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April 30, 2024

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Physical Scientist
Land, Chemicals, and Redevelopment Division – RCRA Branch
U.S. Environmental Protection Agency – Region 9
775 Hawthorne Street, LND-4-2
San Francisco, California 94105

Subject: First Quarter 2024 Groundwater Monitoring Results, Former Thermal Treatment Unit, Nammo Defense Systems Inc., Mesa, Arizona

Dear Ms. Clark:

Pinyon Environmental, Inc. (Pinyon), has prepared the following First Quarter 2024 (Q1 2024) Groundwater Monitoring Report (Report) on behalf of Nammo Defense Systems Inc. (NDS). This report documents field activities and results for groundwater sampling conducted at the NDS Former Thermal Treatment Unit (TTU) located in Mesa, Arizona (the Site; Figure 1). The monitoring activities were planned and executed following the scope of work and requirements outlined in the *Groundwater Water Sampling and Analysis Plan, Former Thermal Treatment Unit, Nammo Defense Systems Inc., Mesa Arizona*, dated September 30, 2022 (TTU SAP); and the *Quality Assurance Project Plan, Nammo Defense Systems Inc. Facility, Mesa, Arizona*, dated April 28, 2022 (NDS Facility QAPP). The finalized TTU SAP was submitted to the U.S. Environmental Protection Agency (EPA) on September 30, 2022. Any changes or deviations from these documents are provided in subsequent sections of this report.

Based on the *Data Management Plan, Nammo Defense Systems Inc., Mesa Arizona*, dated November 2022 (revised May 2023) groundwater data will be provided as electronic data deliverables (EDDs) with spatial data provided as shape files. The EDDs and shape files will be provided as a separate deliverable.

1. SCOPE OF ACTIVITIES

Groundwater monitoring and pumping/extraction wells were sampled between February 12 and 14, 2024. Well construction details are summarized in Table 1. A Site Vicinity Map and well locations are shown on Figures 1 and 2, respectively.

1.1 Deviations from Work Plan

The Q1 2024 groundwater monitoring was conducted in accordance with the TTU SAP and NDS Facility QAPP, with the following exceptions:

- Depth to groundwater was not obtained from well PF-2 as the water level meter hit an obstruction within the well at a depth of 168 feet while the well is 400 feet deep.

- TTU-3 was resampled due to an exceedance of 1,4-dioxane during the routine sampling event. 1,4-dioxane was not detected at a concentration above the laboratory reporting limit in the second sample.The laboratory analytical order will be updated to prevent suspected carryover.

- TTU-15 and TTU-16 had sample results that appeared to have been switched, they were subsequently resampled, and the confirmation analytical results were in line with historical results confirming the lab had switched the samples during analysis. The data presented in the figures and tables represent the results

from resampling. 1,4-dioxane was not included in the analytical suite during resampling; however, based on the results of the other analysis the February sample results were switched and those results have been rejected and are not considered valid for use on Figure 4.

1.2 Groundwater Elevation Measurements

Groundwater depth was measured utilizing an electronic water level indicator. Measurements were collected to the nearest 0.01 foot on the north side, top of casing at each well. Monitoring well TTU-18 is historically dry, and therefore no sample was collected during the Q1 event. Depth to water will be re-checked during the second quarter (Q2) 2024 groundwater monitoring event. A summary of groundwater elevation data for the Q1 2024 sampling event is provided in Table 2.

1.3 Groundwater Sampling

For extraction/pumping wells, the wells were activated and allowed to purge for at least 15 minutes prior to sample collection. Water was taken from the spigot closest to the wellhead. From each sampled well, field parameter measurements were collected utilizing a YSI Professional Plus water quality meter. Parameters collected include water temperature, pH, oxidation reduction potential (ORP), conductivity, and dissolved oxygen (DO). Turbidity was measured utilizing a 2020 LaMotte Turbidimeter. For the production wells TTU-1, TTU-2, TTU-20, and PF-2, field readings were collected every 5 minutes during the minimum 15-minute purging/stabilization period. For non-pumping wells, one round of field parameter measurements were collected at the time of sample collection. Field parameter stabilization forms for the Q1 2024 sampling event are provided in Attachment 1.

Monitoring wells and other non-pumping wells were sampled utilizing HydraSleeve samplers that were deployed by Pinyon following Q4 2024 sample collection at each well. The samplers were suspended inside the wells at the depths summarized in Table 3.

Groundwater samples were collected into laboratory provided and preserved sample containers based on analytical method requirements as described in the TTU SAP. Each groundwater sample was labeled, secured from breakage, and stored on-ice inside an insulated cooler. The samples were transported under chain-of-custody protocol to Pace Analytical and Eurofins-Phoenix for analysis. Pace Analytical and Eurofins Phoenix are Arizona Department of Health Services (ADHS) certified laboratories (#AZ0728 and #AZ0728, respectively).

The groundwater samples were analyzed for volatile organic compounds (VOCs) via USEPA Method 8260B, 1,4-dioxane via USEPA Method 8260B-SIM, and perchlorate via USEPA Method 314.0 Mod by Pace Analytical. The sample collected from PF-2 was analyzed by Eurofins Phoenix for perchlorate via USEPA Method 314.0.

1.4 Sampling Equipment Decontamination

Disposable sampling equipment such as protective gloves and paper towels were containerized and disposed of as non-hazardous commercial or household waste. Reusable equipment such as the YSI meter and the water level indicator were decontaminated prior to use and between each well using an Alconox and distilled water solution followed by a double rinse with distilled water. Reusable equipment was allowed to air dry prior to its next use.

2. GROUNDWATER MONITORING RESULTS

Laboratory reports and chain-of-custody forms are presented in Attachment 2.

2.1 Estimated Groundwater Flow Direction

The horizontal groundwater gradient was measured across the Site at approximately 0.12 feet per foot (ft/ft) to the west for the plane defined between wells TTU-5, TTU-9A, and TTU-10. This gradient is the same as the 0.12 ft/ft reported during the Q4 2023 groundwater sampling event. Groundwater flow appears to be affected by the significant land surface elevation increase at TTU-15, TTU-16, and TTU-17 creating an area of northern groundwater flow (Figure 2).

2.2 Groundwater Laboratory Results

Perchlorate was detected at concentrations above the Arizona Department of Environmental Quality (ADEQ) Health Based Guidance Level (HBGL) of 14 micrograms per liter (µg/L) in 15 of the 25 wells sampled. 1,4-dioxane was detected at concentrations above the interim screening level of 3.5 µg/L in 14 of the 25 wells sampled. 1,1-dichloroethene (DCE) was detected at concentrations above the Arizona Aquifer Water Quality Standard (AWQS) of 7 µg/L in 10 of the 25 wells sampled. Trichloroethene (TCE) was detected at concentrations above the AWQS of 5 µg/L in 13 of the 25 wells sampled. Analytical results for Q1 2024 are summarized in Table 3 for perchlorate and in Table 4 for detected VOCs. Historical 1,4-Dioxane and TCE concentrations are summarized in Table 5. Concentration contours and VOC exceedances are depicted on Figures 3 through 7.

The Q1 2024 monitoring activities were generally consistent with Q4 2023, but also include the addition of newly established trigger levels (TLs) and follow up sampling procedures as outlined in the *Contingency Plan for Groundwater Contamination Limit Exceedances at Primate Facility Well No. 2* (Geosyntec, 2023). Per the Contingency Plan, notification and resampling must be made if the following TLs are exceeded:

- PF-2: if perchlorate exceeds 3.2 µg/L;
- TTU-6: if 1,4-dioxane exceeds 3.5 µg/L and/or other VOCs reach 50% of the AWQS; and
- TTU-1, TTU-2, TTU-3, TTU-4, TTU-6, TTU-7, TTU-8, TTU-9a, TTU-10, TTU-14, PF-1, and PF-2: if an order-of-magnitude increase in the concentration[1] of a COPC that was previously measured at a concentration exceeding the project screening level (e.g., AWQS).

During the February sampling event, 1,4-dioxane was detected at 5.72 µg/L at TTU-3. Historically since 2018, 1,4-dioxane was not observed at TTU-3 above method detection limits (0.26 ug/L - 0.597 ug/L), with the exception of the previous event where concentrations were observed at 9.22 ug/L. Due to the sample concentrations being an order of magnitude1 higher than the average of the three previous sampling events, TTU-3 was subsequently resampled in March 2024. 1,4-dioxane was not detected in the March resampling, thus additional monthly sampling was not necessary.

During a review of analytical data, the concentrations of COCs at TTU-15 and TTU-16 appeared to be inversed. These wells were resampled to confirm the results in April 2024, and based on the resampling concentrations, it was confirmed the February samples were in fact reported incorrectly. The April resampling results were used for COC plume geometry and concentration versus time plots.

2.3 Groundwater Concentration and Elevation versus Time Plots

Concentration and groundwater elevation versus time plots for TCE, 1,1- DCE, perchlorate, and 1,4-dioxane are presented in Attachment 3. Based on a review of the plots, no groundwater elevations or chemical concentrations were outside of their respective historical ranges. A full analysis of the concentration trends will be provided in the 2024 Annual Groundwater Monitoring Report prepared following the Q4 2024 sampling event.

Prior to the Q3 2023 sampling event, TTU-20 was converted to an extraction well and initially the concentrations of COCs decreased following injection of a carbon donor but have rebounded following the cessation of the pilot test. No other significant changes in concentrations at any other wells were noted during Q1 2024 event relative to historical concentrations.

2.4 Discussion

On February 12, 2024, 1,4-dioxane was detected at TTU-3 at a concentration of 5.72 µg/L, which exceeded the established trigger level. TTU-3 was subsequently resampled for 1,4-dioxane on March 18, 2024, and no detectable concentrations of 1,4-dioxane were observed. Based on the confirmation sample and the trend of the last three sampling events at TTU-3, the non-detection was consistent with historical results and was utilized to determine 1,4-dioxane plume geometry.

2.5 Data Validation

A Tier 1A data validation of the laboratory results according to EPA guidance and the laboratory results are qualified as usable for meeting project objectives. A data validation summary is provided in Attachment 4.

3. CLOSING

Overall, the Q1 2024 groundwater monitoring data indicates primarily stable conditions associated with the Site.

Please contact the undersigned at 720-536-4178 or musson@pinyon-env.com if you have any questions.

Sincerely,
Pinyon Environmental, Inc.

Isabella Foster
Environmental Scientist

Reviewed By:

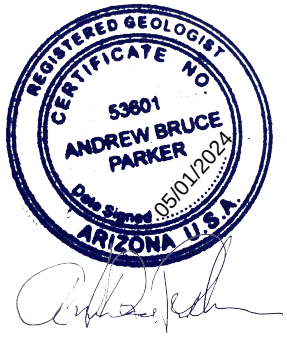

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TABLE 1:
FORMER THERMAL TREATMENT UNIT
GROUNDWATER MONITORING NETWORK
NAMMO DEFENSE SYSTEMS INC.
MESA, ARIZONA

Well ID (Location)	Install Date	Latitude	Longitude	Survey Date	Survey Coordinate Datum	Measuring Point Elevation Top of Casing (ft asml)	Ground Surface Elevation (ft asml)	Well Stickup Height (ft)	ADWR Number	Well Type/Use	Well Name/ Owner	Well Owner Information	Well Const	Well Diameter (in)	Screen Interval (ft bgs)	Casing Depth (ft bgs)	Boring Depth (ft bgs)
Plume Monitoring Wells																	
TTU-3	10/18/2013	33 29 57.98	-111 43 00.91	NP	NAVD 88	1308.03	1305.50	2.50	N/A	Monitoring	Nammo Defense Systems Inc.	P.O. Box 34299 Mesa, AZ 85279	PVC	4	78.1-138.1	143.6	180
TTU-4	10/25/2013	33 30 01.65	-111 42 59.09	NP	NAVD 88	1305.12	1302.50	2.50	N/A	Monitoring	Nammo Defense Systems Inc.	P.O. Box 34299 Mesa, AZ 85280	PVC	4	39.5-99.5	104.9	180
TTU-5	9/20/2014	33 29 52.48	-111 42 58.40	NP	NAVD 88	1314.93	1312.30	3.00	N/A	Monitoring	Nammo Defense Systems Inc.	P.O. Box 34299 Mesa, AZ 85281	PVC	4	59.5-164.5	169.5	174
TTU-6	10/7/2014	33 29 57.57	-111 43 04.79	NP	NAVD 88	1300.84	1299.40	2.37	N/A	Monitoring	Nammo Defense Systems Inc.	P.O. Box 34299 Mesa, AZ 85282	PVC	4	110-175	180	185
TTU-7	10/8/2014	33 29 57.85	-111 43 05.18	NP	NAVD 88	1301.84	1299.30	2.52	N/A	Monitoring	Nammo Defense Systems Inc.	P.O. Box 34299 Mesa, AZ 85282	Steel Open Borehole	8.5 8	282-410	282 None	410
TTU-8	4/18/2016	33 30 01.91	-111 43 05.31	NP	NAVD 88	1310.23	1307.60	2.98	N/A	Monitoring	Nammo Defense Systems Inc.	P.O. Box 34299 Mesa, AZ 85282	PVC	4	135-185	190	204
TTU-9A	6/16/2016	33 30 04.61	-111 42 51.19	NP	NAVD 88	1318.04	1316.00	2.5	N/A	Monitoring	Nammo Defense Systems Inc.	P.O. Box 34299 Mesa, AZ 85282	PVC	4	24-99	104	105
TTU-10	4/18/2016	33 29 54.60	-111 43 07.90	NP	NAVD 88	1302.42	1299.80	3.17	N/A	Monitoring	Nammo Defense Systems Inc.	P.O. Box 34299 Mesa, AZ 85282	PVC	4	115-180	185	204
TTU-12	7/19/2018	33 29 56.03	-111 42 58.38	NP	NP	1312.21	NP	1.33	N/A	Monitoring	Nammo Defense Systems Inc.	P.O. Box 34299 Mesa, AZ 85282	Steel Open Borehole	5.5 5	30-180	30 None	180
TTU-13	7/20/2018	33 29 58.99	-111 42 56.85	NP	NP	1310.79	NP	1.46	N/A	Monitoring	Nammo Defense Systems Inc.	P.O. Box 34299 Mesa, AZ 85283	Steel Open Borehole	5.5 5	30-80	30 None	80
TTU-14	7/19/2018	33 29 57.20	-111 42 57.46	NP	NP	1319.30	1316.80	1.38	N/A	Monitoring	Nammo Defense Systems Inc.	P.O. Box 34299 Mesa, AZ 85284	Steel Open Borehole	5.5 5	45-100	45 None	100
TTU-15	1/25/2018	33 29 56.78	-111 42 47.03	NP	NP	1350.85	NP	1.88	55-228014	Monitoring	Nammo Defense Systems Inc.	P.O. Box 34299 Mesa, AZ 85285	Steel Open Borehole	5 4.5	10-100	10 None	100
TTU-16	1/28/2020	33 29 56.18	-111 42 49.59	NP	NP	1338.55	NP	1.19	55-231730	Monitoring	Nammo Defense Systems Inc.	P.O. Box 34299 Mesa, AZ 85286	Steel Open Borehole	8 8	20-95.6	20 None	95.6
TTU-17	1/28/2020	33 29 58.61	-111 42 45.69	NP	NP	1347.49	NP	0.60	55-231735	Monitoring	Nammo Defense Systems Inc.	P.O. Box 34299 Mesa, AZ 85287	Steel Open Borehole	8 8	20-101	20 None	101
TTU-18	1/25/2020	33 29 47.20	-111 42 58.10	NP	NP	1320.25	NP	NP	55-231737	Monitoring	Nammo Defense Systems Inc.	P.O. Box 34299 Mesa, AZ 85288	Steel Open Borehole	8 8	21-140	21 None	140

TABLE 1:
FORMER THERMAL TREATMENT UNIT
GROUNDWATER MONITORING NETWORK
NAMMO DEFENSE SYSTEMS INC.
MESA, ARIZONA

Well ID (Location)	Install Date	Latitude	Longitude	Survey Date	Survey Coordinate Datum	Measuring Point Elevation Top of Casing (ft asml)	Ground Surface Elevation (ft asml)	Well Stickup Height (ft)	ADWR Number	Well Type/Use	Well Name/ Owner	Well Owner Information	Well Const	Well Diameter (in)	Screen Interval (ft bgs)	Casing Depth (ft bgs)	Boring Depth (ft bgs)
Extraction and Injection Wells																	
TTU-1	6/6/2012	33 29 59.14	-111 42 56.27	NP	NAVD 88	1312.73	1309.70	3.03	55-914440	Extraction	Nammo Defense Systems Inc.	P.O. Box 34299 Mesa, AZ 85277	PVC	4	30-70	75	200
TTU-2	10/17/2013	33 29 55.85	-111 42 57.85	NP	NAVD 88	1314.44	1311.80	2.64	N/A	Extraction	Nammo Defense Systems Inc.	P.O. Box 34299 Mesa, AZ 85278	PVC	4	49.4-179.6	185	187.5
TTU-11	9/11/2015	33 29 55.28	-111 42 51.47	NP	NAVD 88	1339.20	1336.60	2.60	55-918534	Extraction/ Injection [1]	Nammo Defense Systems Inc.	P.O. Box 34299 Mesa, AZ 85282	PVC	4	24.1-89.1	94	136
TTU-20	9/24/2020	33 29 55.17	-111 42 51.58	NP	NP	1336.90	NP	0.85	55-232968	Extraction [4]	Nammo Defense Systems Inc.	P.O. Box 34299 Mesa, AZ 85288	PVC	4	25-95	95	100
TTU-19	9/24/2020	33 29 55.25	-111 42 51.50	NP	NP	1336.67	NP	NP	55-232969	Monitoring/ Injection [2]	Nammo Defense Systems Inc.	P.O. Box 34299 Mesa, AZ 85288	PVC	4	25-90	95	96
TTU-EX-1	1/25/2020	33 29 58.42	-111 42 52.55	NP	NP	1321.69	NP	1.60	55-231733	Extraction	Nammo Defense Systems Inc.	P.O. Box 34299 Mesa, AZ 85288	Steel	8	20-110.7	20	110.7
													Open Borehole	8		None	
TTU-EX-2	1/23/2020	33 29 57.61	-111 42 53.79	NP	NP	1316.40	NP	1.10	55-231734	Extraction	Nammo Defense Systems Inc.	P.O. Box 34299 Mesa, AZ 85289	Steel	8	20-110	20	110
													Open Borehole	8		None	
TTU-EX-3	1/24/2020	33 29 56.29	-111 42 54.12	NP	NP	1316.85	NP	0.58	55-231731	Extraction	Nammo Defense Systems Inc.	P.O. Box 34299 Mesa, AZ 85290	Steel	8	20-101.45	20	111
													Open Borehole	8		None	
TTU-EX-4	1/24/2020	33 29 55.46	-111 42 54.39	NP	NP	1319.96	NP	1.42	55-231732	Extraction	Nammo Defense Systems Inc.	P.O. Box 34299 Mesa, AZ 85291	Steel	8	20-110.7	20	110.7
													Open Borehole	8		None	
TTU-EX-5	1/24/2020	33 29 54.68	-111 42 54.62	NP	NP	1319.50	NP	0.96	55-231736	Extraction	Nammo Defense Systems Inc.	P.O. Box 34299 Mesa, AZ 85292	Steel	8	20-110.8	20	110.8
													Open Borehole	8		None	
Production Wells																	
PF-1	NP	33 29 56.60	-111 43 09.75	NP	NP	1295.99	NP	NP	N/A	Production	University of Washington	4202 N Higley Rd Mesa, AZ 85215	Unknown	Unknown	Unknown	Unknown	Unknown
PF-2	3/27/2013	33 29 56.65	-111 43 09.96	NP	NP	1296.35	NP	NP	N/A	Production	University of Washington	4202 N Higley Rd Mesa, AZ 85215	Steel	6 5/8	300-400	400	400

Notes:

ft amsl = feet above mean sea level (NAVD88)
ADWR = Arizona Department of Water Resources
Const = construction
in = inches
N/A = Not applicable
PVC = polyvinyl chloride
ft bgs = feet below ground surface
TTU = Thermal Treatment Unit
EX = Extraction
PF = Primate Facility
NP = Not Provided
Drill Log TOC Different from Original
Drill Log TOC listed

(1) - TTU-11 was converted from an extraction well to an injection well in October 2020 for a In-Situ Bioremediation Pilot Test.
(2) - TTU-19 was converted from a monitoring well to an injection well in February 2021 for an In-Situ Bioremediation Pilot Test
(3) - Monitoring well stick-up was measured using a tape measure to the top of the protective casing and not to the top of the well casing as no survey equipment was utilized.
(4) - TTU-20 was converted from a monioring well to an extraction well in INSERTHERE 2023 for
TTU-EX-1 through TTU-EX-5 are not currently operating as extraction wells. TTU-11 and TTU-19 are not currently operating as injection wells.

TABLE 2:
GROUNDWATER ELEVATIONS - FIRST QUARTER 2024
FORMER THERMAL TREATMENT UNIT
NAMMO DEFENSE SYSTEMS INC.

Location	Northing (intl ft)	Easting (intl ft)	Top of Casing Elevation (ft asml)	Date Measured	Depth to Water (ft btoc)	Groundwater Elevation (ft asml)
TTU-1	909420.734	761281.203	1312.73	2/13/2024	45.63	1,267.10
TTU-2	909087.852	761148.265	1314.44	2/13/2024	64.01	1,250.43
TTU-3	909303.363	760888.204	1308.03	2/12/2024	88.82	1,219.21
	909303.363	760888.204	1308.03	3/18/2024	89.07	1,218.96
TTU-4	909673.680	761041.975	1305.12	2/13/2024	52.93	1,252.19
TTU-5	908747.636	761102.227	1314.93	2/13/2024	82.77	1,232.16
TTU-6	909260.820	760560.096	1300.84	2/12/2024	129.01	1,171.83
TTU-7	909287.611	760527.269	1301.84	2/12/2024	135.50	1,166.34
TTU-8	909699.266	760514.908	1310.23	2/14/2024	144.25	1,165.98
TTU-9A	909974.490	761710.151	1318.04	2/13/2024	66.96	1,251.08
TTU-10	908960.114	760297.013	1302.42	2/14/2024	166.62	1,135.80
TTU-11	909029.758	761706.470	1339.20	2/13/2024	36.84	1,302.36
TTU-12	909105.990	761103.280	1312.21	2/12/2024	74.17	1,238.04
TTU-13	909405.920	761232.180	1310.79	2/12/2024	43.70	1,267.09
TTU-14	909224.260	761181.230	1316.80	2/12/2024	60.39	1,256.41
TTU-15	909185.100	762065.910	1350.85	2/13/2024	33.82	1,317.03
TTU-16	909124.980	761848.851	1338.55	2/13/2024	26.24	1,312.31
TTU-17	909370.903	762179.168	1347.49	2/12/2024	41.19	1,306.30
TTU-18	908215.829	761130.011	1320.25		DRY	
TTU-19	909030.750	761687.700	1336.81	2/13/2024	34.87	1,301.94
TTU-20	909022.530	761681.990	1336.90	2/13/2024	38.06	1,298.84
TTU-EX-1	909350.574	761597.823	1321.69	2/13/2024	30.80	1,290.89
TTU-EX-2	909268.187	761493.214	1316.40	2/13/2024	39.87	1,276.53
TTU-EX-3	909134.941	761465.507	1316.85	2/12/2024	41.78	1,275.07
TTU-EX-4	909051.298	761442.876	1319.96	2/12/2024	44.74	1,275.22
TTU-EX-5	908971.770	761423.325	1319.50	2/13/2024	41.22	1,278.28
PF-1	909161.578	760140.434	1295.99	2/14/2024	156.10	1,139.89
PF-2	909166.890	760122.250	1296.35	2/14/2024	NM*	--

Notes:

intl ft - international foot

ft amsl - feet above mean sea level

ft btoc - feet below top of casing

"--" - Not enough information to determine groundwater elevation

NM - not measured

** - DTW not obtained due to obstruction in well*

TABLE 3:

SUMMARY OF PERCHLORATE CONCENTRATIONS - FIRST QUARTER 2024

FORMER THERMAL TREATMENT UNIT

NAMMO DEFENSE SYSTEMS INC.

			Analyte	Perchlorate	
			EPA Method	314	314.0
			Units	µg/l	
			HBGL	14	
Location	Sample Depth (ft btoc)	Sample Date	Sample Type	Concentration	
PF-2	400	2/14/2024	Primary	--	<1.0 E8
TTU-1	50	2/13/2024	Primary	**11,700**	--
TTU-2	114	2/13/2024	Primary	**150,000**	--
TTU-3	108	2/12/2024	Primary	**54.0**	--
TTU-4	57	2/13/2024	Primary	<4.00	--
DUP-01			Duplicate	<4.00 M2	--
TTU-5	110	2/13/2024	Primary	**37.7**	--
TTU-6	143	2/12/2024	Primary	**18.1**	--
TTU-7	345	2/12/2024	Primary	<4.00	--
TTU-8	164	2/14/2024	Primary	<4.00	--
TTU-9A	61	2/13/2024	Primary	6.06	--
TTU-10	172	2/14/2024	Primary	<4.00	--
DUP-03			Duplicate	<4.00	--
TTU-11	73	2/13/2024	Primary	<4.00	--
TTU-12	82	2/12/2024	Primary	**139,000**	--
TTU-13	51	2/12/2024	Primary	**24,000**	--
TTU-14	64	2/12/2024	Primary	**147,000**	--
TTU-15	75	4/9/2024	Primary	**11,100**	--
TTU-16	80	4/9/2024	Primary	**817,000**	--
TTU-17	80	2/12/2024	Primary	<4.00	--
TTU-19	73	2/13/2024	Primary	<4.00	--
TTU-20	73	2/13/2024	Primary	**85,900**	--
TTU-EX-1	69	2/13/2024	Primary	**90,000**	--
TTU-EX-2	74	2/13/2024	Primary	**93,200**	--
TTU-EX-3	76	2/12/2024	Primary	**468,000**	--
TTU-EX-4	77	2/12/2024	Primary	**85,600**	--
TTU-EX-5	80	2/13/2024	Primary	<4.00	--
DUP-02			Duplicate	<4.00	--

Notes:

ft btoc - feet below top of casing

µg/l - micrograms per liter

EPA - United States Environmental Protection Agency

HBGL - Health-Based Guidance Level

<Grey - Concentration is below laboratory reporting limits

** - The concentration limit for PF-2 is 6.4 µg/l and the trigger level is 3.2 µg/l*

-- - Not reported

***BOLD** - Concentration exceeds its respective HGBL*

E8 - Analyte reported to MDL per project specification. Target analyte was not detected in the sample.

M2 - Matrix spike recovery was low; the associated blank spike recovery was acceptable.

TABLE 4:
SUMMARY OF DETECTED VOC CONCENTRATIONS - FIRST QUARTER 2024
FORMER THERMAL TREATMENT UNIT
NAMMO DEFENSE SYSTEMS INC.

Location	Sample Depth (ft btoc)	Chemical Name	1,4-Dioxane	Acetone	1,1-dichloroethane	1,1-dichloroethene	1,2-dichloroethane	Benzene	Carbon Disulfide	Chloroform	cis-1,2-dichloroethene	Dichloromethane (methylene chloride)	Isopropylbenzene	Tetrachloroethene	Toluene	trans-1,2-dichloroethene	1,1,2-trichloroethane	Trichloroethene	Vinyl chloride	2-butanone (MEK)	4-methyl-2-pentanone (MIBK)	Xylene Total
		EPA Method	8260B SIM	8260B																		
		Unit	µg/l																			
		Screening Level	3.5(1)	1800(2)	2.8(2)	7	5	5	81(2)	80(3)	70	5	45(2)	5	1000	100	5	5	2	560(2)	630(2)	10000
		Sample Date																				
PF-2	400	2/14/2024	<0.597	<11.3	<0.100	<0.188	<0.0819	<0.0941	0.172 B1;E4	<0.111	<0.126	<0.430	<0.105	<0.3	<0.278	<0.149	<0.158	<0.190	<0.234	<1.19	<0.478 MI	<0.174
TTU-1	50	2/13/2024	**18.2**	<11.3	<0.100	1.21	<0.0819	<0.0941	<0.0962	<0.111	<0.126	<0.430	<0.105	<0.300	<0.278	<0.149	<0.158	**5.44**	<0.234	<1.19	<0.478	<0.174
TTU-2	114	2/13/2024	**207**	<113	<1.00	**31.6**	<0.819	1.14 E4	<0.962	1.65 E4	<1.26	<4.30	<1.05	<3.00	<2.78	<1.49	<1.58	**529**	<2.34	<11.9	<4.78	<1.74
TTU-3	108	2/12/2024	**5.72**	<11.3	<0.100	<0.188	<0.0819	<0.0941	<0.0962	<0.111	<0.126	<0.430	<0.105	<0.300	<0.278	<0.149	<0.158	<0.190	<0.234	<1.19	<0.478	<0.174
TTU-3 *		3/18/2024	<0.597	NS	NS	NS	NS	NS	NS	NS	NS	NS	NS	NS	NS	NS	NS	NS	NS	NS	NS	NS
TTU-4	57	2/13/2024	<0.597	<11.3	<0.100	<0.188	<0.0819	<0.0941	<0.0962	<0.111	<0.126	<0.430	<0.105	<0.300	<0.278	<0.149	<0.158	<0.190	<0.234	<1.19	<0.478	<0.174
DUP-01			<0.597	<11.3	<0.100	<0.188	<0.0819	<0.0941	<0.0962	<0.111	<0.126	<0.430	<0.105	<0.300	<0.278	<0.149	<0.158	<0.190	<0.234	<1.19	<0.478	<0.174
TTU-5	110	2/13/2024	<0.597	<11.3	<0.100	<0.188	<0.0819	<0.0941	<0.0962	<0.111	<0.126	<0.430	<0.105	<0.300	<0.278	<0.149	<0.158	<0.190	<0.234	<1.19	<0.478	<0.174
TTU-6	143	2/12/2024	<0.597	<11.3	<0.100	<0.188	<0.0819	<0.0941	<0.0962	<0.111	<0.126	<0.430	<0.105	<0.300	<0.278	<0.149	<0.158	<0.190	<0.234	<1.19	<0.478	<0.174
TTU-7	345	2/12/2024	<0.597	<11.3	<0.100	<0.188	<0.0819	0.113 E4	<0.0962	<0.111	<0.126	<0.430	<0.105	<0.300	0.540 E4	<0.149	<0.158	<0.190	<0.234	<1.19	<0.478	0.185 E4
TTU-8	164	2/14/2024	<0.597	<11.3	<0.100	<0.188	<0.0819	<0.0941	0.186 B1;E4;R7	<0.111	<0.126	<0.430	<0.105	<0.300	<0.278	<0.149	<0.158	<0.190	<0.234	<1.19	<0.478	<0.174
TTU-9A	61	2/13/2024	<0.597	<11.3	<0.100	<0.188	<0.0819	<0.0941	<0.0962	<0.111	<0.126	<0.430	<0.105	<0.300	<0.278	<0.149	<0.158	<0.190	<0.234	<1.19	<0.478	<0.174
TTU-10	172	2/14/2024	<0.597	<11.3	<0.100	<0.188	<0.0819	<0.0941	0.17 B1;E4;R7	<0.111	<0.126	<0.430	<0.105	<0.300	<0.278	<0.149	<0.158	<0.190	<0.234	<1.19	<0.478	<0.174
DUP-03			<0.597	<11.3	<0.100	<0.188	<0.0819	<0.0941	<0.0962	<0.111	<0.126	<0.430	<0.105	<0.300	<0.278	<0.149	<0.158	<0.190	<0.234	<1.19	<0.478	<0.174
TTU-11	73	2/13/2024	<0.597	**2320**	<1.00	5.47 E4	<0.819	<0.941	<0.962	<1.11	28.8	4.98 E4	3.71 E4	<3.00	<2.78	<1.49	<1.58	**76.6**	<2.34	**1260**	183	<1.74
TTU-12	82	2/12/2024	<0.597	<113	<1.00	**52.6**	<0.819	1.02 E4	<0.962	1.55 E4	<1.26	<4.30	<1.05	<3.00	<2.78	<1.49	<1.58	**511**	<2.34	<11.9	<4.78	<1.74
TTU-13	51	2/12/2024	**31.4**	<11.3	<0.100	4.80	<0.0819	<0.0941	<0.0962	<0.111	<0.126	<0.430	<0.105	<0.300	<0.278	<0.149	<0.158	**13.6**	<0.234	<1.19	<0.478	<0.174
TTU-14	64	2/12/2024	**286**	<113	<1.00	**81.7**	<0.819	1.84 E4	<0.962	2.01 E4	2.28 E4	<4.30	<1.05	<3.00	<2.78	<1.49	2.31 E4	**829**	<2.34	<11.9	<4.78	<1.74
TTU-15	75	4/9/2024	NS	<226	<2	<3.76	<1.64	<1.88	3.98 B1;E4	<2.22	<2.52	<8.60	<2.10	<6.00	<5.56	<2.98	<3.16	<3.8	<4.68	<23.8	<9.56	<3.48
TTU-16	80	4/9/2024	NS	<28200	<250	**2440 E4**	<205	**260 E4**	**494 B1;E4**	<278	<315	**52100**	<263	<750	<695	<373	<395	**68500**	<585	<2980	<1200	<435
TTU-17	80	2/12/2024	<0.597	<11.3	<0.100	<0.188	<0.0819	<0.0941	<0.0962	<0.111	0.507 E4	<0.430	<0.105	<0.300	<0.278	<0.149	<0.158	0.661 E4	<0.234	<1.19	<0.478	<0.174
TTU-19	73	2/13/2024	**192**	<56.5	<0.500	**17.2**	<0.409	3.46 E4	0.496 E4	<0.555	**155**	**5.44 E4**	1.83 E4	<1.50	<1.39	5.90	<0.790	**195**	**3.99 E4**	<5.95	3.61 E4	<0.870
TTU-20	73	2/13/2024	**378**	<113	**3.36 E4**	**154**	<0.819	4.48 E4	<0.962	3.86 E4	18.9	**8.89 E4**	<1.05	<3.00	<2.78	<1.49	3.28 E4	**1860**	<2.34	<11.9	<4.78	<1.74
TTU-EX-1	69	2/13/2024	**326**	<113	<1.00	**88.9**	<0.819	<0.941	<0.962	1.33 E4	<1.26	<4.30	<1.05	<3.00	<2.78	<1.49	<1.58	**241**	<2.34	<11.9	<4.78	<1.74
TTU-EX-2	74	2/13/2024	**256**	<226	<2.00	**83.3**	<1.64	<1.88	<1.92	<2.22	<2.52	<8.60	<2.10	<6.00	<5.56	<2.98	<3.16	**415**	<4.68	<23.8	<9.56	<3.48
TTU-EX-3	76	2/12/2024	**523**	<2820	<25.0	**836**	<20.5	<23.5	<24.1	<27.8	<31.5	**160 E4**	<26.3	<75.0	<69.5	<37.3	<39.5	**5930**	<58.5	<298	<120	<43.5
TTU-EX-4	77	2/12/2024	**15.9**	<282	<2.50	**91.1**	<2.05	<2.35	<2.41	<2.78	<3.15	**16.1 E4**	<2.63	<7.50	<6.95	<3.73	<3.95	**588**	<5.85	<29.8	<12.0	<4.35
TTU-EX-5	80	2/13/2024	**9.87**	<11.3	<0.100	<0.188	<0.0819	<0.0941	<0.0962	<0.111	0.234 E4	<0.430	<0.105	<0.300	<0.278	<0.149	<0.158	3.57	<0.234	<1.19	<0.478	<0.174
DUP-02			<0.597	<11.3	<0.100	<0.188	<0.0819	<0.0941	<0.0962	<0.111	<0.126	<0.430	<0.105	<0.300	<0.278	<0.149	<0.158	3.08	<0.234	<1.19	<0.478	<0.174
			3.50	**1800.00**	**2.80**	**7.00**	**5.00**	**5.00**	**81.00**	**80.00**	**70.00**	**5.00**	**45.00**	**5.00**	**1000.00**	**100.00**	**5.00**	**5.00**	**2.00**	**560.00**	**630.00**	**10000.00**

Notes:
ft btoc - feet below top of casing
µg/l - micrograms per liter
AWQS - Arizona Aquifer Water Quality Standard
Detected VOCs with established screening value presented
EPA - Environmental Protection Agency
SIM - Selected Ion Monitoring
<Gray - Concentration is below laboratory reporting limits
BOLD - Concentration exceeds its respective AWQS or screening level
NS - Not sampled

B1 - Target analyte detected in method blank at or above the method reporting limit.
E4 - Concentration estimated. Analyte was detected below laboratory minimum reporting level (MRL) but above MDL.
R7 - LFB/LFBD RPD exceeded the laboratory acceptance limit. Recovery met acceptance criteria.
(1) - Interim Screening Level
(2) - Tap water regional screening level, AWQS has not been established
(3) - The total trihalomethane (TTHM) standard is exceeded when the sum of the four compounds exceeds 80 µg/L as a rolling average
PF-2, TTU-6, and TTU-7 have a Concentration Limit at the AWQS and a Trigger Level at 1/2 the AWQS
(*) - Confirmation sample per Contigency Plan 2023

TABLE 5:
HISTORICAL 1,4-DIOXANE AND TCE CONCENTRATIONS
FORMER THERMAL TREATMENT UNIT
NAMMO DEFENSE SYSTEMS INC.

		Chemical Name	1,4-Dioxane	Trichloroethene
		EPA Method	8260B SIM	8260B
		Unit	µg/l	
Location	Sample Type	Screening Level	3.5[1]	5
		Sample Date		
Contigency Level/ Trigger Level at PF-1, PF-2, TTU-6, and TTU-7			3.5/1.75	5/2.5
PF-1	Primary	3/27/2018	<0.36 U	<0.18 U
	Primary	6/28/2018	<0.72 U	<0.15 U
	Primary	9/10/2018	<0.26 U	<0.15 U
	Primary	12/10/2018	<0.26 U	<0.15 U
	Primary	3/26/2019	<0.597 U	<0.398 U
PF-2	Primary	3/27/2018	<0.36 U	<0.18 U
	Primary	6/28/2018	<0.72 U	<0.15 U
	Primary	9/10/2018	<0.26 U	<0.15 U
	Primary	12/10/2018	<0.26 U	<0.15 U
	Primary	3/26/2019	<0.597 U	<0.398 U
	Primary	9/16/2019	<0.597 U	<0.398 U
	Field Duplicate	9/16/2019	<0.597 U	<0.398 U
	Primary	12/23/2019	<0.597 U	<0.398 U
	Primary	3/13/2020	<0.597 U	<0.398 U
	Primary	12/4/2020	<0.597 U	<0.190 U
	Field Duplicate	12/4/2020	<0.597 U	<0.190 U
	Primary	3/29/2021	<0.597 U	<0.190 U
	Primary	5/6/2021	<0.597 U	<0.190 U
	Primary	8/6/2021	<0.597 U	<0.190 U
	Primary	11/18/2021	<0.597 U	<0.190 U
	Primary	3/31/2022	<0.597	<0.190 R7
	Field Duplicate	3/31/2022	<0.597	<0.190 R7
	Primary	6/21/2022	<0.597	<0.190 J3
	Field Duplicate	6/21/2022	<0.597	<0.190
	Primary	9/9/2022	<0.597 R7	<0.190
	Primary	11/30/2022	<0.597	<0.190
	Primary	2/27/2023	<0.597	<0.190
	Primary	6/12/2023	<0.597	<0.190
	Field Duplicate	6/12/2023	<0.597	<0.190 L1
	Primary	9/5/2023	<0.597	<0.190
	Primary	11/21/2023	<0.597	<0.190 L1
	Primary	2/14/2024	<0.597	<0.190
TTU-1	Primary	11/18/2014	NA	6.1
	Primary	12/23/2014	NA	8.8
	Primary	2/5/2015	26	10
	Primary	5/18/2015	20	6.1
	Primary	9/9/2015	17	5.2
	Primary	11/23/2015	14	5.1
	Primary	2/25/2016	11	4.6
	Primary	6/1/2016	12.7	3.03
	Primary	8/18/2016	11	3.7
	Primary	11/22/2016	27	5.5
	Primary	2/22/2017	18.4	5.5
	Primary	5/23/2017	14.1	7.2
	Primary	8/29/2017	11	1.4
	Primary	11/27/2017	17.7	7.1
	Field Duplicate	11/27/2017	18.1	7.2
	Primary	3/27/2018	17.1	4.6
	Primary	9/12/2018	31.8	11.2
	Field Duplicate	9/12/2018	29.1	12.4
	Primary	12/4/2018	7.3	4.4
	Primary	9/16/2019	13.9	5.72
	Field Duplicate	9/16/2019	10.8	4.85
	Primary	12/20/2019	5.06	5.19
	Primary	3/12/2020	4.63 J	3.91
	Primary	6/18/2020	17.1	7.6
	Primary	7/20/2020	3.71	6.09
	Primary	12/2/2020	29.9	1.33
	Primary	3/30/2021	18.9 J	6.40
	Primary	5/6/2021	22.0	17.1 J
	Primary	7/29/2021	37.7	14.3
	Primary	12/22/2021	11.1	8.82
	Primary	3/26/2022	18.4	3.72
	Field Duplicate	3/26/2022	19.9	4.46
	Primary	6/16/2022	17.5 Q	4.42
	Field Duplicate	6/16/2022	35.5	4.12
	Primary	10/11/2022	15.1	5.13
	Field Duplicate	10/11/2022	14.5	5.85
	Primary	11/28/2022	11.8 B	4.86
	Primary	2/23/2023	14.9	5.02 M1
	Primary	6/15/2023	9.66	4.18
	Primary	9/14/2023	6.07	1.38
	Primary	11/17/2023	9.45	3.51
	Primary	2/13/2024	18.2	5.44

TABLE 5:
HISTORICAL 1,4-DIOXANE AND TCE CONCENTRATIONS
FORMER THERMAL TREATMENT UNIT
NAMMO DEFENSE SYSTEMS INC.

		Chemical Name	1,4-Dioxane	Trichloroethene
		EPA Method	8260B SIM	8260B
		Unit	µg/l	
Location	Sample Type	Screening Level	3.5[(1)]	5
		Sample Date		
TTU-2	Primary	11/18/2014	NA	370
	Primary	12/23/2014	NA	280
	Primary	2/5/2015	170	280
	Primary	5/18/2015	160	190
	Primary	9/9/2015	170	200
	Primary	11/23/2015	140	150
	Primary	2/25/2016	110	150
	Primary	6/1/2016	88.2	50.3
	Primary	8/18/2016	150	360
	Primary	11/22/2016	260	780
	Primary	2/22/2017	244	727
	Primary	5/23/2017	222	880
	Primary	8/29/2017	241	93.2
	Field Duplicate	8/29/2017	227	89.7
	Primary	11/27/2017	235	353
	Primary	3/27/2018	219	236
	Field Duplicate	3/27/2018	152	274
	Primary	6/28/2018	246	498
	Primary	9/10/2018	246	433
	Primary	12/4/2018	232	288
	Primary	3/25/2019	313	364
	Primary	9/16/2019	295	475
	Primary	12/20/2019	211	711
	Field Duplicate	12/20/2019	215	742
	Primary	3/12/2020	227 J	511
	Primary	6/18/2020	292	824
	Primary	7/20/2020	156	959
	Primary	12/2/2020	329	785
	Primary	3/30/2021	196 J	656
	Field Duplicate	3/30/2021	244 J	720
	Primary	5/6/2021	316	683
	Primary	7/29/2021	373	654
	Primary	12/22/2021	280	627
	Field Duplicate	12/22/2021	281	653
	Primary	3/26/2022	251	823
	Primary	6/16/2022	246 Q	443
	Primary	10/10/2022	170	596 M3
	Primary	11/28/2022	230 V	643 V
	Primary	2/23/2023	362	648
	Primary	6/15/2023	239	406
	Field Duplicate	6/15/2023	269	429
	Primary	9/14/2023	200	501
	Field Duplicate	9/14/2023	234	566
	Primary	11/17/2023	221	561
	Primary	2/13/2024	207	529
TTU-3	Primary	5/23/2017	NA	2.5
	Primary	3/27/2018	<0.36 U	<0.18 U
	Primary	6/28/2018	<0.72 U	<0.15 U
	Primary	9/10/2018	<0.26 U	<0.15 U
	Primary	12/10/2018	<0.26 U	<0.15 U
	Primary	3/26/2019	<0.597 U	<0.398 U
	Primary	6/7/2019	<0.597 U	<0.398 U
	Primary	9/16/2019	<0.597 U	<0.398 U
	Primary	12/23/2019	<0.597 U	<0.398 U
	Primary	3/13/2020	<0.597 U	<0.398 U
	Primary	6/18/2020	<0.597 U	<0.190 U
	Primary	7/21/2020	<0.597 U	<0.190 U
	Primary	12/4/2020	<0.597 U	<0.190 U
	Primary	3/29/2021	<0.597 U	<0.190 U
	Primary	5/6/2021	<0.597 U	<0.190 U
	Field Duplicate	5/6/2021	<0.597 U	<0.190 U
	Primary	7/30/2021	<0.597 U	<0.190 U
	Primary	11/18/2021	<0.597 U	<0.190 U
	Primary	3/22/2022	<0.597	0.454 E4
	Primary	6/14/2022	<0.597 J3	<0.190 J3
	Primary	9/9/2022	<0.597	<0.190
	Primary	11/30/2022	<0.597 J3	<0.190
	Field Duplicate	11/30/2022	<0.597	<0.190
	Primary	2/25/2023	<0.597	0.266 E4
	Field Duplicate	2/25/2023	<0.597	<0.190
	Primary	6/12/2023	<0.597	<0.190
	Primary	9/5/2023	<0.597	<0.190
	Primary	11/17/2023	9.22	<0.190
	Primary	12/13/2023	<0.597	NS
	Primary	2/12/2024	5.72	<0.190

TABLE 5:
HISTORICAL 1,4-DIOXANE AND TCE CONCENTRATIONS
FORMER THERMAL TREATMENT UNIT
NAMMO DEFENSE SYSTEMS INC.

		Chemical Name	1,4-Dioxane	Trichloroethene
		EPA Method	8260B SIM	8260B
		Unit	µg/l	
Location	Sample Type	Screening Level	3.5[(1)]	5
		Sample Date		
TTU-4	Primary	5/23/2017	NA	0.31
	Primary	3/27/2018	<0.36 U	<0.18 U
	Field Duplicate	3/27/2018	<0.36 U	<0.18 U
	Primary	6/28/2018	<0.72 U	<0.15 U
	Field Duplicate	6/28/2018	<0.72 U	<0.15 U
	Primary	9/10/2018	<0.26 U	<0.15 U
	Field Duplicate	9/10/2018	<0.26 U	<0.15 U
	Primary	12/10/2018	<0.26 U	<0.15 U
	Field Duplicate	12/10/2018	<0.26 U	<0.15 U
	Primary	3/26/2019	<0.597 U	<0.398 U
	Primary	6/7/2019	<0.597 U	<0.398 U
	Primary	9/16/2019	<0.597 U	<0.398 U
	Primary	12/23/2019	<0.597 U	<0.398 U
	Primary	3/13/2020	<0.597 U	<0.398 U
	Primary	6/18/2020	<0.597 U	<0.190 U
	Primary	7/21/2020	<0.597 U	<0.190 U
	Field Duplicate	7/21/2020	<0.597 U	<0.190 U
	Primary	12/4/2020	<0.597 U	<0.190 U
	Primary	3/29/2021	<0.597 U	<0.190 U
	Primary	5/6/2021	<0.597 U	<0.190 U
	Field Duplicate	5/6/2021	<0.597 U	<0.190 U
	Primary	7/30/2021	<0.597 U	<0.190 U
	Field Duplicate	7/30/2021	<0.597 U	<0.190 U
	Primary	11/18/2021	<0.597 U	<0.190 U
	Primary	3/22/2022	<0.597	<0.190
	Field Duplicate	3/22/2022	2.59 E4	<0.190
	Primary	6/14/2022	11.1	<0.190J3
	Primary	7/21/2022	<0.597	<0.190
	Field Duplicate	7/21/2022	<0.597	<0.190
	Primary	9/9/2022	<0.597	<0.190
	Primary	11/30/2022	1.84 J	<0.190
	Primary	2/25/2023	<0.597	<0.190
	Primary	6/12/2023	<0.597	<0.190
	Primary	9/5/2023	<0.597	<0.190
	Primary	11/17/2023	<0.597	<0.190
	Primary	2/13/2024	<0.597	<0.190
	Field Duplicate	2/13/2024	<0.597	<0.190
TTU-5	Primary	3/27/2018	<0.36 U	<0.18 U
	Primary	6/28/2018	<0.72 U	<0.15 U
	Primary	9/10/2018	<0.26 U	<0.15 U
	Primary	12/10/2018	<0.26 U	<0.15 U
	Primary	3/26/2019	<0.597 U	<0.398 U
	Primary	6/7/2019	<0.597 U	<0.398 U
	Primary	9/16/2019	<0.597 U	<0.398 U
	Primary	12/20/2019	3.54	<0.398 U
	Primary	3/12/2020	<0.597 U	<0.398 U
	Primary	6/17/2020	<0.597 U	<0.190 U
	Primary	7/20/2020	<0.597 U	<0.190 U
	Primary	12/2/2020	<0.597 U	0.877 J
	Primary	3/30/2021	<0.597 U	<0.190 U
	Primary	5/6/2021	<0.597 U	<0.190 U
	Primary	7/29/2021	<0.597 U	<0.190 U
	Primary	11/17/2021	<0.597 U	<0.190 U
	Primary	3/21/2022	<0.597	0.64 E4
	Primary	6/13/2022	130	<0.190
	Primary	7/21/2022	<0.597	<0.190
	Primary	9/8/2022	<0.597	<0.190
	Primary	11/29/2022	21.5	<0.190
	Primary	2/25/2023	<0.597	<0.190
	Primary	6/13/2023	<0.597	<0.190
	Primary	9/6/2023	<0.597	<0.190
	Primary	11/17/2023	<0.597	<0.190
	Field Duplicate	11/17/2023	<0.597	<0.190
	Primary	2/13/2024	<0.597	<0.190

TABLE 5:
HISTORICAL 1,4-DIOXANE AND TCE CONCENTRATIONS
FORMER THERMAL TREATMENT UNIT
NAMMO DEFENSE SYSTEMS INC.

		Chemical Name	1,4-Dioxane	Trichloroethene
		EPA Method	8260B SIM	8260B
		Unit	µg/l	
Location	Sample Type	Screening Level	3.5[(1)]	5
		Sample Date		
TTU-6	Primary	8/29/2017	NA	0.38
	Primary	3/27/2018	<0.36 U	<0.18 U
	Primary	6/28/2018	<0.72 U	<0.15 U
	Primary	9/10/2018	<0.26 U	<0.15 U
	Primary	12/10/2018	<0.26 U	<0.15 U
	Primary	3/26/2019	<0.597 U	<0.398 U
	Primary	6/7/2019	<0.597 U	<0.398 U
	Field Duplicate	6/7/2019	<0.597 U	<0.398 U
	Primary	9/16/2019	<0.597 U	<0.398 U
	Primary	12/23/2019	<0.597 U	<0.398 U
	Primary	3/13/2020	<0.597 U	<0.398 U
	Primary	6/18/2020	<0.597 U	<0.190 U
	Primary	7/21/2020	<0.597 U	<0.190 U
	Primary	12/4/2020	<0.597 U	<0.190 U
	Primary	3/29/2021	<0.597 U	<0.190 U
	Primary	5/6/2021	<0.597 U	<0.190 U
	Primary	7/30/2021	<0.597 U	<0.190 U
	Primary	11/18/2021	<0.597 U	<0.190 U
	Primary	3/22/2022	<0.597	<0.190
	Primary	6/14/2022	<0.597 J3	<0.190
	Primary	9/9/2022	<0.597 R5	<0.190
	Primary	11/30/2022	<0.597	<0.190
	Primary	2/25/2023	<0.597	0.218 E4
	Primary	6/12/2023	<0.597	<0.190
	Primary	9/5/2023	<0.597	<0.190
	Primary	11/17/2023	<0.597	<0.190
	Primary	2/12/2024	<0.597	<0.190
TTU-7	Primary	3/27/2018	<0.36 U	<0.18 U
	Primary	6/28/2018	<0.72 U	<0.15 U
	Primary	9/10/2018	<0.26 U	<0.15 U
	Primary	12/10/2018	<0.26 U	<0.15 U
	Primary	3/26/2019	<0.597 U	<0.398 U
	Field Duplicate	3/26/2019	<0.597 U	<0.398 U
	Primary	6/7/2019	<0.597 U	<0.398 U
	Primary	9/16/2019	<0.597 U	<0.398 U
	Primary	12/23/2019	<0.597 U	<0.398 U
	Primary	3/13/2020	<0.597 U	<0.398 U
	Primary	6/18/2020	<0.597 U	<0.190 U
	Primary	7/21/2020	<0.597 U	<0.190 U
	Primary	12/4/2020	<0.597 U	<0.190 U
	Primary	3/29/2021	<0.597 U	<0.190 U
	Primary	5/6/2021	<0.597 U	<0.190 U
	Primary	7/30/2021	<0.597 U	<0.190 U
	Primary	11/18/2021	<0.597 U	<0.190 U
	Primary	3/22/2022	<0.597	<0.190
	Primary	6/14/2022	<0.597 J3	<0.190 J3
	Primary	9/9/2022	<0.597	<0.190
	Primary	11/30/2022	<0.597	<0.190
	Primary	2/25/2023	<0.597	<0.190
	Primary	6/12/2023	<0.597	<0.190
	Field Duplicate	6/12/2023	<0.597	<0.190 L1
	Primary	9/5/2023	<0.597	<0.190
	Primary	11/17/2024	<0.597	<0.190
	Primary	2/12/2024	<0.597	<0.190

TABLE 5:
HISTORICAL 1,4-DIOXANE AND TCE CONCENTRATIONS
FORMER THERMAL TREATMENT UNIT
NAMMO DEFENSE SYSTEMS INC.

		Chemical Name	1,4-Dioxane	Trichloroethene
		EPA Method	8260B SIM	8260B
		Unit	µg/l	
Location	Sample Type	Screening Level	3.5[(1)]	5
		Sample Date		
TTU-8	Primary	3/27/2018	<0.36 U	<0.18 U
	Primary	6/28/2018	<0.72 U	<0.15 U
	Primary	9/10/2018	<0.26 U	<0.15 U
	Primary	12/10/2018	<0.26 U	<0.15 U
	Primary	3/26/2019	<0.597 U	<0.398 U
	Primary	6/7/2019	<0.597 U	<0.398 U
	Primary	9/16/2019	<0.597 U	<0.398 U
	Primary	12/23/2019	<0.597 U	<0.398 U
	Primary	3/16/2020	<0.597 U	<0.398 U
	Field Duplicate	3/16/2020	<0.597 U	<0.398 U
	Primary	6/18/2020	<0.597 U	<0.190 U
	Field Duplicate	6/18/2020	<0.597 U	<0.190 U
	Primary	7/21/2020	<0.597 U	<0.190 U
	Primary	12/4/2020	<0.597 U	<0.190 U
	Primary	3/29/2021	<0.597 U	<0.190 U
	Primary	5/6/2021	<0.597 U	<0.190 U
	Primary	7/30/2021	<0.597 U	<0.190 U
	Primary	11/18/2021	<0.597 U	<0.190 U
	Primary	3/22/2022	<0.597	<0.190
	Primary	6/14/2022	<0.597	<0.190
	Primary	9/9/2022	<0.597	<0.190
	Field Duplicate	9/9/2022	<0.597	<0.190
	Primary	11/30/2022	<0.597	<0.190
	Primary	2/25/2023	<0.597	<0.190
	Primary	6/12/2023	<0.597	<0.190
	Primary	9/5/2023	<0.597	<0.190
	Field Duplicate	9/6/2023	28.7	<0.190
	Primary	11/17/2023	<0.597	<0.190
	Field Duplicate	11/17/2023	<0.597	<0.190
	Primary	2/14/2024	<0.597	<0.190
TTU-9A	Primary	3/27/2018	<0.36 U	<0.18 U
	Primary	6/28/2018	<0.72 U	<0.15 U
	Primary	9/10/2018	<0.26 U	<0.15 U
	Primary	12/10/2018	<0.26 U	<0.15 U
	Primary	3/26/2019	<0.597 U	<0.398 U
	Primary	6/7/2019	<0.597 U	<0.398 U
	Primary	9/16/2019	<0.597 U	<0.398 U
	Primary	12/20/2019	1.01 J	<0.398 U
	Primary	3/12/2020	11.9 J	<0.398 U
	Primary	6/17/2020	<0.597 U	<0.190U
	Primary	7/20/2020	<0.597 U	<0.190U
	Primary	12/2/2020	<0.597 U	6.46 J
	Primary	3/30/2021	<0.597 U	7.53
	Primary	5/6/2021	<0.597 U	4.76
	Primary	7/29/2021	<0.597 U	2.75
	Primary	11/17/2021	<0.597 U	0.911 J
	Field Duplicate	11/17/2021	<0.597 U	0.985 J
	Primary	3/22/2022	<0.597	0.944 E4
	Primary	6/13/2022	4.82	<0.190
	Primary	7/21/2022	<0.597	0.221 J
	Primary	9/8/2022	<0.597	<0.190
	Primary	11/29/2022	<0.597	<0.190
	Primary	2/23/2023	18.4	<0.190
	Primary	3/21/2023	<0.597	<0.190
	Primary	6/13/2023	<0.597	<0.190
	Field Duplicate	6/13/2023	<0.597	<0.190 L1
	Primary	9/6/2023	<0.597	<0.190
	Primary	11/17/2023	<0.597	<0.190
	Primary	2/13/2024	<0.597	<0.190

TABLE 5:
HISTORICAL 1,4-DIOXANE AND TCE CONCENTRATIONS
FORMER THERMAL TREATMENT UNIT
NAMMO DEFENSE SYSTEMS INC.

		Chemical Name	1,4-Dioxane	Trichloroethene
		EPA Method	8260B SIM	8260B
		Unit	µg/l	
Location	Sample Type	Screening Level	3.5[(1)]	5
		Sample Date		
TTU-10	Primary	3/27/2018	<0.36 U	<0.18 U
	Primary	6/28/2018	<0.72 U	<0.15 U
	Primary	9/10/2018	<0.26 U	<0.15 U
	Primary	12/10/2018	<0.26 U	<0.15 U
	Primary	3/26/2019	<0.597 U	<0.398 U
	Primary	9/16/2019	<0.597 U	<0.398 U
	Primary	12/23/2019	<0.597 U	<0.398 U
	Primary	3/13/2020	<0.597 U	<0.398 U
	Primary	6/18/2020	<0.597 U	<0.190 U
	Primary	7/21/2020	<0.597 U	<0.190 U
	Primary	12/4/2020	<0.597 U	<0.190 U
	Primary	3/29/2021	<0.597 U	<0.190 U
	Primary	5/6/2021	<0.597 U	<0.190 U
	Primary	8/6/2021	<0.597 U	<0.190 U
	Primary	11/18/2021	<0.597 U	<0.190 U
	Primary	3/22/2022	1.58 E4	<0.190
	Primary	6/14/2022	<0.597 J3	<0.190
	Field Duplicate	6/14/2022	<0.597 J3	<0.190
	Primary	9/9/2022	<0.597	<0.190
	Primary	11/30/2022	<0.597	<0.190
	Primary	2/27/2023	<0.597	<0.190
	Primary	6/12/2023	<0.597	<0.190
	Primary	9/5/2023	<0.597	<0.190 R5
	Primary	11/21/2023	<0.597	<0.190 L1;R7
	Field Duplicate	11/21/2023	<0.597	<0.190 L1;R7
	Primary	2/14/2024	<0.597	<0.190
	Field Duplicate	2/14/2024	<0.597	<0.190
TTU-11	Primary	9/23/2015	380	3100
	Field Duplicate	9/23/2015	400	3100
	Primary	11/23/2015	270	2900
	Primary	2/25/2016	250	2400
	Primary	6/1/2016	282	1600
	Primary	8/18/2016	240	1800
	Primary	11/22/2016	310	2500
	Field Duplicate	11/22/2016	340	2400
	Primary	2/22/2017	222	2010
	Field Duplicate	2/22/2017	224	2080
	Primary	5/23/2017	201	1560
	Field Duplicate	5/23/2017	192	1710
	Primary	8/29/2017	1450	807
	Primary	3/27/2018	671	461
	Primary	9/12/2018	1060	4650
	Primary	12/4/2018	1820	14500
	Field Duplicate	12/4/2018	1840	14800
	Primary	12/10/2018	1820	14500
	Field Duplicate	12/10/2018	1840	14800
	Primary	9/16/2019	1510	11200
	Primary	12/20/2019	855 J-	11500
	Field Duplicate	12/20/2019	907 J-	9400
	Primary	3/12/2020	863	6780
	Primary	6/18/2020	1570	15000
	Primary	7/20/2020	977	17600
	Primary	10/26/2020	358 J	4430
	Primary	10/26/2020	562 J	4870
	Primary	9/23/2021	6.95 J	69.8
	Primary	6/20/2022	<0.597	56.3
	Primary	9/3/2022	<0.597 R7	58.2
	Primary	11/30/2022	<0.597 J3	71.5
	Primary	2/25/2023	11600	67.8
	Primary	6/13/2023	<5.97	45.5
	Primary	9/5/2023	<0.597	78.6
	Primary	11/18/2023	<0.597	67.9
	Primary	2/13/2024	<0.597	76.6

TABLE 5:
HISTORICAL 1,4-DIOXANE AND TCE CONCENTRATIONS
FORMER THERMAL TREATMENT UNIT
NAMMO DEFENSE SYSTEMS INC.

		Chemical Name	1,4-Dioxane	Trichloroethene
		EPA Method	8260B SIM	8260B
		Unit	µg/l	
Location	Sample Type	Screening Level	3.5[(1)]	5
		Sample Date		
TTU-12	Primary	8/29/2017	85.7	335
	Primary	11/27/2017	84.1	301
	Primary	3/27/2018	85.5	484
	Primary	6/28/2018	108	339
	Primary	9/10/2018	91	460
	Primary	12/10/2018	107	454
	Primary	3/25/2019	136	176
	Primary	6/7/2019	120	507
	Primary	9/16/2019	160	543
	Primary	12/20/2019	106	567
	Primary	3/12/2020	94.8 J	407
	Primary	6/17/2020	184	471
	Primary	7/20/2020	82.2	547
	Primary	12/2/2020	159	531
	Primary	3/30/2021	115 J	480
	Primary	5/6/2021	142	540
	Primary	7/29/2021	176	466
	Primary	11/18/2021	133	624
	Field Duplicate	11/18/2021	141	617
	Primary	3/22/2022	149	538
	Primary	6/13/2022	170	487
	Primary	9/9/2022	119	529
	Primary	11/29/2022	117	463
	Primary	2/23/2023	209	452
	Primary	9/7/2023	131	519 M3
	Primary	11/18/2023	153	427
	Primary	12/2/2024	<0.597	511
	Primary	2/12/2024	<0.597	511
TTU-13	Primary	8/29/2017	4	2.6
	Primary	11/27/2017	14.1	5.7
	Primary	3/27/2018	18.3	7.3
	Primary	6/28/2018	33.9	12.6
	Primary	9/10/2018	47.3	24.2
	Primary	12/10/2018	45.2	20.1
	Primary	3/25/2019	55.8	21.7
	Primary	6/7/2019	39.9	22.6
	Primary	9/16/2019	58	18.3
	Primary	12/20/2019	40.2	17
	Primary	3/16/2020	32.2 J	15.4
	Field Duplicate	3/16/2020	33.5 J	14.9
	Primary	6/17/2020	48.5	14.6
	Field Duplicate	6/17/2020	54.1	16.6
	Primary	7/20/2020	29.6	13.3
	Field Duplicate	7/20/2020	27.7	13.8
	Primary	12/3/2020	25.3	11.2 J
	Primary	3/30/2021	37.7 J	17.1
	Primary	5/6/2021	37.9	12.9
	Primary	7/29/2021	58.6	11.1
	Primary	11/18/2021	3.26	1.44 J
	Primary	3/22/2022	9.96	5.76
	Primary	6/13/2022	28.9	5.52
	Primary	9/8/2022	13.7	7.06
	Primary	11/29/2022	33.5	12.7
	Primary	2/23/2023	40.1	12.8
	Primary	6/13/2023	32.5	9.34
	Primary	9/6/2023	37.4	9.79
	Primary	11/17/2023	20.9	9.37
	Primary	2/12/2024	31.4	13.6

TABLE 5:
HISTORICAL 1,4-DIOXANE AND TCE CONCENTRATIONS
FORMER THERMAL TREATMENT UNIT
NAMMO DEFENSE SYSTEMS INC.

		Chemical Name	1,4-Dioxane	Trichloroethene
		EPA Method	8260B SIM	8260B
		Unit	µg/l	
Location	Sample Type	Screening Level	3.5[1]	5
		Sample Date		
TTU-14	Primary	8/29/2017	367	657
	Primary	11/27/2017	356	828
	Primary	3/27/2018	363	1030
	Primary	6/28/2018	381	875
	Primary	9/10/2018	338	689
	Primary	12/17/2018	331	694
	Primary	3/27/2019	356	780
	Primary	9/16/2019	422	921
	Primary	12/20/2019	280	1060
	Primary	3/12/2020	278 J	880
	Primary	6/17/2020	504	891
	Primary	7/20/2020	241	1210
	Primary	12/2/2020	388	917
	Primary	3/30/2021	280 J	990
	Primary	5/6/2021	370	831
	Primary	7/29/2021	493	966
	Primary	11/18/2021	279	917
	Primary	3/22/2022	339	908
	Field Duplicate	3/22/2022	321	879
	Primary	6/14/2022	297 J3	1040
	Primary	9/9/2022	297	1020
	Primary	11/29/2022	288	882
	Primary	2/25/2023	339	807
	Primary	6/13/2023	269	764
	Primary	9/7/2023	246	921
	Primary	11/18/2023	356	712
	Primary	2/12/2024	286	829
TTU-15	Primary	3/27/2019	3.54	<0.398 U
	Primary	9/16/2019	3.95	<0.398 U
	Primary	12/20/2019	6.09	<0.398 U
	Primary	3/12/2020	3.02	<0.398 U
	Primary	6/17/2020	5.32	<0.190U
	Primary	7/20/2020	2.81 J	<0.190U
	Primary	12/2/2020	<0.597 U	3.1
	Primary	3/29/2021	5.33 J	12.9
	Primary	5/5/2021	3.83	11.7
	Primary	7/29/2021	6.26	13
	Primary	11/17/2021	5.9	10.3
	Primary	3/21/2022	6.93	7.89
	Primary	6/13/2022	9.83	6.23
	Primary	9/8/2022	8.21	6.08
	Primary	11/29/2022	27.5	5.13
	Primary	2/25/2023	15.7	4.9
	Primary	6/12/2023	13.5	3.76
	Primary	9/6/2023	4.69	3.05
	Primary	11/18/2023	12.2	3.24
	Field Duplicate	11/18/2023	6.38	3.21
	Primary	4/9/2024	NS	<3.80
TTU-16	Primary	3/13/2020	2470 J	51500
	Primary	6/17/2020	4310	68400
	Field Duplicate	6/17/2020	5610	70200
	Primary	7/20/2020	2220 J-	92200
	Primary	12/2/2020	1730	80000
	Field Duplicate	12/2/2020	1990	96000
	Primary	3/29/2021	2880	76800
	Field Duplicate	3/29/2021	2550	71800
	Primary	5/5/2021	4920	77400 J
	Field Duplicate	5/5/2021	5270	38500 J
	Primary	7/29/2021	5140	86000
	Field Duplicate	7/29/2021	5710	87300
	Primary	11/17/2021	3930	93200
	Primary	3/21/2022	5430	103000
	Primary	6/13/2022	3600 J3	96500
	Primary	9/8/2022	3820 R7	9520
	Primary	11/29/2022	3180	80000
	Primary	2/25/2023	32800	69100
	Field Duplicate	2/25/2023	39600	83600
	Primary	6/13/2023	2300	78000 L1
	Primary	9/7/2023	2880	59400
	Primary	11/18/2023	3470	30800
	Primary	4/9/2024	NS	68500

TABLE 5:
HISTORICAL 1,4-DIOXANE AND TCE CONCENTRATIONS
FORMER THERMAL TREATMENT UNIT
NAMMO DEFENSE SYSTEMS INC.

		Chemical Name	1,4-Dioxane	Trichloroethene
		EPA Method	8260B SIM	8260B
		Unit	µg/l	
Location	Sample Type	Screening Level	3.5[(1)]	5
		Sample Date		
TTU-17	Primary	3/13/2020	<0.0474 U	0.463 J
	Primary	6/17/2020	<0.597 U	0.321 J
	Primary	7/20/2020	<0.597 U	0.367 J
	Primary	12/2/2020	<0.597 U	1.56
	Primary	3/29/2021	<0.597 U	5
	Primary	5/5/2021	<0.597 U	4.13
	Primary	7/29/2021	<0.597 U	3.99
	Primary	11/17/2021	<0.597 U	3.08
	Primary	3/21/2022	4.75	3.51
	Primary	6/13/2022	10.1	2.1
	Primary	9/8/2022	242	2.1
	Primary	11/29/2022	264	1.41
	Field Duplicate	11/29/2022	2.11 B;J	1.57
	Primary	2/25/2023	<0.597	1.28
	Primary	6/12/2023	<0.597	1.15 L1
	Primary	9/6/2023	<0.597	1.20
	Primary	11/18/2023	<0.597	0.716 E4
	Primary	2/12/2024	<0.597	0.661 E4
TTU-19	Primary	10/26/2020	915 J	9990
	Primary	10/26/2020	781 J	12900
	Primary	11/18/2023	<0.597	478
	Primary	6/20/2022	<0.597	189
	Field Duplicate	6/20/2022	<0.597	373
	Primary	9/3/2022	152 H1	293 M3
	Primary	11/30/2022	<0.597	360
	Primary	2/25/2023	318	348
	Primary	6/13/2023	247	310 L1
	Primary	9/7/2023	209	158
	Primary	11/18/2023	279	187
	Primary	2/13/2024	192	195
TTU-20	Primary	10/26/2020	567 J	4480
	Primary	10/26/2020	824 J	6360
	Primary	6/14/2021	1450 J	11200 J
	Primary	9/23/2021	841 J	14300
	Primary	11/18/2021	2140	13400
	Primary	6/16/2022	1540 Q	10800
	Primary	9/3/2022	1140 Q	13200 L1
	Field Duplicate	9/3/2022	1250 Q	10700
	Primary	11/30/2022	1490	12400
	Primary	2/25/2023	19600	12800
	Primary	9/14/2023	499	2.43 E4
	Primary	2/13/2024	378	1860

TABLE 5:
HISTORICAL 1,4-DIOXANE AND TCE CONCENTRATIONS
FORMER THERMAL TREATMENT UNIT
NAMMO DEFENSE SYSTEMS INC.

		Chemical Name	1,4-Dioxane	Trichloroethene
		EPA Method	8260B SIM	8260B
		Unit	µg/l	
Location	Sample Type	Screening Level	3.5[(1)]	5
		Sample Date		
TTU-EX-1	Primary	3/13/2020	24.5	265
	Primary	6/17/2020	284	168
	Primary	7/20/2020	207	163
	Primary	12/2/2020	466	240
	Primary	3/29/2021	340 J	262
	Primary	5/5/2021	258	286
	Primary	7/29/2021	702	372
	Primary	11/17/2021	112	79
	Primary	3/21/2022	244	181
	Primary	6/13/2022	324 J3	174
	Primary	9/8/2022	68.2	75.1
	Primary	11/29/2022	105	59.1
	Primary	2/23/2023	251	127
	Primary	6/13/2023	220	161 L1
	Primary	9/6/2023	163	195
	Field Duplicate	9/6/2023	202	194
	Primary	11/18/2023	351	182
	Primary	2/13/2024	326	241
TTU-EX-2	Primary	3/13/2020	198 J	327
	Primary	6/17/2020	405	549
	Primary	7/20/2020	212	561
	Primary	12/2/2020	424	506
	Primary	3/30/2021	334 J	634
	Primary	5/5/2021	218	536
	Primary	7/29/2021	523	630
	Primary	11/17/2021	158	238
	Primary	3/21/2022	213	234
	Primary	6/13/2022	189 J3	315
	Primary	9/8/2022	74.9	68.1
	Primary	11/29/2022	143	197
	Primary	2/23/2023	162	166
	Field Duplicate	2/23/2023	197	143
	Primary	6/13/2023	220	161 L1
	Primary	9/6/2023	191	433
	Primary	11/18/2023	266	329
	Primary	2/13/2024	256	415
TTU-EX-3	Primary	3/13/2020	175 J	5960
	Primary	6/17/2020	785	6050
	Primary	7/20/2020	610	7390
	Primary	12/2/2020	805 J-	5970 J
	Primary	3/30/2021	697	5560
	Primary	5/5/2021	536	5540
	Primary	7/29/2021	1010	7260
	Primary	11/17/2021	909	8120
	Field Duplicate	11/17/2021	969	8010
	Primary	3/21/2022	885	6560
	Primary	6/13/2022	863 J3	6020
	Primary	9/8/2022	741	7220
	Primary	11/29/2022	735	6620
	Primary	2/23/2023	916	6520
	Primary	6/13/2023	761	7690 L1
	Field Duplicate	6/12/2023	721	7580 L1
	Primary	9/6/2023	598	8100
	Primary	11/18/2023	1490	6890
	Primary	2/12/2024	523	5930

TABLE 5:
HISTORICAL 1,4-DIOXANE AND TCE CONCENTRATIONS
FORMER THERMAL TREATMENT UNIT
NAMMO DEFENSE SYSTEMS INC.

