/22/21 18:14
Surrogates							
n-Triacontane-d62 (surr)	97.3	50-150		%	1		07/22/21 18:14

#### **Batch Information**

Analytical Batch: XFC16015 Analytical Method: AK103

Analyst: IVM

Analytical Date/Time: 07/22/21 18:14 Container ID: 1214128002-A Prep Batch: XXX45210
Prep Method: SW3520C
Prep Date/Time: 07/21/21 18:36
Prep Initial Wt./Vol.: 265 mL
Prep Extract Vol: 1 mL

Print Date: 07/29/2021 10:21:10AM



## Results of 101528-WTB

Client Sample ID: 101528-WTB

Client Project ID: 101528-002 5701 NorthwoodDrive

Lab Sample ID: 1214128003 Lab Project ID: 1214128 Collection Date: 07/09/21 11:00 Received Date: 07/09/21 14:46 Matrix: Water (Surface, Eff., Ground)

Solids (%): Location:

# Results by Volatile GC/MS

Parameter	Result Qual	LOQ/CL	<u>DL</u>	<u>Units</u>	<u>DF</u>	Allowable Limits	Date Analyzed
1,1,1,2-Tetrachloroethane	0.250 U	0.500	0.150	ug/L	1		07/21/21 16:37
1,1,1-Trichloroethane	0.500 U	1.00	0.310	ug/L	1		07/21/21 16:37
1,1,2,2-Tetrachloroethane	0.250 U	0.500	0.150	ug/L	1		07/21/21 16:37
1,1,2-Trichloroethane	0.200 U	0.400	0.120	ug/L	1		07/21/21 16:37
1,1-Dichloroethane	0.500 U	1.00	0.310	ug/L	1		07/21/21 16:37
1,1-Dichloroethene	0.500 U	1.00	0.310	ug/L	1		07/21/21 16:37
1,1-Dichloropropene	0.500 U	1.00	0.310	ug/L	1		07/21/21 16:37
1,2,3-Trichlorobenzene	0.500 U	1.00	0.310	ug/L	1		07/21/21 16:37
1,2,3-Trichloropropane	0.500 U	1.00	0.310	ug/L	1		07/21/21 16:37
1,2,4-Trichlorobenzene	0.500 U	1.00	0.310	ug/L	1		07/21/21 16:37
1,2,4-Trimethylbenzene	0.500 U	1.00	0.310	ug/L	1		07/21/21 16:37
1,2-Dibromo-3-chloropropane	5.00 U	10.0	3.10	ug/L	1		07/21/21 16:37
1,2-Dibromoethane	0.0375 U	0.0750	0.0180	ug/L	1		07/21/21 16:37
1,2-Dichlorobenzene	0.500 U	1.00	0.310	ug/L	1		07/21/21 16:37
1,2-Dichloroethane	0.250 U	0.500	0.150	ug/L	1		07/21/21 16:37
1,2-Dichloropropane	0.500 U	1.00	0.310	ug/L	1		07/21/21 16:37
1,3,5-Trimethylbenzene	0.500 U	1.00	0.310	ug/L	1		07/21/21 16:37
1,3-Dichlorobenzene	0.500 U	1.00	0.310	ug/L	1		07/21/21 16:37
1,3-Dichloropropane	0.250 U	0.500	0.150	ug/L	1		07/21/21 16:37
1,4-Dichlorobenzene	0.250 U	0.500	0.150	ug/L	1		07/21/21 16:37
2,2-Dichloropropane	0.500 U	1.00	0.310	ug/L	1		07/21/21 16:37
2-Butanone (MEK)	5.00 U	10.0	3.10	ug/L	1		07/21/21 16:37
2-Chlorotoluene	0.500 U	1.00	0.310	ug/L	1		07/21/21 16:37
2-Hexanone	5.00 U	10.0	3.10	ug/L	1		07/21/21 16:37
4-Chlorotoluene	0.500 U	1.00	0.310	ug/L	1		07/21/21 16:37
4-Isopropyltoluene	0.500 U	1.00	0.310	ug/L	1		07/21/21 16:37
4-Methyl-2-pentanone (MIBK)	5.00 U	10.0	3.10	ug/L	1		07/21/21 16:37
Benzene	0.200 U	0.400	0.120	ug/L	1		07/21/21 16:37
Bromobenzene	0.500 U	1.00	0.310	ug/L	1		07/21/21 16:37
Bromochloromethane	0.500 U	1.00	0.310	ug/L	1		07/21/21 16:37
Bromodichloromethane	0.250 U	0.500	0.150	ug/L	1		07/21/21 16:37
Bromoform	0.500 U	1.00	0.310	ug/L	1		07/21/21 16:37
Bromomethane	2.50 U	5.00	2.00	ug/L	1		07/21/21 16:37
Carbon disulfide	5.00 U	10.0	3.10	ug/L	1		07/21/21 16:37
Carbon tetrachloride	0.500 U	1.00	0.310	ug/L	1		07/21/21 16:37
Chlorobenzene	0.250 U	0.500	0.150	ug/L	1		07/21/21 16:37
Chloroethane	0.500 U	1.00	0.310	ug/L	1		07/21/21 16:37

Print Date: 07/29/2021 10:21:10AM



## Results of 101528-WTB

Client Sample ID: 101528-WTB

Client Project ID: 101528-002 5701 NorthwoodDrive

Lab Sample ID: 1214128003 Lab Project ID: 1214128 Collection Date: 07/09/21 11:00 Received Date: 07/09/21 14:46 Matrix: Water (Surface, Eff., Ground)

Solids (%): Location:

# Results by Volatile GC/MS

						Allowable	
<u>Parameter</u>	Result Qual	LOQ/CL	<u>DL</u>	<u>Units</u>	<u>DF</u>	<u>Limits</u> <u>Da</u>	<u>te Analyzed</u>
Chloroform	0.500 U	1.00	0.310	ug/L	1	07.	21/21 16:37
Chloromethane	0.500 U	1.00	0.310	ug/L	1	07.	21/21 16:37
cis-1,2-Dichloroethene	0.500 U	1.00	0.310	ug/L	1	07.	21/21 16:37
cis-1,3-Dichloropropene	0.250 U	0.500	0.150	ug/L	1	07.	21/21 16:37
Dibromochloromethane	0.250 U	0.500	0.150	ug/L	1	07.	21/21 16:37
Dibromomethane	0.500 U	1.00	0.310	ug/L	1	07.	21/21 16:37
Dichlorodifluoromethane	0.500 U	1.00	0.310	ug/L	1	07.	21/21 16:37
Ethylbenzene	0.500 U	1.00	0.310	ug/L	1	07.	21/21 16:37
Freon-113	5.00 U	10.0	3.10	ug/L	1	07.	21/21 16:37
Hexachlorobutadiene	0.500 U	1.00	0.310	ug/L	1	07.	21/21 16:37
Isopropylbenzene (Cumene)	0.500 U	1.00	0.310	ug/L	1	07.	21/21 16:37
Methylene chloride	5.00 U	10.0	3.10	ug/L	1	07.	21/21 16:37
Methyl-t-butyl ether	5.00 U	10.0	3.10	ug/L	1	07.	21/21 16:37
Naphthalene	0.500 U	1.00	0.310	ug/L	1	07.	21/21 16:37
n-Butylbenzene	0.500 U	1.00	0.310	ug/L	1	07.	21/21 16:37
n-Propylbenzene	0.500 U	1.00	0.310	ug/L	1	07.	21/21 16:37
o-Xylene	0.500 U	1.00	0.310	ug/L	1	07.	21/21 16:37
P & M -Xylene	1.00 U	2.00	0.620	ug/L	1	07.	21/21 16:37
sec-Butylbenzene	0.500 U	1.00	0.310	ug/L	1	07.	21/21 16:37
Styrene	0.500 U	1.00	0.310	ug/L	1	07.	21/21 16:37
tert-Butylbenzene	0.500 U	1.00	0.310	ug/L	1	07.	21/21 16:37
Tetrachloroethene	0.500 U	1.00	0.310	ug/L	1	07.	21/21 16:37
Toluene	0.500 U	1.00	0.310	ug/L	1	07.	21/21 16:37
trans-1,2-Dichloroethene	0.500 U	1.00	0.310	ug/L	1	07.	21/21 16:37
trans-1,3-Dichloropropene	0.500 U	1.00	0.310	ug/L	1	07.	21/21 16:37
Trichloroethene	0.500 U	1.00	0.310	ug/L	1	07.	21/21 16:37
Trichlorofluoromethane	0.500 U	1.00	0.310	ug/L	1	07.	21/21 16:37
Vinyl acetate	5.00 U	10.0	3.10	ug/L	1	07.	21/21 16:37
Vinyl chloride	0.0750 U	0.150	0.0500	ug/L	1	07.	21/21 16:37
Xylenes (total)	1.50 U	3.00	1.00	ug/L	1	07	21/21 16:37
Surrogates							
1,2-Dichloroethane-D4 (surr)	113	81-118		%	1	07	21/21 16:37
4-Bromofluorobenzene (surr)	97.9	85-114		%	1	07.	21/21 16:37
Toluene-d8 (surr)	102	89-112		%	1	07	21/21 16:37

Print Date: 07/29/2021 10:21:10AM



## Results of 101528-WTB

Client Sample ID: 101528-WTB

Client Project ID: 101528-002 5701 NorthwoodDrive

Lab Sample ID: 1214128003 Lab Project ID: 1214128 Collection Date: 07/09/21 11:00 Received Date: 07/09/21 14:46 Matrix: Water (Surface, Eff., Ground)

Solids (%): Location:

# Results by Volatile GC/MS

# **Batch Information**

Analytical Batch: VMS20950 Analytical Method: SW8260D

Analyst: JMG

Analytical Date/Time: 07/21/21 16:37 Container ID: 1214128003-A Prep Batch: VXX37468
Prep Method: SW5030B
Prep Date/Time: 07/21/21 06:00
Prep Initial Wt./Vol.: 5 mL
Prep Extract Vol: 5 mL

Print Date: 07/29/2021 10:21:10AM J flagging is activated



Blank ID: MB for HBN 1822803 [VXX/37468]

Blank Lab ID: 1624875

QC for Samples: 1214128003

Matrix: Water (Surface, Eff., Ground)

# Results by SW8260D

Doromotor	Populto	LOQ/CL	<u>DL</u>	Units
Parameter 1,1,1,2-Tetrachloroethane	Results 0.250U	0.500	<u>DL</u> 0.150	ug/L
1,1,1-Trichloroethane	0.500U	1.00	0.310	ug/L
1,1,2,2-Tetrachloroethane	0.250U	0.500	0.150	ug/L
1,1,2-Trichloroethane	0.200U	0.400	0.120	ug/L
1,1-Dichloroethane	0.500U	1.00	0.310	ug/L
1,1-Dichloroethene	0.500U	1.00	0.310	ug/L
1,1-Dichloropropene	0.500U	1.00	0.310	ug/L
1,2,3-Trichlorobenzene	0.500U	1.00	0.310	ug/L
1,2,3-Trichloropropane	0.500U	1.00	0.310	ug/L
1,2,4-Trichlorobenzene	0.500U	1.00	0.310	ug/L
1,2,4-Trimethylbenzene	0.500U	1.00	0.310	ug/L
1,2-Dibromo-3-chloropropane	5.00U	10.0	3.10	ug/L
1,2-Dibromoethane	0.0375U	0.0750	0.0180	ug/L
1,2-Dichlorobenzene	0.500U	1.00	0.310	ug/L
1,2-Dichloroethane	0.250U	0.500	0.150	ug/L
1,2-Dichloropropane	0.500U	1.00	0.310	ug/L
1,3,5-Trimethylbenzene	0.500U	1.00	0.310	ug/L
1,3-Dichlorobenzene	0.500U	1.00	0.310	ug/L
1,3-Dichloropropane	0.250U	0.500	0.150	ug/L
1,4-Dichlorobenzene	0.250U	0.500	0.150	ug/L
2,2-Dichloropropane	0.500U	1.00	0.310	ug/L
2-Butanone (MEK)	5.00U	10.0	3.10	ug/L
2-Chlorotoluene	0.500U	1.00	0.310	ug/L
2-Hexanone	5.00U	10.0	3.10	ug/L
4-Chlorotoluene	0.500U	1.00	0.310	ug/L
4-Isopropyltoluene	0.500U	1.00	0.310	ug/L
4-Methyl-2-pentanone (MIBK)	5.00U	10.0	3.10	ug/L
Benzene	0.200U	0.400	0.120	ug/L
Bromobenzene	0.500U	1.00	0.310	ug/L
Bromochloromethane	0.500U	1.00	0.310	ug/L
Bromodichloromethane	0.250U	0.500	0.150	ug/L
Bromoform	0.500U	1.00	0.310	ug/L
Bromomethane	2.50U	5.00	2.00	ug/L
Carbon disulfide	5.00U	10.0	3.10	ug/L
Carbon tetrachloride	0.500U	1.00	0.310	ug/L
Chlorobenzene	0.250U	0.500	0.150	ug/L
Chloroethane	0.500U	1.00	0.310	ug/L
Chloroform	0.500U	1.00	0.310	ug/L

Print Date: 07/29/2021 10:21:12AM



Blank ID: MB for HBN 1822803 [VXX/37468]

Blank Lab ID: 1624875

QC for Samples: 1214128003

Matrix: Water (Surface, Eff., Ground)

# Results by SW8260D

Parameter	Results	LOQ/CL	<u>DL</u>	Units
Chloromethane	0.500U	1.00	0.310	ug/L
cis-1,2-Dichloroethene	0.500U	1.00	0.310	ug/L
cis-1,3-Dichloropropene	0.250U	0.500	0.150	ug/L
Dibromochloromethane	0.250U	0.500	0.150	ug/L
Dibromomethane	0.500U	1.00	0.310	ug/L
Dichlorodifluoromethane	0.500U	1.00	0.310	ug/L
Ethylbenzene	0.500U	1.00	0.310	ug/L
Freon-113	5.00U	10.0	3.10	ug/L
Hexachlorobutadiene	0.500U	1.00	0.310	ug/L
Isopropylbenzene (Cumene)	0.500U	1.00	0.310	ug/L
Methylene chloride	5.00U	10.0	3.10	ug/L
Methyl-t-butyl ether	5.00U	10.0	3.10	ug/L
Naphthalene	0.500U	1.00	0.310	ug/L
n-Butylbenzene	0.500U	1.00	0.310	ug/L
n-Propylbenzene	0.500U	1.00	0.310	ug/L
o-Xylene	0.500U	1.00	0.310	ug/L
P & M -Xylene	1.00U	2.00	0.620	ug/L
sec-Butylbenzene	0.500U	1.00	0.310	ug/L
Styrene	0.500U	1.00	0.310	ug/L
tert-Butylbenzene	0.500U	1.00	0.310	ug/L
Tetrachloroethene	0.500U	1.00	0.310	ug/L
Toluene	0.500U	1.00	0.310	ug/L
trans-1,2-Dichloroethene	0.500U	1.00	0.310	ug/L
trans-1,3-Dichloropropene	0.500U	1.00	0.310	ug/L
Trichloroethene	0.500U	1.00	0.310	ug/L
Trichlorofluoromethane	0.500U	1.00	0.310	ug/L
Vinyl acetate	5.00U	10.0	3.10	ug/L
Vinyl chloride	0.0750U	0.150	0.0500	ug/L
Xylenes (total)	1.50U	3.00	1.00	ug/L
Surrogates				
1,2-Dichloroethane-D4 (surr)	108	81-118		%
4-Bromofluorobenzene (surr)	101	85-114		%
Toluene-d8 (surr)	102	89-112		%

Print Date: 07/29/2021 10:21:12AM



Blank ID: MB for HBN 1822803 [VXX/37468]

Blank Lab ID: 1624875

QC for Samples: 1214128003

Matrix: Water (Surface, Eff., Ground)

#### Results by SW8260D

<u>Parameter</u> <u>Results</u> <u>LOQ/CL</u> <u>DL</u> <u>Units</u>

#### **Batch Information**

Analytical Batch: VMS20950 Analytical Method: SW8260D Instrument: VPA 780/5975 GC/MS

Analyst: JMG

Analytical Date/Time: 7/21/2021 1:38:00PM

Prep Batch: VXX37468 Prep Method: SW5030B

Prep Date/Time: 7/21/2021 6:00:00AM

Prep Initial Wt./Vol.: 5 mL Prep Extract Vol: 5 mL

Print Date: 07/29/2021 10:21:12AM



Blank Spike ID: LCS for HBN 1214128 [VXX37468]

Blank Spike Lab ID: 1624876 Date Analyzed: 07/21/2021 13:53

QC for Samples: 1214128003

Spike Duplicate ID: LCSD for HBN 1214128

[VXX37468]

Spike Duplicate Lab ID: 1624877 Matrix: Water (Surface, Eff., Ground)

# Results by SW8260D

		Blank Spike	e (ug/L)	:	Spike Dupli	cate (ug/L)			
<u>Parameter</u>	<u>Spike</u>	Result	Rec (%)	<u>Spike</u>	Result	Rec (%)	CL	RPD (%)	RPD CL
1,1,1,2-Tetrachloroethane	30	28.6	95	30	28.4	95	(78-124)	0.67	(< 20)
1,1,1-Trichloroethane	30	26.8	89	30	26.8	89	(74-131)	0.10	(< 20)
1,1,2,2-Tetrachloroethane	30	29.1	97	30	29.4	98	(71-121)	1.10	(< 20)
1,1,2-Trichloroethane	30	29.1	97	30	28.8	96	(80-119)	1.20	(< 20)
1,1-Dichloroethane	30	27.0	90	30	26.8	89	(77-125)	0.88	(< 20)
1,1-Dichloroethene	30	28.4	95	30	28.3	94	(71-131)	0.49	(< 20)
1,1-Dichloropropene	30	27.0	90	30	27.1	90	(79-125)	0.16	(< 20)
1,2,3-Trichlorobenzene	30	30.1	100	30	30.8	103	(69-129)	2.40	(< 20)
1,2,3-Trichloropropane	30	29.0	97	30	29.4	98	(73-122)	1.50	(< 20)
1,2,4-Trichlorobenzene	30	30.0	100	30	30.5	102	(69-130)	1.50	(< 20)
1,2,4-Trimethylbenzene	30	28.9	96	30	29.3	98	(79-124)	1.40	(< 20)
1,2-Dibromo-3-chloropropane	30	28.2	94	30	28.1	94	(62-128)	0.25	(< 20)
1,2-Dibromoethane	30	29.3	98	30	29.2	97	(77-121)	0.18	(< 20)
1,2-Dichlorobenzene	30	29.0	97	30	29.1	97	(80-119)	0.32	(< 20)
1,2-Dichloroethane	30	27.7	92	30	27.4	91	(73-128)	1.10	(< 20)
1,2-Dichloropropane	30	27.8	93	30	27.8	93	(78-122)	0.29	(< 20)
1,3,5-Trimethylbenzene	30	28.9	96	30	29.5	98	(75-124)	2.00	(< 20)
1,3-Dichlorobenzene	30	28.7	96	30	28.9	96	(80-119)	0.60	(< 20)
1,3-Dichloropropane	30	29.0	97	30	28.8	96	(80-119)	0.49	(< 20)
1,4-Dichlorobenzene	30	28.5	95	30	29.0	97	(79-118)	1.90	(< 20)
2,2-Dichloropropane	30	25.9	86	30	26.0	87	(60-139)	0.39	(< 20)
2-Butanone (MEK)	90	93.1	103	90	92.1	102	(56-143)	1.10	(< 20)
2-Chlorotoluene	30	28.6	95	30	28.5	95	(79-122)	0.22	(< 20)
2-Hexanone	90	93.6	104	90	93.1	103	(57-139)	0.50	(< 20)
4-Chlorotoluene	30	28.8	96	30	28.9	96	(78-122)	0.40	(< 20)
4-Isopropyltoluene	30	29.4	98	30	30.1	100	(77-127)	2.50	(< 20)
4-Methyl-2-pentanone (MIBK)	90	88.1	98	90	86.8	96	(67-130)	1.50	(< 20)
Benzene	30	27.1	91	30	27.0	90	(79-120)	0.35	(< 20)
Bromobenzene	30	28.1	94	30	28.4	95	(80-120)	0.97	(< 20)
Bromochloromethane	30	27.8	93	30	27.7	92	(78-123)	0.31	(< 20)
Bromodichloromethane	30	27.7	92	30	27.6	92	(79-125)	0.33	(< 20)
Bromoform	30	28.6	95	30	28.0	94	(66-130)	1.80	(< 20)
Bromomethane	30	28.5	95	30	28.6	95	(53-141)	0.35	(< 20)
Carbon disulfide	45	43.0	96	45	42.4	94	(64-133)	1.50	(< 20)

Print Date: 07/29/2021 10:21:14AM



Blank Spike ID: LCS for HBN 1214128 [VXX37468]

Blank Spike Lab ID: 1624876 Date Analyzed: 07/21/2021 13:53

QC for Samples: 1214128003

Spike Duplicate ID: LCSD for HBN 1214128

[VXX37468]

Spike Duplicate Lab ID: 1624877 Matrix: Water (Surface, Eff., Ground)