		Chemical Name	1,4-Dioxane	Trichloroethene
		EPA Method	8260B SIM	8260B
		Unit	µg/l	
Location	Sample Type	Screening Level	3.5[(1)]	5
		Sample Date		
TTU-EX-4	Primary	3/13/2020	16.1	811
	Primary	6/17/2020	23.7	1040
	Primary	7/20/2020	18.1	934
	Primary	12/2/2020	20.7	501
	Primary	3/30/2021	16.3	486
	Primary	5/5/2021	12.8	420
	Primary	7/29/2021	29	461
	Primary	11/17/2021	16.1	755
	Primary	3/21/2022	23.9	909
	Primary	6/13/2022	27.4	579
	Field Duplicate	6/13/2022	26.1	635
	Primary	9/8/2022	41.4	698
	Primary	11/29/2022	51.5	612
	Primary	2/23/2023	16.7	836
	Primary	6/13/2023	13.5	970 L1
	Primary	9/7/2023	19.8	698
	Field Duplicate	9/7/2023	12.1	742
	Primary	11/18/2023	9.81	673
	Primary	2/12/2024	15.9	588
TTU-EX-5	Primary	3/13/2020	<0.0532 U	0.929 J
	Field Duplicate	3/13/2020	<0.055 U	0.775 J
	Primary	6/17/2020	<0.597 U	0.456 J
	Primary	7/20/2020	<0.597 U	0.562 J
	Field Duplicate	7/20/2020	<0.597 U	0.637 J
	Primary	12/2/2020	<0.597 U	4.18 J
	Field Duplicate	12/2/2020	<0.597 U	3.89 J
	Primary	3/30/2021	<0.597 U	6.53
	Primary	5/5/2021	<0.597 U	5.52
	Primary	7/29/2021	<0.597 U	5.51
	Primary	11/17/2021	<0.597 U	6.91
	Primary	3/21/2022	<0.597	5.74
	Field Duplicate	3/21/2022	<0.597	5.98
	Primary	6/13/2022	<0.597	5.58
	Primary	9/8/2022	2.16 E4	4.96
	Field Duplicate	9/8/2022	<0.597	5.06
	Primary	11/29/2022	3.4 B	4.51
	Primary	2/23/2023	<0.597	4.45
	Primary	6/12/2023	<0.597	4.01 L1
	Primary	9/7/2023	<0.597	3.45
	Primary	11/17/2023	<0.597	3.65
	Primary	2/13/2024	9.87	3.57
	Field Duplicate	2/13/2024	<0.597	3.08

Notes:

µg/l - micrograms per liter

EPA - Environmental Protection Agency

NA - Not Analyzed

NS - No sample collected

SIM - Selected Ion Monitoring

< - Concentration is below laboratory reporting limits

Concentration detected above the method detection limit

(1) - Interim Screening Level

V = The sample concentration is too high to evalulate accurate spike recoveries

J = The analyte was positively identified; the associated numerical value is the approximate

J3 = The associated batch QC was outside the established quality control range for precision

R7 = LFB/LFBD RPD exceeded the laboratory acceptance limit. Recovery met acceptance criteria

Q = Sample was prepared and/or analyzed past holding time as defined in the method. Concentration

E4 = Concentration estimated. Analyte was detected below laboratory minimum reporting level (MRL) but above MDL.

U = Not detected at the Reporting Limit (or MDL where applicable).

B = The same analyte is found in the associated blank.

M1 = Matrix spike recovery was high; the associated blank spike recovery was acceptable.

L1 = The associated blank spikerecovery was above laboratory acceptance limits.

Figures

PLOT DATE: 2/2/2024

C:\Users\Hernandez\OneDrive - Pinyon Environmental Inc\DWG\TTU Q4 2023\TTU_GW01.dwg

LEGEND

Approximate Property Boundary

N

Notes:
-NG: Northrop Grumman
-DFR: Desert Firing Range
-NDS: Nammo Defense Systems Inc.
-SRP: Salt River Project
-SRPMIC: Salt River Pima-Maricopa Indian Community
-SDI: Special Devices Inc.
-WBO: Water Bore-Out
-TTU: Thermal Treatment Unit
Drawing not to scale

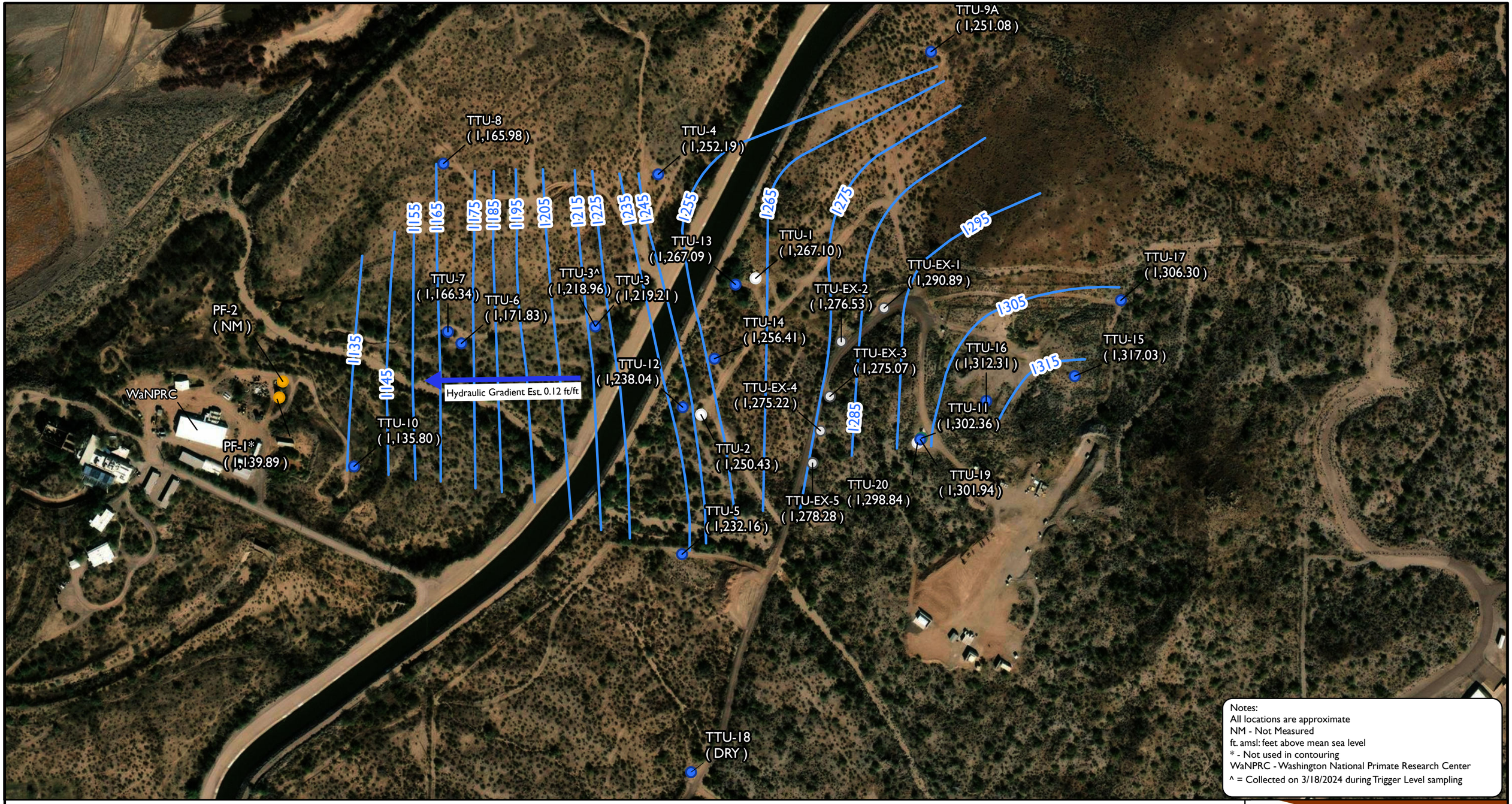
Pinyon Environmental, Inc.

SITE VICINITY MAP

Nammo Defense Systems Inc. Plant #3
Formal Thermal Treatment Unit (TTU)
Mesa, Arizona

Site Location: Section 3, 15 and 27 Township 1N, Range 6E, Gila-Salt River Meridian	Drawn By: SJA	Figure: 1
Pinyon Project Number: 7/22-1522-01.REM001.2	Reviewed By: AP	Date: 2/2/2024

Coordinate System: NAD83 ARIZONA STATE PLANES, CENTRAL ZONE, US FOOT - AZ83-CF



Notes:
All locations are approximate
NM - Not Measured
ft. amsl: feet above mean sea level
* - Not used in contouring
WaNPRC - Washington National Primate Research Center
^ = Collected on 3/18/2024 during Trigger Level sampling

N

Legend

- Extraction Well
- Extraction Well Currently Used for Monitoring
- Extraction/Injection Well
- Monitoring Well
- Private Production Well
- Groundwater Elevation Contour (ft amsl) (Contour Interval: 10ft)
- Estimated Regional Groundwater Flow Direction

TTU-1 (1,145.24) = Monitoring Well Location (Groundwater Elevation (ft. amsl))

0 130 260 Feet

Pinyon Environmental, Inc.

QUARTERLY GROUNDWATER CONTOUR MAP- FIRST QUARTER 2024

Nammo Defense Systems Inc.
Former Thermal Treatment Unit (TTU)
Mesa, Arizona

Site Location: Sections 23, Townships 12 North, Range 6 East, Gila-Salt River Meridian	Drawn By: CJB / Figure: 2
Pinyon Project Number: 7/22-1522-01.REM002.1b	Reviewed By: AP / Date: 4/29/2024

Document Path: C:\Users\bowes\Desktop\Arizona Project Delivery - NDS - WBO_TTU GWM\Figures\ArcPro\TTU\2024\1Q\72215220I_NDS_TTU_2024Q1.aprx

Notes:
µg/l - micrograms per liter
Health-Based Guidance Level for perchlorate is 14 ug/L
BOLD - Concentration exceeds its respective AWQS or screening level
E8 - Analyte reported to MDL per project specification. Target analyte was not detected in the sample
M2 - Matrix spike recovery was low; the associated blank spike recovery was acceptable
< - Less Than
TTU-15 and TTU-16 results appear to have been switched. The results shown are based on resampling

N

Legend

- Extraction Well
- Extraction Well Currently Used for Monitoring
- Extraction/Injection Well
- Monitoring Well
- Private Production Well
- Estimated extent of Perchlorate concentrations above the AWQS of 14 ug/l
- Dashed where Inferred

((Result/Duplicate Result))

0 100 200 Feet

Pinyon Environmental, Inc.

PERCHLORATE DETECTIONS IN GROUNDWATER - FIRST QUARTER 2024

Nammo Defense Systems Inc.
Former Thermal Treatment Unit (TTU)
Mesa, Arizona

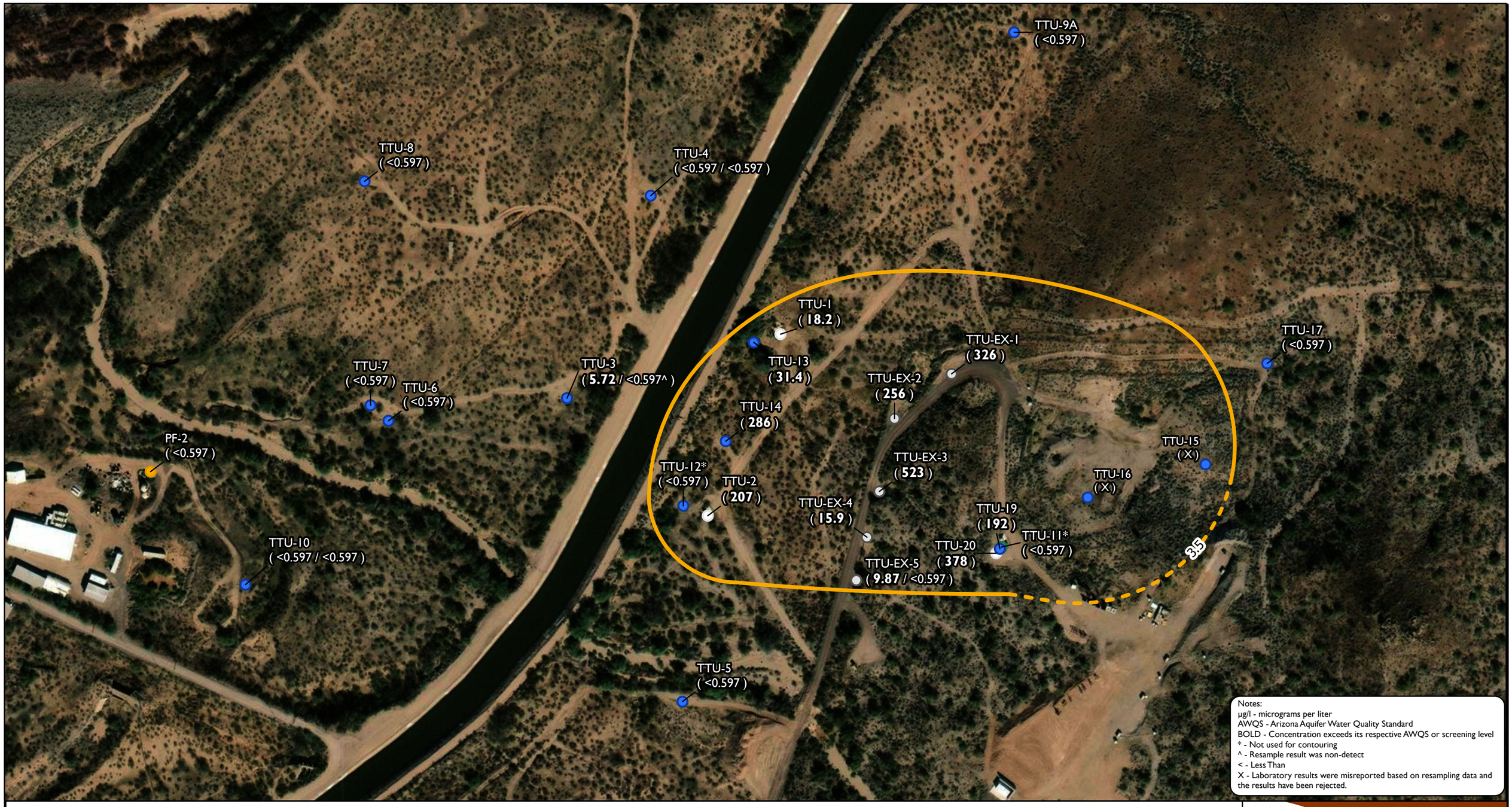
Site Location: Sections 23, Townships 12 North, Range 6 East, Gila-Salt River Meridian

Pinyon Project Number: 7/22-1522-01.REM002.1b

Drawn By: CJB | Figure: 3

Reviewed By: AP | Date: 4/29/2024

Document Path: C:\Users\bowes\Desktop\Arizona Project Delivery - NDS - WBO_TTU GWM\Figures\ArcPro\TTU\2024\1Q\722152201_NDS_TTU_2024Q1.aprx



Notes:
ug/l - micrograms per liter
AWQS - Arizona Aquifer Water Quality Standard
BOLD - Concentration exceeds its respective AWQS or screening level
* - Not used for contouring
^ - Resample result was non-detect
< - Less Than
X - Laboratory results were misreported based on resampling data and the results have been rejected.

N

Legend

- Extraction Well
- Extraction Well Currently Used for Monitoring
- Extraction/Injection Well
- Monitoring Well
- Private Production Well
- Dashed where Inferred
- Estimated extent of 1,4-Dioxane concentrations above the interim screening level of 3.5 ug/l

((Result/Duplicate Result))

0 100 200 Feet

Pinyon Environmental, Inc.

1,4-DIOXANE DETECTIONS IN GROUNDWATER - FIRST QUARTER 2024

Nammo Defense Systems Inc.
Former Thermal Treatment Unit (TTU)
Mesa, Arizona

Site Location: Sections 23, Townships 12 North, Range 6 East, Gila-Salt River Meridian	Drawn By: CJB / Figure: 4
Pinyon Project Number: 7/22-1522-01.REM002.1b	Reviewed By: AP / Date: 4/28/2024

Document Path: C:\Users\bowes\Desktop\Arizona Project Delivery - NDS - WBO_TTU GWM\Figures\ArcPro\TTU\2024\1Q\722152201_NDS_TTU_2024Q1.aprx

TTU-9A
(<0.188)

TTU-8
(<0.188)

TTU-4
(<0.188 / <0.188)

TTU-13
(4.80)

TTU-1
(1.21)

TTU-EX-1
(**88.9**)

TTU-17
(<0.188)

TTU-7
(<0.188)

TTU-6
(<0.188)

TTU-3
(<0.188)

TTU-EX-2
(**83.3**)

TTU-14
(**81.7**)

TTU-15
(<3.76)

PF-2
(<0.188)

TTU-12
(**52.6**)

TTU-2
(**31.6**)

TTU-EX-3
(**836**)

TTU-16
(**2,440 E4**)

TTU-19
(**17.2**)

TTU-10
(<0.188 / <0.188)

TTU-EX-4
(**91.1**)

TTU-11*
(5.47 E4)

TTU-EX-5
(<0.188 / <0.188)

TTU-20
(**154**)

TTU-5
(<0.188)

Notes:
µg/l - micrograms per liter
AWQS - Arizona Aquifer Water Quality Standard
BOLD - Concentration exceeds its respective AWQS or screening level
* - Not used for contouring
E4 - Concentration estimated. Analyte was detected below laboratory minimum reporting level (MRL) but above method detection limit (MDL)
< - Less Than
TTU-15 and TTU-16 results appear to have been switched. The results shown are based on resampling

N

Legend

- Extraction Well
- Extraction Well Currently Used for Monitoring
- Extraction/Injection Well
- Monitoring Well
- Private Production Well
- Dashed where Inferred
- Estimated extent of 1,1 Dichloroethene concentrations above the AWQS of 7 ug/l

((Result/Duplicate Result))

0 100 200 Feet

Pinyon Environmental, Inc.

1,1 DICHLOROETHENE DETECTIONS IN GROUNDWATER - FIRST QUARTER 2024

Nammo Defense Systems Inc.
Former Thermal Treatment Unit (TTU)
Mesa, Arizona

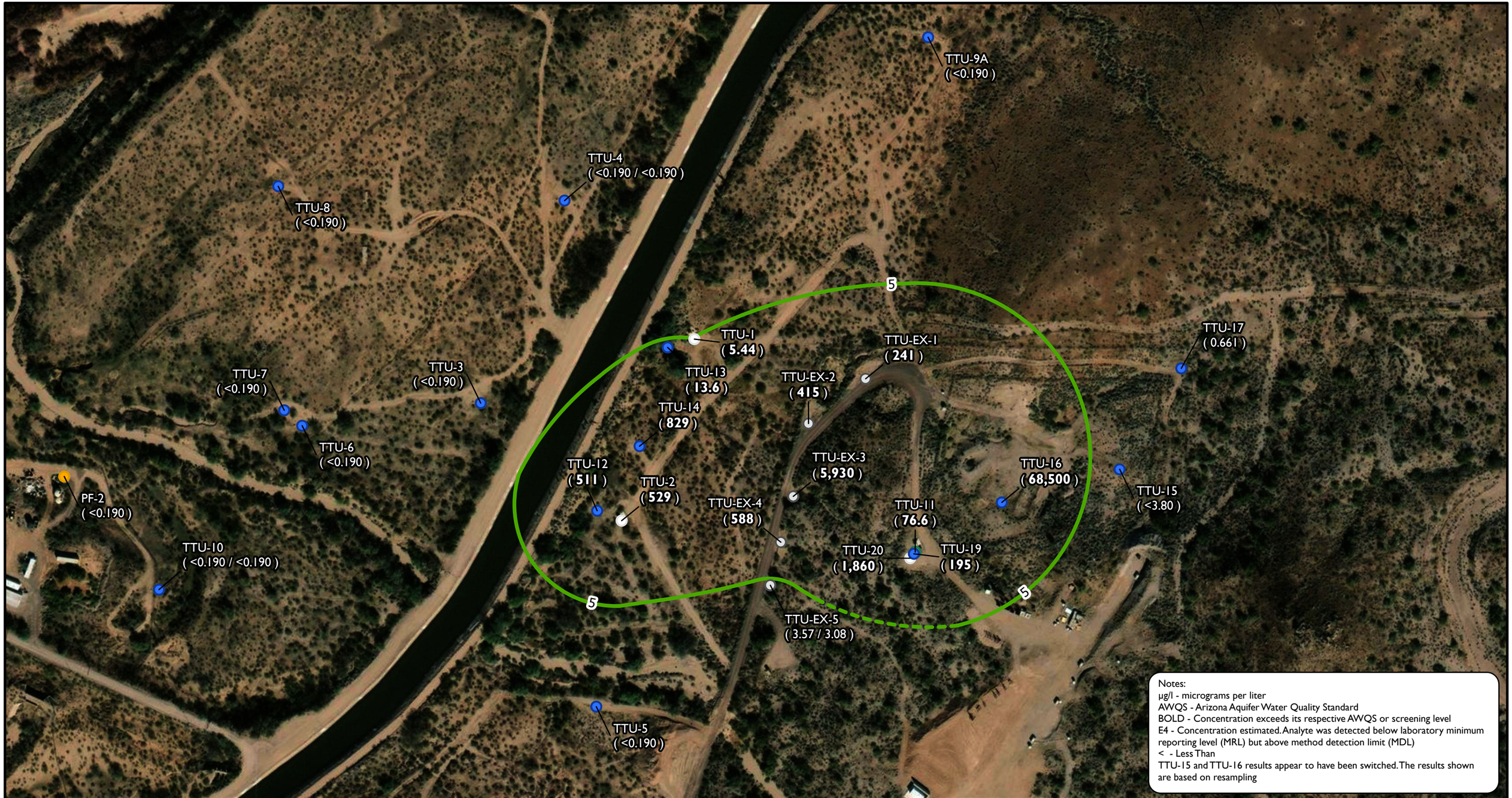
Site Location: Sections 23, Townships 12 North, Range 6 East, Gila-Salt River Meridian

Pinyon Project Number: 7/22-1522-01.REM002.1b

Drawn By: CJB | Figure: 5

Reviewed By: AP | Date: 4/29/2024

Document Path: C:\Users\bowes\Desktop\Arizona Project Delivery - NDS - WBO_TTU GWM\Figures\ArcPro\TTU\2024\1Q\72152201_NDS_TTU_2024Q1.aprx



Notes:
µg/l - micrograms per liter
AWQS - Arizona Aquifer Water Quality Standard
BOLD - Concentration exceeds its respective AWQS or screening level
E4 - Concentration estimated. Analyte was detected below laboratory minimum reporting level (MRL) but above method detection limit (MDL)
< - Less Than
TTU-15 and TTU-16 results appear to have been switched. The results shown are based on resampling

N

Legend

- Extraction Well
- Extraction Well Currently Used for Monitoring
- Extraction/Injection Well
- Monitoring Well
- Private Production Well
- Dashed where Inferred
- Estimated extent of Trichloroethene (TCE) concentrations above the AWQS of 5 ug/l

((Result/Duplicate Result))

0 150 300 Feet

Pinyon Environmental, Inc.

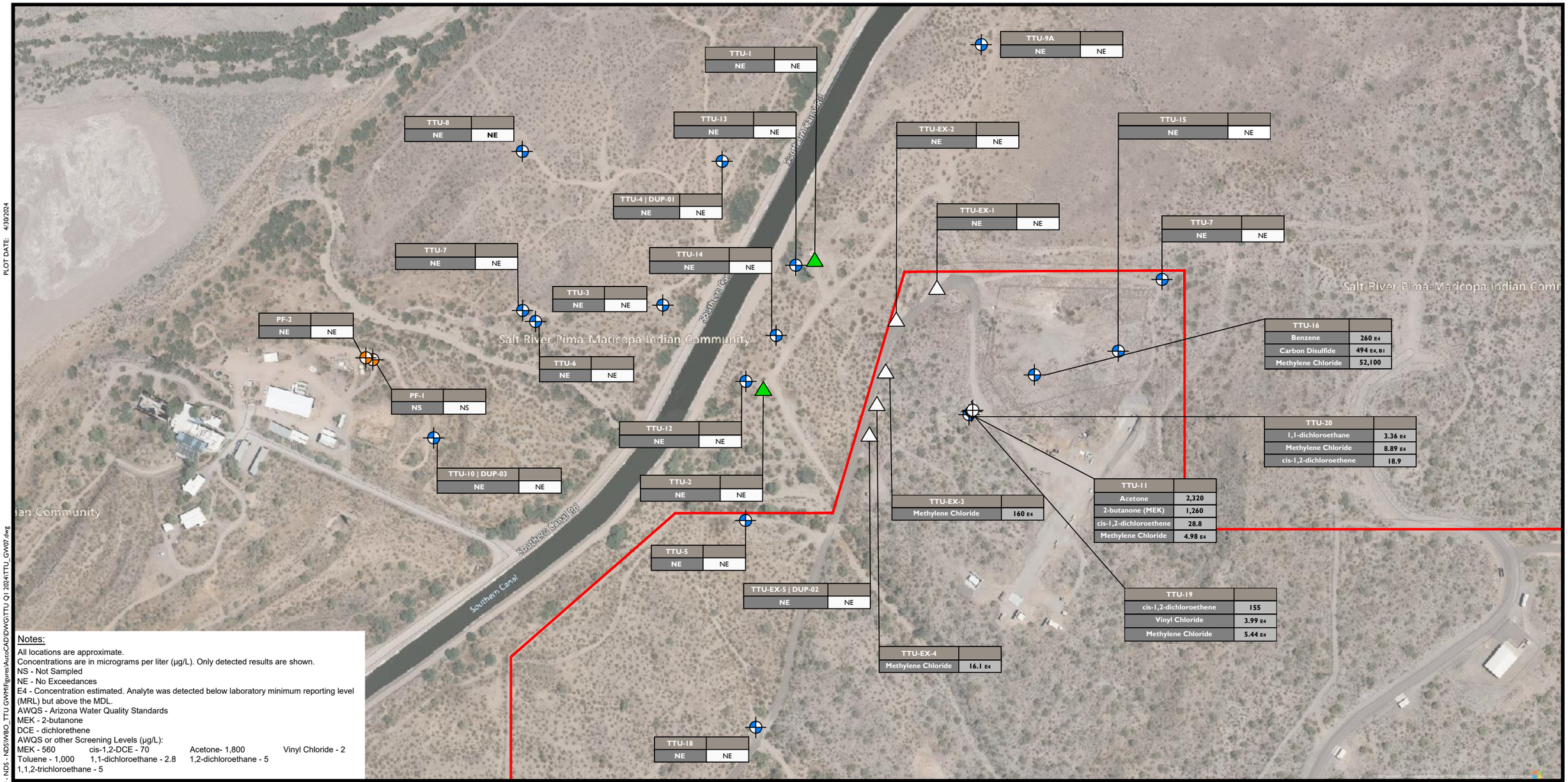
TRICHLOROETHENE DETECTIONS IN GROUNDWATER - FIRST QUARTER 2024

Nammo Defense Systems Inc.
Former Thermal Treatment Unit (TTU)
Mesa, Arizona

Site Location: Sections 23, Townships 12 North, Range 6 East, Gila-Salt River Meridian	Drawn By: CJB / Figure: 6
Pinyon Project Number: 7/22-1522-01.REM002.1b	Reviewed By: AP / Date: 4/29/2024

Document Path: C:\Users\bowes\Desktop\Arizona Project Delivery - NDS - WBO_TTU GWM\Figures\ArcPro\TTU\2024\1Q\722152201_NDS_TTU_2024Q1.aprx

PLOT DATE: 4/30/2024



Notes:

All locations are approximate.
Concentrations are in micrograms per liter (µg/L). Only detected results are shown.
NS - Not Sampled
NE - No Exceedances
E4 - Concentration estimated. Analyte was detected below laboratory minimum reporting level (MRL) but above the MDL.
AWQS - Arizona Water Quality Standards
MEK - 2-butanone
DCE - dichlorethene
AWQS or other Screening Levels (µg/L):

MEK - 560	cis-1,2-DCE - 70	Acetone- 1,800	Vinyl Chloride - 2
Toluene - 1,000	1,1-dichloroethane - 2.8	1,2-dichloroethane - 5	
1,1,2-trichloroethane - 5			

LEGEND

- Extraction Well
- Monitoring Well
- Primate Production Well
- Extraction and Pilot Test Injection Well
- Monitoring / Injection Well
- Extraction Well Currently used for Monitoring
- TTU-1 = Monitoring Well Location
- NDS Leased Property Boundary with SRP-MIC

5.76 Exceeds Aquifer Water Quality Standards or other applicable screening levels

N

0 250' 500'

SCALE: 1" = 250'

Pinyon Environmental, Inc.

VOC EXCEEDANCES IN GROUNDWATER - FIRST QUARTER 2024

Nammo Defense Systems Inc.
Former Thermal Treatment Unit (TTU)
Mesa, Arizona

Site Location: Section 23, Township 12N, Range 6E, Gila-Salt River Meridian

Pinyon Project Number: 7/22-1522-01.REM002.1b

Drawn By: SJA | Figure: 7

Reviewed By: DW | Date: 4/30/2024

Coordinate System: NAD83 ARIZONA STATE PLANES, CENTRAL ZONE, US FOOT - AZ83-CF

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Attachments

Attachment 1 – Field Notes

Location	Measurement Date	Depth to Groundwater (ft btoc)	Measurement Date	Depth to Groundwater (ft btoc)
TTU-1	11/17/2023	43.07	2-13-24	45.63
TTU-2	11/17/2023	64.03	2-13-24	64.01
TTU-3	11/17/2023	92.07	2-12-24	88.82
TTU-4	11/17/2023	52.35	2-13-24	52.93
TTU-5	11/17/2023	81.76	2-13-24	82.77
TTU-6	11/17/2023	123.85	2-12-24	129.01
TTU-7	11/17/2023	131.85	2-12-24	135.5
TTU-8	11/17/2023	143.08	2-14-24	144.25
TTU-9A	11/17/2023	27.35	2-13-24	66.96
TTU-10	11/21/2023	165.18	2-14-24	166.62
TTU-11	11/18/2023	34.33	2-13-24	36.84
TTU-12	11/18/2023	75.62	2-12-24	74.17
TTU-13	11/17/2023	43.27	2-12-24	43.7
TTU-14	11/18/2023	61.81	2-12-24	60.39
TTU-15	11/18/2023	32.70	2-13-24	33.82
TTU-16	11/18/2023	24.67	2-13-24	26.24
TTU-17	11/18/2023	39.98	2-12-24	41.19
TTU-18	11/18/2023	Dry	—	—
TTU-19	11/18/2023	32.56	2-13-24	34.87
TTU-20	11/17/2023	33.65	2-13-24	38.06
TTU-EX-1	11/18/2023	29.27	2-13-24	30.8
TTU-EX-2	11/18/2023	38.04	2-13-24	39.87
TTU-EX-3	11/18/2023	40.18	2-12-24	41.78
TTU-EX-4	11/18/2023	43.78	2-12-24	44.74
TTU-EX-5	11/17/2023	40.71	2-12-24	41.22
PF-1	11/21/2023	No sound	—	—
PF-2	11/21/2023	NM (probe won't fit)	2-14-24	NA

Pinyon
Environmental, Inc.

Well Sampling Record	
Project Name	Nammo TTU
Project Number	722152201.002
Well ID / ADWR #	PF-2
Date Completed	N/A
Casing Material	Steel
Casing Diameter (in)	6 5/8
Screen (ft btoc)	300-400
Well Total Depth (ft btoc)	400
Survey Information	Northing: 909166.890 / Easting: 760122.250
Deployment	
Date / Time	NA
Type of Sampler	Production Well; spigot
Size of Sampler	NA
DTW (ft btoc)	NA
Deployment Depth (ft btoc)	400
Personnel	NA
Notes	Purge 15 minutes. Take parameters (starting and 15 min. after)
Retrieval and/or Sampling	
Date / Time	2/14/24 @ 1500
DTW (ft btoc)	*Collect before turning well on* WLM stopping at 168, no hit
Sampler Integrity	NA
Personnel	MER, IGF, CJW
Notes	

	±1°C	±0.1	**Field Parameters**	±3%	±10%	±10%
Date / Time	Water Temp (°C)	pH (SU)	±10 ORP (mV)	Sp Cond (µS/cm)	DO (mg/L)	Turbidity
2/14/24 1414	23.9	7.22	-223.2	1294	2.64	7.1
1419	23.9	7.20	-224.7	1290	2.69	3.0
1424	24.7	7.18	-104.6	water	spilled	3.1
1429	25.1	7.17	-228.6	1277	3.05	2.3

Sample ID	PF-2-400- 20240214
QAQC Samples	~~DUP-01~~ MS/MSD #3
Containers	(2) 250 mL HDPE (no pres. & filtered) & (~~10~~ 12) 40 mL amber VOAs (10F)
Preservatives	HCl
Analysis	Perchlorate (Method 314.0 at Eurofins) / VOCs / 1,4-Dioxane
Sampler Reset	Yes / No

Notes

Well on at 1410; turned off at 1502.

Make sure to indicate "drinking water" on COC for Eurofins

1434	25.2	7.27	-115.9	1291	3.89	2.2
1439	24.3	6.95	-231.9	1287	3.30	(10F) ~~1.5~~ 2.2
1446	24.1	7.15	-225.1	1288	3.43	1.2
1451	24.0	7.13	-226.1	1290	3.22	1.0

Pinyon Environmental, Inc.

Well Sampling Record	
Project Name	Nammo TTV
Project Number	722152201.002
Well ID / ADWR #	TTV-1 / 55-914440
Date Completed	6/6/2012
Casing Material	PVC
Casing Diameter (in)	4
Screen (ft btoc)	30-70
Well Total Depth (ft btoc)	75
Survey Information	Northing 909420.734 / Easting 761281.203
Deployment	
Date / Time	~~11-17-23 @ 1342~~ CW NA
Type of Sampler	Pump, Production well
Size of Sampler	NA
DTW (ft btoc)	43.07 * Before turning on well on 11-17-23
Deployment Depth (ft btoc)	50
Personnel	16F
Notes	System is odorous
Retrieval and/or Sampling	
Date / Time	2-13-24 @ 1027
DTW (ft btoc)	45.63
Sampler Integrity	Spigot
Personnel	16F, CJW
Notes	-

Field Parameters

Date / Time	~~Water Temp (°C)~~ pH	~~pH (SU)~~ Water temp (°C)	ORP (mV)	Sp Cond (µS/cm)	DO (mg/L)	Turbidity
2-13-24 @ 1003	7.68	18.9	-130.7	1018	5.80	2.5
1008	7.84	21.8	-102.0	1028	5.41	1.6
1013	7.83	22.3	-74.5	1026	5.19	1.4
1018	7.79	22.4	-78.9	1020	5.19	1.6

Sample ID	TTV-1-50-20240213
QAQC Samples	-
Containers	(1) 125 mL HDPE (6) 40mL Amber Vials
Preservatives	HCl
Analysis	Perchlorate / VOCs / 1,4-Dioxane
Sampler Reset	Yes / (No)
Notes	Well off upon arrival, well on @ 0857, well on upon departure.

Pinyon Environmental, Inc.

Well Sampling Record	
Project Name	Nammo TTU
Project Number	~~722152201.001~~
Well ID / ADWR #	~~TTU-2~~ / NA
Date Completed	10/17/2013
Casing Material	PVC
Casing Diameter (in)	4
Screen (ft btoc)	49.4 - 179.6
Well Total Depth (ft btoc)	185
Survey Information	Northing 909087.852 / Easting 761148.265
Deployment	
Date / Time	~~2-13-24 @ 1101~~ NA
Type of Sampler	~~Spigot~~ Production well, Spigot
Size of Sampler	NA
DTW (ft btoc)	~~71.54 on~~ aw 64.03 on 11-17-23
Deployment Depth (ft btoc)	114.5
Personnel	IGF, CJW
Notes	-
Retrieval and/or Sampling	
Date / Time	2-13-24 @ 1101
DTW (ft btoc)	-
Sampler Integrity	Spigot
Personnel	IGF, CJW
Notes	-

Field Parameters

Date / Time	Water Temp (°C)	pH (SU)	ORP (mV)	Sp Cond (µS/cm)	DO (mg/L)	Turbidity
1037	21.7	7.28	-209.9	3419	5.37	3.0
1042	22.6	7.26	-206.6	3472	5.13	1.3
1047	22.9	7.20	-189.1	3471	5.11	1.5
1052	22.7	7.26	-209.2	3466	5.14	1.4

Sample ID	TTU-2-114-20240212
QAQC Samples	-
Containers	(1) 125 mL HDPE, (6) 40mL Amber VOAs
Preservatives	HCl
Analysis	Perchlorate, VOC's, 1,4-Dioxane
Sampler Reset	Yes / (No)

Notes: ORP did not stabalize within 15 minutes, well on upon departure.

Pinyon Environmental, Inc.

Well Sampling Record	
Project Name	Nammo TTU
Project Number	722152201.002
Well ID / ADWR #	TTU-3 / NA
Date Completed	10/18/2013
Casing Material	PVC
Casing Diameter (in)	4
Screen (ft btoc)	78.1-138.1
Well Total Depth (ft btoc)	143.6
Survey Information	Northing: 909303.363 / Easting: 760888.204
Deployment	
Date / Time	11/17/2023 at 11:25
Type of Sampler	HydraSleeve
Size of Sampler	HS-2-1L
DTW (ft btoc)	92.07
Deployment Depth (ft btoc)	108
Personnel	IGF, CJW
Notes	Water a little cloudy
Retrieval and/or Sampling	
Date / Time	2-12-24 @ 1518
DTW (ft btoc)	88.82
Sampler Integrity	Good
Personnel	IGF, CJW
Notes	" "

Field Parameters

Date / Time	Water Temp (°C)	pH (SU)	ORP (mV)	Sp Cond (µS/cm)	DO (mg/L)	Turbidity
2-12-24 @ 1523	21.9	7.08	-163.3	1327	~~4.17~~ 4.71	14

Sample ID	TTU-3-108- 20240212
QAQC Samples	—
Containers	(1) 125 mL HDPE (no pres.) & (6) 40 mL amber VOAs
Preservatives	HCl
Analysis	Perchlorate / VOCs / 1,4-Dioxane
Sampler Reset	Yes / No

Notes

*__Be careful__ of tether when first opening well; is not secured and can fall in easily

Pinyon Environmental, Inc.

Well Sampling Record	
Project Name	Nammo TTU
Project Number	722152201.002
Well ID / ADWR #	TTU-4 / NA
Date Completed	10-25-2013
Casing Material	PVC
Casing Diameter (in)	4
Screen (ft btoc)	39.5-99.5
Well Total Depth (ft btoc)	104.9
Survey Information	Northing 909673.680 / Easting 761041.975
Deployment	
Date / Time	11-17-2023 @ 0954
Type of Sampler	Hydrasleeve
Size of Sampler	HS-2-1L
DTW (ft btoc)	52.35
Deployment Depth (ft btoc)	52.05
Personnel	IGF, CJW
Notes	Water clear
Retrieval and/or Sampling	
Date / Time	~~TTU-4~~ 2-13-24 @ 1528
DTW (ft btoc)	-
Sampler Integrity	Good
Personnel	IGF, CJW
Notes	Water clear ~~parameters~~

Field Parameters

Date / Time	Water Temp (°C)	pH (SU)	ORP (mV)	Sp Cond (µS/cm)	DO (mg/L)	Turbidity
2-13-24 @ 1542	22.7	7.57	-249.7	~~SC~~ 2198	2.49	2.8

Sample ID	TTU-4-57-20240213	
QAQC Samples	Dup-01	
Containers	(2) 125 mL HDPE, (6) 40ml Amber VOAs	
Preservatives	HCl	
Analysis	VOC's, 1,4-Dioxane	
Sampler Reset	Yes	No
Notes		

Pinyon Environmental, Inc.

Well Sampling Record	
Project Name	Nammo TTU
Project Number	722152201.002
Well ID / ADWR #	TTU-5 / NA
Date Completed	9/20/2014
Casing Material	PVC
Casing Diameter (in)	4
Screen (ft btoc)	59.5-164.5
Well Total Depth (ft btoc)	169.5
Survey Information	Northing: 908747.636 / Easting: 761102.227
Deployment	
Date / Time	11/17/2023 at 13:00
Type of Sampler	HydraSleeve
Size of Sampler	HS-2-1L
DTW (ft btoc)	81.76
Deployment Depth (ft btoc)	110
Personnel	IGF & CJW
Notes	Water clear
Retrieval and/or Sampling	
Date / Time	2/13/24 @ 1450
DTW (ft btoc)	82.77
Sampler Integrity	Good
Personnel	IGF, CJW
Notes	" "

Field Parameters

Date / Time	Water Temp (°C)	pH (SU)	ORP (mV)	Sp Cond (µS/cm)	DO (mg/L)	Turbidity
2/13/24 @ 1455	25.3	~~7.9~~ 7.49	-254.2	688	3.00	4.7

Sample ID	TTU-5-110-20240213	
QAQC Samples	—	
Containers	(1) 125 mL HDPE (no pres.) & (6) 40 mL amber VOAs	
Preservatives	HCl	
Analysis	Perchlorate / VOCs / 1,4-Dioxane	
Sampler Reset	(Yes)	No
Notes		

Pinyon Environmental, Inc.

Well Sampling Record	
Project Name	Nammo TTU
Project Number	722152201.002
Well ID / ADWR #	TTU-6 / NA
Date Completed	10/7/2014
Casing Material	PVC
Casing Diameter (in)	4
Screen (ft btoc)	110-175
Well Total Depth (ft btoc)	180
Survey Information	Northing: 909260.820 / Easting: 760560.096
Deployment	
Date / Time	11/17/2023 at 11:05
Type of Sampler	HydraSleeve
Size of Sampler	HS-2-1L
DTW (ft btoc)	123.85
Deployment Depth (ft btoc)	143
Personnel	IGF & CJW
Notes	Water clear
Retrieval and/or Sampling	
Date / Time	2-12-23 @ 1458
DTW (ft btoc)	129.01
Sampler Integrity	IGF, CJW GOOD
Personnel	
Notes	Slightly cloudy

Field Parameters

Date / Time	Water Temp (°C)	pH (SU)	ORP (mV)	Sp Cond (µS/cm)	DO (mg/L)	Turbidity
2-12-24 @ 1504	22.0	7.11	-165.9	2897	1.16	27

Sample ID	TTU-6-143-20240212	
QAQC Samples	—	
Containers	(1) 125 mL HDPE (no pres.) & (6) 40 mL amber VOAs	
Preservatives	HCl	
Analysis	Perchlorate / VOCs / 1,4-Dioxane	
Sampler Reset	Yes	No
Notes		

Pinyon Environmental, Inc.

Well Sampling Record	
Project Name	Nammo TTU
Project Number	722152201.002
Well ID / ADWR #	TTU-7 / NA
Date Completed	10/8/2014
Casing Material	Steel
Casing Diameter (in)	8.5
Screen (ft btoc)	280-410
Well Total Depth (ft btoc)	410
Survey Information	Northing: 909287.611 / Easting: 760527.269
Deployment	
Date / Time	11/17/2023 at 10:48
Type of Sampler	HydraSleeve
Size of Sampler	HS-2-1L
DTW (ft btoc)	131.35
Deployment Depth (ft btoc)	345
Personnel	IGF & CJW
Notes	Black sediment in bottom half of sleeve
Retrieval and/or Sampling	
Date / Time	2/12/24 @ 1435
DTW (ft btoc)	135.50
Sampler Integrity	Good
Personnel	IGF, CJW
Notes	" "

Field Parameters

Date / Time	Water Temp (°C)	pH (SU)	ORP (mV)	Sp Cond (µS/cm)	DO (mg/L)	Turbidity
2/12/24 @ 1445	22.0	6.97	-175.3	3838	1.55	15

Sample ID	TTU-7-345-20240212	
QAQC Samples	-	
Containers	(1) 125 mL HDPE (no pres.) & (4) 40 mL amber VOAs	
Preservatives	HCl	
Analysis	Perchlorate / VOCs / 1,4-Dioxane	
Sampler Reset	(Yes)	No
Notes		

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Well Sampling Record	
Project Name	Nammo TTU
Project Number	722152201.002
Well ID / ADWR #	~~TTU-15 / 55-228014~~ TTU-8 / NA
Date Completed	~~1/25/2018~~ 4-18-2016
Casing Material	~~Steel~~ PVC
Casing Diameter (in)	~~N/A~~ 4
Screen (ft btoc)	~~Open~~ 135-185
Well Total Depth (ft btoc)	~~100~~ 190
Survey Information	~~Northing: 909185.100 / Easting: 762065.910~~ Northing 909699.266 Easting 760514.908
Deployment	
Date / Time	~~11/18/2023 at 12:25~~ 11-17-23 @ 1018
Type of Sampler	~~HydraSleeve~~ Hydrasleeve
Size of Sampler	~~HS-2-1L~~ HS-2-1L
DTW (ft btoc)	~~32.70~~ 143.08
Deployment Depth (ft btoc)	~~75~~ 164
Personnel	~~IGF & CJW~~ IGF, CJW
Notes	~~Slightly cloudy~~ Water clear
Retrieval and/or Sampling	
Date / Time	2/14/24 @ 1520
DTW (ft btoc)	144.25 @ 2-12-24
Sampler Integrity	GOOD
Personnel	MER, IGF, CJW
Notes	Water clear

Field Parameters

Date / Time	Water Temp (°C)	pH (SU)	ORP (mV)	Sp Cond (µS/cm)	DO (mg/L)	Turbidity
2/14/24 @ 1529	25.8	6.91	-54.9	3107	1.67	5.6

Sample ID	~~TTU-15-75-~~ TTU-8-164-20240214	
QAQC Samples	-	
Containers	(1) 125 mL HDPE (no pres.) & (6) 40 mL amber VOAs	
Preservatives	HCl	
Analysis	Perchlorate / VOCs / 1,4-Dioxane	
Sampler Reset	(Yes)	No
Notes		

Pinyon Environmental, Inc.

Well Sampling Record	
Project Name	Nammo TTU
Project Number	722152201.002
Well ID / ADWR #	~~TTU-EX-3 / 55-231731~~ TTU-9A / NA
Date Completed	~~1/24/2020~~ 6-16-2016
Casing Material	PVC
Casing Diameter (in)	~~8~~ 4
Screen (ft btoc)	~~OPEN~~ 24-99
Well Total Depth (ft btoc)	~~111~~ 104
Survey Information	~~Northing: 909134.941 / Easting: 761465.507~~ Northing 909974.490 Easting 761710.151
Deployment	
Date / Time	11/18/2023 at 10:56 11-17-2023 @ 1224
Type of Sampler	HydraSleeve Hydra Sleeve
Size of Sampler	HS-2-1L HS-2-1L
DTW (ft btoc)	38.75 27.35
Deployment Depth (ft btoc)	76 26.61
Personnel	IGF & HMC IGF, CJW
Notes	Rocky sediment on & in bag Water clear
Retrieval and/or Sampling	
Date / Time	2-13-24 @ 0932
DTW (ft btoc)	66.96
Sampler Integrity	Good
Personnel	IGF + CJW
Notes	Water clear

Field Parameters

Date / Time	Water Temp (°C)	pH (SU)	ORP (mV)	Sp Cond (µS/cm)	DO (mg/L)	Turbidity
2-13-24 @ 0942	18.7	7.45	-173.9	1605	4.90 IGF	3.4

Sample ID	~~TTU-EX-3-76~~ TTU-9A-01-20240213	
QAQC Samples	MS/MSD #2	
Containers	(2) 125 mL HDPE (no pres.) & (12) 40 mL amber VOAs	
Preservatives	HCl	
Analysis	Perchlorate / VOCs / 1,4-Dioxane	
Sampler Reset	(Yes)	No
Notes		

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Well Sampling Record	
Project Name	Nammo TTU
Project Number	722152201.002
Well ID / ADWR #	TTU-10 / NA
Date Completed	4/18/2016
Casing Material	PVC
Casing Diameter (in)	4
Screen (ft btoc)	115-180
Well Total Depth (ft btoc)	185
Survey Information	Northing: 908960.114 / Easting: 760297.013
Deployment	
Date / Time	11/21/2023 at 12:40
Type of Sampler	HydraSleeve
Size of Sampler	HS-2-1L
DTW (ft btoc)	165.18
Deployment Depth (ft btoc)	172
Personnel	IGF & CJW
Notes	Water clear, small bit of sediment at bottom
Retrieval and/or Sampling	
Date / Time	2/14/24 @ 1349
DTW (ft btoc)	166.62
Sampler Integrity	GOOD
Personnel	MER, IGF, CJW
Notes	water clear

Field Parameters

Date / Time	Water Temp (°C)	pH (SU)	ORP (mV)	Sp Cond (µS/cm)	DO (mg/L)	Turbidity
2/14/24 @ 1356	23.5	7.37	-205.9	1431	4.47	21

Sample ID	TTU-10-172-20240214
QAQC Samples	DUP-03
Containers	(2) 125 mL HDPE (no pres.) & (12) 40 mL amber VOAs
Preservatives	HCl
Analysis	Perchlorate / ~~VOCs~~ / 1,4-Dioxane
Sampler Reset	(Yes) / No
Notes	

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Well Sampling Record		
Project Name	Nammo TTU	
Project Number	722152201.002	
Well ID / ADWR #	~~TTU-13 / NA~~	TTU-11 / 55-918534
Date Completed	7/20/2018	9-11-2015
Casing Material	Steel	PVC
Casing Diameter (in)	5	4
Screen (ft btoc)	Open to 80	24-89
Well Total Depth (ft btoc)	80	94
Survey Information	~~Northing: 909405.920 / Easting: 761232.180~~	Northing 909029.758 Easting 761706.470
Deployment		
Date / Time	1.1/17/23 at 14:35	11-18-23 @ 1415
Type of Sampler	HydraSleeve	Hydra Sleeve
Size of Sampler	HS-2-1L	HS-2-1L
DTW (ft btoc)	43.27	34.33
Deployment Depth (ft btoc)	51	73
Personnel	IGF & CJW	IGF, CJW
Notes	~~Tree/plant roots on outside of sleeve~~	Odorous effervescent, decrease in gel-like substance
Retrieval and/or Sampling		
Date / Time	2/13/24 @ 1225	
DTW (ft btoc)	36.84	
Sampler Integrity	Good	
Personnel	IGF, CJW	
Notes	Gel like, slightly effervescent, floating precipitate	

Field Parameters

Date / Time	Water Temp (°C)	pH (SU)	ORP (mV)	Sp Cond (µS/cm)	DO (mg/L)	Turbidity
2/13/24 @ 1233	26.4	5.80	-276.9	1533	0.56	140

Sample ID	TTU-~~13-51~~-73-20240213	
QAQC Samples	—	
Containers	(1) 125 mL HDPE (no pres.) & (6) 40 mL amber VOAs	
Preservatives	HCl	
Analysis	Perchlorate / VOCs / 1,4-Dioxane	
Sampler Reset	(Yes)	No
Notes		

Pinyon Environmental, Inc.

Well Sampling Record	
Project Name	Nammo TTU
Project Number	722152201.002
Well ID / ADWR #	TTU-12 / NA
Date Completed	7/31/2018
Casing Material	Steel
Casing Diameter (in)	5
Screen (ft btoc)	Open to 180
Well Total Depth (ft btoc)	180
Survey Information	Northing: 909105.990 / Easting: 761103.280
Deployment	
Date / Time	11/18/23 at 10:15
Type of Sampler	HydraSleeve
Size of Sampler	HS-2-1L
DTW (ft btoc)	75.62
Deployment Depth (ft btoc)	82
Personnel	IGF & CJW
Notes	HS only ½ full
Retrieval and/or Sampling	
Date / Time	2/12/24 @ 1000
DTW (ft btoc)	74.17
Sampler Integrity	Good
Personnel	IGF, CJW
Notes	3/4 full, 1/2" Sediment @ bottom of HS, Sediment on outside of HS.

Field Parameters

Date / Time	Water Temp (°C)	pH (SU)	ORP (mV)	Sp Cond (µS/cm)	DO (mg/L)	Turbidity
2/12/24 @ 1010	19.3	6.87	-207.7	*Over 3000 last time 3232	4.0	7.9

Sample ID	TTU-12-82-20240212	
QAQC Samples	–	
Containers	(1) 125 mL HDPE (no pres.) & (6) 40 mL amber VOAs	
Preservatives	HCl	
Analysis	Perchlorate / ~~VOCs~~ / 1,4-Dioxane	
Sampler Reset	Yes	No
Notes		

Pinyon Environmental, Inc.

Well Sampling Record	
Project Name	Nammo TTU
Project Number	722152201.002
Well ID / ADWR #	TTU-13 / NA
Date Completed	7/20/2018
Casing Material	Steel
Casing Diameter (in)	5
Screen (ft btoc)	Open to 80
Well Total Depth (ft btoc)	80
Survey Information	Northing: 909405.920 / Easting: 761232.180
Deployment	
Date / Time	11/17/23 at 14:35
Type of Sampler	HydraSleeve
Size of Sampler	HS-2-1L
DTW (ft btoc)	43.27
Deployment Depth (ft btoc)	51
Personnel	IGF & CJW
Notes	Tree/plant roots on outside of sleeve
Retrieval and/or Sampling	
Date / Time	~~0~~ 2-12-24 @ 0905
DTW (ft btoc)	43.70
Sampler Integrity	Good
Personnel	IGF, CJW
Notes	" "

Field Parameters

Date / Time	Water Temp (°C)	pH (SU)	ORP (mV)	Sp Cond (µS/cm)	DO (mg/L)	Turbidity
2/12/24 @ 0915	16.1	6.82	-89.0	1302	3.73	6.8

Sample ID	TTU-13-51-20240212
QAQC Samples	MS/MSD #1
Containers	(2) 125 mL HDPE (no pres.) & (2) 40 mL amber VOAs
Preservatives	HCl
Analysis	Perchlorate / VOCs / 1,4-Dioxane
Sampler Reset	(Yes) / No
Notes	

Pinyon Environmental, Inc.

Well Sampling Record	
Project Name	Nammo TTU
Project Number	722152201.002
Well ID / ADWR #	TTU-14 / NA
Date Completed	7/19/2018
Casing Material	Steel
Casing Diameter (in)	5
Screen (ft btoc)	Open to 100
Well Total Depth (ft btoc)	100
Survey Information	Northing: 909224.260 / Easting: 761181.230
Deployment	
Date / Time	11/18/23 at 09:54
Type of Sampler	HydraSleeve
Size of Sampler	HS-2-1L
DTW (ft btoc)	61.81
Deployment Depth (ft btoc)	64
Personnel	IGF & CJW
Notes	Small amount of rocky sediment
Retrieval and/or Sampling	
Date / Time	2-12-24 @ 0935
DTW (ft btoc)	60.39
Sampler Integrity	Good
Personnel	IGF, CJW
Notes	" "

Field Parameters

Date / Time	Water Temp (°C)	pH (SU)	ORP (mV)	Sp Cond (µS/cm)	DO (mg/L)	Turbidity
2-12-24 @ 0942	18.3	6.86	-73.6	2900	3.22	17

Sample ID	TTU-14-64-20240212
QAQC Samples	~~DUP-~~ CW
Containers	(1) ~~(2)~~ 125 mL HDPE (no pres.) & (6) ~~(10)~~ 40 mL amber VOAs
Preservatives	HCl
Analysis	Perchlorate / VOCs / 1,4-Dioxane
Sampler Reset	Yes / No

Notes

**Be careful, tether not hooked into anything

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Well Sampling Record	
Project Name	Nammo TTU
Project Number	722152201.002
Well ID / ADWR #	TTU-15 / 55-228014
Date Completed	1/25/2018
Casing Material	Steel
Casing Diameter (in)	N/A
Screen (ft btoc)	Open
Well Total Depth (ft btoc)	100
Survey Information	Northing: 909185.100 / Easting: 762065.910
Deployment	
Date / Time	11/18/2023 at 12:25
Type of Sampler	HydraSleeve
Size of Sampler	HS-2-1L
DTW (ft btoc)	32.70
Deployment Depth (ft btoc)	75
Personnel	IGF & CJW
Notes	Slightly cloudy
Retrieval and/or Sampling	
Date / Time	2024-13-2 @ 1302
DTW (ft btoc)	33.82
Sampler Integrity	Good
Personnel	IGF, CJW
Notes	" "

Field Parameters

Date / Time	Water Temp (°C)	pH (SU)	ORP (mV)	Sp Cond (µS/cm)	DO (mg/L)	Turbidity
2-13-24 @ 1308	23.4	7.23	-88.4	2452	0.83	M

Sample ID	TTU-15-75-20240213	
QAQC Samples	—	
Containers	(1) 125 mL HDPE (no pres.) & (6) 40 mL amber VOAs	
Preservatives	HCl	
Analysis	Perchlorate / VOCs / 1,4-Dioxane	
Sampler Reset	Yes	No
Notes		

Pinyon Environmental, Inc.

Well Sampling Record	
Project Name	Nammo TTU
Project Number	722152201.002
Well ID / ADWR #	TTU-16 / 55-231730
Date Completed	1/23/2020
Casing Material	Steel
Casing Diameter (in)	8
Screen (ft btoc)	OPEN
Well Total Depth (ft btoc)	96.6
Survey Information	Northing: 909124.980 / Easting: 761848.851
Deployment	
Date / Time	11/18/2023 at 13:04
Type of Sampler	HydraSleeve
Size of Sampler	HS-2-1L
DTW (ft btoc)	24.67
Deployment Depth (ft btoc)	80
Personnel	IGF & CJW
Notes	Water had red tint; slightly effervescent
Retrieval and/or Sampling	
Date / Time	2/13/24 @ 1325
DTW (ft btoc)	26.24
Sampler Integrity	Good
Personnel	IGF CJW
Notes	2-3" Sediment @ bottom of HS. Slightly less red tint.

Field Parameters

Date / Time	Water Temp (°C)	pH (SU)	ORP (mV)	Sp Cond (µS/cm)	DO (mg/L)	Turbidity
2/13/24 @ 1331	23.7	6.54	-287.9	*Was over 8,000 last time 9,530	0.73	50

Sample ID	TTU-16-80-20240213	
QAQC Samples	—	
Containers	(1) 125 mL HDPE (no pres.) & (6) 40 mL amber VOAs	
Preservatives	HCl	
Analysis	Perchlorate / VOCs / 1,4-Dioxane	
Sampler Reset	(Yes)	No

Notes
Known high concentrations

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Well Sampling Record	
Project Name	Nammo TTU
Project Number	722152201.002
Well ID / ADWR #	TTU-17 / 55-23173
Date Completed	1/22/2020
Casing Material	Steel
Casing Diameter (in)	8
Screen (ft btoc)	OPEN
Well Total Depth (ft btoc)	102
Survey Information	Northing: 909370.903 / Easting: 762179.168
Deployment	
Date / Time	11/18/2023 at 11:58
Type of Sampler	HydraSleeve
Size of Sampler	HS-2-1L
DTW (ft btoc)	38.80
Deployment Depth (ft btoc)	80
Personnel	IGF & HMC
Notes	~2" sediment in bottom of HS, bottom 1/3 cloudy, sulfur smell
Retrieval and/or Sampling	
Date / Time	2-12-24 @ 1035
DTW (ft btoc)	41.19
Sampler Integrity	Good
Personnel	IGF, CJW
Notes	" "sulfur-like smell

Field Parameters

Date / Time	Water Temp (°C)	pH (SU)	ORP (mV)	Sp Cond (µS/cm)	DO (mg/L)	Turbidity
2-12-24 @ 1040	20.5	7.06	-72.9	1083	1.14	50

Sample ID	TTU-17-80-20240212	
QAQC Samples	—	
Containers	(1) 125 mL HDPE (no pres.) & (6) 40 mL amber VOAs	
Preservatives	HCl	
Analysis	Perchlorate / ~~VOCs~~ / 1,4-Dioxane	
Sampler Reset	Yes	No
Notes		

Pinyon Environmental, Inc.

Well Sampling Record	
Project Name	Nammo TTU
Project Number	722152201.002
Well ID / ADWR #	TTU-19 / 55-232969
Date Completed	9/24/2020
Casing Material	PVC
Casing Diameter (in)	4
Screen (ft btoc)	25-90
Well Total Depth (ft btoc)	95
Survey Information	Northing: 909030.750 / Easting: 761687.700
Deployment	
Date / Time	11/18/2023 at 13:44
Type of Sampler	HydraSleeve
Size of Sampler	HS-2-1L
DTW (ft btoc)	32.56
Deployment Depth (ft btoc)	73
Personnel	IGF & CJW
Notes	Much black sed. at bottom of HS; odorous; slightly effervescent
Retrieval and/or Sampling	
Date / Time	2/13/24 @ 1200
DTW (ft btoc)	34.87
Sampler Integrity	Good
Personnel	IGF, CJW
Notes	" "

Field Parameters

Date / Time	Water Temp (°C)	pH (SU)	ORP (mV)	Sp Cond (µS/cm)	DO (mg/L)	Turbidity
2/13/24 @ 1210	24.6	6.61	-162.5	2403	0.46	270

Sample ID	TTU-19-73-20240213	
QAQC Samples	—	
Containers	(1) 125 mL HDPE (no pres.) & (6) 40 mL amber VOAs	
Preservatives	HCl	
Analysis	Perchlorate / VOCs / 1,4-Dioxane	
Sampler Reset	(Yes)	No
Notes		

Pinyon Environmental, Inc.

Well Sampling Record	
Project Name	Nammo TTU
Project Number	722152201.002
Well ID / ADWR #	TTU-20 / 55-232968
Date Completed	9/24/2020
Casing Material	PVC
Casing Diameter (in)	4
Screen (ft btoc)	25-90
Well Total Depth (ft btoc)	95
Survey Information	Northing: 909022.530 / Easting: 761681.990
Deployment	
Date / Time	—
Type of Sampler	Dedicated pump
Size of Sampler	HS-2-1L
DTW (ft btoc)	33.65 on 11/17/23
Deployment Depth (ft btoc)	73
Personnel	IGF & CJW
Notes	
Retrieval and/or Sampling	
Date / Time	2/13/24 @ 1140
DTW (ft btoc)	38.06
Sampler Integrity	Spigot
Personnel	IGF, CJW
Notes	Water clear

Field Parameters

Date / Time	Water Temp (°C)	pH (SU)	ORP (mV)	Sp Cond (µS/cm)	DO (mg/L)	Turbidity
2/13/24 1122	22.9	7.08	-233.9	3047	1.17	2.6
1127	24.3	7.03	-251.9	2291	1.32	1.9
1132	24.1	6.98	-183.6	2228	2.15	2.6

Sample ID	TTU-20-73- 20240213
QAQC Samples	—
Containers	(1) 125 mL HDPE (no pres.) & (6) 40 mL amber VOAs
Preservatives	HCl
Analysis	Perchlorate / VOCs / 1,4-Dioxane
Sampler Reset	Yes / (No)
Notes	*Known high concentrations*

Pinyon Environmental, Inc.

Well Sampling Record	
Project Name	Nammo TTU
Project Number	722152201.002
Well ID / ADWR #	TTU-EX-1 / 55-231733
Date Completed	1/29/2020
Casing Material	
Casing Diameter (in)	8
Screen (ft btoc)	OPEN
Well Total Depth (ft btoc)	109
Survey Information	Northing: 909350.574 / Easting: 761597.823
Deployment	
Date / Time	11/18/2023 at 11:32
Type of Sampler	HydraSleeve
Size of Sampler	HS-2-1L
DTW (ft btoc)	29.27
Deployment Depth (ft btoc)	69
Personnel	IGF & CJW
Notes	Bottom ½ of HS cloudy, ~1" silt/sediment at bottom
Retrieval and/or Sampling	
Date / Time	2/13/24 @ 1343
DTW (ft btoc)	30.80
Sampler Integrity	Good
Personnel	IGF, CJW
Notes	

Field Parameters

Date / Time	Water Temp (°C)	pH (SU)	ORP (mV)	Sp Cond (µS/cm)	DO (mg/L)	Turbidity
2/13/24 @ 1349	23.9	7.17	-191.8	2862	2.44	14

Sample ID	TTU-EX-1-69-20240213	
QAQC Samples	—	
Containers	(1) 125 mL HDPE (no pres.) & (6) 40 mL amber VOAs	
Preservatives	HCl	
Analysis	Perchlorate / VOCs / 1,4-Dioxane	
Sampler Reset	Yes	No
Notes		

Pinyon Environmental, Inc.

Well Sampling Record	
Project Name	Nammo TTU
Project Number	722152201.002
Well ID / ADWR #	TTU-EX-2 / 55-231734
Date Completed	1/28/2020
Casing Material	
Casing Diameter (in)	8
Screen (ft btoc)	OPEN
Well Total Depth (ft btoc)	110
Survey Information	Northing: 909268.187 / Easting: 761493.214
Deployment	
Date / Time	11/18/2023 at 11:12
Type of Sampler	HydraSleeve
Size of Sampler	HS-2-1L
DTW (ft btoc)	38.04
Deployment Depth (ft btoc)	74
Personnel	IGF & CJW
Notes	Bottom ½ of sleeve cloudy
Retrieval and/or Sampling	
Date / Time	2-13-24 @ 1359
DTW (ft btoc)	39.87
Sampler Integrity	GOOD
Personnel	IGF, CJW
Notes	"...slightly cloudy"

Field Parameters

Date / Time	Water Temp (°C)	pH (SU)	ORP (mV)	Sp Cond (µS/cm)	DO (mg/L)	Turbidity
2-13-24 @ 1405	23.8	7.20	-202.6	2116	0.90	33

Sample ID	TTU-EX-2-74-20240213	
QAQC Samples	—	
Containers	(1) 125 mL HDPE (no pres.) & (6) 40 mL amber VOAs	
Preservatives	HCl	
Analysis	Perchlorate / VOCs / 1,4-Dioxane	
Sampler Reset	Yes	No
Notes		

Pinyon Environmental, Inc.

Well Sampling Record	
Project Name	Nammo TTU
Project Number	722152201.002
Well ID / ADWR #	TTU-EX-3 / 55-231731
Date Completed	1/24/2020
Casing Material	
Casing Diameter (in)	8
Screen (ft btoc)	OPEN
Well Total Depth (ft btoc)	111
Survey Information	Northing: 909134.941 / Easting: 761465.507
Deployment	
Date / Time	11/18/2023 at 10:56
Type of Sampler	HydraSleeve
Size of Sampler	HS-2-1L
DTW (ft btoc)	38.75
Deployment Depth (ft btoc)	76
Personnel	IGF & HMC
Notes	Rocky sediment on & in bag
Retrieval and/or Sampling	
Date / Time	2/12/24 @ 1100
DTW (ft btoc)	41.78
Sampler Integrity	GOOD
Personnel	IGF, CJW
Notes	" "

Field Parameters

Date / Time	Water Temp (°C)	pH (SU)	ORP (mV)	Sp Cond (µS/cm)	DO (mg/L)	Turbidity
2/12/24 @ 1107	20.5	6.65	-229.0	5661	1.35	12

Sample ID	TTU-EX-3-76-20240212	
QAQC Samples	—	
Containers	(1) 125 mL HDPE (no pres.) & ((6)) 40 mL amber VOAs	
Preservatives	HCl	
Analysis	Perchlorate / VOCs / 1,4-Dioxane	
Sampler Reset	(Yes)	No
Notes		

Pinyon Environmental, Inc.

Well Sampling Record	
Project Name	Nammo TTU
Project Number	722152201.002
Well ID / ADWR #	TTU-EX-4 / 55-231732
Date Completed	1/25/2020
Casing Material	
Casing Diameter (in)	8
Screen (ft btoc)	OPEN
Well Total Depth (ft btoc)	112
Survey Information	Northing: 909051.298 / Easting: 761442.876
Deployment	
Date / Time	11/18/2023 at 10:38
Type of Sampler	HydraSleeve
Size of Sampler	HS-2-1L
DTW (ft btoc)	42.98
Deployment Depth (ft btoc)	77
Personnel	IGF & CJW
Notes	Partially cloudy bottom ½
Retrieval and/or Sampling	
Date / Time	2/12/24 @ 1113
DTW (ft btoc)	44.74
Sampler Integrity	Good
Personnel	IGF. CJW
Notes	" "

Field Parameters

Date / Time	Water Temp (°C)	pH (SU)	ORP (mV)	Sp Cond (µS/cm)	DO (mg/L)	Turbidity
2/12/24 @ 1118	20.4	6.98	-175.0	2157	1.11	55

Sample ID	TTU-EX-4-77-20240212	
QAQC Samples	—	
Containers	(1) 125 mL HDPE (no pres.) & (6) 40 mL amber VOAs	
Preservatives	HCl	
Analysis	Perchlorate / ~~VOCs~~ / 1,4-Dioxane	
Sampler Reset	Yes	No
Notes		

Pinyon Environmental, Inc

Well Sampling Record	
Project Name	Nammo TTU
Project Number	722152201.002
Well ID / ADWR #	TTU-EX-5 / 55-231736
Date Completed	1/24/2020
Casing Material	
Casing Diameter (in)	8
Screen (ft btoc)	OPEN
Well Total Depth (ft btoc)	112.4
Survey Information	Northing: 908971.770 / Easting: 761423.325
Deployment	
Date / Time	11/17/2023 at 15:07
Type of Sampler	HydraSleeve
Size of Sampler	HS-2-1L
DTW (ft btoc)	40.71
Deployment Depth (ft btoc)	80
Personnel	IGF & CJW
Notes	Cloudy towards bottom of HS
Retrieval and/or Sampling	
Date / Time	2024-2-13 @ 1415
DTW (ft btoc)	41.22
Sampler Integrity	Good
Personnel	IGF, CJW
Notes	—

Field Parameters

Date / Time	Water Temp (°C)	pH (SU)	ORP (mV)	Sp Cond (µS/cm)	DO (mg/L)	Turbidity
2/13/24 @ 1425	24.4	7.27	-282.9	1162	0.85	26

Sample ID	TTU-EX-5-80-20240213	
QAQC Samples	DUP-02 ✓	
Containers	(2) 125 mL HDPE (no pres.) & (12) 40 mL amber VOAs	
Preservatives	HCl	
Analysis	Perchlorate / ~~VOCs~~ / 1,4-Dioxane	
Sampler Reset	Yes	No
Notes		

Attachment 2 – Laboratory Analytical Reports

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ANALYTICAL REPORT

PREPARED FOR

Attn: Andrew Parker
Pinyon Environmental Inc
1783 W University Drive
Suite 137
Tempe, Arizona 85281

Generated 3/29/2024 11:43:06 AM Revision 2

JOB DESCRIPTION

Nammo TTU

JOB NUMBER

550-214412-1

Eurofins Phoenix
4625 East Cotton Center Boulevard
Suite #189
Phoenix AZ 85040

See page two for job notes and contact information.

Job Notes

This report may not be reproduced except in full, and with written approval from the laboratory. The results relate only to the samples tested. For questions please contact the Project Manager at the e-mail address or telephone number listed on this page.

The test results in this report relate only to the samples as received by the laboratory and will meet all requirements of the methodology, with any exceptions noted. This report shall not be reproduced except in full, without the express written approval of the laboratory. All questions should be directed to the Eurofins Environment Testing Southwest, LLC Project Manager.

Authorization

Generated
3/29/2024 11:43:06 AM
Revision 2

Authorized for release by
Rachel Sester, Project Manager I
Rachel.Sester@et.eurofinsus.com
(602)659-7615

Table of Contents

Definitions/Glossary

Client: Pinyon Environmental Inc
Project/Site: Nammo TTU

Job ID: 550-214412-1

Qualifiers

HPLC/IC

Qualifier	Qualifier Description
E4	Concentration estimated. Analyte was detected below laboratory minimum reporting level (MRL) but above MDL.
E8	Analyte reported to MDL per project specification. Target analyte was not detected in the sample.

Glossary

Abbreviation	These commonly used abbreviations may or may not be present in this report.
¤	Listed under the "D" column to designate that the result is reported on a dry weight basis
%R	Percent Recovery
CFL	Contains Free Liquid
CFU	Colony Forming Unit
CNF	Contains No Free Liquid
DER	Duplicate Error Ratio (normalized absolute difference)
Dil Fac	Dilution Factor
DL	Detection Limit (DoD/DOE)
DL, RA, RE, IN	Indicates a Dilution, Re-analysis, Re-extraction, or additional Initial metals/anion analysis of the sample
DLC	Decision Level Concentration (Radiochemistry)
EDL	Estimated Detection Limit (Dioxin)
LOD	Limit of Detection (DoD/DOE)
LOQ	Limit of Quantitation (DoD/DOE)
MCL	EPA recommended "Maximum Contaminant Level"
MDA	Minimum Detectable Activity (Radiochemistry)
MDC	Minimum Detectable Concentration (Radiochemistry)
MDL	Method Detection Limit
ML	Minimum Level (Dioxin)
MPN	Most Probable Number
MQL	Method Quantitation Limit
NC	Not Calculated
ND	Not Detected at the reporting limit (or MDL or EDL if shown)
NEG	Negative / Absent
POS	Positive / Present
PQL	Practical Quantitation Limit
PRES	Presumptive
QC	Quality Control
RER	Relative Error Ratio (Radiochemistry)
RL	Reporting Limit or Requested Limit (Radiochemistry)
RPD	Relative Percent Difference, a measure of the relative difference between two points
TEF	Toxicity Equivalent Factor (Dioxin)
TEQ	Toxicity Equivalent Quotient (Dioxin)
TNTC	Too Numerous To Count

Job ID: 550-214412-1 Eurofins Phoenix

Job Narrative
550-214412-1

Analytical test results meet all requirements of the associated regulatory program listed on the Accreditation/Certification Summary Page unless otherwise noted under the individual analysis. Data qualifiers are applied to indicate exceptions. Noncompliant quality control (QC) is further explained in narrative comments.

- Matrix QC may not be reported if insufficient sample or site-specific QC samples were not submitted. In these situations, to demonstrate precision and accuracy at a batch level, a LCS/LCSD may be performed, unless otherwise specified in the method.
- Surrogate and/or isotope dilution analyte recoveries (if applicable) which are outside of the QC window are confirmed unless attributed to a dilution or otherwise noted in the narrative.

Regulated compliance samples (e.g. SDWA, NPDES) must comply with the associated agency requirements/permits.

Revision 2
This report was revised on 3/97/24 to update the certification matrix to potable water. This final report replaces the report that was generated on 2/27/24 at 2:59 PM.

Revision
This report was revised on 3/27/24 to change the name of the project to "Nammo TTU", per client request. This final report replaces the report that was generated on 2/26/24 at 11:35 AM.

Receipt
The sample was received on 2/15/2024 9:35 AM. Unless otherwise noted below, the sample arrived in good condition, and, where required, properly preserved and on ice. The temperature of the cooler at receipt time was 5.1°C.

HPLC/IC
No additional analytical or quality issues were noted, other than those described above or in the Definitions/ Glossary page.

Sample Summary

Client: Pinyon Environmental Inc
Project/Site: Nammo TTU

Job ID: 550-214412-1

Lab Sample ID	Client Sample ID	Matrix	Collected	Received
550-214412-1	PF-2-400-20240214	Water	02/14/24 00:00	02/15/24 09:35

Detection Summary

Client: Pinyon Environmental Inc
Project/Site: Nammo TTU

Job ID: 550-214412-1

Client Sample ID: PF-2-400-20240214 **Lab Sample ID: 550-214412-1**

No Detections.

This Detection Summary does not include radiochemical test results.

Client Sample Results

Client: Pinyon Environmental Inc
Project/Site: Nammo TTU

Job ID: 550-214412-1

Client Sample ID: PF-2-400-20240214
Date Collected: 02/14/24 00:00
Date Received: 02/15/24 09:35

Lab Sample ID: 550-214412-1
Matrix: Water

Method: EPA 314.0 - Perchlorate (IC)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perchlorate	ND	E8	1.0	0.31	ug/L			02/23/24 21:30	1

Method: 314.0 - Perchlorate (IC)

Lab Sample ID: MB 550-316351/1002
Matrix: Water
Analysis Batch: 316351

Client Sample ID: Method Blank
Prep Type: Total/NA

Analyte	MB Result	MB Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perchlorate	ND	E8	1.0	0.31	ug/L			02/23/24 12:20	1

Lab Sample ID: LCS 550-316351/4
Matrix: Water
Analysis Batch: 316351

Client Sample ID: Lab Control Sample
Prep Type: Total/NA

Analyte			Spike Added	LCS Result	LCS Qualifier	Unit	D	%Rec	%Rec Limits		
Perchlorate			25.0	23.4		ug/L		93	85 - 115		

Lab Sample ID: LCSD 550-316351/5
Matrix: Water
Analysis Batch: 316351

Client Sample ID: Lab Control Sample Dup
Prep Type: Total/NA

Analyte			Spike Added	LCSD Result	LCSD Qualifier	Unit	D	%Rec	%Rec Limits	RPD	RPD Limit
Perchlorate			25.0	23.4		ug/L		94	85 - 115	0	15

Lab Sample ID: MRL 550-316351/1003
Matrix: Water
Analysis Batch: 316351

Client Sample ID: Lab Control Sample
Prep Type: Total/NA

Analyte			Spike Added	MRL Result	MRL Qualifier	Unit	D	%Rec	%Rec Limits		
Perchlorate			1.00	0.836	E4	ug/L		84	75 - 125		

Lab Sample ID: 550-214272-A-6 MS ^100
Matrix: Water
Analysis Batch: 316351

Client Sample ID: Matrix Spike
Prep Type: Total/NA

Analyte	Sample Result	Sample Qualifier	Spike Added	MS Result	MS Qualifier	Unit	D	%Rec	%Rec Limits		
Perchlorate	1100		2500	3200		ug/L		85	80 - 120		

Lab Sample ID: 550-214272-A-6 MSD ^100
Matrix: Water
Analysis Batch: 316351

Client Sample ID: Matrix Spike Duplicate
Prep Type: Total/NA

Analyte	Sample Result	Sample Qualifier	Spike Added	MSD Result	MSD Qualifier	Unit	D	%Rec	%Rec Limits	RPD	RPD Limit
Perchlorate	1100		2500	3190		ug/L		85	80 - 120	0	15

QC Association Summary

Client: Pinyon Environmental Inc
Project/Site: Nammo TTU

Job ID: 550-214412-1

HPLC/IC

Analysis Batch: 316351

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
550-214412-1	PF-2-400-20240214	Total/NA	Water	314.0	
MB 550-316351/1002	Method Blank	Total/NA	Water	314.0	
LCS 550-316351/4	Lab Control Sample	Total/NA	Water	314.0	
LCSD 550-316351/5	Lab Control Sample Dup	Total/NA	Water	314.0	
MRL 550-316351/1003	Lab Control Sample	Total/NA	Water	314.0	
550-214272-A-6 MS ^100	Matrix Spike	Total/NA	Water	314.0	
550-214272-A-6 MSD ^100	Matrix Spike Duplicate	Total/NA	Water	314.0	

Lab Chronicle

Client: Pinyon Environmental Inc
Project/Site: Nammo TTU

Job ID: 550-214412-1

Client Sample ID: PF-2-400-20240214

Date Collected: 02/14/24 00:00
Date Received: 02/15/24 09:35

Lab Sample ID: 550-214412-1

Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Analyst	Lab	Prepared or Analyzed
Total/NA	Analysis	314.0		1	316351	RDC	EET PHX	02/23/24 21:30

Laboratory References:

EET PHX = Eurofins Phoenix, 4625 East Cotton Center Boulevard, Suite #189, Phoenix, AZ 85040, TEL (602)437-3340

Accreditation/Certification Summary

Client: Pinyon Environmental Inc
Project/Site: Nammo TTU

Job ID: 550-214412-1

Laboratory: Eurofins Phoenix

The accreditations/certifications listed below are applicable to this report.

Authority	Program	Identification Number	Expiration Date
Arizona	State	AZ0728	06-10-24

Method Summary

Client: Pinyon Environmental Inc
Project/Site: Nammo TTU

Job ID: 550-214412-1

Method	Method Description	Protocol	Laboratory
314.0	Perchlorate (IC)	EPA	EET PHX

Protocol References:

EPA = US Environmental Protection Agency

Laboratory References:

EET PHX = Eurofins Phoenix, 4625 East Cotton Center Boulevard, Suite #189, Phoenix, AZ 85040, TEL (602)437-3340

Eurofins Phoenix
4625 East Cotton Center Boulevard
Suite 189
Phoenix, AZ 85040-4807
phone 602.437.3340

Chain of Custody Record

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2144122

3/29/2024 (Rev. 2)

Regulatory Program: ☐DW ☐NPDES ☐RCRA ☒Other: EPA

Eurofins Environment Testing America

Client Contact	**Project Manager:** Andrew Parker	**Site Contact:** –	**Date:**
Company Name: Pinyon Environmental	**Email:** parker@pinyon-env.com	**Lab Contact:** Jeremy Russell	**Carrier:**
Address: 1783 W. University Dr. Suite 137	**Tel/Fax:** (720)614-5032	**Analysis Turnaround Time**	
City/State/Zip: Tempe, AZ 85281	☐ CALENDAR DAYS ☐ WORKING DAYS		
Phone: (602)274-0533	TAT if different from Below _____		
FAX: –	☐ 2 weeks ☐ 1 week ☐ 2 days ☐ 1 day	Standard TAT	
Project Name: NDS TTU			
Site: –			
PO# 72152201.002			

COC No:	Sampler: Isabella Foster
1 of 1 COCs	**For Lab Use Only:**
TALS Project #:	Walk-in Client:
	Lab Sampling:
	Job / SDG No.:

Sample Identification	Sample Date	Sample Time	Sample Type (C=Comp, G=Grab)	Matrix	# of Cont.	Filtered Sample (Y/N)	Perform MS/MSD (Y/N)	Pecnbrate 314.0	Sample Specific Notes:
PF-2-4000-20240214 -1	2/14/24	~~DW~~ 1617	C	DW	2	N	Y	X	

550-214412 Chain of Custody

Preservation Used: 1= Ice, 2= HCl; 3= H_2SO_4; 4=HNO_3; 5=NaOH; 6= Other _____

Possible Hazard Identification:
Are any samples from a listed EPA Hazardous Waste? Please List any EPA Waste Codes for the sample in the Comments Section if the lab is to dispose of the sample.

☒Non-Hazard ☐Flammable ☐Skin Irritant ☐Poison B ☐Unknown

Sample Disposal (A fee may be assessed if samples are retained longer than 1 month)

☐Return to Client ☒Disposal by Lab ☐Archive for _____ Months

Special Instructions/QC Requirements & Comments:

5.1°C/ICE/cpe

Custody Seals Intact: ☐ Yes ☐ No	Custody Seal No.:		Cooler Temp. (°C): Obs'd: _____ Corr'd: _____ Therm ID No.: _____		
Relinquished by:	Company: Pinyon	Date/Time: 2/15/24 0915	Received by:	Company:	Date/Time:
Relinquished by:	Company: Pinyon	Date/Time: 2/15/24 0935	Received by:	Company:	Date/Time:
Relinquished by:	Company:	Date/Time:	Received in Laboratory by:	Company: EETA-PHX	Date/Time: 2/15/24 0935

Login Sample Receipt Checklist

Client: Pinyon Environmental Inc

Job Number: 550-214412-1

Login Number: 214412
List Number: 1
Creator: Maycock, Lisa

List Source: Eurofins Phoenix

Question	Answer	Comment
Radioactivity wasn't checked or is </= background as measured by a survey meter.	True	
The cooler's custody seal, if present, is intact.	True	
Sample custody seals, if present, are intact.	True	
The cooler or samples do not appear to have been compromised or tampered with.	True	
Samples were received on ice.	True	
Cooler Temperature is acceptable.	True	
Cooler Temperature is recorded.	True	
COC is present.	True	
COC is filled out in ink and legible.	True	
COC is filled out with all pertinent information.	True	
Is the Field Sampler's name present on COC?	True	
There are no discrepancies between the containers received and the COC.	True	
Samples are received within Holding Time (excluding tests with immediate HTs)	True	
Sample containers have legible labels.	True	
Containers are not broken or leaking.	True	
Sample collection date/times are provided.	True	
Appropriate sample containers are used.	True	
Sample bottles are completely filled.	True	
Sample Preservation Verified.	True	
There is sufficient vol. for all requested analyses, incl. any requested MS/MSDs	True	
Containers requiring zero headspace have no headspace or bubble is <6mm (1/4").	True	
Multiphasic samples are not present.	True	
Samples do not require splitting or compositing.	True	
Residual Chlorine Checked.	False	Check done at department level as required.

ANALYTICAL REPORT

March 22, 2024

Revised Report

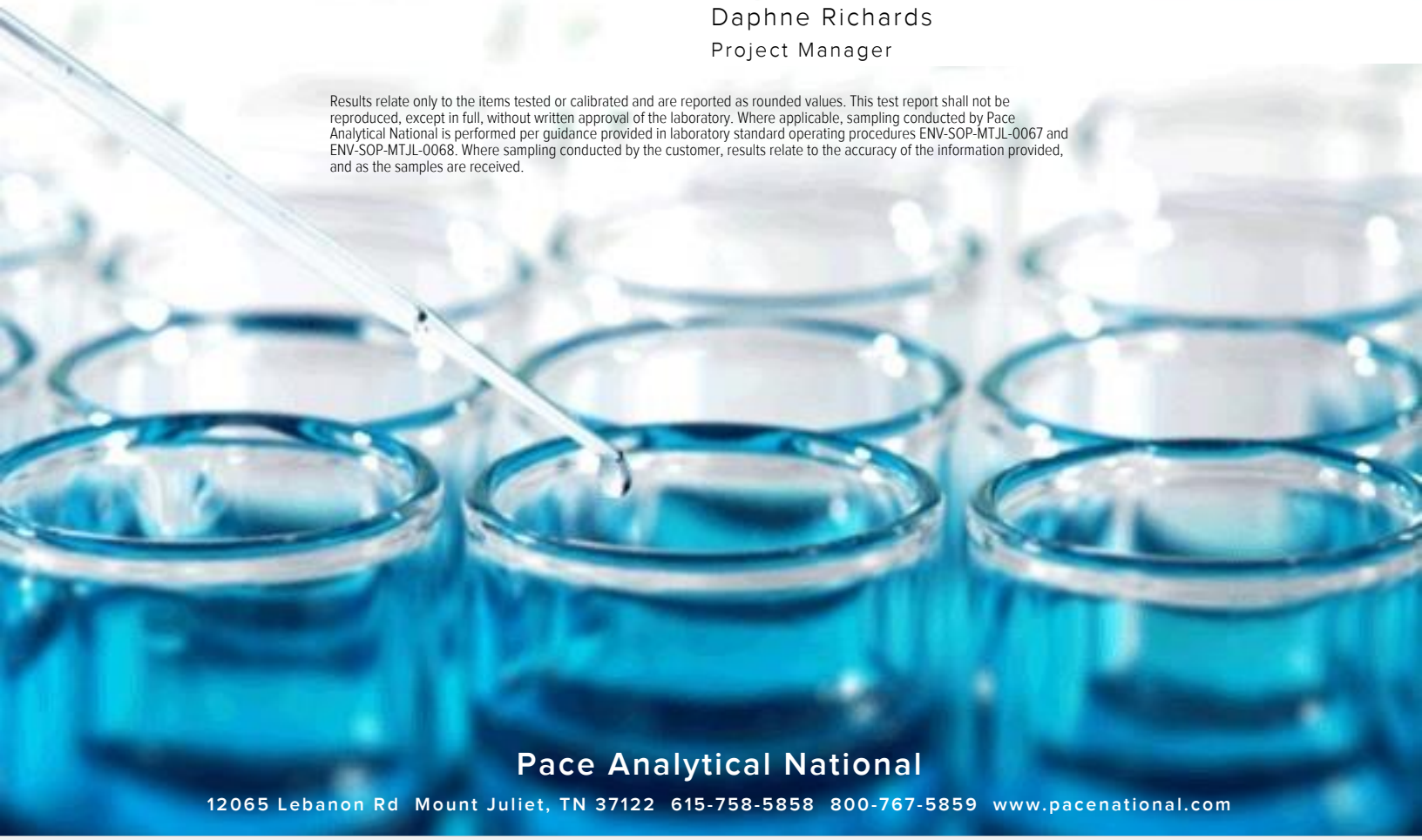
Pinyon Environmental

Sample Delivery Group:	L1705776
Samples Received:	02/14/2024
Project Number:	722152201.002
Description:	Nammo TTU Groundwater Monitoring

Report To: Isbella Foster and Andrew Parker
3222 S. Vance Street Suite 200
Lakewood, CO 80227

Entire Report Reviewed By: Daphne R Richards

Daphne Richards
Project Manager



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SAMPLE SUMMARY

TTU-1-50-20240213 L1705776-01 GW

Collected by: Isabella Foster | Collected date/time: 02/13/24 10:27 | Received date/time: 02/14/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2228552	500	02/24/24 23:29	02/24/24 23:29	JEA	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2228368	1	02/17/24 11:32	02/17/24 11:32	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2227341	1	02/16/24 04:38	02/16/24 04:38	DWR	Mt. Juliet, TN

TTU-2-114-20240213 L1705776-02 GW

Collected by: Isabella Foster | Collected date/time: 02/13/24 11:01 | Received date/time: 02/14/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2228552	5000	02/21/24 01:26	02/21/24 01:26	JEA	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2228368	10	02/17/24 15:18	02/17/24 15:18	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2229520	10	02/20/24 01:23	02/20/24 01:23	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2227341	1	02/16/24 05:00	02/16/24 05:00	DWR	Mt. Juliet, TN

TTU-3-108-20240212 L1705776-03 GW

Collected by: Isabella Foster | Collected date/time: 02/12/24 15:18 | Received date/time: 02/14/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2228552	1	02/21/24 01:54	02/21/24 01:54	JEA	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2228368	1	02/17/24 11:52	02/17/24 11:52	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2227341	1	02/16/24 05:22	02/16/24 05:22	DWR	Mt. Juliet, TN

TTU-4-57-20240213 L1705776-04 GW

Collected by: Isabella Foster | Collected date/time: 02/13/24 15:28 | Received date/time: 02/14/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2228552	1	02/21/24 02:22	02/21/24 02:22	JEA	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2228368	1	02/17/24 12:13	02/17/24 12:13	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2227341	1	02/16/24 05:43	02/16/24 05:43	DWR	Mt. Juliet, TN

TTU-5-110-20240213 L1705776-05 GW

Collected by: Isabella Foster | Collected date/time: 02/13/24 14:50 | Received date/time: 02/14/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2228552	1	02/21/24 02:50	02/21/24 02:50	JEA	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2228368	1	02/17/24 12:33	02/17/24 12:33	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2227341	1	02/16/24 06:05	02/16/24 06:05	DWR	Mt. Juliet, TN

TTU-6-143-20240212 L1705776-06 GW

Collected by: Isabella Foster | Collected date/time: 02/12/24 14:58 | Received date/time: 02/14/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2228552	1	02/21/24 04:14	02/21/24 04:14	JEA	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2228368	1	02/17/24 12:54	02/17/24 12:54	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2227341	1	02/16/24 06:27	02/16/24 06:27	DWR	Mt. Juliet, TN

TTU-7-345-20240212 L1705776-07 GW

Collected by: Isabella Foster | Collected date/time: 02/12/24 14:35 | Received date/time: 02/14/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2228552	1	02/21/24 04:41	02/21/24 04:41	JEA	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2228368	1	02/17/24 13:15	02/17/24 13:15	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2227341	1	02/16/24 06:49	02/16/24 06:49	DWR	Mt. Juliet, TN

TTU-9A-61-20240213 L1705776-08 GW

Collected by: Isabella Foster | Collected date/time: 02/13/24 09:32 | Received date/time: 02/14/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2228552	1	02/21/24 05:09	02/21/24 05:09	JEA	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2228368	1	02/17/24 13:35	02/17/24 13:35	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2230157	1	02/20/24 16:25	02/20/24 16:25	JBE	Mt. Juliet, TN

TTU-11-73-20240213 L1705776-09 GW

Collected by: Isabella Foster | Collected date/time: 02/13/24 12:25 | Received date/time: 02/14/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2228552	1	02/21/24 06:05	02/21/24 06:05	JEA	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2229520	10	02/20/24 01:47	02/20/24 01:47	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2230157	1	02/20/24 16:47	02/20/24 16:47	JBE	Mt. Juliet, TN

TTU-12-82-20240212 L1705776-10 GW

Collected by: Isabella Foster | Collected date/time: 02/12/24 10:00 | Received date/time: 02/14/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2228552	5000	02/21/24 06:33	02/21/24 06:33	JEA	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2228368	10	02/17/24 15:59	02/17/24 15:59	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2229520	10	02/20/24 02:10	02/20/24 02:10	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2230157	1	02/20/24 17:09	02/20/24 17:09	JBE	Mt. Juliet, TN

TTU-13-51-20240212 L1705776-11 GW

Collected by: Isabella Foster | Collected date/time: 02/12/24 09:05 | Received date/time: 02/14/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2228552	1000	02/21/24 07:01	02/21/24 07:01	JEA	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2228368	1	02/17/24 13:56	02/17/24 13:56	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2230157	1	02/20/24 17:31	02/20/24 17:31	JBE	Mt. Juliet, TN

TTU-14-64-20240212 L1705776-12 GW

Collected by: Isabella Foster | Collected date/time: 02/12/24 09:35 | Received date/time: 02/14/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2228552	5000	02/25/24 21:43	02/25/24 21:43	JEA	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2228368	10	02/17/24 16:20	02/17/24 16:20	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2229520	10	02/20/24 02:33	02/20/24 02:33	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2230157	1	02/20/24 17:53	02/20/24 17:53	JBE	Mt. Juliet, TN

TTU-15-75-20240213 L1705776-13 GW

Collected by: Isabella Foster | Collected date/time: 02/13/24 13:02 | Received date/time: 02/14/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2228552	10000	02/25/24 00:24	02/25/24 00:24	JEA	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2228368	1	02/17/24 14:16	02/17/24 14:16	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2229520	2500	02/20/24 02:56	02/20/24 02:56	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2230690	20	02/21/24 15:55	02/21/24 15:55	JBE	Mt. Juliet, TN

TTU-16-80-20240213 L1705776-14 GW

Collected by: Isabella Foster | Collected date/time: 02/13/24 13:25 | Received date/time: 02/14/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2228552	500	02/25/24 00:52	02/25/24 00:52	JEA	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2229520	20	02/20/24 03:19	02/20/24 03:19	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2230690	1	02/21/24 14:29	02/21/24 14:29	JBE	Mt. Juliet, TN

TTU-17-80-20240212 L1705776-15 GW

Collected by: Isabella Foster | Collected date/time: 02/12/24 10:35 | Received date/time: 02/14/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2228552	1	02/25/24 01:20	02/25/24 01:20	JEA	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2228368	1	02/17/24 14:37	02/17/24 14:37	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2229520	1	02/20/24 03:42	02/20/24 03:42	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2230690	1	02/21/24 14:50	02/21/24 14:50	JBE	Mt. Juliet, TN

TTU-19-73-20240213 L1705776-16 GW

Collected by: Isabella Foster | Collected date/time: 02/13/24 12:00 | Received date/time: 02/14/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2228552	1	02/21/24 10:44	02/21/24 10:44	JEA	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2228368	5	02/17/24 17:01	02/17/24 17:01	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2229520	5	02/20/24 04:06	02/20/24 04:06	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2230157	5	02/20/24 21:53	02/20/24 21:53	JBE	Mt. Juliet, TN

TTU-20-73-20240213 L1705776-17 GW

Collected by: Isabella Foster | Collected date/time: 02/13/24 11:40 | Received date/time: 02/14/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2228552	1000	02/25/24 02:44	02/25/24 02:44	JEA	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2228368	10	02/17/24 17:21	02/17/24 17:21	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2229520	10	02/20/24 04:29	02/20/24 04:29	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2230157	1	02/20/24 18:58	02/20/24 18:58	JBE	Mt. Juliet, TN

TTU-EX-5-80-20240213 L1705776-18 GW

Collected by: Isabella Foster | Collected date/time: 02/13/24 14:15 | Received date/time: 02/14/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2228552	1	02/21/24 11:40	02/21/24 11:40	JEA	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2228368	1	02/17/24 14:57	02/17/24 14:57	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2229520	1	02/20/24 04:52	02/20/24 04:52	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2230157	1	02/20/24 19:20	02/20/24 19:20	JBE	Mt. Juliet, TN

TTU-EX-1-69-20240213 L1705776-19 GW

Collected by: Isabella Foster | Collected date/time: 02/13/24 13:43 | Received date/time: 02/14/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2228552	1000	02/25/24 03:12	02/25/24 03:12	JEA	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2228368	10	02/17/24 17:42	02/17/24 17:42	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2229520	10	02/20/24 05:15	02/20/24 05:15	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2230157	10	02/20/24 22:15	02/20/24 22:15	JBE	Mt. Juliet, TN

TTU-EX-2-74-20240213 L1705776-20 GW

Collected by: Isabella Foster | Collected date/time: 02/13/24 13:59 | Received date/time: 02/14/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2228552	1000	02/25/24 03:40	02/25/24 03:40	JEA	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2228368	20	02/17/24 18:02	02/17/24 18:02	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2229520	20	02/20/24 05:38	02/20/24 05:38	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2230690	1	02/21/24 15:12	02/21/24 15:12	JBE	Mt. Juliet, TN

TTU-EX-3-76-20240212 L1705776-21 GW

Collected by: Isabella Foster | Collected date/time: 02/12/24 11:00 | Received date/time: 02/14/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2228553	5000	02/29/24 17:10	02/29/24 17:10	ASM	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2228641	250	02/18/24 04:48	02/18/24 04:48	DYW	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2231600	10	02/22/24 13:45	02/22/24 13:45	JBE	Mt. Juliet, TN

TTU-EX-4-77-20240212 L1705776-22 GW

Collected by: Isabella Foster | Collected date/time: 02/12/24 11:13 | Received date/time: 02/14/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2228553	5000	02/29/24 17:38	02/29/24 17:38	ASM	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2228641	25	02/18/24 05:11	02/18/24 05:11	DYW	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2230157	1	02/20/24 19:42	02/20/24 19:42	JBE	Mt. Juliet, TN

DUP-01 L1705776-23 GW

Collected by: Isabella Foster | Collected date/time: 02/13/24 00:00 | Received date/time: 02/14/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2228553	1	02/29/24 18:06	02/29/24 18:06	ASM	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2228641	1	02/18/24 02:05	02/18/24 02:05	DYW	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2230157	1	02/20/24 20:04	02/20/24 20:04	JBE	Mt. Juliet, TN

DUP-02 L1705776-24 GW

Collected by: Isabella Foster | Collected date/time: 02/13/24 00:00 | Received date/time: 02/14/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2228553	1	02/29/24 18:34	02/29/24 18:34	ASM	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2228641	1	02/18/24 02:28	02/18/24 02:28	DYW	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2230157	1	02/20/24 20:26	02/20/24 20:26	JBE	Mt. Juliet, TN

SAMPLE SUMMARY

TRIP BLANK L1705776-25 GW

Collected by: Isabella Foster
Collected date/time: 02/13/24 00:00
Received date/time: 02/14/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2228641	1	02/17/24 23:23	02/17/24 23:23	DYW	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2230157	1	02/20/24 16:03	02/20/24 16:03	JBE	Mt. Juliet, TN

CASE NARRATIVE

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

Daphne Richards
Project Manager

Report Revision History

Level II Report - Version 1: 03/06/24 14:18

Project Narrative

Sample id correction

Sample Delivery Group (SDG) Narrative

Insufficient sample volume to perform MS/MSD analyses per method QC requirements.

Lab Sample ID	Project Sample ID	Method
L1705776-02	TTU-2-114-20240213	8260B
L1705776-09	TTU-11-73-20240213	8260B
L1705776-10	TTU-12-82-20240212	8260B
L1705776-12	TTU-14-64-20240212	8260B
L1705776-13	TTU-15-75-20240213	8260B, 8260B-SIM
L1705776-14	TTU-16-80-20240213	8260B, 8260B-SIM
L1705776-15	TTU-17-80-20240212	8260B, 8260B-SIM
L1705776-16	TTU-19-73-20240213	8260B
L1705776-17	TTU-20-73-20240213	8260B
L1705776-18	TTU-EX-5-80-20240213	8260B
L1705776-19	TTU-EX-1-69-20240213	8260B
L1705776-20	TTU-EX-2-74-20240213	8260B, 8260B-SIM

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	11700		150	2000	500	02/24/2024 23:29	WG2228552

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		11.3	50.0	1	02/17/2024 11:32	WG2228368
Acrolein	U		2.54	50.0	1	02/17/2024 11:32	WG2228368
Acrylonitrile	U		0.671	10.0	1	02/17/2024 11:32	WG2228368
Benzene	U		0.0941	1.00	1	02/17/2024 11:32	WG2228368
Bromobenzene	U		0.118	1.00	1	02/17/2024 11:32	WG2228368
Bromodichloromethane	U		0.136	1.00	1	02/17/2024 11:32	WG2228368
Bromoform	U		0.129	1.00	1	02/17/2024 11:32	WG2228368
Bromomethane	U		0.605	5.00	1	02/17/2024 11:32	WG2228368
1,3-Butadiene	U		0.299	2.00	1	02/17/2024 11:32	WG2228368
n-Butylbenzene	U		0.157	1.00	1	02/17/2024 11:32	WG2228368
sec-Butylbenzene	U		0.125	1.00	1	02/17/2024 11:32	WG2228368
tert-Butylbenzene	U		0.127	1.00	1	02/17/2024 11:32	WG2228368
Carbon tetrachloride	U		0.128	1.00	1	02/17/2024 11:32	WG2228368
Carbon disulfide	U		0.0962	1.00	1	02/17/2024 11:32	WG2228368
Chlorobenzene	U		0.116	1.00	1	02/17/2024 11:32	WG2228368
Chlorodibromomethane	U		0.140	1.00	1	02/17/2024 11:32	WG2228368
Chloroethane	U		0.192	5.00	1	02/17/2024 11:32	WG2228368
Chloroform	U		0.111	5.00	1	02/17/2024 11:32	WG2228368
Chloromethane	U		0.960	2.50	1	02/17/2024 11:32	WG2228368
Cyclohexane	U		0.188	1.00	1	02/17/2024 11:32	WG2228368
2-Chlorotoluene	U		0.106	1.00	1	02/17/2024 11:32	WG2228368
4-Chlorotoluene	U		0.114	1.00	1	02/17/2024 11:32	WG2228368
1,2-Dibromo-3-Chloropropane	U		0.276	5.00	1	02/17/2024 11:32	WG2228368
1,2-Dibromoethane	U		0.126	1.00	1	02/17/2024 11:32	WG2228368
Dibromomethane	U		0.122	1.00	1	02/17/2024 11:32	WG2228368
1,2-Dichlorobenzene	U		0.107	1.00	1	02/17/2024 11:32	WG2228368
1,3-Dichlorobenzene	U		0.110	1.00	1	02/17/2024 11:32	WG2228368
1,4-Dichlorobenzene	U		0.120	1.00	1	02/17/2024 11:32	WG2228368
Dichlorodifluoromethane	U		0.374	5.00	1	02/17/2024 11:32	WG2228368
1,1-Dichloroethane	U		0.100	1.00	1	02/17/2024 11:32	WG2228368
1,2-Dichloroethane	U		0.0819	1.00	1	02/17/2024 11:32	WG2228368
1,1-Dichloroethene	1.21		0.188	1.00	1	02/17/2024 11:32	WG2228368
cis-1,2-Dichloroethene	U		0.126	1.00	1	02/17/2024 11:32	WG2228368
trans-1,2-Dichloroethene	U		0.149	1.00	1	02/17/2024 11:32	WG2228368
1,2-Dichloropropane	U		0.149	1.00	1	02/17/2024 11:32	WG2228368
1,1-Dichloropropene	U		0.142	1.00	1	02/17/2024 11:32	WG2228368
1,3-Dichloropropane	U		0.110	1.00	1	02/17/2024 11:32	WG2228368
cis-1,3-Dichloropropene	U		0.111	1.00	1	02/17/2024 11:32	WG2228368
trans-1,3-Dichloropropene	U		0.118	1.00	1	02/17/2024 11:32	WG2228368
2,2-Dichloropropane	U		0.161	1.00	1	02/17/2024 11:32	WG2228368
Dicyclopentadiene	U		0.253	1.00	1	02/17/2024 11:32	WG2228368
Di-isopropyl ether	U		0.105	1.00	1	02/17/2024 11:32	WG2228368
Ethylbenzene	U		0.137	1.00	1	02/17/2024 11:32	WG2228368
4-Ethyltoluene	U		0.208	1.00	1	02/17/2024 11:32	WG2228368
Hexachloro-1,3-butadiene	U		0.337	1.00	1	02/17/2024 11:32	WG2228368
n-Hexane	U		0.749	10.0	1	02/17/2024 11:32	WG2228368
Isopropylbenzene	U		0.105	1.00	1	02/17/2024 11:32	WG2228368
p-Isopropyltoluene	U		0.120	1.00	1	02/17/2024 11:32	WG2228368
2-Butanone (MEK)	U		1.19	10.0	1	02/17/2024 11:32	WG2228368
Methyl Cyclohexane	U		0.660	1.00	1	02/17/2024 11:32	WG2228368

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Methylene Chloride	U		0.430	5.00	1	02/17/2024 11:32	WG2228368
4-Methyl-2-pentanone (MIBK)	U		0.478	10.0	1	02/17/2024 11:32	WG2228368
Methyl tert-butyl ether	U		0.101	1.00	1	02/17/2024 11:32	WG2228368
Naphthalene	U		1.00	5.00	1	02/17/2024 11:32	WG2228368
Propene	U		0.936	2.50	1	02/17/2024 11:32	WG2228368
n-Propylbenzene	U		0.0993	1.00	1	02/17/2024 11:32	WG2228368
Styrene	U		0.118	1.00	1	02/17/2024 11:32	WG2228368
1,1,1,2-Tetrachloroethane	U		0.147	1.00	1	02/17/2024 11:32	WG2228368
1,1,2,2-Tetrachloroethane	U		0.133	1.00	1	02/17/2024 11:32	WG2228368
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00	1	02/17/2024 11:32	WG2228368
Tetrachloroethene	U		0.300	1.00	1	02/17/2024 11:32	WG2228368
Toluene	U		0.278	1.00	1	02/17/2024 11:32	WG2228368
1,2,3-Trichlorobenzene	U		0.230	1.00	1	02/17/2024 11:32	WG2228368
1,2,4-Trichlorobenzene	U		0.481	1.00	1	02/17/2024 11:32	WG2228368
1,1,1-Trichloroethane	U		0.149	1.00	1	02/17/2024 11:32	WG2228368
1,1,2-Trichloroethane	U		0.158	1.00	1	02/17/2024 11:32	WG2228368
Trichloroethene	5.44		0.190	1.00	1	02/17/2024 11:32	WG2228368
Trichlorofluoromethane	U		0.160	5.00	1	02/17/2024 11:32	WG2228368
1,2,3-Trichloropropane	U		0.237	2.50	1	02/17/2024 11:32	WG2228368
1,2,4-Trimethylbenzene	U		0.322	1.00	1	02/17/2024 11:32	WG2228368
1,2,3-Trimethylbenzene	U		0.104	1.00	1	02/17/2024 11:32	WG2228368
1,3,5-Trimethylbenzene	U		0.104	1.00	1	02/17/2024 11:32	WG2228368
Vinyl chloride	U		0.234	1.00	1	02/17/2024 11:32	WG2228368
Xylenes, Total	U		0.174	3.00	1	02/17/2024 11:32	WG2228368
(S) Toluene-d8	*102*			*80.0-120*		*02/17/2024 11:32*	WG2228368
(S) 4-Bromofluorobenzene	*92.4*			*77.0-126*		*02/17/2024 11:32*	WG2228368
(S) 1,2-Dichloroethane-d4	*104*			*70.0-130*		*02/17/2024 11:32*	WG2228368

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
1,4-Dioxane	18.2		0.597	3.00	1	02/16/2024 04:38	WG2227341
(S) Toluene-d8	*96.1*			*77.0-127*		*02/16/2024 04:38*	WG2227341

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	150000		1500	20000	5000	02/21/2024 01:26	WG2228552

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		113	500	10	02/17/2024 15:18	WG2228368
Acrolein	U		25.4	500	10	02/17/2024 15:18	WG2228368
Acrylonitrile	U		6.71	100	10	02/17/2024 15:18	WG2228368
Benzene	1.14	E4	0.941	10.0	10	02/17/2024 15:18	WG2228368
Bromobenzene	U		1.18	10.0	10	02/17/2024 15:18	WG2228368
Bromodichloromethane	U		1.36	10.0	10	02/17/2024 15:18	WG2228368
Bromoform	U		1.29	10.0	10	02/17/2024 15:18	WG2228368
Bromomethane	U		6.05	50.0	10	02/17/2024 15:18	WG2228368
1,3-Butadiene	U		2.99	20.0	10	02/17/2024 15:18	WG2228368
n-Butylbenzene	U		1.57	10.0	10	02/17/2024 15:18	WG2228368
sec-Butylbenzene	U		1.25	10.0	10	02/17/2024 15:18	WG2228368
tert-Butylbenzene	U		1.27	10.0	10	02/17/2024 15:18	WG2228368
Carbon tetrachloride	U		1.28	10.0	10	02/17/2024 15:18	WG2228368
Carbon disulfide	U		0.962	10.0	10	02/17/2024 15:18	WG2228368
Chlorobenzene	U		1.16	10.0	10	02/17/2024 15:18	WG2228368
Chlorodibromomethane	U		1.40	10.0	10	02/17/2024 15:18	WG2228368
Chloroethane	U		1.92	50.0	10	02/17/2024 15:18	WG2228368
Chloroform	1.65	E4	1.11	50.0	10	02/17/2024 15:18	WG2228368
Chloromethane	U		9.60	25.0	10	02/17/2024 15:18	WG2228368
Cyclohexane	U		1.88	10.0	10	02/17/2024 15:18	WG2228368
2-Chlorotoluene	U		1.06	10.0	10	02/17/2024 15:18	WG2228368
4-Chlorotoluene	U		1.14	10.0	10	02/17/2024 15:18	WG2228368
1,2-Dibromo-3-Chloropropane	U		2.76	50.0	10	02/17/2024 15:18	WG2228368
1,2-Dibromoethane	U		1.26	10.0	10	02/17/2024 15:18	WG2228368
Dibromomethane	U		1.22	10.0	10	02/17/2024 15:18	WG2228368
1,2-Dichlorobenzene	U		1.07	10.0	10	02/17/2024 15:18	WG2228368
1,3-Dichlorobenzene	U		1.10	10.0	10	02/17/2024 15:18	WG2228368
1,4-Dichlorobenzene	U		1.20	10.0	10	02/17/2024 15:18	WG2228368
Dichlorodifluoromethane	U		3.74	50.0	10	02/17/2024 15:18	WG2228368
1,1-Dichloroethane	U		1.00	10.0	10	02/17/2024 15:18	WG2228368
1,2-Dichloroethane	U		0.819	10.0	10	02/17/2024 15:18	WG2228368
1,1-Dichloroethene	31.6		1.88	10.0	10	02/17/2024 15:18	WG2228368
cis-1,2-Dichloroethene	U		1.26	10.0	10	02/17/2024 15:18	WG2228368
trans-1,2-Dichloroethene	U		1.49	10.0	10	02/17/2024 15:18	WG2228368
1,2-Dichloropropane	U		1.49	10.0	10	02/17/2024 15:18	WG2228368
1,1-Dichloropropene	U		1.42	10.0	10	02/17/2024 15:18	WG2228368
1,3-Dichloropropane	U		1.10	10.0	10	02/17/2024 15:18	WG2228368
cis-1,3-Dichloropropene	U		1.11	10.0	10	02/17/2024 15:18	WG2228368
trans-1,3-Dichloropropene	U		1.18	10.0	10	02/17/2024 15:18	WG2228368
2,2-Dichloropropane	U		1.61	10.0	10	02/17/2024 15:18	WG2228368
Dicyclopentadiene	U		2.53	10.0	10	02/17/2024 15:18	WG2228368
Di-isopropyl ether	U		1.05	10.0	10	02/17/2024 15:18	WG2228368
Ethylbenzene	U		1.37	10.0	10	02/17/2024 15:18	WG2228368
4-Ethyltoluene	U		2.08	10.0	10	02/17/2024 15:18	WG2228368
Hexachloro-1,3-butadiene	U		3.37	10.0	10	02/17/2024 15:18	WG2228368
n-Hexane	U		7.49	100	10	02/17/2024 15:18	WG2228368
Isopropylbenzene	U		1.05	10.0	10	02/17/2024 15:18	WG2228368
p-Isopropyltoluene	U		1.20	10.0	10	02/17/2024 15:18	WG2228368
2-Butanone (MEK)	U		11.9	100	10	02/17/2024 15:18	WG2228368
Methyl Cyclohexane	8.46	E4	6.60	10.0	10	02/17/2024 15:18	WG2228368

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Methylene Chloride	U		4.30	50.0	10	02/20/2024 01:23	WG2229520
4-Methyl-2-pentanone (MIBK)	U		4.78	100	10	02/17/2024 15:18	WG2228368
Methyl tert-butyl ether	U		1.01	10.0	10	02/17/2024 15:18	WG2228368
Naphthalene	U		10.0	50.0	10	02/17/2024 15:18	WG2228368
Propene	U		9.36	25.0	10	02/17/2024 15:18	WG2228368
n-Propylbenzene	U		0.993	10.0	10	02/17/2024 15:18	WG2228368
Styrene	U		1.18	10.0	10	02/17/2024 15:18	WG2228368
1,1,1,2-Tetrachloroethane	U		1.47	10.0	10	02/17/2024 15:18	WG2228368
1,1,2,2-Tetrachloroethane	U		1.33	10.0	10	02/17/2024 15:18	WG2228368
1,1,2-Trichlorotrifluoroethane	U		1.80	10.0	10	02/17/2024 15:18	WG2228368
Tetrachloroethene	U		3.00	10.0	10	02/17/2024 15:18	WG2228368
Toluene	U		2.78	10.0	10	02/17/2024 15:18	WG2228368
1,2,3-Trichlorobenzene	U		2.30	10.0	10	02/17/2024 15:18	WG2228368
1,2,4-Trichlorobenzene	U		4.81	10.0	10	02/17/2024 15:18	WG2228368
1,1,1-Trichloroethane	U		1.49	10.0	10	02/17/2024 15:18	WG2228368
1,1,2-Trichloroethane	U		1.58	10.0	10	02/17/2024 15:18	WG2228368
Trichloroethene	529		1.90	10.0	10	02/20/2024 01:23	WG2229520
Trichlorofluoromethane	U		1.60	50.0	10	02/17/2024 15:18	WG2228368
1,2,3-Trichloropropane	U		2.37	25.0	10	02/17/2024 15:18	WG2228368
1,2,4-Trimethylbenzene	U		3.22	10.0	10	02/17/2024 15:18	WG2228368
1,2,3-Trimethylbenzene	U		1.04	10.0	10	02/17/2024 15:18	WG2228368
1,3,5-Trimethylbenzene	U		1.04	10.0	10	02/17/2024 15:18	WG2228368
Vinyl chloride	U		2.34	10.0	10	02/17/2024 15:18	WG2228368
Xylenes, Total	U		1.74	30.0	10	02/17/2024 15:18	WG2228368
(S) Toluene-d8	*102*			*80.0-120*		*02/17/2024 15:18*	WG2228368
(S) Toluene-d8	*94.0*			*80.0-120*		*02/20/2024 01:23*	WG2229520
(S) 4-Bromofluorobenzene	*89.1*			*77.0-126*		*02/17/2024 15:18*	WG2228368
(S) 4-Bromofluorobenzene	*87.9*			*77.0-126*		*02/20/2024 01:23*	WG2229520
(S) 1,2-Dichloroethane-d4	*90.9*			*70.0-130*		*02/17/2024 15:18*	WG2228368
(S) 1,2-Dichloroethane-d4	*92.7*			*70.0-130*		*02/20/2024 01:23*	WG2229520

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
1,4-Dioxane	207		0.597	3.00	1	02/16/2024 05:00	WG2227341
(S) Toluene-d8	*96.8*			*77.0-127*		*02/16/2024 05:00*	WG2227341

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	54.0		0.300	4.00	1	02/21/2024 01:54	WG2228552

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		11.3	50.0	1	02/17/2024 11:52	WG2228368
Acrolein	U		2.54	50.0	1	02/17/2024 11:52	WG2228368
Acrylonitrile	U		0.671	10.0	1	02/17/2024 11:52	WG2228368
Benzene	U		0.0941	1.00	1	02/17/2024 11:52	WG2228368
Bromobenzene	U		0.118	1.00	1	02/17/2024 11:52	WG2228368
Bromodichloromethane	U		0.136	1.00	1	02/17/2024 11:52	WG2228368
Bromoform	U		0.129	1.00	1	02/17/2024 11:52	WG2228368
Bromomethane	U		0.605	5.00	1	02/17/2024 11:52	WG2228368
1,3-Butadiene	U		0.299	2.00	1	02/17/2024 11:52	WG2228368
n-Butylbenzene	U		0.157	1.00	1	02/17/2024 11:52	WG2228368
sec-Butylbenzene	U		0.125	1.00	1	02/17/2024 11:52	WG2228368
tert-Butylbenzene	U		0.127	1.00	1	02/17/2024 11:52	WG2228368
Carbon tetrachloride	U		0.128	1.00	1	02/17/2024 11:52	WG2228368
Carbon disulfide	U		0.0962	1.00	1	02/17/2024 11:52	WG2228368
Chlorobenzene	U		0.116	1.00	1	02/17/2024 11:52	WG2228368
Chlorodibromomethane	U		0.140	1.00	1	02/17/2024 11:52	WG2228368
Chloroethane	U		0.192	5.00	1	02/17/2024 11:52	WG2228368
Chloroform	U		0.111	5.00	1	02/17/2024 11:52	WG2228368
Chloromethane	U		0.960	2.50	1	02/17/2024 11:52	WG2228368
Cyclohexane	U		0.188	1.00	1	02/17/2024 11:52	WG2228368
2-Chlorotoluene	U		0.106	1.00	1	02/17/2024 11:52	WG2228368
4-Chlorotoluene	U		0.114	1.00	1	02/17/2024 11:52	WG2228368
1,2-Dibromo-3-Chloropropane	U		0.276	5.00	1	02/17/2024 11:52	WG2228368
1,2-Dibromoethane	U		0.126	1.00	1	02/17/2024 11:52	WG2228368
Dibromomethane	U		0.122	1.00	1	02/17/2024 11:52	WG2228368
1,2-Dichlorobenzene	U		0.107	1.00	1	02/17/2024 11:52	WG2228368
1,3-Dichlorobenzene	U		0.110	1.00	1	02/17/2024 11:52	WG2228368
1,4-Dichlorobenzene	U		0.120	1.00	1	02/17/2024 11:52	WG2228368
Dichlorodifluoromethane	U		0.374	5.00	1	02/17/2024 11:52	WG2228368
1,1-Dichloroethane	U		0.100	1.00	1	02/17/2024 11:52	WG2228368
1,2-Dichloroethane	U		0.0819	1.00	1	02/17/2024 11:52	WG2228368
1,1-Dichloroethene	U		0.188	1.00	1	02/17/2024 11:52	WG2228368
cis-1,2-Dichloroethene	U		0.126	1.00	1	02/17/2024 11:52	WG2228368
trans-1,2-Dichloroethene	U		0.149	1.00	1	02/17/2024 11:52	WG2228368
1,2-Dichloropropane	U		0.149	1.00	1	02/17/2024 11:52	WG2228368
1,1-Dichloropropene	U		0.142	1.00	1	02/17/2024 11:52	WG2228368
1,3-Dichloropropane	U		0.110	1.00	1	02/17/2024 11:52	WG2228368
cis-1,3-Dichloropropene	U		0.111	1.00	1	02/17/2024 11:52	WG2228368
trans-1,3-Dichloropropene	U		0.118	1.00	1	02/17/2024 11:52	WG2228368
2,2-Dichloropropane	U		0.161	1.00	1	02/17/2024 11:52	WG2228368
Dicyclopentadiene	U		0.253	1.00	1	02/17/2024 11:52	WG2228368
Di-isopropyl ether	U		0.105	1.00	1	02/17/2024 11:52	WG2228368
Ethylbenzene	U		0.137	1.00	1	02/17/2024 11:52	WG2228368
4-Ethyltoluene	U		0.208	1.00	1	02/17/2024 11:52	WG2228368
Hexachloro-1,3-butadiene	U		0.337	1.00	1	02/17/2024 11:52	WG2228368
n-Hexane	U		0.749	10.0	1	02/17/2024 11:52	WG2228368
Isopropylbenzene	U		0.105	1.00	1	02/17/2024 11:52	WG2228368
p-Isopropyltoluene	U		0.120	1.00	1	02/17/2024 11:52	WG2228368
2-Butanone (MEK)	U		1.19	10.0	1	02/17/2024 11:52	WG2228368
Methyl Cyclohexane	U		0.660	1.00	1	02/17/2024 11:52	WG2228368

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Methylene Chloride	U		0.430	5.00	1	02/17/2024 11:52	WG2228368
4-Methyl-2-pentanone (MIBK)	U		0.478	10.0	1	02/17/2024 11:52	WG2228368
Methyl tert-butyl ether	U		0.101	1.00	1	02/17/2024 11:52	WG2228368
Naphthalene	U		1.00	5.00	1	02/17/2024 11:52	WG2228368
Propene	U		0.936	2.50	1	02/17/2024 11:52	WG2228368
n-Propylbenzene	U		0.0993	1.00	1	02/17/2024 11:52	WG2228368
Styrene	U		0.118	1.00	1	02/17/2024 11:52	WG2228368
1,1,1,2-Tetrachloroethane	U		0.147	1.00	1	02/17/2024 11:52	WG2228368
1,1,2,2-Tetrachloroethane	U		0.133	1.00	1	02/17/2024 11:52	WG2228368
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00	1	02/17/2024 11:52	WG2228368
Tetrachloroethene	U		0.300	1.00	1	02/17/2024 11:52	WG2228368
Toluene	U		0.278	1.00	1	02/17/2024 11:52	WG2228368
1,2,3-Trichlorobenzene	U		0.230	1.00	1	02/17/2024 11:52	WG2228368
1,2,4-Trichlorobenzene	U		0.481	1.00	1	02/17/2024 11:52	WG2228368
1,1,1-Trichloroethane	U		0.149	1.00	1	02/17/2024 11:52	WG2228368
1,1,2-Trichloroethane	U		0.158	1.00	1	02/17/2024 11:52	WG2228368
Trichloroethene	U		0.190	1.00	1	02/17/2024 11:52	WG2228368
Trichlorofluoromethane	U		0.160	5.00	1	02/17/2024 11:52	WG2228368
1,2,3-Trichloropropane	U		0.237	2.50	1	02/17/2024 11:52	WG2228368
1,2,4-Trimethylbenzene	U		0.322	1.00	1	02/17/2024 11:52	WG2228368
1,2,3-Trimethylbenzene	U		0.104	1.00	1	02/17/2024 11:52	WG2228368
1,3,5-Trimethylbenzene	U		0.104	1.00	1	02/17/2024 11:52	WG2228368
Vinyl chloride	U		0.234	1.00	1	02/17/2024 11:52	WG2228368
Xylenes, Total	U		0.174	3.00	1	02/17/2024 11:52	WG2228368
(S) Toluene-d8	*101*			*80.0-120*		*02/17/2024 11:52*	WG2228368
(S) 4-Bromofluorobenzene	*88.4*			*77.0-126*		*02/17/2024 11:52*	WG2228368
(S) 1,2-Dichloroethane-d4	*101*			*70.0-130*		*02/17/2024 11:52*	WG2228368

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
1,4-Dioxane	5.72		0.597	3.00	1	02/16/2024 05:22	WG2227341
(S) Toluene-d8	*93.2*			*77.0-127*		*02/16/2024 05:22*	WG2227341

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	U		0.300	4.00	1	02/21/2024 02:22	WG2228552

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		11.3	50.0	1	02/17/2024 12:13	WG2228368
Acrolein	U		2.54	50.0	1	02/17/2024 12:13	WG2228368
Acrylonitrile	U		0.671	10.0	1	02/17/2024 12:13	WG2228368
Benzene	U		0.0941	1.00	1	02/17/2024 12:13	WG2228368
Bromobenzene	U		0.118	1.00	1	02/17/2024 12:13	WG2228368
Bromodichloromethane	U		0.136	1.00	1	02/17/2024 12:13	WG2228368
Bromoform	U		0.129	1.00	1	02/17/2024 12:13	WG2228368
Bromomethane	U		0.605	5.00	1	02/17/2024 12:13	WG2228368
1,3-Butadiene	U		0.299	2.00	1	02/17/2024 12:13	WG2228368
n-Butylbenzene	U		0.157	1.00	1	02/17/2024 12:13	WG2228368
sec-Butylbenzene	U		0.125	1.00	1	02/17/2024 12:13	WG2228368
tert-Butylbenzene	U		0.127	1.00	1	02/17/2024 12:13	WG2228368
Carbon tetrachloride	U		0.128	1.00	1	02/17/2024 12:13	WG2228368
Carbon disulfide	U		0.0962	1.00	1	02/17/2024 12:13	WG2228368
Chlorobenzene	U		0.116	1.00	1	02/17/2024 12:13	WG2228368
Chlorodibromomethane	U		0.140	1.00	1	02/17/2024 12:13	WG2228368
Chloroethane	U		0.192	5.00	1	02/17/2024 12:13	WG2228368
Chloroform	U		0.111	5.00	1	02/17/2024 12:13	WG2228368
Chloromethane	U		0.960	2.50	1	02/17/2024 12:13	WG2228368
Cyclohexane	U		0.188	1.00	1	02/17/2024 12:13	WG2228368
2-Chlorotoluene	U		0.106	1.00	1	02/17/2024 12:13	WG2228368
4-Chlorotoluene	U		0.114	1.00	1	02/17/2024 12:13	WG2228368
1,2-Dibromo-3-Chloropropane	U		0.276	5.00	1	02/17/2024 12:13	WG2228368
1,2-Dibromoethane	U		0.126	1.00	1	02/17/2024 12:13	WG2228368
Dibromomethane	U		0.122	1.00	1	02/17/2024 12:13	WG2228368
1,2-Dichlorobenzene	U		0.107	1.00	1	02/17/2024 12:13	WG2228368
1,3-Dichlorobenzene	U		0.110	1.00	1	02/17/2024 12:13	WG2228368
1,4-Dichlorobenzene	U		0.120	1.00	1	02/17/2024 12:13	WG2228368
Dichlorodifluoromethane	U		0.374	5.00	1	02/17/2024 12:13	WG2228368
1,1-Dichloroethane	U		0.100	1.00	1	02/17/2024 12:13	WG2228368
1,2-Dichloroethane	U		0.0819	1.00	1	02/17/2024 12:13	WG2228368
1,1-Dichloroethene	U		0.188	1.00	1	02/17/2024 12:13	WG2228368
cis-1,2-Dichloroethene	U		0.126	1.00	1	02/17/2024 12:13	WG2228368
trans-1,2-Dichloroethene	U		0.149	1.00	1	02/17/2024 12:13	WG2228368
1,2-Dichloropropane	U		0.149	1.00	1	02/17/2024 12:13	WG2228368
1,1-Dichloropropene	U		0.142	1.00	1	02/17/2024 12:13	WG2228368
1,3-Dichloropropane	U		0.110	1.00	1	02/17/2024 12:13	WG2228368
cis-1,3-Dichloropropene	U		0.111	1.00	1	02/17/2024 12:13	WG2228368
trans-1,3-Dichloropropene	U		0.118	1.00	1	02/17/2024 12:13	WG2228368
2,2-Dichloropropane	U		0.161	1.00	1	02/17/2024 12:13	WG2228368
Dicyclopentadiene	U		0.253	1.00	1	02/17/2024 12:13	WG2228368
Di-isopropyl ether	U		0.105	1.00	1	02/17/2024 12:13	WG2228368
Ethylbenzene	U		0.137	1.00	1	02/17/2024 12:13	WG2228368
4-Ethyltoluene	U		0.208	1.00	1	02/17/2024 12:13	WG2228368
Hexachloro-1,3-butadiene	U		0.337	1.00	1	02/17/2024 12:13	WG2228368
n-Hexane	U		0.749	10.0	1	02/17/2024 12:13	WG2228368
Isopropylbenzene	U		0.105	1.00	1	02/17/2024 12:13	WG2228368
p-Isopropyltoluene	U		0.120	1.00	1	02/17/2024 12:13	WG2228368
2-Butanone (MEK)	U		1.19	10.0	1	02/17/2024 12:13	WG2228368
Methyl Cyclohexane	U		0.660	1.00	1	02/17/2024 12:13	WG2228368

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Methylene Chloride	U		0.430	5.00	1	02/17/2024 12:13	WG2228368
4-Methyl-2-pentanone (MIBK)	U		0.478	10.0	1	02/17/2024 12:13	WG2228368
Methyl tert-butyl ether	U		0.101	1.00	1	02/17/2024 12:13	WG2228368
Naphthalene	U		1.00	5.00	1	02/17/2024 12:13	WG2228368
Propene	U		0.936	2.50	1	02/17/2024 12:13	WG2228368
n-Propylbenzene	U		0.0993	1.00	1	02/17/2024 12:13	WG2228368
Styrene	U		0.118	1.00	1	02/17/2024 12:13	WG2228368
1,1,1,2-Tetrachloroethane	U		0.147	1.00	1	02/17/2024 12:13	WG2228368
1,1,2,2-Tetrachloroethane	U		0.133	1.00	1	02/17/2024 12:13	WG2228368
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00	1	02/17/2024 12:13	WG2228368
Tetrachloroethene	U		0.300	1.00	1	02/17/2024 12:13	WG2228368
Toluene	U		0.278	1.00	1	02/17/2024 12:13	WG2228368
1,2,3-Trichlorobenzene	U		0.230	1.00	1	02/17/2024 12:13	WG2228368
1,2,4-Trichlorobenzene	U		0.481	1.00	1	02/17/2024 12:13	WG2228368
1,1,1-Trichloroethane	U		0.149	1.00	1	02/17/2024 12:13	WG2228368
1,1,2-Trichloroethane	U		0.158	1.00	1	02/17/2024 12:13	WG2228368
Trichloroethene	U		0.190	1.00	1	02/17/2024 12:13	WG2228368
Trichlorofluoromethane	U		0.160	5.00	1	02/17/2024 12:13	WG2228368
1,2,3-Trichloropropane	U		0.237	2.50	1	02/17/2024 12:13	WG2228368
1,2,4-Trimethylbenzene	U		0.322	1.00	1	02/17/2024 12:13	WG2228368
1,2,3-Trimethylbenzene	U		0.104	1.00	1	02/17/2024 12:13	WG2228368
1,3,5-Trimethylbenzene	U		0.104	1.00	1	02/17/2024 12:13	WG2228368
Vinyl chloride	U		0.234	1.00	1	02/17/2024 12:13	WG2228368
Xylenes, Total	U		0.174	3.00	1	02/17/2024 12:13	WG2228368
(S) Toluene-d8	*102*			*80.0-120*		*02/17/2024 12:13*	WG2228368
(S) 4-Bromofluorobenzene	*90.4*			*77.0-126*		*02/17/2024 12:13*	WG2228368
(S) 1,2-Dichloroethane-d4	*105*			*70.0-130*		*02/17/2024 12:13*	WG2228368

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
1,4-Dioxane	U		0.597	3.00	1	02/16/2024 05:43	WG2227341
(S) Toluene-d8	*93.3*			*77.0-127*		*02/16/2024 05:43*	WG2227341

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	37.7		0.300	4.00	1	02/21/2024 02:50	WG2228552

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		11.3	50.0	1	02/17/2024 12:33	WG2228368
Acrolein	U		2.54	50.0	1	02/17/2024 12:33	WG2228368
Acrylonitrile	U		0.671	10.0	1	02/17/2024 12:33	WG2228368
Benzene	U		0.0941	1.00	1	02/17/2024 12:33	WG2228368
Bromobenzene	U		0.118	1.00	1	02/17/2024 12:33	WG2228368
Bromodichloromethane	U		0.136	1.00	1	02/17/2024 12:33	WG2228368
Bromoform	U		0.129	1.00	1	02/17/2024 12:33	WG2228368
Bromomethane	U		0.605	5.00	1	02/17/2024 12:33	WG2228368
1,3-Butadiene	U		0.299	2.00	1	02/17/2024 12:33	WG2228368
n-Butylbenzene	U		0.157	1.00	1	02/17/2024 12:33	WG2228368
sec-Butylbenzene	U		0.125	1.00	1	02/17/2024 12:33	WG2228368
tert-Butylbenzene	U		0.127	1.00	1	02/17/2024 12:33	WG2228368
Carbon tetrachloride	U		0.128	1.00	1	02/17/2024 12:33	WG2228368
Carbon disulfide	U		0.0962	1.00	1	02/17/2024 12:33	WG2228368
Chlorobenzene	U		0.116	1.00	1	02/17/2024 12:33	WG2228368
Chlorodibromomethane	U		0.140	1.00	1	02/17/2024 12:33	WG2228368
Chloroethane	U		0.192	5.00	1	02/17/2024 12:33	WG2228368
Chloroform	U		0.111	5.00	1	02/17/2024 12:33	WG2228368
Chloromethane	U		0.960	2.50	1	02/17/2024 12:33	WG2228368
Cyclohexane	U		0.188	1.00	1	02/17/2024 12:33	WG2228368
2-Chlorotoluene	U		0.106	1.00	1	02/17/2024 12:33	WG2228368
4-Chlorotoluene	U		0.114	1.00	1	02/17/2024 12:33	WG2228368
1,2-Dibromo-3-Chloropropane	U		0.276	5.00	1	02/17/2024 12:33	WG2228368
1,2-Dibromoethane	U		0.126	1.00	1	02/17/2024 12:33	WG2228368
Dibromomethane	U		0.122	1.00	1	02/17/2024 12:33	WG2228368
1,2-Dichlorobenzene	U		0.107	1.00	1	02/17/2024 12:33	WG2228368
1,3-Dichlorobenzene	U		0.110	1.00	1	02/17/2024 12:33	WG2228368
1,4-Dichlorobenzene	U		0.120	1.00	1	02/17/2024 12:33	WG2228368
Dichlorodifluoromethane	U		0.374	5.00	1	02/17/2024 12:33	WG2228368
1,1-Dichloroethane	U		0.100	1.00	1	02/17/2024 12:33	WG2228368
1,2-Dichloroethane	U		0.0819	1.00	1	02/17/2024 12:33	WG2228368
1,1-Dichloroethene	U		0.188	1.00	1	02/17/2024 12:33	WG2228368
cis-1,2-Dichloroethene	U		0.126	1.00	1	02/17/2024 12:33	WG2228368
trans-1,2-Dichloroethene	U		0.149	1.00	1	02/17/2024 12:33	WG2228368
1,2-Dichloropropane	U		0.149	1.00	1	02/17/2024 12:33	WG2228368
1,1-Dichloropropene	U		0.142	1.00	1	02/17/2024 12:33	WG2228368
1,3-Dichloropropane	U		0.110	1.00	1	02/17/2024 12:33	WG2228368
cis-1,3-Dichloropropene	U		0.111	1.00	1	02/17/2024 12:33	WG2228368
trans-1,3-Dichloropropene	U		0.118	1.00	1	02/17/2024 12:33	WG2228368
2,2-Dichloropropane	U		0.161	1.00	1	02/17/2024 12:33	WG2228368
Dicyclopentadiene	U		0.253	1.00	1	02/17/2024 12:33	WG2228368
Di-isopropyl ether	U		0.105	1.00	1	02/17/2024 12:33	WG2228368
Ethylbenzene	U		0.137	1.00	1	02/17/2024 12:33	WG2228368
4-Ethyltoluene	U		0.208	1.00	1	02/17/2024 12:33	WG2228368
Hexachloro-1,3-butadiene	U		0.337	1.00	1	02/17/2024 12:33	WG2228368
n-Hexane	U		0.749	10.0	1	02/17/2024 12:33	WG2228368
Isopropylbenzene	U		0.105	1.00	1	02/17/2024 12:33	WG2228368
p-Isopropyltoluene	U		0.120	1.00	1	02/17/2024 12:33	WG2228368
2-Butanone (MEK)	U		1.19	10.0	1	02/17/2024 12:33	WG2228368
Methyl Cyclohexane	U		0.660	1.00	1	02/17/2024 12:33	WG2228368

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Methylene Chloride	U		0.430	5.00	1	02/17/2024 12:33	WG2228368
4-Methyl-2-pentanone (MIBK)	U		0.478	10.0	1	02/17/2024 12:33	WG2228368
Methyl tert-butyl ether	U		0.101	1.00	1	02/17/2024 12:33	WG2228368
Naphthalene	U		1.00	5.00	1	02/17/2024 12:33	WG2228368
Propene	U		0.936	2.50	1	02/17/2024 12:33	WG2228368
n-Propylbenzene	U		0.0993	1.00	1	02/17/2024 12:33	WG2228368
Styrene	U		0.118	1.00	1	02/17/2024 12:33	WG2228368
1,1,1,2-Tetrachloroethane	U		0.147	1.00	1	02/17/2024 12:33	WG2228368
1,1,2,2-Tetrachloroethane	U		0.133	1.00	1	02/17/2024 12:33	WG2228368
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00	1	02/17/2024 12:33	WG2228368
Tetrachloroethene	U		0.300	1.00	1	02/17/2024 12:33	WG2228368
Toluene	U		0.278	1.00	1	02/17/2024 12:33	WG2228368
1,2,3-Trichlorobenzene	U		0.230	1.00	1	02/17/2024 12:33	WG2228368
1,2,4-Trichlorobenzene	U		0.481	1.00	1	02/17/2024 12:33	WG2228368
1,1,1-Trichloroethane	U		0.149	1.00	1	02/17/2024 12:33	WG2228368
1,1,2-Trichloroethane	U		0.158	1.00	1	02/17/2024 12:33	WG2228368
Trichloroethene	U		0.190	1.00	1	02/17/2024 12:33	WG2228368
Trichlorofluoromethane	U		0.160	5.00	1	02/17/2024 12:33	WG2228368
1,2,3-Trichloropropane	U		0.237	2.50	1	02/17/2024 12:33	WG2228368
1,2,4-Trimethylbenzene	U		0.322	1.00	1	02/17/2024 12:33	WG2228368
1,2,3-Trimethylbenzene	U		0.104	1.00	1	02/17/2024 12:33	WG2228368
1,3,5-Trimethylbenzene	U		0.104	1.00	1	02/17/2024 12:33	WG2228368
Vinyl chloride	U		0.234	1.00	1	02/17/2024 12:33	WG2228368
Xylenes, Total	U		0.174	3.00	1	02/17/2024 12:33	WG2228368
(S) Toluene-d8	*100*			*80.0-120*		*02/17/2024 12:33*	WG2228368
(S) 4-Bromofluorobenzene	*88.6*			*77.0-126*		*02/17/2024 12:33*	WG2228368
(S) 1,2-Dichloroethane-d4	*107*			*70.0-130*		*02/17/2024 12:33*	WG2228368

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
1,4-Dioxane	U		0.597	3.00	1	02/16/2024 06:05	WG2227341
(S) Toluene-d8	*93.1*			*77.0-127*		*02/16/2024 06:05*	WG2227341

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	18.1		0.300	4.00	1	02/21/2024 04:14	WG2228552

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		11.3	50.0	1	02/17/2024 12:54	WG2228368
Acrolein	U		2.54	50.0	1	02/17/2024 12:54	WG2228368
Acrylonitrile	U		0.671	10.0	1	02/17/2024 12:54	WG2228368
Benzene	U		0.0941	1.00	1	02/17/2024 12:54	WG2228368
Bromobenzene	U		0.118	1.00	1	02/17/2024 12:54	WG2228368
Bromodichloromethane	U		0.136	1.00	1	02/17/2024 12:54	WG2228368
Bromoform	U		0.129	1.00	1	02/17/2024 12:54	WG2228368
Bromomethane	U		0.605	5.00	1	02/17/2024 12:54	WG2228368
1,3-Butadiene	U		0.299	2.00	1	02/17/2024 12:54	WG2228368
n-Butylbenzene	U		0.157	1.00	1	02/17/2024 12:54	WG2228368
sec-Butylbenzene	U		0.125	1.00	1	02/17/2024 12:54	WG2228368
tert-Butylbenzene	U		0.127	1.00	1	02/17/2024 12:54	WG2228368
Carbon tetrachloride	U		0.128	1.00	1	02/17/2024 12:54	WG2228368
Carbon disulfide	U		0.0962	1.00	1	02/17/2024 12:54	WG2228368
Chlorobenzene	U		0.116	1.00	1	02/17/2024 12:54	WG2228368
Chlorodibromomethane	U		0.140	1.00	1	02/17/2024 12:54	WG2228368
Chloroethane	U		0.192	5.00	1	02/17/2024 12:54	WG2228368
Chloroform	U		0.111	5.00	1	02/17/2024 12:54	WG2228368
Chloromethane	U		0.960	2.50	1	02/17/2024 12:54	WG2228368
Cyclohexane	U		0.188	1.00	1	02/17/2024 12:54	WG2228368
2-Chlorotoluene	U		0.106	1.00	1	02/17/2024 12:54	WG2228368
4-Chlorotoluene	U		0.114	1.00	1	02/17/2024 12:54	WG2228368
1,2-Dibromo-3-Chloropropane	U		0.276	5.00	1	02/17/2024 12:54	WG2228368
1,2-Dibromoethane	U		0.126	1.00	1	02/17/2024 12:54	WG2228368
Dibromomethane	U		0.122	1.00	1	02/17/2024 12:54	WG2228368
1,2-Dichlorobenzene	U		0.107	1.00	1	02/17/2024 12:54	WG2228368
1,3-Dichlorobenzene	U		0.110	1.00	1	02/17/2024 12:54	WG2228368
1,4-Dichlorobenzene	U		0.120	1.00	1	02/17/2024 12:54	WG2228368
Dichlorodifluoromethane	U		0.374	5.00	1	02/17/2024 12:54	WG2228368
1,1-Dichloroethane	U		0.100	1.00	1	02/17/2024 12:54	WG2228368
1,2-Dichloroethane	U		0.0819	1.00	1	02/17/2024 12:54	WG2228368
1,1-Dichloroethene	U		0.188	1.00	1	02/17/2024 12:54	WG2228368
cis-1,2-Dichloroethene	U		0.126	1.00	1	02/17/2024 12:54	WG2228368
trans-1,2-Dichloroethene	U		0.149	1.00	1	02/17/2024 12:54	WG2228368
1,2-Dichloropropane	U		0.149	1.00	1	02/17/2024 12:54	WG2228368
1,1-Dichloropropene	U		0.142	1.00	1	02/17/2024 12:54	WG2228368
1,3-Dichloropropane	U		0.110	1.00	1	02/17/2024 12:54	WG2228368
cis-1,3-Dichloropropene	U		0.111	1.00	1	02/17/2024 12:54	WG2228368
trans-1,3-Dichloropropene	U		0.118	1.00	1	02/17/2024 12:54	WG2228368
2,2-Dichloropropane	U		0.161	1.00	1	02/17/2024 12:54	WG2228368
Dicyclopentadiene	U		0.253	1.00	1	02/17/2024 12:54	WG2228368
Di-isopropyl ether	U		0.105	1.00	1	02/17/2024 12:54	WG2228368
Ethylbenzene	U		0.137	1.00	1	02/17/2024 12:54	WG2228368
4-Ethyltoluene	U		0.208	1.00	1	02/17/2024 12:54	WG2228368
Hexachloro-1,3-butadiene	U		0.337	1.00	1	02/17/2024 12:54	WG2228368
n-Hexane	U		0.749	10.0	1	02/17/2024 12:54	WG2228368
Isopropylbenzene	U		0.105	1.00	1	02/17/2024 12:54	WG2228368
p-Isopropyltoluene	U		0.120	1.00	1	02/17/2024 12:54	WG2228368
2-Butanone (MEK)	U		1.19	10.0	1	02/17/2024 12:54	WG2228368
Methyl Cyclohexane	U		0.660	1.00	1	02/17/2024 12:54	WG2228368

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Methylene Chloride	U		0.430	5.00	1	02/17/2024 12:54	WG2228368
4-Methyl-2-pentanone (MIBK)	U		0.478	10.0	1	02/17/2024 12:54	WG2228368
Methyl tert-butyl ether	U		0.101	1.00	1	02/17/2024 12:54	WG2228368
Naphthalene	U		1.00	5.00	1	02/17/2024 12:54	WG2228368
Propene	U		0.936	2.50	1	02/17/2024 12:54	WG2228368
n-Propylbenzene	U		0.0993	1.00	1	02/17/2024 12:54	WG2228368
Styrene	U		0.118	1.00	1	02/17/2024 12:54	WG2228368
1,1,1,2-Tetrachloroethane	U		0.147	1.00	1	02/17/2024 12:54	WG2228368
1,1,2,2-Tetrachloroethane	U		0.133	1.00	1	02/17/2024 12:54	WG2228368
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00	1	02/17/2024 12:54	WG2228368
Tetrachloroethene	U		0.300	1.00	1	02/17/2024 12:54	WG2228368
Toluene	U		0.278	1.00	1	02/17/2024 12:54	WG2228368
1,2,3-Trichlorobenzene	U		0.230	1.00	1	02/17/2024 12:54	WG2228368
1,2,4-Trichlorobenzene	U		0.481	1.00	1	02/17/2024 12:54	WG2228368
1,1,1-Trichloroethane	U		0.149	1.00	1	02/17/2024 12:54	WG2228368
1,1,2-Trichloroethane	U		0.158	1.00	1	02/17/2024 12:54	WG2228368
Trichloroethene	U		0.190	1.00	1	02/17/2024 12:54	WG2228368
Trichlorofluoromethane	U		0.160	5.00	1	02/17/2024 12:54	WG2228368
1,2,3-Trichloropropane	U		0.237	2.50	1	02/17/2024 12:54	WG2228368
1,2,4-Trimethylbenzene	U		0.322	1.00	1	02/17/2024 12:54	WG2228368
1,2,3-Trimethylbenzene	U		0.104	1.00	1	02/17/2024 12:54	WG2228368
1,3,5-Trimethylbenzene	U		0.104	1.00	1	02/17/2024 12:54	WG2228368
Vinyl chloride	U		0.234	1.00	1	02/17/2024 12:54	WG2228368
Xylenes, Total	U		0.174	3.00	1	02/17/2024 12:54	WG2228368
(S) Toluene-d8	*103*			*80.0-120*		*02/17/2024 12:54*	WG2228368
(S) 4-Bromofluorobenzene	*91.6*			*77.0-126*		*02/17/2024 12:54*	WG2228368
(S) 1,2-Dichloroethane-d4	*102*			*70.0-130*		*02/17/2024 12:54*	WG2228368