# Results by SW8260D

Parameter         Spike         Result         Rec (%)         Spike         Result         Rec (%)         Spike         Result         Rec (%)         Clambor           Carbon tetrachloride         30         26.4         88         30         26.9         90         (72-136)         1.90           Chlorobenzene         30         28.6         99         30         27.0         90         30         (79-124)         0.21           Chloromethane         30         27.0         90         30         25.8         85         (50-139)         0.87           cis-1,2-Dichloroethene         30         27.0         90         30         25.8         89         (78-123)         0.53           cis-1,3-Dichloropropene         30         27.5         92         30         28.3         94         (78-126)         0.03           Dibromoethlane         30         27.8         93         30         28.0         93         (79-123)         0.65           Dibromoethlane         30         27.8         93         30         28.0         94         (79-124)         1.00           Dibromoethlane         30         27.8         92         30         27.4 <t< th=""><th></th><th>Blank Sp</th><th>ke (ug/L)</th><th></th><th>Spike Dupli</th><th>cate (ug/L)</th><th></th><th></th><th></th></t<>		Blank Sp	ke (ug/L)		Spike Dupli	cate (ug/L)			
Chlorobenzene         30         28.3         94         30         28.0         93         (82.118)         0.87           Chloroethane         30         29.6         99         30         27.9         93         (60-138)         5.70           Chloroform         30         27.2         91         30         27.1         90         (79-124)         0.21           Chloromethane         30         25.8         86         30         25.8         85         (50-139)         0.83           cis-1,2-Dichloroptopene         30         27.5         92         30         27.5         92         (75-124)         0.07           Dibromochloromethane         30         27.6         92         30         28.0         93         (79-123)         0.65           Dibromochloromethane         30         27.6         92         30         28.0         93         (79-124)         1.00           Dibromochloromethane         30         27.6         92         30         28.1         94         (79-121)         1.60           Ethylbenzene         30         28.6         95         30         28.1         94         (79-121)         1.60 <th< th=""><th><u>er</u></th><th><u>ike</u> <u>Result</u></th><th>Rec (%)</th><th>Spike</th><th>Result</th><th>Rec (%)</th><th>CL</th><th>RPD (%)</th><th>RPD CL</th></th<>	<u>er</u>	<u>ike</u> <u>Result</u>	Rec (%)	Spike	Result	Rec (%)	CL	RPD (%)	RPD CL
Chloroethane   30   29.6   99   30   27.9   93   (60-138)   5.70	etrachloride	26.4	88	30	26.9	90	(72-136)	1.90	(< 20)
Chloroform	enzene	28.3	94	30	28.0	93	(82-118)	0.87	(< 20)
Chloromethane         30         25.8         86         30         25.6         85         (50-139)         0.87           cis-1,2-Dichloroethene         30         27.0         90         30         26.8         89         (78-123)         0.53           cis-1,3-Dichloropropene         30         27.5         92         30         27.5         92         (75-124)         0.07           Dibromochloromethane         30         28.6         95         30         28.3         94         (74-126)         1.00           Dibromomethane         30         27.6         92         30         27.4         91         (32-152)         0.66           Ethylbenzene         30         28.6         95         30         28.1         94         (79-121)         1.60           Freon-113         45         43.8         97         45         43.6         97         (70-136)         0.51           Hexachlorobutadiene         30         29.2         97         30         28.8         96         (72-131)         1.20           Methylene chloride         30         28.3         94         30         28.1         94         (74-124)         0.39	hane	29.6	99	30	27.9	93	(60-138)	5.70	(< 20)
cis-1,2-Dichloroethene         30         27.0         90         30         26.8         89         (78-123)         0.53           cis-1,3-Dichloropropene         30         27.5         92         30         27.5         92         (75-124)         0.07           Dibromochloromethane         30         28.6         95         30         28.3         94         (74-126)         1.00           Dibromochloromethane         30         27.8         93         30         28.0         93         (79-123)         0.65           Dichlorodifluoromethane         30         27.6         92         30         27.4         91         (32-152)         0.46           Ethylbenzene         30         28.6         95         30         28.1         94         (79-121)         1.60           Freon-113         45         43.8         97         45         43.6         97         (70-136)         0.51           Hexachlorobutadiene         30         29.4         98         30         30.1         100         (66-134)         22.0           Isopropylbenzene (Cumene)         30         29.3         94         30         28.1         94         (74-124)         0.69	rm	27.2	91	30	27.1	90	(79-124)	0.21	(< 20)
cis-1,3-Dichloropropene         30         27.5         92         30         27.5         92         (75-124)         0.07           Dibromochloromethane         30         28.6         95         30         28.3         94         (74-126)         1.00           Dibromomethane         30         27.8         93         30         28.0         93         (79-123)         0.65           Dichlorodifluoromethane         30         27.6         92         30         27.4         91         (32-152)         0.46           Ethylbenzene         30         28.6         95         30         28.1         94         (79-121)         1.60           Freon-113         45         43.8         97         45         43.6         97         (70-136)         0.51           Hexachlorobutadiene         30         29.4         98         30         30.1         100         (66-134)         2.20           Isopropylbenzene (Cumene)         30         29.2         97         30         28.1         94         (74-124)         0.69           Methyl-t-butyl ether         45         42.1         94         45         41.9         93         (71-124)         0.38 <td>ethane</td> <td>25.8</td> <td>86</td> <td>30</td> <td>25.6</td> <td>85</td> <td>(50-139)</td> <td>0.87</td> <td>(&lt; 20)</td>	ethane	25.8	86	30	25.6	85	(50-139)	0.87	(< 20)
Dibromochloromethane   30	oichloroethene	27.0	90	30	26.8	89	(78-123)	0.53	(< 20)
Dibromomethane   30   27.8   93   30   28.0   93   (79-123)   0.65	ichloropropene	27.5	92	30	27.5	92	(75-124)	0.07	(< 20)
Dichlorodifluoromethane   30   27.6   92   30   27.4   91   (32-152)   0.46	chloromethane	28.6	95	30	28.3	94	(74-126)	1.00	(< 20)
Ethylbenzene         30         28.6         95         30         28.1         94         (79-121)         1.60           Freon-113         45         43.8         97         45         43.6         97         (70-136)         0.51           Hexachlorobutadiene         30         29.4         98         30         30.1         100         (66-134)         2.20           Isopropylbenzene (Cumene)         30         29.2         97         30         28.8         96         (72-131)         1.20           Methylene chloride         30         28.3         94         30         28.1         94         (74-124)         0.69           Methyl-t-butyl ether         45         42.1         94         45         41.9         93         (71-124)         0.38           Naphthalene         30         28.9         96         30         30.0         100         (61-128)         3.60           n-Butylbenzene         30         29.0         97         30         29.0         97         (76-126)         0.16           o-Xylene         30         28.2         94         30         27.9         93         (78-122)         1.20           P &	methane	27.8	93	30	28.0	93	(79-123)	0.65	(< 20)
Freon-113         45         43.8         97         45         43.6         97         (70-136)         0.51           Hexachlorobutadiene         30         29.4         98         30         30.1         100         (66-134)         2.20           Isopropylbenzene (Cumene)         30         29.2         97         30         28.8         96         (72-131)         1.20           Methylene chloride         30         28.3         94         30         28.1         94         (74-124)         0.69           Methyl-t-butyl ether         45         42.1         94         45         41.9         93         (71-124)         0.38           Naphthalene         30         28.9         96         30         30.0         100         (61-128)         3.60           n-Butylbenzene         30         30.7         102         30         31.2         104         (75-128)         1.50           n-Propylbenzene         30         29.0         97         30         29.0         97         (76-126)         0.16           o-Xylene         30         28.2         94         30         27.9         93         (78-122)         1.20 <td< td=""><td>difluoromethane</td><td>27.6</td><td>92</td><td>30</td><td>27.4</td><td>91</td><td>(32-152)</td><td>0.46</td><td>(&lt; 20)</td></td<>	difluoromethane	27.6	92	30	27.4	91	(32-152)	0.46	(< 20)
Hexachlorobutadiene         30         29.4         98         30         30.1         100         (66-134)         2.20           Isopropylbenzene (Cumene)         30         29.2         97         30         28.8         96         (72-131)         1.20           Methylene chloride         30         28.3         94         30         28.1         94         (74-124)         0.69           Methyl-t-butyl ether         45         42.1         94         45         41.9         93         (71-124)         0.38           Naphthalene         30         28.9         96         30         30.0         100         (61-128)         3.60           n-Butylbenzene         30         29.0         97         30         29.0         97         (76-126)         0.16           o-Xylene         30         29.0         97         30         29.0         97         (76-126)         0.16           o-Xylene         30         28.2         94         30         27.9         93         (78-122)         1.20           P & M-Xylene         30         29.7         99         30         30.3         101         (77-126)         1.90           Styr	zene	28.6	95	30	28.1	94	(79-121)	1.60	(< 20)
Sopropylbenzene (Cumene)   30   29.2   97   30   28.8   96   (72-131)   1.20	13	43.8	97	45	43.6	97	(70-136)	0.51	(< 20)
Methylene chloride         30         28.3         94         30         28.1         94         (74-124)         0.69           Methyl-t-butyl ether         45         42.1         94         45         41.9         93         (71-124)         0.38           Naphthalene         30         28.9         96         30         30.0         100         (61-128)         3.60           n-Butylbenzene         30         30.7         102         30         31.2         104         (75-128)         1.50           n-Propylbenzene         30         29.0         97         30         29.0         97         (76-126)         0.16           o-Xylene         30         28.2         94         30         27.9         93         (78-122)         1.20           P & M -Xylene         60         57.1         95         60         55.8         93         (80-121)         2.20           sec-Butylbenzene         30         29.7         99         30         30.3         101         (77-126)         1.90           Styrene         30         29.6         99         30         29.3         98         (78-123)         1.10           tert-Butylbenzen	orobutadiene	29.4	98	30	30.1	100	(66-134)	2.20	(< 20)
Methyl-t-butyl ether         45         42.1         94         45         41.9         93         (71-124)         0.38           Naphthalene         30         28.9         96         30         30.0         100         (61-128)         3.60           n-Butylbenzene         30         30.7         102         30         31.2         104         (75-128)         1.50           n-Propylbenzene         30         29.0         97         30         29.0         97         (76-126)         0.16           o-Xylene         30         28.2         94         30         27.9         93         (78-122)         1.20           P & M -Xylene         60         57.1         95         60         55.8         93         (80-121)         2.20           sec-Butylbenzene         30         29.7         99         30         30.3         101         (77-126)         1.90           Styrene         30         29.6         99         30         29.3         98         (78-123)         1.10           tert-Butylbenzene         30         27.6         92         30         27.6         92         (74-129)         0.06           Tetrachloroethene	lbenzene (Cumene)	29.2	97	30	28.8	96	(72-131)	1.20	(< 20)
Naphthalene         30         28.9         96         30         30.0         100         (61-128)         3.60           n-Butylbenzene         30         30.7         102         30         31.2         104         (75-128)         1.50           n-Propylbenzene         30         29.0         97         30         29.0         97         (76-126)         0.16           o-Xylene         30         28.2         94         30         27.9         93         (78-122)         1.20           P & M -Xylene         60         57.1         95         60         55.8         93         (80-121)         2.20           sec-Butylbenzene         30         29.7         99         30         30.3         101         (77-126)         1.90           Styrene         30         29.6         99         30         29.3         98         (78-123)         1.10           tert-Butylbenzene         30         28.2         94         30         29.0         97         (78-124)         2.80           Tetrachloroethene         30         27.6         92         30         27.6         92         (74-129)         0.06           Toluene	ne chloride	28.3	94	30	28.1	94	(74-124)	0.69	(< 20)
n-Butylbenzene 30 30.7 102 30 31.2 104 (75-128) 1.50 n-Propylbenzene 30 29.0 97 30 29.0 97 (76-126) 0.16 o-Xylene 30 28.2 94 30 27.9 93 (78-122) 1.20 P & M -Xylene 60 57.1 95 60 55.8 93 (80-121) 2.20 sec-Butylbenzene 30 29.7 99 30 30.3 101 (77-126) 1.90 Styrene 30 29.6 99 30 29.3 98 (78-123) 1.10 tert-Butylbenzene 30 28.2 94 30 29.0 97 (78-124) 2.80 Tetrachloroethene 30 27.6 92 30 27.6 92 (74-129) 0.06 Toluene 30 27.2 91 30 27.0 90 (80-121) 0.99 trans-1,2-Dichloroethene 30 27.1 90 30 26.9 90 (75-124) 0.76 trans-1,3-Dichloropropene 30 29.4 98 30 29.4 98 (73-127) 0.18 Trichloroethene 30 27.0 90 30 26.9 90 (79-123) 0.37 Trichloroethene 30 29.1 97 30 28.8 96 (54-146) 0.85 Vinyl chloride 30 27.9 93 30 27.7 92 (58-137) 0.79	butyl ether	42.1	94	45	41.9	93	(71-124)	0.38	(< 20)
n-Propylbenzene         30         29.0         97         30         29.0         97         (76-126)         0.16           o-Xylene         30         28.2         94         30         27.9         93         (78-122)         1.20           P & M -Xylene         60         57.1         95         60         55.8         93         (80-121)         2.20           sec-Butylbenzene         30         29.7         99         30         30.3         101         (77-126)         1.90           Styrene         30         29.6         99         30         29.3         98         (78-123)         1.10           tert-Butylbenzene         30         28.2         94         30         29.0         97         (78-124)         2.80           Tetrachloroethene         30         27.6         92         30         27.6         92         (74-129)         0.06           Toluene         30         27.2         91         30         27.0         90         (80-121)         0.99           trans-1,2-Dichloroethene         30         27.1         90         30         26.9         90         (75-124)         0.18           Trichloroethene </td <td>lene</td> <td>28.9</td> <td>96</td> <td>30</td> <td>30.0</td> <td>100</td> <td>(61-128)</td> <td>3.60</td> <td>(&lt; 20)</td>	lene	28.9	96	30	30.0	100	(61-128)	3.60	(< 20)
o-Xylene         30         28.2         94         30         27.9         93         (78-122)         1.20           P & M -Xylene         60         57.1         95         60         55.8         93         (80-121)         2.20           sec-Butylbenzene         30         29.7         99         30         30.3         101         (77-126)         1.90           Styrene         30         29.6         99         30         29.3         98         (78-123)         1.10           tert-Butylbenzene         30         29.6         99         30         29.0         97         (78-124)         2.80           Tetrachloroethene         30         27.6         92         30         27.6         92         (74-129)         0.06           Toluene         30         27.2         91         30         27.0         90         (80-121)         0.99           trans-1,2-Dichloroethene         30         27.1         90         30         26.9         90         (75-124)         0.76           trans-1,3-Dichloropropene         30         27.0         90         30         26.9         90         (79-123)         0.37           Trichlo	enzene	30.7	102	30	31.2	104	(75-128)	1.50	(< 20)
P & M -Xylene       60       57.1       95       60       55.8       93       (80-121)       2.20         sec-Butylbenzene       30       29.7       99       30       30.3       101       (77-126)       1.90         Styrene       30       29.6       99       30       29.3       98       (78-123)       1.10         tert-Butylbenzene       30       28.2       94       30       29.0       97       (78-124)       2.80         Tetrachloroethene       30       27.6       92       30       27.6       92       (74-129)       0.06         Toluene       30       27.2       91       30       27.0       90       (80-121)       0.99         trans-1,2-Dichloroethene       30       27.1       90       30       26.9       90       (75-124)       0.76         trans-1,3-Dichloropropene       30       27.0       90       30       29.4       98       (73-127)       0.18         Trichloroethene       30       27.0       90       30       26.9       90       (79-123)       0.37         Trichlorofluoromethane       30       29.2       97       30       28.9       96       (65-	benzene	29.0	97	30	29.0	97	(76-126)	0.16	(< 20 )
sec-Butylbenzene         30         29.7         99         30         30.3         101         (77-126)         1.90           Styrene         30         29.6         99         30         29.3         98         (78-123)         1.10           tert-Butylbenzene         30         28.2         94         30         29.0         97         (78-124)         2.80           Tetrachloroethene         30         27.6         92         30         27.6         92         (74-129)         0.06           Toluene         30         27.2         91         30         27.0         90         (80-121)         0.99           trans-1,2-Dichloroethene         30         27.1         90         30         26.9         90         (75-124)         0.76           trans-1,3-Dichloropropene         30         29.4         98         30         29.4         98         (73-127)         0.18           Trichloroethene         30         27.0         90         30         26.9         90         (79-123)         0.37           Trichlorofluoromethane         30         29.2         97         30         28.9         96         (65-141)         1.10	:	28.2	94	30	27.9	93	(78-122)	1.20	(< 20)
Styrene         30         29.6         99         30         29.3         98         (78-123)         1.10           tert-Butylbenzene         30         28.2         94         30         29.0         97         (78-124)         2.80           Tetrachloroethene         30         27.6         92         30         27.6         92         (74-129)         0.06           Toluene         30         27.2         91         30         27.0         90         (80-121)         0.99           trans-1,2-Dichloroethene         30         27.1         90         30         26.9         90         (75-124)         0.76           trans-1,3-Dichloropropene         30         29.4         98         30         29.4         98         (73-127)         0.18           Trichloroethene         30         27.0         90         30         26.9         90         (79-123)         0.37           Trichlorofluoromethane         30         29.2         97         30         28.9         96         (65-141)         1.10           Vinyl acetate         30         29.1         97         30         28.8         96         (54-146)         0.85	(ylene	57.1	95	60	55.8	93	(80-121)	2.20	(< 20)
tert-Butylbenzene         30         28.2         94         30         29.0         97         (78-124)         2.80           Tetrachloroethene         30         27.6         92         30         27.6         92         (74-129)         0.06           Toluene         30         27.2         91         30         27.0         90         (80-121)         0.99           trans-1,2-Dichloroethene         30         27.1         90         30         26.9         90         (75-124)         0.76           trans-1,3-Dichloropropene         30         29.4         98         30         29.4         98         (73-127)         0.18           Trichloroethene         30         27.0         90         30         26.9         90         (79-123)         0.37           Trichlorofluoromethane         30         29.2         97         30         28.9         96         (65-141)         1.10           Vinyl acetate         30         27.9         93         30         27.7         92         (58-137)         0.79	lbenzene	29.7	99	30	30.3	101	(77-126)	1.90	(< 20)
Tetrachloroethene         30         27.6         92         30         27.6         92         (74-129)         0.06           Toluene         30         27.2         91         30         27.0         90         (80-121)         0.99           trans-1,2-Dichloroethene         30         27.1         90         30         26.9         90         (75-124)         0.76           trans-1,3-Dichloropropene         30         29.4         98         30         29.4         98         (73-127)         0.18           Trichloroethene         30         27.0         90         30         26.9         90         (79-123)         0.37           Trichlorofluoromethane         30         29.2         97         30         28.9         96         (65-141)         1.10           Vinyl acetate         30         29.1         97         30         28.8         96         (54-146)         0.85           Vinyl chloride         30         27.9         93         30         27.7         92         (58-137)         0.79	;	29.6	99	30	29.3	98	(78-123)	1.10	(< 20)
Toluene         30         27.2         91         30         27.0         90         (80-121)         0.99           trans-1,2-Dichloroethene         30         27.1         90         30         26.9         90         (75-124)         0.76           trans-1,3-Dichloropropene         30         29.4         98         30         29.4         98         (73-127)         0.18           Trichloroethene         30         27.0         90         30         26.9         90         (79-123)         0.37           Trichlorofluoromethane         30         29.2         97         30         28.9         96         (65-141)         1.10           Vinyl acetate         30         29.1         97         30         28.8         96         (54-146)         0.85           Vinyl chloride         30         27.9         93         30         27.7         92         (58-137)         0.79	Ibenzene	28.2	94	30	29.0	97	(78-124)	2.80	(< 20 )
trans-1,2-Dichloroethene       30       27.1       90       30       26.9       90       (75-124)       0.76         trans-1,3-Dichloropropene       30       29.4       98       30       29.4       98       (73-127)       0.18         Trichloroethene       30       27.0       90       30       26.9       90       (79-123)       0.37         Trichlorofluoromethane       30       29.2       97       30       28.9       96       (65-141)       1.10         Vinyl acetate       30       29.1       97       30       28.8       96       (54-146)       0.85         Vinyl chloride       30       27.9       93       30       27.7       92       (58-137)       0.79	proethene	27.6	92	30	27.6	92	(74-129)	0.06	(< 20 )
trans-1,3-Dichloropropene       30       29.4       98       30       29.4       98       (73-127)       0.18         Trichloroethene       30       27.0       90       30       26.9       90       (79-123)       0.37         Trichlorofluoromethane       30       29.2       97       30       28.9       96       (65-141)       1.10         Vinyl acetate       30       29.1       97       30       28.8       96       (54-146)       0.85         Vinyl chloride       30       27.9       93       30       27.7       92       (58-137)       0.79	;	27.2	91	30	27.0	90	(80-121)	0.99	(< 20)
Trichloroethene         30         27.0         90         30         26.9         90         (79-123)         0.37           Trichlorofluoromethane         30         29.2         97         30         28.9         96         (65-141)         1.10           Vinyl acetate         30         29.1         97         30         28.8         96         (54-146)         0.85           Vinyl chloride         30         27.9         93         30         27.7         92         (58-137)         0.79	-Dichloroethene	27.1	90	30	26.9	90	(75-124)	0.76	(< 20)
Trichlorofluoromethane       30       29.2       97       30       28.9       96       (65-141)       1.10         Vinyl acetate       30       29.1       97       30       28.8       96       (54-146)       0.85         Vinyl chloride       30       27.9       93       30       27.7       92       (58-137)       0.79	-Dichloropropene	29.4	98	30	29.4	98	(73-127)	0.18	(< 20)
Vinyl acetate       30       29.1       97       30       28.8       96       (54-146)       0.85         Vinyl chloride       30       27.9       93       30       27.7       92       (58-137)       0.79	ethene	27.0	90	30	26.9	90	(79-123)	0.37	(< 20 )
Vinyl chloride         30         27.9         93         30         27.7         92         (58-137)         0.79	fluoromethane	29.2	97	30	28.9	96	(65-141)	1.10	(< 20 )
	etate	29.1	97	30	28.8	96	(54-146)	0.85	(< 20 )
	oride	27.9	93	30	27.7	92	(58-137)	0.79	(< 20 )
Xylenes (total) 90 85.3 95 90 83.7 93 (79-121) 1.90	(total)	85.3	95	90	83.7	93	(79-121)	1.90	(< 20 )

Print Date: 07/29/2021 10:21:14AM



Blank Spike ID: LCS for HBN 1214128 [VXX37468]

Blank Spike Lab ID: 1624876 Date Analyzed: 07/21/2021 13:53

QC for Samples: 1214128003

Spike Duplicate ID: LCSD for HBN 1214128

[VXX37468]

Spike Duplicate Lab ID: 1624877 Matrix: Water (Surface, Eff., Ground)

#### Results by SW8260D

		Blank Spil	ke (%)		Spike Dup	licate (%)			
<u>Parameter</u>	<u>Spike</u>	Result	Rec (%)	<u>Spike</u>	Result	Rec (%)	CL	RPD (%)	RPD CL
Surrogates									
1,2-Dichloroethane-D4 (surr)	30		103	30		102	(81-118)	1.60	
4-Bromofluorobenzene (surr)	30		98	30		99	(85-114)	0.90	
Toluene-d8 (surr)	30		103	30		103	(89-112)	0.09	

#### **Batch Information**

Analytical Batch: VMS20950 Analytical Method: SW8260D Instrument: VPA 780/5975 GC/MS

Analyst: JMG

Prep Batch: VXX37468
Prep Method: SW5030B

Prep Date/Time: 07/21/2021 06:00

Spike Init Wt./Vol.: 30 ug/L Extract Vol: 5 mL Dupe Init Wt./Vol.: 30 ug/L Extract Vol: 5 mL

Print Date: 07/29/2021 10:21:14AM



Blank ID: MB for HBN 1822812 [VXX/37469]

Blank Lab ID: 1624913

QC for Samples: 1214128001

Matrix: Water (Surface, Eff., Ground)