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
1,4-Dioxane	U		0.597	3.00	1	02/16/2024 06:27	WG2227341
(S) Toluene-d8	*93.7*			*77.0-127*		*02/16/2024 06:27*	WG2227341

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	U		0.300	4.00	1	02/21/2024 04:41	WG2228552

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		11.3	50.0	1	02/17/2024 13:15	WG2228368
Acrolein	U		2.54	50.0	1	02/17/2024 13:15	WG2228368
Acrylonitrile	U		0.671	10.0	1	02/17/2024 13:15	WG2228368
Benzene	0.113	E4	0.0941	1.00	1	02/17/2024 13:15	WG2228368
Bromobenzene	U		0.118	1.00	1	02/17/2024 13:15	WG2228368
Bromodichloromethane	U		0.136	1.00	1	02/17/2024 13:15	WG2228368
Bromoform	U		0.129	1.00	1	02/17/2024 13:15	WG2228368
Bromomethane	U		0.605	5.00	1	02/17/2024 13:15	WG2228368
1,3-Butadiene	U		0.299	2.00	1	02/17/2024 13:15	WG2228368
n-Butylbenzene	U		0.157	1.00	1	02/17/2024 13:15	WG2228368
sec-Butylbenzene	U		0.125	1.00	1	02/17/2024 13:15	WG2228368
tert-Butylbenzene	U		0.127	1.00	1	02/17/2024 13:15	WG2228368
Carbon tetrachloride	U		0.128	1.00	1	02/17/2024 13:15	WG2228368
Carbon disulfide	U		0.0962	1.00	1	02/17/2024 13:15	WG2228368
Chlorobenzene	U		0.116	1.00	1	02/17/2024 13:15	WG2228368
Chlorodibromomethane	U		0.140	1.00	1	02/17/2024 13:15	WG2228368
Chloroethane	U		0.192	5.00	1	02/17/2024 13:15	WG2228368
Chloroform	U		0.111	5.00	1	02/17/2024 13:15	WG2228368
Chloromethane	U		0.960	2.50	1	02/17/2024 13:15	WG2228368
Cyclohexane	U		0.188	1.00	1	02/17/2024 13:15	WG2228368
2-Chlorotoluene	U		0.106	1.00	1	02/17/2024 13:15	WG2228368
4-Chlorotoluene	U		0.114	1.00	1	02/17/2024 13:15	WG2228368
1,2-Dibromo-3-Chloropropane	U		0.276	5.00	1	02/17/2024 13:15	WG2228368
1,2-Dibromoethane	U		0.126	1.00	1	02/17/2024 13:15	WG2228368
Dibromomethane	U		0.122	1.00	1	02/17/2024 13:15	WG2228368
1,2-Dichlorobenzene	U		0.107	1.00	1	02/17/2024 13:15	WG2228368
1,3-Dichlorobenzene	U		0.110	1.00	1	02/17/2024 13:15	WG2228368
1,4-Dichlorobenzene	U		0.120	1.00	1	02/17/2024 13:15	WG2228368
Dichlorodifluoromethane	U		0.374	5.00	1	02/17/2024 13:15	WG2228368
1,1-Dichloroethane	U		0.100	1.00	1	02/17/2024 13:15	WG2228368
1,2-Dichloroethane	U		0.0819	1.00	1	02/17/2024 13:15	WG2228368
1,1-Dichloroethene	U		0.188	1.00	1	02/17/2024 13:15	WG2228368
cis-1,2-Dichloroethene	U		0.126	1.00	1	02/17/2024 13:15	WG2228368
trans-1,2-Dichloroethene	U		0.149	1.00	1	02/17/2024 13:15	WG2228368
1,2-Dichloropropane	U		0.149	1.00	1	02/17/2024 13:15	WG2228368
1,1-Dichloropropene	U		0.142	1.00	1	02/17/2024 13:15	WG2228368
1,3-Dichloropropane	U		0.110	1.00	1	02/17/2024 13:15	WG2228368
cis-1,3-Dichloropropene	U		0.111	1.00	1	02/17/2024 13:15	WG2228368
trans-1,3-Dichloropropene	U		0.118	1.00	1	02/17/2024 13:15	WG2228368
2,2-Dichloropropane	U		0.161	1.00	1	02/17/2024 13:15	WG2228368
Dicyclopentadiene	U		0.253	1.00	1	02/17/2024 13:15	WG2228368
Di-isopropyl ether	U		0.105	1.00	1	02/17/2024 13:15	WG2228368
Ethylbenzene	U		0.137	1.00	1	02/17/2024 13:15	WG2228368
4-Ethyltoluene	U		0.208	1.00	1	02/17/2024 13:15	WG2228368
Hexachloro-1,3-butadiene	U		0.337	1.00	1	02/17/2024 13:15	WG2228368
n-Hexane	U		0.749	10.0	1	02/17/2024 13:15	WG2228368
Isopropylbenzene	U		0.105	1.00	1	02/17/2024 13:15	WG2228368
p-Isopropyltoluene	U		0.120	1.00	1	02/17/2024 13:15	WG2228368
2-Butanone (MEK)	U		1.19	10.0	1	02/17/2024 13:15	WG2228368
Methyl Cyclohexane	U		0.660	1.00	1	02/17/2024 13:15	WG2228368

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Methylene Chloride	U		0.430	5.00	1	02/17/2024 13:15	WG2228368
4-Methyl-2-pentanone (MIBK)	U		0.478	10.0	1	02/17/2024 13:15	WG2228368
Methyl tert-butyl ether	U		0.101	1.00	1	02/17/2024 13:15	WG2228368
Naphthalene	U		1.00	5.00	1	02/17/2024 13:15	WG2228368
Propene	7.32		0.936	2.50	1	02/17/2024 13:15	WG2228368
n-Propylbenzene	U		0.0993	1.00	1	02/17/2024 13:15	WG2228368
Styrene	U		0.118	1.00	1	02/17/2024 13:15	WG2228368
1,1,1,2-Tetrachloroethane	U		0.147	1.00	1	02/17/2024 13:15	WG2228368
1,1,2,2-Tetrachloroethane	U		0.133	1.00	1	02/17/2024 13:15	WG2228368
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00	1	02/17/2024 13:15	WG2228368
Tetrachloroethene	U		0.300	1.00	1	02/17/2024 13:15	WG2228368
Toluene	0.540	E4	0.278	1.00	1	02/17/2024 13:15	WG2228368
1,2,3-Trichlorobenzene	U		0.230	1.00	1	02/17/2024 13:15	WG2228368
1,2,4-Trichlorobenzene	U		0.481	1.00	1	02/17/2024 13:15	WG2228368
1,1,1-Trichloroethane	U		0.149	1.00	1	02/17/2024 13:15	WG2228368
1,1,2-Trichloroethane	U		0.158	1.00	1	02/17/2024 13:15	WG2228368
Trichloroethene	U		0.190	1.00	1	02/17/2024 13:15	WG2228368
Trichlorofluoromethane	U		0.160	5.00	1	02/17/2024 13:15	WG2228368
1,2,3-Trichloropropane	U		0.237	2.50	1	02/17/2024 13:15	WG2228368
1,2,4-Trimethylbenzene	U		0.322	1.00	1	02/17/2024 13:15	WG2228368
1,2,3-Trimethylbenzene	U		0.104	1.00	1	02/17/2024 13:15	WG2228368
1,3,5-Trimethylbenzene	U		0.104	1.00	1	02/17/2024 13:15	WG2228368
Vinyl chloride	U		0.234	1.00	1	02/17/2024 13:15	WG2228368
Xylenes, Total	0.185	E4	0.174	3.00	1	02/17/2024 13:15	WG2228368
(S) Toluene-d8	*102*			*80.0-120*		*02/17/2024 13:15*	WG2228368
(S) 4-Bromofluorobenzene	*88.3*			*77.0-126*		*02/17/2024 13:15*	WG2228368
(S) 1,2-Dichloroethane-d4	*101*			*70.0-130*		*02/17/2024 13:15*	WG2228368

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
1,4-Dioxane	U		0.597	3.00	1	02/16/2024 06:49	WG2227341
(S) Toluene-d8	*93.1*			*77.0-127*		*02/16/2024 06:49*	WG2227341

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	6.06		0.300	4.00	1	02/21/2024 05:09	WG2228552

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		11.3	50.0	1	02/17/2024 13:35	WG2228368
Acrolein	U		2.54	50.0	1	02/17/2024 13:35	WG2228368
Acrylonitrile	U		0.671	10.0	1	02/17/2024 13:35	WG2228368
Benzene	U		0.0941	1.00	1	02/17/2024 13:35	WG2228368
Bromobenzene	U		0.118	1.00	1	02/17/2024 13:35	WG2228368
Bromodichloromethane	U		0.136	1.00	1	02/17/2024 13:35	WG2228368
Bromoform	U		0.129	1.00	1	02/17/2024 13:35	WG2228368
Bromomethane	U		0.605	5.00	1	02/17/2024 13:35	WG2228368
1,3-Butadiene	U		0.299	2.00	1	02/17/2024 13:35	WG2228368
n-Butylbenzene	U		0.157	1.00	1	02/17/2024 13:35	WG2228368
sec-Butylbenzene	U		0.125	1.00	1	02/17/2024 13:35	WG2228368
tert-Butylbenzene	U		0.127	1.00	1	02/17/2024 13:35	WG2228368
Carbon tetrachloride	U		0.128	1.00	1	02/17/2024 13:35	WG2228368
Carbon disulfide	U		0.0962	1.00	1	02/17/2024 13:35	WG2228368
Chlorobenzene	U		0.116	1.00	1	02/17/2024 13:35	WG2228368
Chlorodibromomethane	U		0.140	1.00	1	02/17/2024 13:35	WG2228368
Chloroethane	U		0.192	5.00	1	02/17/2024 13:35	WG2228368
Chloroform	U		0.111	5.00	1	02/17/2024 13:35	WG2228368
Chloromethane	U		0.960	2.50	1	02/17/2024 13:35	WG2228368
Cyclohexane	U		0.188	1.00	1	02/17/2024 13:35	WG2228368
2-Chlorotoluene	U		0.106	1.00	1	02/17/2024 13:35	WG2228368
4-Chlorotoluene	U		0.114	1.00	1	02/17/2024 13:35	WG2228368
1,2-Dibromo-3-Chloropropane	U		0.276	5.00	1	02/17/2024 13:35	WG2228368
1,2-Dibromoethane	U		0.126	1.00	1	02/17/2024 13:35	WG2228368
Dibromomethane	U		0.122	1.00	1	02/17/2024 13:35	WG2228368
1,2-Dichlorobenzene	U		0.107	1.00	1	02/17/2024 13:35	WG2228368
1,3-Dichlorobenzene	U		0.110	1.00	1	02/17/2024 13:35	WG2228368
1,4-Dichlorobenzene	U		0.120	1.00	1	02/17/2024 13:35	WG2228368
Dichlorodifluoromethane	U		0.374	5.00	1	02/17/2024 13:35	WG2228368
1,1-Dichloroethane	U		0.100	1.00	1	02/17/2024 13:35	WG2228368
1,2-Dichloroethane	U		0.0819	1.00	1	02/17/2024 13:35	WG2228368
1,1-Dichloroethene	U		0.188	1.00	1	02/17/2024 13:35	WG2228368
cis-1,2-Dichloroethene	U		0.126	1.00	1	02/17/2024 13:35	WG2228368
trans-1,2-Dichloroethene	U		0.149	1.00	1	02/17/2024 13:35	WG2228368
1,2-Dichloropropane	U		0.149	1.00	1	02/17/2024 13:35	WG2228368
1,1-Dichloropropene	U		0.142	1.00	1	02/17/2024 13:35	WG2228368
1,3-Dichloropropane	U		0.110	1.00	1	02/17/2024 13:35	WG2228368
cis-1,3-Dichloropropene	U		0.111	1.00	1	02/17/2024 13:35	WG2228368
trans-1,3-Dichloropropene	U		0.118	1.00	1	02/17/2024 13:35	WG2228368
2,2-Dichloropropane	U		0.161	1.00	1	02/17/2024 13:35	WG2228368
Dicyclopentadiene	U		0.253	1.00	1	02/17/2024 13:35	WG2228368
Di-isopropyl ether	U		0.105	1.00	1	02/17/2024 13:35	WG2228368
Ethylbenzene	U		0.137	1.00	1	02/17/2024 13:35	WG2228368
4-Ethyltoluene	U		0.208	1.00	1	02/17/2024 13:35	WG2228368
Hexachloro-1,3-butadiene	U		0.337	1.00	1	02/17/2024 13:35	WG2228368
n-Hexane	U		0.749	10.0	1	02/17/2024 13:35	WG2228368
Isopropylbenzene	U		0.105	1.00	1	02/17/2024 13:35	WG2228368
p-Isopropyltoluene	U		0.120	1.00	1	02/17/2024 13:35	WG2228368
2-Butanone (MEK)	U		1.19	10.0	1	02/17/2024 13:35	WG2228368
Methyl Cyclohexane	U		0.660	1.00	1	02/17/2024 13:35	WG2228368

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Methylene Chloride	U		0.430	5.00	1	02/17/2024 13:35	WG2228368
4-Methyl-2-pentanone (MIBK)	U		0.478	10.0	1	02/17/2024 13:35	WG2228368
Methyl tert-butyl ether	U		0.101	1.00	1	02/17/2024 13:35	WG2228368
Naphthalene	U		1.00	5.00	1	02/17/2024 13:35	WG2228368
Propene	U		0.936	2.50	1	02/17/2024 13:35	WG2228368
n-Propylbenzene	U		0.0993	1.00	1	02/17/2024 13:35	WG2228368
Styrene	U		0.118	1.00	1	02/17/2024 13:35	WG2228368
1,1,1,2-Tetrachloroethane	U		0.147	1.00	1	02/17/2024 13:35	WG2228368
1,1,2,2-Tetrachloroethane	U		0.133	1.00	1	02/17/2024 13:35	WG2228368
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00	1	02/17/2024 13:35	WG2228368
Tetrachloroethene	U		0.300	1.00	1	02/17/2024 13:35	WG2228368
Toluene	U		0.278	1.00	1	02/17/2024 13:35	WG2228368
1,2,3-Trichlorobenzene	U		0.230	1.00	1	02/17/2024 13:35	WG2228368
1,2,4-Trichlorobenzene	U		0.481	1.00	1	02/17/2024 13:35	WG2228368
1,1,1-Trichloroethane	U		0.149	1.00	1	02/17/2024 13:35	WG2228368
1,1,2-Trichloroethane	U		0.158	1.00	1	02/17/2024 13:35	WG2228368
Trichloroethene	U		0.190	1.00	1	02/17/2024 13:35	WG2228368
Trichlorofluoromethane	U		0.160	5.00	1	02/17/2024 13:35	WG2228368
1,2,3-Trichloropropane	U		0.237	2.50	1	02/17/2024 13:35	WG2228368
1,2,4-Trimethylbenzene	U		0.322	1.00	1	02/17/2024 13:35	WG2228368
1,2,3-Trimethylbenzene	U		0.104	1.00	1	02/17/2024 13:35	WG2228368
1,3,5-Trimethylbenzene	U		0.104	1.00	1	02/17/2024 13:35	WG2228368
Vinyl chloride	U		0.234	1.00	1	02/17/2024 13:35	WG2228368
Xylenes, Total	U		0.174	3.00	1	02/17/2024 13:35	WG2228368
(S) Toluene-d8	*99.2*			*80.0-120*		*02/17/2024 13:35*	WG2228368
(S) 4-Bromofluorobenzene	*90.1*			*77.0-126*		*02/17/2024 13:35*	WG2228368
(S) 1,2-Dichloroethane-d4	*103*			*70.0-130*		*02/17/2024 13:35*	WG2228368

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
1,4-Dioxane	U		0.597	3.00	1	02/20/2024 16:25	WG2230157
(S) Toluene-d8	*94.6*			*77.0-127*		*02/20/2024 16:25*	WG2230157

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	U		0.300	4.00	1	02/21/2024 06:05	WG2228552

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	2320		113	500	10	02/20/2024 01:47	WG2229520
Acrolein	U		25.4	500	10	02/20/2024 01:47	WG2229520
Acrylonitrile	U		6.71	100	10	02/20/2024 01:47	WG2229520
Benzene	U		0.941	10.0	10	02/20/2024 01:47	WG2229520
Bromobenzene	U		1.18	10.0	10	02/20/2024 01:47	WG2229520
Bromodichloromethane	U		1.36	10.0	10	02/20/2024 01:47	WG2229520
Bromoform	U	L2	1.29	10.0	10	02/20/2024 01:47	WG2229520
Bromomethane	U		6.05	50.0	10	02/20/2024 01:47	WG2229520
1,3-Butadiene	U		2.99	20.0	10	02/20/2024 01:47	WG2229520
n-Butylbenzene	U		1.57	10.0	10	02/20/2024 01:47	WG2229520
sec-Butylbenzene	U		1.25	10.0	10	02/20/2024 01:47	WG2229520
tert-Butylbenzene	U		1.27	10.0	10	02/20/2024 01:47	WG2229520
Carbon tetrachloride	U		1.28	10.0	10	02/20/2024 01:47	WG2229520
Carbon disulfide	U		0.962	10.0	10	02/20/2024 01:47	WG2229520
Chlorobenzene	U		1.16	10.0	10	02/20/2024 01:47	WG2229520
Chlorodibromomethane	U		1.40	10.0	10	02/20/2024 01:47	WG2229520
Chloroethane	U		1.92	50.0	10	02/20/2024 01:47	WG2229520
Chloroform	U		1.11	50.0	10	02/20/2024 01:47	WG2229520
Chloromethane	U		9.60	25.0	10	02/20/2024 01:47	WG2229520
Cyclohexane	U		1.88	10.0	10	02/20/2024 01:47	WG2229520
2-Chlorotoluene	U		1.06	10.0	10	02/20/2024 01:47	WG2229520
4-Chlorotoluene	U		1.14	10.0	10	02/20/2024 01:47	WG2229520
1,2-Dibromo-3-Chloropropane	U		2.76	50.0	10	02/20/2024 01:47	WG2229520
1,2-Dibromoethane	U		1.26	10.0	10	02/20/2024 01:47	WG2229520
Dibromomethane	U		1.22	10.0	10	02/20/2024 01:47	WG2229520
1,2-Dichlorobenzene	U	L2	1.07	10.0	10	02/20/2024 01:47	WG2229520
1,3-Dichlorobenzene	U	L2	1.10	10.0	10	02/20/2024 01:47	WG2229520
1,4-Dichlorobenzene	U	L2	1.20	10.0	10	02/20/2024 01:47	WG2229520
Dichlorodifluoromethane	U		3.74	50.0	10	02/20/2024 01:47	WG2229520
1,1-Dichloroethane	U		1.00	10.0	10	02/20/2024 01:47	WG2229520
1,2-Dichloroethane	U		0.819	10.0	10	02/20/2024 01:47	WG2229520
1,1-Dichloroethene	5.47	E4	1.88	10.0	10	02/20/2024 01:47	WG2229520
cis-1,2-Dichloroethene	28.8		1.26	10.0	10	02/20/2024 01:47	WG2229520
trans-1,2-Dichloroethene	U		1.49	10.0	10	02/20/2024 01:47	WG2229520
1,2-Dichloropropane	U		1.49	10.0	10	02/20/2024 01:47	WG2229520
1,1-Dichloropropene	U		1.42	10.0	10	02/20/2024 01:47	WG2229520
1,3-Dichloropropane	U		1.10	10.0	10	02/20/2024 01:47	WG2229520
cis-1,3-Dichloropropene	U		1.11	10.0	10	02/20/2024 01:47	WG2229520
trans-1,3-Dichloropropene	U		1.18	10.0	10	02/20/2024 01:47	WG2229520
2,2-Dichloropropane	U		1.61	10.0	10	02/20/2024 01:47	WG2229520
Dicyclopentadiene	U		2.53	10.0	10	02/20/2024 01:47	WG2229520
Di-isopropyl ether	U		1.05	10.0	10	02/20/2024 01:47	WG2229520
Ethylbenzene	U		1.37	10.0	10	02/20/2024 01:47	WG2229520
4-Ethyltoluene	U		2.08	10.0	10	02/20/2024 01:47	WG2229520
Hexachloro-1,3-butadiene	U		3.37	10.0	10	02/20/2024 01:47	WG2229520
n-Hexane	U		7.49	100	10	02/20/2024 01:47	WG2229520
Isopropylbenzene	3.71	E4	1.05	10.0	10	02/20/2024 01:47	WG2229520
p-Isopropyltoluene	U		1.20	10.0	10	02/20/2024 01:47	WG2229520
2-Butanone (MEK)	1260		11.9	100	10	02/20/2024 01:47	WG2229520
Methyl Cyclohexane	U		6.60	10.0	10	02/20/2024 01:47	WG2229520

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Methylene Chloride	4.98	E4	4.30	50.0	10	02/20/2024 01:47	WG2229520
4-Methyl-2-pentanone (MIBK)	183		4.78	100	10	02/20/2024 01:47	WG2229520
Methyl tert-butyl ether	U		1.01	10.0	10	02/20/2024 01:47	WG2229520
Naphthalene	U		10.0	50.0	10	02/20/2024 01:47	WG2229520
Propene	U		9.36	25.0	10	02/20/2024 01:47	WG2229520
n-Propylbenzene	U		0.993	10.0	10	02/20/2024 01:47	WG2229520
Styrene	U	L2	1.18	10.0	10	02/20/2024 01:47	WG2229520
1,1,1,2-Tetrachloroethane	U		1.47	10.0	10	02/20/2024 01:47	WG2229520
1,1,2,2-Tetrachloroethane	U		1.33	10.0	10	02/20/2024 01:47	WG2229520
1,1,2-Trichlorotrifluoroethane	U		1.80	10.0	10	02/20/2024 01:47	WG2229520
Tetrachloroethene	U		3.00	10.0	10	02/20/2024 01:47	WG2229520
Toluene	U		2.78	10.0	10	02/20/2024 01:47	WG2229520
1,2,3-Trichlorobenzene	U		2.30	10.0	10	02/20/2024 01:47	WG2229520
1,2,4-Trichlorobenzene	U		4.81	10.0	10	02/20/2024 01:47	WG2229520
1,1,1-Trichloroethane	U		1.49	10.0	10	02/20/2024 01:47	WG2229520
1,1,2-Trichloroethane	U		1.58	10.0	10	02/20/2024 01:47	WG2229520
Trichloroethene	76.6		1.90	10.0	10	02/20/2024 01:47	WG2229520
Trichlorofluoromethane	U		1.60	50.0	10	02/20/2024 01:47	WG2229520
1,2,3-Trichloropropane	U		2.37	25.0	10	02/20/2024 01:47	WG2229520
1,2,4-Trimethylbenzene	U		3.22	10.0	10	02/20/2024 01:47	WG2229520
1,2,3-Trimethylbenzene	U		1.04	10.0	10	02/20/2024 01:47	WG2229520
1,3,5-Trimethylbenzene	U		1.04	10.0	10	02/20/2024 01:47	WG2229520
Vinyl chloride	U		2.34	10.0	10	02/20/2024 01:47	WG2229520
Xylenes, Total	U		1.74	30.0	10	02/20/2024 01:47	WG2229520
(S) Toluene-d8	*124*	S10		*80.0-120*		*02/20/2024 01:47*	WG2229520
(S) 4-Bromofluorobenzene	*114*			*77.0-126*		*02/20/2024 01:47*	WG2229520
(S) 1,2-Dichloroethane-d4	*88.6*			*70.0-130*		*02/20/2024 01:47*	WG2229520

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
1,4-Dioxane	U		0.597	3.00	1	02/20/2024 16:47	WG2230157
(S) Toluene-d8	*105*			*77.0-127*		*02/20/2024 16:47*	WG2230157

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	139000		1500	20000	5000	02/21/2024 06:33	WG2228552

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		113	500	10	02/17/2024 15:59	WG2228368
Acrolein	U		25.4	500	10	02/17/2024 15:59	WG2228368
Acrylonitrile	U		6.71	100	10	02/17/2024 15:59	WG2228368
Benzene	1.02	E4	0.941	10.0	10	02/17/2024 15:59	WG2228368
Bromobenzene	U		1.18	10.0	10	02/17/2024 15:59	WG2228368
Bromodichloromethane	U		1.36	10.0	10	02/17/2024 15:59	WG2228368
Bromoform	U		1.29	10.0	10	02/17/2024 15:59	WG2228368
Bromomethane	U		6.05	50.0	10	02/17/2024 15:59	WG2228368
1,3-Butadiene	U		2.99	20.0	10	02/17/2024 15:59	WG2228368
n-Butylbenzene	U		1.57	10.0	10	02/17/2024 15:59	WG2228368
sec-Butylbenzene	U		1.25	10.0	10	02/17/2024 15:59	WG2228368
tert-Butylbenzene	U		1.27	10.0	10	02/17/2024 15:59	WG2228368
Carbon tetrachloride	U		1.28	10.0	10	02/17/2024 15:59	WG2228368
Carbon disulfide	U		0.962	10.0	10	02/17/2024 15:59	WG2228368
Chlorobenzene	U		1.16	10.0	10	02/17/2024 15:59	WG2228368
Chlorodibromomethane	U		1.40	10.0	10	02/17/2024 15:59	WG2228368
Chloroethane	U		1.92	50.0	10	02/17/2024 15:59	WG2228368
Chloroform	1.55	E4	1.11	50.0	10	02/17/2024 15:59	WG2228368
Chloromethane	U		9.60	25.0	10	02/17/2024 15:59	WG2228368
Cyclohexane	U		1.88	10.0	10	02/17/2024 15:59	WG2228368
2-Chlorotoluene	U		1.06	10.0	10	02/17/2024 15:59	WG2228368
4-Chlorotoluene	U		1.14	10.0	10	02/17/2024 15:59	WG2228368
1,2-Dibromo-3-Chloropropane	U		2.76	50.0	10	02/17/2024 15:59	WG2228368
1,2-Dibromoethane	U		1.26	10.0	10	02/17/2024 15:59	WG2228368
Dibromomethane	U		1.22	10.0	10	02/17/2024 15:59	WG2228368
1,2-Dichlorobenzene	U		1.07	10.0	10	02/17/2024 15:59	WG2228368
1,3-Dichlorobenzene	U		1.10	10.0	10	02/17/2024 15:59	WG2228368
1,4-Dichlorobenzene	U		1.20	10.0	10	02/17/2024 15:59	WG2228368
Dichlorodifluoromethane	U		3.74	50.0	10	02/17/2024 15:59	WG2228368
1,1-Dichloroethane	U		1.00	10.0	10	02/17/2024 15:59	WG2228368
1,2-Dichloroethane	U		0.819	10.0	10	02/17/2024 15:59	WG2228368
1,1-Dichloroethene	52.6		1.88	10.0	10	02/17/2024 15:59	WG2228368
cis-1,2-Dichloroethene	U		1.26	10.0	10	02/17/2024 15:59	WG2228368
trans-1,2-Dichloroethene	U		1.49	10.0	10	02/17/2024 15:59	WG2228368
1,2-Dichloropropane	U		1.49	10.0	10	02/17/2024 15:59	WG2228368
1,1-Dichloropropene	U		1.42	10.0	10	02/17/2024 15:59	WG2228368
1,3-Dichloropropane	U		1.10	10.0	10	02/17/2024 15:59	WG2228368
cis-1,3-Dichloropropene	U		1.11	10.0	10	02/17/2024 15:59	WG2228368
trans-1,3-Dichloropropene	U		1.18	10.0	10	02/17/2024 15:59	WG2228368
2,2-Dichloropropane	U		1.61	10.0	10	02/17/2024 15:59	WG2228368
Dicyclopentadiene	U		2.53	10.0	10	02/17/2024 15:59	WG2228368
Di-isopropyl ether	U		1.05	10.0	10	02/17/2024 15:59	WG2228368
Ethylbenzene	U		1.37	10.0	10	02/17/2024 15:59	WG2228368
4-Ethyltoluene	U		2.08	10.0	10	02/17/2024 15:59	WG2228368
Hexachloro-1,3-butadiene	U		3.37	10.0	10	02/17/2024 15:59	WG2228368
n-Hexane	U		7.49	100	10	02/17/2024 15:59	WG2228368
Isopropylbenzene	U		1.05	10.0	10	02/17/2024 15:59	WG2228368
p-Isopropyltoluene	U		1.20	10.0	10	02/17/2024 15:59	WG2228368
2-Butanone (MEK)	U		11.9	100	10	02/17/2024 15:59	WG2228368
Methyl Cyclohexane	U		6.60	10.0	10	02/17/2024 15:59	WG2228368

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Methylene Chloride	U		4.30	50.0	10	02/20/2024 02:10	WG2229520
4-Methyl-2-pentanone (MIBK)	U		4.78	100	10	02/17/2024 15:59	WG2228368
Methyl tert-butyl ether	U		1.01	10.0	10	02/17/2024 15:59	WG2228368
Naphthalene	U		10.0	50.0	10	02/17/2024 15:59	WG2228368
Propene	U		9.36	25.0	10	02/17/2024 15:59	WG2228368
n-Propylbenzene	U		0.993	10.0	10	02/17/2024 15:59	WG2228368
Styrene	U		1.18	10.0	10	02/17/2024 15:59	WG2228368
1,1,1,2-Tetrachloroethane	U		1.47	10.0	10	02/17/2024 15:59	WG2228368
1,1,2,2-Tetrachloroethane	U		1.33	10.0	10	02/17/2024 15:59	WG2228368
1,1,2-Trichlorotrifluoroethane	U		1.80	10.0	10	02/17/2024 15:59	WG2228368
Tetrachloroethene	U		3.00	10.0	10	02/17/2024 15:59	WG2228368
Toluene	U		2.78	10.0	10	02/17/2024 15:59	WG2228368
1,2,3-Trichlorobenzene	U		2.30	10.0	10	02/17/2024 15:59	WG2228368
1,2,4-Trichlorobenzene	U		4.81	10.0	10	02/17/2024 15:59	WG2228368
1,1,1-Trichloroethane	U		1.49	10.0	10	02/17/2024 15:59	WG2228368
1,1,2-Trichloroethane	U		1.58	10.0	10	02/17/2024 15:59	WG2228368
Trichloroethene	511		1.90	10.0	10	02/17/2024 15:59	WG2228368
Trichlorofluoromethane	U		1.60	50.0	10	02/17/2024 15:59	WG2228368
1,2,3-Trichloropropane	U		2.37	25.0	10	02/17/2024 15:59	WG2228368
1,2,4-Trimethylbenzene	U		3.22	10.0	10	02/17/2024 15:59	WG2228368
1,2,3-Trimethylbenzene	U		1.04	10.0	10	02/17/2024 15:59	WG2228368
1,3,5-Trimethylbenzene	U		1.04	10.0	10	02/17/2024 15:59	WG2228368
Vinyl chloride	U		2.34	10.0	10	02/17/2024 15:59	WG2228368
Xylenes, Total	U		1.74	30.0	10	02/17/2024 15:59	WG2228368
(S) Toluene-d8	*103*			*80.0-120*		*02/17/2024 15:59*	WG2228368
(S) Toluene-d8	*98.9*			*80.0-120*		*02/20/2024 02:10*	WG2229520
(S) 4-Bromofluorobenzene	*93.0*			*77.0-126*		*02/17/2024 15:59*	WG2228368
(S) 4-Bromofluorobenzene	*94.3*			*77.0-126*		*02/20/2024 02:10*	WG2229520
(S) 1,2-Dichloroethane-d4	*102*			*70.0-130*		*02/17/2024 15:59*	WG2228368
(S) 1,2-Dichloroethane-d4	*90.9*			*70.0-130*		*02/20/2024 02:10*	WG2229520

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
1,4-Dioxane	U		0.597	3.00	1	02/20/2024 17:09	WG2230157
(S) Toluene-d8	*99.4*			*77.0-127*		*02/20/2024 17:09*	WG2230157

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	24000		300	4000	1000	02/21/2024 07:01	WG2228552

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		11.3	50.0	1	02/17/2024 13:56	WG2228368
Acrolein	U		2.54	50.0	1	02/17/2024 13:56	WG2228368
Acrylonitrile	U		0.671	10.0	1	02/17/2024 13:56	WG2228368
Benzene	U		0.0941	1.00	1	02/17/2024 13:56	WG2228368
Bromobenzene	U		0.118	1.00	1	02/17/2024 13:56	WG2228368
Bromodichloromethane	U		0.136	1.00	1	02/17/2024 13:56	WG2228368
Bromoform	0.241	E4	0.129	1.00	1	02/17/2024 13:56	WG2228368
Bromomethane	U		0.605	5.00	1	02/17/2024 13:56	WG2228368
1,3-Butadiene	U		0.299	2.00	1	02/17/2024 13:56	WG2228368
n-Butylbenzene	U		0.157	1.00	1	02/17/2024 13:56	WG2228368
sec-Butylbenzene	U		0.125	1.00	1	02/17/2024 13:56	WG2228368
tert-Butylbenzene	U		0.127	1.00	1	02/17/2024 13:56	WG2228368
Carbon tetrachloride	U		0.128	1.00	1	02/17/2024 13:56	WG2228368
Carbon disulfide	U		0.0962	1.00	1	02/17/2024 13:56	WG2228368
Chlorobenzene	U		0.116	1.00	1	02/17/2024 13:56	WG2228368
Chlorodibromomethane	U		0.140	1.00	1	02/17/2024 13:56	WG2228368
Chloroethane	U		0.192	5.00	1	02/17/2024 13:56	WG2228368
Chloroform	U		0.111	5.00	1	02/17/2024 13:56	WG2228368
Chloromethane	U		0.960	2.50	1	02/17/2024 13:56	WG2228368
Cyclohexane	U		0.188	1.00	1	02/17/2024 13:56	WG2228368
2-Chlorotoluene	U		0.106	1.00	1	02/17/2024 13:56	WG2228368
4-Chlorotoluene	U		0.114	1.00	1	02/17/2024 13:56	WG2228368
1,2-Dibromo-3-Chloropropane	U		0.276	5.00	1	02/17/2024 13:56	WG2228368
1,2-Dibromoethane	U		0.126	1.00	1	02/17/2024 13:56	WG2228368
Dibromomethane	U		0.122	1.00	1	02/17/2024 13:56	WG2228368
1,2-Dichlorobenzene	U		0.107	1.00	1	02/17/2024 13:56	WG2228368
1,3-Dichlorobenzene	U		0.110	1.00	1	02/17/2024 13:56	WG2228368
1,4-Dichlorobenzene	U		0.120	1.00	1	02/17/2024 13:56	WG2228368
Dichlorodifluoromethane	U		0.374	5.00	1	02/17/2024 13:56	WG2228368
1,1-Dichloroethane	U		0.100	1.00	1	02/17/2024 13:56	WG2228368
1,2-Dichloroethane	U		0.0819	1.00	1	02/17/2024 13:56	WG2228368
1,1-Dichloroethene	4.80		0.188	1.00	1	02/17/2024 13:56	WG2228368
cis-1,2-Dichloroethene	U		0.126	1.00	1	02/17/2024 13:56	WG2228368
trans-1,2-Dichloroethene	U		0.149	1.00	1	02/17/2024 13:56	WG2228368
1,2-Dichloropropane	U		0.149	1.00	1	02/17/2024 13:56	WG2228368
1,1-Dichloropropene	U		0.142	1.00	1	02/17/2024 13:56	WG2228368
1,3-Dichloropropane	U		0.110	1.00	1	02/17/2024 13:56	WG2228368
cis-1,3-Dichloropropene	U		0.111	1.00	1	02/17/2024 13:56	WG2228368
trans-1,3-Dichloropropene	U		0.118	1.00	1	02/17/2024 13:56	WG2228368
2,2-Dichloropropane	U		0.161	1.00	1	02/17/2024 13:56	WG2228368
Dicyclopentadiene	U		0.253	1.00	1	02/17/2024 13:56	WG2228368
Di-isopropyl ether	U		0.105	1.00	1	02/17/2024 13:56	WG2228368
Ethylbenzene	U		0.137	1.00	1	02/17/2024 13:56	WG2228368
4-Ethyltoluene	U		0.208	1.00	1	02/17/2024 13:56	WG2228368
Hexachloro-1,3-butadiene	U		0.337	1.00	1	02/17/2024 13:56	WG2228368
n-Hexane	U		0.749	10.0	1	02/17/2024 13:56	WG2228368
Isopropylbenzene	U		0.105	1.00	1	02/17/2024 13:56	WG2228368
p-Isopropyltoluene	U		0.120	1.00	1	02/17/2024 13:56	WG2228368
2-Butanone (MEK)	U		1.19	10.0	1	02/17/2024 13:56	WG2228368
Methyl Cyclohexane	U		0.660	1.00	1	02/17/2024 13:56	WG2228368

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Methylene Chloride	U		0.430	5.00	1	02/17/2024 13:56	WG2228368
4-Methyl-2-pentanone (MIBK)	U		0.478	10.0	1	02/17/2024 13:56	WG2228368
Methyl tert-butyl ether	U		0.101	1.00	1	02/17/2024 13:56	WG2228368
Naphthalene	U		1.00	5.00	1	02/17/2024 13:56	WG2228368
Propene	U		0.936	2.50	1	02/17/2024 13:56	WG2228368
n-Propylbenzene	U		0.0993	1.00	1	02/17/2024 13:56	WG2228368
Styrene	U		0.118	1.00	1	02/17/2024 13:56	WG2228368
1,1,1,2-Tetrachloroethane	U		0.147	1.00	1	02/17/2024 13:56	WG2228368
1,1,2,2-Tetrachloroethane	U		0.133	1.00	1	02/17/2024 13:56	WG2228368
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00	1	02/17/2024 13:56	WG2228368
Tetrachloroethene	U		0.300	1.00	1	02/17/2024 13:56	WG2228368
Toluene	U		0.278	1.00	1	02/17/2024 13:56	WG2228368
1,2,3-Trichlorobenzene	U		0.230	1.00	1	02/17/2024 13:56	WG2228368
1,2,4-Trichlorobenzene	U		0.481	1.00	1	02/17/2024 13:56	WG2228368
1,1,1-Trichloroethane	U		0.149	1.00	1	02/17/2024 13:56	WG2228368
1,1,2-Trichloroethane	U		0.158	1.00	1	02/17/2024 13:56	WG2228368
Trichloroethene	13.6		0.190	1.00	1	02/17/2024 13:56	WG2228368
Trichlorofluoromethane	U		0.160	5.00	1	02/17/2024 13:56	WG2228368
1,2,3-Trichloropropane	U		0.237	2.50	1	02/17/2024 13:56	WG2228368
1,2,4-Trimethylbenzene	U		0.322	1.00	1	02/17/2024 13:56	WG2228368
1,2,3-Trimethylbenzene	U		0.104	1.00	1	02/17/2024 13:56	WG2228368
1,3,5-Trimethylbenzene	U		0.104	1.00	1	02/17/2024 13:56	WG2228368
Vinyl chloride	U		0.234	1.00	1	02/17/2024 13:56	WG2228368
Xylenes, Total	U		0.174	3.00	1	02/17/2024 13:56	WG2228368
(S) Toluene-d8	*101*			*80.0-120*		*02/17/2024 13:56*	WG2228368
(S) 4-Bromofluorobenzene	*87.0*			*77.0-126*		*02/17/2024 13:56*	WG2228368
(S) 1,2-Dichloroethane-d4	*107*			*70.0-130*		*02/17/2024 13:56*	WG2228368

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
1,4-Dioxane	31.4		0.597	3.00	1	02/20/2024 17:31	WG2230157
(S) Toluene-d8	*98.9*			*77.0-127*		*02/20/2024 17:31*	WG2230157

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	147000		1500	20000	5000	02/25/2024 21:43	WG2228552

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		113	500	10	02/17/2024 16:20	WG2228368
Acrolein	U		25.4	500	10	02/17/2024 16:20	WG2228368
Acrylonitrile	U		6.71	100	10	02/17/2024 16:20	WG2228368
Benzene	1.84	E4	0.941	10.0	10	02/17/2024 16:20	WG2228368
Bromobenzene	U		1.18	10.0	10	02/17/2024 16:20	WG2228368
Bromodichloromethane	U		1.36	10.0	10	02/17/2024 16:20	WG2228368
Bromoform	U		1.29	10.0	10	02/17/2024 16:20	WG2228368
Bromomethane	U		6.05	50.0	10	02/17/2024 16:20	WG2228368
1,3-Butadiene	U		2.99	20.0	10	02/17/2024 16:20	WG2228368
n-Butylbenzene	U		1.57	10.0	10	02/17/2024 16:20	WG2228368
sec-Butylbenzene	U		1.25	10.0	10	02/17/2024 16:20	WG2228368
tert-Butylbenzene	U		1.27	10.0	10	02/17/2024 16:20	WG2228368
Carbon tetrachloride	U		1.28	10.0	10	02/17/2024 16:20	WG2228368
Carbon disulfide	U		0.962	10.0	10	02/17/2024 16:20	WG2228368
Chlorobenzene	U		1.16	10.0	10	02/17/2024 16:20	WG2228368
Chlorodibromomethane	U		1.40	10.0	10	02/17/2024 16:20	WG2228368
Chloroethane	U		1.92	50.0	10	02/17/2024 16:20	WG2228368
Chloroform	2.01	E4	1.11	50.0	10	02/17/2024 16:20	WG2228368
Chloromethane	U		9.60	25.0	10	02/17/2024 16:20	WG2228368
Cyclohexane	U		1.88	10.0	10	02/17/2024 16:20	WG2228368
2-Chlorotoluene	U		1.06	10.0	10	02/17/2024 16:20	WG2228368
4-Chlorotoluene	U		1.14	10.0	10	02/17/2024 16:20	WG2228368
1,2-Dibromo-3-Chloropropane	U		2.76	50.0	10	02/17/2024 16:20	WG2228368
1,2-Dibromoethane	U		1.26	10.0	10	02/17/2024 16:20	WG2228368
Dibromomethane	U		1.22	10.0	10	02/17/2024 16:20	WG2228368
1,2-Dichlorobenzene	U		1.07	10.0	10	02/17/2024 16:20	WG2228368
1,3-Dichlorobenzene	U		1.10	10.0	10	02/17/2024 16:20	WG2228368
1,4-Dichlorobenzene	U		1.20	10.0	10	02/17/2024 16:20	WG2228368
Dichlorodifluoromethane	U		3.74	50.0	10	02/17/2024 16:20	WG2228368
1,1-Dichloroethane	U		1.00	10.0	10	02/17/2024 16:20	WG2228368
1,2-Dichloroethane	U		0.819	10.0	10	02/17/2024 16:20	WG2228368
1,1-Dichloroethene	81.7		1.88	10.0	10	02/17/2024 16:20	WG2228368
cis-1,2-Dichloroethene	2.28	E4	1.26	10.0	10	02/17/2024 16:20	WG2228368
trans-1,2-Dichloroethene	U		1.49	10.0	10	02/17/2024 16:20	WG2228368
1,2-Dichloropropane	U		1.49	10.0	10	02/17/2024 16:20	WG2228368
1,1-Dichloropropene	U		1.42	10.0	10	02/17/2024 16:20	WG2228368
1,3-Dichloropropane	U		1.10	10.0	10	02/17/2024 16:20	WG2228368
cis-1,3-Dichloropropene	U		1.11	10.0	10	02/17/2024 16:20	WG2228368
trans-1,3-Dichloropropene	U		1.18	10.0	10	02/17/2024 16:20	WG2228368
2,2-Dichloropropane	U		1.61	10.0	10	02/17/2024 16:20	WG2228368
Dicyclopentadiene	U		2.53	10.0	10	02/17/2024 16:20	WG2228368
Di-isopropyl ether	U		1.05	10.0	10	02/17/2024 16:20	WG2228368
Ethylbenzene	U		1.37	10.0	10	02/17/2024 16:20	WG2228368
4-Ethyltoluene	U		2.08	10.0	10	02/17/2024 16:20	WG2228368
Hexachloro-1,3-butadiene	U		3.37	10.0	10	02/17/2024 16:20	WG2228368
n-Hexane	U		7.49	100	10	02/17/2024 16:20	WG2228368
Isopropylbenzene	U		1.05	10.0	10	02/17/2024 16:20	WG2228368
p-Isopropyltoluene	U		1.20	10.0	10	02/17/2024 16:20	WG2228368
2-Butanone (MEK)	U		11.9	100	10	02/17/2024 16:20	WG2228368
Methyl Cyclohexane	U		6.60	10.0	10	02/17/2024 16:20	WG2228368

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

Analyte	Result ug/l	Qualifier	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	Batch
Methylene Chloride	U		4.30	50.0	10	02/20/2024 02:33	WG2229520
4-Methyl-2-pentanone (MIBK)	U		4.78	100	10	02/17/2024 16:20	WG2228368
Methyl tert-butyl ether	U		1.01	10.0	10	02/17/2024 16:20	WG2228368
Naphthalene	U		10.0	50.0	10	02/17/2024 16:20	WG2228368
Propene	U		9.36	25.0	10	02/17/2024 16:20	WG2228368
n-Propylbenzene	U		0.993	10.0	10	02/17/2024 16:20	WG2228368
Styrene	U		1.18	10.0	10	02/17/2024 16:20	WG2228368
1,1,1,2-Tetrachloroethane	U		1.47	10.0	10	02/17/2024 16:20	WG2228368
1,1,2,2-Tetrachloroethane	U		1.33	10.0	10	02/17/2024 16:20	WG2228368
1,1,2-Trichlorotrifluoroethane	U		1.80	10.0	10	02/17/2024 16:20	WG2228368
Tetrachloroethene	U		3.00	10.0	10	02/17/2024 16:20	WG2228368
Toluene	U		2.78	10.0	10	02/17/2024 16:20	WG2228368
1,2,3-Trichlorobenzene	U		2.30	10.0	10	02/17/2024 16:20	WG2228368
1,2,4-Trichlorobenzene	U		4.81	10.0	10	02/17/2024 16:20	WG2228368
1,1,1-Trichloroethane	U		1.49	10.0	10	02/17/2024 16:20	WG2228368
1,1,2-Trichloroethane	2.31	E4	1.58	10.0	10	02/17/2024 16:20	WG2228368
Trichloroethene	829		1.90	10.0	10	02/17/2024 16:20	WG2228368
Trichlorofluoromethane	U		1.60	50.0	10	02/17/2024 16:20	WG2228368
1,2,3-Trichloropropane	U		2.37	25.0	10	02/17/2024 16:20	WG2228368
1,2,4-Trimethylbenzene	U		3.22	10.0	10	02/17/2024 16:20	WG2228368
1,2,3-Trimethylbenzene	U		1.04	10.0	10	02/17/2024 16:20	WG2228368
1,3,5-Trimethylbenzene	U		1.04	10.0	10	02/17/2024 16:20	WG2228368
Vinyl chloride	U		2.34	10.0	10	02/17/2024 16:20	WG2228368
Xylenes, Total	U		1.74	30.0	10	02/17/2024 16:20	WG2228368
(S) Toluene-d8	*103*			*80.0-120*		*02/17/2024 16:20*	WG2228368
(S) Toluene-d8	*106*			*80.0-120*		*02/20/2024 02:33*	WG2229520
(S) 4-Bromofluorobenzene	*88.7*			*77.0-126*		*02/17/2024 16:20*	WG2228368
(S) 4-Bromofluorobenzene	*106*			*77.0-126*		*02/20/2024 02:33*	WG2229520
(S) 1,2-Dichloroethane-d4	*102*			*70.0-130*		*02/17/2024 16:20*	WG2228368
(S) 1,2-Dichloroethane-d4	*93.4*			*70.0-130*		*02/20/2024 02:33*	WG2229520

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result ug/l	Qualifier	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	Batch
1,4-Dioxane	286		0.597	3.00	1	02/20/2024 17:53	WG2230157
(S) Toluene-d8	*98.8*			*77.0-127*		*02/20/2024 17:53*	WG2230157

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	800000		3000	40000	10000	02/25/2024 00:24	WG2228552

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		11.3	50.0	1	02/17/2024 14:16	WG2228368
Acrolein	U		2.54	50.0	1	02/17/2024 14:16	WG2228368
Acrylonitrile	U		0.671	10.0	1	02/17/2024 14:16	WG2228368
Benzene	U		235	2500	2500	02/20/2024 02:56	WG2229520
Bromobenzene	U		0.118	1.00	1	02/17/2024 14:16	WG2228368
Bromodichloromethane	U		0.136	1.00	1	02/17/2024 14:16	WG2228368
Bromoform	2.35		0.129	1.00	1	02/17/2024 14:16	WG2228368
Bromomethane	U		0.605	5.00	1	02/17/2024 14:16	WG2228368
1,3-Butadiene	U		0.299	2.00	1	02/17/2024 14:16	WG2228368
n-Butylbenzene	U		0.157	1.00	1	02/17/2024 14:16	WG2228368
sec-Butylbenzene	U		0.125	1.00	1	02/17/2024 14:16	WG2228368
tert-Butylbenzene	U		0.127	1.00	1	02/17/2024 14:16	WG2228368
Carbon tetrachloride	U		0.128	1.00	1	02/17/2024 14:16	WG2228368
Carbon disulfide	0.264	E4	0.0962	1.00	1	02/17/2024 14:16	WG2228368
Chlorobenzene	2.09		0.116	1.00	1	02/17/2024 14:16	WG2228368
Chlorodibromomethane	0.573	E4	0.140	1.00	1	02/17/2024 14:16	WG2228368
Chloroethane	U		0.192	5.00	1	02/17/2024 14:16	WG2228368
Chloroform	79.2		0.111	5.00	1	02/17/2024 14:16	WG2228368
Chloromethane	U		0.960	2.50	1	02/17/2024 14:16	WG2228368
Cyclohexane	U		0.188	1.00	1	02/17/2024 14:16	WG2228368
2-Chlorotoluene	U		0.106	1.00	1	02/17/2024 14:16	WG2228368
4-Chlorotoluene	U		0.114	1.00	1	02/17/2024 14:16	WG2228368
1,2-Dibromo-3-Chloropropane	U		0.276	5.00	1	02/17/2024 14:16	WG2228368
1,2-Dibromoethane	U		0.126	1.00	1	02/17/2024 14:16	WG2228368
Dibromomethane	U		0.122	1.00	1	02/17/2024 14:16	WG2228368
1,2-Dichlorobenzene	3.02		0.107	1.00	1	02/17/2024 14:16	WG2228368
1,3-Dichlorobenzene	0.287	E4	0.110	1.00	1	02/17/2024 14:16	WG2228368
1,4-Dichlorobenzene	0.719	E4	0.120	1.00	1	02/17/2024 14:16	WG2228368
Dichlorodifluoromethane	U		0.374	5.00	1	02/17/2024 14:16	WG2228368
1,1-Dichloroethane	53.2		0.100	1.00	1	02/17/2024 14:16	WG2228368
1,2-Dichloroethane	27.6		0.0819	1.00	1	02/17/2024 14:16	WG2228368
1,1-Dichloroethene	5240		470	2500	2500	02/20/2024 02:56	WG2229520
cis-1,2-Dichloroethene	13.4		0.126	1.00	1	02/17/2024 14:16	WG2228368
trans-1,2-Dichloroethene	9.46		0.149	1.00	1	02/17/2024 14:16	WG2228368
1,2-Dichloropropane	U		0.149	1.00	1	02/17/2024 14:16	WG2228368
1,1-Dichloropropene	U		0.142	1.00	1	02/17/2024 14:16	WG2228368
1,3-Dichloropropane	U		0.110	1.00	1	02/17/2024 14:16	WG2228368
cis-1,3-Dichloropropene	U		0.111	1.00	1	02/17/2024 14:16	WG2228368
trans-1,3-Dichloropropene	U		0.118	1.00	1	02/17/2024 14:16	WG2228368
2,2-Dichloropropane	U		0.161	1.00	1	02/17/2024 14:16	WG2228368
Dicyclopentadiene	U		0.253	1.00	1	02/17/2024 14:16	WG2228368
Di-isopropyl ether	0.432	E4	0.105	1.00	1	02/17/2024 14:16	WG2228368
Ethylbenzene	0.721	E4	0.137	1.00	1	02/17/2024 14:16	WG2228368
4-Ethyltoluene	0.253	E4	0.208	1.00	1	02/17/2024 14:16	WG2228368
Hexachloro-1,3-butadiene	U		0.337	1.00	1	02/17/2024 14:16	WG2228368
n-Hexane	U		0.749	10.0	1	02/17/2024 14:16	WG2228368
Isopropylbenzene	0.531	E4	0.105	1.00	1	02/17/2024 14:16	WG2228368
p-Isopropyltoluene	U		0.120	1.00	1	02/17/2024 14:16	WG2228368
2-Butanone (MEK)	U		1.19	10.0	1	02/17/2024 14:16	WG2228368
Methyl Cyclohexane	U		0.660	1.00	1	02/17/2024 14:16	WG2228368

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Methylene Chloride	83000		1080	12500	2500	02/20/2024 02:56	WG2229520
4-Methyl-2-pentanone (MIBK)	U		0.478	10.0	1	02/17/2024 14:16	WG2228368
Methyl tert-butyl ether	U		0.101	1.00	1	02/17/2024 14:16	WG2228368
Naphthalene	1.17	E4	1.00	5.00	1	02/17/2024 14:16	WG2228368
Propene	3.71		0.936	2.50	1	02/17/2024 14:16	WG2228368
n-Propylbenzene	U		0.0993	1.00	1	02/17/2024 14:16	WG2228368
Styrene	U		0.118	1.00	1	02/17/2024 14:16	WG2228368
1,1,1,2-Tetrachloroethane	U		0.147	1.00	1	02/17/2024 14:16	WG2228368
1,1,2,2-Tetrachloroethane	U		0.133	1.00	1	02/17/2024 14:16	WG2228368
1,1,2-Trichlorotrifluoroethane	8.85		0.180	1.00	1	02/17/2024 14:16	WG2228368
Tetrachloroethene	57.4		0.300	1.00	1	02/17/2024 14:16	WG2228368
Toluene	17.1		0.278	1.00	1	02/17/2024 14:16	WG2228368
1,2,3-Trichlorobenzene	U		0.230	1.00	1	02/17/2024 14:16	WG2228368
1,2,4-Trichlorobenzene	U		0.481	1.00	1	02/17/2024 14:16	WG2228368
1,1,1-Trichloroethane	U		0.149	1.00	1	02/17/2024 14:16	WG2228368
1,1,2-Trichloroethane	67.7		0.158	1.00	1	02/17/2024 14:16	WG2228368
Trichloroethene	77500		475	2500	2500	02/20/2024 02:56	WG2229520
Trichlorofluoromethane	U		0.160	5.00	1	02/17/2024 14:16	WG2228368
1,2,3-Trichloropropane	0.276	E4	0.237	2.50	1	02/17/2024 14:16	WG2228368
1,2,4-Trimethylbenzene	0.675	E4	0.322	1.00	1	02/17/2024 14:16	WG2228368
1,2,3-Trimethylbenzene	1.57		0.104	1.00	1	02/17/2024 14:16	WG2228368
1,3,5-Trimethylbenzene	0.220	E4	0.104	1.00	1	02/17/2024 14:16	WG2228368
Vinyl chloride	0.941	E4	0.234	1.00	1	02/17/2024 14:16	WG2228368
Xylenes, Total	99.1		0.174	3.00	1	02/17/2024 14:16	WG2228368
(S) Toluene-d8	*95.8*			*80.0-120*		*02/17/2024 14:16*	WG2228368
(S) Toluene-d8	*107*			*80.0-120*		*02/20/2024 02:56*	WG2229520
(S) 4-Bromofluorobenzene	*93.9*			*77.0-126*		*02/17/2024 14:16*	WG2228368
(S) 4-Bromofluorobenzene	*104*			*77.0-126*		*02/20/2024 02:56*	WG2229520
(S) 1,2-Dichloroethane-d4	*105*			*70.0-130*		*02/17/2024 14:16*	WG2228368
(S) 1,2-Dichloroethane-d4	*86.3*			*70.0-130*		*02/20/2024 02:56*	WG2229520

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
1,4-Dioxane	4020		11.9	60.0	20	02/21/2024 15:55	WG2230690
(S) Toluene-d8	*99.0*			*77.0-127*		*02/21/2024 15:55*	WG2230690

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	11600		150	2000	500	02/25/2024 00:52	WG2228552

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		226	1000	20	02/20/2024 03:19	WG2229520
Acrolein	U		50.8	1000	20	02/20/2024 03:19	WG2229520
Acrylonitrile	U		13.4	200	20	02/20/2024 03:19	WG2229520
Benzene	U		1.88	20.0	20	02/20/2024 03:19	WG2229520
Bromobenzene	U		2.36	20.0	20	02/20/2024 03:19	WG2229520
Bromodichloromethane	U		2.72	20.0	20	02/20/2024 03:19	WG2229520
Bromoform	U	L2	2.58	20.0	20	02/20/2024 03:19	WG2229520
Bromomethane	U		12.1	100	20	02/20/2024 03:19	WG2229520
1,3-Butadiene	U		5.98	40.0	20	02/20/2024 03:19	WG2229520
n-Butylbenzene	U		3.14	20.0	20	02/20/2024 03:19	WG2229520
sec-Butylbenzene	U		2.50	20.0	20	02/20/2024 03:19	WG2229520
tert-Butylbenzene	U		2.54	20.0	20	02/20/2024 03:19	WG2229520
Carbon tetrachloride	U		2.56	20.0	20	02/20/2024 03:19	WG2229520
Carbon disulfide	U		1.92	20.0	20	02/20/2024 03:19	WG2229520
Chlorobenzene	U		2.32	20.0	20	02/20/2024 03:19	WG2229520
Chlorodibromomethane	U		2.80	20.0	20	02/20/2024 03:19	WG2229520
Chloroethane	U		3.84	100	20	02/20/2024 03:19	WG2229520
Chloroform	U		2.22	100	20	02/20/2024 03:19	WG2229520
Chloromethane	U		19.2	50.0	20	02/20/2024 03:19	WG2229520
Cyclohexane	U		3.76	20.0	20	02/20/2024 03:19	WG2229520
2-Chlorotoluene	U		2.12	20.0	20	02/20/2024 03:19	WG2229520
4-Chlorotoluene	U		2.28	20.0	20	02/20/2024 03:19	WG2229520
1,2-Dibromo-3-Chloropropane	U		5.52	100	20	02/20/2024 03:19	WG2229520
1,2-Dibromoethane	U		2.52	20.0	20	02/20/2024 03:19	WG2229520
Dibromomethane	U		2.44	20.0	20	02/20/2024 03:19	WG2229520
1,2-Dichlorobenzene	U	L2	2.14	20.0	20	02/20/2024 03:19	WG2229520
1,3-Dichlorobenzene	U	L2	2.20	20.0	20	02/20/2024 03:19	WG2229520
1,4-Dichlorobenzene	U	L2	2.40	20.0	20	02/20/2024 03:19	WG2229520
Dichlorodifluoromethane	U		7.48	100	20	02/20/2024 03:19	WG2229520
1,1-Dichloroethane	U		2.00	20.0	20	02/20/2024 03:19	WG2229520
1,2-Dichloroethane	U		1.64	20.0	20	02/20/2024 03:19	WG2229520
1,1-Dichloroethene	U		3.76	20.0	20	02/20/2024 03:19	WG2229520
cis-1,2-Dichloroethene	U		2.52	20.0	20	02/20/2024 03:19	WG2229520
trans-1,2-Dichloroethene	U		2.98	20.0	20	02/20/2024 03:19	WG2229520
1,2-Dichloropropane	U		2.98	20.0	20	02/20/2024 03:19	WG2229520
1,1-Dichloropropene	U		2.84	20.0	20	02/20/2024 03:19	WG2229520
1,3-Dichloropropane	U		2.20	20.0	20	02/20/2024 03:19	WG2229520
cis-1,3-Dichloropropene	U		2.22	20.0	20	02/20/2024 03:19	WG2229520
trans-1,3-Dichloropropene	U		2.36	20.0	20	02/20/2024 03:19	WG2229520
2,2-Dichloropropane	U		3.22	20.0	20	02/20/2024 03:19	WG2229520
Dicyclopentadiene	U		5.06	20.0	20	02/20/2024 03:19	WG2229520
Di-isopropyl ether	U		2.10	20.0	20	02/20/2024 03:19	WG2229520
Ethylbenzene	U		2.74	20.0	20	02/20/2024 03:19	WG2229520
4-Ethyltoluene	U		4.16	20.0	20	02/20/2024 03:19	WG2229520
Hexachloro-1,3-butadiene	U		6.74	20.0	20	02/20/2024 03:19	WG2229520
n-Hexane	U		15.0	200	20	02/20/2024 03:19	WG2229520
Isopropylbenzene	U		2.10	20.0	20	02/20/2024 03:19	WG2229520
p-Isopropyltoluene	U		2.40	20.0	20	02/20/2024 03:19	WG2229520
2-Butanone (MEK)	U		23.8	200	20	02/20/2024 03:19	WG2229520
Methyl Cyclohexane	U		13.2	20.0	20	02/20/2024 03:19	WG2229520

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

Analyte	Result ug/l	Qualifier	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	Batch
Methylene Chloride	U		8.60	100	20	02/20/2024 03:19	WG2229520
4-Methyl-2-pentanone (MIBK)	U		9.56	200	20	02/20/2024 03:19	WG2229520
Methyl tert-butyl ether	U		2.02	20.0	20	02/20/2024 03:19	WG2229520
Naphthalene	U		20.0	100	20	02/20/2024 03:19	WG2229520
Propene	U		18.7	50.0	20	02/20/2024 03:19	WG2229520
n-Propylbenzene	U		1.99	20.0	20	02/20/2024 03:19	WG2229520
Styrene	U	L2	2.36	20.0	20	02/20/2024 03:19	WG2229520
1,1,1,2-Tetrachloroethane	U		2.94	20.0	20	02/20/2024 03:19	WG2229520
1,1,2,2-Tetrachloroethane	U		2.66	20.0	20	02/20/2024 03:19	WG2229520
1,1,2-Trichlorotrifluoroethane	U		3.60	20.0	20	02/20/2024 03:19	WG2229520
Tetrachloroethene	U		6.00	20.0	20	02/20/2024 03:19	WG2229520
Toluene	U		5.56	20.0	20	02/20/2024 03:19	WG2229520
1,2,3-Trichlorobenzene	U		4.60	20.0	20	02/20/2024 03:19	WG2229520
1,2,4-Trichlorobenzene	U		9.62	20.0	20	02/20/2024 03:19	WG2229520
1,1,1-Trichloroethane	U		2.98	20.0	20	02/20/2024 03:19	WG2229520
1,1,2-Trichloroethane	U		3.16	20.0	20	02/20/2024 03:19	WG2229520
Trichloroethene	U		3.80	20.0	20	02/20/2024 03:19	WG2229520
Trichlorofluoromethane	U		3.20	100	20	02/20/2024 03:19	WG2229520
1,2,3-Trichloropropane	U		4.74	50.0	20	02/20/2024 03:19	WG2229520
1,2,4-Trimethylbenzene	U		6.44	20.0	20	02/20/2024 03:19	WG2229520
1,2,3-Trimethylbenzene	U		2.08	20.0	20	02/20/2024 03:19	WG2229520
1,3,5-Trimethylbenzene	U		2.08	20.0	20	02/20/2024 03:19	WG2229520
Vinyl chloride	U		4.68	20.0	20	02/20/2024 03:19	WG2229520
Xylenes, Total	U		3.48	60.0	20	02/20/2024 03:19	WG2229520
(S) Toluene-d8	*104*			*80.0-120*		*02/20/2024 03:19*	WG2229520
(S) 4-Bromofluorobenzene	*101*			*77.0-126*		*02/20/2024 03:19*	WG2229520
(S) 1,2-Dichloroethane-d4	*91.4*			*70.0-130*		*02/20/2024 03:19*	WG2229520

Sample Narrative:

L1705776-14 WG2229520: Elevated RI due to sample matrix/historical data.

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result ug/l	Qualifier	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	Batch
1,4-Dioxane	7.27		0.597	3.00	1	02/21/2024 14:29	WG2230690
(S) Toluene-d8	*98.3*			*77.0-127*		*02/21/2024 14:29*	WG2230690

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	U		0.300	4.00	1	02/25/2024 01:20	WG2228552

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		11.3	50.0	1	02/17/2024 14:37	WG2228368
Acrolein	U		2.54	50.0	1	02/17/2024 14:37	WG2228368
Acrylonitrile	U		0.671	10.0	1	02/17/2024 14:37	WG2228368
Benzene	U		0.0941	1.00	1	02/17/2024 14:37	WG2228368
Bromobenzene	U		0.118	1.00	1	02/17/2024 14:37	WG2228368
Bromodichloromethane	U		0.136	1.00	1	02/17/2024 14:37	WG2228368
Bromoform	U		0.129	1.00	1	02/17/2024 14:37	WG2228368
Bromomethane	U		0.605	5.00	1	02/17/2024 14:37	WG2228368
1,3-Butadiene	U		0.299	2.00	1	02/17/2024 14:37	WG2228368
n-Butylbenzene	U		0.157	1.00	1	02/17/2024 14:37	WG2228368
sec-Butylbenzene	U		0.125	1.00	1	02/17/2024 14:37	WG2228368
tert-Butylbenzene	U		0.127	1.00	1	02/17/2024 14:37	WG2228368
Carbon tetrachloride	U		0.128	1.00	1	02/17/2024 14:37	WG2228368
Carbon disulfide	U		0.0962	1.00	1	02/17/2024 14:37	WG2228368
Chlorobenzene	U		0.116	1.00	1	02/17/2024 14:37	WG2228368
Chlorodibromomethane	U		0.140	1.00	1	02/17/2024 14:37	WG2228368
Chloroethane	U		0.192	5.00	1	02/17/2024 14:37	WG2228368
Chloroform	U		0.111	5.00	1	02/17/2024 14:37	WG2228368
Chloromethane	U		0.960	2.50	1	02/17/2024 14:37	WG2228368
Cyclohexane	U		0.188	1.00	1	02/17/2024 14:37	WG2228368
2-Chlorotoluene	U		0.106	1.00	1	02/17/2024 14:37	WG2228368
4-Chlorotoluene	U		0.114	1.00	1	02/17/2024 14:37	WG2228368
1,2-Dibromo-3-Chloropropane	U		0.276	5.00	1	02/17/2024 14:37	WG2228368
1,2-Dibromoethane	U		0.126	1.00	1	02/17/2024 14:37	WG2228368
Dibromomethane	U		0.122	1.00	1	02/17/2024 14:37	WG2228368
1,2-Dichlorobenzene	U		0.107	1.00	1	02/17/2024 14:37	WG2228368
1,3-Dichlorobenzene	U		0.110	1.00	1	02/17/2024 14:37	WG2228368
1,4-Dichlorobenzene	U		0.120	1.00	1	02/17/2024 14:37	WG2228368
Dichlorodifluoromethane	U		0.374	5.00	1	02/17/2024 14:37	WG2228368
1,1-Dichloroethane	U		0.100	1.00	1	02/17/2024 14:37	WG2228368
1,2-Dichloroethane	U		0.0819	1.00	1	02/17/2024 14:37	WG2228368
1,1-Dichloroethene	U		0.188	1.00	1	02/20/2024 03:42	WG2229520
cis-1,2-Dichloroethene	0.507	E4	0.126	1.00	1	02/17/2024 14:37	WG2228368
trans-1,2-Dichloroethene	U		0.149	1.00	1	02/17/2024 14:37	WG2228368
1,2-Dichloropropane	U		0.149	1.00	1	02/17/2024 14:37	WG2228368
1,1-Dichloropropene	U		0.142	1.00	1	02/17/2024 14:37	WG2228368
1,3-Dichloropropane	U		0.110	1.00	1	02/17/2024 14:37	WG2228368
cis-1,3-Dichloropropene	U		0.111	1.00	1	02/17/2024 14:37	WG2228368
trans-1,3-Dichloropropene	U		0.118	1.00	1	02/17/2024 14:37	WG2228368
2,2-Dichloropropane	U		0.161	1.00	1	02/17/2024 14:37	WG2228368
Dicyclopentadiene	U		0.253	1.00	1	02/17/2024 14:37	WG2228368
Di-isopropyl ether	U		0.105	1.00	1	02/17/2024 14:37	WG2228368
Ethylbenzene	U		0.137	1.00	1	02/17/2024 14:37	WG2228368
4-Ethyltoluene	U		0.208	1.00	1	02/17/2024 14:37	WG2228368
Hexachloro-1,3-butadiene	U		0.337	1.00	1	02/17/2024 14:37	WG2228368
n-Hexane	U		0.749	10.0	1	02/17/2024 14:37	WG2228368
Isopropylbenzene	U		0.105	1.00	1	02/17/2024 14:37	WG2228368
p-Isopropyltoluene	U		0.120	1.00	1	02/17/2024 14:37	WG2228368
2-Butanone (MEK)	U		1.19	10.0	1	02/17/2024 14:37	WG2228368
Methyl Cyclohexane	U		0.660	1.00	1	02/17/2024 14:37	WG2228368

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Methylene Chloride	U		0.430	5.00	1	02/20/2024 03:42	WG2229520
4-Methyl-2-pentanone (MIBK)	U		0.478	10.0	1	02/17/2024 14:37	WG2228368
Methyl tert-butyl ether	U		0.101	1.00	1	02/17/2024 14:37	WG2228368
Naphthalene	U		1.00	5.00	1	02/17/2024 14:37	WG2228368
Propene	U		0.936	2.50	1	02/17/2024 14:37	WG2228368
n-Propylbenzene	U		0.0993	1.00	1	02/17/2024 14:37	WG2228368
Styrene	U		0.118	1.00	1	02/17/2024 14:37	WG2228368
1,1,1,2-Tetrachloroethane	U		0.147	1.00	1	02/17/2024 14:37	WG2228368
1,1,2,2-Tetrachloroethane	U		0.133	1.00	1	02/17/2024 14:37	WG2228368
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00	1	02/17/2024 14:37	WG2228368
Tetrachloroethene	U		0.300	1.00	1	02/17/2024 14:37	WG2228368
Toluene	U		0.278	1.00	1	02/17/2024 14:37	WG2228368
1,2,3-Trichlorobenzene	U		0.230	1.00	1	02/17/2024 14:37	WG2228368
1,2,4-Trichlorobenzene	U		0.481	1.00	1	02/17/2024 14:37	WG2228368
1,1,1-Trichloroethane	U		0.149	1.00	1	02/17/2024 14:37	WG2228368
1,1,2-Trichloroethane	U		0.158	1.00	1	02/17/2024 14:37	WG2228368
Trichloroethene	0.661	E4	0.190	1.00	1	02/20/2024 03:42	WG2229520
Trichlorofluoromethane	U		0.160	5.00	1	02/17/2024 14:37	WG2228368
1,2,3-Trichloropropane	U		0.237	2.50	1	02/17/2024 14:37	WG2228368
1,2,4-Trimethylbenzene	U		0.322	1.00	1	02/17/2024 14:37	WG2228368
1,2,3-Trimethylbenzene	U		0.104	1.00	1	02/17/2024 14:37	WG2228368
1,3,5-Trimethylbenzene	U		0.104	1.00	1	02/17/2024 14:37	WG2228368
Vinyl chloride	U		0.234	1.00	1	02/17/2024 14:37	WG2228368
Xylenes, Total	U		0.174	3.00	1	02/17/2024 14:37	WG2228368
(S) Toluene-d8	*90.1*			*80.0-120*		*02/17/2024 14:37*	WG2228368
(S) Toluene-d8	*109*			*80.0-120*		*02/20/2024 03:42*	WG2229520
(S) 4-Bromofluorobenzene	*79.9*			*77.0-126*		*02/17/2024 14:37*	WG2228368
(S) 4-Bromofluorobenzene	*102*			*77.0-126*		*02/20/2024 03:42*	WG2229520
(S) 1,2-Dichloroethane-d4	*101*			*70.0-130*		*02/17/2024 14:37*	WG2228368
(S) 1,2-Dichloroethane-d4	*84.6*			*70.0-130*		*02/20/2024 03:42*	WG2229520

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
1,4-Dioxane	U		0.597	3.00	1	02/21/2024 14:50	WG2230690
(S) Toluene-d8	*93.9*			*77.0-127*		*02/21/2024 14:50*	WG2230690

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	U		0.300	4.00	1	02/21/2024 10:44	WG2228552

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		56.5	250	5	02/17/2024 17:01	WG2228368
Acrolein	U		12.7	250	5	02/17/2024 17:01	WG2228368
Acrylonitrile	U		3.36	50.0	5	02/17/2024 17:01	WG2228368
Benzene	3.46	E4	0.471	5.00	5	02/17/2024 17:01	WG2228368
Bromobenzene	U		0.590	5.00	5	02/17/2024 17:01	WG2228368
Bromodichloromethane	U		0.680	5.00	5	02/17/2024 17:01	WG2228368
Bromoform	U		0.645	5.00	5	02/17/2024 17:01	WG2228368
Bromomethane	U		3.03	25.0	5	02/17/2024 17:01	WG2228368
1,3-Butadiene	U		1.49	10.0	5	02/17/2024 17:01	WG2228368
n-Butylbenzene	U		0.785	5.00	5	02/17/2024 17:01	WG2228368
sec-Butylbenzene	U		0.625	5.00	5	02/17/2024 17:01	WG2228368
tert-Butylbenzene	U		0.635	5.00	5	02/17/2024 17:01	WG2228368
Carbon tetrachloride	U		0.640	5.00	5	02/17/2024 17:01	WG2228368
Carbon disulfide	0.496	E4	0.481	5.00	5	02/17/2024 17:01	WG2228368
Chlorobenzene	U		0.580	5.00	5	02/17/2024 17:01	WG2228368
Chlorodibromomethane	U		0.700	5.00	5	02/17/2024 17:01	WG2228368
Chloroethane	U		0.960	25.0	5	02/17/2024 17:01	WG2228368
Chloroform	U		0.555	25.0	5	02/17/2024 17:01	WG2228368
Chloromethane	U		4.80	12.5	5	02/17/2024 17:01	WG2228368
Cyclohexane	U		0.940	5.00	5	02/17/2024 17:01	WG2228368
2-Chlorotoluene	U		0.530	5.00	5	02/17/2024 17:01	WG2228368
4-Chlorotoluene	U		0.570	5.00	5	02/17/2024 17:01	WG2228368
1,2-Dibromo-3-Chloropropane	U		1.38	25.0	5	02/17/2024 17:01	WG2228368
1,2-Dibromoethane	U		0.630	5.00	5	02/17/2024 17:01	WG2228368
Dibromomethane	U		0.610	5.00	5	02/17/2024 17:01	WG2228368
1,2-Dichlorobenzene	U		0.535	5.00	5	02/17/2024 17:01	WG2228368
1,3-Dichlorobenzene	U		0.550	5.00	5	02/17/2024 17:01	WG2228368
1,4-Dichlorobenzene	U		0.600	5.00	5	02/17/2024 17:01	WG2228368
Dichlorodifluoromethane	U		1.87	25.0	5	02/17/2024 17:01	WG2228368
1,1-Dichloroethane	U		0.500	5.00	5	02/17/2024 17:01	WG2228368
1,2-Dichloroethane	U		0.409	5.00	5	02/17/2024 17:01	WG2228368
1,1-Dichloroethene	17.2		0.940	5.00	5	02/17/2024 17:01	WG2228368
cis-1,2-Dichloroethene	155		0.630	5.00	5	02/17/2024 17:01	WG2228368
trans-1,2-Dichloroethene	5.90		0.745	5.00	5	02/17/2024 17:01	WG2228368
1,2-Dichloropropane	U		0.745	5.00	5	02/17/2024 17:01	WG2228368
1,1-Dichloropropene	U		0.710	5.00	5	02/17/2024 17:01	WG2228368
1,3-Dichloropropane	U		0.550	5.00	5	02/17/2024 17:01	WG2228368
cis-1,3-Dichloropropene	U		0.555	5.00	5	02/17/2024 17:01	WG2228368
trans-1,3-Dichloropropene	U		0.590	5.00	5	02/17/2024 17:01	WG2228368
2,2-Dichloropropane	U		0.805	5.00	5	02/17/2024 17:01	WG2228368
Dicyclopentadiene	U		1.27	5.00	5	02/17/2024 17:01	WG2228368
Di-isopropyl ether	U		0.525	5.00	5	02/17/2024 17:01	WG2228368
Ethylbenzene	U		0.685	5.00	5	02/17/2024 17:01	WG2228368
4-Ethyltoluene	U		1.04	5.00	5	02/17/2024 17:01	WG2228368
Hexachloro-1,3-butadiene	U		1.69	5.00	5	02/17/2024 17:01	WG2228368
n-Hexane	U		3.74	50.0	5	02/17/2024 17:01	WG2228368
Isopropylbenzene	1.83	E4	0.525	5.00	5	02/17/2024 17:01	WG2228368
p-Isopropyltoluene	U		0.600	5.00	5	02/17/2024 17:01	WG2228368
2-Butanone (MEK)	U		5.95	50.0	5	02/17/2024 17:01	WG2228368
Methyl Cyclohexane	U		3.30	5.00	5	02/17/2024 17:01	WG2228368

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Methylene Chloride	5.44	E4	2.15	25.0	5	02/20/2024 04:06	WG2229520
4-Methyl-2-pentanone (MIBK)	3.61	E4	2.39	50.0	5	02/17/2024 17:01	WG2228368
Methyl tert-butyl ether	U		0.505	5.00	5	02/17/2024 17:01	WG2228368
Naphthalene	U		5.00	25.0	5	02/17/2024 17:01	WG2228368
Propene	U		4.68	12.5	5	02/17/2024 17:01	WG2228368
n-Propylbenzene	U		0.497	5.00	5	02/17/2024 17:01	WG2228368
Styrene	U		0.590	5.00	5	02/17/2024 17:01	WG2228368
1,1,1,2-Tetrachloroethane	U		0.735	5.00	5	02/17/2024 17:01	WG2228368
1,1,2,2-Tetrachloroethane	U		0.665	5.00	5	02/17/2024 17:01	WG2228368
1,1,2-Trichlorotrifluoroethane	U		0.900	5.00	5	02/17/2024 17:01	WG2228368
Tetrachloroethene	U		1.50	5.00	5	02/17/2024 17:01	WG2228368
Toluene	U		1.39	5.00	5	02/17/2024 17:01	WG2228368
1,2,3-Trichlorobenzene	U		1.15	5.00	5	02/17/2024 17:01	WG2228368
1,2,4-Trichlorobenzene	U		2.41	5.00	5	02/17/2024 17:01	WG2228368
1,1,1-Trichloroethane	U		0.745	5.00	5	02/17/2024 17:01	WG2228368
1,1,2-Trichloroethane	U		0.790	5.00	5	02/17/2024 17:01	WG2228368
Trichloroethene	195		0.950	5.00	5	02/17/2024 17:01	WG2228368
Trichlorofluoromethane	U		0.800	25.0	5	02/17/2024 17:01	WG2228368
1,2,3-Trichloropropane	U		1.19	12.5	5	02/17/2024 17:01	WG2228368
1,2,4-Trimethylbenzene	U		1.61	5.00	5	02/17/2024 17:01	WG2228368
1,2,3-Trimethylbenzene	U		0.520	5.00	5	02/17/2024 17:01	WG2228368
1,3,5-Trimethylbenzene	U		0.520	5.00	5	02/17/2024 17:01	WG2228368
Vinyl chloride	3.99	E4	1.17	5.00	5	02/17/2024 17:01	WG2228368
Xylenes, Total	U		0.870	15.0	5	02/17/2024 17:01	WG2228368
(S) Toluene-d8	*100*			*80.0-120*		*02/17/2024 17:01*	WG2228368
(S) Toluene-d8	*107*			*80.0-120*		*02/20/2024 04:06*	WG2229520
(S) 4-Bromofluorobenzene	*100*			*77.0-126*		*02/17/2024 17:01*	WG2228368
(S) 4-Bromofluorobenzene	*100*			*77.0-126*		*02/20/2024 04:06*	WG2229520
(S) 1,2-Dichloroethane-d4	*101*			*70.0-130*		*02/17/2024 17:01*	WG2228368
(S) 1,2-Dichloroethane-d4	*92.3*			*70.0-130*		*02/20/2024 04:06*	WG2229520

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
1,4-Dioxane	192		2.99	15.0	5	02/20/2024 21:53	WG2230157
(S) Toluene-d8	*104*			*77.0-127*		*02/20/2024 21:53*	WG2230157

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	85900		300	4000	1000	02/25/2024 02:44	WG2228552

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		113	500	10	02/17/2024 17:21	WG2228368
Acrolein	U		25.4	500	10	02/17/2024 17:21	WG2228368
Acrylonitrile	U		6.71	100	10	02/17/2024 17:21	WG2228368
Benzene	4.48	E4	0.941	10.0	10	02/17/2024 17:21	WG2228368
Bromobenzene	U		1.18	10.0	10	02/17/2024 17:21	WG2228368
Bromodichloromethane	U		1.36	10.0	10	02/17/2024 17:21	WG2228368
Bromoform	U		1.29	10.0	10	02/17/2024 17:21	WG2228368
Bromomethane	U		6.05	50.0	10	02/17/2024 17:21	WG2228368
1,3-Butadiene	U		2.99	20.0	10	02/17/2024 17:21	WG2228368
n-Butylbenzene	U		1.57	10.0	10	02/17/2024 17:21	WG2228368
sec-Butylbenzene	U		1.25	10.0	10	02/17/2024 17:21	WG2228368
tert-Butylbenzene	U		1.27	10.0	10	02/17/2024 17:21	WG2228368
Carbon tetrachloride	U		1.28	10.0	10	02/17/2024 17:21	WG2228368
Carbon disulfide	U		0.962	10.0	10	02/17/2024 17:21	WG2228368
Chlorobenzene	U		1.16	10.0	10	02/17/2024 17:21	WG2228368
Chlorodibromomethane	U		1.40	10.0	10	02/17/2024 17:21	WG2228368
Chloroethane	U		1.92	50.0	10	02/17/2024 17:21	WG2228368
Chloroform	3.86	E4	1.11	50.0	10	02/17/2024 17:21	WG2228368
Chloromethane	U		9.60	25.0	10	02/17/2024 17:21	WG2228368
Cyclohexane	U		1.88	10.0	10	02/17/2024 17:21	WG2228368
2-Chlorotoluene	U		1.06	10.0	10	02/17/2024 17:21	WG2228368
4-Chlorotoluene	U		1.14	10.0	10	02/17/2024 17:21	WG2228368
1,2-Dibromo-3-Chloropropane	U		2.76	50.0	10	02/17/2024 17:21	WG2228368
1,2-Dibromoethane	U		1.26	10.0	10	02/17/2024 17:21	WG2228368
Dibromomethane	U		1.22	10.0	10	02/17/2024 17:21	WG2228368
1,2-Dichlorobenzene	U		1.07	10.0	10	02/17/2024 17:21	WG2228368
1,3-Dichlorobenzene	U		1.10	10.0	10	02/17/2024 17:21	WG2228368
1,4-Dichlorobenzene	U		1.20	10.0	10	02/17/2024 17:21	WG2228368
Dichlorodifluoromethane	U		3.74	50.0	10	02/17/2024 17:21	WG2228368
1,1-Dichloroethane	3.36	E4	1.00	10.0	10	02/17/2024 17:21	WG2228368
1,2-Dichloroethane	U		0.819	10.0	10	02/17/2024 17:21	WG2228368
1,1-Dichloroethene	154		1.88	10.0	10	02/17/2024 17:21	WG2228368
cis-1,2-Dichloroethene	18.9		1.26	10.0	10	02/17/2024 17:21	WG2228368
trans-1,2-Dichloroethene	U		1.49	10.0	10	02/17/2024 17:21	WG2228368
1,2-Dichloropropane	U		1.49	10.0	10	02/17/2024 17:21	WG2228368
1,1-Dichloropropene	U		1.42	10.0	10	02/17/2024 17:21	WG2228368
1,3-Dichloropropane	U		1.10	10.0	10	02/17/2024 17:21	WG2228368
cis-1,3-Dichloropropene	U		1.11	10.0	10	02/17/2024 17:21	WG2228368
trans-1,3-Dichloropropene	U		1.18	10.0	10	02/17/2024 17:21	WG2228368
2,2-Dichloropropane	U		1.61	10.0	10	02/17/2024 17:21	WG2228368
Dicyclopentadiene	U		2.53	10.0	10	02/17/2024 17:21	WG2228368
Di-isopropyl ether	U		1.05	10.0	10	02/17/2024 17:21	WG2228368
Ethylbenzene	U		1.37	10.0	10	02/17/2024 17:21	WG2228368
4-Ethyltoluene	U		2.08	10.0	10	02/17/2024 17:21	WG2228368
Hexachloro-1,3-butadiene	U		3.37	10.0	10	02/17/2024 17:21	WG2228368
n-Hexane	U		7.49	100	10	02/17/2024 17:21	WG2228368
Isopropylbenzene	U		1.05	10.0	10	02/17/2024 17:21	WG2228368
p-Isopropyltoluene	U		1.20	10.0	10	02/17/2024 17:21	WG2228368
2-Butanone (MEK)	U		11.9	100	10	02/17/2024 17:21	WG2228368
Methyl Cyclohexane	U		6.60	10.0	10	02/17/2024 17:21	WG2228368

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Methylene Chloride	8.89	E4	4.30	50.0	10	02/20/2024 04:29	WG2229520
4-Methyl-2-pentanone (MIBK)	U		4.78	100	10	02/17/2024 17:21	WG2228368
Methyl tert-butyl ether	U		1.01	10.0	10	02/17/2024 17:21	WG2228368
Naphthalene	U		10.0	50.0	10	02/17/2024 17:21	WG2228368
Propene	U		9.36	25.0	10	02/17/2024 17:21	WG2228368
n-Propylbenzene	U		0.993	10.0	10	02/17/2024 17:21	WG2228368
Styrene	U		1.18	10.0	10	02/17/2024 17:21	WG2228368
1,1,1,2-Tetrachloroethane	U		1.47	10.0	10	02/17/2024 17:21	WG2228368
1,1,2,2-Tetrachloroethane	U		1.33	10.0	10	02/17/2024 17:21	WG2228368
1,1,2-Trichlorotrifluoroethane	U		1.80	10.0	10	02/17/2024 17:21	WG2228368
Tetrachloroethene	U		3.00	10.0	10	02/17/2024 17:21	WG2228368
Toluene	U		2.78	10.0	10	02/17/2024 17:21	WG2228368
1,2,3-Trichlorobenzene	U		2.30	10.0	10	02/17/2024 17:21	WG2228368
1,2,4-Trichlorobenzene	U		4.81	10.0	10	02/17/2024 17:21	WG2228368
1,1,1-Trichloroethane	U		1.49	10.0	10	02/17/2024 17:21	WG2228368
1,1,2-Trichloroethane	3.28	E4	1.58	10.0	10	02/17/2024 17:21	WG2228368
Trichloroethene	1860		1.90	10.0	10	02/17/2024 17:21	WG2228368
Trichlorofluoromethane	U		1.60	50.0	10	02/17/2024 17:21	WG2228368
1,2,3-Trichloropropane	U		2.37	25.0	10	02/17/2024 17:21	WG2228368
1,2,4-Trimethylbenzene	U		3.22	10.0	10	02/17/2024 17:21	WG2228368
1,2,3-Trimethylbenzene	U		1.04	10.0	10	02/17/2024 17:21	WG2228368
1,3,5-Trimethylbenzene	U		1.04	10.0	10	02/17/2024 17:21	WG2228368
Vinyl chloride	U		2.34	10.0	10	02/17/2024 17:21	WG2228368
Xylenes, Total	U		1.74	30.0	10	02/17/2024 17:21	WG2228368
(S) Toluene-d8	*106*			*80.0-120*		*02/17/2024 17:21*	WG2228368
(S) Toluene-d8	*108*			*80.0-120*		*02/20/2024 04:29*	WG2229520
(S) 4-Bromofluorobenzene	*92.5*			*77.0-126*		*02/17/2024 17:21*	WG2228368
(S) 4-Bromofluorobenzene	*101*			*77.0-126*		*02/20/2024 04:29*	WG2229520
(S) 1,2-Dichloroethane-d4	*98.6*			*70.0-130*		*02/17/2024 17:21*	WG2228368
(S) 1,2-Dichloroethane-d4	*87.2*			*70.0-130*		*02/20/2024 04:29*	WG2229520

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
1,4-Dioxane	378		0.597	3.00	1	02/20/2024 18:58	WG2230157
(S) Toluene-d8	*99.6*			*77.0-127*		*02/20/2024 18:58*	WG2230157

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	U		0.300	4.00	1	02/21/2024 11:40	WG2228552

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		11.3	50.0	1	02/17/2024 14:57	WG2228368
Acrolein	U		2.54	50.0	1	02/17/2024 14:57	WG2228368
Acrylonitrile	U		0.671	10.0	1	02/17/2024 14:57	WG2228368
Benzene	U		0.0941	1.00	1	02/17/2024 14:57	WG2228368
Bromobenzene	U		0.118	1.00	1	02/17/2024 14:57	WG2228368
Bromodichloromethane	U		0.136	1.00	1	02/17/2024 14:57	WG2228368
Bromoform	U		0.129	1.00	1	02/17/2024 14:57	WG2228368
Bromomethane	U		0.605	5.00	1	02/17/2024 14:57	WG2228368
1,3-Butadiene	U		0.299	2.00	1	02/17/2024 14:57	WG2228368
n-Butylbenzene	U		0.157	1.00	1	02/17/2024 14:57	WG2228368
sec-Butylbenzene	U		0.125	1.00	1	02/17/2024 14:57	WG2228368
tert-Butylbenzene	U		0.127	1.00	1	02/17/2024 14:57	WG2228368
Carbon tetrachloride	U		0.128	1.00	1	02/17/2024 14:57	WG2228368
Carbon disulfide	U		0.0962	1.00	1	02/17/2024 14:57	WG2228368
Chlorobenzene	U		0.116	1.00	1	02/17/2024 14:57	WG2228368
Chlorodibromomethane	U		0.140	1.00	1	02/17/2024 14:57	WG2228368
Chloroethane	U		0.192	5.00	1	02/17/2024 14:57	WG2228368
Chloroform	U		0.111	5.00	1	02/17/2024 14:57	WG2228368
Chloromethane	U		0.960	2.50	1	02/17/2024 14:57	WG2228368
Cyclohexane	U		0.188	1.00	1	02/17/2024 14:57	WG2228368
2-Chlorotoluene	U		0.106	1.00	1	02/17/2024 14:57	WG2228368
4-Chlorotoluene	U		0.114	1.00	1	02/17/2024 14:57	WG2228368
1,2-Dibromo-3-Chloropropane	U		0.276	5.00	1	02/17/2024 14:57	WG2228368
1,2-Dibromoethane	U		0.126	1.00	1	02/17/2024 14:57	WG2228368
Dibromomethane	U		0.122	1.00	1	02/17/2024 14:57	WG2228368
1,2-Dichlorobenzene	U		0.107	1.00	1	02/17/2024 14:57	WG2228368
1,3-Dichlorobenzene	U		0.110	1.00	1	02/17/2024 14:57	WG2228368
1,4-Dichlorobenzene	U		0.120	1.00	1	02/17/2024 14:57	WG2228368
Dichlorodifluoromethane	U		0.374	5.00	1	02/17/2024 14:57	WG2228368
1,1-Dichloroethane	U		0.100	1.00	1	02/17/2024 14:57	WG2228368
1,2-Dichloroethane	U		0.0819	1.00	1	02/17/2024 14:57	WG2228368
1,1-Dichloroethene	U		0.188	1.00	1	02/20/2024 04:52	WG2229520
cis-1,2-Dichloroethene	0.234	E4	0.126	1.00	1	02/17/2024 14:57	WG2228368
trans-1,2-Dichloroethene	U		0.149	1.00	1	02/17/2024 14:57	WG2228368
1,2-Dichloropropane	U		0.149	1.00	1	02/17/2024 14:57	WG2228368
1,1-Dichloropropene	U		0.142	1.00	1	02/17/2024 14:57	WG2228368
1,3-Dichloropropane	U		0.110	1.00	1	02/17/2024 14:57	WG2228368
cis-1,3-Dichloropropene	U		0.111	1.00	1	02/17/2024 14:57	WG2228368
trans-1,3-Dichloropropene	U		0.118	1.00	1	02/17/2024 14:57	WG2228368
2,2-Dichloropropane	U		0.161	1.00	1	02/17/2024 14:57	WG2228368
Dicyclopentadiene	U		0.253	1.00	1	02/17/2024 14:57	WG2228368
Di-isopropyl ether	U		0.105	1.00	1	02/17/2024 14:57	WG2228368
Ethylbenzene	U		0.137	1.00	1	02/17/2024 14:57	WG2228368
4-Ethyltoluene	U		0.208	1.00	1	02/17/2024 14:57	WG2228368
Hexachloro-1,3-butadiene	U		0.337	1.00	1	02/17/2024 14:57	WG2228368
n-Hexane	U		0.749	10.0	1	02/17/2024 14:57	WG2228368
Isopropylbenzene	U		0.105	1.00	1	02/17/2024 14:57	WG2228368
p-Isopropyltoluene	U		0.120	1.00	1	02/17/2024 14:57	WG2228368
2-Butanone (MEK)	U		1.19	10.0	1	02/17/2024 14:57	WG2228368
Methyl Cyclohexane	U		0.660	1.00	1	02/17/2024 14:57	WG2228368

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Methylene Chloride	U		0.430	5.00	1	02/20/2024 04:52	WG2229520
4-Methyl-2-pentanone (MIBK)	U		0.478	10.0	1	02/17/2024 14:57	WG2228368
Methyl tert-butyl ether	U		0.101	1.00	1	02/17/2024 14:57	WG2228368
Naphthalene	U		1.00	5.00	1	02/17/2024 14:57	WG2228368
Propene	U		0.936	2.50	1	02/17/2024 14:57	WG2228368
n-Propylbenzene	U		0.0993	1.00	1	02/17/2024 14:57	WG2228368
Styrene	U		0.118	1.00	1	02/17/2024 14:57	WG2228368
1,1,1,2-Tetrachloroethane	U		0.147	1.00	1	02/17/2024 14:57	WG2228368
1,1,2,2-Tetrachloroethane	U		0.133	1.00	1	02/17/2024 14:57	WG2228368
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00	1	02/17/2024 14:57	WG2228368
Tetrachloroethene	U		0.300	1.00	1	02/17/2024 14:57	WG2228368
Toluene	U		0.278	1.00	1	02/17/2024 14:57	WG2228368
1,2,3-Trichlorobenzene	U		0.230	1.00	1	02/17/2024 14:57	WG2228368
1,2,4-Trichlorobenzene	U		0.481	1.00	1	02/17/2024 14:57	WG2228368
1,1,1-Trichloroethane	U		0.149	1.00	1	02/17/2024 14:57	WG2228368
1,1,2-Trichloroethane	U		0.158	1.00	1	02/17/2024 14:57	WG2228368
Trichloroethene	3.57		0.190	1.00	1	02/20/2024 04:52	WG2229520
Trichlorofluoromethane	U		0.160	5.00	1	02/17/2024 14:57	WG2228368
1,2,3-Trichloropropane	U		0.237	2.50	1	02/17/2024 14:57	WG2228368
1,2,4-Trimethylbenzene	U		0.322	1.00	1	02/17/2024 14:57	WG2228368
1,2,3-Trimethylbenzene	U		0.104	1.00	1	02/17/2024 14:57	WG2228368
1,3,5-Trimethylbenzene	U		0.104	1.00	1	02/17/2024 14:57	WG2228368
Vinyl chloride	U		0.234	1.00	1	02/17/2024 14:57	WG2228368
Xylenes, Total	U		0.174	3.00	1	02/17/2024 14:57	WG2228368
(S) Toluene-d8	*101*			*80.0-120*		*02/17/2024 14:57*	WG2228368
(S) Toluene-d8	*103*			*80.0-120*		*02/20/2024 04:52*	WG2229520
(S) 4-Bromofluorobenzene	*91.4*			*77.0-126*		*02/17/2024 14:57*	WG2228368
(S) 4-Bromofluorobenzene	*99.4*			*77.0-126*		*02/20/2024 04:52*	WG2229520
(S) 1,2-Dichloroethane-d4	*101*			*70.0-130*		*02/17/2024 14:57*	WG2228368
(S) 1,2-Dichloroethane-d4	*90.9*			*70.0-130*		*02/20/2024 04:52*	WG2229520

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
1,4-Dioxane	9.87		0.597	3.00	1	02/20/2024 19:20	WG2230157
(S) Toluene-d8	*100*			*77.0-127*		*02/20/2024 19:20*	WG2230157

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	90000		300	4000	1000	02/25/2024 03:12	WG2228552

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		113	500	10	02/17/2024 17:42	WG2228368
Acrolein	U		25.4	500	10	02/17/2024 17:42	WG2228368
Acrylonitrile	U		6.71	100	10	02/17/2024 17:42	WG2228368
Benzene	U		0.941	10.0	10	02/17/2024 17:42	WG2228368
Bromobenzene	U		1.18	10.0	10	02/17/2024 17:42	WG2228368
Bromodichloromethane	U		1.36	10.0	10	02/17/2024 17:42	WG2228368
Bromoform	U		1.29	10.0	10	02/17/2024 17:42	WG2228368
Bromomethane	U		6.05	50.0	10	02/17/2024 17:42	WG2228368
1,3-Butadiene	U		2.99	20.0	10	02/17/2024 17:42	WG2228368
n-Butylbenzene	U		1.57	10.0	10	02/17/2024 17:42	WG2228368
sec-Butylbenzene	U		1.25	10.0	10	02/17/2024 17:42	WG2228368
tert-Butylbenzene	U		1.27	10.0	10	02/17/2024 17:42	WG2228368
Carbon tetrachloride	U		1.28	10.0	10	02/17/2024 17:42	WG2228368
Carbon disulfide	U		0.962	10.0	10	02/17/2024 17:42	WG2228368
Chlorobenzene	U		1.16	10.0	10	02/17/2024 17:42	WG2228368
Chlorodibromomethane	U		1.40	10.0	10	02/17/2024 17:42	WG2228368
Chloroethane	U		1.92	50.0	10	02/17/2024 17:42	WG2228368
Chloroform	1.33	E4	1.11	50.0	10	02/17/2024 17:42	WG2228368
Chloromethane	U		9.60	25.0	10	02/17/2024 17:42	WG2228368
Cyclohexane	U		1.88	10.0	10	02/17/2024 17:42	WG2228368
2-Chlorotoluene	U		1.06	10.0	10	02/17/2024 17:42	WG2228368
4-Chlorotoluene	U		1.14	10.0	10	02/17/2024 17:42	WG2228368
1,2-Dibromo-3-Chloropropane	U		2.76	50.0	10	02/17/2024 17:42	WG2228368
1,2-Dibromoethane	U		1.26	10.0	10	02/17/2024 17:42	WG2228368
Dibromomethane	U		1.22	10.0	10	02/17/2024 17:42	WG2228368
1,2-Dichlorobenzene	U		1.07	10.0	10	02/17/2024 17:42	WG2228368
1,3-Dichlorobenzene	U		1.10	10.0	10	02/17/2024 17:42	WG2228368
1,4-Dichlorobenzene	U		1.20	10.0	10	02/17/2024 17:42	WG2228368
Dichlorodifluoromethane	U		3.74	50.0	10	02/17/2024 17:42	WG2228368
1,1-Dichloroethane	U		1.00	10.0	10	02/17/2024 17:42	WG2228368
1,2-Dichloroethane	U		0.819	10.0	10	02/17/2024 17:42	WG2228368
1,1-Dichloroethene	88.9		1.88	10.0	10	02/17/2024 17:42	WG2228368
cis-1,2-Dichloroethene	U		1.26	10.0	10	02/17/2024 17:42	WG2228368
trans-1,2-Dichloroethene	U		1.49	10.0	10	02/17/2024 17:42	WG2228368
1,2-Dichloropropane	U		1.49	10.0	10	02/17/2024 17:42	WG2228368
1,1-Dichloropropene	U		1.42	10.0	10	02/17/2024 17:42	WG2228368
1,3-Dichloropropane	U		1.10	10.0	10	02/17/2024 17:42	WG2228368
cis-1,3-Dichloropropene	U		1.11	10.0	10	02/17/2024 17:42	WG2228368
trans-1,3-Dichloropropene	U		1.18	10.0	10	02/17/2024 17:42	WG2228368
2,2-Dichloropropane	U		1.61	10.0	10	02/17/2024 17:42	WG2228368
Dicyclopentadiene	U		2.53	10.0	10	02/17/2024 17:42	WG2228368
Di-isopropyl ether	U		1.05	10.0	10	02/17/2024 17:42	WG2228368
Ethylbenzene	U		1.37	10.0	10	02/17/2024 17:42	WG2228368
4-Ethyltoluene	U		2.08	10.0	10	02/17/2024 17:42	WG2228368
Hexachloro-1,3-butadiene	U		3.37	10.0	10	02/17/2024 17:42	WG2228368
n-Hexane	U		7.49	100	10	02/17/2024 17:42	WG2228368
Isopropylbenzene	U		1.05	10.0	10	02/17/2024 17:42	WG2228368
p-Isopropyltoluene	U		1.20	10.0	10	02/17/2024 17:42	WG2228368
2-Butanone (MEK)	U		11.9	100	10	02/17/2024 17:42	WG2228368
Methyl Cyclohexane	U		6.60	10.0	10	02/17/2024 17:42	WG2228368

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Methylene Chloride	U		4.30	50.0	10	02/20/2024 05:15	WG2229520
4-Methyl-2-pentanone (MIBK)	U		4.78	100	10	02/17/2024 17:42	WG2228368
Methyl tert-butyl ether	U		1.01	10.0	10	02/17/2024 17:42	WG2228368
Naphthalene	U		10.0	50.0	10	02/17/2024 17:42	WG2228368
Propene	U		9.36	25.0	10	02/17/2024 17:42	WG2228368
n-Propylbenzene	U		0.993	10.0	10	02/17/2024 17:42	WG2228368
Styrene	U		1.18	10.0	10	02/17/2024 17:42	WG2228368
1,1,1,2-Tetrachloroethane	U		1.47	10.0	10	02/17/2024 17:42	WG2228368
1,1,2,2-Tetrachloroethane	U		1.33	10.0	10	02/17/2024 17:42	WG2228368
1,1,2-Trichlorotrifluoroethane	U		1.80	10.0	10	02/17/2024 17:42	WG2228368
Tetrachloroethene	U		3.00	10.0	10	02/17/2024 17:42	WG2228368
Toluene	U		2.78	10.0	10	02/17/2024 17:42	WG2228368
1,2,3-Trichlorobenzene	U		2.30	10.0	10	02/17/2024 17:42	WG2228368
1,2,4-Trichlorobenzene	U		4.81	10.0	10	02/17/2024 17:42	WG2228368
1,1,1-Trichloroethane	U		1.49	10.0	10	02/17/2024 17:42	WG2228368
1,1,2-Trichloroethane	U		1.58	10.0	10	02/17/2024 17:42	WG2228368
Trichloroethene	241		1.90	10.0	10	02/17/2024 17:42	WG2228368
Trichlorofluoromethane	U		1.60	50.0	10	02/17/2024 17:42	WG2228368
1,2,3-Trichloropropane	U		2.37	25.0	10	02/17/2024 17:42	WG2228368
1,2,4-Trimethylbenzene	U		3.22	10.0	10	02/17/2024 17:42	WG2228368
1,2,3-Trimethylbenzene	U		1.04	10.0	10	02/17/2024 17:42	WG2228368
1,3,5-Trimethylbenzene	U		1.04	10.0	10	02/17/2024 17:42	WG2228368
Vinyl chloride	U		2.34	10.0	10	02/17/2024 17:42	WG2228368
Xylenes, Total	U		1.74	30.0	10	02/17/2024 17:42	WG2228368
(S) Toluene-d8	*103*			*80.0-120*		*02/17/2024 17:42*	WG2228368
(S) Toluene-d8	*106*			*80.0-120*		*02/20/2024 05:15*	WG2229520
(S) 4-Bromofluorobenzene	*91.4*			*77.0-126*		*02/17/2024 17:42*	WG2228368
(S) 4-Bromofluorobenzene	*99.5*			*77.0-126*		*02/20/2024 05:15*	WG2229520
(S) 1,2-Dichloroethane-d4	*105*			*70.0-130*		*02/17/2024 17:42*	WG2228368
(S) 1,2-Dichloroethane-d4	*85.6*			*70.0-130*		*02/20/2024 05:15*	WG2229520

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
1,4-Dioxane	326		5.97	30.0	10	02/20/2024 22:15	WG2230157
(S) Toluene-d8	*99.1*			*77.0-127*		*02/20/2024 22:15*	WG2230157

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	93200		300	4000	1000	02/25/2024 03:40	WG2228552

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		226	1000	20	02/17/2024 18:02	WG2228368
Acrolein	U		50.8	1000	20	02/17/2024 18:02	WG2228368
Acrylonitrile	U		13.4	200	20	02/17/2024 18:02	WG2228368
Benzene	U		1.88	20.0	20	02/17/2024 18:02	WG2228368
Bromobenzene	U		2.36	20.0	20	02/17/2024 18:02	WG2228368
Bromodichloromethane	U		2.72	20.0	20	02/17/2024 18:02	WG2228368
Bromoform	U		2.58	20.0	20	02/17/2024 18:02	WG2228368
Bromomethane	U		12.1	100	20	02/17/2024 18:02	WG2228368
1,3-Butadiene	U		5.98	40.0	20	02/17/2024 18:02	WG2228368
n-Butylbenzene	U		3.14	20.0	20	02/17/2024 18:02	WG2228368
sec-Butylbenzene	U		2.50	20.0	20	02/17/2024 18:02	WG2228368
tert-Butylbenzene	U		2.54	20.0	20	02/17/2024 18:02	WG2228368
Carbon tetrachloride	U		2.56	20.0	20	02/17/2024 18:02	WG2228368
Carbon disulfide	U		1.92	20.0	20	02/17/2024 18:02	WG2228368
Chlorobenzene	U		2.32	20.0	20	02/17/2024 18:02	WG2228368
Chlorodibromomethane	U		2.80	20.0	20	02/17/2024 18:02	WG2228368
Chloroethane	U		3.84	100	20	02/17/2024 18:02	WG2228368
Chloroform	U		2.22	100	20	02/17/2024 18:02	WG2228368
Chloromethane	U		19.2	50.0	20	02/17/2024 18:02	WG2228368
Cyclohexane	U		3.76	20.0	20	02/17/2024 18:02	WG2228368
2-Chlorotoluene	U		2.12	20.0	20	02/17/2024 18:02	WG2228368
4-Chlorotoluene	U		2.28	20.0	20	02/17/2024 18:02	WG2228368
1,2-Dibromo-3-Chloropropane	U		5.52	100	20	02/17/2024 18:02	WG2228368
1,2-Dibromoethane	U		2.52	20.0	20	02/17/2024 18:02	WG2228368
Dibromomethane	U		2.44	20.0	20	02/17/2024 18:02	WG2228368
1,2-Dichlorobenzene	U		2.14	20.0	20	02/17/2024 18:02	WG2228368
1,3-Dichlorobenzene	U		2.20	20.0	20	02/17/2024 18:02	WG2228368
1,4-Dichlorobenzene	U		2.40	20.0	20	02/17/2024 18:02	WG2228368
Dichlorodifluoromethane	U		7.48	100	20	02/17/2024 18:02	WG2228368
1,1-Dichloroethane	U		2.00	20.0	20	02/17/2024 18:02	WG2228368
1,2-Dichloroethane	U		1.64	20.0	20	02/17/2024 18:02	WG2228368
1,1-Dichloroethene	83.3		3.76	20.0	20	02/17/2024 18:02	WG2228368
cis-1,2-Dichloroethene	U		2.52	20.0	20	02/17/2024 18:02	WG2228368
trans-1,2-Dichloroethene	U		2.98	20.0	20	02/17/2024 18:02	WG2228368
1,2-Dichloropropane	U		2.98	20.0	20	02/17/2024 18:02	WG2228368
1,1-Dichloropropene	U		2.84	20.0	20	02/17/2024 18:02	WG2228368
1,3-Dichloropropane	U		2.20	20.0	20	02/17/2024 18:02	WG2228368
cis-1,3-Dichloropropene	U		2.22	20.0	20	02/17/2024 18:02	WG2228368
trans-1,3-Dichloropropene	U		2.36	20.0	20	02/17/2024 18:02	WG2228368
2,2-Dichloropropane	U		3.22	20.0	20	02/17/2024 18:02	WG2228368
Dicyclopentadiene	U		5.06	20.0	20	02/17/2024 18:02	WG2228368
Di-isopropyl ether	U		2.10	20.0	20	02/17/2024 18:02	WG2228368
Ethylbenzene	U		2.74	20.0	20	02/17/2024 18:02	WG2228368
4-Ethyltoluene	U		4.16	20.0	20	02/17/2024 18:02	WG2228368
Hexachloro-1,3-butadiene	U		6.74	20.0	20	02/17/2024 18:02	WG2228368
n-Hexane	U		15.0	200	20	02/17/2024 18:02	WG2228368
Isopropylbenzene	U		2.10	20.0	20	02/17/2024 18:02	WG2228368
p-Isopropyltoluene	U		2.40	20.0	20	02/17/2024 18:02	WG2228368
2-Butanone (MEK)	U		23.8	200	20	02/17/2024 18:02	WG2228368
Methyl Cyclohexane	U		13.2	20.0	20	02/17/2024 18:02	WG2228368

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Methylene Chloride	U		8.60	100	20	02/20/2024 05:38	WG2229520
4-Methyl-2-pentanone (MIBK)	U		9.56	200	20	02/17/2024 18:02	WG2228368
Methyl tert-butyl ether	U		2.02	20.0	20	02/17/2024 18:02	WG2228368
Naphthalene	U		20.0	100	20	02/17/2024 18:02	WG2228368
Propene	U		18.7	50.0	20	02/17/2024 18:02	WG2228368
n-Propylbenzene	U		1.99	20.0	20	02/17/2024 18:02	WG2228368
Styrene	U		2.36	20.0	20	02/17/2024 18:02	WG2228368
1,1,1,2-Tetrachloroethane	U		2.94	20.0	20	02/17/2024 18:02	WG2228368
1,1,2,2-Tetrachloroethane	U		2.66	20.0	20	02/17/2024 18:02	WG2228368
1,1,2-Trichlorotrifluoroethane	U		3.60	20.0	20	02/17/2024 18:02	WG2228368
Tetrachloroethene	U		6.00	20.0	20	02/17/2024 18:02	WG2228368
Toluene	U		5.56	20.0	20	02/17/2024 18:02	WG2228368
1,2,3-Trichlorobenzene	U		4.60	20.0	20	02/17/2024 18:02	WG2228368
1,2,4-Trichlorobenzene	U		9.62	20.0	20	02/17/2024 18:02	WG2228368
1,1,1-Trichloroethane	U		2.98	20.0	20	02/17/2024 18:02	WG2228368
1,1,2-Trichloroethane	U		3.16	20.0	20	02/17/2024 18:02	WG2228368
Trichloroethene	415		3.80	20.0	20	02/17/2024 18:02	WG2228368
Trichlorofluoromethane	U		3.20	100	20	02/17/2024 18:02	WG2228368
1,2,3-Trichloropropane	U		4.74	50.0	20	02/17/2024 18:02	WG2228368
1,2,4-Trimethylbenzene	U		6.44	20.0	20	02/17/2024 18:02	WG2228368
1,2,3-Trimethylbenzene	U		2.08	20.0	20	02/17/2024 18:02	WG2228368
1,3,5-Trimethylbenzene	U		2.08	20.0	20	02/17/2024 18:02	WG2228368
Vinyl chloride	U		4.68	20.0	20	02/17/2024 18:02	WG2228368
Xylenes, Total	U		3.48	60.0	20	02/17/2024 18:02	WG2228368
(S) Toluene-d8	*102*			*80.0-120*		*02/17/2024 18:02*	WG2228368
(S) Toluene-d8	*109*			*80.0-120*		*02/20/2024 05:38*	WG2229520
(S) 4-Bromofluorobenzene	*90.1*			*77.0-126*		*02/17/2024 18:02*	WG2228368
(S) 4-Bromofluorobenzene	*99.1*			*77.0-126*		*02/20/2024 05:38*	WG2229520
(S) 1,2-Dichloroethane-d4	*106*			*70.0-130*		*02/17/2024 18:02*	WG2228368
(S) 1,2-Dichloroethane-d4	*88.4*			*70.0-130*		*02/20/2024 05:38*	WG2229520

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
1,4-Dioxane	256		0.597	3.00	1	02/21/2024 15:12	WG2230690
(S) Toluene-d8	*98.4*			*77.0-127*		*02/21/2024 15:12*	WG2230690

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	468000		1500	20000	5000	02/29/2024 17:10	WG2228553

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		2820	12500	250	02/18/2024 04:48	WG2228641
Acrolein	U		635	12500	250	02/18/2024 04:48	WG2228641
Acrylonitrile	U		168	2500	250	02/18/2024 04:48	WG2228641
Benzene	U		23.5	250	250	02/18/2024 04:48	WG2228641
Bromobenzene	U		29.5	250	250	02/18/2024 04:48	WG2228641
Bromodichloromethane	U		34.0	250	250	02/18/2024 04:48	WG2228641
Bromoform	U		32.3	250	250	02/18/2024 04:48	WG2228641
Bromomethane	U		151	1250	250	02/18/2024 04:48	WG2228641
1,3-Butadiene	U		74.8	500	250	02/18/2024 04:48	WG2228641
n-Butylbenzene	U		39.3	250	250	02/18/2024 04:48	WG2228641
sec-Butylbenzene	U		31.3	250	250	02/18/2024 04:48	WG2228641
tert-Butylbenzene	U		31.8	250	250	02/18/2024 04:48	WG2228641
Carbon tetrachloride	U		32.0	250	250	02/18/2024 04:48	WG2228641
Carbon disulfide	U		24.1	250	250	02/18/2024 04:48	WG2228641
Chlorobenzene	U		29.0	250	250	02/18/2024 04:48	WG2228641
Chlorodibromomethane	U		35.0	250	250	02/18/2024 04:48	WG2228641
Chloroethane	U	L1	48.0	1250	250	02/18/2024 04:48	WG2228641
Chloroform	U		27.8	1250	250	02/18/2024 04:48	WG2228641
Chloromethane	U		240	625	250	02/18/2024 04:48	WG2228641
Cyclohexane	U		47.0	250	250	02/18/2024 04:48	WG2228641
2-Chlorotoluene	U		26.5	250	250	02/18/2024 04:48	WG2228641
4-Chlorotoluene	U		28.5	250	250	02/18/2024 04:48	WG2228641
1,2-Dibromo-3-Chloropropane	U		69.0	1250	250	02/18/2024 04:48	WG2228641
1,2-Dibromoethane	U		31.5	250	250	02/18/2024 04:48	WG2228641
Dibromomethane	U		30.5	250	250	02/18/2024 04:48	WG2228641
1,2-Dichlorobenzene	U	L2	26.8	250	250	02/18/2024 04:48	WG2228641
1,3-Dichlorobenzene	U		27.5	250	250	02/18/2024 04:48	WG2228641
1,4-Dichlorobenzene	U		30.0	250	250	02/18/2024 04:48	WG2228641
Dichlorodifluoromethane	U		93.5	1250	250	02/18/2024 04:48	WG2228641
1,1-Dichloroethane	U		25.0	250	250	02/18/2024 04:48	WG2228641
1,2-Dichloroethane	U		20.5	250	250	02/18/2024 04:48	WG2228641
1,1-Dichloroethene	836		47.0	250	250	02/18/2024 04:48	WG2228641
cis-1,2-Dichloroethene	U		31.5	250	250	02/18/2024 04:48	WG2228641
trans-1,2-Dichloroethene	U		37.3	250	250	02/18/2024 04:48	WG2228641
1,2-Dichloropropane	U		37.3	250	250	02/18/2024 04:48	WG2228641
1,1-Dichloropropene	U		35.5	250	250	02/18/2024 04:48	WG2228641
1,3-Dichloropropane	U		27.5	250	250	02/18/2024 04:48	WG2228641
cis-1,3-Dichloropropene	U		27.8	250	250	02/18/2024 04:48	WG2228641
trans-1,3-Dichloropropene	U		29.5	250	250	02/18/2024 04:48	WG2228641
2,2-Dichloropropane	U		40.3	250	250	02/18/2024 04:48	WG2228641
Dicyclopentadiene	U		63.3	250	250	02/18/2024 04:48	WG2228641
Di-isopropyl ether	U		26.3	250	250	02/18/2024 04:48	WG2228641
Ethylbenzene	U		34.3	250	250	02/18/2024 04:48	WG2228641
4-Ethyltoluene	U		52.0	250	250	02/18/2024 04:48	WG2228641
Hexachloro-1,3-butadiene	U		84.3	250	250	02/18/2024 04:48	WG2228641
n-Hexane	U		187	2500	250	02/18/2024 04:48	WG2228641
Isopropylbenzene	U		26.3	250	250	02/18/2024 04:48	WG2228641
p-Isopropyltoluene	U		30.0	250	250	02/18/2024 04:48	WG2228641
2-Butanone (MEK)	U		298	2500	250	02/18/2024 04:48	WG2228641
Methyl Cyclohexane	U		165	250	250	02/18/2024 04:48	WG2228641

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

Analyte	Result ug/l	Qualifier	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	Batch
Methylene Chloride	160	E4	108	1250	250	02/18/2024 04:48	WG2228641
4-Methyl-2-pentanone (MIBK)	U		120	2500	250	02/18/2024 04:48	WG2228641
Methyl tert-butyl ether	U		25.3	250	250	02/18/2024 04:48	WG2228641
Naphthalene	U	L2	250	1250	250	02/18/2024 04:48	WG2228641
Propene	U		234	625	250	02/18/2024 04:48	WG2228641
n-Propylbenzene	U		24.8	250	250	02/18/2024 04:48	WG2228641
Styrene	U		29.5	250	250	02/18/2024 04:48	WG2228641
1,1,1,2-Tetrachloroethane	U		36.8	250	250	02/18/2024 04:48	WG2228641
1,1,2,2-Tetrachloroethane	U		33.3	250	250	02/18/2024 04:48	WG2228641
1,1,2-Trichlorotrifluoroethane	U		45.0	250	250	02/18/2024 04:48	WG2228641
Tetrachloroethene	U		75.0	250	250	02/18/2024 04:48	WG2228641
Toluene	U		69.5	250	250	02/18/2024 04:48	WG2228641
1,2,3-Trichlorobenzene	U		57.5	250	250	02/18/2024 04:48	WG2228641
1,2,4-Trichlorobenzene	U		120	250	250	02/18/2024 04:48	WG2228641
1,1,1-Trichloroethane	U		37.3	250	250	02/18/2024 04:48	WG2228641
1,1,2-Trichloroethane	U		39.5	250	250	02/18/2024 04:48	WG2228641
Trichloroethene	5930		47.5	250	250	02/18/2024 04:48	WG2228641
Trichlorofluoromethane	U	L1	40.0	1250	250	02/18/2024 04:48	WG2228641
1,2,3-Trichloropropane	U		59.3	625	250	02/18/2024 04:48	WG2228641
1,2,4-Trimethylbenzene	U		80.5	250	250	02/18/2024 04:48	WG2228641
1,2,3-Trimethylbenzene	U		26.0	250	250	02/18/2024 04:48	WG2228641
1,3,5-Trimethylbenzene	U		26.0	250	250	02/18/2024 04:48	WG2228641
Vinyl chloride	U		58.5	250	250	02/18/2024 04:48	WG2228641
Xylenes, Total	U		43.5	750	250	02/18/2024 04:48	WG2228641
(S) Toluene-d8	*108*			*80.0-120*		*02/18/2024 04:48*	WG2228641
(S) 4-Bromofluorobenzene	*121*			*77.0-126*		*02/18/2024 04:48*	WG2228641
(S) 1,2-Dichloroethane-d4	*87.9*			*70.0-130*		*02/18/2024 04:48*	WG2228641

Sample Narrative:

L1705776-21 WG2228641: Target compounds too high to run at a lower dilution.

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result ug/l	Qualifier	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	Batch
1,4-Dioxane	523		5.97	30.0	10	02/22/2024 13:45	WG2231600
(S) Toluene-d8	*96.8*			*77.0-127*		*02/22/2024 13:45*	WG2231600

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	85600		1500	20000	5000	02/29/2024 17:38	WG2228553

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		282	1250	25	02/18/2024 05:11	WG2228641
Acrolein	U		63.5	1250	25	02/18/2024 05:11	WG2228641
Acrylonitrile	U		16.8	250	25	02/18/2024 05:11	WG2228641
Benzene	U		2.35	25.0	25	02/18/2024 05:11	WG2228641
Bromobenzene	U		2.95	25.0	25	02/18/2024 05:11	WG2228641
Bromodichloromethane	U		3.40	25.0	25	02/18/2024 05:11	WG2228641
Bromoform	U		3.22	25.0	25	02/18/2024 05:11	WG2228641
Bromomethane	U		15.1	125	25	02/18/2024 05:11	WG2228641
1,3-Butadiene	U		7.48	50.0	25	02/18/2024 05:11	WG2228641
n-Butylbenzene	U		3.93	25.0	25	02/18/2024 05:11	WG2228641
sec-Butylbenzene	U		3.13	25.0	25	02/18/2024 05:11	WG2228641
tert-Butylbenzene	U		3.18	25.0	25	02/18/2024 05:11	WG2228641
Carbon tetrachloride	U		3.20	25.0	25	02/18/2024 05:11	WG2228641
Carbon disulfide	U		2.41	25.0	25	02/18/2024 05:11	WG2228641
Chlorobenzene	U		2.90	25.0	25	02/18/2024 05:11	WG2228641
Chlorodibromomethane	U		3.50	25.0	25	02/18/2024 05:11	WG2228641
Chloroethane	U	L1	4.80	125	25	02/18/2024 05:11	WG2228641
Chloroform	U		2.78	125	25	02/18/2024 05:11	WG2228641
Chloromethane	U		24.0	62.5	25	02/18/2024 05:11	WG2228641
Cyclohexane	U		4.70	25.0	25	02/18/2024 05:11	WG2228641
2-Chlorotoluene	U		2.65	25.0	25	02/18/2024 05:11	WG2228641
4-Chlorotoluene	U		2.85	25.0	25	02/18/2024 05:11	WG2228641
1,2-Dibromo-3-Chloropropane	U		6.90	125	25	02/18/2024 05:11	WG2228641
1,2-Dibromoethane	U		3.15	25.0	25	02/18/2024 05:11	WG2228641
Dibromomethane	U		3.05	25.0	25	02/18/2024 05:11	WG2228641
1,2-Dichlorobenzene	U	L2	2.68	25.0	25	02/18/2024 05:11	WG2228641
1,3-Dichlorobenzene	U		2.75	25.0	25	02/18/2024 05:11	WG2228641
1,4-Dichlorobenzene	U		3.00	25.0	25	02/18/2024 05:11	WG2228641
Dichlorodifluoromethane	U		9.35	125	25	02/18/2024 05:11	WG2228641
1,1-Dichloroethane	U		2.50	25.0	25	02/18/2024 05:11	WG2228641
1,2-Dichloroethane	U		2.05	25.0	25	02/18/2024 05:11	WG2228641
1,1-Dichloroethene	91.1		4.70	25.0	25	02/18/2024 05:11	WG2228641
cis-1,2-Dichloroethene	U		3.15	25.0	25	02/18/2024 05:11	WG2228641
trans-1,2-Dichloroethene	U		3.73	25.0	25	02/18/2024 05:11	WG2228641
1,2-Dichloropropane	U		3.73	25.0	25	02/18/2024 05:11	WG2228641
1,1-Dichloropropene	U		3.55	25.0	25	02/18/2024 05:11	WG2228641
1,3-Dichloropropane	U		2.75	25.0	25	02/18/2024 05:11	WG2228641
cis-1,3-Dichloropropene	U		2.78	25.0	25	02/18/2024 05:11	WG2228641
trans-1,3-Dichloropropene	U		2.95	25.0	25	02/18/2024 05:11	WG2228641
2,2-Dichloropropane	U		4.03	25.0	25	02/18/2024 05:11	WG2228641
Dicyclopentadiene	U		6.33	25.0	25	02/18/2024 05:11	WG2228641
Di-isopropyl ether	U		2.63	25.0	25	02/18/2024 05:11	WG2228641
Ethylbenzene	U		3.43	25.0	25	02/18/2024 05:11	WG2228641
4-Ethyltoluene	U		5.20	25.0	25	02/18/2024 05:11	WG2228641
Hexachloro-1,3-butadiene	U		8.43	25.0	25	02/18/2024 05:11	WG2228641
n-Hexane	U		18.7	250	25	02/18/2024 05:11	WG2228641
Isopropylbenzene	U		2.63	25.0	25	02/18/2024 05:11	WG2228641
p-Isopropyltoluene	U		3.00	25.0	25	02/18/2024 05:11	WG2228641
2-Butanone (MEK)	U		29.8	250	25	02/18/2024 05:11	WG2228641
Methyl Cyclohexane	U		16.5	25.0	25	02/18/2024 05:11	WG2228641

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Methylene Chloride	16.1	E4	10.7	125	25	02/18/2024 05:11	WG2228641
4-Methyl-2-pentanone (MIBK)	U		12.0	250	25	02/18/2024 05:11	WG2228641
Methyl tert-butyl ether	U		2.53	25.0	25	02/18/2024 05:11	WG2228641
Naphthalene	U	L2	25.0	125	25	02/18/2024 05:11	WG2228641
Propene	U		23.4	62.5	25	02/18/2024 05:11	WG2228641
n-Propylbenzene	U		2.48	25.0	25	02/18/2024 05:11	WG2228641
Styrene	U		2.95	25.0	25	02/18/2024 05:11	WG2228641
1,1,1,2-Tetrachloroethane	U		3.68	25.0	25	02/18/2024 05:11	WG2228641
1,1,2,2-Tetrachloroethane	U		3.33	25.0	25	02/18/2024 05:11	WG2228641
1,1,2-Trichlorotrifluoroethane	U		4.50	25.0	25	02/18/2024 05:11	WG2228641
Tetrachloroethene	U		7.50	25.0	25	02/18/2024 05:11	WG2228641
Toluene	U		6.95	25.0	25	02/18/2024 05:11	WG2228641
1,2,3-Trichlorobenzene	U		5.75	25.0	25	02/18/2024 05:11	WG2228641
1,2,4-Trichlorobenzene	U		12.0	25.0	25	02/18/2024 05:11	WG2228641
1,1,1-Trichloroethane	U		3.73	25.0	25	02/18/2024 05:11	WG2228641
1,1,2-Trichloroethane	U		3.95	25.0	25	02/18/2024 05:11	WG2228641
Trichloroethene	588		4.75	25.0	25	02/18/2024 05:11	WG2228641
Trichlorofluoromethane	U	L1	4.00	125	25	02/18/2024 05:11	WG2228641
1,2,3-Trichloropropane	U		5.93	62.5	25	02/18/2024 05:11	WG2228641
1,2,4-Trimethylbenzene	U		8.05	25.0	25	02/18/2024 05:11	WG2228641
1,2,3-Trimethylbenzene	U		2.60	25.0	25	02/18/2024 05:11	WG2228641
1,3,5-Trimethylbenzene	U		2.60	25.0	25	02/18/2024 05:11	WG2228641
Vinyl chloride	U		5.85	25.0	25	02/18/2024 05:11	WG2228641
Xylenes, Total	U		4.35	75.0	25	02/18/2024 05:11	WG2228641
(S) Toluene-d8	*108*			*80.0-120*		*02/18/2024 05:11*	WG2228641
(S) 4-Bromofluorobenzene	*128*	S10		*77.0-126*		*02/18/2024 05:11*	WG2228641
(S) 1,2-Dichloroethane-d4	*86.3*			*70.0-130*		*02/18/2024 05:11*	WG2228641

Sample Narrative:

L1705776-22 WG2228641: Non-target compounds too high to run at a lower dilution.

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
1,4-Dioxane	15.9		0.597	3.00	1	02/20/2024 19:42	WG2230157
(S) Toluene-d8	*99.9*			*77.0-127*		*02/20/2024 19:42*	WG2230157

Wet Chemistry by Method 314.0 Mod

Analyte	Result ug/l	Qualifier	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	Batch
Perchlorate	U	M2	0.300	4.00	1	02/29/2024 18:06	WG2228553

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

Analyte	Result ug/l	Qualifier	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	Batch
Acetone	U		11.3	50.0	1	02/18/2024 02:05	WG2228641
Acrolein	U		2.54	50.0	1	02/18/2024 02:05	WG2228641
Acrylonitrile	U		0.671	10.0	1	02/18/2024 02:05	WG2228641
Benzene	U		0.0941	1.00	1	02/18/2024 02:05	WG2228641
Bromobenzene	U		0.118	1.00	1	02/18/2024 02:05	WG2228641
Bromodichloromethane	U		0.136	1.00	1	02/18/2024 02:05	WG2228641
Bromoform	U		0.129	1.00	1	02/18/2024 02:05	WG2228641
Bromomethane	U		0.605	5.00	1	02/18/2024 02:05	WG2228641
1,3-Butadiene	U		0.299	2.00	1	02/18/2024 02:05	WG2228641
n-Butylbenzene	U		0.157	1.00	1	02/18/2024 02:05	WG2228641
sec-Butylbenzene	U		0.125	1.00	1	02/18/2024 02:05	WG2228641
tert-Butylbenzene	U		0.127	1.00	1	02/18/2024 02:05	WG2228641
Carbon tetrachloride	U		0.128	1.00	1	02/18/2024 02:05	WG2228641
Carbon disulfide	U		0.0962	1.00	1	02/18/2024 02:05	WG2228641
Chlorobenzene	U		0.116	1.00	1	02/18/2024 02:05	WG2228641
Chlorodibromomethane	U		0.140	1.00	1	02/18/2024 02:05	WG2228641
Chloroethane	U	L1	0.192	5.00	1	02/18/2024 02:05	WG2228641
Chloroform	U		0.111	5.00	1	02/18/2024 02:05	WG2228641
Chloromethane	U		0.960	2.50	1	02/18/2024 02:05	WG2228641
Cyclohexane	U		0.188	1.00	1	02/18/2024 02:05	WG2228641
2-Chlorotoluene	U		0.106	1.00	1	02/18/2024 02:05	WG2228641
4-Chlorotoluene	U		0.114	1.00	1	02/18/2024 02:05	WG2228641
1,2-Dibromo-3-Chloropropane	U		0.276	5.00	1	02/18/2024 02:05	WG2228641
1,2-Dibromoethane	U		0.126	1.00	1	02/18/2024 02:05	WG2228641
Dibromomethane	U		0.122	1.00	1	02/18/2024 02:05	WG2228641
1,2-Dichlorobenzene	U	L2	0.107	1.00	1	02/18/2024 02:05	WG2228641
1,3-Dichlorobenzene	U		0.110	1.00	1	02/18/2024 02:05	WG2228641
1,4-Dichlorobenzene	U		0.120	1.00	1	02/18/2024 02:05	WG2228641
Dichlorodifluoromethane	U		0.374	5.00	1	02/18/2024 02:05	WG2228641
1,1-Dichloroethane	U		0.100	1.00	1	02/18/2024 02:05	WG2228641
1,2-Dichloroethane	U		0.0819	1.00	1	02/18/2024 02:05	WG2228641
1,1-Dichloroethene	U		0.188	1.00	1	02/18/2024 02:05	WG2228641
cis-1,2-Dichloroethene	U		0.126	1.00	1	02/18/2024 02:05	WG2228641
trans-1,2-Dichloroethene	U		0.149	1.00	1	02/18/2024 02:05	WG2228641
1,2-Dichloropropane	U		0.149	1.00	1	02/18/2024 02:05	WG2228641
1,1-Dichloropropene	U		0.142	1.00	1	02/18/2024 02:05	WG2228641
1,3-Dichloropropane	U		0.110	1.00	1	02/18/2024 02:05	WG2228641
cis-1,3-Dichloropropene	U		0.111	1.00	1	02/18/2024 02:05	WG2228641
trans-1,3-Dichloropropene	U		0.118	1.00	1	02/18/2024 02:05	WG2228641
2,2-Dichloropropane	U		0.161	1.00	1	02/18/2024 02:05	WG2228641
Dicyclopentadiene	U		0.253	1.00	1	02/18/2024 02:05	WG2228641
Di-isopropyl ether	U		0.105	1.00	1	02/18/2024 02:05	WG2228641
Ethylbenzene	U		0.137	1.00	1	02/18/2024 02:05	WG2228641
4-Ethyltoluene	U		0.208	1.00	1	02/18/2024 02:05	WG2228641
Hexachloro-1,3-butadiene	U		0.337	1.00	1	02/18/2024 02:05	WG2228641
n-Hexane	U		0.749	10.0	1	02/18/2024 02:05	WG2228641
Isopropylbenzene	U		0.105	1.00	1	02/18/2024 02:05	WG2228641
p-Isopropyltoluene	U		0.120	1.00	1	02/18/2024 02:05	WG2228641
2-Butanone (MEK)	U		1.19	10.0	1	02/18/2024 02:05	WG2228641
Methyl Cyclohexane	U		0.660	1.00	1	02/18/2024 02:05	WG2228641

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Methylene Chloride	U		0.430	5.00	1	02/18/2024 02:05	WG2228641
4-Methyl-2-pentanone (MIBK)	U		0.478	10.0	1	02/18/2024 02:05	WG2228641
Methyl tert-butyl ether	U		0.101	1.00	1	02/18/2024 02:05	WG2228641
Naphthalene	U	L2	1.00	5.00	1	02/18/2024 02:05	WG2228641
Propene	U		0.936	2.50	1	02/18/2024 02:05	WG2228641
n-Propylbenzene	U		0.0993	1.00	1	02/18/2024 02:05	WG2228641
Styrene	U		0.118	1.00	1	02/18/2024 02:05	WG2228641
1,1,1,2-Tetrachloroethane	U		0.147	1.00	1	02/18/2024 02:05	WG2228641
1,1,2,2-Tetrachloroethane	U		0.133	1.00	1	02/18/2024 02:05	WG2228641
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00	1	02/18/2024 02:05	WG2228641
Tetrachloroethene	U		0.300	1.00	1	02/18/2024 02:05	WG2228641
Toluene	U		0.278	1.00	1	02/18/2024 02:05	WG2228641
1,2,3-Trichlorobenzene	U		0.230	1.00	1	02/18/2024 02:05	WG2228641
1,2,4-Trichlorobenzene	U		0.481	1.00	1	02/18/2024 02:05	WG2228641
1,1,1-Trichloroethane	U		0.149	1.00	1	02/18/2024 02:05	WG2228641
1,1,2-Trichloroethane	U		0.158	1.00	1	02/18/2024 02:05	WG2228641
Trichloroethene	U		0.190	1.00	1	02/18/2024 02:05	WG2228641
Trichlorofluoromethane	U	L1	0.160	5.00	1	02/18/2024 02:05	WG2228641
1,2,3-Trichloropropane	U		0.237	2.50	1	02/18/2024 02:05	WG2228641
1,2,4-Trimethylbenzene	U		0.322	1.00	1	02/18/2024 02:05	WG2228641
1,2,3-Trimethylbenzene	U		0.104	1.00	1	02/18/2024 02:05	WG2228641
1,3,5-Trimethylbenzene	U		0.104	1.00	1	02/18/2024 02:05	WG2228641
Vinyl chloride	U		0.234	1.00	1	02/18/2024 02:05	WG2228641
Xylenes, Total	U		0.174	3.00	1	02/18/2024 02:05	WG2228641
(S) Toluene-d8	*108*			*80.0-120*		*02/18/2024 02:05*	WG2228641
(S) 4-Bromofluorobenzene	*99.8*			*77.0-126*		*02/18/2024 02:05*	WG2228641
(S) 1,2-Dichloroethane-d4	*86.3*			*70.0-130*		*02/18/2024 02:05*	WG2228641

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
1,4-Dioxane	U		0.597	3.00	1	02/20/2024 20:04	WG2230157
(S) Toluene-d8	*98.8*			*77.0-127*		*02/20/2024 20:04*	WG2230157

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	U		0.300	4.00	1	02/29/2024 18:34	WG2228553

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		11.3	50.0	1	02/18/2024 02:28	WG2228641
Acrolein	U		2.54	50.0	1	02/18/2024 02:28	WG2228641
Acrylonitrile	U		0.671	10.0	1	02/18/2024 02:28	WG2228641
Benzene	U		0.0941	1.00	1	02/18/2024 02:28	WG2228641
Bromobenzene	U		0.118	1.00	1	02/18/2024 02:28	WG2228641
Bromodichloromethane	U		0.136	1.00	1	02/18/2024 02:28	WG2228641
Bromoform	U		0.129	1.00	1	02/18/2024 02:28	WG2228641
Bromomethane	U		0.605	5.00	1	02/18/2024 02:28	WG2228641
1,3-Butadiene	U		0.299	2.00	1	02/18/2024 02:28	WG2228641
n-Butylbenzene	U		0.157	1.00	1	02/18/2024 02:28	WG2228641
sec-Butylbenzene	U		0.125	1.00	1	02/18/2024 02:28	WG2228641
tert-Butylbenzene	U		0.127	1.00	1	02/18/2024 02:28	WG2228641
Carbon tetrachloride	U		0.128	1.00	1	02/18/2024 02:28	WG2228641
Carbon disulfide	U		0.0962	1.00	1	02/18/2024 02:28	WG2228641
Chlorobenzene	U		0.116	1.00	1	02/18/2024 02:28	WG2228641
Chlorodibromomethane	U		0.140	1.00	1	02/18/2024 02:28	WG2228641
Chloroethane	U	L1	0.192	5.00	1	02/18/2024 02:28	WG2228641
Chloroform	U		0.111	5.00	1	02/18/2024 02:28	WG2228641
Chloromethane	U		0.960	2.50	1	02/18/2024 02:28	WG2228641
Cyclohexane	U		0.188	1.00	1	02/18/2024 02:28	WG2228641
2-Chlorotoluene	U		0.106	1.00	1	02/18/2024 02:28	WG2228641
4-Chlorotoluene	U		0.114	1.00	1	02/18/2024 02:28	WG2228641
1,2-Dibromo-3-Chloropropane	U		0.276	5.00	1	02/18/2024 02:28	WG2228641
1,2-Dibromoethane	U		0.126	1.00	1	02/18/2024 02:28	WG2228641
Dibromomethane	U		0.122	1.00	1	02/18/2024 02:28	WG2228641
1,2-Dichlorobenzene	U	L2	0.107	1.00	1	02/18/2024 02:28	WG2228641
1,3-Dichlorobenzene	U		0.110	1.00	1	02/18/2024 02:28	WG2228641
1,4-Dichlorobenzene	U		0.120	1.00	1	02/18/2024 02:28	WG2228641
Dichlorodifluoromethane	U		0.374	5.00	1	02/18/2024 02:28	WG2228641
1,1-Dichloroethane	U		0.100	1.00	1	02/18/2024 02:28	WG2228641
1,2-Dichloroethane	U		0.0819	1.00	1	02/18/2024 02:28	WG2228641
1,1-Dichloroethene	U		0.188	1.00	1	02/18/2024 02:28	WG2228641
cis-1,2-Dichloroethene	U		0.126	1.00	1	02/18/2024 02:28	WG2228641
trans-1,2-Dichloroethene	U		0.149	1.00	1	02/18/2024 02:28	WG2228641
1,2-Dichloropropane	U		0.149	1.00	1	02/18/2024 02:28	WG2228641
1,1-Dichloropropene	U		0.142	1.00	1	02/18/2024 02:28	WG2228641
1,3-Dichloropropane	U		0.110	1.00	1	02/18/2024 02:28	WG2228641
cis-1,3-Dichloropropene	U		0.111	1.00	1	02/18/2024 02:28	WG2228641
trans-1,3-Dichloropropene	U		0.118	1.00	1	02/18/2024 02:28	WG2228641
2,2-Dichloropropane	U		0.161	1.00	1	02/18/2024 02:28	WG2228641
Dicyclopentadiene	U		0.253	1.00	1	02/18/2024 02:28	WG2228641
Di-isopropyl ether	U		0.105	1.00	1	02/18/2024 02:28	WG2228641
Ethylbenzene	U		0.137	1.00	1	02/18/2024 02:28	WG2228641
4-Ethyltoluene	U		0.208	1.00	1	02/18/2024 02:28	WG2228641
Hexachloro-1,3-butadiene	U		0.337	1.00	1	02/18/2024 02:28	WG2228641
n-Hexane	U		0.749	10.0	1	02/18/2024 02:28	WG2228641
Isopropylbenzene	U		0.105	1.00	1	02/18/2024 02:28	WG2228641
p-Isopropyltoluene	U		0.120	1.00	1	02/18/2024 02:28	WG2228641
2-Butanone (MEK)	U		1.19	10.0	1	02/18/2024 02:28	WG2228641
Methyl Cyclohexane	U		0.660	1.00	1	02/18/2024 02:28	WG2228641

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Methylene Chloride	U		0.430	5.00	1	02/18/2024 02:28	WG2228641
4-Methyl-2-pentanone (MIBK)	U		0.478	10.0	1	02/18/2024 02:28	WG2228641
Methyl tert-butyl ether	U		0.101	1.00	1	02/18/2024 02:28	WG2228641
Naphthalene	U	L2	1.00	5.00	1	02/18/2024 02:28	WG2228641
Propene	U		0.936	2.50	1	02/18/2024 02:28	WG2228641
n-Propylbenzene	U		0.0993	1.00	1	02/18/2024 02:28	WG2228641
Styrene	U		0.118	1.00	1	02/18/2024 02:28	WG2228641
1,1,1,2-Tetrachloroethane	U		0.147	1.00	1	02/18/2024 02:28	WG2228641
1,1,2,2-Tetrachloroethane	U		0.133	1.00	1	02/18/2024 02:28	WG2228641
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00	1	02/18/2024 02:28	WG2228641
Tetrachloroethene	U		0.300	1.00	1	02/18/2024 02:28	WG2228641
Toluene	U		0.278	1.00	1	02/18/2024 02:28	WG2228641
1,2,3-Trichlorobenzene	U		0.230	1.00	1	02/18/2024 02:28	WG2228641
1,2,4-Trichlorobenzene	U		0.481	1.00	1	02/18/2024 02:28	WG2228641
1,1,1-Trichloroethane	U		0.149	1.00	1	02/18/2024 02:28	WG2228641
1,1,2-Trichloroethane	U		0.158	1.00	1	02/18/2024 02:28	WG2228641
Trichloroethene	3.08		0.190	1.00	1	02/18/2024 02:28	WG2228641
Trichlorofluoromethane	U	L1	0.160	5.00	1	02/18/2024 02:28	WG2228641
1,2,3-Trichloropropane	U		0.237	2.50	1	02/18/2024 02:28	WG2228641
1,2,4-Trimethylbenzene	U		0.322	1.00	1	02/18/2024 02:28	WG2228641
1,2,3-Trimethylbenzene	U		0.104	1.00	1	02/18/2024 02:28	WG2228641
1,3,5-Trimethylbenzene	U		0.104	1.00	1	02/18/2024 02:28	WG2228641
Vinyl chloride	U		0.234	1.00	1	02/18/2024 02:28	WG2228641
Xylenes, Total	U		0.174	3.00	1	02/18/2024 02:28	WG2228641
(S) Toluene-d8	*109*			*80.0-120*		*02/18/2024 02:28*	WG2228641
(S) 4-Bromofluorobenzene	*99.9*			*77.0-126*		*02/18/2024 02:28*	WG2228641
(S) 1,2-Dichloroethane-d4	*85.6*			*70.0-130*		*02/18/2024 02:28*	WG2228641

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
1,4-Dioxane	U		0.597	3.00	1	02/20/2024 20:26	WG2230157
(S) Toluene-d8	*99.1*			*77.0-127*		*02/20/2024 20:26*	WG2230157

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		11.3	50.0	1	02/17/2024 23:23	WG2228641
Acrolein	U		2.54	50.0	1	02/17/2024 23:23	WG2228641
Acrylonitrile	U		0.671	10.0	1	02/17/2024 23:23	WG2228641
Benzene	U		0.0941	1.00	1	02/17/2024 23:23	WG2228641
Bromobenzene	U		0.118	1.00	1	02/17/2024 23:23	WG2228641
Bromodichloromethane	U		0.136	1.00	1	02/17/2024 23:23	WG2228641
Bromoform	U		0.129	1.00	1	02/17/2024 23:23	WG2228641
Bromomethane	U		0.605	5.00	1	02/17/2024 23:23	WG2228641
1,3-Butadiene	U		0.299	2.00	1	02/17/2024 23:23	WG2228641
n-Butylbenzene	U		0.157	1.00	1	02/17/2024 23:23	WG2228641
sec-Butylbenzene	U		0.125	1.00	1	02/17/2024 23:23	WG2228641
tert-Butylbenzene	U		0.127	1.00	1	02/17/2024 23:23	WG2228641
Carbon tetrachloride	U		0.128	1.00	1	02/17/2024 23:23	WG2228641
Carbon disulfide	U		0.0962	1.00	1	02/17/2024 23:23	WG2228641
Chlorobenzene	U		0.116	1.00	1	02/17/2024 23:23	WG2228641
Chlorodibromomethane	U		0.140	1.00	1	02/17/2024 23:23	WG2228641
Chloroethane	U	L1	0.192	5.00	1	02/17/2024 23:23	WG2228641
Chloroform	U		0.111	5.00	1	02/17/2024 23:23	WG2228641
Chloromethane	U		0.960	2.50	1	02/17/2024 23:23	WG2228641
Cyclohexane	U		0.188	1.00	1	02/17/2024 23:23	WG2228641
2-Chlorotoluene	U		0.106	1.00	1	02/17/2024 23:23	WG2228641
4-Chlorotoluene	U		0.114	1.00	1	02/17/2024 23:23	WG2228641
1,2-Dibromo-3-Chloropropane	U		0.276	5.00	1	02/17/2024 23:23	WG2228641
1,2-Dibromoethane	U		0.126	1.00	1	02/17/2024 23:23	WG2228641
Dibromomethane	U		0.122	1.00	1	02/17/2024 23:23	WG2228641
1,2-Dichlorobenzene	U	L2	0.107	1.00	1	02/17/2024 23:23	WG2228641
1,3-Dichlorobenzene	U		0.110	1.00	1	02/17/2024 23:23	WG2228641
1,4-Dichlorobenzene	U		0.120	1.00	1	02/17/2024 23:23	WG2228641
Dichlorodifluoromethane	U		0.374	5.00	1	02/17/2024 23:23	WG2228641
1,1-Dichloroethane	U		0.100	1.00	1	02/17/2024 23:23	WG2228641
1,2-Dichloroethane	U		0.0819	1.00	1	02/17/2024 23:23	WG2228641
1,1-Dichloroethene	U		0.188	1.00	1	02/17/2024 23:23	WG2228641
cis-1,2-Dichloroethene	U		0.126	1.00	1	02/17/2024 23:23	WG2228641
trans-1,2-Dichloroethene	U		0.149	1.00	1	02/17/2024 23:23	WG2228641
1,2-Dichloropropane	U		0.149	1.00	1	02/17/2024 23:23	WG2228641
1,1-Dichloropropene	U		0.142	1.00	1	02/17/2024 23:23	WG2228641
1,3-Dichloropropane	U		0.110	1.00	1	02/17/2024 23:23	WG2228641
cis-1,3-Dichloropropene	U		0.111	1.00	1	02/17/2024 23:23	WG2228641
trans-1,3-Dichloropropene	U		0.118	1.00	1	02/17/2024 23:23	WG2228641
2,2-Dichloropropane	U		0.161	1.00	1	02/17/2024 23:23	WG2228641
Dicyclopentadiene	U		0.253	1.00	1	02/17/2024 23:23	WG2228641
Di-isopropyl ether	U		0.105	1.00	1	02/17/2024 23:23	WG2228641
Ethylbenzene	U		0.137	1.00	1	02/17/2024 23:23	WG2228641
4-Ethyltoluene	U		0.208	1.00	1	02/17/2024 23:23	WG2228641
Hexachloro-1,3-butadiene	U		0.337	1.00	1	02/17/2024 23:23	WG2228641
n-Hexane	U		0.749	10.0	1	02/17/2024 23:23	WG2228641
Isopropylbenzene	U		0.105	1.00	1	02/17/2024 23:23	WG2228641
p-Isopropyltoluene	U		0.120	1.00	1	02/17/2024 23:23	WG2228641
2-Butanone (MEK)	U		1.19	10.0	1	02/17/2024 23:23	WG2228641
Methyl Cyclohexane	U		0.660	1.00	1	02/17/2024 23:23	WG2228641
Methylene Chloride	U		0.430	5.00	1	02/17/2024 23:23	WG2228641
4-Methyl-2-pentanone (MIBK)	U		0.478	10.0	1	02/17/2024 23:23	WG2228641
Methyl tert-butyl ether	U		0.101	1.00	1	02/17/2024 23:23	WG2228641
Naphthalene	U	L2	1.00	5.00	1	02/17/2024 23:23	WG2228641
Propene	U		0.936	2.50	1	02/17/2024 23:23	WG2228641
n-Propylbenzene	U		0.0993	1.00	1	02/17/2024 23:23	WG2228641