# Results by SW8260D

<u>Parameter</u>	Results	LOQ/CL	<u>DL</u>	<u>Units</u>
1,1,1,2-Tetrachloroethane	0.250U	0.500	0.150	ug/L
1,1,1-Trichloroethane	0.500U	1.00	0.310	ug/L
1,1,2,2-Tetrachloroethane	0.250U	0.500	0.150	ug/L
1,1,2-Trichloroethane	0.200U	0.400	0.120	ug/L
1,1-Dichloroethane	0.500U	1.00	0.310	ug/L
1,1-Dichloroethene	0.500U	1.00	0.310	ug/L
1,1-Dichloropropene	0.500U	1.00	0.310	ug/L
1,2,3-Trichlorobenzene	0.500U	1.00	0.310	ug/L
1,2,3-Trichloropropane	0.500U	1.00	0.310	ug/L
1,2,4-Trichlorobenzene	0.500U	1.00	0.310	ug/L
1,2,4-Trimethylbenzene	0.500U	1.00	0.310	ug/L
1,2-Dibromo-3-chloropropane	5.00U	10.0	3.10	ug/L
1,2-Dibromoethane	0.0375U	0.0750	0.0180	ug/L
1,2-Dichlorobenzene	0.500U	1.00	0.310	ug/L
1,2-Dichloroethane	0.250U	0.500	0.150	ug/L
1,2-Dichloropropane	0.500U	1.00	0.310	ug/L
1,3,5-Trimethylbenzene	0.500U	1.00	0.310	ug/L
1,3-Dichlorobenzene	0.500U	1.00	0.310	ug/L
1,3-Dichloropropane	0.250U	0.500	0.150	ug/L
1,4-Dichlorobenzene	0.250U	0.500	0.150	ug/L
2,2-Dichloropropane	0.500U	1.00	0.310	ug/L
2-Butanone (MEK)	5.00U	10.0	3.10	ug/L
2-Chlorotoluene	0.500U	1.00	0.310	ug/L
2-Hexanone	5.00U	10.0	3.10	ug/L
4-Chlorotoluene	0.500U	1.00	0.310	ug/L
4-Isopropyltoluene	0.500U	1.00	0.310	ug/L
4-Methyl-2-pentanone (MIBK)	5.00U	10.0	3.10	ug/L
Benzene	0.200U	0.400	0.120	ug/L
Bromobenzene	0.500U	1.00	0.310	ug/L
Bromochloromethane	0.500U	1.00	0.310	ug/L
Bromodichloromethane	0.250U	0.500	0.150	ug/L
Bromoform	0.500U	1.00	0.310	ug/L
Bromomethane	2.50U	5.00	2.00	ug/L
Carbon disulfide	5.00U	10.0	3.10	ug/L
Carbon tetrachloride	0.500U	1.00	0.310	ug/L
Chlorobenzene	0.250U	0.500	0.150	ug/L
Chloroethane	0.500U	1.00	0.310	ug/L
Chloroform	0.500U	1.00	0.310	ug/L

Print Date: 07/29/2021 10:21:17AM



Blank ID: MB for HBN 1822812 [VXX/37469]

Blank Lab ID: 1624913

QC for Samples: 1214128001

Matrix: Water (Surface, Eff., Ground)

# Results by SW8260D

Parameter	Results	LOQ/CL	<u>DL</u>	Units
Chloromethane	0.500U	1.00	0.310	ug/L
cis-1,2-Dichloroethene	0.500U	1.00	0.310	ug/L
cis-1,3-Dichloropropene	0.250U	0.500	0.150	ug/L
Dibromochloromethane	0.250U	0.500	0.150	ug/L
Dibromomethane	0.500U	1.00	0.310	ug/L
Dichlorodifluoromethane	0.500U	1.00	0.310	ug/L
Ethylbenzene	0.500U	1.00	0.310	ug/L
Freon-113	5.00U	10.0	3.10	ug/L
Hexachlorobutadiene	0.500U	1.00	0.310	ug/L
Isopropylbenzene (Cumene)	0.500U	1.00	0.310	ug/L
Methylene chloride	5.00U	10.0	3.10	ug/L
Methyl-t-butyl ether	5.00U	10.0	3.10	ug/L
Naphthalene	0.500U	1.00	0.310	ug/L
n-Butylbenzene	0.500U	1.00	0.310	ug/L
n-Propylbenzene	0.500U	1.00	0.310	ug/L
o-Xylene	0.500U	1.00	0.310	ug/L
P & M -Xylene	1.00U	2.00	0.620	ug/L
sec-Butylbenzene	0.500U	1.00	0.310	ug/L
Styrene	0.500U	1.00	0.310	ug/L
tert-Butylbenzene	0.500U	1.00	0.310	ug/L
Tetrachloroethene	0.500U	1.00	0.310	ug/L
Toluene	0.500U	1.00	0.310	ug/L
trans-1,2-Dichloroethene	0.500U	1.00	0.310	ug/L
trans-1,3-Dichloropropene	0.500U	1.00	0.310	ug/L
Trichloroethene	0.500U	1.00	0.310	ug/L
Trichlorofluoromethane	0.500U	1.00	0.310	ug/L
Vinyl acetate	5.00U	10.0	3.10	ug/L
Vinyl chloride	0.0750U	0.150	0.0500	ug/L
Xylenes (total)	1.50U	3.00	1.00	ug/L
Surrogates				
1,2-Dichloroethane-D4 (surr)	101	81-118		%
4-Bromofluorobenzene (surr)	102	85-114		%
Toluene-d8 (surr)	101	89-112		%

Print Date: 07/29/2021 10:21:17AM



Blank ID: MB for HBN 1822812 [VXX/37469]

Blank Lab ID: 1624913

QC for Samples: 1214128001

Matrix: Water (Surface, Eff., Ground)

# Results by SW8260D

<u>Parameter</u> <u>Results</u> <u>LOQ/CL</u> <u>DL</u> <u>Units</u>

#### **Batch Information**

Analytical Batch: VMS20951 Analytical Method: SW8260D Instrument: Agilent 7890-75MS

Analyst: JMG

Analytical Date/Time: 7/21/2021 1:37:00PM

Prep Batch: VXX37469 Prep Method: SW5030B

Prep Date/Time: 7/21/2021 6:00:00AM

Prep Initial Wt./Vol.: 5 mL Prep Extract Vol: 5 mL

Print Date: 07/29/2021 10:21:17AM



Blank Spike ID: LCS for HBN 1214128 [VXX37469]

Blank Spike Lab ID: 1624914 Date Analyzed: 07/21/2021 13:51

QC for Samples: 1214128001

Spike Duplicate ID: LCSD for HBN 1214128

[VXX37469]

Spike Duplicate Lab ID: 1624915 Matrix: Water (Surface, Eff., Ground)

# Results by SW8260D

		Blank Spike	e (ug/L)		Spike Dupli	cate (ug/L)			
<u>Parameter</u>	<u>Spike</u>	Result	Rec (%)	Spike	Result	Rec (%)	CL	RPD (%)	RPD CL
1,1,1,2-Tetrachloroethane	30	29.1	97	30	29.4	98	(78-124)	1.00	(< 20)
1,1,1-Trichloroethane	30	29.3	98	30	28.6	95	(74-131)	2.50	(< 20)
1,1,2,2-Tetrachloroethane	30	29.0	97	30	29.8	99	(71-121)	2.70	(< 20)
1,1,2-Trichloroethane	30	29.3	98	30	30.0	100	(80-119)	2.40	(< 20)
1,1-Dichloroethane	30	28.7	96	30	28.2	94	(77-125)	1.80	(< 20 )
1,1-Dichloroethene	30	29.9	100	30	28.8	96	(71-131)	3.70	(< 20 )
1,1-Dichloropropene	30	30.1	100	30	29.1	97	(79-125)	3.30	(< 20 )
1,2,3-Trichlorobenzene	30	28.0	93	30	29.0	97	(69-129)	3.50	(< 20 )
1,2,3-Trichloropropane	30	28.6	96	30	29.7	99	(73-122)	3.60	(< 20 )
1,2,4-Trichlorobenzene	30	28.8	96	30	29.4	98	(69-130)	1.80	(< 20 )
1,2,4-Trimethylbenzene	30	29.7	99	30	29.3	98	(79-124)	1.20	(< 20 )
1,2-Dibromo-3-chloropropane	30	27.4	91	30	28.9	96	(62-128)	5.20	(< 20 )
1,2-Dibromoethane	30	28.6	95	30	29.3	98	(77-121)	2.60	(< 20 )
1,2-Dichlorobenzene	30	29.4	98	30	29.6	99	(80-119)	0.75	(< 20 )
1,2-Dichloroethane	30	27.2	91	30	27.3	91	(73-128)	0.56	(< 20 )
1,2-Dichloropropane	30	28.8	96	30	28.7	96	(78-122)	0.42	(< 20 )
1,3,5-Trimethylbenzene	30	30.4	101	30	29.8	99	(75-124)	1.90	(< 20 )
1,3-Dichlorobenzene	30	29.6	99	30	29.8	99	(80-119)	0.48	(< 20 )
1,3-Dichloropropane	30	29.1	97	30	29.7	99	(80-119)	1.90	(< 20 )
1,4-Dichlorobenzene	30	29.3	98	30	29.6	99	(79-118)	1.10	(< 20 )
2,2-Dichloropropane	30	29.1	97	30	28.1	94	(60-139)	3.60	(< 20 )
2-Butanone (MEK)	90	79.4	88	90	81.4	90	(56-143)	2.50	(< 20 )
2-Chlorotoluene	30	30.0	100	30	29.6	99	(79-122)	1.40	(< 20 )
2-Hexanone	90	81.5	91	90	84.8	94	(57-139)	4.00	(< 20 )
4-Chlorotoluene	30	30.0	100	30	29.9	100	(78-122)	0.42	(< 20 )
4-Isopropyltoluene	30	30.8	103	30	30.3	101	(77-127)	1.60	(< 20 )
4-Methyl-2-pentanone (MIBK)	90	80.3	89	90	82.7	92	(67-130)	2.90	(< 20 )
Benzene	30	28.9	97	30	28.1	94	(79-120)	3.00	(< 20 )
Bromobenzene	30	29.6	99	30	29.5	98	(80-120)	0.39	(< 20 )
Bromochloromethane	30	27.9	93	30	28.0	93	(78-123)	0.40	(< 20 )
Bromodichloromethane	30	28.6	95	30	28.5	95	(79-125)	0.09	(< 20 )
Bromoform	30	28.5	95	30	29.8	100	(66-130)	4.50	(< 20 )
Bromomethane	30	27.9	93	30	28.8	96	(53-141)	3.00	(< 20 )
Carbon disulfide	45	45.0	100	45	43.3	96	(64-133)	3.90	(< 20 )

Print Date: 07/29/2021 10:21:18AM



Blank Spike ID: LCS for HBN 1214128 [VXX37469]

Blank Spike Lab ID: 1624914 Date Analyzed: 07/21/2021 13:51

QC for Samples: 1214128001

Spike Duplicate ID: LCSD for HBN 1214128

[VXX37469]

Spike Duplicate Lab ID: 1624915 Matrix: Water (Surface, Eff., Ground)

# Results by SW8260D

Parameter         Spike         Result         Rec (%)         Spike         Result         Rec (%)         CL           Carbon tetrachloride         30         29.9         100         30         28.7         96         (72-136)           Chlorobenzene         30         28.8         96         30         28.6         95         (82-118)           Chloroethane         30         35.3         118         30         30.8         103         (60-138)           Chloroform         30         27.9         93         30         27.6         92         (79-124)           Chloromethane         30         27.9         93         30         27.5         92         (50-139)           cis-1,2-Dichloroethene         30         28.3         94         30         27.8         93         (78-124)           Dibromochloromethane         30         28.5         95         30         28.7         96         (75-124)           Dibromomethane         30         27.9         93         30         28.3         94         (79-123)           Dichlorodifluoromethane         30         31.6         105         30         30.6         102         (32-152)		RPD CL (< 20)
Chlorobenzene         30         28.8         96         30         28.6         95         (82-118)           Chloroethane         30         35.3         118         30         30.8         103         (60-138)           Chloroform         30         27.9         93         30         27.6         92         (79-124)           Chloromethane         30         27.9         93         30         27.5         92         (50-139)           cis-1,2-Dichloroethene         30         28.3         94         30         27.8         93         (78-123)           cis-1,3-Dichloropropene         30         28.5         95         30         28.7         96         (75-124)           Dibromochloromethane         30         29.0         97         30         29.7         99         (74-126)           Dibromomethane         30         27.9         93         30         28.3         94         (79-123)           Dichlorodifluoromethane         30         31.6         105         30         30.6         102         (32-152)           Ethylbenzene         30         29.3         98         30         28.9         96         (79-121)		(< 20)
Chloroethane       30       35.3       118       30       30.8       103       (60-138)         Chloroform       30       27.9       93       30       27.6       92       (79-124)         Chloromethane       30       27.9       93       30       27.5       92       (50-139)         cis-1,2-Dichloroethene       30       28.3       94       30       27.8       93       (78-123)         cis-1,3-Dichloropropene       30       28.5       95       30       28.7       96       (75-124)         Dibromochloromethane       30       29.0       97       30       29.7       99       (74-126)         Dibromomethane       30       27.9       93       30       28.3       94       (79-123)         Dichlorodifluoromethane       30       31.6       105       30       30.6       102       (32-152)         Ethylbenzene       30       29.3       98       30       28.9       96       (79-121)         Freon-113       45       46.0       102       45       44.3       99       (70-136)         Hexachlorobutadiene       30       30.9       103       30       30.1       100	) 0.74	
Chloroform       30       27.9       93       30       27.6       92       (79-124)         Chloromethane       30       27.9       93       30       27.5       92       (50-139)         cis-1,2-Dichloroethene       30       28.3       94       30       27.8       93       (78-123)         cis-1,3-Dichloropropene       30       28.5       95       30       28.7       96       (75-124)         Dibromochloromethane       30       29.0       97       30       29.7       99       (74-126)         Dibromomethane       30       27.9       93       30       28.3       94       (79-123)         Dichlorodifluoromethane       30       31.6       105       30       30.6       102       (32-152)         Ethylbenzene       30       29.3       98       30       28.9       96       (79-121)         Freon-113       45       46.0       102       45       44.3       99       (70-136)         Hexachlorobutadiene       30       30.9       103       30       30.1       100       (66-134)         Isopropylbenzene (Cumene)       30       29.9       100       30       29.6 <td< td=""><td></td><td>(&lt; 20)</td></td<>		(< 20)
Chloromethane         30         27.9         93         30         27.5         92         (50-139)           cis-1,2-Dichloroethene         30         28.3         94         30         27.8         93         (78-123)           cis-1,3-Dichloropropene         30         28.5         95         30         28.7         96         (75-124)           Dibromochloromethane         30         29.0         97         30         29.7         99         (74-126)           Dibromomethane         30         27.9         93         30         28.3         94         (79-123)           Dichlorodifluoromethane         30         31.6         105         30         30.6         102         (32-152)           Ethylbenzene         30         29.3         98         30         28.9         96         (79-121)           Freon-113         45         46.0         102         45         44.3         99         (70-136)           Hexachlorobutadiene         30         30.9         103         30         30.1         100         (66-134)           Isopropylbenzene (Cumene)         30         29.9         100         30         29.6         99         (72-131) <td>) 13.70</td> <td>(&lt; 20)</td>	) 13.70	(< 20)
cis-1,2-Dichloroethene       30       28.3       94       30       27.8       93       (78-123)         cis-1,3-Dichloropropene       30       28.5       95       30       28.7       96       (75-124)         Dibromochloromethane       30       29.0       97       30       29.7       99       (74-126)         Dibromomethane       30       27.9       93       30       28.3       94       (79-123)         Dichlorodifluoromethane       30       31.6       105       30       30.6       102       (32-152)         Ethylbenzene       30       29.3       98       30       28.9       96       (79-121)         Freon-113       45       46.0       102       45       44.3       99       (70-136)         Hexachlorobutadiene       30       30.9       103       30       30.1       100       (66-134)         Isopropylbenzene (Cumene)       30       29.9       100       30       29.6       99       (72-131)	) 1.20	(< 20)
cis-1,3-Dichloropropene       30       28.5       95       30       28.7       96       (75-124)         Dibromochloromethane       30       29.0       97       30       29.7       99       (74-126)         Dibromomethane       30       27.9       93       30       28.3       94       (79-123)         Dichlorodifluoromethane       30       31.6       105       30       30.6       102       (32-152)         Ethylbenzene       30       29.3       98       30       28.9       96       (79-121)         Freon-113       45       46.0       102       45       44.3       99       (70-136)         Hexachlorobutadiene       30       30.9       103       30       30.1       100       (66-134)         Isopropylbenzene (Cumene)       30       29.9       100       30       29.6       99       (72-131)	) 1.50	(< 20)
Dibromochloromethane         30         29.0         97         30         29.7         99         ( 74-126           Dibromomethane         30         27.9         93         30         28.3         94         ( 79-123           Dichlorodifluoromethane         30         31.6         105         30         30.6         102         ( 32-152           Ethylbenzene         30         29.3         98         30         28.9         96         ( 79-121           Freon-113         45         46.0         102         45         44.3         99         ( 70-136           Hexachlorobutadiene         30         30.9         103         30         30.1         100         ( 66-134           Isopropylbenzene (Cumene)         30         29.9         100         30         29.6         99         ( 72-131	) 1.60	(< 20)
Dibromomethane       30       27.9       93       30       28.3       94       ( 79-123         Dichlorodifluoromethane       30       31.6       105       30       30.6       102       ( 32-152         Ethylbenzene       30       29.3       98       30       28.9       96       ( 79-121         Freon-113       45       46.0       102       45       44.3       99       ( 70-136         Hexachlorobutadiene       30       30.9       103       30       30.1       100       ( 66-134         Isopropylbenzene (Cumene)       30       29.9       100       30       29.6       99       ( 72-131	) 0.59	(< 20)
Dichlorodifluoromethane       30       31.6       105       30       30.6       102       ( 32-152)         Ethylbenzene       30       29.3       98       30       28.9       96       ( 79-121)         Freon-113       45       46.0       102       45       44.3       99       ( 70-136)         Hexachlorobutadiene       30       30.9       103       30       30.1       100       ( 66-134)         Isopropylbenzene (Cumene)       30       29.9       100       30       29.6       99       ( 72-131)	) 2.40	(< 20)
Ethylbenzene       30       29.3       98       30       28.9       96       (79-121)         Freon-113       45       46.0       102       45       44.3       99       (70-136)         Hexachlorobutadiene       30       30.9       103       30       30.1       100       (66-134)         Isopropylbenzene (Cumene)       30       29.9       100       30       29.6       99       (72-131)	) 1.60	(< 20)
Freon-113       45       46.0       102       45       44.3       99       (70-136)         Hexachlorobutadiene       30       30.9       103       30       30.1       100       (66-134)         Isopropylbenzene (Cumene)       30       29.9       100       30       29.6       99       (72-131)	) 3.50	(< 20 )
Hexachlorobutadiene       30       30.9       103       30       30.1       100       (66-134)         Isopropylbenzene (Cumene)       30       29.9       100       30       29.6       99       (72-131)	) 1.40	(< 20)
<b>Isopropylbenzene (Cumene)</b> 30 29.9 <b>100</b> 30 29.6 <b>99</b> (72-131	) 3.70	(< 20)
	) 2.60	(< 20 )
Methylene chloride         30         28.5         95         30         28.7         96         ( 74-124	) 1.10	(< 20 )
	) 0.62	(< 20)
Methyl-t-butyl ether 45 42.7 95 45 43.4 97 (71-124	) 1.60	(< 20 )
Naphthalene 30 26.4 <b>88</b> 30 28.1 <b>94</b> (61-128	) 6.30	(< 20 )
<b>n-Butylbenzene</b> 30 31.6 <b>105</b> 30 31.2 <b>104</b> (75-128	) 1.40	(< 20 )
<b>n-Propylbenzene</b> 30 30.9 <b>103</b> 30 30.1 <b>100</b> (76-126	) 2.50	(< 20 )
<b>o-Xylene</b> 30 29.0 <b>97</b> 30 28.7 <b>96</b> (78-122	) 0.92	(< 20 )
P & M -Xylene 60 58.0 97 60 57.1 95 (80-121	) 1.60	(< 20 )
sec-Butylbenzene         30         31.0         103         30         30.4         101         ( 77-126	) 2.10	(< 20 )
<b>Styrene</b> 30 29.2 <b>98</b> 30 29.2 <b>97</b> (78-123	) 0.11	(< 20 )
tert-Butylbenzene 30 30.6 <b>102</b> 30 29.9 <b>100</b> (78-124	) 2.20	(< 20 )
<b>Tetrachloroethene</b> 30 30.0 <b>100</b> 30 29.3 <b>98</b> (74-129	) 2.60	(< 20 )
Toluene 30 28.6 <b>95</b> 30 28.3 <b>94</b> (80-121	) 1.00	(< 20 )
trans-1,2-Dichloroethene 30 29.1 <b>97</b> 30 28.3 <b>94</b> (75-124	) 2.70	(< 20 )
trans-1,3-Dichloropropene 30 29.3 98 30 30.0 100 (73-127	) 2.30	(< 20 )
Trichloroethene 30 29.2 <b>97</b> 30 28.4 <b>95</b> (79-123	) 2.80	(< 20 )
<b>Trichlorofluoromethane</b> 30 30.7 <b>102</b> 30 29.1 <b>97</b> (65-141	) 5.30	(< 20 )
Vinyl acetate         30         28.6         95         30         29.2         97         ( 54-146)	) 2.20	(< 20 )
Vinyl chloride         30         29.3         98         30         28.3         94         ( 58-137)	) 3.40	(< 20)
<b>Xylenes (total)</b> 90 86.9 <b>97</b> 90 85.8 <b>95</b> (79-121	) 3.40	( - 20 )

Print Date: 07/29/2021 10:21:18AM



Blank Spike ID: LCS for HBN 1214128 [VXX37469]

Blank Spike Lab ID: 1624914 Date Analyzed: 07/21/2021 13:51

QC for Samples: 1214128001

Spike Duplicate ID: LCSD for HBN 1214128

[VXX37469]

Spike Duplicate Lab ID: 1624915 Matrix: Water (Surface, Eff., Ground)

#### Results by SW8260D

		Blank Spik	(e (%)		Spike Dup	licate (%)			
<u>Parameter</u>	<u>Spike</u>	Result	Rec (%)	Spike	Result	Rec (%)	CL	RPD (%)	RPD CL
Surrogates									
1,2-Dichloroethane-D4 (surr)	30		99	30		99	(81-118)	0.04	
4-Bromofluorobenzene (surr)	30		101	30		101	(85-114)	0.11	
Toluene-d8 (surr)	30		100	30		101	(89-112)	0.65	

#### **Batch Information**

Analytical Batch: VMS20951 Analytical Method: SW8260D Instrument: Agilent 7890-75MS

Analyst: JMG

Prep Batch: VXX37469
Prep Method: SW5030B

Prep Date/Time: 07/21/2021 06:00

Spike Init Wt./Vol.: 30 ug/L Extract Vol: 5 mL Dupe Init Wt./Vol.: 30 ug/L Extract Vol: 5 mL

Print Date: 07/29/2021 10:21:18AM



Blank ID: MB for HBN 1822240 [XXX/45148]

Blank Lab ID: 1622573

QC for Samples: 1214128001

Matrix: Water (Surface, Eff., Ground)