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

Analyte	Result ug/l	Qualifier	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	Batch
Styrene	U		0.118	1.00	1	02/17/2024 23:23	WG2228641
1,1,1,2-Tetrachloroethane	U		0.147	1.00	1	02/17/2024 23:23	WG2228641
1,1,2,2-Tetrachloroethane	U		0.133	1.00	1	02/17/2024 23:23	WG2228641
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00	1	02/17/2024 23:23	WG2228641
Tetrachloroethene	U		0.300	1.00	1	02/17/2024 23:23	WG2228641
Toluene	U		0.278	1.00	1	02/17/2024 23:23	WG2228641
1,2,3-Trichlorobenzene	U		0.230	1.00	1	02/17/2024 23:23	WG2228641
1,2,4-Trichlorobenzene	U		0.481	1.00	1	02/17/2024 23:23	WG2228641
1,1,1-Trichloroethane	U		0.149	1.00	1	02/17/2024 23:23	WG2228641
1,1,2-Trichloroethane	U		0.158	1.00	1	02/17/2024 23:23	WG2228641
Trichloroethene	U		0.190	1.00	1	02/17/2024 23:23	WG2228641
Trichlorofluoromethane	U	L1	0.160	5.00	1	02/17/2024 23:23	WG2228641
1,2,3-Trichloropropane	U		0.237	2.50	1	02/17/2024 23:23	WG2228641
1,2,4-Trimethylbenzene	U		0.322	1.00	1	02/17/2024 23:23	WG2228641
1,2,3-Trimethylbenzene	U		0.104	1.00	1	02/17/2024 23:23	WG2228641
1,3,5-Trimethylbenzene	U		0.104	1.00	1	02/17/2024 23:23	WG2228641
Vinyl chloride	U		0.234	1.00	1	02/17/2024 23:23	WG2228641
Xylenes, Total	U		0.174	3.00	1	02/17/2024 23:23	WG2228641
(S) Toluene-d8	*108*			*80.0-120*		*02/17/2024 23:23*	WG2228641
(S) 4-Bromofluorobenzene	*99.7*			*77.0-126*		*02/17/2024 23:23*	WG2228641
(S) 1,2-Dichloroethane-d4	*87.7*			*70.0-130*		*02/17/2024 23:23*	WG2228641

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result ug/l	Qualifier	MDL ug/l	RDL ug/l	Dilution	Analysis date / time	Batch
1,4-Dioxane	U		0.597	3.00	1	02/20/2024 16:03	WG2230157
(S) Toluene-d8	*93.3*			*77.0-127*		*02/20/2024 16:03*	WG2230157

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

Method Blank (MB)

(MB) R4038381-1 02/20/24 23:34

Analyte	MB Result	MB Qualifier	MB MDL	MB RDL
	ug/l		ug/l	ug/l
Perchlorate	U		0.300	4.00

Method Blank (MB)

(MB) R4038995-1 02/24/24 22:05

Analyte	MB Result	MB Qualifier	MB MDL	MB RDL
	ug/l		ug/l	ug/l
Perchlorate	U		0.300	4.00

L1705776-08 Original Sample (OS) • Duplicate (DUP)

(OS) L1705776-08 02/21/24 05:09 • (DUP) R4038381-3 02/21/24 05:37

Analyte	Original Result	DUP Result	Dilution	DUP RPD	DUP Qualifier	DUP RPD Limits
	ug/l	ug/l		%		%
Perchlorate	6.06	5.84	1	3.68		15

L1705776-11 Original Sample (OS) • Duplicate (DUP)

(OS) L1705776-11 02/21/24 07:01 • (DUP) R4038381-4 02/21/24 07:29

Analyte	Original Result	DUP Result	Dilution	DUP RPD	DUP Qualifier	DUP RPD Limits
	ug/l	ug/l		%		%
Perchlorate	24000	22400	1000	6.68		15

L1705776-08 Original Sample (OS) • Duplicate (DUP)

(OS) L1705776-08 02/27/24 17:15 • (DUP) R4038995-7 02/27/24 17:43

Analyte	Original Result	DUP Result	Dilution	DUP RPD	DUP Qualifier	DUP RPD Limits
	ug/l	ug/l		%		%
Perchlorate	5.34	5.26	1	1.55		15

L1705776-11 Original Sample (OS) • Duplicate (DUP)

(OS) L1705776-11 02/27/24 18:11 • (DUP) R4038995-8 02/27/24 18:39

Analyte	Original Result ug/l	DUP Result ug/l	Dilution	DUP RPD %	DUP Qualifier	DUP RPD Limits %
Perchlorate	24100	25700	1000	6.46		15

Laboratory Control Sample (LCS)

(LCS) R4038381-2 02/21/24 00:30

Analyte	Spike Amount ug/l	LCS Result ug/l	LCS Rec. %	Rec. Limits %	LCS Qualifier
Perchlorate	10.0	9.43	94.3	90.0-110	

Laboratory Control Sample (LCS)

(LCS) R4038995-2 02/24/24 23:01

Analyte	Spike Amount ug/l	LCS Result ug/l	LCS Rec. %	Rec. Limits %	LCS Qualifier
Perchlorate	10.0	9.62	96.2	90.0-110	

L1705776-08 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1705776-08 02/27/24 17:15 • (MS) R4038995-3 02/25/24 04:08 • (MSD) R4038995-4 02/25/24 04:36

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MSD Result ug/l	MS Rec. %	MSD Rec. %	Dilution	Rec. Limits %	MS Qualifier	MSD Qualifier	RPD %	RPD Limits %
Perchlorate	10.0	5.34	14.8	16.1	94.2	108	1	80.0-120			8.89	15

L1705776-11 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1705776-11 02/27/24 18:11 • (MS) R4038995-5 02/25/24 05:04 • (MSD) R4038995-6 02/25/24 05:31

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MSD Result ug/l	MS Rec. %	MSD Rec. %	Dilution	Rec. Limits %	MS Qualifier	MSD Qualifier	RPD %	RPD Limits %
Perchlorate	10000	24100	32600	33700	85.1	96.9	1000	80.0-120			3.55	15

Method Blank (MB)

(MB) R4042116-1 02/29/24 13:52

Analyte	MB Result ug/l	MB Qualifier	MB MDL ug/l	MB RDL ug/l
Perchlorate	U		0.300	4.00

Method Blank (MB)

(MB) R4042116-2 02/29/24 14:50

Analyte	MB Result ug/l	MB Qualifier	MB MDL ug/l	MB RDL ug/l
Perchlorate	U		0.300	4.00

Method Blank (MB)

(MB) R4042116-4 02/29/24 16:14

Analyte	MB Result ug/l	MB Qualifier	MB MDL ug/l	MB RDL ug/l
Perchlorate	U		0.300	4.00

Method Blank (MB)

(MB) R4042116-14 03/01/24 10:15

Analyte	MB Result ug/l	MB Qualifier	MB MDL ug/l	MB RDL ug/l
Perchlorate	U		0.300	4.00

L1706680-03 Original Sample (OS) • Duplicate (DUP)

(OS) L1706680-03 03/01/24 00:29 • (DUP) R4042116-5 03/01/24 00:57

Analyte	Original Result ug/l	DUP Result ug/l	Dilution	DUP RPD %	DUP Qualifier	DUP RPD Limits %
Perchlorate	10100	12700	1000	22.6	R8	15

L1706680-06 Original Sample (OS) • Duplicate (DUP)

(OS) L1706680-06 03/01/24 03:16 • (DUP) R4042116-6 03/01/24 03:44

Analyte	Original Result ug/l	DUP Result ug/l	Dilution	DUP RPD %	DUP Qualifier	DUP RPD Limits %
Perchlorate	5.54	5.53	1	0.255		15

Laboratory Control Sample (LCS)

(LCS) R4042116-3 02/29/24 15:46

Analyte	Spike Amount ug/l	LCS Result ug/l	LCS Rec. %	Rec. Limits %	LCS Qualifier
Perchlorate	10.0	9.65	96.5	90.0-110	

L1706682-02 Original Sample (OS) • Matrix Spike (MS)

(OS) L1706682-02 03/01/24 05:36 • (MS) R4042116-7 03/01/24 06:04

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MS Rec. %	Dilution	Rec. Limits %	MS Qualifier
Perchlorate	10.0	U	10.2	102	1	80.0-120	

L1705776-23 Original Sample (OS) • Matrix Spike (MS)

(OS) L1705776-23 02/29/24 18:06 • (MS) R4042116-8 03/01/24 06:32

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MS Rec. %	Dilution	Rec. Limits %	MS Qualifier
Perchlorate	10.0	U	7.94	79.4	1	80.0-120	M2

L1706680-03 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1706680-03 03/01/24 00:29 • (MS) R4042116-9 03/01/24 07:00 • (MSD) R4042116-10 03/01/24 08:23

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MSD Result ug/l	MS Rec. %	MSD Rec. %	Dilution	Rec. Limits %	MS Qualifier	MSD Qualifier	RPD %	RPD Limits %
Perchlorate	10.0	10100	20100	22600	99800	125000	1000	80.0-120	M3	M3	11.9	15

L1706680-06 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1706680-06 03/01/24 03:16 • (MS) R4042116-11 03/01/24 08:51 • (MSD) R4042116-12 03/01/24 09:19

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MSD Result ug/l	MS Rec. %	MSD Rec. %	Dilution	Rec. Limits %	MS Qualifier	MSD Qualifier	RPD %	RPD Limits %
Perchlorate	10.0	5.54	14.9	15.5	93.3	99.5	1	80.0-120			4.09	15

L1706680-08 Original Sample (OS) • Matrix Spike (MS)

(OS) L1706680-08 03/01/24 04:40 • (MS) R4042116-13 03/01/24 09:47

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MS Rec. %	Dilution	Rec. Limits %	MS Qualifier
Perchlorate	10.0	4.55	13.7	91.8	1	80.0-120	

Method Blank (MB)

(MB) R4035581-3 02/17/24 08:23

Analyte	MB Result ug/l	MB Qualifier	MB MDL ug/l	MB RDL ug/l
Acetone	U		11.3	50.0
Acrolein	U		2.54	50.0
Acrylonitrile	U		0.671	10.0
Benzene	U		0.0941	1.00
Bromobenzene	U		0.118	1.00
Bromodichloromethane	U		0.136	1.00
Bromoform	U		0.129	1.00
Bromomethane	U		0.605	5.00
1,3-Butadiene	U		0.299	2.00
n-Butylbenzene	U		0.157	1.00
sec-Butylbenzene	U		0.125	1.00
tert-Butylbenzene	U		0.127	1.00
Carbon tetrachloride	U		0.128	1.00
Carbon disulfide	U		0.0962	1.00
Chlorobenzene	U		0.116	1.00
Chlorodibromomethane	U		0.140	1.00
Chloroethane	U		0.192	5.00
Chloroform	U		0.111	5.00
Chloromethane	U		0.960	2.50
Cyclohexane	U		0.188	1.00
2-Chlorotoluene	U		0.106	1.00
4-Chlorotoluene	U		0.114	1.00
1,2-Dibromo-3-Chloropropane	U		0.276	5.00
1,2-Dibromoethane	U		0.126	1.00
Dibromomethane	U		0.122	1.00
1,2-Dichlorobenzene	U		0.107	1.00
1,3-Dichlorobenzene	U		0.110	1.00
1,4-Dichlorobenzene	U		0.120	1.00
Dichlorodifluoromethane	U		0.374	5.00
1,1-Dichloroethane	U		0.100	1.00
1,2-Dichloroethane	U		0.0819	1.00
1,1-Dichloroethene	U		0.188	1.00
cis-1,2-Dichloroethene	U		0.126	1.00
trans-1,2-Dichloroethene	U		0.149	1.00
1,2-Dichloropropane	U		0.149	1.00
1,1-Dichloropropene	U		0.142	1.00
1,3-Dichloropropane	U		0.110	1.00
cis-1,3-Dichloropropene	U		0.111	1.00
trans-1,3-Dichloropropene	U		0.118	1.00
2,2-Dichloropropane	U		0.161	1.00

Method Blank (MB)

(MB) R4035581-3 02/17/24 08:23

Analyte	MB Result ug/l	MB Qualifier	MB MDL ug/l	MB RDL ug/l
Dicyclopentadiene	U		0.253	1.00
Di-isopropyl ether	U		0.105	1.00
Ethylbenzene	U		0.137	1.00
4-Ethyltoluene	U		0.208	1.00
Hexachloro-1,3-butadiene	U		0.337	1.00
n-Hexane	U		0.749	10.0
Isopropylbenzene	U		0.105	1.00
p-Isopropyltoluene	U		0.120	1.00
2-Butanone (MEK)	U		1.19	10.0
Methyl Cyclohexane	U		0.660	1.00
Methylene Chloride	U		0.430	5.00
4-Methyl-2-pentanone (MIBK)	U		0.478	10.0
Methyl tert-butyl ether	U		0.101	1.00
Naphthalene	U		1.00	5.00
Propene	U		0.936	2.50
n-Propylbenzene	U		0.0993	1.00
Styrene	U		0.118	1.00
1,1,1,2-Tetrachloroethane	U		0.147	1.00
1,1,2,2-Tetrachloroethane	U		0.133	1.00
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00
Tetrachloroethene	U		0.300	1.00
Toluene	U		0.278	1.00
1,2,3-Trichlorobenzene	U		0.230	1.00
1,2,4-Trichlorobenzene	U		0.481	1.00
1,1,1-Trichloroethane	U		0.149	1.00
1,1,2-Trichloroethane	U		0.158	1.00
Trichloroethene	U		0.190	1.00
Trichlorofluoromethane	U		0.160	5.00
1,2,3-Trichloropropane	U		0.237	2.50
1,2,4-Trimethylbenzene	U		0.322	1.00
1,2,3-Trimethylbenzene	U		0.104	1.00
1,3,5-Trimethylbenzene	U		0.104	1.00
Vinyl chloride	U		0.234	1.00
Xylenes, Total	U		0.174	3.00
(S) Toluene-d8	*102*			*80.0-120*
(S) 4-Bromofluorobenzene	*91.3*			*77.0-126*
(S) 1,2-Dichloroethane-d4	*105*			*70.0-130*

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R4035581-1 02/17/24 07:22 • (LCSD) R4035581-2 02/17/24 07:42

Analyte	Spike Amount ug/l	LCS Result ug/l	LCSD Result ug/l	LCS Rec. %	LCSD Rec. %	Rec. Limits %	LCS Qualifier	LCSD Qualifier	RPD %	RPD Limits %
Acetone	25.0	26.2	21.4	105	85.6	19.0-160			20.2	27
Acrolein	25.0	25.4	27.2	102	109	30.0-160			6.84	26
Acrylonitrile	25.0	24.7	25.0	98.8	100	55.0-149			1.21	20
Benzene	5.00	5.25	5.42	105	108	70.0-123			3.19	20
Bromobenzene	5.00	5.27	5.58	105	112	73.0-121			5.71	20
Bromodichloromethane	5.00	5.33	5.53	107	111	75.0-120			3.68	20
Bromoform	5.00	5.16	5.70	103	114	68.0-132			9.94	20
Bromomethane	5.00	4.64	5.12	92.8	102	30.0-160			9.84	25
1,3-Butadiene	5.00	5.49	5.45	110	109	45.0-147			0.731	20
n-Butylbenzene	5.00	4.46	4.68	89.2	93.6	73.0-125			4.81	20
sec-Butylbenzene	5.00	4.59	4.75	91.8	95.0	75.0-125			3.43	20
tert-Butylbenzene	5.00	4.84	5.17	96.8	103	76.0-124			6.59	20
Carbon tetrachloride	5.00	5.81	5.68	116	114	68.0-126			2.26	20
Carbon disulfide	5.00	5.29	5.71	106	114	61.0-128			7.64	20
Chlorobenzene	5.00	5.28	5.69	106	114	80.0-121			7.47	20
Chlorodibromomethane	5.00	5.41	5.51	108	110	77.0-125			1.83	20
Chloroethane	5.00	5.17	5.47	103	109	47.0-150			5.64	20
Chloroform	5.00	5.15	5.31	103	106	73.0-120			3.06	20
Chloromethane	5.00	5.15	5.96	103	119	41.0-142			14.6	20
Cyclohexane	5.00	4.63	4.79	92.6	95.8	71.0-124			3.40	20
2-Chlorotoluene	5.00	5.11	5.42	102	108	76.0-123			5.89	20
4-Chlorotoluene	5.00	5.14	5.24	103	105	75.0-122			1.93	20
1,2-Dibromo-3-Chloropropane	5.00	4.64	4.73	92.8	94.6	58.0-134			1.92	20
1,2-Dibromoethane	5.00	5.31	5.33	106	107	80.0-122			0.376	20
Dibromomethane	5.00	5.44	5.30	109	106	80.0-120			2.61	20
1,2-Dichlorobenzene	5.00	5.15	5.53	103	111	79.0-121			7.12	20
1,3-Dichlorobenzene	5.00	5.15	5.33	103	107	79.0-120			3.44	20
1,4-Dichlorobenzene	5.00	5.28	5.58	106	112	79.0-120			5.52	20
Dichlorodifluoromethane	5.00	4.49	5.03	89.8	101	51.0-149			11.3	20
1,1-Dichloroethane	5.00	5.22	5.53	104	111	70.0-126			5.77	20
1,2-Dichloroethane	5.00	5.41	5.46	108	109	70.0-128			0.920	20
1,1-Dichloroethene	5.00	4.93	5.16	98.6	103	71.0-124			4.56	20
cis-1,2-Dichloroethene	5.00	5.25	5.57	105	111	73.0-120			5.91	20
trans-1,2-Dichloroethene	5.00	5.27	5.53	105	111	73.0-120			4.81	20
1,2-Dichloropropane	5.00	5.06	5.38	101	108	77.0-125			6.13	20
1,1-Dichloropropene	5.00	5.40	5.78	108	116	74.0-126			6.80	20
1,3-Dichloropropane	5.00	5.19	5.25	104	105	80.0-120			1.15	20
cis-1,3-Dichloropropene	5.00	5.16	5.26	103	105	80.0-123			1.92	20
trans-1,3-Dichloropropene	5.00	5.08	5.63	102	113	78.0-124			10.3	20
2,2-Dichloropropane	5.00	5.78	5.79	116	116	58.0-130			0.173	20

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R4035581-1 02/17/24 07:22 • (LCSD) R4035581-2 02/17/24 07:42

Analyte	Spike Amount ug/l	LCS Result ug/l	LCSD Result ug/l	LCS Rec. %	LCSD Rec. %	Rec. Limits %	LCS Qualifier	LCSD Qualifier	RPD %	RPD Limits %
Dicyclopentadiene	5.00	5.24	5.63	105	113	74.0-126			7.18	20
Di-isopropyl ether	5.00	5.10	5.31	102	106	58.0-138			4.03	20
Ethylbenzene	5.00	4.74	5.44	94.8	109	79.0-123			13.8	20
4-Ethyltoluene	5.00	4.84	5.32	96.8	106	74.0-127			9.45	20
Hexachloro-1,3-butadiene	5.00	5.65	5.44	113	109	54.0-138			3.79	20
n-Hexane	5.00	5.49	5.04	110	101	57.0-133			8.55	20
Isopropylbenzene	5.00	5.09	5.37	102	107	76.0-127			5.35	20
p-Isopropyltoluene	5.00	4.55	4.71	91.0	94.2	76.0-125			3.46	20
2-Butanone (MEK)	25.0	25.4	24.1	102	96.4	44.0-160			5.25	20
Methyl Cyclohexane	5.00	4.54	4.46	90.8	89.2	68.0-126			1.78	20
Methylene Chloride	5.00	5.56	5.47	111	109	67.0-120			1.63	20
4-Methyl-2-pentanone (MIBK)	25.0	25.4	26.2	102	105	68.0-142			3.10	20
Methyl tert-butyl ether	5.00	5.26	5.35	105	107	68.0-125			1.70	20
Naphthalene	5.00	3.86	4.50	77.2	90.0	54.0-135			15.3	20
Propene	5.00	4.67	4.56	93.4	91.2	30.0-160			2.38	20
n-Propylbenzene	5.00	4.89	5.10	97.8	102	77.0-124			4.20	20
Styrene	5.00	5.02	5.52	100	110	73.0-130			9.49	20
1,1,1,2-Tetrachloroethane	5.00	5.12	5.50	102	110	75.0-125			7.16	20
1,1,2,2-Tetrachloroethane	5.00	5.42	5.69	108	114	65.0-130			4.86	20
1,1,2-Trichlorotrifluoroethane	5.00	5.24	5.04	105	101	69.0-132			3.89	20
Tetrachloroethene	5.00	5.44	5.59	109	112	72.0-132			2.72	20
Toluene	5.00	5.07	5.48	101	110	79.0-120			7.77	20
1,2,3-Trichlorobenzene	5.00	5.01	5.69	100	114	50.0-138			12.7	20
1,2,4-Trichlorobenzene	5.00	4.88	5.84	97.6	117	57.0-137			17.9	20
1,1,1-Trichloroethane	5.00	5.48	5.57	110	111	73.0-124			1.63	20
1,1,2-Trichloroethane	5.00	5.41	5.91	108	118	80.0-120			8.83	20
Trichloroethene	5.00	5.13	5.26	103	105	78.0-124			2.50	20
Trichlorofluoromethane	5.00	5.67	5.71	113	114	59.0-147			0.703	20
1,2,3-Trichloropropane	5.00	5.18	5.24	104	105	73.0-130			1.15	20
1,2,4-Trimethylbenzene	5.00	4.66	5.12	93.2	102	76.0-121			9.41	20
1,2,3-Trimethylbenzene	5.00	5.01	5.26	100	105	77.0-120			4.87	20
1,3,5-Trimethylbenzene	5.00	4.88	5.12	97.6	102	76.0-122			4.80	20
Vinyl chloride	5.00	5.08	5.39	102	108	67.0-131			5.92	20
Xylenes, Total	15.0	14.7	16.2	98.0	108	79.0-123			9.71	20
(S) Toluene-d8				*97.2*	*100*	*80.0-120*				
(S) 4-Bromofluorobenzene				*91.0*	*94.6*	*77.0-126*				
(S) 1,2-Dichloroethane-d4				*102*	*99.5*	*70.0-130*				

L1705776-08 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1705776-08 02/17/24 13:35 • (MS) R4035581-4 02/17/24 18:23 • (MSD) R4035581-5 02/17/24 18:44

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MSD Result ug/l	MS Rec. %	MSD Rec. %	Dilution	Rec. Limits %	MS Qualifier	MSD Qualifier	RPD %	RPD Limits %
Acetone	25.0	U	17.9	17.2	71.6	68.8	1	10.0-160			3.99	35
Acrolein	25.0	U	17.8	19.3	71.2	77.2	1	10.0-160			8.09	39
Acrylonitrile	25.0	U	19.3	19.9	77.2	79.6	1	21.0-160			3.06	32
Benzene	5.00	U	4.11	4.19	82.2	83.8	1	17.0-158			1.93	27
Bromobenzene	5.00	U	4.26	4.27	85.2	85.4	1	30.0-149			0.234	28
Bromodichloromethane	5.00	U	4.22	4.57	84.4	91.4	1	31.0-150			7.96	27
Bromoform	5.00	U	4.19	4.28	83.8	85.6	1	29.0-150			2.13	29
Bromomethane	5.00	U	3.66	4.08	73.2	81.6	1	10.0-160			10.9	38
1,3-Butadiene	5.00	U	4.26	4.94	85.2	98.8	1	10.0-160			14.8	22
n-Butylbenzene	5.00	U	3.39	3.50	67.8	70.0	1	31.0-150			3.19	30
sec-Butylbenzene	5.00	U	3.63	3.93	72.6	78.6	1	33.0-155			7.94	29
tert-Butylbenzene	5.00	U	3.89	4.03	77.8	80.6	1	34.0-153			3.54	28
Carbon tetrachloride	5.00	U	4.90	5.06	98.0	101	1	23.0-159			3.21	28
Carbon disulfide	5.00	U	3.56	3.65	71.2	73.0	1	10.0-156			2.50	28
Chlorobenzene	5.00	U	4.05	4.32	81.0	86.4	1	33.0-152			6.45	27
Chlorodibromomethane	5.00	U	4.14	4.19	82.8	83.8	1	37.0-149			1.20	27
Chloroethane	5.00	U	5.07	4.51	101	90.2	1	10.0-160			11.7	30
Chloroform	5.00	U	4.09	4.42	81.8	88.4	1	29.0-154			7.76	28
Chloromethane	5.00	U	4.15	4.48	83.0	89.6	1	10.0-160			7.65	29
Cyclohexane	5.00	U	3.91	3.93	78.2	78.6	1	19.0-160			0.510	23
2-Chlorotoluene	5.00	U	4.04	4.14	80.8	82.8	1	32.0-153			2.44	28
4-Chlorotoluene	5.00	U	3.93	4.19	78.6	83.8	1	32.0-150			6.40	28
1,2-Dibromo-3-Chloropropane	5.00	U	3.52	3.86	70.4	77.2	1	22.0-151			9.21	34
1,2-Dibromoethane	5.00	U	3.89	4.14	77.8	82.8	1	34.0-147			6.23	27
Dibromomethane	5.00	U	4.25	4.34	85.0	86.8	1	30.0-151			2.10	27
1,2-Dichlorobenzene	5.00	U	4.09	4.43	81.8	88.6	1	34.0-149			7.98	28
1,3-Dichlorobenzene	5.00	U	4.10	4.25	82.0	85.0	1	36.0-146			3.59	27
1,4-Dichlorobenzene	5.00	U	4.32	4.37	86.4	87.4	1	35.0-142			1.15	27
Dichlorodifluoromethane	5.00	U	4.00	3.98	80.0	79.6	1	10.0-160			0.501	29
1,1-Dichloroethane	5.00	U	4.33	4.40	86.6	88.0	1	25.0-158			1.60	27
1,2-Dichloroethane	5.00	U	4.15	4.41	83.0	88.2	1	29.0-151			6.07	27
1,1-Dichloroethene	5.00	U	4.12	4.17	82.4	83.4	1	11.0-160			1.21	29
cis-1,2-Dichloroethene	5.00	U	4.16	4.30	83.2	86.0	1	10.0-160			3.31	27
trans-1,2-Dichloroethene	5.00	U	4.18	4.17	83.6	83.4	1	17.0-153			0.240	27
1,2-Dichloropropane	5.00	U	4.05	4.26	81.0	85.2	1	30.0-156			5.05	27
1,1-Dichloropropene	5.00	U	4.26	4.66	85.2	93.2	1	25.0-158			8.97	27
1,3-Dichloropropane	5.00	U	3.98	4.34	79.6	86.8	1	38.0-147			8.65	27
cis-1,3-Dichloropropene	5.00	U	3.71	3.66	74.2	73.2	1	34.0-149			1.36	28
trans-1,3-Dichloropropene	5.00	U	3.88	4.07	77.6	81.4	1	32.0-149			4.78	28
2,2-Dichloropropane	5.00	U	4.29	4.42	85.8	88.4	1	24.0-152			2.99	29

L1705776-08 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1705776-08 02/17/24 13:35 • (MS) R4035581-4 02/17/24 18:23 • (MSD) R4035581-5 02/17/24 18:44

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MSD Result ug/l	MS Rec. %	MSD Rec. %	Dilution	Rec. Limits %	MS Qualifier	MSD Qualifier	RPD %	RPD Limits %
Dicyclopentadiene	5.00	U	2.80	2.83	56.0	56.6	1	51.0-139			1.07	20
Di-isopropyl ether	5.00	U	3.91	4.15	78.2	83.0	1	21.0-160			5.96	28
Ethylbenzene	5.00	U	3.86	4.09	77.2	81.8	1	30.0-155			5.79	27
4-Ethyltoluene	5.00	U	3.89	4.00	77.8	80.0	1	10.0-160			2.79	20
Hexachloro-1,3-butadiene	5.00	U	4.26	4.32	85.2	86.4	1	20.0-154			1.40	34
n-Hexane	5.00	U	4.28	4.51	85.6	90.2	1	10.0-153			5.23	28
Isopropylbenzene	5.00	U	3.99	4.22	79.8	84.4	1	28.0-157			5.60	27
p-Isopropyltoluene	5.00	U	3.62	3.64	72.4	72.8	1	30.0-154			0.551	29
2-Butanone (MEK)	25.0	U	17.8	18.9	71.2	75.6	1	10.0-160			5.99	32
Methyl Cyclohexane	5.00	U	3.75	3.97	75.0	79.4	1	11.0-160			5.70	24
Methylene Chloride	5.00	U	4.26	4.40	85.2	88.0	1	23.0-144			3.23	28
4-Methyl-2-pentanone (MIBK)	25.0	U	18.7	19.6	74.8	78.4	1	29.0-160			4.70	29
Methyl tert-butyl ether	5.00	U	3.92	4.18	78.4	83.6	1	28.0-150			6.42	29
Naphthalene	5.00	U	3.05	3.23	61.0	64.6	1	12.0-156			5.73	35
Propene	5.00	U	3.09	3.34	61.8	66.8	1	10.0-160			7.78	29
n-Propylbenzene	5.00	U	3.98	4.15	79.6	83.0	1	31.0-154			4.18	28
Styrene	5.00	U	3.70	3.86	74.0	77.2	1	33.0-155			4.23	28
1,1,1,2-Tetrachloroethane	5.00	U	4.13	4.43	82.6	88.6	1	36.0-151			7.01	29
1,1,2,2-Tetrachloroethane	5.00	U	4.39	4.65	87.8	93.0	1	33.0-150			5.75	28
1,1,2-Trichlorotrifluoroethane	5.00	U	4.45	4.58	89.0	91.6	1	23.0-160			2.88	30
Tetrachloroethene	5.00	U	4.28	4.36	85.6	87.2	1	10.0-160			1.85	27
Toluene	5.00	U	4.01	4.13	80.2	82.6	1	26.0-154			2.95	28
1,2,3-Trichlorobenzene	5.00	U	3.75	3.76	75.0	75.2	1	17.0-150			0.266	36
1,2,4-Trichlorobenzene	5.00	U	3.59	3.88	71.8	77.6	1	24.0-150			7.76	33
1,1,1-Trichloroethane	5.00	U	4.56	4.80	91.2	96.0	1	23.0-160			5.13	28
1,1,2-Trichloroethane	5.00	U	4.30	4.31	86.0	86.2	1	35.0-147			0.232	27
Trichloroethene	5.00	U	4.63	4.74	92.6	94.8	1	10.0-160			2.35	25
Trichlorofluoromethane	5.00	U	4.86	5.03	97.2	101	1	17.0-160			3.44	31
1,2,3-Trichloropropane	5.00	U	4.18	4.15	83.6	83.0	1	34.0-151			0.720	29
1,2,4-Trimethylbenzene	5.00	U	3.83	4.05	76.6	81.0	1	26.0-154			5.58	27
1,2,3-Trimethylbenzene	5.00	U	3.70	4.13	74.0	82.6	1	32.0-149			11.0	28
1,3,5-Trimethylbenzene	5.00	U	3.71	3.92	74.2	78.4	1	28.0-153			5.50	27
Vinyl chloride	5.00	U	4.18	4.52	83.6	90.4	1	10.0-160			7.82	27
Xylenes, Total	15.0	U	11.4	11.9	76.0	79.3	1	29.0-154			4.29	28
(S) Toluene-d8					*93.6*	*92.1*		*80.0-120*				
(S) 4-Bromofluorobenzene					*89.9*	*89.8*		*77.0-126*				
(S) 1,2-Dichloroethane-d4					*101*	*101*		*70.0-130*				

L1705776-11 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1705776-11 02/17/24 13:56 • (MS) R4035581-6 02/17/24 19:04 • (MSD) R4035581-7 02/17/24 19:25

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MSD Result ug/l	MS Rec. %	MSD Rec. %	Dilution	Rec. Limits %	MS Qualifier	MSD Qualifier	RPD %	RPD Limits %
Acetone	25.0	U	19.3	19.4	77.2	77.6	1	10.0-160			0.517	35
Acrolein	25.0	U	22.3	21.7	89.2	86.8	1	10.0-160			2.73	39
Acrylonitrile	25.0	U	21.7	22.3	86.8	89.2	1	21.0-160			2.73	32
Benzene	5.00	U	4.55	4.68	91.0	93.6	1	17.0-158			2.82	27
Bromobenzene	5.00	U	4.47	4.87	89.4	97.4	1	30.0-149			8.57	28
Bromodichloromethane	5.00	U	4.79	4.93	95.8	98.6	1	31.0-150			2.88	27
Bromoform	5.00	0.241	5.21	5.27	99.4	101	1	29.0-150			1.15	29
Bromomethane	5.00	U	4.34	4.55	86.8	91.0	1	10.0-160			4.72	38
1,3-Butadiene	5.00	U	4.60	4.63	92.0	92.6	1	10.0-160			0.650	22
n-Butylbenzene	5.00	U	3.65	3.71	73.0	74.2	1	31.0-150			1.63	30
sec-Butylbenzene	5.00	U	3.93	4.45	78.6	89.0	1	33.0-155			12.4	29
tert-Butylbenzene	5.00	U	4.22	4.67	84.4	93.4	1	34.0-153			10.1	28
Carbon tetrachloride	5.00	U	5.30	5.28	106	106	1	23.0-159			0.378	28
Carbon disulfide	5.00	U	3.82	3.96	76.4	79.2	1	10.0-156			3.60	28
Chlorobenzene	5.00	U	4.69	4.81	93.8	96.2	1	33.0-152			2.53	27
Chlorodibromomethane	5.00	U	4.79	4.90	95.8	98.0	1	37.0-149			2.27	27
Chloroethane	5.00	U	4.98	5.21	99.6	104	1	10.0-160			4.51	30
Chloroform	5.00	U	4.93	4.98	98.6	99.6	1	29.0-154			1.01	28
Chloromethane	5.00	U	4.77	4.72	95.4	94.4	1	10.0-160			1.05	29
Cyclohexane	5.00	U	4.15	4.35	83.0	87.0	1	19.0-160			4.71	23
2-Chlorotoluene	5.00	U	4.47	4.86	89.4	97.2	1	32.0-153			8.36	28
4-Chlorotoluene	5.00	U	4.37	4.61	87.4	92.2	1	32.0-150			5.35	28
1,2-Dibromo-3-Chloropropane	5.00	U	3.78	4.40	75.6	88.0	1	22.0-151			15.2	34
1,2-Dibromoethane	5.00	U	4.45	4.53	89.0	90.6	1	34.0-147			1.78	27
Dibromomethane	5.00	U	4.65	4.68	93.0	93.6	1	30.0-151			0.643	27
1,2-Dichlorobenzene	5.00	U	4.36	4.68	87.2	93.6	1	34.0-149			7.08	28
1,3-Dichlorobenzene	5.00	U	4.57	4.70	91.4	94.0	1	36.0-146			2.80	27
1,4-Dichlorobenzene	5.00	U	4.49	4.92	89.8	98.4	1	35.0-142			9.14	27
Dichlorodifluoromethane	5.00	U	4.30	4.19	86.0	83.8	1	10.0-160			2.59	29
1,1-Dichloroethane	5.00	U	4.66	4.81	93.2	96.2	1	25.0-158			3.17	27
1,2-Dichloroethane	5.00	U	4.62	4.73	92.4	94.6	1	29.0-151			2.35	27
1,1-Dichloroethene	5.00	4.80	8.77	8.60	79.4	76.0	1	11.0-160			1.96	29
cis-1,2-Dichloroethene	5.00	U	4.56	4.83	91.2	96.6	1	10.0-160			5.75	27
trans-1,2-Dichloroethene	5.00	U	4.31	4.66	86.2	93.2	1	17.0-153			7.80	27
1,2-Dichloropropane	5.00	U	4.51	4.68	90.2	93.6	1	30.0-156			3.70	27
1,1-Dichloropropene	5.00	U	4.75	4.97	95.0	99.4	1	25.0-158			4.53	27
1,3-Dichloropropane	5.00	U	4.65	4.79	93.0	95.8	1	38.0-147			2.97	27
cis-1,3-Dichloropropene	5.00	U	4.13	4.25	82.6	85.0	1	34.0-149			2.86	28
trans-1,3-Dichloropropene	5.00	U	4.46	4.65	89.2	93.0	1	32.0-149			4.17	28
2,2-Dichloropropane	5.00	U	4.79	4.76	95.8	95.2	1	24.0-152			0.628	29

L1705776-11 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1705776-11 02/17/24 13:56 • (MS) R4035581-6 02/17/24 19:04 • (MSD) R4035581-7 02/17/24 19:25

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MSD Result ug/l	MS Rec. %	MSD Rec. %	Dilution	Rec. Limits %	MS Qualifier	MSD Qualifier	RPD %	RPD Limits %
Dicyclopentadiene	5.00	U	3.34	3.08	66.8	61.6	1	51.0-139			8.10	20
Di-isopropyl ether	5.00	U	4.54	4.56	90.8	91.2	1	21.0-160			0.440	28
Ethylbenzene	5.00	U	4.45	4.69	89.0	93.8	1	30.0-155			5.25	27
4-Ethyltoluene	5.00	U	4.23	4.66	84.6	93.2	1	10.0-160			9.67	20
Hexachloro-1,3-butadiene	5.00	U	4.81	4.85	96.2	97.0	1	20.0-154			0.828	34
n-Hexane	5.00	U	4.64	4.57	92.8	91.4	1	10.0-153			1.52	28
Isopropylbenzene	5.00	U	4.57	4.71	91.4	94.2	1	28.0-157			3.02	27
p-Isopropyltoluene	5.00	U	3.80	4.07	76.0	81.4	1	30.0-154			6.86	29
2-Butanone (MEK)	25.0	U	21.2	21.7	84.8	86.8	1	10.0-160			2.33	32
Methyl Cyclohexane	5.00	U	4.04	4.15	80.8	83.0	1	11.0-160			2.69	24
Methylene Chloride	5.00	U	4.75	4.75	95.0	95.0	1	23.0-144			0.000	28
4-Methyl-2-pentanone (MIBK)	25.0	U	22.6	22.8	90.4	91.2	1	29.0-160			0.881	29
Methyl tert-butyl ether	5.00	U	4.41	4.69	88.2	93.8	1	28.0-150			6.15	29
Naphthalene	5.00	U	3.92	4.07	78.4	81.4	1	12.0-156			3.75	35
Propene	5.00	U	3.22	3.20	64.4	64.0	1	10.0-160			0.623	29
n-Propylbenzene	5.00	U	4.22	4.75	84.4	95.0	1	31.0-154			11.8	28
Styrene	5.00	U	4.24	4.23	84.8	84.6	1	33.0-155			0.236	28
1,1,1,2-Tetrachloroethane	5.00	U	4.74	5.31	94.8	106	1	36.0-151			11.3	29
1,1,2,2-Tetrachloroethane	5.00	U	4.84	5.16	96.8	103	1	33.0-150			6.40	28
1,1,2-Trichlorotrifluoroethane	5.00	U	5.03	4.84	101	96.8	1	23.0-160			3.85	30
Tetrachloroethene	5.00	U	4.67	5.08	93.4	102	1	10.0-160			8.41	27
Toluene	5.00	U	4.60	4.75	92.0	95.0	1	26.0-154			3.21	28
1,2,3-Trichlorobenzene	5.00	U	4.67	5.05	93.4	101	1	17.0-150			7.82	36
1,2,4-Trichlorobenzene	5.00	U	4.61	4.56	92.2	91.2	1	24.0-150			1.09	33
1,1,1-Trichloroethane	5.00	U	5.09	5.21	102	104	1	23.0-160			2.33	28
1,1,2-Trichloroethane	5.00	U	4.96	4.87	99.2	97.4	1	35.0-147			1.83	27
Trichloroethene	5.00	13.6	18.2	17.7	92.0	82.0	1	10.0-160			2.79	25
Trichlorofluoromethane	5.00	U	5.43	5.17	109	103	1	17.0-160			4.91	31
1,2,3-Trichloropropane	5.00	U	4.77	4.88	95.4	97.6	1	34.0-151			2.28	29
1,2,4-Trimethylbenzene	5.00	U	4.10	4.37	82.0	87.4	1	26.0-154			6.38	27
1,2,3-Trimethylbenzene	5.00	U	4.28	4.56	85.6	91.2	1	32.0-149			6.33	28
1,3,5-Trimethylbenzene	5.00	U	4.05	4.43	81.0	88.6	1	28.0-153			8.96	27
Vinyl chloride	5.00	U	4.58	4.65	91.6	93.0	1	10.0-160			1.52	27
Xylenes, Total	15.0	U	13.2	13.8	88.0	92.0	1	29.0-154			4.44	28
(S) Toluene-d8					*95.1*	*96.1*		*80.0-120*				
(S) 4-Bromofluorobenzene					*91.3*	*93.5*		*77.0-126*				
(S) 1,2-Dichloroethane-d4					*102*	*100*		*70.0-130*				

Method Blank (MB)

(MB) R4035528-3 02/17/24 21:24

Analyte	MB Result ug/l	MB Qualifier	MB MDL ug/l	MB RDL ug/l
Acetone	U		11.3	50.0
Acrolein	U		2.54	50.0
Acrylonitrile	U		0.671	10.0
Benzene	U		0.0941	1.00
Bromobenzene	U		0.118	1.00
Bromodichloromethane	U		0.136	1.00
Bromoform	U		0.129	1.00
Bromomethane	U		0.605	5.00
1,3-Butadiene	U		0.299	2.00
n-Butylbenzene	U		0.157	1.00
sec-Butylbenzene	U		0.125	1.00
tert-Butylbenzene	U		0.127	1.00
Carbon tetrachloride	U		0.128	1.00
Carbon disulfide	U		0.0962	1.00
Chlorobenzene	U		0.116	1.00
Chlorodibromomethane	U		0.140	1.00
Chloroethane	U		0.192	5.00
Chloroform	U		0.111	5.00
Chloromethane	U		0.960	2.50
Cyclohexane	U		0.188	1.00
2-Chlorotoluene	U		0.106	1.00
4-Chlorotoluene	U		0.114	1.00
1,2-Dibromo-3-Chloropropane	U		0.276	5.00
1,2-Dibromoethane	U		0.126	1.00
Dibromomethane	U		0.122	1.00
1,2-Dichlorobenzene	U		0.107	1.00
1,3-Dichlorobenzene	U		0.110	1.00
1,4-Dichlorobenzene	U		0.120	1.00
Dichlorodifluoromethane	U		0.374	5.00
1,1-Dichloroethane	U		0.100	1.00
1,2-Dichloroethane	U		0.0819	1.00
1,1-Dichloroethene	U		0.188	1.00
cis-1,2-Dichloroethene	U		0.126	1.00
trans-1,2-Dichloroethene	U		0.149	1.00
1,2-Dichloropropane	U		0.149	1.00
1,1-Dichloropropene	U		0.142	1.00
1,3-Dichloropropane	U		0.110	1.00
cis-1,3-Dichloropropene	U		0.111	1.00
trans-1,3-Dichloropropene	U		0.118	1.00
2,2-Dichloropropane	U		0.161	1.00

Method Blank (MB)

(MB) R4035528-3 02/17/24 21:24

Analyte	MB Result ug/l	MB Qualifier	MB MDL ug/l	MB RDL ug/l
Dicyclopentadiene	U		0.253	1.00
Di-isopropyl ether	U		0.105	1.00
Ethylbenzene	U		0.137	1.00
4-Ethyltoluene	U		0.208	1.00
Hexachloro-1,3-butadiene	U		0.337	1.00
n-Hexane	U		0.749	10.0
Isopropylbenzene	U		0.105	1.00
p-Isopropyltoluene	U		0.120	1.00
2-Butanone (MEK)	U		1.19	10.0
Methyl Cyclohexane	U		0.660	1.00
Methylene Chloride	U		0.430	5.00
4-Methyl-2-pentanone (MIBK)	U		0.478	10.0
Methyl tert-butyl ether	U		0.101	1.00
Naphthalene	U		1.00	5.00
Propene	U		0.936	2.50
n-Propylbenzene	U		0.0993	1.00
Styrene	U		0.118	1.00
1,1,1,2-Tetrachloroethane	U		0.147	1.00
1,1,2,2-Tetrachloroethane	U		0.133	1.00
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00
Tetrachloroethene	U		0.300	1.00
Toluene	U		0.278	1.00
1,2,3-Trichlorobenzene	U		0.230	1.00
1,2,4-Trichlorobenzene	U		0.481	1.00
1,1,1-Trichloroethane	U		0.149	1.00
1,1,2-Trichloroethane	U		0.158	1.00
Trichloroethene	U		0.190	1.00
Trichlorofluoromethane	U		0.160	5.00
1,2,3-Trichloropropane	U		0.237	2.50
1,2,4-Trimethylbenzene	U		0.322	1.00
1,2,3-Trimethylbenzene	U		0.104	1.00
1,3,5-Trimethylbenzene	U		0.104	1.00
Vinyl chloride	U		0.234	1.00
Xylenes, Total	U		0.174	3.00
(S) Toluene-d8	*107*			*80.0-120*
(S) 4-Bromofluorobenzene	*99.1*			*77.0-126*
(S) 1,2-Dichloroethane-d4	*85.6*			*70.0-130*

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R4035528-1 02/17/24 20:15 • (LCSD) R4035528-2 02/17/24 20:38

Analyte	Spike Amount ug/l	LCS Result ug/l	LCSD Result ug/l	LCS Rec. %	LCSD Rec. %	Rec. Limits %	LCS Qualifier	LCSD Qualifier	RPD %	RPD Limits %
Acetone	25.0	21.7	21.2	86.8	84.8	19.0-160			2.33	27
Acrolein	25.0	20.2	19.7	80.8	78.8	30.0-160			2.51	26
Acrylonitrile	25.0	20.0	19.2	80.0	76.8	55.0-149			4.08	20
Benzene	5.00	5.02	5.23	100	105	70.0-123			4.10	20
Bromobenzene	5.00	4.20	4.35	84.0	87.0	73.0-121			3.51	20
Bromodichloromethane	5.00	4.78	5.05	95.6	101	75.0-120			5.49	20
Bromoform	5.00	3.43	3.47	68.6	69.4	68.0-132			1.16	20
Bromomethane	5.00	5.87	7.37	117	147	30.0-160			22.7	25
1,3-Butadiene	5.00	4.84	5.10	96.8	102	45.0-147			5.23	20
n-Butylbenzene	5.00	3.73	3.98	74.6	79.6	73.0-125			6.49	20
sec-Butylbenzene	5.00	4.32	4.61	86.4	92.2	75.0-125			6.49	20
tert-Butylbenzene	5.00	4.21	4.42	84.2	88.4	76.0-124			4.87	20
Carbon tetrachloride	5.00	4.89	5.04	97.8	101	68.0-126			3.02	20
Carbon disulfide	5.00	4.72	4.87	94.4	97.4	61.0-128			3.13	20
Chlorobenzene	5.00	4.34	4.60	86.8	92.0	80.0-121			5.82	20
Chlorodibromomethane	5.00	4.09	4.03	81.8	80.6	77.0-125			1.48	20
Chloroethane	5.00	8.56	9.30	171	186	47.0-150	L1	L1	8.29	20
Chloroform	5.00	4.86	5.00	97.2	100	73.0-120			2.84	20
Chloromethane	5.00	4.29	4.68	85.8	93.6	41.0-142			8.70	20
Cyclohexane	5.00	4.80	5.39	96.0	108	71.0-124			11.6	20
2-Chlorotoluene	5.00	4.20	4.38	84.0	87.6	76.0-123			4.20	20
4-Chlorotoluene	5.00	4.09	4.34	81.8	86.8	75.0-122			5.93	20
1,2-Dibromo-3-Chloropropane	5.00	3.30	3.47	66.0	69.4	58.0-134			5.02	20
1,2-Dibromoethane	5.00	4.26	4.36	85.2	87.2	80.0-122			2.32	20
Dibromomethane	5.00	5.03	4.89	101	97.8	80.0-120			2.82	20
1,2-Dichlorobenzene	5.00	3.85	3.96	77.0	79.2	79.0-121	L2		2.82	20
1,3-Dichlorobenzene	5.00	4.06	4.30	81.2	86.0	79.0-120			5.74	20
1,4-Dichlorobenzene	5.00	3.99	4.09	79.8	81.8	79.0-120			2.48	20
Dichlorodifluoromethane	5.00	3.93	4.36	78.6	87.2	51.0-149			10.4	20
1,1-Dichloroethane	5.00	4.64	4.81	92.8	96.2	70.0-126			3.60	20
1,2-Dichloroethane	5.00	4.34	4.51	86.8	90.2	70.0-128			3.84	20
1,1-Dichloroethene	5.00	5.77	5.99	115	120	71.0-124			3.74	20
cis-1,2-Dichloroethene	5.00	4.98	5.00	99.6	100	73.0-120			0.401	20
trans-1,2-Dichloroethene	5.00	4.94	5.27	98.8	105	73.0-120			6.46	20
1,2-Dichloropropane	5.00	4.70	4.83	94.0	96.6	77.0-125			2.73	20
1,1-Dichloropropene	5.00	5.38	5.39	108	108	74.0-126			0.186	20
1,3-Dichloropropane	5.00	4.14	4.47	82.8	89.4	80.0-120			7.67	20
cis-1,3-Dichloropropene	5.00	4.58	4.60	91.6	92.0	80.0-123			0.436	20
trans-1,3-Dichloropropene	5.00	4.16	4.10	83.2	82.0	78.0-124			1.45	20
2,2-Dichloropropane	5.00	4.66	4.76	93.2	95.2	58.0-130			2.12	20

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R4035528-1 02/17/24 20:15 • (LCSD) R4035528-2 02/17/24 20:38

Analyte	Spike Amount ug/l	LCS Result ug/l	LCSD Result ug/l	LCS Rec. %	LCSD Rec. %	Rec. Limits %	LCS Qualifier	LCSD Qualifier	RPD %	RPD Limits %
Dicyclopentadiene	5.00	4.30	4.60	86.0	92.0	74.0-126			6.74	20
Di-isopropyl ether	5.00	3.98	4.04	79.6	80.8	58.0-138			1.50	20
Ethylbenzene	5.00	4.35	4.60	87.0	92.0	79.0-123			5.59	20
4-Ethyltoluene	5.00	4.22	4.51	84.4	90.2	74.0-127			6.64	20
Hexachloro-1,3-butadiene	5.00	3.56	3.75	71.2	75.0	54.0-138			5.20	20
n-Hexane	5.00	4.20	4.17	84.0	83.4	57.0-133			0.717	20
Isopropylbenzene	5.00	4.11	4.38	82.2	87.6	76.0-127			6.36	20
p-Isopropyltoluene	5.00	4.07	4.33	81.4	86.6	76.0-125			6.19	20
2-Butanone (MEK)	25.0	19.2	20.0	76.8	80.0	44.0-160			4.08	20
Methyl Cyclohexane	5.00	4.93	5.06	98.6	101	68.0-126			2.60	20
Methylene Chloride	5.00	5.19	5.36	104	107	67.0-120			3.22	20
4-Methyl-2-pentanone (MIBK)	25.0	17.4	17.6	69.6	70.4	68.0-142			1.14	20
Methyl tert-butyl ether	5.00	4.55	4.69	91.0	93.8	68.0-125			3.03	20
Naphthalene	5.00	2.93	2.67	58.6	53.4	54.0-135		L2	9.29	20
Propene	5.00	1.78	2.06	35.6	41.2	30.0-160			14.6	20
n-Propylbenzene	5.00	4.36	4.72	87.2	94.4	77.0-124			7.93	20
Styrene	5.00	3.70	3.87	74.0	77.4	73.0-130			4.49	20
1,1,1,2-Tetrachloroethane	5.00	4.11	4.33	82.2	86.6	75.0-125			5.21	20
1,1,2,2-Tetrachloroethane	5.00	3.75	3.79	75.0	75.8	65.0-130			1.06	20
1,1,2-Trichlorotrifluoroethane	5.00	4.54	4.76	90.8	95.2	69.0-132			4.73	20
Tetrachloroethene	5.00	4.67	5.01	93.4	100	72.0-132			7.02	20
Toluene	5.00	4.54	4.73	90.8	94.6	79.0-120			4.10	20
1,2,3-Trichlorobenzene	5.00	3.92	3.34	78.4	66.8	50.0-138			16.0	20
1,2,4-Trichlorobenzene	5.00	3.27	3.51	65.4	70.2	57.0-137			7.08	20
1,1,1-Trichloroethane	5.00	5.12	5.30	102	106	73.0-124			3.45	20
1,1,2-Trichloroethane	5.00	4.39	4.64	87.8	92.8	80.0-120			5.54	20
Trichloroethene	5.00	5.15	5.37	103	107	78.0-124			4.18	20
Trichlorofluoromethane	5.00	7.28	7.66	146	153	59.0-147		L1	5.09	20
1,2,3-Trichloropropane	5.00	4.30	4.34	86.0	86.8	73.0-130			0.926	20
1,2,4-Trimethylbenzene	5.00	3.99	4.24	79.8	84.8	76.0-121			6.08	20
1,2,3-Trimethylbenzene	5.00	3.96	4.18	79.2	83.6	77.0-120			5.41	20
1,3,5-Trimethylbenzene	5.00	4.14	4.31	82.8	86.2	76.0-122			4.02	20
Vinyl chloride	5.00	5.79	5.84	116	117	67.0-131			0.860	20
Xylenes, Total	15.0	12.6	13.2	84.0	88.0	79.0-123			4.65	20
(S) Toluene-d8				*106*	*106*	*80.0-120*				
(S) 4-Bromofluorobenzene				*101*	*100*	*77.0-126*				
(S) 1,2-Dichloroethane-d4				*86.1*	*84.2*	*70.0-130*				

L1705404-01 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1705404-01 02/18/24 00:33 • (MS) R4035528-4 02/18/24 06:21 • (MSD) R4035528-5 02/18/24 06:44

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MSD Result ug/l	MS Rec. %	MSD Rec. %	Dilution	Rec. Limits %	MS Qualifier	MSD Qualifier	RPD %	RPD Limits %
Acetone	25.0	U	18.8	19.4	75.2	77.6	1	10.0-160			3.14	35
Acrolein	25.0	U	31.1	32.4	124	130	1	10.0-160			4.09	39
Acrylonitrile	25.0	U	26.1	20.4	104	81.6	1	21.0-160			24.5	32
Benzene	5.00	U	7.46	5.63	149	113	1	17.0-158		R5	28.0	27
Bromobenzene	5.00	U	5.43	4.97	109	99.4	1	30.0-149			8.85	28
Bromodichloromethane	5.00	U	5.10	5.48	102	110	1	31.0-150			7.18	27
Bromoform	5.00	U	3.48	3.37	69.6	67.4	1	29.0-150			3.21	29
Bromomethane	5.00	U	2.70	4.09	54.0	81.8	1	10.0-160		R5	40.9	38
1,3-Butadiene	5.00	U	6.00	6.90	120	138	1	10.0-160			14.0	22
n-Butylbenzene	5.00	U	4.62	4.59	92.4	91.8	1	31.0-150			0.651	30
sec-Butylbenzene	5.00	U	4.70	5.14	94.0	103	1	33.0-155			8.94	29
tert-Butylbenzene	5.00	U	5.32	5.41	106	108	1	34.0-153			1.68	28
Carbon tetrachloride	5.00	U	5.76	5.80	115	116	1	23.0-159			0.692	28
Carbon disulfide	5.00	U	4.98	5.18	99.6	104	1	10.0-156			3.94	28
Chlorobenzene	5.00	U	4.59	4.51	91.8	90.2	1	33.0-152			1.76	27
Chlorodibromomethane	5.00	U	5.32	3.86	106	77.2	1	37.0-149		R5	31.8	27
Chloroethane	5.00	U	6.91	7.81	138	156	1	10.0-160			12.2	30
Chloroform	5.00	3.49	9.21	8.96	184	179	1	29.0-154	M1	M1	2.75	28
Chloromethane	5.00	U	4.86	5.00	97.2	100	1	10.0-160			2.84	29
Cyclohexane	5.00	U	6.40	6.33	128	127	1	19.0-160			1.10	23
2-Chlorotoluene	5.00	U	4.57	5.42	91.4	108	1	32.0-153			17.0	28
4-Chlorotoluene	5.00	U	5.68	4.88	114	97.6	1	32.0-150			15.2	28
1,2-Dibromo-3-Chloropropane	5.00	U	4.03	3.86	80.6	77.2	1	22.0-151			4.31	34
1,2-Dibromoethane	5.00	U	4.84	4.04	96.8	80.8	1	34.0-147			18.0	27
Dibromomethane	5.00	U	5.79	5.32	116	106	1	30.0-151			8.46	27
1,2-Dichlorobenzene	5.00	U	3.94	4.16	78.8	83.2	1	34.0-149			5.43	28
1,3-Dichlorobenzene	5.00	U	4.08	4.38	81.6	87.6	1	36.0-146			7.09	27
1,4-Dichlorobenzene	5.00	U	3.97	4.26	79.4	85.2	1	35.0-142			7.05	27
Dichlorodifluoromethane	5.00	U	5.12	5.66	102	113	1	10.0-160			10.0	29
1,1-Dichloroethane	5.00	U	6.70	5.31	134	106	1	25.0-158			23.1	27
1,2-Dichloroethane	5.00	U	5.42	5.12	108	102	1	29.0-151			5.69	27
1,1-Dichloroethene	5.00	U	6.85	6.82	137	136	1	11.0-160			0.439	29
cis-1,2-Dichloroethene	5.00	U	5.35	5.35	107	107	1	10.0-160			0.000	27
trans-1,2-Dichloroethene	5.00	U	6.02	5.70	120	114	1	17.0-153			5.46	27
1,2-Dichloropropane	5.00	U	5.66	5.23	113	105	1	30.0-156			7.90	27
1,1-Dichloropropene	5.00	U	6.95	6.14	139	123	1	25.0-158			12.4	27
1,3-Dichloropropane	5.00	U	5.80	4.23	116	84.6	1	38.0-147		R5	31.3	27
cis-1,3-Dichloropropene	5.00	U	4.51	4.93	90.2	98.6	1	34.0-149			8.90	28
trans-1,3-Dichloropropene	5.00	U	4.30	4.30	86.0	86.0	1	32.0-149			0.000	28
2,2-Dichloropropane	5.00	U	5.88	5.71	118	114	1	24.0-152			2.93	29

L1705404-01 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1705404-01 02/18/24 00:33 • (MS) R4035528-4 02/18/24 06:21 • (MSD) R4035528-5 02/18/24 06:44

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MSD Result ug/l	MS Rec. %	MSD Rec. %	Dilution	Rec. Limits %	MS Qualifier	MSD Qualifier	RPD %	RPD Limits %
Dicyclopentadiene	5.00	U	3.18	4.45	63.6	89.0	1	51.0-139		R5	33.3	20
Di-isopropyl ether	5.00	U	4.76	4.54	95.2	90.8	1	21.0-160			4.73	28
Ethylbenzene	5.00	U	4.67	4.64	93.4	92.8	1	30.0-155			0.644	27
4-Ethyltoluene	5.00	U	4.70	5.11	94.0	102	1	10.0-160			8.36	20
Hexachloro-1,3-butadiene	5.00	U	5.33	4.13	107	82.6	1	20.0-154			25.4	34
n-Hexane	5.00	U	6.04	5.15	121	103	1	10.0-153			15.9	28
Isopropylbenzene	5.00	U	4.82	4.59	96.4	91.8	1	28.0-157			4.89	27
p-Isopropyltoluene	5.00	U	4.53	4.81	90.6	96.2	1	30.0-154			6.00	29
2-Butanone (MEK)	25.0	U	20.1	18.9	80.4	75.6	1	10.0-160			6.15	32
Methyl Cyclohexane	5.00	U	6.25	6.50	125	130	1	11.0-160			3.92	24
Methylene Chloride	5.00	U	5.70	5.57	114	111	1	23.0-144			2.31	28
4-Methyl-2-pentanone (MIBK)	25.0	U	19.6	18.1	78.4	72.4	1	29.0-160			7.96	29
Methyl tert-butyl ether	5.00	U	5.62	5.05	112	101	1	28.0-150			10.7	29
Naphthalene	5.00	U	4.74	4.00	94.8	80.0	1	12.0-156			16.9	35
Propene	5.00	U	2.87	3.19	57.4	63.8	1	10.0-160			10.6	29
n-Propylbenzene	5.00	U	5.62	5.30	112	106	1	31.0-154			5.86	28
Styrene	5.00	U	3.89	4.40	77.8	88.0	1	33.0-155			12.3	28
1,1,1,2-Tetrachloroethane	5.00	U	4.39	4.38	87.8	87.6	1	36.0-151			0.228	29
1,1,2,2-Tetrachloroethane	5.00	U	5.47	5.00	109	100	1	33.0-150			8.98	28
1,1,2-Trichlorotrifluoroethane	5.00	U	6.13	6.06	123	121	1	23.0-160			1.15	30
Tetrachloroethene	5.00	U	5.35	4.86	107	97.2	1	10.0-160			9.60	27
Toluene	5.00	U	6.82	4.55	136	91.0	1	26.0-154		R5	39.9	28
1,2,3-Trichlorobenzene	5.00	U	12.1	4.73	242	94.6	1	17.0-150	M1	R5	87.6	36
1,2,4-Trichlorobenzene	5.00	U	5.05	4.34	101	86.8	1	24.0-150			15.1	33
1,1,1-Trichloroethane	5.00	U	6.25	5.82	125	116	1	23.0-160			7.13	28
1,1,2-Trichloroethane	5.00	U	5.06	4.46	101	89.2	1	35.0-147			12.6	27
Trichloroethene	5.00	U	5.14	5.05	103	101	1	10.0-160			1.77	25
Trichlorofluoromethane	5.00	U	7.23	8.46	145	169	1	17.0-160		M1	15.7	31
1,2,3-Trichloropropane	5.00	U	5.07	5.02	101	100	1	34.0-151			0.991	29
1,2,4-Trimethylbenzene	5.00	U	4.38	4.76	87.6	95.2	1	26.0-154			8.32	27
1,2,3-Trimethylbenzene	5.00	U	4.25	4.60	85.0	92.0	1	32.0-149			7.91	28
1,3,5-Trimethylbenzene	5.00	U	4.54	5.35	90.8	107	1	28.0-153			16.4	27
Vinyl chloride	5.00	U	6.36	7.11	127	142	1	10.0-160			11.1	27
Xylenes, Total	15.0	U	13.6	14.3	90.7	95.3	1	29.0-154			5.02	28
(S) Toluene-d8					*129*	*95.7*		*80.0-120*	S10			
(S) 4-Bromofluorobenzene					*116*	*95.4*		*77.0-126*				
(S) 1,2-Dichloroethane-d4					*113*	*94.1*		*70.0-130*				

Method Blank (MB)

(MB) R4035816-3 02/19/24 19:40

Analyte	MB Result ug/l	MB Qualifier	MB MDL ug/l	MB RDL ug/l
Acetone	U		11.3	50.0
Acrolein	U		2.54	50.0
Acrylonitrile	U		0.671	10.0
Benzene	U		0.0941	1.00
Bromobenzene	U		0.118	1.00
Bromodichloromethane	U		0.136	1.00
Bromoform	U		0.129	1.00
Bromomethane	U		0.605	5.00
1,3-Butadiene	U		0.299	2.00
n-Butylbenzene	U		0.157	1.00
sec-Butylbenzene	U		0.125	1.00
tert-Butylbenzene	U		0.127	1.00
Carbon tetrachloride	U		0.128	1.00
Carbon disulfide	U		0.0962	1.00
Chlorobenzene	U		0.116	1.00
Chlorodibromomethane	U		0.140	1.00
Chloroethane	U		0.192	5.00
Chloroform	U		0.111	5.00
Chloromethane	U		0.960	2.50
Cyclohexane	U		0.188	1.00
2-Chlorotoluene	U		0.106	1.00
4-Chlorotoluene	U		0.114	1.00
1,2-Dibromo-3-Chloropropane	U		0.276	5.00
1,2-Dibromoethane	U		0.126	1.00
Dibromomethane	U		0.122	1.00
1,2-Dichlorobenzene	U		0.107	1.00
1,3-Dichlorobenzene	U		0.110	1.00
1,4-Dichlorobenzene	U		0.120	1.00
Dichlorodifluoromethane	U		0.374	5.00
1,1-Dichloroethane	U		0.100	1.00
1,2-Dichloroethane	U		0.0819	1.00
1,1-Dichloroethene	U		0.188	1.00
cis-1,2-Dichloroethene	U		0.126	1.00
trans-1,2-Dichloroethene	U		0.149	1.00
1,2-Dichloropropane	U		0.149	1.00
1,1-Dichloropropene	U		0.142	1.00
1,3-Dichloropropane	U		0.110	1.00
cis-1,3-Dichloropropene	U		0.111	1.00
trans-1,3-Dichloropropene	U		0.118	1.00
2,2-Dichloropropane	U		0.161	1.00

Method Blank (MB)

(MB) R4035816-3 02/19/24 19:40

Analyte	MB Result ug/l	MB Qualifier	MB MDL ug/l	MB RDL ug/l
Dicyclopentadiene	U		0.253	1.00
Di-isopropyl ether	U		0.105	1.00
Ethylbenzene	U		0.137	1.00
4-Ethyltoluene	U		0.208	1.00
Hexachloro-1,3-butadiene	U		0.337	1.00
n-Hexane	U		0.749	10.0
Isopropylbenzene	U		0.105	1.00
p-Isopropyltoluene	U		0.120	1.00
2-Butanone (MEK)	U		1.19	10.0
Methyl Cyclohexane	U		0.660	1.00
Methylene Chloride	U		0.430	5.00
4-Methyl-2-pentanone (MIBK)	U		0.478	10.0
Methyl tert-butyl ether	U		0.101	1.00
Naphthalene	U		1.00	5.00
Propene	U		0.936	2.50
n-Propylbenzene	U		0.0993	1.00
Styrene	U		0.118	1.00
1,1,1,2-Tetrachloroethane	U		0.147	1.00
1,1,2,2-Tetrachloroethane	U		0.133	1.00
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00
Tetrachloroethene	U		0.300	1.00
Toluene	U		0.278	1.00
1,2,3-Trichlorobenzene	U		0.230	1.00
1,2,4-Trichlorobenzene	U		0.481	1.00
1,1,1-Trichloroethane	U		0.149	1.00
1,1,2-Trichloroethane	U		0.158	1.00
Trichloroethene	U		0.190	1.00
Trichlorofluoromethane	U		0.160	5.00
1,2,3-Trichloropropane	U		0.237	2.50
1,2,4-Trimethylbenzene	U		0.322	1.00
1,2,3-Trimethylbenzene	U		0.104	1.00
1,3,5-Trimethylbenzene	U		0.104	1.00
Vinyl chloride	U		0.234	1.00
Xylenes, Total	U		0.174	3.00
(S) Toluene-d8	*106*			*80.0-120*
(S) 4-Bromofluorobenzene	*99.5*			*77.0-126*
(S) 1,2-Dichloroethane-d4	*79.7*			*70.0-130*

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R4035816-1 02/19/24 18:31 • (LCSD) R4035816-2 02/19/24 18:54

Analyte	Spike Amount ug/l	LCS Result ug/l	LCSD Result ug/l	LCS Rec. %	LCSD Rec. %	Rec. Limits %	LCS Qualifier	LCSD Qualifier	RPD %	RPD Limits %
Acetone	25.0	19.7	19.4	78.8	77.6	19.0-160			1.53	27
Acrolein	25.0	22.9	26.6	91.6	106	30.0-160			14.9	26
Acrylonitrile	25.0	18.0	19.3	72.0	77.2	55.0-149			6.97	20
Benzene	5.00	5.02	5.19	100	104	70.0-123			3.33	20
Bromobenzene	5.00	4.99	4.54	99.8	90.8	73.0-121			9.44	20
Bromodichloromethane	5.00	4.76	5.18	95.2	104	75.0-120			8.45	20
Bromoform	5.00	3.28	3.47	65.6	69.4	68.0-132	L2		5.63	20
Bromomethane	5.00	6.95	7.14	139	143	30.0-160			2.70	25
1,3-Butadiene	5.00	3.98	3.82	79.6	76.4	45.0-147			4.10	20
n-Butylbenzene	5.00	4.05	4.34	81.0	86.8	73.0-125			6.91	20
sec-Butylbenzene	5.00	4.33	4.58	86.6	91.6	75.0-125			5.61	20
tert-Butylbenzene	5.00	4.19	4.29	83.8	85.8	76.0-124			2.36	20
Carbon tetrachloride	5.00	4.69	4.90	93.8	98.0	68.0-126			4.38	20
Carbon disulfide	5.00	4.04	4.15	80.8	83.0	61.0-128			2.69	20
Chlorobenzene	5.00	4.32	4.55	86.4	91.0	80.0-121			5.19	20
Chlorodibromomethane	5.00	4.08	4.10	81.6	82.0	77.0-125			0.489	20
Chloroethane	5.00	6.66	6.72	133	134	47.0-150			0.897	20
Chloroform	5.00	4.93	5.29	98.6	106	73.0-120			7.05	20
Chloromethane	5.00	3.18	3.35	63.6	67.0	41.0-142			5.21	20
Cyclohexane	5.00	4.94	4.84	98.8	96.8	71.0-124			2.04	20
2-Chlorotoluene	5.00	4.25	4.46	85.0	89.2	76.0-123			4.82	20
4-Chlorotoluene	5.00	4.21	4.31	84.2	86.2	75.0-122			2.35	20
1,2-Dibromo-3-Chloropropane	5.00	3.11	3.59	62.2	71.8	58.0-134			14.3	20
1,2-Dibromoethane	5.00	4.06	4.27	81.2	85.4	80.0-122			5.04	20
Dibromomethane	5.00	4.83	4.86	96.6	97.2	80.0-120			0.619	20
1,2-Dichlorobenzene	5.00	3.72	4.04	74.4	80.8	79.0-121	L2		8.25	20
1,3-Dichlorobenzene	5.00	3.92	4.23	78.4	84.6	79.0-120	L2		7.61	20
1,4-Dichlorobenzene	5.00	3.91	4.24	78.2	84.8	79.0-120	L2		8.10	20
Dichlorodifluoromethane	5.00	3.88	3.99	77.6	79.8	51.0-149			2.80	20
1,1-Dichloroethane	5.00	4.73	4.97	94.6	99.4	70.0-126			4.95	20
1,2-Dichloroethane	5.00	4.57	4.61	91.4	92.2	70.0-128			0.871	20
1,1-Dichloroethene	5.00	5.46	5.92	109	118	71.0-124			8.08	20
cis-1,2-Dichloroethene	5.00	4.82	4.94	96.4	98.8	73.0-120			2.46	20
trans-1,2-Dichloroethene	5.00	4.67	5.28	93.4	106	73.0-120			12.3	20
1,2-Dichloropropane	5.00	4.70	4.84	94.0	96.8	77.0-125			2.94	20
1,1-Dichloropropene	5.00	4.92	5.21	98.4	104	74.0-126			5.73	20
1,3-Dichloropropane	5.00	4.25	4.77	85.0	95.4	80.0-120			11.5	20
cis-1,3-Dichloropropene	5.00	4.59	4.71	91.8	94.2	80.0-123			2.58	20
trans-1,3-Dichloropropene	5.00	4.14	4.37	82.8	87.4	78.0-124			5.41	20
2,2-Dichloropropane	5.00	5.56	5.76	111	115	58.0-130			3.53	20

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R4035816-1 02/19/24 18:31 • (LCSD) R4035816-2 02/19/24 18:54

Analyte	Spike Amount ug/l	LCS Result ug/l	LCSD Result ug/l	LCS Rec. %	LCSD Rec. %	Rec. Limits %	LCS Qualifier	LCSD Qualifier	RPD %	RPD Limits %
Dicyclopentadiene	5.00	4.81	4.45	96.2	89.0	74.0-126			7.78	20
Di-isopropyl ether	5.00	4.18	4.34	83.6	86.8	58.0-138			3.76	20
Ethylbenzene	5.00	4.30	4.52	86.0	90.4	79.0-123			4.99	20
4-Ethyltoluene	5.00	4.57	4.54	91.4	90.8	74.0-127			0.659	20
Hexachloro-1,3-butadiene	5.00	3.82	4.36	76.4	87.2	54.0-138			13.2	20
n-Hexane	5.00	3.79	3.93	75.8	78.6	57.0-133			3.63	20
Isopropylbenzene	5.00	4.18	4.41	83.6	88.2	76.0-127			5.36	20
p-Isopropyltoluene	5.00	4.62	4.59	92.4	91.8	76.0-125			0.651	20
2-Butanone (MEK)	25.0	17.5	18.2	70.0	72.8	44.0-160			3.92	20
Methyl Cyclohexane	5.00	4.83	4.74	96.6	94.8	68.0-126			1.88	20
Methylene Chloride	5.00	5.39	5.64	108	113	67.0-120			4.53	20
4-Methyl-2-pentanone (MIBK)	25.0	17.3	18.0	69.2	72.0	68.0-142			3.97	20
Methyl tert-butyl ether	5.00	4.54	4.72	90.8	94.4	68.0-125			3.89	20
Naphthalene	5.00	2.97	3.14	59.4	62.8	54.0-135			5.56	20
Propene	5.00	1.90	1.71	38.0	34.2	30.0-160			10.5	20
n-Propylbenzene	5.00	5.11	5.01	102	100	77.0-124			1.98	20
Styrene	5.00	3.64	4.03	72.8	80.6	73.0-130	L2		10.2	20
1,1,1,2-Tetrachloroethane	5.00	4.16	4.49	83.2	89.8	75.0-125			7.63	20
1,1,2,2-Tetrachloroethane	5.00	4.65	4.37	93.0	87.4	65.0-130			6.21	20
1,1,2-Trichlorotrifluoroethane	5.00	4.57	4.88	91.4	97.6	69.0-132			6.56	20
Tetrachloroethene	5.00	4.59	4.74	91.8	94.8	72.0-132			3.22	20
Toluene	5.00	4.43	4.74	88.6	94.8	79.0-120			6.76	20
1,2,3-Trichlorobenzene	5.00	3.32	3.66	66.4	73.2	50.0-138			9.74	20
1,2,4-Trichlorobenzene	5.00	3.52	3.98	70.4	79.6	57.0-137			12.3	20
1,1,1-Trichloroethane	5.00	4.95	5.27	99.0	105	73.0-124			6.26	20
1,1,2-Trichloroethane	5.00	4.27	4.35	85.4	87.0	80.0-120			1.86	20
Trichloroethene	5.00	4.56	4.84	91.2	96.8	78.0-124			5.96	20
Trichlorofluoromethane	5.00	5.76	5.54	115	111	59.0-147			3.89	20
1,2,3-Trichloropropane	5.00	4.09	4.28	81.8	85.6	73.0-130			4.54	20
1,2,4-Trimethylbenzene	5.00	3.94	4.34	78.8	86.8	76.0-121			9.66	20
1,2,3-Trimethylbenzene	5.00	3.98	4.28	79.6	85.6	77.0-120			7.26	20
1,3,5-Trimethylbenzene	5.00	4.26	4.38	85.2	87.6	76.0-122			2.78	20
Vinyl chloride	5.00	4.62	4.49	92.4	89.8	67.0-131			2.85	20
Xylenes, Total	15.0	12.4	13.5	82.7	90.0	79.0-123			8.49	20
(S) Toluene-d8				*106*	*106*	*80.0-120*				
(S) 4-Bromofluorobenzene				*123*	*100*	*77.0-126*				
(S) 1,2-Dichloroethane-d4				*85.1*	*86.7*	*70.0-130*				

Method Blank (MB)

(MB) R4035347-3 02/15/24 22:44

Analyte	MB Result	MB Qualifier	MB MDL	MB RDL
	ug/l		ug/l	ug/l
1,4-Dioxane	U		0.597	3.00
(S) Toluene-d8	*92.9*			*77.0-127*

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R4035347-1 02/15/24 21:38 • (LCSD) R4035347-2 02/15/24 22:00

Analyte	Spike Amount	LCS Result	LCSD Result	LCS Rec.	LCSD Rec.	Rec. Limits	LCS Qualifier	LCSD Qualifier	RPD	RPD Limits
	ug/l	ug/l	ug/l	%	%	%			%	%
1,4-Dioxane	50.0	46.7	46.5	93.4	93.0	55.0-138			0.429	24
(S) Toluene-d8				*92.3*	*92.9*	*77.0-127*				

L1705409-01 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1705409-01 02/16/24 02:07 • (MS) R4035347-4 02/16/24 07:10 • (MSD) R4035347-5 02/16/24 07:32

Analyte	Spike Amount	Original Result	MS Result	MSD Result	MS Rec.	MSD Rec.	Dilution	Rec. Limits	MS Qualifier	MSD Qualifier	RPD	RPD Limits
	ug/l	ug/l	ug/l	ug/l	%	%		%			%	%
1,4-Dioxane	50.0	34.8	68.3	65.1	67.0	60.6	1	13.0-160			4.80	31
(S) Toluene-d8					*96.6*	*97.4*		*77.0-127*				

Method Blank (MB)

(MB) R4036334-3 02/20/24 11:58

Analyte	MB Result	MB Qualifier	MB MDL	MB RDL
	ug/l		ug/l	ug/l
1,4-Dioxane	U		0.597	3.00
(S) Toluene-d8	*94.9*			*77.0-127*

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R4036334-1 02/20/24 10:53 • (LCSD) R4036334-2 02/20/24 11:15

Analyte	Spike Amount	LCS Result	LCSD Result	LCS Rec.	LCSD Rec.	Rec. Limits	LCS Qualifier	LCSD Qualifier	RPD	RPD Limits
	ug/l	ug/l	ug/l	%	%	%			%	%
1,4-Dioxane	50.0	43.2	44.8	86.4	89.6	55.0-138			3.64	24
(S) Toluene-d8				*95.2*	*94.9*	*77.0-127*				

L1705776-08 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1705776-08 02/20/24 16:25 • (MS) R4036334-4 02/20/24 23:20 • (MSD) R4036334-5 02/20/24 23:42

Analyte	Spike Amount	Original Result	MS Result	MSD Result	MS Rec.	MSD Rec.	Dilution	Rec. Limits	MS Qualifier	MSD Qualifier	RPD	RPD Limits
	ug/l	ug/l	ug/l	ug/l	%	%		%			%	%
1,4-Dioxane	50.0	U	43.6	52.2	87.2	104	1	13.0-160			18.0	31
(S) Toluene-d8					*95.4*	*97.9*		*77.0-127*				

L1705776-11 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1705776-11 02/20/24 17:31 • (MS) R4036334-6 02/21/24 00:04 • (MSD) R4036334-7 02/21/24 00:26

Analyte	Spike Amount	Original Result	MS Result	MSD Result	MS Rec.	MSD Rec.	Dilution	Rec. Limits	MS Qualifier	MSD Qualifier	RPD	RPD Limits
	ug/l	ug/l	ug/l	ug/l	%	%		%			%	%
1,4-Dioxane	50.0	31.4	81.6	78.5	100	94.2	1	13.0-160			3.87	31
(S) Toluene-d8					*98.6*	*98.3*		*77.0-127*				

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

Method Blank (MB)

(MB) R4036886-3 02/21/24 13:02

Analyte	MB Result ug/l	MB Qualifier	MB MDL ug/l	MB RDL ug/l
1,4-Dioxane	U		0.597	3.00
(S) Toluene-d8	*95.4*			*77.0-127*

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R4036886-1 02/21/24 11:57 • (LCSD) R4036886-2 02/21/24 12:19

Analyte	Spike Amount ug/l	LCS Result ug/l	LCSD Result ug/l	LCS Rec. %	LCSD Rec. %	Rec. Limits %	LCS Qualifier	LCSD Qualifier	RPD %	RPD Limits %
1,4-Dioxane	50.0	53.2	52.6	106	105	55.0-138			1.13	24
(S) Toluene-d8				*95.3*	*95.4*	*77.0-127*				

Method Blank (MB)

(MB) R4037174-3 02/22/24 13:02

Analyte	MB Result	MB Qualifier	MB MDL	MB RDL
	ug/l		ug/l	ug/l
1,4-Dioxane	U		0.597	3.00
(S) Toluene-d8	*93.8*			*77.0-127*

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R4037174-1 02/22/24 11:40 • (LCSD) R4037174-2 02/22/24 12:02

Analyte	Spike Amount	LCS Result	LCSD Result	LCS Rec.	LCSD Rec.	Rec. Limits	LCS Qualifier	LCSD Qualifier	RPD	RPD Limits
	ug/l	ug/l	ug/l	%	%	%			%	%
1,4-Dioxane	50.0	43.1	44.7	86.2	89.4	55.0-138			3.64	24
(S) Toluene-d8				*93.8*	*93.9*	*77.0-127*				

L1707469-01 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1707469-01 02/22/24 13:24 • (MS) R4037174-4 02/22/24 14:07 • (MSD) R4037174-5 02/22/24 14:29

Analyte	Spike Amount	Original Result	MS Result	MSD Result	MS Rec.	MSD Rec.	Dilution	Rec. Limits	MS Qualifier	MSD Qualifier	RPD	RPD Limits
	ug/l	ug/l	ug/l	ug/l	%	%		%			%	%
1,4-Dioxane	50.0	U	48.8	50.7	97.6	101	1	13.0-160			3.82	31
(S) Toluene-d8					*93.5*	*93.8*		*77.0-127*				

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

INTERNAL STANDARD SUMMARY

Instrument: VOCMS20 • File ID: 0217_02

02/17/24 07:22

Sample ID	File ID	8260-FLUOROBENZENE	8260-CHLOROBENZENE-D5	8260-1,4-DICHLOROBENZENE-D4
		Response	Response	Response
Standard	0217_02	178985	82056	73389
Upper Limit		357970	164112	146778
Lower Limit		89493	41028	36695
LCS R4035581-1 WG2228368 1x	0217_02LCS	178985	82056	73389
LCSD R4035581-2 WG2228368 1x	0217_03	189400	83365	74691
BLANK R4035581-3 WG2228368 1x	0217_05	169332	72597	61265
L1705776-01 WG2228368 1x	0217_06	149886	63248	55448
L1705776-03 WG2228368 1x	0217_07	169600	72527	60412
L1705776-04 WG2228368 1x	0217_08	165955	70111	60000
L1705776-05 WG2228368 1x	0217_09	159842	69451	59424
L1705776-06 WG2228368 1x	0217_10	164016	68367	57134
L1705776-07 WG2228368 1x	0217_11	166442	71494	59286
L1705776-08 WG2228368 1x	0217_12	167705	73223	60160
L1705776-11 WG2228368 1x	0217_13	155025	68709	56707
L1705776-13 WG2228368 1x	0217_14	200108	86372	86945
L1705776-15 WG2228368 1x	0217_15	161913	78879	58858
L1705776-18 WG2228368 1x	0217_16	165878	71000	60023
L1705776-02 WG2228368 10x	0217_17	163942	70856	57154
L1705776-10 WG2228368 10x	0217_19	165497	69827	61667
L1705776-12 WG2228368 10x	0217_20	165379	71930	61037
L1705776-16 WG2228368 5x	0217_22	164066	72914	69665
L1705776-17 WG2228368 10x	0217_23	167644	68885	59335
L1705776-19 WG2228368 10x	0217_24	159083	67988	58704
L1705776-20 WG2228368 20x	0217_25	155534	67176	55885
MS R4035581-4 WG2228368 1x	0217_26	171094	78521	67647
MSD R4035581-5 WG2228368 1x	0217_27	171698	79067	67358
MS R4035581-6 WG2228368 1x	0217_28	172607	77609	71066
MSD R4035581-7 WG2228368 1x	0217_29	178632	79432	69181

INTERNAL STANDARD SUMMARY

Instrument: VOCMS30 • File ID: 0217_27

02/17/24 20:15

Sample ID	File ID	8260-FLUOROBENZENE	8260-CHLOROBENZENE-D5	8260-1,4-DICHLOROBENZENE-D4
		Response	Response	Response
Standard	0217_27	241910	125515	124898
Upper Limit		483820	251030	249796
Lower Limit		120955	62758	62449
LCS R4035528-1 WG2228641 1x	0217_27LCS	241910	125515	124898
LCSD R4035528-2 WG2228641 1x	0217_28	241632	123524	121913
BLANK R4035528-3 WG2228641 1x	0217_30	218894	111895	109016
L1705776-25 WG2228641 1x	0217_33	214964	109465	106702
L1705776-23 WG2228641 1x	0217_40	207761	105528	103053
L1705776-24 WG2228641 1x	0217_41	210866	105878	103752
L1705776-21 WG2228641 250x	0217_47	243158	125378	124335
L1705776-22 WG2228641 25x	0217_48	226908	115206	114749
MS R4035528-4 WG2228641 1x	0217_51	293553	134798	143816
MSD R4035528-5 WG2228641 1x	0217_52	262487	146770	130529

Instrument: VOCMS30 • File ID: 0219_14

02/19/24 18:31

Sample ID	File ID	8260-FLUOROBENZENE	8260-CHLOROBENZENE-D5	8260-1,4-DICHLOROBENZENE-D4
		Response	Response	Response
Standard	0219_14	252425	128569	128185
Upper Limit		504850	257138	256370
Lower Limit		126213	64285	64093
LCS R4035816-1 WG2229520 1x	0219_14LCSB	252425	128569	128185
LCSD R4035816-2 WG2229520 1x	0219_15B	254832	129414	126380
BLANK R4035816-3 WG2229520 1x	0219_17B	252006	120609	117134
L1705776-02 WG2229520 10x	0219_30	237841	143495	121905
L1705776-09 WG2229520 10x	0219_31	228466	125685	135152
L1705776-10 WG2229520 10x	0219_32	215773	121945	116025
L1705776-12 WG2229520 10x	0219_33	226221	122632	124867
L1705776-13 WG2229520 2500x	0219_34	215854	114014	115757
L1705776-14 WG2229520 20x	0219_35	244033	130363	125757
L1705776-15 WG2229520 1x	0219_36	227649	114226	115815
L1705776-16 WG2229520 5x	0219_37	246567	130636	126107
L1705776-17 WG2229520 10x	0219_38	221003	116042	116726
L1705776-18 WG2229520 1x	0219_39	253943	134458	124858
L1705776-19 WG2229520 10x	0219_40	218266	114755	113614

INTERNAL STANDARD SUMMARY

Instrument: VOCMS30 • File ID: 0219_14

02/19/24 18:31

Sample ID	File ID	8260-FLUOROBENZENE Response	8260-CHLOROBENZENE-D5 Response	8260-1,4-DICHLOROBENZENE-D4 Response
L1705776-20 WG2229520 20x	0219_41	214621	114978	113634

INTERNAL STANDARD SUMMARY

Instrument: VOCMS27 • File ID: 0215A_03

02/15/24 20:57

Sample ID	File ID	8260-FLUOROBENZENE Response
Standard	0215A_03	329629
Upper Limit		659258
Lower Limit		164815
LCS R4035347-1 WG2227341 1x	0215A_04	310737
LCSD R4035347-2 WG2227341 1x	0215A_05	315477
BLANK R4035347-3 WG2227341 1x	0215A_07	297814
L1705776-01 WG2227341 1x	0215A_21	316092
L1705776-02 WG2227341 1x	0215A_22	332011
L1705776-03 WG2227341 1x	0215A_23	328428
L1705776-04 WG2227341 1x	0215A_24	328059
L1705776-05 WG2227341 1x	0215A_25	334712
L1705776-06 WG2227341 1x	0215A_26	329664
L1705776-07 WG2227341 1x	0215A_27	332564
MS R4035347-4 WG2227341 1x	0215A_28	293066
MSD R4035347-5 WG2227341 1x	0215A_29	313329

Instrument: VOCMS27 • File ID: 0220_03

02/20/24 10:32

Sample ID	File ID	8260-FLUOROBENZENE Response
Standard	0220_03	344527
Upper Limit		689054
Lower Limit		172264
LCS R4036334-1 WG2230157 1x	0220_04	320440
LCSD R4036334-2 WG2230157 1x	0220_05	314803
BLANK R4036334-3 WG2230157 1x	0220_07	292457
L1705776-25 WG2230157 1x	0220_08	350811
L1705776-08 WG2230157 1x	0220_09	330385
L1705776-09 WG2230157 1x	0220_10	305544
L1705776-10 WG2230157 1x	0220_11	327147
L1705776-11 WG2230157 1x	0220_12	323713
L1705776-12 WG2230157 1x	0220_13	330858
L1705776-17 WG2230157 1x	0220_16	343497
L1705776-18 WG2230157 1x	0220_17	322569
L1705776-22 WG2230157 1x	0220_18	326896

INTERNAL STANDARD SUMMARY

Instrument: VOCMS27 • File ID: 0220_03

02/20/24 10:32

Sample ID	File ID	8260-FLUOROBENZENE Response
L1705776-23 WG2230157 1x	0220_19	308527
L1705776-24 WG2230157 1x	0220_20	318609
L1705776-16 WG2230157 5x	0220_24	310905
L1705776-19 WG2230157 10x	0220_25	318869
MS R4036334-4 WG2230157 1x	0220_28	321121
MSD R4036334-5 WG2230157 1x	0220_29	300584
MS R4036334-6 WG2230157 1x	0220_30	307253
MSD R4036334-7 WG2230157 1x	0220_31	314153

Instrument: VOCMS27 • File ID: 0221_03

02/21/24 11:35

Sample ID	File ID	8260-FLUOROBENZENE Response
Standard	0221_03	303435
Upper Limit		606870
Lower Limit		151718
LCS R4036886-1 WG2230690 1x	0221_04	295880
LCSD R4036886-2 WG2230690 1x	0221_05	287985
BLANK R4036886-3 WG2230690 1x	0221_07	280046
L1705776-14 WG2230690 1x	0221_11	310135
L1705776-15 WG2230690 1x	0221_12	304526
L1705776-20 WG2230690 1x	0221_13	307065
L1705776-13 WG2230690 20x	0221_15	307787

Instrument: VOCMS27 • File ID: 0222_02

02/22/24 11:18

Sample ID	File ID	8260-FLUOROBENZENE Response
Standard	0222_02	312476
Upper Limit		624952
Lower Limit		156238
LCS R4037174-1 WG2231600 1x	0222_03	318906

INTERNAL STANDARD SUMMARY

Instrument: VOCMS27 • File ID: 0222_02

02/22/24 11:18

Sample ID	File ID	8260-FLUOROBENZENE Response
LCSD R4037174-2 WG2231600 1x	0222_04	310746
BLANK R4037174-3 WG2231600 1x	0222_06	309059
L1705776-21 WG2231600 10x	0222_08	304679
MS R4037174-4 WG2231600 1x	0222_09	310642
MSD R4037174-5 WG2231600 1x	0222_10	322726

GLOSSARY OF TERMS

Guide to Reading and Understanding Your Laboratory Report

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

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

Abbreviations and Definitions

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

Qualifier	Description
E4	Concentration estimated. Analyte was detected below laboratory minimum reporting level (MRL) but above MDL.
L1	The associated blank spike recovery was above laboratory acceptance limits.
L2	The associated blank spike recovery was below laboratory acceptance limits.
M1	Matrix spike recovery was high, the method control sample recovery was acceptable.
M2	Matrix spike recovery was low, the method control sample recovery was acceptable.
M3	The spike recovery value is unusable since the analyte concentration in the sample is disproportionate to the spike level. The associated blank spike recovery was acceptable.
R5	MS/MSD RPD exceeded the laboratory acceptance limit. Recovery met acceptance criteria.
R8	Sample RPD exceeded the method acceptance limit.
S10	Surrogate recovery was above laboratory and method acceptance limits.

ACCREDITATIONS & LOCATIONS

Pace Analytical National 12065 Lebanon Rd Mount Juliet, TN 37122

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

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

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

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

Company Name/Address:
Pinyon Environmental

3222 S. Vance Street Suite 200
Lakewood, CO 80227

Billing Information:
Accounts Payable
3222 S Vance Street
Suite 200
Lakewood, CO 80227

Pres Chk

Report to:
Isbella Foster and Andrew Parker

Email To: **foster@pinyon-env.com;parker@pinyon-**

Project Description:
Nammo TTU Groundwater Monitoring

City/State Collected: Mesa, AZ

Please Circle: PT (MT) CT ET

Phone: **303-785-7697**

Client Project #
722152201.002

Lab Project #
PINYONMAZ-722152201

Collected by (print):
Isabella Foster

Site/Facility ID #

P.O. #

Collected by (signature):
Immediately
Packed on Ice N ___ Y X

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

Quote #

Date Results Needed: Standard TAT

Analysis / Container / Preservative

Sample ID	Comp/Grab	Matrix *	Depth	Date	Time	No. of Cntrs	PERCHLORATE 125mlHDPE-NoPres	V8260AZ 40mlAmb-HCl	V8260LL14D 40mlAmb-HCl							Remarks	Sample # (lab only)
TTU-1-50-20240213	C	GW	50	2/13/24	1027	7	X	X	X								-01
TTU-2-114-20240213	C	GW	114	2/13/24	1101	7	X	X	X								-02
TTU-3-108-20240212	G	GW	108	2/12/24	1518	7	X	X	X								-03
TTU-4-57-20240213	G	GW	57	2/13/24	1528	7	X	X	X								-04
TTU-5-110-20240213	G	GW	110	2/13/24	1450	7	X	X	X								-05
TTU-6-143-20240212	G	GW	143	2/12/24	1458	7	X	X	X								-06
TTU-7-345-20240212	G	GW	345	2/12/24	1435	7	X	X	X								-07
~~TTU-8~~		GW															
TTU-9a-61-20240213	G	GW	61	2/13/24	0932	14	X	X	X							MS/MSD	-08
~~TTU-10~~	G	GW															

Chain of Custody Page 1 of 3

Pace
PEOPLE ADVANCING SCIENCE

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

SDG # L1705776

Table #

Acctnum: **PINYONMAZ**
Template: **T205653**
Prelogin: **P1054553**
PM: **288 - Daphne Richards**
PB:

Shipped Via:

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

Remarks:

pH _____ Temp _____
Flow _____ Other _____

Samples returned via:
__ UPS __ FedEx __ Courier _____

Tracking #

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

Relinquished by : (Signature)	Date:	Time:	Received by: (Signature)	Trip Blank Received: (Yes / No) HCL / MeoH TBR	If preservation required by Login: Date/Time
	2/13/24	1550		1	
Relinquished by : (Signature)	Date:	Time:	Received by: (Signature)	Temp: °C Bottles Received:	
	2/13/24	1658			
Relinquished by : (Signature)	Date:	Time:	Received for lab by: (Signature)	Date:	Time:	Hold:	Condition: NCF / OK
	2-13-24	1805		2/14/24	0800		

Company Name/Address:
Pinyon Environmental

3222 S. Vance Street Suite 200
Lakewood, CO 80227

Billing Information:
Accounts Payable
3222 S Vance Street
Suite 200
Lakewood, CO 80227

Pres Chk

Report to:
Isbella Foster and Andrew Parker

Email To: **foster@pinyon-env.com;parker@pinyon-**

Project Description:
Nammo TTU Groundwater Monitoring

City/State Collected: Mesa, AZ

Please Circle: PT (MT) CT ET

Phone: **303-785-7697**

Client Project #
722152201.002

Lab Project #
PINYONMAZ-722152201

Collected by (print):
Isabella Foster

Site/Facility ID #

P.O. #

Collected by (signature):

Immediately Packed on Ice N ___ Y X

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

Quote #

Date Results Needed
Standard TAT

Analysis / Container / Preservative

Chain of Custody Page 1 of 3

Pace
PEOPLE ADVANCING SCIENCE

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

SDG # L1205726

Table #

Acctnum: **PINYONMAZ**
Template: **T205653**
Prelogin: **P1054553**
PM: **288 - Daphne Richards**
PB:

Shipped Via:

Sample ID	Comp/Grab	Matrix *	Depth	Date	Time	No. of Cntrs	PERCHLORATE 125mlHDPE-NoPres	V8260AZ 40mlAmb-HCl	V8260LL14D 40mlAmb-HCl	Remarks	Sample # (lab only)
TTU-11-73-20240213	G	GW	73	2/13/24	1225	7	X	X	X		-09
TTU-12-82-20240212	G	GW	82	2/12/24	1000	7	X	X	X		-10
TTU-13-51-20240212	G	GW	51	2/12/24	0905	14	X	X	X	MS/MSD	-11
TTU-14-64-20240212	G	GW	64	2/12/24	0935	7	X	X	X		-12
TTU-15-75-20240213	G	GW	75	2/13/24	1302	7	X	X	X		-13
TTU-16-80-20240213	G	GW	80	2/13/24	1325	7	X	X	X		-14
TTU-17-80-20240212	G	GW	80	2/12/24	1035	7	X	X	X		-15
TTU-19-73-20240213	G	GW	73	2/13/24	1200	7	X	X	X		-16
TTU-20-73-20240213	G	GW	73	2/13/24	1140	7	X	X	X		-17
TTU-EX-5-80-20240213	G	GW	80	2/13/24	1415	7	X	X	X		-18

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

Remarks:

pH _____ Temp _____

Flow _____ Other _____

Samples returned via:
__ UPS __ FedEx __ Courier _____

Tracking #

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

Relinquished by : (Signature)	Date:	Time:	Received by: (Signature)	Trip Blank Received: Yes / No HCl / MeoH TBR	If preservation required by Login: Date/Time
	2/13/24	1556		1	
Relinquished by : (Signature)	Date:	Time:	Received by: (Signature)	Temp: °C Bottles Received:	
	2/13/24	1658			
Relinquished by : (Signature)	Date:	Time:	Received for lab by: (Signature)	Date:	Time:	Hold:	Condition: NCF / OK
	2-13-24	1855		2/14/24	0800		

Company Name/Address:
Pinyon Environmental

3222 S. Vance Street Suite 200
Lakewood, CO 80227

Billing Information:
Accounts Payable
3222 S Vance Street
Suite 200
Lakewood, CO 80227

Pres Chk

Report to:
Isbella Foster and Andrew Parker

Email To: **foster@pinyon-env.com;parker@pinyon-**

Project Description:
Nammo TTU Groundwater Monitoring

City/State Collected: Mesa, Az

Please Circle: PT (MT) CT ET

Phone: **303-785-7697**

Client Project #
722152201.002

Lab Project #
PINYONMAZ-722152201

Collected by (print):
Isabella Foster

Site/Facility ID #

P.O. #

Collected by (signature):

Immediately Packed on Ice N ___ Y X

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

Quote #

Date Results Needed: Standard TAT

No. of Cntrs

Analysis / Container / Preservative

Sample ID	Comp/Grab	Matrix *	Depth	Date	Time	No. of Cntrs	PERCHLORATE 125mlHDPE-NoPres	V8260AZ 40mlAmb-HCl	V8260LL14D 40mlAmb-HCl	Remarks	Sample # (lab only)
TTU-EX-1-69-20240213	G	GW	69	2-13-24	1343	7	X	X	X		-19
TTU-EX-2-74-20240213	G	GW	74	2-13-24	1359	7	X	X	X		-20
TTU-EX-3-76-20240212	G	GW	76	2-12-24	1100	7	X	X	X		-21
TTU-EX-4-77-20240212	G	GW	77	2-12-24	1113	7	X	X	X		-22
Dup-01	—	GW	—	—	—	7	X	X	X		-23
Dup-02	—	GW	—	—	—	7	X	X	X		-24
Trip Blank	—	GW	—	—	—	1		X	X		-25
		GW									
		GW									
		GW									

Chain of Custody Page 3 of 3

Pace
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MT JULIET, TN
12065 Lebanon Rd Mount Juliet, TN 37122
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SDG # L1705776

Table #

Acctnum: **PINYONMAZ**
Template: **T205653**
Prelogin: **P1054553**
PM: **288 - Daphne Richards**
PB:

Shipped Via:

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

Remarks:

pH _____ Temp _____

Flow _____ Other _____

Samples returned via:
__ UPS __ FedEx __ Courier _____

Tracking #

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

Relinquished by : (Signature)	Date:	Time:	Received by: (Signature)	Trip Blank Received: (Yes) / No HCL / MeoH TBR	If preservation required by Login: Date/Time
	2/13/24	1550			
Relinquished by : (Signature)	Date:	Time:	Received by: (Signature)	Temp: °C Bottles Received:	
	2/13/24	1658			
Relinquished by : (Signature)	Date:	Time:	Received for lab by: (Signature)	Date:	Time:
	2-13-24	1800 SLM		2/14/24	0800

Hold: | Condition: NCF / (OK)

LT05776

Tracking Numbers	Temperature
SWA	DFA 0.9 to
SWA	DFA 1.5 to

Name

Date

ANALYTICAL REPORT

March 27, 2024

Revised Report

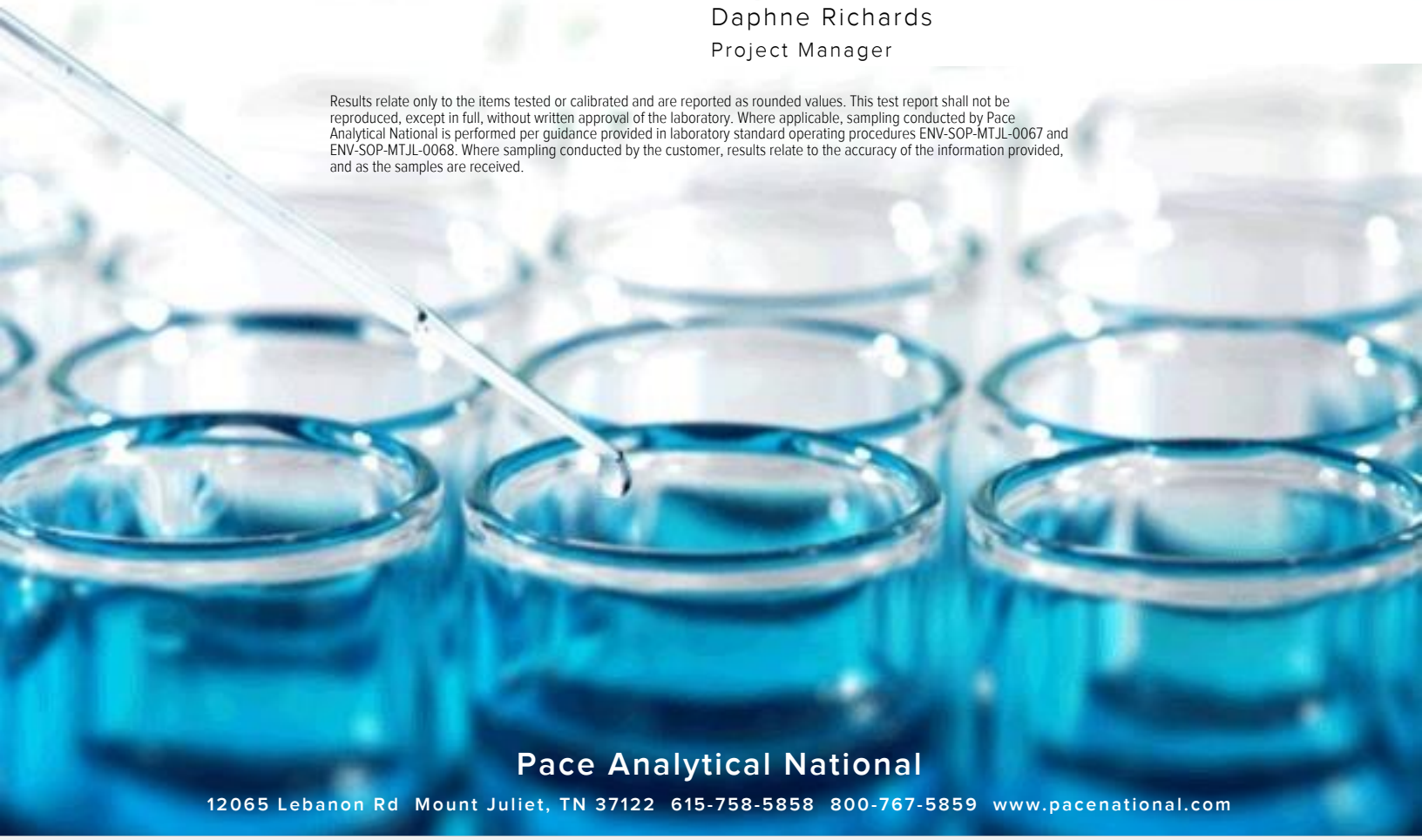
Pinyon Environmental

Sample Delivery Group:	L1706682
Samples Received:	02/16/2024
Project Number:	722152201.002
Description:	Nammo TTU

Report To: Isbella Foster and Andrew Parker
3222 S. Vance Street Suite 200
Lakewood, CO 80227

Entire Report Reviewed By: Daphne R Richards

Daphne Richards
Project Manager



Pace Analytical National
12065 Lebanon Rd Mount Juliet, TN 37122 615-758-5858 800-767-5859 www.pacenational.com

TABLE OF CONTENTS

PF-2-400-20240214 L1706682-01 GW

Collected by: Isabella Foster | Collected date/time: 02/14/24 15:00 | Received date/time: 02/16/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2230378	1	02/21/24 05:43	02/21/24 05:43	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2231965	1	02/25/24 15:16	02/25/24 15:16	JCP	Mt. Juliet, TN

TTU-8-164-20240214 L1706682-02 GW

Collected by: Isabella Foster | Collected date/time: 02/14/24 15:20 | Received date/time: 02/16/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2228553	1	03/01/24 05:36	03/01/24 05:36	ASM	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2230378	1	02/21/24 06:05	02/21/24 06:05	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2231221	1	02/22/24 05:29	02/22/24 05:29	ACG	Mt. Juliet, TN

TTU-10-172-20240214 L1706682-03 GW

Collected by: Isabella Foster | Collected date/time: 02/14/24 13:49 | Received date/time: 02/16/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2234696	1	03/03/24 22:32	03/03/24 22:32	ASM	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2230378	1	02/21/24 06:26	02/21/24 06:26	JCP	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2231221	1	02/22/24 05:51	02/22/24 05:51	ACG	Mt. Juliet, TN

DUP-03 L1706682-04 GW

Collected by: Isabella Foster | Collected date/time: 02/14/24 00:00 | Received date/time: 02/16/24 08:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2234696	1	03/03/24 23:00	03/03/24 23:00	ASM	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2230965	1	02/22/24 05:07	02/22/24 05:07	ADM	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2231221	1	02/22/24 06:13	02/22/24 06:13	ACG	Mt. Juliet, TN

CASE NARRATIVE

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

Daphne Richards
Project Manager

Report Revision History

Level II Report - Version 1: 03/07/24 09:14

Project Narrative

Sample description and project number update

Sample Delivery Group (SDG) Narrative

Insufficient sample volume to perform MS/MSD analyses per method QC requirements.

Lab Sample ID	Project Sample ID	Method
L1706682-04	DUP-03	8260B

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		11.3	50.0	1	02/21/2024 05:43	WG2230378
Acrolein	U		2.54	50.0	1	02/21/2024 05:43	WG2230378
Acrylonitrile	U		0.671	10.0	1	02/21/2024 05:43	WG2230378
Benzene	U		0.0941	1.00	1	02/21/2024 05:43	WG2230378
Bromobenzene	U		0.118	1.00	1	02/21/2024 05:43	WG2230378
Bromodichloromethane	U		0.136	1.00	1	02/21/2024 05:43	WG2230378
Bromoform	U		0.129	1.00	1	02/21/2024 05:43	WG2230378
Bromomethane	U		0.605	5.00	1	02/21/2024 05:43	WG2230378
1,3-Butadiene	U		0.299	2.00	1	02/21/2024 05:43	WG2230378
n-Butylbenzene	U		0.157	1.00	1	02/21/2024 05:43	WG2230378
sec-Butylbenzene	U		0.125	1.00	1	02/21/2024 05:43	WG2230378
tert-Butylbenzene	U		0.127	1.00	1	02/21/2024 05:43	WG2230378
Carbon tetrachloride	U		0.128	1.00	1	02/21/2024 05:43	WG2230378
Carbon disulfide	0.172	B1 E4	0.0962	1.00	1	02/21/2024 05:43	WG2230378
Chlorobenzene	U		0.116	1.00	1	02/21/2024 05:43	WG2230378
Chlorodibromomethane	U		0.140	1.00	1	02/21/2024 05:43	WG2230378
Chloroethane	U		0.192	5.00	1	02/21/2024 05:43	WG2230378
Chloroform	U		0.111	5.00	1	02/21/2024 05:43	WG2230378
Chloromethane	U		0.960	2.50	1	02/21/2024 05:43	WG2230378
Cyclohexane	U		0.188	1.00	1	02/21/2024 05:43	WG2230378
2-Chlorotoluene	U		0.106	1.00	1	02/21/2024 05:43	WG2230378
4-Chlorotoluene	U		0.114	1.00	1	02/21/2024 05:43	WG2230378
1,2-Dibromo-3-Chloropropane	U		0.276	5.00	1	02/21/2024 05:43	WG2230378
1,2-Dibromoethane	U		0.126	1.00	1	02/21/2024 05:43	WG2230378
Dibromomethane	U		0.122	1.00	1	02/21/2024 05:43	WG2230378
1,2-Dichlorobenzene	U		0.107	1.00	1	02/21/2024 05:43	WG2230378
1,3-Dichlorobenzene	U		0.110	1.00	1	02/21/2024 05:43	WG2230378
1,4-Dichlorobenzene	U		0.120	1.00	1	02/21/2024 05:43	WG2230378
Dichlorodifluoromethane	U		0.374	5.00	1	02/21/2024 05:43	WG2230378
1,1-Dichloroethane	U		0.100	1.00	1	02/21/2024 05:43	WG2230378
1,2-Dichloroethane	U		0.0819	1.00	1	02/21/2024 05:43	WG2230378
1,1-Dichloroethene	U		0.188	1.00	1	02/21/2024 05:43	WG2230378
cis-1,2-Dichloroethene	U		0.126	1.00	1	02/21/2024 05:43	WG2230378
trans-1,2-Dichloroethene	U		0.149	1.00	1	02/21/2024 05:43	WG2230378
1,2-Dichloropropane	U		0.149	1.00	1	02/21/2024 05:43	WG2230378
1,1-Dichloropropene	U		0.142	1.00	1	02/21/2024 05:43	WG2230378
1,3-Dichloropropane	U		0.110	1.00	1	02/21/2024 05:43	WG2230378
cis-1,3-Dichloropropene	U		0.111	1.00	1	02/21/2024 05:43	WG2230378
trans-1,3-Dichloropropene	U		0.118	1.00	1	02/21/2024 05:43	WG2230378
2,2-Dichloropropane	U		0.161	1.00	1	02/21/2024 05:43	WG2230378
Dicyclopentadiene	U		0.253	1.00	1	02/21/2024 05:43	WG2230378
Di-isopropyl ether	U	M1	0.105	1.00	1	02/21/2024 05:43	WG2230378
Ethylbenzene	U		0.137	1.00	1	02/21/2024 05:43	WG2230378
4-Ethyltoluene	U		0.208	1.00	1	02/21/2024 05:43	WG2230378
Hexachloro-1,3-butadiene	U		0.337	1.00	1	02/21/2024 05:43	WG2230378
n-Hexane	U		0.749	10.0	1	02/21/2024 05:43	WG2230378
Isopropylbenzene	U		0.105	1.00	1	02/21/2024 05:43	WG2230378
p-Isopropyltoluene	U		0.120	1.00	1	02/21/2024 05:43	WG2230378
2-Butanone (MEK)	U		1.19	10.0	1	02/21/2024 05:43	WG2230378
Methyl Cyclohexane	U		0.660	1.00	1	02/21/2024 05:43	WG2230378
Methylene Chloride	U		0.430	5.00	1	02/21/2024 05:43	WG2230378
4-Methyl-2-pentanone (MIBK)	U	M1	0.478	10.0	1	02/21/2024 05:43	WG2230378
Methyl tert-butyl ether	U		0.101	1.00	1	02/21/2024 05:43	WG2230378
Naphthalene	U		1.00	5.00	1	02/21/2024 05:43	WG2230378
Propene	U		0.936	2.50	1	02/21/2024 05:43	WG2230378
n-Propylbenzene	U		0.0993	1.00	1	02/21/2024 05:43	WG2230378

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Styrene	U		0.118	1.00	1	02/21/2024 05:43	WG2230378
1,1,1,2-Tetrachloroethane	U		0.147	1.00	1	02/21/2024 05:43	WG2230378
1,1,2,2-Tetrachloroethane	U		0.133	1.00	1	02/21/2024 05:43	WG2230378
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00	1	02/21/2024 05:43	WG2230378
Tetrachloroethene	U		0.300	1.00	1	02/21/2024 05:43	WG2230378
Toluene	U		0.278	1.00	1	02/21/2024 05:43	WG2230378
1,2,3-Trichlorobenzene	U		0.230	1.00	1	02/21/2024 05:43	WG2230378
1,2,4-Trichlorobenzene	U		0.481	1.00	1	02/21/2024 05:43	WG2230378
1,1,1-Trichloroethane	U		0.149	1.00	1	02/21/2024 05:43	WG2230378
1,1,2-Trichloroethane	U		0.158	1.00	1	02/21/2024 05:43	WG2230378
Trichloroethene	U		0.190	1.00	1	02/21/2024 05:43	WG2230378
Trichlorofluoromethane	U		0.160	5.00	1	02/21/2024 05:43	WG2230378
1,2,3-Trichloropropane	U		0.237	2.50	1	02/21/2024 05:43	WG2230378
1,2,4-Trimethylbenzene	U		0.322	1.00	1	02/21/2024 05:43	WG2230378
1,2,3-Trimethylbenzene	U		0.104	1.00	1	02/21/2024 05:43	WG2230378
1,3,5-Trimethylbenzene	U		0.104	1.00	1	02/21/2024 05:43	WG2230378
Vinyl chloride	U		0.234	1.00	1	02/21/2024 05:43	WG2230378
Xylenes, Total	U		0.174	3.00	1	02/21/2024 05:43	WG2230378
(S) Toluene-d8	*114*			*80.0-120*		*02/21/2024 05:43*	WG2230378
(S) 4-Bromofluorobenzene	*107*			*77.0-126*		*02/21/2024 05:43*	WG2230378
(S) 1,2-Dichloroethane-d4	*124*			*70.0-130*		*02/21/2024 05:43*	WG2230378

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
1,4-Dioxane	U		0.597	3.00	1	02/25/2024 15:16	WG2231965
(S) Toluene-d8	*94.2*			*77.0-127*		*02/25/2024 15:16*	WG2231965

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	U		0.300	4.00	1	03/01/2024 05:36	WG2228553

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		11.3	50.0	1	02/21/2024 06:05	WG2230378
Acrolein	U		2.54	50.0	1	02/21/2024 06:05	WG2230378
Acrylonitrile	U		0.671	10.0	1	02/21/2024 06:05	WG2230378
Benzene	U		0.0941	1.00	1	02/21/2024 06:05	WG2230378
Bromobenzene	U		0.118	1.00	1	02/21/2024 06:05	WG2230378
Bromodichloromethane	U		0.136	1.00	1	02/21/2024 06:05	WG2230378
Bromoform	U		0.129	1.00	1	02/21/2024 06:05	WG2230378
Bromomethane	U	R7	0.605	5.00	1	02/21/2024 06:05	WG2230378
1,3-Butadiene	U	R7	0.299	2.00	1	02/21/2024 06:05	WG2230378
n-Butylbenzene	U		0.157	1.00	1	02/21/2024 06:05	WG2230378
sec-Butylbenzene	U		0.125	1.00	1	02/21/2024 06:05	WG2230378
tert-Butylbenzene	U		0.127	1.00	1	02/21/2024 06:05	WG2230378
Carbon tetrachloride	U		0.128	1.00	1	02/21/2024 06:05	WG2230378
Carbon disulfide	0.186	B1 E4 R7	0.0962	1.00	1	02/21/2024 06:05	WG2230378
Chlorobenzene	U		0.116	1.00	1	02/21/2024 06:05	WG2230378
Chlorodibromomethane	U		0.140	1.00	1	02/21/2024 06:05	WG2230378
Chloroethane	U		0.192	5.00	1	02/21/2024 06:05	WG2230378
Chloroform	U		0.111	5.00	1	02/21/2024 06:05	WG2230378
Chloromethane	U	R7	0.960	2.50	1	02/21/2024 06:05	WG2230378
Cyclohexane	U		0.188	1.00	1	02/21/2024 06:05	WG2230378
2-Chlorotoluene	U		0.106	1.00	1	02/21/2024 06:05	WG2230378
4-Chlorotoluene	U		0.114	1.00	1	02/21/2024 06:05	WG2230378
1,2-Dibromo-3-Chloropropane	U		0.276	5.00	1	02/21/2024 06:05	WG2230378
1,2-Dibromoethane	U		0.126	1.00	1	02/21/2024 06:05	WG2230378
Dibromomethane	U		0.122	1.00	1	02/21/2024 06:05	WG2230378
1,2-Dichlorobenzene	U		0.107	1.00	1	02/21/2024 06:05	WG2230378
1,3-Dichlorobenzene	U		0.110	1.00	1	02/21/2024 06:05	WG2230378
1,4-Dichlorobenzene	U		0.120	1.00	1	02/21/2024 06:05	WG2230378
Dichlorodifluoromethane	U	R7	0.374	5.00	1	02/21/2024 06:05	WG2230378
1,1-Dichloroethane	U		0.100	1.00	1	02/21/2024 06:05	WG2230378
1,2-Dichloroethane	U		0.0819	1.00	1	02/21/2024 06:05	WG2230378
1,1-Dichloroethene	U		0.188	1.00	1	02/21/2024 06:05	WG2230378
cis-1,2-Dichloroethene	U		0.126	1.00	1	02/21/2024 06:05	WG2230378
trans-1,2-Dichloroethene	U		0.149	1.00	1	02/21/2024 06:05	WG2230378
1,2-Dichloropropane	U		0.149	1.00	1	02/21/2024 06:05	WG2230378
1,1-Dichloropropene	U		0.142	1.00	1	02/21/2024 06:05	WG2230378
1,3-Dichloropropane	U		0.110	1.00	1	02/21/2024 06:05	WG2230378
cis-1,3-Dichloropropene	U		0.111	1.00	1	02/21/2024 06:05	WG2230378
trans-1,3-Dichloropropene	U		0.118	1.00	1	02/21/2024 06:05	WG2230378
2,2-Dichloropropane	U		0.161	1.00	1	02/21/2024 06:05	WG2230378
Dicyclopentadiene	U		0.253	1.00	1	02/21/2024 06:05	WG2230378
Di-isopropyl ether	U		0.105	1.00	1	02/21/2024 06:05	WG2230378
Ethylbenzene	U		0.137	1.00	1	02/21/2024 06:05	WG2230378
4-Ethyltoluene	U		0.208	1.00	1	02/21/2024 06:05	WG2230378
Hexachloro-1,3-butadiene	U		0.337	1.00	1	02/21/2024 06:05	WG2230378
n-Hexane	U		0.749	10.0	1	02/21/2024 06:05	WG2230378
Isopropylbenzene	U		0.105	1.00	1	02/21/2024 06:05	WG2230378
p-Isopropyltoluene	U		0.120	1.00	1	02/21/2024 06:05	WG2230378
2-Butanone (MEK)	U		1.19	10.0	1	02/21/2024 06:05	WG2230378
Methyl Cyclohexane	U		0.660	1.00	1	02/21/2024 06:05	WG2230378

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Methylene Chloride	U		0.430	5.00	1	02/21/2024 06:05	WG2230378
4-Methyl-2-pentanone (MIBK)	U		0.478	10.0	1	02/21/2024 06:05	WG2230378
Methyl tert-butyl ether	U		0.101	1.00	1	02/21/2024 06:05	WG2230378
Naphthalene	U		1.00	5.00	1	02/21/2024 06:05	WG2230378
Propene	U	R7	0.936	2.50	1	02/21/2024 06:05	WG2230378
n-Propylbenzene	U		0.0993	1.00	1	02/21/2024 06:05	WG2230378
Styrene	U		0.118	1.00	1	02/21/2024 06:05	WG2230378
1,1,1,2-Tetrachloroethane	U		0.147	1.00	1	02/21/2024 06:05	WG2230378
1,1,2,2-Tetrachloroethane	U		0.133	1.00	1	02/21/2024 06:05	WG2230378
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00	1	02/21/2024 06:05	WG2230378
Tetrachloroethene	U		0.300	1.00	1	02/21/2024 06:05	WG2230378
Toluene	U		0.278	1.00	1	02/21/2024 06:05	WG2230378
1,2,3-Trichlorobenzene	U		0.230	1.00	1	02/21/2024 06:05	WG2230378
1,2,4-Trichlorobenzene	U		0.481	1.00	1	02/21/2024 06:05	WG2230378
1,1,1-Trichloroethane	U		0.149	1.00	1	02/21/2024 06:05	WG2230378
1,1,2-Trichloroethane	U		0.158	1.00	1	02/21/2024 06:05	WG2230378
Trichloroethene	U		0.190	1.00	1	02/21/2024 06:05	WG2230378
Trichlorofluoromethane	U		0.160	5.00	1	02/21/2024 06:05	WG2230378
1,2,3-Trichloropropane	U		0.237	2.50	1	02/21/2024 06:05	WG2230378
1,2,4-Trimethylbenzene	U		0.322	1.00	1	02/21/2024 06:05	WG2230378
1,2,3-Trimethylbenzene	U		0.104	1.00	1	02/21/2024 06:05	WG2230378
1,3,5-Trimethylbenzene	U		0.104	1.00	1	02/21/2024 06:05	WG2230378
Vinyl chloride	U		0.234	1.00	1	02/21/2024 06:05	WG2230378
Xylenes, Total	U		0.174	3.00	1	02/21/2024 06:05	WG2230378
(S) Toluene-d8	*114*			*80.0-120*		*02/21/2024 06:05*	WG2230378
(S) 4-Bromofluorobenzene	*106*			*77.0-126*		*02/21/2024 06:05*	WG2230378
(S) 1,2-Dichloroethane-d4	*124*			*70.0-130*		*02/21/2024 06:05*	WG2230378

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
1,4-Dioxane	U		0.597	3.00	1	02/22/2024 05:29	WG2231221
(S) Toluene-d8	*96.0*			*77.0-127*		*02/22/2024 05:29*	WG2231221

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	U		0.300	4.00	1	03/03/2024 22:32	WG2234696

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		11.3	50.0	1	02/21/2024 06:26	WG2230378
Acrolein	U		2.54	50.0	1	02/21/2024 06:26	WG2230378
Acrylonitrile	U		0.671	10.0	1	02/21/2024 06:26	WG2230378
Benzene	U		0.0941	1.00	1	02/21/2024 06:26	WG2230378
Bromobenzene	U		0.118	1.00	1	02/21/2024 06:26	WG2230378
Bromodichloromethane	U		0.136	1.00	1	02/21/2024 06:26	WG2230378
Bromoform	U		0.129	1.00	1	02/21/2024 06:26	WG2230378
Bromomethane	U	R7	0.605	5.00	1	02/21/2024 06:26	WG2230378
1,3-Butadiene	U	R7	0.299	2.00	1	02/21/2024 06:26	WG2230378
n-Butylbenzene	U		0.157	1.00	1	02/21/2024 06:26	WG2230378
sec-Butylbenzene	U		0.125	1.00	1	02/21/2024 06:26	WG2230378
tert-Butylbenzene	U		0.127	1.00	1	02/21/2024 06:26	WG2230378
Carbon tetrachloride	U		0.128	1.00	1	02/21/2024 06:26	WG2230378
Carbon disulfide	0.170	B1 E4 R7	0.0962	1.00	1	02/21/2024 06:26	WG2230378
Chlorobenzene	U		0.116	1.00	1	02/21/2024 06:26	WG2230378
Chlorodibromomethane	U		0.140	1.00	1	02/21/2024 06:26	WG2230378
Chloroethane	U		0.192	5.00	1	02/21/2024 06:26	WG2230378
Chloroform	U		0.111	5.00	1	02/21/2024 06:26	WG2230378
Chloromethane	U	R7	0.960	2.50	1	02/21/2024 06:26	WG2230378
Cyclohexane	U		0.188	1.00	1	02/21/2024 06:26	WG2230378
2-Chlorotoluene	U		0.106	1.00	1	02/21/2024 06:26	WG2230378
4-Chlorotoluene	U		0.114	1.00	1	02/21/2024 06:26	WG2230378
1,2-Dibromo-3-Chloropropane	U		0.276	5.00	1	02/21/2024 06:26	WG2230378
1,2-Dibromoethane	U		0.126	1.00	1	02/21/2024 06:26	WG2230378
Dibromomethane	U		0.122	1.00	1	02/21/2024 06:26	WG2230378
1,2-Dichlorobenzene	U		0.107	1.00	1	02/21/2024 06:26	WG2230378
1,3-Dichlorobenzene	U		0.110	1.00	1	02/21/2024 06:26	WG2230378
1,4-Dichlorobenzene	U		0.120	1.00	1	02/21/2024 06:26	WG2230378
Dichlorodifluoromethane	U	R7	0.374	5.00	1	02/21/2024 06:26	WG2230378
1,1-Dichloroethane	U		0.100	1.00	1	02/21/2024 06:26	WG2230378
1,2-Dichloroethane	U		0.0819	1.00	1	02/21/2024 06:26	WG2230378
1,1-Dichloroethene	U		0.188	1.00	1	02/21/2024 06:26	WG2230378
cis-1,2-Dichloroethene	U		0.126	1.00	1	02/21/2024 06:26	WG2230378
trans-1,2-Dichloroethene	U		0.149	1.00	1	02/21/2024 06:26	WG2230378
1,2-Dichloropropane	U		0.149	1.00	1	02/21/2024 06:26	WG2230378
1,1-Dichloropropene	U		0.142	1.00	1	02/21/2024 06:26	WG2230378
1,3-Dichloropropane	U		0.110	1.00	1	02/21/2024 06:26	WG2230378
cis-1,3-Dichloropropene	U		0.111	1.00	1	02/21/2024 06:26	WG2230378
trans-1,3-Dichloropropene	U		0.118	1.00	1	02/21/2024 06:26	WG2230378
2,2-Dichloropropane	U		0.161	1.00	1	02/21/2024 06:26	WG2230378
Dicyclopentadiene	U		0.253	1.00	1	02/21/2024 06:26	WG2230378
Di-isopropyl ether	U		0.105	1.00	1	02/21/2024 06:26	WG2230378
Ethylbenzene	U		0.137	1.00	1	02/21/2024 06:26	WG2230378
4-Ethyltoluene	U		0.208	1.00	1	02/21/2024 06:26	WG2230378
Hexachloro-1,3-butadiene	U		0.337	1.00	1	02/21/2024 06:26	WG2230378
n-Hexane	U		0.749	10.0	1	02/21/2024 06:26	WG2230378
Isopropylbenzene	U		0.105	1.00	1	02/21/2024 06:26	WG2230378
p-Isopropyltoluene	U		0.120	1.00	1	02/21/2024 06:26	WG2230378
2-Butanone (MEK)	U		1.19	10.0	1	02/21/2024 06:26	WG2230378
Methyl Cyclohexane	U		0.660	1.00	1	02/21/2024 06:26	WG2230378

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Methylene Chloride	U		0.430	5.00	1	02/21/2024 06:26	WG2230378
4-Methyl-2-pentanone (MIBK)	U		0.478	10.0	1	02/21/2024 06:26	WG2230378
Methyl tert-butyl ether	U		0.101	1.00	1	02/21/2024 06:26	WG2230378
Naphthalene	U		1.00	5.00	1	02/21/2024 06:26	WG2230378
Propene	U	R7	0.936	2.50	1	02/21/2024 06:26	WG2230378
n-Propylbenzene	U		0.0993	1.00	1	02/21/2024 06:26	WG2230378
Styrene	U		0.118	1.00	1	02/21/2024 06:26	WG2230378
1,1,1,2-Tetrachloroethane	U		0.147	1.00	1	02/21/2024 06:26	WG2230378
1,1,2,2-Tetrachloroethane	U		0.133	1.00	1	02/21/2024 06:26	WG2230378
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00	1	02/21/2024 06:26	WG2230378
Tetrachloroethene	U		0.300	1.00	1	02/21/2024 06:26	WG2230378
Toluene	U		0.278	1.00	1	02/21/2024 06:26	WG2230378
1,2,3-Trichlorobenzene	U		0.230	1.00	1	02/21/2024 06:26	WG2230378
1,2,4-Trichlorobenzene	U		0.481	1.00	1	02/21/2024 06:26	WG2230378
1,1,1-Trichloroethane	U		0.149	1.00	1	02/21/2024 06:26	WG2230378
1,1,2-Trichloroethane	U		0.158	1.00	1	02/21/2024 06:26	WG2230378
Trichloroethene	U		0.190	1.00	1	02/21/2024 06:26	WG2230378
Trichlorofluoromethane	U		0.160	5.00	1	02/21/2024 06:26	WG2230378
1,2,3-Trichloropropane	U		0.237	2.50	1	02/21/2024 06:26	WG2230378
1,2,4-Trimethylbenzene	U		0.322	1.00	1	02/21/2024 06:26	WG2230378
1,2,3-Trimethylbenzene	U		0.104	1.00	1	02/21/2024 06:26	WG2230378
1,3,5-Trimethylbenzene	U		0.104	1.00	1	02/21/2024 06:26	WG2230378
Vinyl chloride	U		0.234	1.00	1	02/21/2024 06:26	WG2230378
Xylenes, Total	U		0.174	3.00	1	02/21/2024 06:26	WG2230378
(S) Toluene-d8	*113*			*80.0-120*		*02/21/2024 06:26*	WG2230378
(S) 4-Bromofluorobenzene	*105*			*77.0-126*		*02/21/2024 06:26*	WG2230378
(S) 1,2-Dichloroethane-d4	*127*			*70.0-130*		*02/21/2024 06:26*	WG2230378

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
1,4-Dioxane	U		0.597	3.00	1	02/22/2024 05:51	WG2231221
(S) Toluene-d8	*95.8*			*77.0-127*		*02/22/2024 05:51*	WG2231221

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	U		0.300	4.00	1	03/03/2024 23:00	WG2234696

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		11.3	50.0	1	02/22/2024 05:07	WG2230965
Acrolein	U	L1	2.54	50.0	1	02/22/2024 05:07	WG2230965
Acrylonitrile	U		0.671	10.0	1	02/22/2024 05:07	WG2230965
Benzene	U		0.0941	1.00	1	02/22/2024 05:07	WG2230965
Bromobenzene	U		0.118	1.00	1	02/22/2024 05:07	WG2230965
Bromodichloromethane	U		0.136	1.00	1	02/22/2024 05:07	WG2230965
Bromoform	U		0.129	1.00	1	02/22/2024 05:07	WG2230965
Bromomethane	U	R7	0.605	5.00	1	02/22/2024 05:07	WG2230965
1,3-Butadiene	U		0.299	2.00	1	02/22/2024 05:07	WG2230965
n-Butylbenzene	U		0.157	1.00	1	02/22/2024 05:07	WG2230965
sec-Butylbenzene	U		0.125	1.00	1	02/22/2024 05:07	WG2230965
tert-Butylbenzene	U		0.127	1.00	1	02/22/2024 05:07	WG2230965
Carbon tetrachloride	U		0.128	1.00	1	02/22/2024 05:07	WG2230965
Carbon disulfide	U		0.0962	1.00	1	02/22/2024 05:07	WG2230965
Chlorobenzene	U		0.116	1.00	1	02/22/2024 05:07	WG2230965
Chlorodibromomethane	U		0.140	1.00	1	02/22/2024 05:07	WG2230965
Chloroethane	U		0.192	5.00	1	02/22/2024 05:07	WG2230965
Chloroform	U		0.111	5.00	1	02/22/2024 05:07	WG2230965
Chloromethane	U		0.960	2.50	1	02/22/2024 05:07	WG2230965
Cyclohexane	U		0.188	1.00	1	02/22/2024 05:07	WG2230965
2-Chlorotoluene	U		0.106	1.00	1	02/22/2024 05:07	WG2230965
4-Chlorotoluene	U		0.114	1.00	1	02/22/2024 05:07	WG2230965
1,2-Dibromo-3-Chloropropane	U		0.276	5.00	1	02/22/2024 05:07	WG2230965
1,2-Dibromoethane	U		0.126	1.00	1	02/22/2024 05:07	WG2230965
Dibromomethane	U		0.122	1.00	1	02/22/2024 05:07	WG2230965
1,2-Dichlorobenzene	U		0.107	1.00	1	02/22/2024 05:07	WG2230965
1,3-Dichlorobenzene	U		0.110	1.00	1	02/22/2024 05:07	WG2230965
1,4-Dichlorobenzene	U		0.120	1.00	1	02/22/2024 05:07	WG2230965
Dichlorodifluoromethane	U		0.374	5.00	1	02/22/2024 05:07	WG2230965
1,1-Dichloroethane	U		0.100	1.00	1	02/22/2024 05:07	WG2230965
1,2-Dichloroethane	U		0.0819	1.00	1	02/22/2024 05:07	WG2230965
1,1-Dichloroethene	U		0.188	1.00	1	02/22/2024 05:07	WG2230965
cis-1,2-Dichloroethene	U		0.126	1.00	1	02/22/2024 05:07	WG2230965
trans-1,2-Dichloroethene	U		0.149	1.00	1	02/22/2024 05:07	WG2230965
1,2-Dichloropropane	U		0.149	1.00	1	02/22/2024 05:07	WG2230965
1,1-Dichloropropene	U		0.142	1.00	1	02/22/2024 05:07	WG2230965
1,3-Dichloropropane	U		0.110	1.00	1	02/22/2024 05:07	WG2230965
cis-1,3-Dichloropropene	U		0.111	1.00	1	02/22/2024 05:07	WG2230965
trans-1,3-Dichloropropene	U		0.118	1.00	1	02/22/2024 05:07	WG2230965
2,2-Dichloropropane	U		0.161	1.00	1	02/22/2024 05:07	WG2230965
Dicyclopentadiene	U		0.253	1.00	1	02/22/2024 05:07	WG2230965
Di-isopropyl ether	U		0.105	1.00	1	02/22/2024 05:07	WG2230965
Ethylbenzene	U		0.137	1.00	1	02/22/2024 05:07	WG2230965
4-Ethyltoluene	U		0.208	1.00	1	02/22/2024 05:07	WG2230965
Hexachloro-1,3-butadiene	U		0.337	1.00	1	02/22/2024 05:07	WG2230965
n-Hexane	U		0.749	10.0	1	02/22/2024 05:07	WG2230965
Isopropylbenzene	U		0.105	1.00	1	02/22/2024 05:07	WG2230965
p-Isopropyltoluene	U		0.120	1.00	1	02/22/2024 05:07	WG2230965
2-Butanone (MEK)	U		1.19	10.0	1	02/22/2024 05:07	WG2230965
Methyl Cyclohexane	U		0.660	1.00	1	02/22/2024 05:07	WG2230965

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Methylene Chloride	U		0.430	5.00	1	02/22/2024 05:07	WG2230965
4-Methyl-2-pentanone (MIBK)	U		0.478	10.0	1	02/22/2024 05:07	WG2230965
Methyl tert-butyl ether	U		0.101	1.00	1	02/22/2024 05:07	WG2230965
Naphthalene	1.20	E4	1.00	5.00	1	02/22/2024 05:07	WG2230965
Propene	U		0.936	2.50	1	02/22/2024 05:07	WG2230965
n-Propylbenzene	U		0.0993	1.00	1	02/22/2024 05:07	WG2230965
Styrene	U		0.118	1.00	1	02/22/2024 05:07	WG2230965
1,1,1,2-Tetrachloroethane	U		0.147	1.00	1	02/22/2024 05:07	WG2230965
1,1,2,2-Tetrachloroethane	U		0.133	1.00	1	02/22/2024 05:07	WG2230965
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00	1	02/22/2024 05:07	WG2230965
Tetrachloroethene	U		0.300	1.00	1	02/22/2024 05:07	WG2230965
Toluene	U		0.278	1.00	1	02/22/2024 05:07	WG2230965
1,2,3-Trichlorobenzene	U		0.230	1.00	1	02/22/2024 05:07	WG2230965
1,2,4-Trichlorobenzene	U		0.481	1.00	1	02/22/2024 05:07	WG2230965
1,1,1-Trichloroethane	U		0.149	1.00	1	02/22/2024 05:07	WG2230965
1,1,2-Trichloroethane	U		0.158	1.00	1	02/22/2024 05:07	WG2230965
Trichloroethene	U		0.190	1.00	1	02/22/2024 05:07	WG2230965
Trichlorofluoromethane	U		0.160	5.00	1	02/22/2024 05:07	WG2230965
1,2,3-Trichloropropane	U		0.237	2.50	1	02/22/2024 05:07	WG2230965
1,2,4-Trimethylbenzene	U		0.322	1.00	1	02/22/2024 05:07	WG2230965
1,2,3-Trimethylbenzene	U		0.104	1.00	1	02/22/2024 05:07	WG2230965
1,3,5-Trimethylbenzene	U		0.104	1.00	1	02/22/2024 05:07	WG2230965
Vinyl chloride	U		0.234	1.00	1	02/22/2024 05:07	WG2230965
Xylenes, Total	U		0.174	3.00	1	02/22/2024 05:07	WG2230965
(S) Toluene-d8	*105*			*80.0-120*		*02/22/2024 05:07*	WG2230965
(S) 4-Bromofluorobenzene	*96.4*			*77.0-126*		*02/22/2024 05:07*	WG2230965
(S) 1,2-Dichloroethane-d4	*97.2*			*70.0-130*		*02/22/2024 05:07*	WG2230965

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
1,4-Dioxane	U		0.597	3.00	1	02/22/2024 06:13	WG2231221
(S) Toluene-d8	*96.1*			*77.0-127*		*02/22/2024 06:13*	WG2231221

Method Blank (MB)

(MB) R4042116-1 02/29/24 13:52

Analyte	MB Result	MB Qualifier	MB MDL	MB RDL
	ug/l		ug/l	ug/l
Perchlorate	U		0.300	4.00

Method Blank (MB)

(MB) R4042116-2 02/29/24 14:50

Analyte	MB Result	MB Qualifier	MB MDL	MB RDL
	ug/l		ug/l	ug/l
Perchlorate	U		0.300	4.00

Method Blank (MB)

(MB) R4042116-4 02/29/24 16:14

Analyte	MB Result	MB Qualifier	MB MDL	MB RDL
	ug/l		ug/l	ug/l
Perchlorate	U		0.300	4.00

Method Blank (MB)

(MB) R4042116-14 03/01/24 10:15

Analyte	MB Result	MB Qualifier	MB MDL	MB RDL
	ug/l		ug/l	ug/l
Perchlorate	U		0.300	4.00

L1706680-03 Original Sample (OS) • Duplicate (DUP)

(OS) L1706680-03 03/01/24 00:29 • (DUP) R4042116-5 03/01/24 00:57

Analyte	Original Result	DUP Result	Dilution	DUP RPD	DUP Qualifier	DUP RPD Limits
	ug/l	ug/l		%		%
Perchlorate	10100	12700	1000	22.6	R8	15

L1706680-06 Original Sample (OS) • Duplicate (DUP)

(OS) L1706680-06 03/01/24 03:16 • (DUP) R4042116-6 03/01/24 03:44

Analyte	Original Result	DUP Result	Dilution	DUP RPD	DUP Qualifier	DUP RPD Limits
	ug/l	ug/l		%		%
Perchlorate	5.54	5.53	1	0.255		15

Laboratory Control Sample (LCS)

(LCS) R4042116-3 02/29/24 15:46

Analyte	Spike Amount ug/l	LCS Result ug/l	LCS Rec. %	Rec. Limits %	LCS Qualifier
Perchlorate	10.0	9.65	96.5	90.0-110	

L1706682-02 Original Sample (OS) • Matrix Spike (MS)

(OS) L1706682-02 03/01/24 05:36 • (MS) R4042116-7 03/01/24 06:04

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MS Rec. %	Dilution	Rec. Limits %	MS Qualifier
Perchlorate	10.0	U	10.2	102	1	80.0-120	

L1705776-23 Original Sample (OS) • Matrix Spike (MS)

(OS) L1705776-23 02/29/24 18:06 • (MS) R4042116-8 03/01/24 06:32

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MS Rec. %	Dilution	Rec. Limits %	MS Qualifier
Perchlorate	10.0	U	7.94	79.4	1	80.0-120	M2

L1706680-03 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1706680-03 03/01/24 00:29 • (MS) R4042116-9 03/01/24 07:00 • (MSD) R4042116-10 03/01/24 08:23

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MSD Result ug/l	MS Rec. %	MSD Rec. %	Dilution	Rec. Limits %	MS Qualifier	MSD Qualifier	RPD %	RPD Limits %
Perchlorate	10.0	10100	20100	22600	99800	125000	1000	80.0-120	M3	M3	11.9	15

L1706680-06 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1706680-06 03/01/24 03:16 • (MS) R4042116-11 03/01/24 08:51 • (MSD) R4042116-12 03/01/24 09:19

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MSD Result ug/l	MS Rec. %	MSD Rec. %	Dilution	Rec. Limits %	MS Qualifier	MSD Qualifier	RPD %	RPD Limits %
Perchlorate	10.0	5.54	14.9	15.5	93.3	99.5	1	80.0-120			4.09	15

L1706680-08 Original Sample (OS) • Matrix Spike (MS)

(OS) L1706680-08 03/01/24 04:40 • (MS) R4042116-13 03/01/24 09:47

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MS Rec. %	Dilution	Rec. Limits %	MS Qualifier
Perchlorate	10.0	4.55	13.7	91.8	1	80.0-120	

Method Blank (MB)

(MB) R4042154-1 03/03/24 20:12

Analyte	MB Result ug/l	MB Qualifier	MB MDL ug/l	MB RDL ug/l
Perchlorate	U		0.300	4.00

Method Blank (MB)

(MB) R4042154-3 03/03/24 21:36

Analyte	MB Result ug/l	MB Qualifier	MB MDL ug/l	MB RDL ug/l
Perchlorate	U		0.300	4.00

L1708086-06 Original Sample (OS) • Duplicate (DUP)

(OS) L1708086-06 03/04/24 05:58 • (DUP) R4042154-4 03/04/24 07:22

Analyte	Original Result ug/l	DUP Result ug/l	Dilution	DUP RPD %	DUP Qualifier	DUP RPD Limits %
Perchlorate	U	U	1	0.000		15

L1708086-07 Original Sample (OS) • Duplicate (DUP)

(OS) L1708086-07 03/04/24 08:46 • (DUP) R4042154-7 03/04/24 09:14

Analyte	Original Result ug/l	DUP Result ug/l	Dilution	DUP RPD %	DUP Qualifier	DUP RPD Limits %
Perchlorate	U	U	1	0.000		15

Laboratory Control Sample (LCS)

(LCS) R4042154-2 03/03/24 21:08

Analyte	Spike Amount ug/l	LCS Result ug/l	LCS Rec. %	Rec. Limits %	LCS Qualifier
Perchlorate	10.0	9.38	93.8	90.0-110	

L1708086-06 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1708086-06 03/04/24 05:58 • (MS) R4042154-5 03/04/24 07:50 • (MSD) R4042154-6 03/04/24 08:18

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MSD Result ug/l	MS Rec. %	MSD Rec. %	Dilution	Rec. Limits %	MS Qualifier	MSD Qualifier	RPD %	RPD Limits %
Perchlorate	10.0	U	4.42	6.73	44.2	67.3	1	80.0-120	M2	M2 R5	41.4	15

L1708086-06 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1708086-06 03/04/24 05:58 • (MS) R4042154-5 03/04/24 07:50 • (MSD) R4042154-6 03/04/24 08:18

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MSD Result ug/l	MS Rec. %	MSD Rec. %	Dilution	Rec. Limits %	MS Qualifier	MSD Qualifier	RPD %	RPD Limits %

Sample Narrative:

MS: spike failed due to sample matrix

MSD: spike failed due to sample matrix

L1708086-07 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1708086-07 03/04/24 08:46 • (MS) R4042154-8 03/04/24 09:42 • (MSD) R4042154-9 03/04/24 10:10

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MSD Result ug/l	MS Rec. %	MSD Rec. %	Dilution	Rec. Limits %	MS Qualifier	MSD Qualifier	RPD %	RPD Limits %
Perchlorate	10.0	U	8.65	9.01	86.5	90.1	1	80.0-120			4.05	15

Method Blank (MB)