# Results by 8270D SIM LV (PAH)

<u>Parameter</u>	Results	LOQ/CL	<u>DL</u>	<u>Units</u>
1-Methylnaphthalene	0.0250U	0.0500	0.0150	ug/L
2-Methylnaphthalene	0.0250U	0.0500	0.0150	ug/L
Acenaphthene	0.0250U	0.0500	0.0150	ug/L
Acenaphthylene	0.0250U	0.0500	0.0150	ug/L
Anthracene	0.0250U	0.0500	0.0150	ug/L
Benzo(a)Anthracene	0.0250U	0.0500	0.0150	ug/L
Benzo[a]pyrene	0.0100U	0.0200	0.00620	ug/L
Benzo[b]Fluoranthene	0.0250U	0.0500	0.0150	ug/L
Benzo[g,h,i]perylene	0.0250U	0.0500	0.0150	ug/L
Benzo[k]fluoranthene	0.0250U	0.0500	0.0150	ug/L
Chrysene	0.0250U	0.0500	0.0150	ug/L
Dibenzo[a,h]anthracene	0.0100U	0.0200	0.00620	ug/L
Fluoranthene	0.0250U	0.0500	0.0150	ug/L
Fluorene	0.0171J	0.0500	0.0150	ug/L
Indeno[1,2,3-c,d] pyrene	0.0250U	0.0500	0.0150	ug/L
Naphthalene	0.0500U	0.100	0.0310	ug/L
Phenanthrene	0.0243J	0.0500	0.0150	ug/L
Pyrene	0.0250U	0.0500	0.0150	ug/L
Surrogates				
2-Methylnaphthalene-d10 (surr)	58.8	42-86		%
Fluoranthene-d10 (surr)	77.1	50-97		%

#### **Batch Information**

Analytical Batch: XMS12765

Analytical Method: 8270D SIM LV (PAH)

Instrument: Agilent GC 7890B/5977A SWA

Analyst: LAW

Analytical Date/Time: 7/21/2021 6:58:00PM

Prep Batch: XXX45148 Prep Method: SW3535A

Prep Date/Time: 7/13/2021 7:51:14AM

Prep Initial Wt./Vol.: 250 mL Prep Extract Vol: 1 mL

Print Date: 07/29/2021 10:21:21AM



Blank Spike ID: LCS for HBN 1214128 [XXX45148]

Blank Spike Lab ID: 1622574 Date Analyzed: 07/21/2021 19:19

QC for Samples: 1214128001

Spike Duplicate ID: LCSD for HBN 1214128

[XXX45148]

Spike Duplicate Lab ID: 1622575 Matrix: Water (Surface, Eff., Ground)

#### Results by 8270D SIM LV (PAH)

,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,	•	Blank Spike	e (ug/L)		Spike Dupli	cate (ug/L)			
<u>Parameter</u>	<u>Spike</u>	Result	Rec (%)	<u>Spike</u>	Result	Rec (%)	CL	RPD (%)	RPD CL
1-Methylnaphthalene	2	1.42	71	2	1.54	77	(41-115)	8.20	(< 20)
2-Methylnaphthalene	2	1.39	69	2	1.59	79	(39-114)	13.30	(< 20)
Acenaphthene	2	1.58	79	2	1.76	88	(48-114)	10.70	(< 20)
Acenaphthylene	2	1.57	78	2	1.68	84	(35-121)	6.70	(< 20)
Anthracene	2	1.53	77	2	1.61	81	(53-119)	5.20	(< 20)
Benzo(a)Anthracene	2	1.52	76	2	1.60	80	(59-120)	4.80	(< 20)
Benzo[a]pyrene	2	1.61	81	2	1.69	85	(53-120)	4.50	(< 20)
Benzo[b]Fluoranthene	2	1.55	77	2	1.65	83	(53-126)	6.80	(< 20)
Benzo[g,h,i]perylene	2	1.78	89	2	1.82	91	(44-128)	2.00	(< 20)
Benzo[k]fluoranthene	2	1.79	90	2	1.83	92	(54-125)	2.20	(< 20)
Chrysene	2	1.67	84	2	1.73	87	(57-120)	3.70	(< 20)
Dibenzo[a,h]anthracene	2	1.74	87	2	1.81	91	(44-131)	3.80	(< 20)
Fluoranthene	2	1.56	78	2	1.66	83	(58-120)	6.10	(< 20)
Fluorene	2	1.66	83	2	1.75	87	(50-118)	4.90	(< 20)
Indeno[1,2,3-c,d] pyrene	2	1.73	86	2	1.79	90	(48-130)	3.60	(< 20)
Naphthalene	2	1.37	69	2	1.50	75	(43-114)	9.00	(< 20)
Phenanthrene	2	1.67	83	2	1.80	90	(53-115)	7.60	(< 20)
Pyrene	2	1.52	76	2	1.63	81	(53-121)	7.00	(< 20 )
Surrogates									
2-Methylnaphthalene-d10 (surr)	2		68	2		78	(42-86)	12.50	
Fluoranthene-d10 (surr)	2		85	2		92	(50-97)	7.00	

#### **Batch Information**

Analytical Batch: XMS12765

Analytical Method: 8270D SIM LV (PAH)
Instrument: Agilent GC 7890B/5977A SWA

Analyst: LAW

Prep Batch: XXX45148
Prep Method: SW3535A

Prep Date/Time: 07/13/2021 07:51

Spike Init Wt./Vol.: 2 ug/L Extract Vol: 1 mL Dupe Init Wt./Vol.: 2 ug/L Extract Vol: 1 mL

Print Date: 07/29/2021 10:21:23AM



#### **Matrix Spike Summary**

Original Sample ID: 1214101019 MS Sample ID: 1622576 MS MSD Sample ID: 1622577 MSD

QC for Samples: 1214128001

Analysis Date: 07/21/2021 23:46 Analysis Date: 07/22/2021 0:06 Analysis Date: 07/22/2021 0:27 Matrix: Water (Surface, Eff., Ground)

#### Results by 8270D SIM LV (PAH)

		Matrix Spike (ug/L)		Spike Duplicate (ug/L)						
<u>Parameter</u>	<u>Sample</u>	Spike	Result	Rec (%)	Spike	Result	Rec (%)	CL	RPD (%	RPD CL
1-Methylnaphthalene	0.0255U	2.00	1.45	72	2.00	1.18	59	41-115	20.00	(< 20 )
2-Methylnaphthalene	0.0255U	2.00	1.45	72	2.00	1.16	58	39-114	21.80	* (< 20 )
Acenaphthene	0.0255U	2.00	1.56	78	2.00	1.31	65	48-114	17.20	(< 20)
Acenaphthylene	0.0255U	2.00	1.58	79	2.00	1.31	65	35-121	18.70	(< 20)
Anthracene	0.0255U	2.00	1.53	77	2.00	1.38	69	53-119	10.80	(< 20)
Benzo(a)Anthracene	0.0255U	2.00	1.56	78	2.00	1.61	80	59-120	2.90	(< 20)
Benzo[a]pyrene	0.0102U	2.00	1.65	83	2.00	1.65	83	53-120	0.06	(< 20)
Benzo[b]Fluoranthene	0.0255U	2.00	1.69	85	2.00	1.69	85	53-126	0.10	(< 20)
Benzo[g,h,i]perylene	0.0255U	2.00	1.78	89	2.00	1.79	89	44-128	0.68	(< 20)
Benzo[k]fluoranthene	0.0255U	2.00	1.7	85	2.00	1.76	88	54-125	3.40	(< 20)
Chrysene	0.0255U	2.00	1.72	86	2.00	1.69	85	57-120	1.60	(< 20)
Dibenzo[a,h]anthracene	0.0102U	2.00	1.73	87	2.00	1.79	89	44-131	3.10	(< 20)
Fluoranthene	0.0255U	2.00	1.54	77	2.00	1.56	78	58-120	1.20	(< 20)
Fluorene	0.0255U	2.00	1.61	80	2.00	1.44	72	50-118	11.00	(< 20)
Indeno[1,2,3-c,d] pyrene	0.0255U	2.00	1.72	86	2.00	1.76	88	48-130	2.40	(< 20)
Naphthalene	0.0510U	2.00	1.46	73	2.00	1.19	59	43-114	20.30	* (< 20 )
Phenanthrene	0.0255U	2.00	1.61	80	2.00	1.50	75	53-115	6.80	(< 20)
Pyrene	0.0255U	2.00	1.54	77	2.00	1.49	75	53-121	3.20	(< 20 )
Surrogates										
2-Methylnaphthalene-d10 (surr)		2.00	1.36	68	2.00	1.13	56	42-86	18.90	
Fluoranthene-d10 (surr)		2.00	1.68	84	2.00	1.66	83	50-97	1.10	

#### **Batch Information**

Analytical Batch: XMS12765

Analytical Method: 8270D SIM LV (PAH) Instrument: Agilent GC 7890B/5977A SWA

Analyst: LAW

Analytical Date/Time: 7/22/2021 12:06:00AM

Prep Batch: XXX45148

Prep Method: 3535 Solid Phase Ext for 8270 PAH SIM LV

Prep Date/Time: 7/13/2021 7:51:14AM

Prep Initial Wt./Vol.: 250.00mL Prep Extract Vol: 1.00mL

Print Date: 07/29/2021 10:21:24AM



Blank ID: MB for HBN 1822789 [XXX/45210]

Blank Lab ID: 1624808

QC for Samples:

1214128001, 1214128002

Matrix: Water (Surface, Eff., Ground)

#### Results by AK102

 Parameter
 Results
 LOQ/CL
 DL
 Units

 Diesel Range Organics
 0.337J
 0.600
 0.180
 mg/L

**Surrogates** 

5a Androstane (surr) 87.8 60-120 %

#### **Batch Information**

Analytical Batch: XFC16015 Prep Batch: XXX45210
Analytical Method: AK102 Prep Method: SW3520C

Instrument: Agilent 7890B R Prep Date/Time: 7/21/2021 6:36:22PM

Analyst: IVM Prep Initial Wt./Vol.: 250 mL Analytical Date/Time: 7/22/2021 3:34:00PM Prep Extract Vol: 1 mL

Print Date: 07/29/2021 10:21:26AM



Blank Spike ID: LCS for HBN 1214128 [XXX45210]

Blank Spike Lab ID: 1624809 Date Analyzed: 07/22/2021 15:44 Spike Duplicate ID: LCSD for HBN 1214128

[XXX45210]

Spike Duplicate Lab ID: 1624810 Matrix: Water (Surface, Eff., Ground)

QC for Samples: 1214128001, 1214128002

#### Results by AK102

	E	Blank Spike	e (mg/L)	5	Spike Dupli	cate (mg/L)			
<u>Parameter</u>	<u>Spike</u>	Result	Rec (%)	<u>Spike</u>	Result	Rec (%)	<u>CL</u>	RPD (%)	RPD CL
Diesel Range Organics	20	19.1	96	20	20.2	101	(75-125)	5.40	(< 20 )
Surrogates									
5a Androstane (surr)	0.4		105	0.4		112	(60-120)	6.60	

#### **Batch Information**

Analytical Batch: XFC16015 Analytical Method: AK102

Instrument: Agilent 7890B R

Analyst: IVM

Prep Batch: XXX45210 Prep Method: SW3520C

Prep Date/Time: 07/21/2021 18:36

Spike Init Wt./Vol.: 20 mg/L Extract Vol: 1 mL Dupe Init Wt./Vol.: 20 mg/L Extract Vol: 1 mL

Print Date: 07/29/2021 10:21:28AM



Blank ID: MB for HBN 1822789 [XXX/45210]

Blank Lab ID: 1624808

QC for Samples:

1214128001, 1214128002

Matrix: Water (Surface, Eff., Ground)

#### Results by AK103

 Parameter
 Results
 LOQ/CL
 DL
 Units

 Residual Range Organics
 0.250U
 0.500
 0.150
 mg/L

**Surrogates** 

n-Triacontane-d62 (surr) 97.3 60-120 %

#### **Batch Information**

Analytical Batch: XFC16015 Analytical Method: AK103

Instrument: Agilent 7890B R Analyst: IVM

Analytical Date/Time: 7/22/2021 3:34:00PM

Prep Batch: XXX45210 Prep Method: SW3520C

Prep Date/Time: 7/21/2021 6:36:22PM

Prep Initial Wt./Vol.: 250 mL Prep Extract Vol: 1 mL

Print Date: 07/29/2021 10:21:30AM



Blank Spike ID: LCS for HBN 1214128 [XXX45210]

Blank Spike Lab ID: 1624809 Date Analyzed: 07/22/2021 15:44

1214128001, 1214128002

Spike Duplicate ID: LCSD for HBN 1214128

[XXX45210]

Spike Duplicate Lab ID: 1624810 Matrix: Water (Surface, Eff., Ground)

# Results by AK103

QC for Samples:

-/										
l		E	Blank Spike	(mg/L)	5	Spike Duplic	cate (mg/L)			
l	<u>Parameter</u>	<u>Spike</u>	Result	Rec (%)	Spike	Result	Rec (%)	CL	RPD (%)	RPD CL
l	Residual Range Organics	20	19.0	95	20	19.9	99	(60-120)	4.50	(< 20 )
l	Surrogates									
ı	n-Triacontane-d62 (surr)	0.4		96	0.4		106	(60-120)	9.40	

#### **Batch Information**

Analytical Batch: XFC16015 Analytical Method: AK103 Instrument: Agilent 7890B R

Analyst: IVM

Prep Batch: XXX45210
Prep Method: SW3520C

Prep Date/Time: 07/21/2021 18:36

Spike Init Wt./Vol.: 20 mg/L Extract Vol: 1 mL Dupe Init Wt./Vol.: 20 mg/L Extract Vol: 1 mL

Print Date: 07/29/2021 10:21:32AM

# 1214128



P#365300XO

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Lake Oswego, OR 97035 Denve	Bannock Street, Suite 200 er, CO 80204 325-3800 Lab No.	Time	Date Sampled	Silver		20 20 20 20 20 20 20 20 20 20 20 20 20 2	S. A. S. T.	O CON S			10 h	Remark	s/Matrix
101528 - B1mw	(JAG)	12:30	7/9/21		X	X	X				7	Water	
101528 - BIIMW	(2AB)	14:30			X						2	Water	
101528 - WTB	(3AC)	11:00	↓			X					tox!	Trip Blan	<u>k</u>
Project Information	n Samj	ole Recei	pt		inquished				uished E			Relinquishe	
Project Number: 101528 - 00				Signature:		Time: <b>  </b>	<b>(</b> ∕ Sigr	nature:	Time	ə:	_ Sign	ature:	Time:
Project Name 5701 Mathematin	Received Goo			Printed N	ime:	Date: 7/9/	Print	ted Name:	Date	e:	_ Print	red Name:	Date:
Ongoing Project? Yes No				Company:	a Head	7	Com	npany:			Com	npany:	
Sampler: SAH	(attach shipping	bill, if any)			on + Wil								
Requested Turnaround Time:	structions			Rec Signature:	eived By	Time:		Receiv	red By:	<b>2.</b>		Received By ature:	7: 3. Time: <u>/4/46</u>
Special Instructions:	J IF TO LAKE			Printed Na	me:	Date:	Drint	ted Name:	——————————————————————————————————————	2'		ulle Olle ed Name:	Date: <u>7/4/21</u>
						aic.				··	Mi	chelle Albers	au Absent
Distribution: White - w/shipment - re Yellow - w/shipment - 1 Pink - Shannon & Wilse		ilson w/ labora	atory report	Company			Com	npany:			S6	ipany:	87 N65



e-Sample Receipt Form

SGS Workorder #:

1214128

1214128

Review Criteria	Condition (Yes,						
Chain of Custody / Temperature Requi	<u>irements</u>	Υ	'es	Exemption permit	ted if sample	er hand carries/deliv	ers.
Were Custody Seals intact? Note # &	location N/A	absent					
COC accompanied sa	amples? Yes						
DOD: Were samples received in COC corresponding of	coolers? N/A						
Yes **Exemption permitted if		cted <8 hou	ırs a	ago or for sample	s where chill	ing is not required	
Temperature blank compliant* (i.e., 0-6 °C after		Cooler ID:	_	1	@	8.7 °C Therm. ID:	D65
remperature blank compilant (i.e., 0-0 C and	er CF)!		-	•			<b>D</b> 00
	U. b. a	Cooler ID:			@	°C Therm. ID:	
If samples received without a temperature blank, the "cooler temperature" wil documented instead & "COOLER TEMP" will be noted to the right. "ambient" or "ch		Cooler ID:			@	°C Therm. ID:	
be noted if neither is available.		Cooler ID:			@	°C Therm. ID:	
		Cooler ID:			@	°C Therm. ID:	
*If >6°C, were samples collected <8 hours	s ago? Yes						
	<u></u>						
If <0°C, were sample containers ice	e free? N/A						
Note: Identify containers received at non-compliant tempe	rature						
Use form FS-0029 if more space is n							
·							
Holding Time / Documentation / Sample Condition Re	oguirom ente	Nata Dafan		F 000   C - 0		fin In a Latin an Airean	
Were samples received within holding		Note: Refer	10 101	rm F-083 Sample G	ulde" for speci	tic notaing times.	
were samples received within nording	g time r						
Do samples match COC** (i.e.,sample IDs,dates/times colle							
**Note: If times differ <1hr, record details & login per C	COC.						
***Note: If sample information on containers differs from COC, SGS will default to	COC information						
Were analytical requests clear? (i.e., method is specified for ar	nalyses Yes						
with multiple option for analysis (Ex: BTEX,	Metals)						
		N	I/A	***Exemption per	mitted for me	etals (e.g,200.8/602	0A).
Were proper containers (type/mass/volume/preservative***	*)used? Yes						
(7)	,	ľ					
Volatile / LL-Hg Reg	uirements						
Were Trip Blanks (i.e., VOAs, LL-Hg) in cooler with sa							
Were all water VOA vials free of headspace (i.e., bubbles ≤							
Were all soil VOAs field extracted with MeOH							
Note to Client: Any "No", answer above indicates no	on-compliance	with standa	rd p	rocedures and ma	ay impact da	ta quality.	
Additions	al notes (if a	pplicable	).				
Additions		PPHOUDIC	/•				



# **Sample Containers and Preservatives**

Container Id	<u>Preservative</u>	<u>Container</u> <u>Condition</u>	Container Id	<u>Preservative</u>	Container Condition
1214128001-A	HCL to pH < 2	ОК			
1214128001-B	HCL to pH < 2	OK			
1214128001-C	No Preservative Required	OK			
1214128001-D	No Preservative Required	OK			
1214128001-E	HCL to pH < 2	OK			
1214128001-F	HCL to pH < 2	OK			
1214128001-G	HCL to pH < 2	OK			
1214128002-A	HCL to pH < 2	OK			
1214128002-B	HCL to pH < 2	OK			
1214128003-A	HCL to pH < 2	OK			
1214128003-B	HCL to pH < 2	OK			
1214128003-C	HCL to pH < 2	ОК			

#### **Container Condition Glossary**

Containers for bacteriological, low level mercury and VOA vials are not opened prior to analysis and will be assigned condition code OK unless evidence indicates than an inappropriate container was submitted.

- OK The container was received at an acceptable pH for the analysis requested.
- BU The container was received with headspace greater than 6mm.
- DM The container was received damaged.
- FR The container was received frozen and not usable for Bacteria or BOD analyses.
- IC The container provided for microbiology analysis was not a laboratory-supplied, pre-sterilized container and therefore was not suitable for analysis.
- NC- The container provided was not preserved or was under-preserved. The method does not allow for additional preservative added after collection.
- PA The container was received outside of the acceptable pH for the analysis requested. Preservative was added upon receipt and the container is now at the correct pH. See the Sample Receipt Form for details on the amount and lot # of the preservative added.
- PH The container was received outside of the acceptable pH for the analysis requested. Preservative was added upon receipt, but was insufficient to bring the container to the correct pH for the analysis requested. See the Sample Receipt Form for details on the amount and lot # of the preservative added. QN Insufficient sample quantity provided.

### LABORATORY DATA REVIEW CHECKLIST

Completed by: Chris Pepe
Title: Environmental Scientist

Date: December 2021

Consultant Firm: Shannon & Wilson, Inc.

Laboratory Name: SGS North America Inc. Laboratory Report Number: 1214128 Laboratory Report Date: 7/29/2021

Contaminated Site Name: 5701 Northwood Drive

**ADEC File Number:** 2100.38.536 **Hazard Identification Number:** 25942

(**NOTE**: *NA* = not applicable; Text in *italics* added by Shannon & Wilson, Inc.)

### 1. <u>Laboratory</u>

a. Did an ADEC CS approved laboratory receive and <u>perform</u> all of the submitted sample analyses? Yes/ No / NA
 Comments:

**b.** If the samples were transferred to another "network" laboratory or sub-contracted to an alternate laboratory, was the laboratory performing the analyses ADEC CS approved? **Yes / No (NA)** 

Comments: The samples were not transferred to another "network" laboratory or subcontracted to an alternate laboratory.

# 2. Chain of Custody (COC)

a. COC information completed, signed, and dated (including released/received by)?
 Yes/ No / NA
 Comments:

b. Correct analyses requested? Yes / No / NA Comments:

# 3. Laboratory Sample Receipt Documentation

a. Sample/cooler temperature documented and within range at receipt (0° to 6° C)? Yes (No)/ NA

Comments: *The cooler temperature blank was received by the laboratory at* 8.7° *Celsius.* 

- **b.** Sample preservation acceptable acidified waters, Methanol preserved VOC soil (GRO, BTEX, VOCs, etc.)? **Yes**/ **No** / **NA** Comments:
- c. Sample condition documented broken, leaking (MeOH), zero headspace (VOC vials)?
   Yes/ No / NA
   Comments:
- **d.** If there were any discrepancies, were they documented? For example, incorrect sample containers/preservation, sample temperature outside of acceptable range, insufficient or missing samples, etc.? **Yes / No / NA**Comments: *No discrepancies were noted*.
- **e.** Data quality or usability affected? Comments: *It is our opinion that the slight temperature exceedance does not affect the useability of the data because the samples were delivered to the laboratory within 2.5 hours from the collection of the samples.*

# 4. Case Narrative

- a. Present and understandable? Yes/ No / NA Comments:
- **b.** Discrepancies, errors or QC failures noted by the lab? **Yes**/**No**/**NA** Comments: *The case narrative noted the following:* 
  - AK102 DRO is detect in the MB greater than one-half the LOQ, but less than the LOQ.
  - 8270D SIM PAH MS/MSD RPD for 2-methylnaphthalene does not meet QC criteria. This analyte was not detected above the LOQ in the parent sample.
- c. Were all corrective actions documented? **Yes** No / NA Comments:
- **d.** What is the effect on data quality/usability, according to the case narrative? Comments: *See above*.

# 5. Sample Results

a. Correct analyses performed/reported as requested on COC? Ves/ No / NA Comments:

**b.** All applicable holding times met? **Yes** / **No** / **NA** Comments:

c. All soils reported on a dry weight basis? Yes / No / NA Comments:

**d.** Are the reported LOQs less than the Cleanup Level or the minimum required detection level for the project? **Yes** (No) NA Comments: *The LOQ for 1,2,3-trichloropropane are greater than the ADEC cleanup level*.

**e.** Data quality or usability affected? Comments: *There is a potential that the target analyte is present at concentrations greater than the ADEC cleanup level, but less than the LOQ.* 

# 6. QC Samples

#### a. Method Blank

i. One method blank reported per matrix, analysis, and 20 samples?Yes/ No / NAComments:

**ii.** All method blank results less than limit of quantitation (LOQ) or project specified objectives?

(Yes)/ No / NA

Comments: The method blanks contained estimated concentrations of DRO (337 J  $\mu$ g/L), fluorene (0.0171 J  $\mu$ g/L), and phenanthrene (0.0243 J  $\mu$ g/L) less than the LOQs.

iii. If above LOQ or project specified objectives, what samples are affected? Comments: DRO: Samples B1MW and B11MW.

Phenanthrene and fluorene: Sample B1MW

iv. Do the affected sample(s) have data flags? If so, are the data flags clearly defined? Yes / No (NA)

Comments: Samples are flagged "B" in Table 2A when the reported sample concentration is within 10x the reported method blank concentration. The concentrations of DRO detected in Samples B1MW (859 µg/L) and B11MW (1,020 µg/L) are within 5x the reported method blank concentration. The sample concentrations are reported and flagged "B" on Table 2A. Fluorene and phenanthrene were not detected in Sample B1MW, therefore, the sample concentrations are reported as non-detect and flagging is not required.

v. Data quality or usability affected? Comments: *See above*.

# b. Laboratory Control Sample/Duplicate (LCS/LCSD)

- i. Organics One LCS/LCSD reported per matrix, analysis, and 20 samples?
   (LCS/LCSD required per AK methods, LCS required per SW846) (Yes) No / NA Comments:
- ii. Metals/Inorganics One LCS and one sample duplicate reported per matrix, analysis and 20 samples? Yes / No NA Comments:
- iii. Accuracy All percent recoveries (%R) reported and within method or laboratory limits and project specified objectives, if applicable. (AK petroleum methods: AK 101 60%-120%, AK 102 75%-125%, AK 103 60%-120%; all other analyses see the laboratory QC pages) Yes/No/NA Comments:
- iv. Precision All relative percent differences (RPDs) reported and less than method or laboratory limits and project specified objectives, if applicable. RPD reported from LCS/LCSD, and/or sample/sample duplicate. (AK Petroleum methods 20%; all other analyses see the laboratory QC pages) Yes/No/NA
  Comments:
- **v.** If %R or RPD is outside of acceptable limits, what samples are affected? Comments: *See above*.
- vi. Do the affected samples(s) have data flags? If so, are the data flags clearly defined?

  Yes / No / NA

  Comments:
- **vii.** Data quality or usability affected? Comments: *See above*.

# c. Matrix Spike/Matrix Spike Duplicate (MS/MSD) Note: Leave blank if not required for project

i. Organics - One MS/MSD reported per matrix, analysis, and 20 samples?
 Yes/ No / NA
 Comments:

- ii. Metals/Inorganics One MS and one MSD reported per matrix, analysis and 20 samples? Yes / No NA Comments:
- iii. Accuracy All percent recoveries (%R) reported and within method or laboratory limits and project specified objectives, if applicable. (AK petroleum methods: AK 101 60%-120%, AK 102 75%-125%, AK 103 60%-120%; all other analyses see the laboratory QC pages) Yes/No/NA Comments:
- iv. Precision All relative percent differences (RPDs) reported and less than method or laboratory limits and project specified objectives, if applicable. RPD reported from MS/MSD, and/or sample/sample duplicate. (AK Petroleum methods 20%; all other analyses see the laboratory QC pages) Yes (No) NA Comments: The RPD for 2-methylnaphthalene does not meet QC criteria.
- **v.** If %R or RPD is outside of acceptable limits, what samples are affected? Comments: *Sample B1MW*
- vi. Do the affected samples(s) have data flags? If so, are the data flags clearly defined? Yes / No NA

  Comments: 2-methylnaphthalene was not detected above the LOQ in Sample B1MW,

therefore, flagging is not required.

**vii.** Data quality or usability affected? Comments: *See above*.

# d. Surrogates - Organics Only or Isotope Dilution Analytes (IDA) - Isotope Dilution Methods Only

- i. Are surrogate/IDA recoveries reported for organic analyses field, QC, and laboratory samples? Yes/ No / NA
  Comments:
- ii. Accuracy All percent recoveries (%R) reported and within method or laboratory limits and project specified objectives, if applicable. (AK Petroleum methods 50-150 %R; all other analyses see the laboratory report pages) Yes / No / NA Comments:

- iii. Do the sample results with failed surrogate/IDA recoveries have data flags? If so, are the data flags clearly defined? Yes / No / NA Comments:
- **iv.** Data quality or usability affected? Comments: *No, see above.*
- e. Trip Blank Volatile analyses only (GRO, BTEX, VOCs, etc.)
  - i. One trip blank reported per matrix, analysis and for each cooler containing volatile samples? Yes/No/NA
     Comments:
  - ii. Is the cooler used to transport the trip blank and volatile samples clearly indicated on the COC? Yes (No) NA

Comments: *Only one cooler was used to transport the samples.* 

- iii. All results less than LOQ and project specified objectives? Yes/ No / NA Comments:
- iv. If above LOQ or project specified DQOs, what samples are affected? Comments:
- **v.** Data quality or usability affected? Comments: *No, see above.*

# f. Field Duplicate

- i. One field duplicate submitted per matrix, analysis and 10 project samples?

  Yes/ No / NA

  Comments: Duplicate Sample B11MW (duplicate of B1MW) was submitted to the laboratory.
- ii. Were the field duplicates submitted blind to the lab? Yes/ No / NA Comments:
- iii. Precision All relative percent differences (RPDs) less than specified project objectives? (Recommended: 30% for water, 50% for soil Yes / No / NA Comments:
- **iv.** Data quality or usability affected? Comments:

**g. Decontamination or Equipment Blank** (if not applicable, a comment stating why must be entered below).

Yes /No NA

Comments: A decontamination blank or equipment blank was not included in our ADEC-approved workplan.

All results less than LOQ and project specified objectives?
 Yes / No NA
 Comments:

- **ii.** If above LOQ or project specified objectives, what samples are affected? Comments:
- **iii.** Data quality or usability affected? Comments:

# 7. Other Data Flags/Qualifiers (ACOE, AFCEE, Lab Specific, etc.)

**a.** Defined and appropriate? **Yes** / **No** / **NA**Comments: A key is provided on Page 3 of the SGS Laboratory Report.



#### **Laboratory Report of Analysis**

To: Shannon & Wilson, Inc.

5430 Fairbanks Street, Suite 3 Anchorage, AK 99518 (907)433-3223

Report Number: 1216680

Client Project: 101528-002 5701 NorthwoodDrive

Dear Dan McMahon,

Enclosed are the results of the analytical services performed under the referenced project for the received samples and associated QC as applicable. The samples are certified to meet the requirements of the National Environmental Laboratory Accreditation Conference Standards. Copies of this report and supporting data will be retained in our files for a period of ten years in the event they are required for future reference. All results are intended to be used in their entirety and SGS is not responsible for use of less than the complete report. Any samples submitted to our laboratory will be retained for a maximum of fourteen (14) days from the date of this report unless other archiving requirements were included in the quote.

If there are any questions about the report or services performed during this project, please call Justin at (907) 562-2343. We will be happy to answer any questions or concerns which you may have.

Thank you for using SGS North America Inc. for your analytical services. We look forward to working with you again on any additional analytical needs.

Sincerely,

SGS North America Inc.

Justin Nelson 2021.10.20

15:17:01 -08'00'

Justin Nelson Project Manager Justin.Nelson@sgs.com Date

Print Date: 10/20/2021 9:08:14AM Results via Engage



#### **Case Narrative**

SGS Client: **Shannon & Wilson, Inc.**SGS Project: **1216680** 

Project Name/Site: 101528-002 5701 NorthwoodDrive

Project Contact: Dan McMahon

Refer to sample receipt form for information on sample condition.

#### LCSD for HBN 1826750 [XXX/4569 (1640957) LCSD

8270D SIM - PAH LCS/LCSD RPDs for indeno[1,2,3-c,d]pyrene and dibenzo[a,h]anthracene do not meet QC criteria. These analytes were not detected above the LOQ in associated samples.

\*QC comments may be associated with the field samples found in this report. When applicable, comments will be applied to associated field samples.

Print Date: 10/20/2021 9:08:16AM



#### **Report of Manual Integrations**

<u>Laboratory ID</u> <u>Client Sample ID</u> <u>Analytical Batch</u> <u>Analyte</u> <u>Reason</u>

8270D SIM LV (PAH)

1640955 MB for HBN 1826750 [XXX/45697] XMS12943 2-Methylnaphthalene SP

#### Manual Integration Reason Code Descriptions

Code Description

O Original Chromatogram
M Modified Chromatogram
SS Skimmed surrogate
BLG Closed baseline gap
RP Reassign peak name
PIR Pattern integration required

IT Included tail SP Split peak

RSP Removed split peak
FPS Forced peak start/stop
BLC Baseline correction

PNF Peak not found by software

All DRO/RRO analysis are integrated per SOP.

Print Date: 10/20/2021 9:08:17AM



#### **Laboratory Qualifiers**

Enclosed are the analytical results associated with the above work order. The results apply to the samples as received. All results are intended to be used in their entirety and SGS is not responsible for use of less than the complete report. This document is issued by the Company under its General Conditions of Service accessible at <a href="http://www.sgs.com/en/Terms-and-Conditions.aspx">http://www.sgs.com/en/Terms-and-Conditions.aspx</a>. Attention is drawn to the limitation of liability, indenmification and jurisdiction issues defined therein.

Any holder of this document is advised that information contained hereon reflects the Company's findings at the time of its intervention only and within the limits of Client's instructions, if any. The Company's sole responsibility is to its Client and this document does not exonerate parties to a transaction from exercising all their rights and obligations under the transaction documents. Any unauthorized alteration, forgery or falsification of the context or appearance of this document is unlawful and offenders may be prosecuted to the fullest extent of the law.

SGS maintains a formal Quality Assurance/Quality Control (QA/QC) program. A copy of our Quality Assurance Plan (QAP), which outlines this program, is available at your request. The laboratory certification numbers are AK00971 (DW Chemistry & Microbiology) & 17-021 (CS) for ADEC and 2944.01 for DOD ELAP/ISO17025 (RCRA methods: 1020B, 1311, 3010A, 3050B, 3520C, 3550C, 5030B, 5035A, 6020B, 7470A, 7471B, 8015C, 8021B, 8082A, 8260D, 8270D, 8270D-SIM, 9040C, 9045D, 9056A, 9060A, AK101 and AK102/103). SGS is only certified for the analytes listed on our Drinking Water Certification (DW methods: 200.8, 2130B, 2320B, 2510B, 300.0, 4500-CN-C,E, 4500-H-B, 4500-NO3-F, 4500-P-E and 524.2) and only those analytes will be reported to the State of Alaska for compliance. Except as specifically noted, all statements and data in this report are in conformance to the provisions set forth by the SGS QAP and, when applicable, other regulatory authorities.

The following descriptors or qualifiers may be found in your report:

\* The analyte has exceeded allowable regulatory or control limits.

! Surrogate out of control limits.

B Indicates the analyte is found in a blank associated with the sample.

CCV/CVA/CVB Continuing Calibration Verification
CCCV/CVC/CVCA/CVCB Closing Continuing Calibration Verification

CL Control Limit

DF Analytical Dilution Factor

DL Detection Limit (i.e., maximum method detection limit)
E The analyte result is above the calibrated range.

GT Greater Than
IB Instrument Blank

ICV Initial Calibration Verification
J The quantitation is an estimation.
LCS(D) Laboratory Control Spike (Duplicate)
LLQC/LLIQC Low Level Quantitation Check

LOD Limit of Detection (i.e., 1/2 of the LOQ)

LOQ Limit of Quantitation (i.e., reporting or practical quantitation limit)

LT Less Than MB Method Blank

MS(D) Matrix Spike (Duplicate)

ND Indicates the analyte is not detected.

RPD Relative Percent Difference
TNTC Too Numerous To Count

U Indicates the analyte was analyzed for but not detected.

Note: Sample summaries which include a result for "Total Solids" have already been adjusted for moisture content.

All DRO/RRO analyses are integrated per SOP.

Print Date: 10/20/2021 9:08:19AM

SGS North America Inc.

200 West Potter Drive, Anchorage, AK 99518 t 907.562.2343 f 907.561.5301 www.us.sgs.com



# **Sample Summary**

Client Sample ID	Lab Sample ID	Collected	Received	<u>Matrix</u>
101528-B1MW	1216680001	10/07/2021	10/07/2021	Water (Surface, Eff., Ground)
10528-B11MW	1216680002	10/07/2021	10/07/2021	Water (Surface, Eff., Ground)
101528-WTB	1216680003	10/07/2021	10/07/2021	Water (Surface, Eff., Ground)

MethodMethod Description8270D SIM LV (PAH)8270 PAH SIM GC/MS LVAK102DRO/RRO Low Volume WaterAK103DRO/RRO Low Volume Water

SW8260D Volatile Organic Compounds (W) FULL

Print Date: 10/20/2021 9:08:20AM



# **Detectable Results Summary**

Client Sample ID: 101528-B1MW			
Lab Sample ID: 1216680001	<u>Parameter</u>	Result	<u>Units</u>
Polynuclear Aromatics GC/MS	2-Methylnaphthalene	0.0163J	ug/L
	Phenanthrene	0.0228J	ug/L
Semivolatile Organic Fuels	Diesel Range Organics	0.701	mg/L
<del>-</del>	Residual Range Organics	0.752	mg/L
Volatile GC/MS	Benzene	0.921	ug/L
	cis-1,2-Dichloroethene	0.317J	ug/L
	Vinyl chloride	0.418	ug/L
Client Sample ID: 10528-B11MW			
Lab Sample ID: 1216680002	<u>Parameter</u>	Result	<u>Units</u>
Semivolatile Organic Fuels	Diesel Range Organics	0.609	mg/L
	Residual Range Organics	0.760	mg/L

Print Date: 10/20/2021 9:08:22AM



#### Results of 101528-B1MW

Client Sample ID: 101528-B1MW

Client Project ID: 101528-002 5701 NorthwoodDrive

Lab Sample ID: 1216680001 Lab Project ID: 1216680 Collection Date: 10/07/21 12:30 Received Date: 10/07/21 13:52 Matrix: Water (Surface, Eff., Ground)

Solids (%): Location:

# Results by Polynuclear Aromatics GC/MS

						<u>Allowable</u>	
<u>Parameter</u>	Result Qual	LOQ/CL	<u>DL</u>	<u>Units</u>	<u>DF</u>	<u>Limits</u>	Date Analyzed
1-Methylnaphthalene	0.0240 U	0.0481	0.0144	ug/L	1		10/11/21 22:08
2-Methylnaphthalene	0.0163 J	0.0481	0.0144	ug/L	1		10/11/21 22:08
Acenaphthene	0.0240 U	0.0481	0.0144	ug/L	1		10/11/21 22:08
Acenaphthylene	0.0240 U	0.0481	0.0144	ug/L	1		10/11/21 22:08
Anthracene	0.0240 U	0.0481	0.0144	ug/L	1		10/11/21 22:08
Benzo(a)Anthracene	0.0240 U	0.0481	0.0144	ug/L	1		10/11/21 22:08
Benzo[a]pyrene	0.00960 U	0.0192	0.00596	ug/L	1		10/11/21 22:08
Benzo[b]Fluoranthene	0.0240 U	0.0481	0.0144	ug/L	1		10/11/21 22:08
Benzo[g,h,i]perylene	0.0240 U	0.0481	0.0144	ug/L	1		10/11/21 22:08
Benzo[k]fluoranthene	0.0240 U	0.0481	0.0144	ug/L	1		10/11/21 22:08
Chrysene	0.0240 U	0.0481	0.0144	ug/L	1		10/11/21 22:08
Dibenzo[a,h]anthracene	0.00960 U	0.0192	0.00596	ug/L	1		10/11/21 22:08
Fluoranthene	0.0240 U	0.0481	0.0144	ug/L	1		10/11/21 22:08
Fluorene	0.0240 U	0.0481	0.0144	ug/L	1		10/11/21 22:08
Indeno[1,2,3-c,d] pyrene	0.0240 U	0.0481	0.0144	ug/L	1		10/11/21 22:08
Naphthalene	0.0481 U	0.0962	0.0298	ug/L	1		10/11/21 22:08
Phenanthrene	0.0228 J	0.0481	0.0144	ug/L	1		10/11/21 22:08
Pyrene	0.0240 U	0.0481	0.0144	ug/L	1		10/11/21 22:08
Surrogates							
2-Methylnaphthalene-d10 (surr)	73.2	42-86		%	1		10/11/21 22:08
Fluoranthene-d10 (surr)	76.8	50-97		%	1		10/11/21 22:08

#### **Batch Information**

Analytical Batch: XMS12943

Analytical Method: 8270D SIM LV (PAH)

Analyst: LAW

Analytical Date/Time: 10/11/21 22:08 Container ID: 1216680001-C Prep Batch: XXX45697 Prep Method: SW3535A Prep Date/Time: 10/09/21 09:14

Prep Initial Wt./Vol.: 260 mL Prep Extract Vol: 1 mL

Print Date: 10/20/2021 9:08:24AM

J flagging is activated



Client Sample ID: 101528-B1MW

Client Project ID: 101528-002 5701 NorthwoodDrive

Lab Sample ID: 1216680001 Lab Project ID: 1216680 Collection Date: 10/07/21 12:30 Received Date: 10/07/21 13:52 Matrix: Water (Surface, Eff., Ground)

Solids (%): Location:

#### Results by Semivolatile Organic Fuels

						<u>Allowable</u>	
<u>Parameter</u>	Result Qual	LOQ/CL	<u>DL</u>	<u>Units</u>	<u>DF</u>	<u>Limits</u>	Date Analyzed
Diesel Range Organics	0.701	0.588	0.196	mg/L	1		10/13/21 22:16
Surrogates							
5a Androstane (surr)	86.5	50-150		%	1		10/13/21 22:16

#### **Batch Information**

Analytical Batch: XFC16112 Analytical Method: AK102

Analyst: IVM

Analytical Date/Time: 10/13/21 22:16 Container ID: 1216680001-A Prep Batch: XXX45702 Prep Method: SW3520C Prep Date/Time: 10/09/21 15:45 Prep Initial Wt./Vol.: 255 mL Prep Extract Vol: 1 mL

						<u>Allowable</u>	
<u>Parameter</u>	Result Qual	LOQ/CL	<u>DL</u>	<u>Units</u>	<u>DF</u>	<u>Limits</u>	Date Analyzed
Residual Range Organics	0.752	0.490	0.196	mg/L	1		10/13/21 22:16
Surrogates							
n-Triacontane-d62 (surr)	85.3	50-150		%	1		10/13/21 22:16

#### **Batch Information**

Analytical Batch: XFC16112 Analytical Method: AK103

Analyst: IVM

Analytical Date/Time: 10/13/21 22:16 Container ID: 1216680001-A Prep Batch: XXX45702 Prep Method: SW3520C Prep Date/Time: 10/09/21 15:45 Prep Initial Wt./Vol.: 255 mL Prep Extract Vol: 1 mL

Print Date: 10/20/2021 9:08:24AM J flagging is activated



Client Sample ID: 101528-B1MW

Client Project ID: 101528-002 5701 NorthwoodDrive

Lab Sample ID: 1216680001 Lab Project ID: 1216680 Collection Date: 10/07/21 12:30 Received Date: 10/07/21 13:52 Matrix: Water (Surface, Eff., Ground)

Solids (%): Location:

#### Results by Volatile GC/MS

<u>Parameter</u>	Result Qual	LOQ/CL	<u>DL</u>	<u>Units</u>	<u>DF</u>	Allowable <u>Limits</u>	Date Analyzed
1,1,1,2-Tetrachloroethane	0.250 U	0.500	0.150	ug/L	1		10/16/21 16:11
1,1,1-Trichloroethane	0.500 U	1.00	0.310	ug/L	1		10/16/21 16:11
1,1,2,2-Tetrachloroethane	0.250 U	0.500	0.150	ug/L	1		10/16/21 16:11
1,1,2-Trichloroethane	0.200 U	0.400	0.120	ug/L	1		10/16/21 16:11
1,1-Dichloroethane	0.500 U	1.00	0.310	ug/L	1		10/16/21 16:11
1,1-Dichloroethene	0.500 U	1.00	0.310	ug/L	1		10/16/21 16:11
1,1-Dichloropropene	0.500 U	1.00	0.310	ug/L	1		10/16/21 16:11
1,2,3-Trichlorobenzene	0.500 U	1.00	0.310	ug/L	1		10/16/21 16:11
1,2,3-Trichloropropane	0.500 U	1.00	0.310	ug/L	1		10/16/21 16:11
1,2,4-Trichlorobenzene	0.500 U	1.00	0.310	ug/L	1		10/16/21 16:11
1,2,4-Trimethylbenzene	0.500 U	1.00	0.310	ug/L	1		10/16/21 16:11
1,2-Dibromo-3-chloropropane	5.00 U	10.0	3.10	ug/L	1		10/16/21 16:11
1,2-Dibromoethane	0.0375 U	0.0750	0.0180	ug/L	1		10/16/21 16:11
1,2-Dichlorobenzene	0.500 U	1.00	0.310	ug/L	1		10/16/21 16:11
1,2-Dichloroethane	0.250 U	0.500	0.200	ug/L	1		10/16/21 16:11
1,2-Dichloropropane	0.500 U	1.00	0.310	ug/L	1		10/16/21 16:11
1,3,5-Trimethylbenzene	0.500 U	1.00	0.310	ug/L	1		10/16/21 16:11
1,3-Dichlorobenzene	0.500 U	1.00	0.310	ug/L	1		10/16/21 16:11
1,3-Dichloropropane	0.250 U	0.500	0.150	ug/L	1		10/16/21 16:11
1,4-Dichlorobenzene	0.250 U	0.500	0.150	ug/L	1		10/16/21 16:11
2,2-Dichloropropane	0.500 U	1.00	0.310	ug/L	1		10/16/21 16:11
2-Butanone (MEK)	5.00 U	10.0	3.10	ug/L	1		10/16/21 16:11
2-Chlorotoluene	0.500 U	1.00	0.310	ug/L	1		10/16/21 16:11
2-Hexanone	5.00 U	10.0	3.10	ug/L	1		10/16/21 16:11
4-Chlorotoluene	0.500 U	1.00	0.310	ug/L	1		10/16/21 16:11
4-Isopropyltoluene	0.500 U	1.00	0.310	ug/L	1		10/16/21 16:11
4-Methyl-2-pentanone (MIBK)	5.00 U	10.0	3.10	ug/L	1		10/16/21 16:11
Benzene	0.921	0.400	0.120	ug/L	1		10/16/21 16:11
Bromobenzene	0.500 U	1.00	0.310	ug/L	1		10/16/21 16:11
Bromochloromethane	0.500 U	1.00	0.310	ug/L	1		10/16/21 16:11
Bromodichloromethane	0.250 U	0.500	0.150	ug/L	1		10/16/21 16:11
Bromoform	0.500 U	1.00	0.310	ug/L	1		10/16/21 16:11
Bromomethane	3.00 U	6.00	3.00	ug/L	1		10/16/21 16:11
Carbon disulfide	5.00 U	10.0	3.10	ug/L	1		10/16/21 16:11
Carbon tetrachloride	0.500 U	1.00	0.310	ug/L	1		10/16/21 16:11
Chlorobenzene	0.250 U	0.500	0.150	ug/L	1		10/16/21 16:11
Chloroethane	0.500 U	1.00	0.310	ug/L	1		10/16/21 16:11

Print Date: 10/20/2021 9:08:24AM

J flagging is activated



Client Sample ID: 101528-B1MW

Client Project ID: 101528-002 5701 NorthwoodDrive

Lab Sample ID: 1216680001 Lab Project ID: 1216680 Collection Date: 10/07/21 12:30 Received Date: 10/07/21 13:52 Matrix: Water (Surface, Eff., Ground)

Solids (%): Location:

#### Results by Volatile GC/MS

Parameter	Result Qual	LOQ/CL	DL	Units	<u>DF</u>	Allowable Limits Date Analy
<u>-arameter</u> Chloroform	0.500 U	1.00	<u>DL</u> 0.310	ug/L	<u>DF</u> 1	10/16/21 1
Chloromethane	0.500 U	1.00	0.310	ug/L	1	10/16/21 1
cis-1,2-Dichloroethene	0.300 J 0.317 J	1.00	0.310	ug/L ug/L	1	10/16/21 1
cis-1,3-Dichloropropene	0.317 J 0.250 U	0.500	0.310	ug/L ug/L	1	10/16/21 1
Dibromochloromethane	0.250 U	0.500	0.150	ug/L ug/L	1	10/16/21 1
Dibromomethane	0.500 U	1.00	0.130	ŭ	1	10/16/21 1
Dibromomemane Dichlorodifluoromethane	0.500 U	1.00	0.310	ug/L	1	10/16/21 1
				ug/L		
Ethylbenzene	0.500 U	1.00	0.310	ug/L	1	10/16/21 1
Freon-113	5.00 U	10.0	3.10	ug/L	1	10/16/21 1
Hexachlorobutadiene	0.500 U	1.00	0.310	ug/L	1	10/16/21 1
sopropylbenzene (Cumene)	0.500 U	1.00	0.310	ug/L	1	10/16/21 1
Methylene chloride	5.00 U	10.0	3.10	ug/L	1	10/16/21 1
Methyl-t-butyl ether	5.00 U	10.0	3.10	ug/L	1	10/16/21 1
Naphthalene	0.500 U	1.00	0.310	ug/L	1	10/16/21 1
n-Butylbenzene	0.500 U	1.00	0.310	ug/L	1	10/16/21 1
n-Propylbenzene	0.500 U	1.00	0.310	ug/L	1	10/16/21 1
o-Xylene	0.500 U	1.00	0.310	ug/L	1	10/16/21 1
P & M -Xylene	1.00 U	2.00	0.620	ug/L	1	10/16/21 1
sec-Butylbenzene	0.500 U	1.00	0.310	ug/L	1	10/16/21 1
Styrene	0.500 U	1.00	0.310	ug/L	1	10/16/21 1
ert-Butylbenzene	0.500 U	1.00	0.310	ug/L	1	10/16/21 1
Tetrachloroethene	0.500 U	1.00	0.310	ug/L	1	10/16/21 1
Toluene	0.500 U	1.00	0.310	ug/L	1	10/16/21 1
rans-1,2-Dichloroethene	0.500 U	1.00	0.310	ug/L	1	10/16/21 1
rans-1,3-Dichloropropene	0.500 U	1.00	0.310	ug/L	1	10/16/21 1
Trichloroethene	0.500 U	1.00	0.310	ug/L	1	10/16/21 1
Frichlorofluoromethane	0.500 U	1.00	0.310	ug/L	1	10/16/21 1
√inyl acetate	5.00 U	10.0	3.10	ug/L	1	10/16/21 1
/inyl chloride	0.418	0.150	0.0500	ug/L	1	10/16/21 1
Kylenes (total)	1.50 U	3.00	1.00	ug/L	1	10/16/21 1
urrogates						
1,2-Dichloroethane-D4 (surr)	101	81-118		%	1	10/16/21 1
4-Bromofluorobenzene (surr)	99.5	85-114		%	1	10/16/21 1
Toluene-d8 (surr)	99.3	89-112		%	1	10/16/21 1

Print Date: 10/20/2021 9:08:24AM

J flagging is activated



Client Sample ID: 101528-B1MW

Client Project ID: 101528-002 5701 NorthwoodDrive

Lab Sample ID: 1216680001 Lab Project ID: 1216680 Collection Date: 10/07/21 12:30 Received Date: 10/07/21 13:52 Matrix: Water (Surface, Eff., Ground)

Solids (%): Location:

#### Results by Volatile GC/MS

#### **Batch Information**

Analytical Batch: VMS21286 Analytical Method: SW8260D

Analyst: JMG

Analytical Date/Time: 10/16/21 16:11 Container ID: 1216680001-E Prep Batch: VXX38039 Prep Method: SW5030B Prep Date/Time: 10/16/21 06:00 Prep Initial Wt./Vol.: 5 mL Prep Extract Vol: 5 mL

Print Date: 10/20/2021 9:08:24AM J flagging is activated



Client Sample ID: 10528-B11MW

Client Project ID: 101528-002 5701 NorthwoodDrive

Lab Sample ID: 1216680002 Lab Project ID: 1216680 Collection Date: 10/07/21 13:00 Received Date: 10/07/21 13:52 Matrix: Water (Surface, Eff., Ground)

Solids (%): Location:

#### Results by Semivolatile Organic Fuels

						<u>Allowable</u>	
<u>Parameter</u>	Result Qual	LOQ/CL	<u>DL</u>	<u>Units</u>	<u>DF</u>	<u>Limits</u>	Date Analyzed
Diesel Range Organics	0.609	0.577	0.192	mg/L	1		10/13/21 22:26
Surrogates							
5a Androstane (surr)	90.8	50-150		%	1		10/13/21 22:26

#### **Batch Information**

Analytical Batch: XFC16112 Analytical Method: AK102

Analyst: IVM

Analytical Date/Time: 10/13/21 22:26 Container ID: 1216680002-A Prep Batch: XXX45702 Prep Method: SW3520C Prep Date/Time: 10/09/21 15:45 Prep Initial Wt./Vol.: 260 mL Prep Extract Vol: 1 mL

						<u>Allowable</u>	
<u>Parameter</u>	Result Qual	LOQ/CL	<u>DL</u>	<u>Units</u>	<u>DF</u>	<u>Limits</u>	Date Analyzed
Residual Range Organics	0.760	0.481	0.192	mg/L	1		10/13/21 22:26
Surrogates							
n-Triacontane-d62 (surr)	86.6	50-150		%	1		10/13/21 22:26

#### **Batch Information**

Analytical Batch: XFC16112 Analytical Method: AK103

Analyst: IVM

Analytical Date/Time: 10/13/21 22:26 Container ID: 1216680002-A Prep Batch: XXX45702 Prep Method: SW3520C Prep Date/Time: 10/09/21 15:45 Prep Initial Wt./Vol.: 260 mL Prep Extract Vol: 1 mL

Print Date: 10/20/2021 9:08:24AM

200 West Potter Drive Anchorage, AK 95518 t 907.562.2343 f 907.561.5301 www.us.sgs.com



#### Results of 101528-WTB

Client Sample ID: 101528-WTB

Client Project ID: 101528-002 5701 NorthwoodDrive

Lab Sample ID: 1216680003 Lab Project ID: 1216680 Collection Date: 10/07/21 09:00 Received Date: 10/07/21 13:52 Matrix: Water (Surface, Eff., Ground)

Solids (%): Location:

#### Results by Volatile GC/MS

<u>Parameter</u>	Result Qual	LOQ/CL	<u>DL</u>	<u>Units</u>	<u>DF</u>	Allowable Limits Date Analyzed
1,1,1,2-Tetrachloroethane	0.250 U	0.500	0.150	ug/L	1	10/16/21 15:27
1,1,1-Trichloroethane	0.500 U	1.00	0.310	ug/L	1	10/16/21 15:27
1,1,2,2-Tetrachloroethane	0.250 U	0.500	0.150	ug/L	1	10/16/21 15:27
1,1,2-Trichloroethane	0.200 U	0.400	0.120	ug/L	1	10/16/21 15:27
1,1-Dichloroethane	0.500 U	1.00	0.310	ug/L	1	10/16/21 15:27
1,1-Dichloroethene	0.500 U	1.00	0.310	ug/L	1	10/16/21 15:27
1,1-Dichloropropene	0.500 U	1.00	0.310	ug/L	1	10/16/21 15:27
1,2,3-Trichlorobenzene	0.500 U	1.00	0.310	ug/L	1	10/16/21 15:27
1,2,3-Trichloropropane	0.500 U	1.00	0.310	ug/L	1	10/16/21 15:27
1,2,4-Trichlorobenzene	0.500 U	1.00	0.310	ug/L	1	10/16/21 15:27
1,2,4-Trimethylbenzene	0.500 U	1.00	0.310	ug/L	1	10/16/21 15:27
1,2-Dibromo-3-chloropropane	5.00 U	10.0	3.10	ug/L	1	10/16/21 15:27
1,2-Dibromoethane	0.0375 U	0.0750	0.0180	ug/L	1	10/16/21 15:27
1,2-Dichlorobenzene	0.500 U	1.00	0.310	ug/L	1	10/16/21 15:27
1,2-Dichloroethane	0.250 U	0.500	0.200	ug/L	1	10/16/21 15:27
1,2-Dichloropropane	0.500 U	1.00	0.310	ug/L	1	10/16/21 15:27
1,3,5-Trimethylbenzene	0.500 U	1.00	0.310	ug/L	1	10/16/21 15:27
1,3-Dichlorobenzene	0.500 U	1.00	0.310	ug/L	1	10/16/21 15:27
1,3-Dichloropropane	0.250 U	0.500	0.150	ug/L	1	10/16/21 15:27
1,4-Dichlorobenzene	0.250 U	0.500	0.150	ug/L	1	10/16/21 15:27
2,2-Dichloropropane	0.500 U	1.00	0.310	ug/L	1	10/16/21 15:27
2-Butanone (MEK)	5.00 U	10.0	3.10	ug/L	1	10/16/21 15:27
2-Chlorotoluene	0.500 U	1.00	0.310	ug/L	1	10/16/21 15:27
2-Hexanone	5.00 U	10.0	3.10	ug/L	1	10/16/21 15:27
4-Chlorotoluene	0.500 U	1.00	0.310	ug/L	1	10/16/21 15:27
4-Isopropyltoluene	0.500 U	1.00	0.310	ug/L	1	10/16/21 15:27
4-Methyl-2-pentanone (MIBK)	5.00 U	10.0	3.10	ug/L	1	10/16/21 15:27
Benzene	0.200 U	0.400	0.120	ug/L	1	10/16/21 15:27
Bromobenzene	0.500 U	1.00	0.310	ug/L	1	10/16/21 15:27
Bromochloromethane	0.500 U	1.00	0.310	ug/L	1	10/16/21 15:27
Bromodichloromethane	0.250 U	0.500	0.150	ug/L	1	10/16/21 15:27
Bromoform	0.500 U	1.00	0.310	ug/L	1	10/16/21 15:27
Bromomethane	3.00 U	6.00	3.00	ug/L	1	10/16/21 15:27
Carbon disulfide	5.00 U	10.0	3.10	ug/L	1	10/16/21 15:27
Carbon tetrachloride	0.500 U	1.00	0.310	ug/L	1	10/16/21 15:27
Chlorobenzene	0.250 U	0.500	0.150	ug/L	1	10/16/21 15:27
Chloroethane	0.500 U	1.00	0.310	ug/L	1	10/16/21 15:27

Print Date: 10/20/2021 9:08:24AM

J flagging is activated



#### Results of 101528-WTB

Client Sample ID: 101528-WTB

Client Project ID: 101528-002 5701 NorthwoodDrive

Lab Sample ID: 1216680003 Lab Project ID: 1216680 Collection Date: 10/07/21 09:00 Received Date: 10/07/21 13:52 Matrix: Water (Surface, Eff., Ground)

Solids (%): Location:

#### Results by Volatile GC/MS

<u>Parameter</u>	Result Qual	LOQ/CL	<u>DL</u>	Units	<u>DF</u>	Allowable Limits	Date Analyzed
Chloroform	0.500 U	1.00	0.310	ug/L	1		10/16/21 15:27
Chloromethane	0.500 U	1.00	0.310	ug/L	1		10/16/21 15:27
cis-1,2-Dichloroethene	0.500 U	1.00	0.310	ug/L	1		10/16/21 15:27
cis-1,3-Dichloropropene	0.250 U	0.500	0.150	ug/L	1		10/16/21 15:27
Dibromochloromethane	0.250 U	0.500	0.150	ug/L	1		10/16/21 15:27
Dibromomethane	0.500 U	1.00	0.310	ug/L	1		10/16/21 15:27
Dichlorodifluoromethane	0.500 U	1.00	0.310	ug/L	1		10/16/21 15:27
Ethylbenzene	0.500 U	1.00	0.310	ug/L	1		10/16/21 15:27
Freon-113	5.00 U	10.0	3.10	ug/L	1		10/16/21 15:27
Hexachlorobutadiene	0.500 U	1.00	0.310	ug/L	1		10/16/21 15:27
Isopropylbenzene (Cumene)	0.500 U	1.00	0.310	ug/L	1		10/16/21 15:27
Methylene chloride	5.00 U	10.0	3.10	ug/L	1		10/16/21 15:27
Methyl-t-butyl ether	5.00 U	10.0	3.10	ug/L	1		10/16/21 15:27
Naphthalene	0.500 U	1.00	0.310	ug/L	1		10/16/21 15:27
n-Butylbenzene	0.500 U	1.00	0.310	ug/L	1		10/16/21 15:27
n-Propylbenzene	0.500 U	1.00	0.310	ug/L	1		10/16/21 15:27
o-Xylene	0.500 U	1.00	0.310	ug/L	1		10/16/21 15:27
P & M -Xylene	1.00 U	2.00	0.620	ug/L	1		10/16/21 15:27
sec-Butylbenzene	0.500 U	1.00	0.310	ug/L	1		10/16/21 15:27
Styrene	0.500 U	1.00	0.310	ug/L	1		10/16/21 15:27
tert-Butylbenzene	0.500 U	1.00	0.310	ug/L	1		10/16/21 15:27
Tetrachloroethene	0.500 U	1.00	0.310	ug/L	1		10/16/21 15:27
Toluene	0.500 U	1.00	0.310	ug/L	1		10/16/21 15:27
trans-1,2-Dichloroethene	0.500 U	1.00	0.310	ug/L	1		10/16/21 15:27
trans-1,3-Dichloropropene	0.500 U	1.00	0.310	ug/L	1		10/16/21 15:27
Trichloroethene	0.500 U	1.00	0.310	ug/L	1		10/16/21 15:27
Trichlorofluoromethane	0.500 U	1.00	0.310	ug/L	1		10/16/21 15:27
Vinyl acetate	5.00 U	10.0	3.10	ug/L	1		10/16/21 15:27
Vinyl chloride	0.0750 U	0.150	0.0500	ug/L	1		10/16/21 15:27
Xylenes (total)	1.50 U	3.00	1.00	ug/L	1		10/16/21 15:27
Surrogates							
1,2-Dichloroethane-D4 (surr)	100	81-118		%	1		10/16/21 15:27
4-Bromofluorobenzene (surr)	103	85-114		%	1		10/16/21 15:27
Toluene-d8 (surr)	100	89-112		%	1		10/16/21 15:27

Print Date: 10/20/2021 9:08:24AM

J flagging is activated



#### Results of 101528-WTB

Client Sample ID: 101528-WTB

Client Project ID: 101528-002 5701 NorthwoodDrive

Lab Sample ID: 1216680003 Lab Project ID: 1216680 Collection Date: 10/07/21 09:00 Received Date: 10/07/21 13:52 Matrix: Water (Surface, Eff., Ground)

Solids (%): Location:

#### Results by Volatile GC/MS

#### **Batch Information**

Analytical Batch: VMS21286 Analytical Method: SW8260D

Analyst: JMG

Analytical Date/Time: 10/16/21 15:27 Container ID: 1216680003-A Prep Batch: VXX38039 Prep Method: SW5030B Prep Date/Time: 10/16/21 06:00 Prep Initial Wt./Vol.: 5 mL Prep Extract Vol: 5 mL

Print Date: 10/20/2021 9:08:24AM J flagging is activated



Blank ID: MB for HBN 1827292 [VXX/38039]

Blank Lab ID: 1642870

QC for Samples:

1216680001, 1216680003

Matrix: Water (Surface, Eff., Ground)

#### Results by SW8260D

Parameter	Results	LOQ/CL	<u>DL</u>	Units
1,1,1,2-Tetrachloroethane	0.250U	0.500	<u>DL</u> 0.150	ug/L
1,1,1-Trichloroethane	0.500U	1.00	0.310	ug/L
1,1,2,2-Tetrachloroethane	0.250U	0.500	0.150	ug/L
1,1,2-Trichloroethane	0.200U	0.400	0.120	ug/L
1,1-Dichloroethane	0.500U	1.00	0.310	ug/L
1,1-Dichloroethene	0.500U	1.00	0.310	ug/L
1,1-Dichloropropene	0.500U	1.00	0.310	ug/L
1,2,3-Trichlorobenzene	0.500U	1.00	0.310	ug/L
1,2,3-Trichloropropane	0.500U	1.00	0.310	ug/L
1,2,4-Trichlorobenzene	0.500U	1.00	0.310	ug/L
1,2,4-Trimethylbenzene	0.500U	1.00	0.310	ug/L
1,2-Dibromo-3-chloropropane	5.00U	10.0	3.10	ug/L
1,2-Dibromoethane	0.0375U	0.0750	0.0180	ug/L
1,2-Dichlorobenzene	0.500U	1.00	0.310	ug/L
1,2-Dichloroethane	0.250U	0.500	0.200	ug/L
1,2-Dichloropropane	0.500U	1.00	0.310	ug/L
1,3,5-Trimethylbenzene	0.500U	1.00	0.310	ug/L
1,3-Dichlorobenzene	0.500U	1.00	0.310	ug/L
1,3-Dichloropropane	0.250U	0.500	0.150	ug/L
1,4-Dichlorobenzene	0.250U	0.500	0.150	ug/L
2,2-Dichloropropane	0.500U	1.00	0.310	ug/L
2-Butanone (MEK)	5.00U	10.0	3.10	ug/L
2-Chlorotoluene	0.500U	1.00	0.310	ug/L
2-Hexanone	5.00U	10.0	3.10	ug/L
4-Chlorotoluene	0.500U	1.00	0.310	ug/L
4-Isopropyltoluene	0.500U	1.00	0.310	ug/L
4-Methyl-2-pentanone (MIBK)	5.00U	10.0	3.10	ug/L
Benzene	0.200U	0.400	0.120	ug/L
Bromobenzene	0.500U	1.00	0.310	ug/L
Bromochloromethane	0.500U	1.00	0.310	ug/L
Bromodichloromethane	0.250U	0.500	0.150	ug/L
Bromoform	0.500U	1.00	0.310	ug/L
Bromomethane	3.00U	6.00	3.00	ug/L
Carbon disulfide	5.00U	10.0	3.10	ug/L
Carbon tetrachloride	0.500U	1.00	0.310	ug/L
Chlorobenzene	0.250U	0.500	0.150	ug/L
Chloroethane	0.500U	1.00	0.310	ug/L
Chloroform	0.500U	1.00	0.310	ug/L

Print Date: 10/20/2021 9:08:25AM



Blank ID: MB for HBN 1827292 [VXX/38039]

Blank Lab ID: 1642870

QC for Samples:

1216680001, 1216680003

Matrix: Water (Surface, Eff., Ground)

#### Results by SW8260D

Parameter	Results	LOQ/CL	<u>DL</u>	Units
Chloromethane	0.500U	1.00	0.310	ug/L
cis-1,2-Dichloroethene	0.500U	1.00	0.310	ug/L
cis-1,3-Dichloropropene	0.250U	0.500	0.150	ug/L
Dibromochloromethane	0.250U	0.500	0.150	ug/L
Dibromomethane	0.500U	1.00	0.310	ug/L
Dichlorodifluoromethane	0.500U	1.00	0.310	ug/L
Ethylbenzene	0.500U	1.00	0.310	ug/L
Freon-113	5.00U	10.0	3.10	ug/L
Hexachlorobutadiene	0.500U	1.00	0.310	ug/L
Isopropylbenzene (Cumene)	0.500U	1.00	0.310	ug/L
Methylene chloride	5.00U	10.0	3.10	ug/L
Methyl-t-butyl ether	5.00U	10.0	3.10	ug/L
Naphthalene	0.500U	1.00	0.310	ug/L
n-Butylbenzene	0.500U	1.00	0.310	ug/L
n-Propylbenzene	0.500U	1.00	0.310	ug/L
o-Xylene	0.500U	1.00	0.310	ug/L
P & M -Xylene	1.00U	2.00	0.620	ug/L
sec-Butylbenzene	0.500U	1.00	0.310	ug/L
Styrene	0.500U	1.00	0.310	ug/L
tert-Butylbenzene	0.500U	1.00	0.310	ug/L
Tetrachloroethene	0.500U	1.00	0.310	ug/L
Toluene	0.500U	1.00	0.310	ug/L
trans-1,2-Dichloroethene	0.500U	1.00	0.310	ug/L
trans-1,3-Dichloropropene	0.500U	1.00	0.310	ug/L
Trichloroethene	0.500U	1.00	0.310	ug/L
Trichlorofluoromethane	0.500U	1.00	0.310	ug/L
Vinyl acetate	5.00U	10.0	3.10	ug/L
Vinyl chloride	0.0750U	0.150	0.0500	ug/L
Xylenes (total)	1.50U	3.00	1.00	ug/L
Surrogates				
1,2-Dichloroethane-D4 (surr)	100	81-118		%
4-Bromofluorobenzene (surr)	102	85-114		%
Toluene-d8 (surr)	99.7	89-112		%

Print Date: 10/20/2021 9:08:25AM



Blank ID: MB for HBN 1827292 [VXX/38039]

Blank Lab ID: 1642870

QC for Samples:

1216680001, 1216680003

Matrix: Water (Surface, Eff., Ground)

#### Results by SW8260D

Parameter Results LOQ/CL DL Units

#### **Batch Information**

Analytical Batch: VMS21286 Analytical Method: SW8260D Instrument: Agilent 7890-75MS

Analyst: JMG

Analytical Date/Time: 10/16/2021 1:36:00PM

Prep Batch: VXX38039 Prep Method: SW5030B

Prep Date/Time: 10/16/2021 6:00:00AM

Prep Initial Wt./Vol.: 5 mL Prep Extract Vol: 5 mL

Print Date: 10/20/2021 9:08:25AM



Blank Spike ID: LCS for HBN 1216680 [VXX38039]

Blank Spike Lab ID: 1642871 Date Analyzed: 10/16/2021 13:51

QC for Samples: 1216680001, 1216680003

Spike Duplicate ID: LCSD for HBN 1216680

[VXX38039]

Spike Duplicate Lab ID: 1642872 Matrix: Water (Surface, Eff., Ground)

#### Results by SW8260D

		Blank Spike	e (ug/L)	Spike Dupli	cate (ug/L)				
<u>Parameter</u>	Spike	Result	Rec (%)	Spike	Result	Rec (%)	<u>CL</u>	RPD (%)	RPD CL
1,1,1,2-Tetrachloroethane	30	30.5	102	30	29.8	100	(78-124)	2.10	(< 20 )
1,1,1-Trichloroethane	30	30.8	103	30	30.1	100	(74-131)	2.10	(< 20 )
1,1,2,2-Tetrachloroethane	30	30.6	102	30	30.1	100	(71-121)	1.50	(< 20 )
1,1,2-Trichloroethane	30	30.8	103	30	30.3	101	(80-119)	1.70	(< 20 )
1,1-Dichloroethane	30	30.1	100	30	29.6	99	(77-125)	1.60	(< 20 )
1,1-Dichloroethene	30	30.1	100	30	29.6	99	(71-131)	1.50	(< 20 )
1,1-Dichloropropene	30	31.1	104	30	30.4	101	(79-125)	2.10	(< 20 )
1,2,3-Trichlorobenzene	30	31.0	103	30	32.3	108	(69-129)	4.00	(< 20 )
1,2,3-Trichloropropane	30	29.8	99	30	29.7	99	(73-122)	0.54	(< 20 )
1,2,4-Trichlorobenzene	30	30.9	103	30	31.5	105	(69-130)	1.90	(< 20 )
1,2,4-Trimethylbenzene	30	30.5	102	30	30.1	100	(79-124)	1.40	(< 20 )
1,2-Dibromo-3-chloropropane	30	28.9	97	30	29.8	99	(62-128)	3.00	(< 20 )
1,2-Dibromoethane	30	30.6	102	30	30.3	101	(77-121)	1.20	(< 20 )
1,2-Dichlorobenzene	30	30.7	102	30	30.2	101	(80-119)	1.80	(< 20 )
1,2-Dichloroethane	30	29.7	99	30	29.3	98	(73-128)	1.40	(< 20 )
1,2-Dichloropropane	30	30.8	103	30	30.2	101	(78-122)	2.00	(< 20 )
1,3,5-Trimethylbenzene	30	30.3	101	30	29.9	100	(75-124)	1.50	(< 20 )
1,3-Dichlorobenzene	30	30.8	103	30	30.4	101	(80-119)	1.50	(< 20 )
1,3-Dichloropropane	30	30.9	103	30	30.1	100	(80-119)	2.40	(< 20 )
1,4-Dichlorobenzene	30	30.9	103	30	30.1	100	(79-118)	2.70	(< 20 )
2,2-Dichloropropane	30	31.2	104	30	30.5	102	(60-139)	2.20	(< 20 )
2-Butanone (MEK)	90	77.5	86	90	84.0	93	(56-143)	8.00	(< 20 )
2-Chlorotoluene	30	29.9	100	30	29.2	98	(79-122)	2.10	(< 20 )
2-Hexanone	90	85.8	95	90	87.9	98	(57-139)	2.40	(< 20 )
4-Chlorotoluene	30	30.6	102	30	29.9	100	(78-122)	2.30	(< 20 )
4-Isopropyltoluene	30	31.2	104	30	30.6	102	(77-127)	1.90	(< 20 )
4-Methyl-2-pentanone (MIBK)	90	88.6	98	90	89.6	100	(67-130)	1.20	(< 20 )
Benzene	30	30.8	103	30	29.9	100	(79-120)	3.00	(< 20 )
Bromobenzene	30	30.8	103	30	30.0	100	(80-120)	2.60	(< 20 )
Bromochloromethane	30	29.9	100	30	29.7	99	(78-123)	0.82	(< 20 )
Bromodichloromethane	30	30.5	102	30	30.1	100	(79-125)	1.50	(< 20 )
Bromoform	30	28.8	96	30	28.7	96	(66-130)	0.49	(< 20 )
Bromomethane	30	28.3	95	30	28.4	95	(53-141)	0.26	(< 20 )
Carbon disulfide	45	45.1	100	45	44.1	98	(64-133)	2.20	(< 20 )

Print Date: 10/20/2021 9:08:28AM



Blank Spike ID: LCS for HBN 1216680 [VXX38039]

Blank Spike Lab ID: 1642871 Date Analyzed: 10/16/2021 13:51

QC for Samples: 1216680001, 1216680003

Spike Duplicate ID: LCSD for HBN 1216680

[VXX38039]

Spike Duplicate Lab ID: 1642872 Matrix: Water (Surface, Eff., Ground)

#### Results by SW8260D

	Blank Spike	e (ug/L)	ıg/L) Spike Duplicate (ug/L)						
<u>Parameter</u>	Spike	Result	Rec (%)	<u>Spike</u>	Result	Rec (%)	CL	RPD (%)	RPD CL
Carbon tetrachloride	30	30.8	103	30	30.1	100	(72-136)	2.10	(< 20)
Chlorobenzene	30	29.9	100	30	29.4	98	(82-118)	1.80	(< 20)
Chloroethane	30	29.0	97	30	28.7	96	(60-138)	0.95	(< 20)
Chloroform	30	28.5	95	30	28.0	93	(79-124)	1.50	(< 20)
Chloromethane	30	28.2	94	30	27.9	93	(50-139)	0.75	(< 20)
cis-1,2-Dichloroethene	30	30.1	100	30	29.8	99	(78-123)	1.20	(< 20)
cis-1,3-Dichloropropene	30	30.8	103	30	30.2	101	(75-124)	2.00	(< 20 )
Dibromochloromethane	30	30.8	103	30	30.2	101	(74-126)	1.90	(< 20 )
Dibromomethane	30	30.5	102	30	30.1	100	(79-123)	1.50	(< 20 )
Dichlorodifluoromethane	30	27.9	93	30	27.4	91	(32-152)	2.00	(< 20 )
Ethylbenzene	30	29.8	99	30	29.3	98	(79-121)	1.70	(< 20)
Freon-113	45	46.0	102	45	45.0	100	(70-136)	2.10	(< 20 )
Hexachlorobutadiene	30	31.0	103	30	31.0	103	(66-134)	0.07	(< 20 )
Isopropylbenzene (Cumene)	30	30.5	102	30	29.9	100	(72-131)	2.00	(< 20 )
Methylene chloride	30	29.0	97	30	28.7	96	(74-124)	1.20	(< 20 )
Methyl-t-butyl ether	45	45.1	100	45	44.7	99	(71-124)	0.91	(< 20 )
Naphthalene	30	30.0	100	30	32.0	107	(61-128)	6.30	(< 20 )
n-Butylbenzene	30	30.1	100	30	29.6	99	(75-128)	1.70	(< 20 )
n-Propylbenzene	30	31.1	104	30	30.0	100	(76-126)	3.40	(< 20 )
o-Xylene	30	29.9	100	30	29.3	98	(78-122)	2.20	(< 20 )
P & M -Xylene	60	59.3	99	60	58.0	97	(80-121)	2.30	(< 20 )
sec-Butylbenzene	30	31.2	104	30	30.3	101	(77-126)	2.90	(< 20 )
Styrene	30	29.2	97	30	28.5	95	(78-123)	2.20	(< 20 )
tert-Butylbenzene	30	30.9	103	30	30.2	101	(78-124)	2.60	(< 20 )
Tetrachloroethene	30	31.3	104	30	30.7	102	(74-129)	2.10	(< 20 )
Toluene	30	30.3	101	30	29.7	99	(80-121)	1.80	(< 20 )
trans-1,2-Dichloroethene	30	30.5	102	30	30.0	100	(75-124)	1.70	(< 20 )
trans-1,3-Dichloropropene	30	31.0	103	30	30.3	101	(73-127)	2.30	(< 20 )
Trichloroethene	30	30.9	103	30	30.3	101	(79-123)	2.10	(< 20 )
Trichlorofluoromethane	30	29.1	97	30	28.8	96	(65-141)	1.00	(< 20 )
Vinyl acetate	30	30.0	100	30	29.8	100	(54-146)	0.69	(< 20 )
Vinyl chloride	30	29.3	98	30	28.8	96	(58-137)	2.00	(< 20 )
Xylenes (total)	90	89.2	99	90	87.2	97	(79-121)	2.30	(< 20 )

Print Date: 10/20/2021 9:08:28AM



Blank Spike ID: LCS for HBN 1216680 [VXX38039]

Blank Spike Lab ID: 1642871 Date Analyzed: 10/16/2021 13:51

QC for Samples: 1216680001, 1216680003

Spike Duplicate ID: LCSD for HBN 1216680

[VXX38039]

Spike Duplicate Lab ID: 1642872 Matrix: Water (Surface, Eff., Ground)

#### Results by SW8260D

	Blank Spike (%)			Spike Duplicate (%)					
<u>Parameter</u>	<u>Spike</u>	Result	Rec (%)	Spike	Result	Rec (%)	CL	RPD (%)	RPD CL
Surrogates									
1,2-Dichloroethane-D4 (surr)	30		98	30		98	(81-118)	0.19	
4-Bromofluorobenzene (surr)	30		100	30		99	(85-114)	1.10	
Toluene-d8 (surr)	30		100	30		100	(89-112)	0.02	

#### **Batch Information**

Analytical Batch: VMS21286 Analytical Method: SW8260D Instrument: Agilent 7890-75MS

Analyst: JMG

Prep Batch: VXX38039
Prep Method: SW5030B

Prep Date/Time: 10/16/2021 06:00

Spike Init Wt./Vol.: 30 ug/L Extract Vol: 5 mL Dupe Init Wt./Vol.: 30 ug/L Extract Vol: 5 mL

Print Date: 10/20/2021 9:08:28AM



Blank ID: MB for HBN 1826750 [XXX/45697]

Blank Lab ID: 1640955

QC for Samples: 1216680001

Matrix: Water (Surface, Eff., Ground)

#### Results by 8270D SIM LV (PAH)

Parameter	Results	LOQ/CL	<u>DL</u>	Units
1-Methylnaphthalene	0.0250U	0.0500	0.0150	ug/L
2-Methylnaphthalene	0.0163J	0.0500	0.0150	ug/L
Acenaphthene	0.0250U	0.0500	0.0150	ug/L
Acenaphthylene	0.0250U	0.0500	0.0150	ug/L
Anthracene	0.0250U	0.0500	0.0150	ug/L
Benzo(a)Anthracene	0.0250U	0.0500	0.0150	ug/L
Benzo[a]pyrene	0.0100U	0.0200	0.00620	ug/L
Benzo[b]Fluoranthene	0.0250U	0.0500	0.0150	ug/L
Benzo[g,h,i]perylene	0.0250U	0.0500	0.0150	ug/L
Benzo[k]fluoranthene	0.0250U	0.0500	0.0150	ug/L
Chrysene	0.0250U	0.0500	0.0150	ug/L
Dibenzo[a,h]anthracene	0.0100U	0.0200	0.00620	ug/L
Fluoranthene	0.0250U	0.0500	0.0150	ug/L
Fluorene	0.0250U	0.0500	0.0150	ug/L
Indeno[1,2,3-c,d] pyrene	0.0250U	0.0500	0.0150	ug/L
Naphthalene	0.0500U	0.100	0.0310	ug/L
Phenanthrene	0.0214J	0.0500	0.0150	ug/L
Pyrene	0.0250U	0.0500	0.0150	ug/L
Surrogates				
2-Methylnaphthalene-d10 (surr)	59	42-86		%
Fluoranthene-d10 (surr)	69.2	50-97		%

#### **Batch Information**

Analytical Batch: XMS12943

Analytical Method: 8270D SIM LV (PAH)

Instrument: Agilent GC 7890B/5977A SWA

Analyst: LAW

Analytical Date/Time: 10/11/2021 9:07:00PM

Prep Batch: XXX45697 Prep Method: SW3535A

Prep Date/Time: 10/9/2021 9:14:19AM

Prep Initial Wt./Vol.: 250 mL Prep Extract Vol: 1 mL

Print Date: 10/20/2021 9:08:30AM



Blank Spike ID: LCS for HBN 1216680 [XXX45697]

Blank Spike Lab ID: 1640956 Date Analyzed: 10/11/2021 21:27

QC for Samples: 1216680001

Spike Duplicate ID: LCSD for HBN 1216680

[XXX45697]

Spike Duplicate Lab ID: 1640957 Matrix: Water (Surface, Eff., Ground)

#### Results by 8270D SIM LV (PAH)

,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,	•	Blank Spike	e (ug/L)	;	Spike Duplicate (ug/L)				
<u>Parameter</u>	Spike	Result	Rec (%)	<u>Spike</u>	Result	Rec (%)	CL	RPD (%)	RPD CL
1-Methylnaphthalene	2	1.23	61	2	1.35	68	(41-115)	9.60	(< 20 )
2-Methylnaphthalene	2	1.22	61	2	1.33	67	(39-114)	8.70	(< 20 )
Acenaphthene	2	1.43	72	2	1.59	80	(48-114)	10.60	(< 20 )
Acenaphthylene	2	1.46	73	2	1.65	82	(35-121)	11.90	(< 20 )
Anthracene	2	1.60	80	2	1.73	87	(53-119)	7.80	(< 20 )
Benzo(a)Anthracene	2	1.37	68	2	1.51	75	(59-120)	9.70	(< 20 )
Benzo[a]pyrene	2	1.86	93	2	2.09	105	(53-120)	11.70	(< 20 )
Benzo[b]Fluoranthene	2	1.72	86	2	1.81	91	(53-126)	5.50	(< 20 )
Benzo[g,h,i]perylene	2	1.99	99	2	2.38	119	(44-128)	18.20	(< 20 )
Benzo[k]fluoranthene	2	1.88	94	2	1.94	97	(54-125)	3.40	(< 20 )
Chrysene	2	1.54	77	2	1.65	83	(57-120)	7.30	(< 20 )
Dibenzo[a,h]anthracene	2	1.77	89	2	2.40	120	(44-131)	29.90	* (< 20 )
Fluoranthene	2	1.45	73	2	1.51	76	(58-120)	4.20	(< 20 )
Fluorene	2	1.55	78	2	1.73	86	(50-118)	10.70	(< 20 )
Indeno[1,2,3-c,d] pyrene	2	1.71	86	2	2.33	117	(48-130)	30.80	* (< 20 )
Naphthalene	2	1.19	59	2	1.32	66	(43-114)	10.80	(< 20)
Phenanthrene	2	1.58	79	2	1.73	87	(53-115)	9.00	(< 20 )
Pyrene	2	1.46	73	2	1.54	77	(53-121)	5.40	(< 20 )
Surrogates									
2-Methylnaphthalene-d10 (surr)	2		55	2		63	(42-86)	13.50	
Fluoranthene-d10 (surr)	2		68	2		71	(50-97)	4.90	

#### **Batch Information**

Analytical Batch: XMS12943

Analytical Method: 8270D SIM LV (PAH)
Instrument: Agilent GC 7890B/5977A SWA

Analyst: LAW

Prep Batch: XXX45697 Prep Method: SW3535A

Prep Date/Time: 10/09/2021 09:14

Spike Init Wt./Vol.: 2 ug/L Extract Vol: 1 mL Dupe Init Wt./Vol.: 2 ug/L Extract Vol: 1 mL

Print Date: 10/20/2021 9:08:32AM



Blank ID: MB for HBN 1826768 [XXX/45702]

Blank Lab ID: 1641067

QC for Samples:

1216680001, 1216680002

Matrix: Water (Surface, Eff., Ground)

#### Results by AK102

 Parameter
 Results
 LOQ/CL
 DL
 Units

 Diesel Range Organics
 0.300U
 0.600
 0.200
 mg/L

**Surrogates** 

5a Androstane (surr) 99.5 60-120 %

#### **Batch Information**

Analytical Batch: XFC16112 Analytical Method: AK102 Instrument: Agilent 7890B R

Analyst: IVM

Analytical Date/Time: 10/13/2021 6:16:00PM

Prep Batch: XXX45702 Prep Method: SW3520C

Prep Date/Time: 10/9/2021 3:45:03PM

Prep Initial Wt./Vol.: 250 mL Prep Extract Vol: 1 mL

Print Date: 10/20/2021 9:08:35AM



Blank Spike ID: LCS for HBN 1216680 [XXX45702]

Blank Spike Lab ID: 1641068 Date Analyzed: 10/13/2021 18:26 Spike Duplicate ID: LCSD for HBN 1216680

[XXX45702]

Spike Duplicate Lab ID: 1641069 Matrix: Water (Surface, Eff., Ground)

QC for Samples: 1216680001, 1216680002

#### Results by AK102

		Blank Spike	e (mg/L)	5	Spike Dupli	cate (mg/L)			
<u>Parameter</u>	<u>Spike</u>	Result	Rec (%)	Spike	Result	Rec (%)	CL	RPD (%)	RPD CL
Diesel Range Organics	20	19.7	98	20	19.9	100	(75-125)	1.40	(< 20 )
Surrogates									
5a Androstane (surr)	0.4		106	0.4		110	(60-120)	3.00	

#### **Batch Information**

Analytical Batch: XFC16112 Analytical Method: AK102

Instrument: Agilent 7890B R

Analyst: IVM

Prep Batch: XXX45702 Prep Method: SW3520C

Prep Date/Time: 10/09/2021 15:45

Spike Init Wt./Vol.: 20 mg/L Extract Vol: 1 mL Dupe Init Wt./Vol.: 20 mg/L Extract Vol: 1 mL

Print Date: 10/20/2021 9:08:36AM



Blank ID: MB for HBN 1826768 [XXX/45702]

Blank Lab ID: 1641067

QC for Samples:

1216680001, 1216680002

Matrix: Water (Surface, Eff., Ground)

#### Results by AK103

 Parameter
 Results
 LOQ/CL
 DL
 Units

 Residual Range Organics
 0.250U
 0.500
 0.200
 mg/L

**Surrogates** 

n-Triacontane-d62 (surr) 103 60-120 %

#### **Batch Information**

Analytical Batch: XFC16112 Analytical Method: AK103

Instrument: Agilent 7890B R

Analyst: IVM

Analytical Date/Time: 10/13/2021 6:16:00PM

Prep Batch: XXX45702 Prep Method: SW3520C

Prep Date/Time: 10/9/2021 3:45:03PM

Prep Initial Wt./Vol.: 250 mL Prep Extract Vol: 1 mL

Print Date: 10/20/2021 9:08:38AM



Blank Spike ID: LCS for HBN 1216680 [XXX45702]

Blank Spike Lab ID: 1641068 Date Analyzed: 10/13/2021 18:26

QC for Samples: 1216680001, 1216680002 Spike Duplicate ID: LCSD for HBN 1216680

[XXX45702]

Spike Duplicate Lab ID: 1641069 Matrix: Water (Surface, Eff., Ground)

#### Results by AK103

		Blank Spike	(mg/L)	5	Spike Duplic	cate (mg/L)			
<u>Parameter</u>	<u>Spike</u>	Result	Rec (%)	Spike	Result	Rec (%)	CL	RPD (%)	RPD CL
Residual Range Organics	20	19.3	97	20	19.7	99	(60-120)	1.90	(< 20 )
Surrogates									
n-Triacontane-d62 (surr)	0.4		90	0.4		97	(60-120)	7.80	

#### **Batch Information**

Analytical Batch: XFC16112 Analytical Method: AK103 Instrument: Agilent 7890B R

Analyst: IVM

Prep Batch: XXX45702 Prep Method: SW3520C

Prep Date/Time: 10/09/2021 15:45

Spike Init Wt./Vol.: 20 mg/L Extract Vol: 1 mL Dupe Init Wt./Vol.: 20 mg/L Extract Vol: 1 mL

Print Date: 10/20/2021 9:08:41AM

						SGS	North A	merica In	ic.		
Shannon & Wilson, Inc. 5430 Fairbanks Street, Suite 3 Anchorage, Alaska 99518 (907) 561-2120 Fax (206) 695-6777					VOCs - EPA 8260D	PAHs - EPA 8270D SIM		12 <sup>-</sup>	166 <b>:</b>		
Date	Time	Sample ID	Total Containers	-							
10-7-21	12:30	101528-B1MW	7	X	Х	Х					1
10-7-21	13:00	101528-B11MW	2	Х							
10-7-21	9106	101528-WTB	1 BOX		X						
Relinquished	By:	Relinquished By:	The state of the s		Project Info	rmation	dir vije diegrees die die große				
Signature:	144	Signature:			Project Num	ber: 10152	8-002				
Print Name:	Zaen +	Print Name:		Project Name: 5701 Northwood Drive							
Company: Sh	annon & Wilson	, Inc. Company:			Contact: Da	n McMahon					
Date: 10		Date:			<del>                                     </del>	ZJT					
Гіте: 13°	51	Time:			Special Instr	ON MADE AND A STATE OF THE STAT	26 240 E B	73 (F - 5188) (17 <sup>18</sup> (F - 17 <sup></sup>		SIISKELL I	FE 270 - 10 - 10
Received By:		Received By:	a RSC		Sample Re					<b>4</b> 0 50 70 70 10 10 10 10 10 10 10 10 10 10 10 10 10	
Signature:		Signature: Office Class			Shipped V	a: Hand D	elivered	•			
Print Name: Print Name: Ryan Conon					Cooler Ten	anaratura I	Inon Amire	1.			
Company:		Company: 5'G5  Date: 10/7/2\									
1,260					Sample Matrix: Water  10 Working DAY TAT						
Time:		Time' 1 1 2.			110 WOLKINS	DAIJAI					









#### SGS North America Inc.

200 W. Potter Dr., 3180 Peger Rd. Ste. Anchorage, AK 99518 (ph) 190, Fairbanks, AK 907-562-2343, (fax) 907-561-5301 907-474-

# **Sample Kit Request**

Client pickup Date:

10/1/2021

Time:

12:00

	***				Be sure to ask if clien	t will ship by groui	nd (DOT) or air	carrier (IATA)
Does a Pr	rofile exist in LIMS? If not, ple	ase send a request for new profile build.			Deliver to client:			
Client Name:	Shannon d	& Wilson Inc	-		Ship by/Air Carrier:			
Ordered By:	Alec Rizzo		-		Airbill Number:			
Email:	AJR@s	nanwil.com	_		Date to ship by:			
Project Name:	5701 Northwood Drive		-		Notes:			
Quote #:			-		Kit request taken by:	JAN	Date:	September 30, 2021
_			•		Kit prepared by:	EBH	Date: _	9/30/21
Delivery Address:			Kit <i>(in</i>	cluding lid tightness for	r pres'd bottles) checked by:	-15	Date: _	9/30/21
					Kit packed & shipped by:	ીંક	Date: _	9/30/21
Filename: SK	IT_Shannon & Wilson Inc_5701 Northwood Driv	e_2 *Required Items						1 -
No.					Preservative	Hold	# QC	Total
Samples Matrix	Analysis	Container Size & Type	Pres.	Bottle Lot #	Lot#	Time	Bottles	Bottles

No.							Preservative	Hold	#QC	Total
Samples	Matrix	Analysis	Container	Size & Type	Pres.	Bottle Lot #	Lot#	Time	Bottles	Bottles
2	Water	DRO/RRO	2 x 250 mL	Amber	HCI			14 d	0	4
1	Water	VOC	3 x 40 mL	VOA	HCI			14 d	0	3
1	Water	PAH	2 x 250 mL	Amber	None			7 d	0	2

Note: The first 10 Analysis and Preservative columns will auto-fill up to the capacity of the associated COC.

	Note. The mist to Analysis and Freservative column	s will auto-fill up to the capacity of the associate	u 000.
Addit	ional Information	Notes for Kit Prep	Attention Client/Sampler:
Pack for Shipment via:	N/A		Do not rinse container, be aware of any acid preservative.
Temperature Blank:	Yes - Small (125 mL)		Fill container, but do not overfill (except volatiles).
Trip Blank:	Yes - Water (8260, AK101, 8021, 624)	· · · · · · · · · · · · · · · · · · ·	3. Label the container with your sample ID and date/time of collection
Coolers:	Yes		4. Fill out the Chain of Custody.
Gel Ice:	Yes		5. Add frozen gel packs to your cooler and pack to prevent breakage
Labels:	Yes		If you have any questions please contact your Project Manager.
Custody Seals:	Yes		
Paper Chain of Custody:	No - Electronic or Client will Provide COC		
Lot Number Tracking (Required for DOD):	No		



e-Sample Receipt Form

SGS Workorder #:

1216680

1216680

<u> </u>						
Review Criteria	Condition (Yes	, No, N/A		Exception	ns Noted below	
Chain of Custody / Temperature Require	rements		N/A	Exemption permitted	d if sampler hand carries/delive	ers.
Were Custody Seals intact? Note # & I						
COC accompanied sa						
·						
DOD: Were samples received in COC corresponding c						
N/A **Exemption permitted if	chilled & colle	ected <8 h	ours	ago, or for samples w	here chilling is not required	
Temperature blank compliant* (i.e., 0-6 °C afte		Cooler I	_		@ 2.7 °C Therm. ID:	D58
(iiii)	N/A				@ °C Therm. ID:	
			_			
If samples received without a temperature blank, the "cooler temperature" will documented instead & "COOLER TEMP" will be noted to the right. "ambient" or "chi		Cooler I	D:		@ °C Therm. ID:	
be noted if neither is available.		Cooler I	D:		@ C Therm. ID:	
		Cooler I	D:		@ °C Therm. ID:	
*If >6°C, were samples collected <8 hours	ago? N/A					
	1.471	4				
W 202						
If <0°C, were sample containers ice	ree? Yes	Į				
		Ī				
Note: Identify containers received at non-compliant temper	rature .					
Use form FS-0029 if more space is no						
Holding Time / Documentation / Sample Condition Re		Note: Refe	r to fo	orm F-083 "Sample Guide	e" for specific holding times.	
Were samples received within holding	g time? Yes					
		Ĭ				
Do samples match COC** (i.e.,sample IDs,dates/times colle	cted)? Ves					
		1				
**Note: If times differ <1hr, record details & login per Co						
***Note: If sample information on containers differs from COC, SGS will default to C	COC information	ì				
Were analytical requests clear? (i.e., method is specified for an	alyses Yes					
with multiple option for analysis (Ex: BTEX, N	Metals)	Ī				
			N/A	***Evamption normitt	ted for metals (e.g,200.8/6020	וםו
M	10		N/A	Exemption permitt	teu foi filetais (e.g,200.6/6020	<u>וטו.</u>
Were proper containers (type/mass/volume/preservative***)	)used? Yes	Ų				
		1				
Volatile / LL-Hg Req	uirements					
Were Trip Blanks (i.e., VOAs, LL-Hg) in cooler with sar	mples? Yes					
Were all water VOA vials free of headspace (i.e., bubbles ≤ 6		1				
Were all soil VOAs field extracted with MeOH-	+BFB! N/A	<u> </u>				
Note to Client: Any "No", answer above indicates nor	n-compliance	with stand	dard p	procedures and may i	mpact data quality.	
A 1 100			- \			
Additiona	I <mark>l notes (if a</mark>	applicable	e):			



#### **Sample Containers and Preservatives**

Container Id	<u>Preservative</u>	Container Condition	Container Id	<u>Preservative</u>	Container Condition
1216680001-A	HCL to pH < 2	OK			
1216680001-B	HCL to pH < 2	ОК			
1216680001-C	No Preservative Required	OK			
1216680001-D	No Preservative Required	ОК			
1216680001-E	HCL to pH < 2	ОК			
1216680001-F	HCL to pH < 2	OK			
1216680001-G	HCL to pH < 2	OK			
1216680002-A	HCL to pH < 2	ОК			
1216680002-B	HCL to pH < 2	OK			
1216680003-A	HCL to pH < 2	OK			
1216680003-B	HCL to pH < 2	OK			
1216680003-C	HCL to pH < 2	ОК			

#### **Container Condition Glossary**

Containers for bacteriological, low level mercury and VOA vials are not opened prior to analysis and will be assigned condition code OK unless evidence indicates than an inappropriate container was submitted.

- OK The container was received at an acceptable pH for the analysis requested.
- BU The container was received with headspace greater than 6mm.
- DM The container was received damaged.
- FR The container was received frozen and not usable for Bacteria or BOD analyses.
- IC The container provided for microbiology analysis was not a laboratory-supplied, pre-sterilized container and therefore was not suitable for analysis.
- NC- The container provided was not preserved or was under-preserved. The method does not allow for additional preservative added after collection.
- PA The container was received outside of the acceptable pH for the analysis requested. Preservative was added upon receipt and the container is now at the correct pH. See the Sample Receipt Form for details on the amount and lot # of the preservative added.
- PH The container was received outside of the acceptable pH for the analysis requested. Preservative was added upon receipt, but was insufficient to bring the container to the correct pH for the analysis requested. See the Sample Receipt Form for details on the amount and lot # of the preservative added. QN Insufficient sample quantity provided.

31 of 31

#### LABORATORY DATA REVIEW CHECKLIST

Completed by: Chris Pepe
Title: Environmental Scientist

Date: December 2021

Consultant Firm: Shannon & Wilson, Inc.

Laboratory Name: SGS North America Inc. Laboratory Report Number: 1216680 Laboratory Report Date: 10/20/2021

Contaminated Site Name: 5701 Northwood Drive

**ADEC File Number:** 2100.38.536 **Hazard Identification Number:** 25942

(**NOTE**: *NA* = not applicable; Text in *italics* added by Shannon & Wilson, Inc.)

#### 1. <u>Laboratory</u>

a. Did an ADEC CS approved laboratory receive and <u>perform</u> all of the submitted sample analyses? Yes/ No / NA
 Comments:

**b.** If the samples were transferred to another "network" laboratory or sub-contracted to an alternate laboratory, was the laboratory performing the analyses ADEC CS approved? **Yes / No (NA)** 

Comments: The samples were not transferred to another "network" laboratory or subcontracted to an alternate laboratory.

#### 2. Chain of Custody (COC)

a. COC information completed, signed, and dated (including released/received by)?
 Yes/ No / NA
 Comments:

b. Correct analyses requested? Yes / No / NA Comments:

#### 3. Laboratory Sample Receipt Documentation

a. Sample/cooler temperature documented and within range at receipt (0° to 6° C)? (Yes)/ No / NA

Comments: *The cooler temperature blank was 2.7° Celsius.* 

b. Sample preservation acceptable - acidified waters, Methanol preserved VOC soil (GRO, BTEX, VOCs, etc.)? Yes/ No / NA Comments:

c. Sample condition documented - broken, leaking (MeOH), zero headspace (VOC vials)?
Yes/ No / NA

Comments:

**d.** If there were any discrepancies, were they documented? For example, incorrect sample containers/preservation, sample temperature outside of acceptable range, insufficient or missing samples, etc.? **Yes / No /NA** 

Comments: No discrepancies were noted

e. Data quality or usability affected?

Comments: See above.

#### 4. Case Narrative

- a. Present and understandable? Yes/ No / NA Comments:
- **b.** Discrepancies, errors or QC failures noted by the lab? **Yes**/**No**/**NA** Comments: *The case narrative noted the following:* 
  - 8270D SIM PAH LCS/LCSD RPDs for ideno[1,2,3-c,d]pyrene and dibenzo[a,h]anthracene do not meet QC criteria. These analytes were not detected above the LOQ in associated samples.
- c. Were all corrective actions documented? **Yes** No / NA Comments:
- **d.** What is the effect on data quality/usability, according to the case narrative? Comments: *See above*.

#### 5. Sample Results

- a. Correct analyses performed/reported as requested on COC? Ves/No/NA Comments:
- **b.** All applicable holding times met? **Yes** / **No** / **NA** Comments:
- c. All soils reported on a dry weight basis? Yes / No / NA Comments:

d. Are the reported LOQs less than the Cleanup Level or the minimum required detection level for the project? Yes No NA Comments: The LOQ for 1,2,3-trichloropropane are greater than the ADEC cleanup level.

**e.** Data quality or usability affected? Comments: There is a potential that the target analyte is present at concentrations greater than the ADEC cleanup level, but less than the LOO.

#### 6. QC Samples

#### a. Method Blank

One method blank reported per matrix, analysis, and 20 samples?
 Yes/ No / NA
 Comments:

**ii.** All method blank results less than limit of quantitation (LOQ) or project specified objectives?

(Yes)/ No / NA

Comments: The method blank associated with Sample B1MW contained estimated concentrations of 2-methylnaphthalene (0.0163 J  $\mu$ g/L) and phenanthrene (0.0214 J  $\mu$ g/L) less than the LOQ.

- **iii.** If above LOQ or project specified objectives, what samples are affected? Comments: *See above*.
- iv. Do the affected sample(s) have data flags? If so, are the data flags clearly defined?

  Yes / No (NA)

Comments: Samples are flagged "B" in Table 2B when the reported sample concentration is within 10x the reported method blank concentration. Estimated concentrations of 2-methylnaphthalene and phenanthrene were detected in Sample B1MW and the method blank, therefore, the sample concentrations are reported as non-detect at the LOQ and flagged "B".

v. Data quality or usability affected? Comments: *See above*.

#### b. Laboratory Control Sample/Duplicate (LCS/LCSD)

i. Organics - One LCS/LCSD reported per matrix, analysis, and 20 samples?
 (LCS/LCSD required per AK methods, LCS required per SW846) (Ves) No / NA Comments:

- ii. Metals/Inorganics One LCS and one sample duplicate reported per matrix, analysis and 20 samples? Yes / No NA Comments:
- iii. Accuracy All percent recoveries (%R) reported and within method or laboratory limits and project specified objectives, if applicable. (AK petroleum methods: AK 101 60%-120%, AK 102 75%-125%, AK 103 60%-120%; all other analyses see the laboratory QC pages) Yes/No/NA Comments:
- iv. Precision All relative percent differences (RPDs) reported and less than method or laboratory limits and project specified objectives, if applicable. RPD reported from LCS/LCSD, and/or sample/sample duplicate. (AK Petroleum methods 20%; all other analyses see the laboratory QC pages) Yes No/NA Comments: RPDs for indeno[1,2,3-c,d]pyrene and dibenzo[a,h]anthracene do not meet QC criteria.
- **v.** If %R or RPD is outside of acceptable limits, what samples are affected? Comments: *Sample B1MW*
- vi. Do the affected samples(s) have data flags? If so, are the data flags clearly defined? Yes No NA

Comments: *Indeno*[1,2,3-c,d]*pyrene and dibenzo*[a,h]*anthracene were not detected above the LOQ in Sample B1MW, therefore, flagging is not required.* 

**vii.** Data quality or usability affected? Comments: *No, see above.* 

- c. Matrix Spike/Matrix Spike Duplicate (MS/MSD)
  Note: Leave blank if not required for project
  - i. Organics One MS/MSD reported per matrix, analysis, and 20 samples?Yes/ No / NAComments:
  - ii. Metals/Inorganics One MS and one MSD reported per matrix, analysis and 20 samples? Yes / No NA Comments:
  - iii. Accuracy All percent recoveries (%R) reported and within method or laboratory limits and project specified objectives, if applicable. (AK petroleum methods: AK 101 60%-120%, AK 102 75%-125%, AK 103 60%-120%; all other analyses see the laboratory QC pages) Yes / No / NA Comments:

- iv. Precision All relative percent differences (RPDs) reported and less than method or laboratory limits and project specified objectives, if applicable. RPD reported from MS/MSD, and/or sample/sample duplicate. (AK Petroleum methods 20%; all other analyses see the laboratory QC pages) Yes / No / NA Comments:
- v. If %R or RPD is outside of acceptable limits, what samples are affected? Comments:
- vi. Do the affected samples(s) have data flags? If so, are the data flags clearly defined? Yes / No (NA)

Comments: No, see above.

vii. Data quality or usability affected?

Comments: No, see above.

# d. Surrogates - Organics Only or Isotope Dilution Analytes (IDA) - Isotope Dilution Methods Only

- i. Are surrogate/IDA recoveries reported for organic analyses field, QC, and laboratory samples? Yes/ No / NA
  Comments:
- ii. Accuracy All percent recoveries (%R) reported and within method or laboratory limits and project specified objectives, if applicable. (AK Petroleum methods 50-150 %R; all other analyses see the laboratory report pages) Yes / No / NA Comments:
- iii. Do the sample results with failed surrogate/IDA recoveries have data flags? If so, are the data flags clearly defined? Yes / No / NA Comments:
- **iv.** Data quality or usability affected? Comments: *No. see above.*
- e. Trip Blank Volatile analyses only (GRO, BTEX, VOCs, etc.)
  - i. One trip blank reported per matrix, analysis and for each cooler containing volatile samples? Yes/No/NA
     Comments:
  - ii. Is the cooler used to transport the trip blank and volatile samples clearly indicated on the COC? Yes No NA
    - Comments: *Only one cooler was used to transport the samples.*
  - iii. All results less than LOQ and project specified objectives? Yes/ No / NA Comments:

- iv. If above LOQ or project specified DQOs, what samples are affected? Comments:
- **v.** Data quality or usability affected? Comments: *No, see above.*

### f. Field Duplicate

i. One field duplicate submitted per matrix, analysis and 10 project samples? Yes/ No / NA

Comments: Duplicate Sample B11MW (duplicate of B1MW) was submitted to the laboratory.

- ii. Were the field duplicates submitted blind to the lab? Yes/ No / NA Comments:
- iii. Precision All relative percent differences (RPDs) less than specified project objectives? (Recommended: 30% for water, 50% for soil Yes / No / NA Comments:
- **iv.** Data quality or usability affected? Comments:
- **g. Decontamination or Equipment Blank** (if not applicable, a comment stating why must be entered below).

Yes /No NA

Comments: A decontamination blank or equipment blank was not included in our ADEC-approved workplan.

i. All results less than LOQ and project specified objectives?

Yes / No NA Comments:

- **ii.** If above LOQ or project specified objectives, what samples are affected? Comments:
- **iii.** Data quality or usability affected? Comments:

## 7. Other Data Flags/Qualifiers (ACOE, AFCEE, Lab Specific, etc.)

**a.** Defined and appropriate? Yes / No / NA Comments: A key is provided on Page 3 of the SGS Laboratory Report.

# SHANNON & WILSON, INC.

# **ATTACHMENT 4**

# IMPORTANT INFORMATION ABOUT YOUR GEOTECHNICAL/ENVIRONMENTAL REPORT

Attachment to and part of Report 101528-002

Date: January 2022

To: 5701 Northwood Drive

# IMPORTANT INFORMATION ABOUT YOUR GEOTECHNICAL/ENVIRONMENTAL REPORT

#### CONSULTING SERVICES ARE PERFORMED FOR SPECIFIC PURPOSES AND FOR SPECIFIC CLIENTS.

Consultants prepare reports to meet the specific needs of specific individuals. A report prepared for a civil engineer may not be adequate for a construction contractor or even another civil engineer. Unless indicated otherwise, your consultant prepared your report expressly for you and expressly for the purposes you indicated. No one other than you should apply this report for its intended purpose without first conferring with the consultant. No party should apply this report for any purpose other than that originally contemplated without first conferring with the consultant.

#### THE CONSULTANT'S REPORT IS BASED ON PROJECT-SPECIFIC FACTORS.

A geotechnical/environmental report is based on a subsurface exploration plan designed to consider a unique set of project-specific factors. Depending on the project, these may include: the general nature of the structure and property involved; its size and configuration; its historical use and practice; the location of the structure on the site and its orientation; other improvements such as access roads, parking lots, and underground utilities; and the additional risk created by scope-of-service limitations imposed by the client. To help avoid costly problems, ask the consultant to evaluate how any factors that change subsequent to the date of the report may affect the recommendations. Unless your consultant indicates otherwise, your report should not be used: (1) when the nature of the proposed project is changed (for example, if an office building will be erected instead of a parking garage, or if a refrigerated warehouse will be built instead of an unrefrigerated one, or chemicals are discovered on or near the site); (2) when the size, elevation, or configuration of the proposed project is altered; (3) when the location or orientation of the proposed project is modified; (4) when there is a change of ownership; or (5) for application to an adjacent site. Consultants cannot accept responsibility for problems that may occur if they are not consulted after factors which were considered in the development of the report have changed.

#### SUBSURFACE CONDITIONS CAN CHANGE.

Subsurface conditions may be affected as a result of natural processes or human activity. Because a geotechnical/environmental report is based on conditions that existed at the time of subsurface exploration, construction decisions should not be based on a report whose adequacy may have been affected by time. Ask the consultant to advise if additional tests are desirable before construction starts; for example, groundwater conditions commonly vary seasonally.

Construction operations at or adjacent to the site and natural events such as floods, earthquakes, or groundwater fluctuations may also affect subsurface conditions and, thus, the continuing adequacy of a geotechnical/environmental report. The consultant should be kept apprised of any such events, and should be consulted to determine if additional tests are necessary.

#### MOST RECOMMENDATIONS ARE PROFESSIONAL JUDGMENTS.

Site exploration and testing identifies actual surface and subsurface conditions only at those points where samples are taken. The data were extrapolated by your consultant, who then applied judgment to render an opinion about overall subsurface conditions. The actual interface between materials may be far more gradual or abrupt than your report indicates. Actual conditions in areas not sampled may differ from those predicted in your report. While nothing can be done to prevent such situations, you and your consultant can work together to help reduce their impacts. Retaining your consultant to observe subsurface construction operations can be particularly beneficial in this respect.

Page 1 of 2 1/2016

#### A REPORT'S CONCLUSIONS ARE PRELIMINARY.

The conclusions contained in your consultant's report are preliminary because they must be based on the assumption that conditions revealed through selective exploratory sampling are indicative of actual conditions throughout a site. Actual subsurface conditions can be discerned only during earthwork; therefore, you should retain your consultant to observe actual conditions and to provide conclusions. Only the consultant who prepared the report is fully familiar with the background information needed to determine whether or not the report's recommendations based on those conclusions are valid and whether or not the contractor is abiding by applicable recommendations. The consultant who developed your report cannot assume responsibility or liability for the adequacy of the report's recommendations if another party is retained to observe construction.

#### THE CONSULTANT'S REPORT IS SUBJECT TO MISINTERPRETATION.

Costly problems can occur when other design professionals develop their plans based on misinterpretation of a geotechnical/environmental report. To help avoid these problems, the consultant should be retained to work with other project design professionals to explain relevant geotechnical, geological, hydrogeological, and environmental findings, and to review the adequacy of their plans and specifications relative to these issues.

#### BORING LOGS AND/OR MONITORING WELL DATA SHOULD NOT BE SEPARATED FROM THE REPORT.

Final boring logs developed by the consultant are based upon interpretation of field logs (assembled by site personnel), field test results, and laboratory and/or office evaluation of field samples and data. Only final boring logs and data are customarily included in geotechnical/environmental reports. These final logs should not, under any circumstances, be redrawn for inclusion in architectural or other design drawings, because drafters may commit errors or omissions in the transfer process.

To reduce the likelihood of boring log or monitoring well misinterpretation, contractors should be given ready access to the complete geotechnical engineering/environmental report prepared or authorized for their use. If access is provided only to the report prepared for you, you should advise contractors of the report's limitations, assuming that a contractor was not one of the specific persons for whom the report was prepared, and that developing construction cost estimates was not one of the specific purposes for which it was prepared. While a contractor may gain important knowledge from a report prepared for another party, the contractor should discuss the report with your consultant and perform the additional or alternative work believed necessary to obtain the data specifically appropriate for construction cost estimating purposes. Some clients hold the mistaken impression that simply disclaiming responsibility for the accuracy of subsurface information always insulates them from attendant liability. Providing the best available information to contractors helps prevent costly construction problems and the adversarial attitudes that aggravate them to a disproportionate scale.

#### READ RESPONSIBILITY CLAUSES CLOSELY.

Because geotechnical/environmental engineering is based extensively on judgment and opinion, it is far less exact than other design disciplines. This situation has resulted in wholly unwarranted claims being lodged against consultants. To help prevent this problem, consultants have developed a number of clauses for use in their contracts, reports, and other documents. These responsibility clauses are not exculpatory clauses designed to transfer the consultant's liabilities to other parties; rather, they are definitive clauses that identify where the consultant's responsibilities begin and end. Their use helps all parties involved recognize their individual responsibilities and take appropriate action. Some of these definitive clauses are likely to appear in your report, and you are encouraged to read them closely. Your consultant will be pleased to give full and frank answers to your questions.

The preceding paragraphs are based on information provided by the ASFE/Association of Engineering Firms Practicing in the Geosciences, Silver Spring, Maryland

Page 2 of 2 1/2016