(MB) R4036492-4 02/20/24 22:46

Analyte	MB Result ug/l	MB Qualifier	MB MDL ug/l	MB RDL ug/l
Acetone	U		11.3	50.0
Acrolein	U		2.54	50.0
Acrylonitrile	U		0.671	10.0
Benzene	U		0.0941	1.00
Bromobenzene	U		0.118	1.00
Bromodichloromethane	U		0.136	1.00
Bromoform	U		0.129	1.00
Bromomethane	U		0.605	5.00
1,3-Butadiene	U		0.299	2.00
n-Butylbenzene	U		0.157	1.00
sec-Butylbenzene	U		0.125	1.00
tert-Butylbenzene	U		0.127	1.00
Carbon tetrachloride	U		0.128	1.00
Carbon disulfide	0.156	E4	0.0962	1.00
Chlorobenzene	U		0.116	1.00
Chlorodibromomethane	U		0.140	1.00
Chloroethane	U		0.192	5.00
Chloroform	U		0.111	5.00
Chloromethane	U		0.960	2.50
Cyclohexane	U		0.188	1.00
2-Chlorotoluene	U		0.106	1.00
4-Chlorotoluene	U		0.114	1.00
1,2-Dibromo-3-Chloropropane	U		0.276	5.00
1,2-Dibromoethane	U		0.126	1.00
Dibromomethane	U		0.122	1.00
1,2-Dichlorobenzene	U		0.107	1.00
1,3-Dichlorobenzene	U		0.110	1.00
1,4-Dichlorobenzene	U		0.120	1.00
Dichlorodifluoromethane	U		0.374	5.00
1,1-Dichloroethane	U		0.100	1.00
1,2-Dichloroethane	U		0.0819	1.00
1,1-Dichloroethene	U		0.188	1.00
cis-1,2-Dichloroethene	U		0.126	1.00
trans-1,2-Dichloroethene	U		0.149	1.00
1,2-Dichloropropane	U		0.149	1.00
1,1-Dichloropropene	U		0.142	1.00
1,3-Dichloropropane	U		0.110	1.00
cis-1,3-Dichloropropene	U		0.111	1.00
trans-1,3-Dichloropropene	U		0.118	1.00
2,2-Dichloropropane	U		0.161	1.00

Method Blank (MB)

(MB) R4036492-4 02/20/24 22:46

Analyte	MB Result ug/l	MB Qualifier	MB MDL ug/l	MB RDL ug/l
Dicyclopentadiene	U		0.253	1.00
Di-isopropyl ether	U		0.105	1.00
Ethylbenzene	U		0.137	1.00
4-Ethyltoluene	U		0.208	1.00
Hexachloro-1,3-butadiene	U		0.337	1.00
n-Hexane	U		0.749	10.0
Isopropylbenzene	U		0.105	1.00
p-Isopropyltoluene	U		0.120	1.00
2-Butanone (MEK)	U		1.19	10.0
Methyl Cyclohexane	U		0.660	1.00
Methylene Chloride	U		0.430	5.00
4-Methyl-2-pentanone (MIBK)	U		0.478	10.0
Methyl tert-butyl ether	U		0.101	1.00
Naphthalene	U		1.00	5.00
Propene	U		0.936	2.50
n-Propylbenzene	U		0.0993	1.00
Styrene	U		0.118	1.00
1,1,1,2-Tetrachloroethane	U		0.147	1.00
1,1,2,2-Tetrachloroethane	U		0.133	1.00
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00
Tetrachloroethene	U		0.300	1.00
Toluene	U		0.278	1.00
1,2,3-Trichlorobenzene	U		0.230	1.00
1,2,4-Trichlorobenzene	U		0.481	1.00
1,1,1-Trichloroethane	U		0.149	1.00
1,1,2-Trichloroethane	U		0.158	1.00
Trichloroethene	U		0.190	1.00
Trichlorofluoromethane	U		0.160	5.00
1,2,3-Trichloropropane	U		0.237	2.50
1,2,4-Trimethylbenzene	U		0.322	1.00
1,2,3-Trimethylbenzene	U		0.104	1.00
1,3,5-Trimethylbenzene	U		0.104	1.00
Vinyl chloride	U		0.234	1.00
Xylenes, Total	U		0.174	3.00
(S) Toluene-d8	*111*			*80.0-120*
(S) 4-Bromofluorobenzene	*105*			*77.0-126*
(S) 1,2-Dichloroethane-d4	*122*			*70.0-130*

QUALITY CONTROL SUMMARY

L1706682-01,02,03

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R4036492-1 02/20/24 21:20 • (LCSD) R4036492-2 02/20/24 21:42

Analyte	Spike Amount ug/l	LCS Result ug/l	LCSD Result ug/l	LCS Rec. %	LCSD Rec. %	Rec. Limits %	LCS Qualifier	LCSD Qualifier	RPD %	RPD Limits %
Acetone	25.0	21.5	16.4	86.0	65.6	19.0-160			26.9	27
Acrolein	25.0	21.8	24.6	87.2	98.4	30.0-160			12.1	26
Acrylonitrile	25.0	29.9	27.8	120	111	55.0-149			7.28	20
Benzene	5.00	4.64	4.62	92.8	92.4	70.0-123			0.432	20
Bromobenzene	5.00	4.53	4.47	90.6	89.4	73.0-121			1.33	20
Bromodichloromethane	5.00	4.64	4.28	92.8	85.6	75.0-120			8.07	20
Bromoform	5.00	3.73	3.49	74.6	69.8	68.0-132			6.65	20
Bromomethane	5.00	3.38	4.38	67.6	87.6	30.0-160		R7	25.8	25
1,3-Butadiene	5.00	3.74	4.78	74.8	95.6	45.0-147		R7	24.4	20
n-Butylbenzene	5.00	4.71	4.67	94.2	93.4	73.0-125			0.853	20
sec-Butylbenzene	5.00	4.58	4.33	91.6	86.6	75.0-125			5.61	20
tert-Butylbenzene	5.00	4.79	4.50	95.8	90.0	76.0-124			6.24	20
Carbon tetrachloride	5.00	4.84	4.78	96.8	95.6	68.0-126			1.25	20
Carbon disulfide	5.00	3.14	3.96	62.8	79.2	61.0-128		R7	23.1	20
Chlorobenzene	5.00	4.94	4.58	98.8	91.6	80.0-121			7.56	20
Chlorodibromomethane	5.00	4.38	4.00	87.6	80.0	77.0-125			9.07	20
Chloroethane	5.00	4.73	5.14	94.6	103	47.0-150			8.31	20
Chloroform	5.00	4.76	4.74	95.2	94.8	73.0-120			0.421	20
Chloromethane	5.00	3.85	4.96	77.0	99.2	41.0-142		R7	25.2	20
Cyclohexane	5.00	4.07	4.40	81.4	88.0	71.0-124			7.79	20
2-Chlorotoluene	5.00	4.64	4.38	92.8	87.6	76.0-123			5.76	20
4-Chlorotoluene	5.00	4.55	4.46	91.0	89.2	75.0-122			2.00	20
1,2-Dibromo-3-Chloropropane	5.00	3.27	3.05	65.4	61.0	58.0-134			6.96	20
1,2-Dibromoethane	5.00	4.42	4.41	88.4	88.2	80.0-122			0.227	20
Dibromomethane	5.00	4.50	4.37	90.0	87.4	80.0-120			2.93	20
1,2-Dichlorobenzene	5.00	4.82	4.56	96.4	91.2	79.0-121			5.54	20
1,3-Dichlorobenzene	5.00	4.94	4.68	98.8	93.6	79.0-120			5.41	20
1,4-Dichlorobenzene	5.00	4.85	4.78	97.0	95.6	79.0-120			1.45	20
Dichlorodifluoromethane	5.00	3.67	5.31	73.4	106	51.0-149		R7	36.5	20
1,1-Dichloroethane	5.00	5.30	4.98	106	99.6	70.0-126			6.23	20
1,2-Dichloroethane	5.00	5.19	5.28	104	106	70.0-128			1.72	20
1,1-Dichloroethene	5.00	4.17	4.28	83.4	85.6	71.0-124			2.60	20
cis-1,2-Dichloroethene	5.00	4.71	4.60	94.2	92.0	73.0-120			2.36	20
trans-1,2-Dichloroethene	5.00	4.39	4.46	87.8	89.2	73.0-120			1.58	20
1,2-Dichloropropane	5.00	5.40	5.26	108	105	77.0-125			2.63	20
1,1-Dichloropropene	5.00	4.72	4.76	94.4	95.2	74.0-126			0.844	20
1,3-Dichloropropane	5.00	4.92	4.75	98.4	95.0	80.0-120			3.52	20
cis-1,3-Dichloropropene	5.00	5.00	4.67	100	93.4	80.0-123			6.83	20
trans-1,3-Dichloropropene	5.00	5.05	4.79	101	95.8	78.0-124			5.28	20
2,2-Dichloropropane	5.00	5.21	4.79	104	95.8	58.0-130			8.40	20

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R4036492-1 02/20/24 21:20 • (LCSD) R4036492-2 02/20/24 21:42

Analyte	Spike Amount ug/l	LCS Result ug/l	LCSD Result ug/l	LCS Rec. %	LCSD Rec. %	Rec. Limits %	LCS Qualifier	LCSD Qualifier	RPD %	RPD Limits %
Dicyclopentadiene	5.00	4.70	4.43	94.0	88.6	74.0-126			5.91	20
Di-isopropyl ether	5.00	6.38	6.08	128	122	58.0-138			4.82	20
Ethylbenzene	5.00	4.77	4.74	95.4	94.8	79.0-123			0.631	20
4-Ethyltoluene	5.00	4.75	4.51	95.0	90.2	74.0-127			5.18	20
Hexachloro-1,3-butadiene	5.00	4.38	4.22	87.6	84.4	54.0-138			3.72	20
n-Hexane	5.00	4.25	5.09	85.0	102	57.0-133			18.0	20
Isopropylbenzene	5.00	4.95	4.63	99.0	92.6	76.0-127			6.68	20
p-Isopropyltoluene	5.00	4.79	4.52	95.8	90.4	76.0-125			5.80	20
2-Butanone (MEK)	25.0	24.6	22.7	98.4	90.8	44.0-160			8.03	20
Methyl Cyclohexane	5.00	3.82	4.32	76.4	86.4	68.0-126			12.3	20
Methylene Chloride	5.00	4.96	4.34	99.2	86.8	67.0-120			13.3	20
4-Methyl-2-pentanone (MIBK)	25.0	32.5	30.7	130	123	68.0-142			5.70	20
Methyl tert-butyl ether	5.00	4.80	4.57	96.0	91.4	68.0-125			4.91	20
Naphthalene	5.00	2.99	2.88	59.8	57.6	54.0-135			3.75	20
Propene	5.00	2.38	4.23	47.6	84.6	30.0-160		R7	56.0	20
n-Propylbenzene	5.00	4.63	4.40	92.6	88.0	77.0-124			5.09	20
Styrene	5.00	4.65	4.30	93.0	86.0	73.0-130			7.82	20
1,1,1,2-Tetrachloroethane	5.00	4.68	4.38	93.6	87.6	75.0-125			6.62	20
1,1,2,2-Tetrachloroethane	5.00	4.37	4.38	87.4	87.6	65.0-130			0.229	20
1,1,2-Trichlorotrifluoroethane	5.00	4.61	4.75	92.2	95.0	69.0-132			2.99	20
Tetrachloroethene	5.00	4.92	4.88	98.4	97.6	72.0-132			0.816	20
Toluene	5.00	4.89	4.77	97.8	95.4	79.0-120			2.48	20
1,2,3-Trichlorobenzene	5.00	3.65	3.64	73.0	72.8	50.0-138			0.274	20
1,2,4-Trichlorobenzene	5.00	3.80	3.56	76.0	71.2	57.0-137			6.52	20
1,1,1-Trichloroethane	5.00	4.82	4.48	96.4	89.6	73.0-124			7.31	20
1,1,2-Trichloroethane	5.00	4.62	4.40	92.4	88.0	80.0-120			4.88	20
Trichloroethene	5.00	5.14	4.71	103	94.2	78.0-124			8.73	20
Trichlorofluoromethane	5.00	4.92	5.22	98.4	104	59.0-147			5.92	20
1,2,3-Trichloropropane	5.00	4.56	4.23	91.2	84.6	73.0-130			7.51	20
1,2,4-Trimethylbenzene	5.00	4.46	4.23	89.2	84.6	76.0-121			5.29	20
1,2,3-Trimethylbenzene	5.00	4.31	4.24	86.2	84.8	77.0-120			1.64	20
1,3,5-Trimethylbenzene	5.00	4.60	4.36	92.0	87.2	76.0-122			5.36	20
Vinyl chloride	5.00	4.42	5.20	88.4	104	67.0-131			16.2	20
Xylenes, Total	15.0	14.3	13.7	95.3	91.3	79.0-123			4.29	20
(S) Toluene-d8				*111*	*111*	*80.0-120*				
(S) 4-Bromofluorobenzene				*105*	*105*	*77.0-126*				
(S) 1,2-Dichloroethane-d4				*124*	*126*	*70.0-130*				

L1706682-01 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1706682-01 02/21/24 05:43 • (MS) R4036492-6 02/21/24 08:36 • (MSD) R4036492-7 02/21/24 08:58

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MSD Result ug/l	MS Rec. %	MSD Rec. %	Dilution	Rec. Limits %	MS Qualifier	MSD Qualifier	RPD %	RPD Limits %
Acetone	25.0	U	21.5	21.9	86.0	87.6	1	10.0-160			1.84	35
Acrolein	25.0	U	34.9	34.1	140	136	1	10.0-160			2.32	39
Acrylonitrile	25.0	U	38.1	36.8	152	147	1	21.0-160			3.47	32
Benzene	5.00	U	6.15	5.84	123	117	1	17.0-158			5.17	27
Bromobenzene	5.00	U	5.98	5.67	120	113	1	30.0-149			5.32	28
Bromodichloromethane	5.00	U	5.66	5.73	113	115	1	31.0-150			1.23	27
Bromoform	5.00	U	4.73	4.23	94.6	84.6	1	29.0-150			11.2	29
Bromomethane	5.00	U	4.70	4.87	94.0	97.4	1	10.0-160			3.55	38
1,3-Butadiene	5.00	U	6.16	5.85	123	117	1	10.0-160			5.16	22
n-Butylbenzene	5.00	U	6.52	5.96	130	119	1	31.0-150			8.97	30
sec-Butylbenzene	5.00	U	6.30	6.03	126	121	1	33.0-155			4.38	29
tert-Butylbenzene	5.00	U	6.40	6.13	128	123	1	34.0-153			4.31	28
Carbon tetrachloride	5.00	U	6.40	6.16	128	123	1	23.0-159			3.82	28
Carbon disulfide	5.00	0.172	3.61	3.41	68.8	64.8	1	10.0-156			5.70	28
Chlorobenzene	5.00	U	6.14	6.07	123	121	1	33.0-152			1.15	27
Chlorodibromomethane	5.00	U	5.40	5.18	108	104	1	37.0-149			4.16	27
Chloroethane	5.00	U	6.23	5.90	125	118	1	10.0-160			5.44	30
Chloroform	5.00	U	5.94	5.90	119	118	1	29.0-154			0.676	28
Chloromethane	5.00	U	4.74	4.31	94.8	86.2	1	10.0-160			9.50	29
Cyclohexane	5.00	U	5.61	5.30	112	106	1	19.0-160			5.68	23
2-Chlorotoluene	5.00	U	6.24	5.86	125	117	1	32.0-153			6.28	28
4-Chlorotoluene	5.00	U	6.15	5.79	123	116	1	32.0-150			6.03	28
1,2-Dibromo-3-Chloropropane	5.00	U	4.34	4.21	86.8	84.2	1	22.0-151			3.04	34
1,2-Dibromoethane	5.00	U	5.65	5.36	113	107	1	34.0-147			5.27	27
Dibromomethane	5.00	U	5.60	5.54	112	111	1	30.0-151			1.08	27
1,2-Dichlorobenzene	5.00	U	6.44	6.16	129	123	1	34.0-149			4.44	28
1,3-Dichlorobenzene	5.00	U	6.49	6.19	130	124	1	36.0-146			4.73	27
1,4-Dichlorobenzene	5.00	U	6.54	6.21	131	124	1	35.0-142			5.18	27
Dichlorodifluoromethane	5.00	U	5.48	4.97	110	99.4	1	10.0-160			9.76	29
1,1-Dichloroethane	5.00	U	6.68	6.58	134	132	1	25.0-158			1.51	27
1,2-Dichloroethane	5.00	U	6.67	6.58	133	132	1	29.0-151			1.36	27
1,1-Dichloroethene	5.00	U	5.46	5.15	109	103	1	11.0-160			5.84	29
cis-1,2-Dichloroethene	5.00	U	5.96	5.86	119	117	1	10.0-160			1.69	27
trans-1,2-Dichloroethene	5.00	U	5.33	5.14	107	103	1	17.0-153			3.63	27
1,2-Dichloropropane	5.00	U	6.91	6.55	138	131	1	30.0-156			5.35	27
1,1-Dichloropropene	5.00	U	6.06	5.83	121	117	1	25.0-158			3.87	27
1,3-Dichloropropane	5.00	U	6.25	6.05	125	121	1	38.0-147			3.25	27
cis-1,3-Dichloropropene	5.00	U	5.94	5.66	119	113	1	34.0-149			4.83	28
trans-1,3-Dichloropropene	5.00	U	6.04	5.92	121	118	1	32.0-149			2.01	28
2,2-Dichloropropane	5.00	U	6.20	6.03	124	121	1	24.0-152			2.78	29

L1706682-01 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1706682-01 02/21/24 05:43 • (MS) R4036492-6 02/21/24 08:36 • (MSD) R4036492-7 02/21/24 08:58

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MSD Result ug/l	MS Rec. %	MSD Rec. %	Dilution	Rec. Limits %	MS Qualifier	MSD Qualifier	RPD %	RPD Limits %
Dicyclopentadiene	5.00	U	6.01	5.77	120	115	1	51.0-139			4.07	20
Di-isopropyl ether	5.00	U	8.32	8.08	166	162	1	21.0-160	M1	M1	2.93	28
Ethylbenzene	5.00	U	6.42	6.16	128	123	1	30.0-155			4.13	27
4-Ethyltoluene	5.00	U	6.78	6.09	136	122	1	10.0-160			10.7	20
Hexachloro-1,3-butadiene	5.00	U	5.70	5.52	114	110	1	20.0-154			3.21	34
n-Hexane	5.00	U	5.45	5.28	109	106	1	10.0-153			3.17	28
Isopropylbenzene	5.00	U	6.52	6.36	130	127	1	28.0-157			2.48	27
p-Isopropyltoluene	5.00	U	6.53	6.05	131	121	1	30.0-154			7.63	29
2-Butanone (MEK)	25.0	U	31.2	30.5	125	122	1	10.0-160			2.27	32
Methyl Cyclohexane	5.00	U	5.41	5.16	108	103	1	11.0-160			4.73	24
Methylene Chloride	5.00	U	5.66	5.50	113	110	1	23.0-144			2.87	28
4-Methyl-2-pentanone (MIBK)	25.0	U	42.3	40.2	169	161	1	29.0-160	M1	M1	5.09	29
Methyl tert-butyl ether	5.00	U	6.16	6.04	123	121	1	28.0-150			1.97	29
Naphthalene	5.00	U	5.11	4.14	102	82.8	1	12.0-156			21.0	35
Propene	5.00	U	3.84	3.95	76.8	79.0	1	10.0-160			2.82	29
n-Propylbenzene	5.00	U	6.17	5.82	123	116	1	31.0-154			5.84	28
Styrene	5.00	U	6.00	5.73	120	115	1	33.0-155			4.60	28
1,1,1,2-Tetrachloroethane	5.00	U	6.00	5.70	120	114	1	36.0-151			5.13	29
1,1,2,2-Tetrachloroethane	5.00	U	5.81	5.67	116	113	1	33.0-150			2.44	28
1,1,2-Trichlorotrifluoroethane	5.00	U	6.48	6.13	130	123	1	23.0-160			5.55	30
Tetrachloroethene	5.00	U	6.33	6.11	127	122	1	10.0-160			3.54	27
Toluene	5.00	U	6.60	6.12	132	122	1	26.0-154			7.55	28
1,2,3-Trichlorobenzene	5.00	U	5.13	4.88	103	97.6	1	17.0-150			5.00	36
1,2,4-Trichlorobenzene	5.00	U	5.17	4.96	103	99.2	1	24.0-150			4.15	33
1,1,1-Trichloroethane	5.00	U	6.24	6.01	125	120	1	23.0-160			3.76	28
1,1,2-Trichloroethane	5.00	U	5.78	5.51	116	110	1	35.0-147			4.78	27
Trichloroethene	5.00	U	6.02	5.90	120	118	1	10.0-160			2.01	25
Trichlorofluoromethane	5.00	U	6.89	6.55	138	131	1	17.0-160			5.06	31
1,2,3-Trichloropropane	5.00	U	5.51	5.27	110	105	1	34.0-151			4.45	29
1,2,4-Trimethylbenzene	5.00	U	7.02	6.02	140	120	1	26.0-154			15.3	27
1,2,3-Trimethylbenzene	5.00	U	5.97	5.50	119	110	1	32.0-149			8.20	28
1,3,5-Trimethylbenzene	5.00	U	6.36	5.82	127	116	1	28.0-153			8.87	27
Vinyl chloride	5.00	U	6.24	5.92	125	118	1	10.0-160			5.26	27
Xylenes, Total	15.0	U	19.7	18.3	131	122	1	29.0-154			7.37	28
(S) Toluene-d8					*107*	*108*		*80.0-120*				
(S) 4-Bromofluorobenzene					*106*	*108*		*77.0-126*				
(S) 1,2-Dichloroethane-d4					*124*	*125*		*70.0-130*				

Method Blank (MB)

(MB) R4036917-4 02/21/24 21:14

Analyte	MB Result ug/l	MB Qualifier	MB MDL ug/l	MB RDL ug/l
Acetone	U		11.3	50.0
Acrolein	U		2.54	50.0
Acrylonitrile	U		0.671	10.0
Benzene	U		0.0941	1.00
Bromobenzene	U		0.118	1.00
Bromodichloromethane	U		0.136	1.00
Bromoform	U		0.129	1.00
Bromomethane	U		0.605	5.00
1,3-Butadiene	U		0.299	2.00
n-Butylbenzene	U		0.157	1.00
sec-Butylbenzene	U		0.125	1.00
tert-Butylbenzene	U		0.127	1.00
Carbon tetrachloride	U		0.128	1.00
Carbon disulfide	U		0.0962	1.00
Chlorobenzene	U		0.116	1.00
Chlorodibromomethane	U		0.140	1.00
Chloroethane	U		0.192	5.00
Chloroform	U		0.111	5.00
Chloromethane	U		0.960	2.50
Cyclohexane	U		0.188	1.00
2-Chlorotoluene	U		0.106	1.00
4-Chlorotoluene	U		0.114	1.00
1,2-Dibromo-3-Chloropropane	U		0.276	5.00
1,2-Dibromoethane	U		0.126	1.00
Dibromomethane	U		0.122	1.00
1,2-Dichlorobenzene	U		0.107	1.00
1,3-Dichlorobenzene	U		0.110	1.00
1,4-Dichlorobenzene	U		0.120	1.00
Dichlorodifluoromethane	U		0.374	5.00
1,1-Dichloroethane	U		0.100	1.00
1,2-Dichloroethane	U		0.0819	1.00
1,1-Dichloroethene	U		0.188	1.00
cis-1,2-Dichloroethene	U		0.126	1.00
trans-1,2-Dichloroethene	U		0.149	1.00
1,2-Dichloropropane	U		0.149	1.00
1,1-Dichloropropene	U		0.142	1.00
1,3-Dichloropropane	U		0.110	1.00
cis-1,3-Dichloropropene	U		0.111	1.00
trans-1,3-Dichloropropene	U		0.118	1.00
2,2-Dichloropropane	U		0.161	1.00

Method Blank (MB)

(MB) R4036917-4 02/21/24 21:14

Analyte	MB Result ug/l	MB Qualifier	MB MDL ug/l	MB RDL ug/l
Dicyclopentadiene	U		0.253	1.00
Di-isopropyl ether	U		0.105	1.00
Ethylbenzene	U		0.137	1.00
4-Ethyltoluene	U		0.208	1.00
Hexachloro-1,3-butadiene	U		0.337	1.00
n-Hexane	U		0.749	10.0
Isopropylbenzene	U		0.105	1.00
p-Isopropyltoluene	U		0.120	1.00
2-Butanone (MEK)	U		1.19	10.0
Methyl Cyclohexane	U		0.660	1.00
Methylene Chloride	U		0.430	5.00
4-Methyl-2-pentanone (MIBK)	U		0.478	10.0
Methyl tert-butyl ether	U		0.101	1.00
Naphthalene	U		1.00	5.00
Propene	U		0.936	2.50
n-Propylbenzene	U		0.0993	1.00
Styrene	U		0.118	1.00
1,1,1,2-Tetrachloroethane	U		0.147	1.00
1,1,2,2-Tetrachloroethane	U		0.133	1.00
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00
Tetrachloroethene	U		0.300	1.00
Toluene	U		0.278	1.00
1,2,3-Trichlorobenzene	U		0.230	1.00
1,2,4-Trichlorobenzene	U		0.481	1.00
1,1,1-Trichloroethane	U		0.149	1.00
1,1,2-Trichloroethane	U		0.158	1.00
Trichloroethene	U		0.190	1.00
Trichlorofluoromethane	U		0.160	5.00
1,2,3-Trichloropropane	U		0.237	2.50
1,2,4-Trimethylbenzene	U		0.322	1.00
1,2,3-Trimethylbenzene	U		0.104	1.00
1,3,5-Trimethylbenzene	U		0.104	1.00
Vinyl chloride	U		0.234	1.00
Xylenes, Total	U		0.174	3.00
(S) Toluene-d8	*104*			*80.0-120*
(S) 4-Bromofluorobenzene	*95.6*			*77.0-126*
(S) 1,2-Dichloroethane-d4	*101*			*70.0-130*

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R4036917-1 02/21/24 19:33 • (LCSD) R4036917-2 02/21/24 19:53

Analyte	Spike Amount ug/l	LCS Result ug/l	LCSD Result ug/l	LCS Rec. %	LCSD Rec. %	Rec. Limits %	LCS Qualifier	LCSD Qualifier	RPD %	RPD Limits %
Acetone	25.0	23.4	25.0	93.6	100	19.0-160			6.61	27
Acrolein	25.0	102	99.6	408	398	30.0-160	L1	L1	2.38	26
Acrylonitrile	25.0	25.3	27.3	101	109	55.0-149			7.60	20
Benzene	5.00	4.93	5.21	98.6	104	70.0-123			5.52	20
Bromobenzene	5.00	4.67	4.81	93.4	96.2	73.0-121			2.95	20
Bromodichloromethane	5.00	4.72	4.88	94.4	97.6	75.0-120			3.33	20
Bromoform	5.00	3.87	4.17	77.4	83.4	68.0-132			7.46	20
Bromomethane	5.00	3.34	4.42	66.8	88.4	30.0-160		R7	27.8	25
1,3-Butadiene	5.00	4.13	4.58	82.6	91.6	45.0-147			10.3	20
n-Butylbenzene	5.00	3.86	4.05	77.2	81.0	73.0-125			4.80	20
sec-Butylbenzene	5.00	4.46	4.63	89.2	92.6	75.0-125			3.74	20
tert-Butylbenzene	5.00	4.31	4.48	86.2	89.6	76.0-124			3.87	20
Carbon tetrachloride	5.00	4.69	5.42	93.8	108	68.0-126			14.4	20
Carbon disulfide	5.00	4.61	5.22	92.2	104	61.0-128			12.4	20
Chlorobenzene	5.00	4.42	4.68	88.4	93.6	80.0-121			5.71	20
Chlorodibromomethane	5.00	4.08	4.38	81.6	87.6	77.0-125			7.09	20
Chloroethane	5.00	5.13	5.38	103	108	47.0-150			4.76	20
Chloroform	5.00	4.98	5.12	99.6	102	73.0-120			2.77	20
Chloromethane	5.00	5.31	6.04	106	121	41.0-142			12.9	20
Cyclohexane	5.00	4.66	4.93	93.2	98.6	71.0-124			5.63	20
2-Chlorotoluene	5.00	4.65	4.88	93.0	97.6	76.0-123			4.83	20
4-Chlorotoluene	5.00	4.30	4.52	86.0	90.4	75.0-122			4.99	20
1,2-Dibromo-3-Chloropropane	5.00	3.62	3.84	72.4	76.8	58.0-134			5.90	20
1,2-Dibromoethane	5.00	4.39	4.52	87.8	90.4	80.0-122			2.92	20
Dibromomethane	5.00	4.72	5.03	94.4	101	80.0-120			6.36	20
1,2-Dichlorobenzene	5.00	4.39	4.54	87.8	90.8	79.0-121			3.36	20
1,3-Dichlorobenzene	5.00	4.35	4.49	87.0	89.8	79.0-120			3.17	20
1,4-Dichlorobenzene	5.00	4.35	4.55	87.0	91.0	79.0-120			4.49	20
Dichlorodifluoromethane	5.00	5.83	6.20	117	124	51.0-149			6.15	20
1,1-Dichloroethane	5.00	4.94	5.26	98.8	105	70.0-126			6.27	20
1,2-Dichloroethane	5.00	5.15	5.38	103	108	70.0-128			4.37	20
1,1-Dichloroethene	5.00	4.96	5.30	99.2	106	71.0-124			6.63	20
cis-1,2-Dichloroethene	5.00	4.63	5.04	92.6	101	73.0-120			8.48	20
trans-1,2-Dichloroethene	5.00	4.77	5.09	95.4	102	73.0-120			6.49	20
1,2-Dichloropropane	5.00	4.95	5.31	99.0	106	77.0-125			7.02	20
1,1-Dichloropropene	5.00	4.93	5.42	98.6	108	74.0-126			9.47	20
1,3-Dichloropropane	5.00	4.58	4.87	91.6	97.4	80.0-120			6.14	20
cis-1,3-Dichloropropene	5.00	4.60	4.85	92.0	97.0	80.0-123			5.29	20
trans-1,3-Dichloropropene	5.00	4.18	4.60	83.6	92.0	78.0-124			9.57	20
2,2-Dichloropropane	5.00	5.03	5.14	101	103	58.0-130			2.16	20

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R4036917-1 02/21/24 19:33 • (LCSD) R4036917-2 02/21/24 19:53

Analyte	Spike Amount ug/l	LCS Result ug/l	LCSD Result ug/l	LCS Rec. %	LCSD Rec. %	Rec. Limits %	LCS Qualifier	LCSD Qualifier	RPD %	RPD Limits %
Dicyclopentadiene	5.00	4.57	4.86	91.4	97.2	74.0-126			6.15	20
Di-isopropyl ether	5.00	4.90	5.17	98.0	103	58.0-138			5.36	20
Ethylbenzene	5.00	4.39	4.65	87.8	93.0	79.0-123			5.75	20
4-Ethyltoluene	5.00	4.39	4.53	87.8	90.6	74.0-127			3.14	20
Hexachloro-1,3-butadiene	5.00	4.31	4.50	86.2	90.0	54.0-138			4.31	20
n-Hexane	5.00	5.07	5.17	101	103	57.0-133			1.95	20
Isopropylbenzene	5.00	4.25	4.59	85.0	91.8	76.0-127			7.69	20
p-Isopropyltoluene	5.00	4.35	4.60	87.0	92.0	76.0-125			5.59	20
2-Butanone (MEK)	25.0	23.2	24.8	92.8	99.2	44.0-160			6.67	20
Methyl Cyclohexane	5.00	4.52	5.02	90.4	100	68.0-126			10.5	20
Methylene Chloride	5.00	4.98	5.16	99.6	103	67.0-120			3.55	20
4-Methyl-2-pentanone (MIBK)	25.0	24.2	25.7	96.8	103	68.0-142			6.01	20
Methyl tert-butyl ether	5.00	4.84	4.99	96.8	99.8	68.0-125			3.05	20
Naphthalene	5.00	3.36	3.84	67.2	76.8	54.0-135			13.3	20
Propene	5.00	3.26	3.79	65.2	75.8	30.0-160			15.0	20
n-Propylbenzene	5.00	4.51	4.79	90.2	95.8	77.0-124			6.02	20
Styrene	5.00	3.97	4.30	79.4	86.0	73.0-130			7.98	20
1,1,1,2-Tetrachloroethane	5.00	4.42	4.56	88.4	91.2	75.0-125			3.12	20
1,1,2,2-Tetrachloroethane	5.00	4.77	4.70	95.4	94.0	65.0-130			1.48	20
1,1,2-Trichlorotrifluoroethane	5.00	4.63	4.73	92.6	94.6	69.0-132			2.14	20
Tetrachloroethene	5.00	4.63	4.82	92.6	96.4	72.0-132			4.02	20
Toluene	5.00	4.43	4.73	88.6	94.6	79.0-120			6.55	20
1,2,3-Trichlorobenzene	5.00	4.35	5.12	87.0	102	50.0-138			16.3	20
1,2,4-Trichlorobenzene	5.00	4.11	4.51	82.2	90.2	57.0-137			9.28	20
1,1,1-Trichloroethane	5.00	4.97	5.28	99.4	106	73.0-124			6.05	20
1,1,2-Trichloroethane	5.00	4.64	4.85	92.8	97.0	80.0-120			4.43	20
Trichloroethene	5.00	4.66	5.31	93.2	106	78.0-124			13.0	20
Trichlorofluoromethane	5.00	5.21	5.47	104	109	59.0-147			4.87	20
1,2,3-Trichloropropane	5.00	4.66	4.89	93.2	97.8	73.0-130			4.82	20
1,2,4-Trimethylbenzene	5.00	4.27	4.47	85.4	89.4	76.0-121			4.58	20
1,2,3-Trimethylbenzene	5.00	4.45	4.63	89.0	92.6	77.0-120			3.96	20
1,3,5-Trimethylbenzene	5.00	4.43	4.60	88.6	92.0	76.0-122			3.77	20
Vinyl chloride	5.00	5.22	5.81	104	116	67.0-131			10.7	20
Xylenes, Total	15.0	12.6	13.5	84.0	90.0	79.0-123			6.90	20
(S) Toluene-d8				*101*	*100*	*80.0-120*				
(S) 4-Bromofluorobenzene				*94.8*	*97.0*	*77.0-126*				
(S) 1,2-Dichloroethane-d4				*100*	*99.5*	*70.0-130*				

Method Blank (MB)

(MB) R4037088-3 02/21/24 22:25

Analyte	MB Result	MB Qualifier	MB MDL	MB RDL
	ug/l		ug/l	ug/l
1,4-Dioxane	U		0.597	3.00
(S) Toluene-d8	*93.1*			*77.0-127*

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R4037088-1 02/21/24 21:20 • (LCSD) R4037088-2 02/21/24 21:42

Analyte	Spike Amount	LCS Result	LCSD Result	LCS Rec.	LCSD Rec.	Rec. Limits	LCS Qualifier	LCSD Qualifier	RPD	RPD Limits
	ug/l	ug/l	ug/l	%	%	%			%	%
1,4-Dioxane	50.0	45.5	43.9	91.0	87.8	55.0-138			3.58	24
(S) Toluene-d8				*96.1*	*95.9*	*77.0-127*				

L1706680-06 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1706680-06 02/22/24 04:24 • (MS) R4037088-4 02/22/24 09:28 • (MSD) R4037088-5 02/22/24 09:50

Analyte	Spike Amount	Original Result	MS Result	MSD Result	MS Rec.	MSD Rec.	Dilution	Rec. Limits	MS Qualifier	MSD Qualifier	RPD	RPD Limits
	ug/l	ug/l	ug/l	ug/l	%	%		%			%	%
1,4-Dioxane	50.0	U	37.5	48.3	75.0	96.6	1	13.0-160			25.2	31
(S) Toluene-d8					*96.3*	*96.1*		*77.0-127*				

Method Blank (MB)

(MB) R4038326-3 02/25/24 13:01

Analyte	MB Result	MB Qualifier	MB MDL	MB RDL
	ug/l		ug/l	ug/l
1,4-Dioxane	U		0.597	3.00
(S) Toluene-d8	*94.3*			*77.0-127*

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R4038326-1 02/25/24 11:56 • (LCSD) R4038326-2 02/25/24 12:18

Analyte	Spike Amount	LCS Result	LCSD Result	LCS Rec.	LCSD Rec.	Rec. Limits	LCS Qualifier	LCSD Qualifier	RPD	RPD Limits
	ug/l	ug/l	ug/l	%	%	%			%	%
1,4-Dioxane	50.0	45.3	44.0	90.6	88.0	55.0-138			2.91	24
(S) Toluene-d8				*94.3*	*94.6*	*77.0-127*				

L1706682-01 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1706682-01 02/25/24 15:16 • (MS) R4038326-4 02/26/24 00:34 • (MSD) R4038326-5 02/26/24 00:56

Analyte	Spike Amount	Original Result	MS Result	MSD Result	MS Rec.	MSD Rec.	Dilution	Rec. Limits	MS Qualifier	MSD Qualifier	RPD	RPD Limits
	ug/l	ug/l	ug/l	ug/l	%	%		%			%	%
1,4-Dioxane	50.0	U	39.2	43.1	78.4	86.2	1	13.0-160			9.48	31
(S) Toluene-d8					*94.5*	*94.6*		*77.0-127*				

INTERNAL STANDARD SUMMARY

Instrument: VOCMS23 • File ID: 0220_28

02/20/24 21:20

Sample ID	File ID	8260-FLUOROBENZENE	8260-CHLOROBENZENE-D5	8260-1,4-DICHLOROBENZENE-D4
		Response	Response	Response
Standard	0220_28	451930	211496	200886
Upper Limit		903860	422992	401772
Lower Limit		225965	105748	100443
LCS R4036492-1 WG2230378 1x	0220_28LCSA	451930	211496	200886
LCSD R4036492-2 WG2230378 1x	0220_29A	461579	214257	202421
BLANK R4036492-4 WG2230378 1x	0220_32	454490	208042	189610
L1706682-01 WG2230378 1x	0220_50	411383	184436	179916
L1706682-02 WG2230378 1x	0220_51	394696	175771	167965
L1706682-03 WG2230378 1x	0220_52	392083	176422	165080
MS R4036492-6 WG2230378 1x	0220_58	457483	214925	204309
MSD R4036492-7 WG2230378 1x	0220_59	466129	220202	214877

Instrument: VOCMS26 • File ID: 0221_34

02/21/24 19:33

Sample ID	File ID	8260-FLUOROBENZENE	8260-CHLOROBENZENE-D5	8260-1,4-DICHLOROBENZENE-D4
		Response	Response	Response
Standard	0221_34	219583	106973	90711
Upper Limit		439166	213946	181422
Lower Limit		109792	53487	45356
LCS R4036917-1 WG2230965 1x	0221_34LCSA	219583	106973	90711
LCSD R4036917-2 WG2230965 1x	0221_35A	214492	104605	92123
BLANK R4036917-4 WG2230965 1x	0221_39	227009	106124	89131
L1706682-04 WG2230965 1x	0221_61	232976	108419	94516

INTERNAL STANDARD SUMMARY

Instrument: VOCMS27 • File ID: 0221_29

02/21/24 20:58

Sample ID	File ID	8260-FLUOROBENZENE Response
Standard	0221_29	300793
Upper Limit		601586
Lower Limit		150397
LCS R4037088-1 WG2231221 1x	0221_30	304185
LCSD R4037088-2 WG2231221 1x	0221_31	309267
BLANK R4037088-3 WG2231221 1x	0221_33	292020
L1706682-02 WG2231221 1x	0221_45	314572
L1706682-03 WG2231221 1x	0221_46	313268
L1706682-04 WG2231221 1x	0221_47	318978
MS R4037088-4 WG2231221 1x	0221_56	312611
MSD R4037088-5 WG2231221 1x	0221_57	301721

Instrument: VOCMS27 • File ID: 0225_03

02/25/24 11:35

Sample ID	File ID	8260-FLUOROBENZENE Response
Standard	0225_03	311101
Upper Limit		622202
Lower Limit		155551
LCS R4038326-1 WG2231965 1x	0225_04A	301989
LCSD R4038326-2 WG2231965 1x	0225_05A	298077
BLANK R4038326-3 WG2231965 1x	0225_07A	288668
L1706682-01 WG2231965 1x	0225_10	296690
MS R4038326-4 WG2231965 1x	0225_30	321017
MSD R4038326-5 WG2231965 1x	0225_31	302706

GLOSSARY OF TERMS

Guide to Reading and Understanding Your Laboratory Report

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

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

Abbreviations and Definitions

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

Qualifier	Description
B1	Target analyte detected in method blank at or above the method reporting limit.
E4	Concentration estimated. Analyte was detected below laboratory minimum reporting level (MRL) but above MDL.
L1	The associated blank spike recovery was above laboratory acceptance limits.
M1	Matrix spike recovery was high, the method control sample recovery was acceptable.
M2	Matrix spike recovery was low, the method control sample recovery was acceptable.
M3	The spike recovery value is unusable since the analyte concentration in the sample is disproportionate to the spike level. The associated blank spike recovery was acceptable.
R5	MS/MSD RPD exceeded the laboratory acceptance limit. Recovery met acceptance criteria.
R7	LFB/LFBD RPD exceeded the laboratory acceptance limit. Recovery met acceptance criteria.
R8	Sample RPD exceeded the method acceptance limit.

ACCREDITATIONS & LOCATIONS

Pace Analytical National 12065 Lebanon Rd Mount Juliet, TN 37122

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

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

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

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

Company Name/Address:
Pinyon Environmental
3222 S. Vance Street Suite 200
Lakewood, CO 80227

Billing Information:
Accounts Payable
3222 S Vance Street
Suite 200
Lakewood, CO 80227

Pres Chk

Analysis / Container / Preservative

Chain of Custody Page 1 of 1

Pace PEOPLE ADVANCING SCIENCE

MT JULIET, TN

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

Report to:
Isabella Foster and Andrew Parker

Email To: foster@pinyon-env.com;parker@pinyon-

Project Description: Nammo WBO Groundwater Monitoring

City/State Collected: Mesa, AZ

Please Circle: PT (MT) CT ET

Phone: 303-785-7697

Client Project # 722152201

Lab Project # PINYONMAZ-722152201

SDG # 1706682

B130

Collected by (print): Isabella Foster

Site/Facility ID #

P.O. #

Acctnum: PINYONMAZ
Template: T205686
Prelogin: P1054555
PM: 288 - Daphne Richards
PB:

Collected by (signature): [signature]

Immediately Packed on Ice N____ Y X

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

Quote #

Date Results Needed: Standard TAT

Shipped Via:

Sample ID	Comp/Grab	Matrix *	Depth	Date	Time	No. of Cntrs	PERCHLORATE 125mlHDPE-NoPres	V8260AZ 40mlAmb-HCl	V8260LL14D 40mlAmb-HCl	Remarks	Sample # (lab only)
PF-2-400-20240214	C	GW	400	2-14-24	1500	12 ~~7~~	X	X	X	* MS/MSD	-01
TU-8-164-20240214	G	GW	164	2-14-24	1520	7	X	X	X		-02
TU-10-172-20240214	G	GW	172	2-14-24	1349	7	X	X	X		-03
DUP-03	–	GW	–	–	–	7	X	X	X		-04
Trip Blank	–	–	–	–	–	1					-05

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

Remarks:

pH _______ Temp _______
Flow _______ Other _______

Samples returned via:
__ UPS __ FedEx __ Courier ________

Tracking #

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

Relinquished by : (Signature)	Date:	Time:	Received by: (Signature)	
[signature]	2-15-24	0943	[signature]	Trip Blank Received: (Yes) / No HCL / MeoH TBR
[signature]	2-15-24	1800	SWA	Temp: DPA8 4.4+0=4.4 °C Bottles Received:
			Received for lab by: (Signature) [signature]	Date: 2-16-24 Time: 8:00

If preservation required by Login: Date/Time

Hold:

Condition: NCF / (OK)

PNPAZ

ANALYTICAL REPORT

March 20, 2024

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

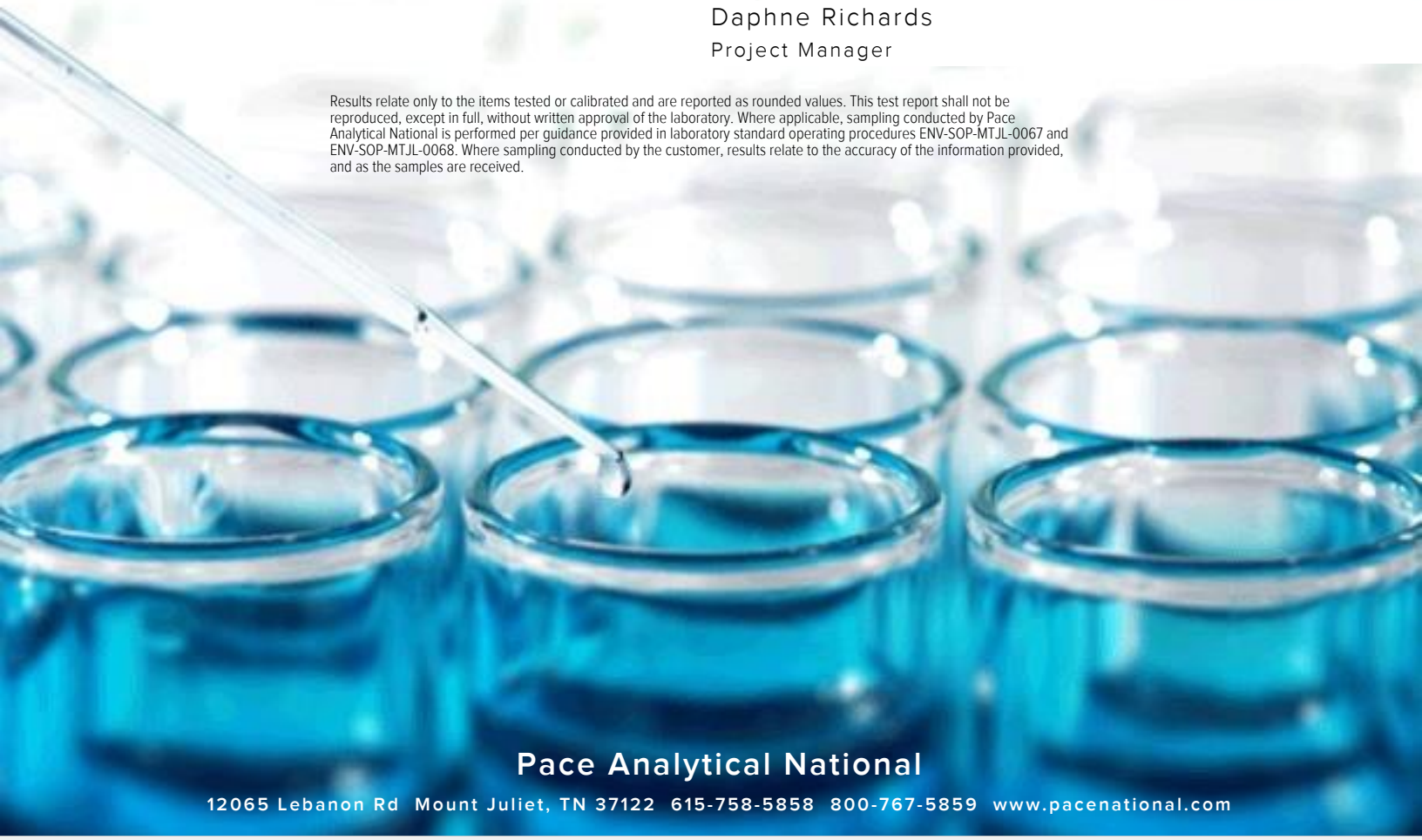
Pinyon Environmental

Sample Delivery Group:	L1716428
Samples Received:	03/19/2024
Project Number:	722152201.002
Description:	Nammo TTU Groundwater Monitoring

Report To: Isbella Foster and Andrew Parker
3222 S. Vance Street Suite 200
Lakewood, CO 80227

Entire Report Reviewed By: Daphne R Richards

Daphne Richards
Project Manager



Pace Analytical National
12065 Lebanon Rd Mount Juliet, TN 37122 615-758-5858 800-767-5859 www.pacenational.com

TABLE OF CONTENTS

SAMPLE SUMMARY

TTU-3-108-20240318 L1716428-01 GW

			Collected by	Collected date/time	Received date/time
				03/18/24 14:22	03/19/24 09:45

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Volatile Organic Compounds (GC/MS) by Method 8260B-SIM	WG2249950	1	03/20/24 01:05	03/20/24 01:05	JAH	Mt. Juliet, TN

CASE NARRATIVE

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

Daphne Richards
Project Manager

Sample Delivery Group (SDG) Narrative

Insufficient sample volume to perform MS/MSD analyses per method QC requirements.

Lab Sample ID	Project Sample ID	Method
L1716428-01	TTU-3-108-20240318	8260B-SIM

Volatile Organic Compounds (GC/MS) by Method 8260B-SIM

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
1,4-Dioxane	U		0.597	3.00	1	03/20/2024 01:05	WG2249950
(S) Toluene-d8	*95.9*			*77.0-127*		*03/20/2024 01:05*	WG2249950

Method Blank (MB)

(MB) R4047904-3 03/20/24 00:30

Analyte	MB Result ug/l	MB Qualifier	MB MDL ug/l	MB RDL ug/l
1,4-Dioxane	U		0.597	3.00
(S) Toluene-d8	*96.5*			*77.0-127*

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R4047904-1 03/19/24 23:25 • (LCSD) R4047904-2 03/19/24 23:46

Analyte	Spike Amount ug/l	LCS Result ug/l	LCSD Result ug/l	LCS Rec. %	LCSD Rec. %	Rec. Limits %	LCS Qualifier	LCSD Qualifier	RPD %	RPD Limits %
1,4-Dioxane	50.0	43.3	42.3	86.6	84.6	55.0-138			2.34	24
(S) Toluene-d8				*96.4*	*96.4*	*77.0-127*				

INTERNAL STANDARD SUMMARY

Instrument: VOCMS27 • File ID: 0319_04

03/19/24 23:03

Sample ID	File ID	8260-FLUOROBENZENE Response
Standard	0319_04	347261
Upper Limit		694522
Lower Limit		173631
LCS R4047904-1 WG2249950 1x	0319_05	344162
LCSD R4047904-2 WG2249950 1x	0319_06	361034
BLANK R4047904-3 WG2249950 1x	0319_08	366776
L1716428-01 WG2249950 1x	0319_09	310130

GLOSSARY OF TERMS

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Abbreviations and Definitions

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

Qualifier	Description

The remainder of this page intentionally left blank, there are no qualifiers applied to this SDG.

ACCREDITATIONS & LOCATIONS

Pace Analytical National 12065 Lebanon Rd Mount Juliet, TN 37122

Alabama	40660	Nebraska	NE-OS-15-05
Alaska	17-026	Nevada	TN000032021-1
Arizona	AZ0612	New Hampshire	2975
Arkansas	88-0469	New Jersey–NELAP	TN002
California	2932	New Mexico [1]	TN00003
Colorado	TN00003	New York	11742
Connecticut	PH-0197	North Carolina	Env375
Florida	E87487	North Carolina [1]	DW21704
Georgia	NELAP	North Carolina [3]	41
Georgia [1]	923	North Dakota	R-140
Idaho	TN00003	Ohio–VAP	CL0069
Illinois	200008	Oklahoma	9915
Indiana	C-TN-01	Oregon	TN200002
Iowa	364	Pennsylvania	68-02979
Kansas	E-10277	Rhode Island	LAO000356
Kentucky [1 6]	KY90010	South Carolina	84004002
Kentucky [2]	16	South Dakota	n/a
Louisiana	AI30792	Tennessee [1 4]	2006
Louisiana	LA018	Texas	T104704245-20-18
Maine	TN00003	Texas [5]	LAB0152
Maryland	324	Utah	TN000032021-11
Massachusetts	M-TN003	Vermont	VT2006
Michigan	9958	Virginia	110033
Minnesota	047-999-395	Washington	C847
Mississippi	TN00003	West Virginia	233
Missouri	340	Wisconsin	998093910
Montana	CERT0086	Wyoming	A2LA
A2LA – ISO 17025	1461.01	AIHA-LAP,LLC EMLAP	100789
A2LA – ISO 17025 [5]	1461.02	DOD	1461.01
Canada	1461.01	USDA	P330-15-00234
EPA–Crypto	TN00003		

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

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

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

Company Name/Address:
Pinyon Environmental

3222 S. Vance Street Suite 200
Lakewood, CO 80227

Billing Information:
Accounts Payable
3222 S Vance Street
Suite 200
Lakewood, CO 80227

Pres Chk

Report to:
Isbella Foster and Andrew Parker

Email To: **foster@pinyon-env.com;parker@pinyon-**

Project Description:
Nammo TTU Groundwater Monitoring

City/State Collected: Mesa, AZ

Please Circle: (PT) MT CT ET

Phone: **303-785-7697**

Client Project #
722152201.002

Lab Project #
PINYONMAZ-722152201

Collected by (print):
Isabella Foster

Site/Facility ID #

P.O. #

Collected by (signature):

Immediately Packed on Ice N ___ Y X

Rush? (Lab MUST Be Notified)
___ Same Day ___ Five Day
___ Next Day ___ 5 Day (Rad Only)
X Two Day ___ 10 Day (Rad Only)
___ Three Day

Quote #

Date Results Needed: 48 hours

Sample ID	Comp/Grab	Matrix *	Depth	Date	Time	No. of Cntrs	PERCHLORATE 125mlHDPE-NoPres	V8260AZ 40mlAmb-HCl	V8260LL14D 40mlAmb-HCl								
TTU-3-108-20240318	G	GW	108	3/18/24	1422	3			X								-01
		GW															
		GW															
		GW															
		GW															
		GW															
		GW															
		GW															
		GW															
		GW															

Analysis / Container / Preservative

Chain of Custody Page 1 of 1

Pace
PEOPLE ADVANCING SCIENCE

MT JULIET, TN

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

SDG # L1716426

F052

Acctnum: **PINYONMAZ**
Template: **T205653**
Prelogin: **P1054553**
PM: **288 - Daphne Richards**
PB:

Shipped Via:

Remarks | Sample # (lab only)

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

Remarks:

pH _____ Temp _____

Flow _____ Other _____

Samples returned via:
__ UPS __ FedEx __ Courier _____

Tracking #

Sample Receipt Checklist

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

	Date:	Time:		
Relinquished by : (Signature)	3-18-24	1523	Received by: (Signature)	Trip Blank Received: Yes / No HCL / MeoH TBR
Relinquished by : (Signature)	3-18-24	1800	Received by: (Signature)	Temp: 0.2 °C 1.2 Bottles Received: 3
Relinquished by : (Signature)	Date:	Time:	Received for lab by: (Signature) Christopher L. Gallin	Date: 3/19/24 Time: 0945

If preservation required by Login: Date/Time

Hold:

Condition: NCF / OK

ANALYTICAL REPORT

April 15, 2024

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

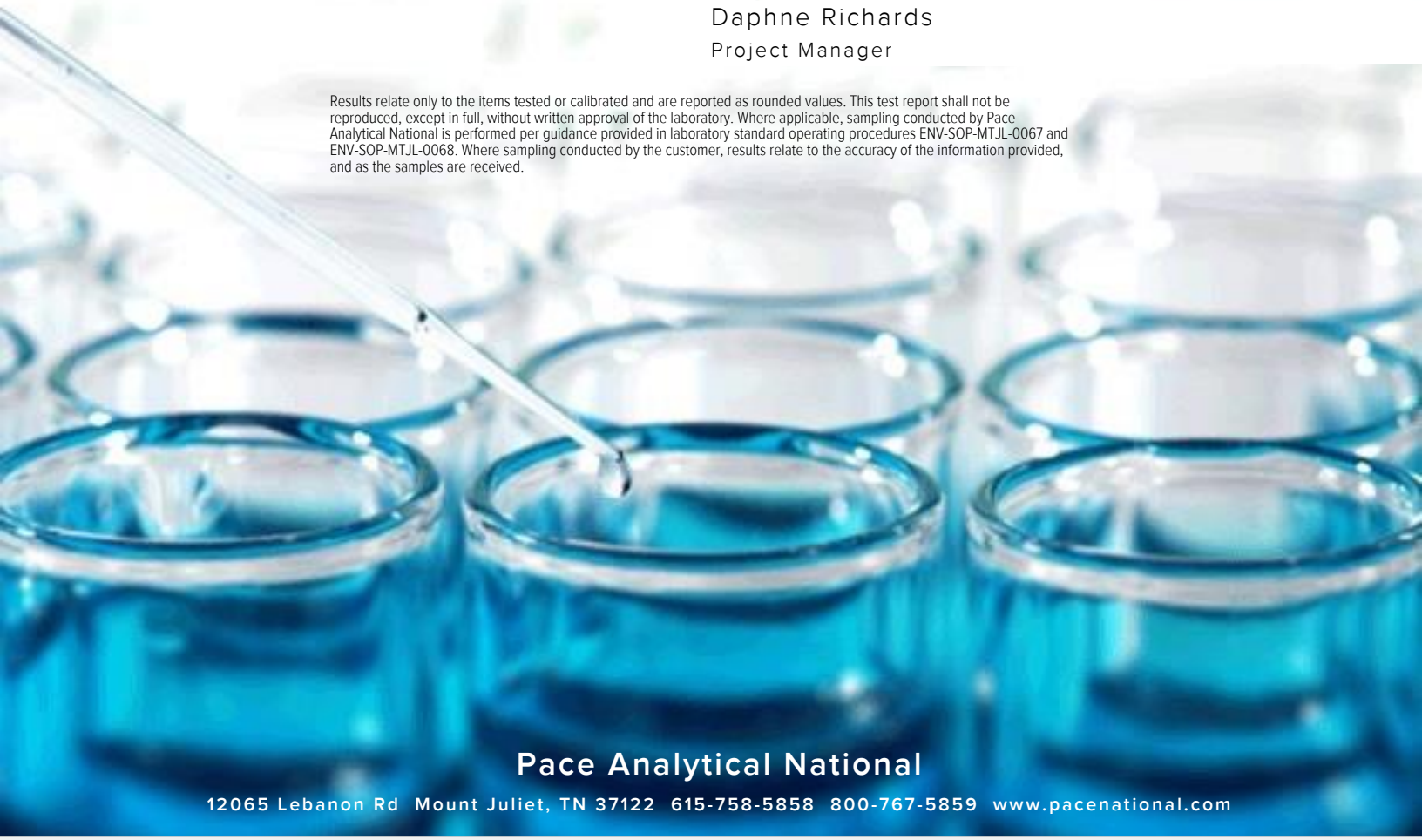
Pinyon Environmental

Sample Delivery Group:	L1724110
Samples Received:	04/10/2024
Project Number:	722152201.002
Description:	Nammo TTU Groundwater

Report To: Andrew Parker
4815 E. Carefree Highway
#108-274
Cave Creek, AZ 85331

Entire Report Reviewed By: Daphne R Richards

Daphne Richards
Project Manager



Pace Analytical National
12065 Lebanon Rd Mount Juliet, TN 37122 615-758-5858 800-767-5859 www.pacenational.com

TABLE OF CONTENTS

SAMPLE SUMMARY

TTU-15-75-20240409 L1724110-01 GW

Collected by: Crissa Wolf
Collected date/time: 04/09/24 10:46
Received date/time: 04/10/24 09:15

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2264996	500	04/13/24 22:25	04/13/24 22:25	ASM	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2264870	20	04/11/24 19:17	04/11/24 19:17	DYW	Mt. Juliet, TN

TTU-16-80-20240409 L1724110-02 GW

Collected by: Crissa Wolf
Collected date/time: 04/09/24 10:26
Received date/time: 04/10/24 09:15

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Wet Chemistry by Method 314.0 Mod	WG2264996	50000	04/13/24 23:24	04/13/24 23:24	ASM	Mt. Juliet, TN
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2264870	2500	04/11/24 19:37	04/11/24 19:37	DYW	Mt. Juliet, TN

TRIP BLANK L1724110-03 GW

Collected by: Crissa Wolf
Collected date/time: 04/09/24 00:00
Received date/time: 04/10/24 09:15

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst	Location
Volatile Organic Compounds (GC/MS) by Method 8260B	WG2264870	1	04/11/24 15:51	04/11/24 15:51	DYW	Mt. Juliet, TN

CASE NARRATIVE

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

Daphne Richards
Project Manager

Sample Delivery Group (SDG) Narrative

Insufficient sample volume to perform MS/MSD analyses per method QC requirements.

Lab Sample ID	Project Sample ID	Method
L1724110-01	TTU-15-75-20240409	8260B
L1724110-02	TTU-16-80-20240409	8260B
L1724110-03	TRIP BLANK	8260B

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	11100		150	2000	500	04/13/2024 22:25	WG2264996

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		226	1000	20	04/11/2024 19:17	WG2264870
Acrolein	U		50.8	1000	20	04/11/2024 19:17	WG2264870
Acrylonitrile	U		13.4	200	20	04/11/2024 19:17	WG2264870
Benzene	U		1.88	20.0	20	04/11/2024 19:17	WG2264870
Bromobenzene	U		2.36	20.0	20	04/11/2024 19:17	WG2264870
Bromodichloromethane	U		2.72	20.0	20	04/11/2024 19:17	WG2264870
Bromoform	U		2.58	20.0	20	04/11/2024 19:17	WG2264870
Bromomethane	U		12.1	100	20	04/11/2024 19:17	WG2264870
1,3-Butadiene	U		5.98	40.0	20	04/11/2024 19:17	WG2264870
n-Butylbenzene	U		3.14	20.0	20	04/11/2024 19:17	WG2264870
sec-Butylbenzene	U		2.50	20.0	20	04/11/2024 19:17	WG2264870
tert-Butylbenzene	U		2.54	20.0	20	04/11/2024 19:17	WG2264870
Carbon tetrachloride	U		2.56	20.0	20	04/11/2024 19:17	WG2264870
Carbon disulfide	3.98	B1 E4	1.92	20.0	20	04/11/2024 19:17	WG2264870
Chlorobenzene	U		2.32	20.0	20	04/11/2024 19:17	WG2264870
Chlorodibromomethane	U		2.80	20.0	20	04/11/2024 19:17	WG2264870
Chloroethane	U		3.84	100	20	04/11/2024 19:17	WG2264870
Chloroform	U		2.22	100	20	04/11/2024 19:17	WG2264870
Chloromethane	U		19.2	50.0	20	04/11/2024 19:17	WG2264870
Cyclohexane	U		3.76	20.0	20	04/11/2024 19:17	WG2264870
2-Chlorotoluene	U		2.12	20.0	20	04/11/2024 19:17	WG2264870
4-Chlorotoluene	U		2.28	20.0	20	04/11/2024 19:17	WG2264870
1,2-Dibromo-3-Chloropropane	U		5.52	100	20	04/11/2024 19:17	WG2264870
1,2-Dibromoethane	U		2.52	20.0	20	04/11/2024 19:17	WG2264870
Dibromomethane	U		2.44	20.0	20	04/11/2024 19:17	WG2264870
1,2-Dichlorobenzene	U		2.14	20.0	20	04/11/2024 19:17	WG2264870
1,3-Dichlorobenzene	U		2.20	20.0	20	04/11/2024 19:17	WG2264870
1,4-Dichlorobenzene	U		2.40	20.0	20	04/11/2024 19:17	WG2264870
Dichlorodifluoromethane	U		7.48	100	20	04/11/2024 19:17	WG2264870
1,1-Dichloroethane	U		2.00	20.0	20	04/11/2024 19:17	WG2264870
1,2-Dichloroethane	U		1.64	20.0	20	04/11/2024 19:17	WG2264870
1,1-Dichloroethene	U		3.76	20.0	20	04/11/2024 19:17	WG2264870
cis-1,2-Dichloroethene	U		2.52	20.0	20	04/11/2024 19:17	WG2264870
trans-1,2-Dichloroethene	U		2.98	20.0	20	04/11/2024 19:17	WG2264870
1,2-Dichloropropane	U		2.98	20.0	20	04/11/2024 19:17	WG2264870
1,1-Dichloropropene	U		2.84	20.0	20	04/11/2024 19:17	WG2264870
1,3-Dichloropropane	U		2.20	20.0	20	04/11/2024 19:17	WG2264870
cis-1,3-Dichloropropene	U		2.22	20.0	20	04/11/2024 19:17	WG2264870
trans-1,3-Dichloropropene	U		2.36	20.0	20	04/11/2024 19:17	WG2264870
2,2-Dichloropropane	U		3.22	20.0	20	04/11/2024 19:17	WG2264870
Dicyclopentadiene	U		5.06	20.0	20	04/11/2024 19:17	WG2264870
Di-isopropyl ether	U		2.10	20.0	20	04/11/2024 19:17	WG2264870
Ethylbenzene	U		2.74	20.0	20	04/11/2024 19:17	WG2264870
4-Ethyltoluene	U		4.16	20.0	20	04/11/2024 19:17	WG2264870
Hexachloro-1,3-butadiene	U		6.74	20.0	20	04/11/2024 19:17	WG2264870
n-Hexane	U		15.0	200	20	04/11/2024 19:17	WG2264870
Isopropylbenzene	U		2.10	20.0	20	04/11/2024 19:17	WG2264870
p-Isopropyltoluene	U		2.40	20.0	20	04/11/2024 19:17	WG2264870
2-Butanone (MEK)	U		23.8	200	20	04/11/2024 19:17	WG2264870
Methyl Cyclohexane	U		13.2	20.0	20	04/11/2024 19:17	WG2264870

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Methylene Chloride	U		8.60	100	20	04/11/2024 19:17	WG2264870
4-Methyl-2-pentanone (MIBK)	U		9.56	200	20	04/11/2024 19:17	WG2264870
Methyl tert-butyl ether	U		2.02	20.0	20	04/11/2024 19:17	WG2264870
Naphthalene	U		20.0	100	20	04/11/2024 19:17	WG2264870
Propene	U		18.7	50.0	20	04/11/2024 19:17	WG2264870
n-Propylbenzene	U		1.99	20.0	20	04/11/2024 19:17	WG2264870
Styrene	U		2.36	20.0	20	04/11/2024 19:17	WG2264870
1,1,1,2-Tetrachloroethane	U		2.94	20.0	20	04/11/2024 19:17	WG2264870
1,1,2,2-Tetrachloroethane	U		2.66	20.0	20	04/11/2024 19:17	WG2264870
1,1,2-Trichlorotrifluoroethane	U		3.60	20.0	20	04/11/2024 19:17	WG2264870
Tetrachloroethene	U		6.00	20.0	20	04/11/2024 19:17	WG2264870
Toluene	U		5.56	20.0	20	04/11/2024 19:17	WG2264870
1,2,3-Trichlorobenzene	U		4.60	20.0	20	04/11/2024 19:17	WG2264870
1,2,4-Trichlorobenzene	U		9.62	20.0	20	04/11/2024 19:17	WG2264870
1,1,1-Trichloroethane	U		2.98	20.0	20	04/11/2024 19:17	WG2264870
1,1,2-Trichloroethane	U		3.16	20.0	20	04/11/2024 19:17	WG2264870
Trichloroethene	U		3.80	20.0	20	04/11/2024 19:17	WG2264870
Trichlorofluoromethane	U		3.20	100	20	04/11/2024 19:17	WG2264870
1,2,3-Trichloropropane	U		4.74	50.0	20	04/11/2024 19:17	WG2264870
1,2,4-Trimethylbenzene	U		6.44	20.0	20	04/11/2024 19:17	WG2264870
1,2,3-Trimethylbenzene	U		2.08	20.0	20	04/11/2024 19:17	WG2264870
1,3,5-Trimethylbenzene	U		2.08	20.0	20	04/11/2024 19:17	WG2264870
Vinyl chloride	U		4.68	20.0	20	04/11/2024 19:17	WG2264870
Xylenes, Total	U		3.48	60.0	20	04/11/2024 19:17	WG2264870
(S) Toluene-d8	*108*			*80.0-120*		*04/11/2024 19:17*	WG2264870
(S) 4-Bromofluorobenzene	*94.1*			*77.0-126*		*04/11/2024 19:17*	WG2264870
(S) 1,2-Dichloroethane-d4	*86.3*			*70.0-130*		*04/11/2024 19:17*	WG2264870

Wet Chemistry by Method 314.0 Mod

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Perchlorate	817000		15000	200000	50000	04/13/2024 23:24	WG2264996

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		28200	125000	2500	04/11/2024 19:37	WG2264870
Acrolein	U		6350	125000	2500	04/11/2024 19:37	WG2264870
Acrylonitrile	U		1680	25000	2500	04/11/2024 19:37	WG2264870
Benzene	260	E4	235	2500	2500	04/11/2024 19:37	WG2264870
Bromobenzene	U		295	2500	2500	04/11/2024 19:37	WG2264870
Bromodichloromethane	U		340	2500	2500	04/11/2024 19:37	WG2264870
Bromoform	U		322	2500	2500	04/11/2024 19:37	WG2264870
Bromomethane	U		1510	12500	2500	04/11/2024 19:37	WG2264870
1,3-Butadiene	U		748	5000	2500	04/11/2024 19:37	WG2264870
n-Butylbenzene	U		393	2500	2500	04/11/2024 19:37	WG2264870
sec-Butylbenzene	U		313	2500	2500	04/11/2024 19:37	WG2264870
tert-Butylbenzene	U		318	2500	2500	04/11/2024 19:37	WG2264870
Carbon tetrachloride	U		320	2500	2500	04/11/2024 19:37	WG2264870
Carbon disulfide	494	B1 E4	241	2500	2500	04/11/2024 19:37	WG2264870
Chlorobenzene	U		290	2500	2500	04/11/2024 19:37	WG2264870
Chlorodibromomethane	U		350	2500	2500	04/11/2024 19:37	WG2264870
Chloroethane	U		480	12500	2500	04/11/2024 19:37	WG2264870
Chloroform	U		278	12500	2500	04/11/2024 19:37	WG2264870
Chloromethane	U		2400	6250	2500	04/11/2024 19:37	WG2264870
Cyclohexane	U		470	2500	2500	04/11/2024 19:37	WG2264870
2-Chlorotoluene	U		265	2500	2500	04/11/2024 19:37	WG2264870
4-Chlorotoluene	U		285	2500	2500	04/11/2024 19:37	WG2264870
1,2-Dibromo-3-Chloropropane	U		690	12500	2500	04/11/2024 19:37	WG2264870
1,2-Dibromoethane	U		315	2500	2500	04/11/2024 19:37	WG2264870
Dibromomethane	U		305	2500	2500	04/11/2024 19:37	WG2264870
1,2-Dichlorobenzene	U		268	2500	2500	04/11/2024 19:37	WG2264870
1,3-Dichlorobenzene	U		275	2500	2500	04/11/2024 19:37	WG2264870
1,4-Dichlorobenzene	U		300	2500	2500	04/11/2024 19:37	WG2264870
Dichlorodifluoromethane	U		935	12500	2500	04/11/2024 19:37	WG2264870
1,1-Dichloroethane	U		250	2500	2500	04/11/2024 19:37	WG2264870
1,2-Dichloroethane	U		205	2500	2500	04/11/2024 19:37	WG2264870
1,1-Dichloroethene	2440	E4	470	2500	2500	04/11/2024 19:37	WG2264870
cis-1,2-Dichloroethene	U		315	2500	2500	04/11/2024 19:37	WG2264870
trans-1,2-Dichloroethene	U		373	2500	2500	04/11/2024 19:37	WG2264870
1,2-Dichloropropane	U		373	2500	2500	04/11/2024 19:37	WG2264870
1,1-Dichloropropene	U		355	2500	2500	04/11/2024 19:37	WG2264870
1,3-Dichloropropane	U		275	2500	2500	04/11/2024 19:37	WG2264870
cis-1,3-Dichloropropene	U		278	2500	2500	04/11/2024 19:37	WG2264870
trans-1,3-Dichloropropene	U		295	2500	2500	04/11/2024 19:37	WG2264870
2,2-Dichloropropane	U		403	2500	2500	04/11/2024 19:37	WG2264870
Dicyclopentadiene	U		633	2500	2500	04/11/2024 19:37	WG2264870
Di-isopropyl ether	U		263	2500	2500	04/11/2024 19:37	WG2264870
Ethylbenzene	U		343	2500	2500	04/11/2024 19:37	WG2264870
4-Ethyltoluene	U		520	2500	2500	04/11/2024 19:37	WG2264870
Hexachloro-1,3-butadiene	U		843	2500	2500	04/11/2024 19:37	WG2264870
n-Hexane	U		1870	25000	2500	04/11/2024 19:37	WG2264870
Isopropylbenzene	U		263	2500	2500	04/11/2024 19:37	WG2264870
p-Isopropyltoluene	U		300	2500	2500	04/11/2024 19:37	WG2264870
2-Butanone (MEK)	U		2980	25000	2500	04/11/2024 19:37	WG2264870
Methyl Cyclohexane	U		1650	2500	2500	04/11/2024 19:37	WG2264870

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Methylene Chloride	52100		1080	12500	2500	04/11/2024 19:37	WG2264870
4-Methyl-2-pentanone (MIBK)	U		1200	25000	2500	04/11/2024 19:37	WG2264870
Methyl tert-butyl ether	U		253	2500	2500	04/11/2024 19:37	WG2264870
Naphthalene	U		2500	12500	2500	04/11/2024 19:37	WG2264870
Propene	U		2340	6250	2500	04/11/2024 19:37	WG2264870
n-Propylbenzene	U		248	2500	2500	04/11/2024 19:37	WG2264870
Styrene	U		295	2500	2500	04/11/2024 19:37	WG2264870
1,1,1,2-Tetrachloroethane	U		368	2500	2500	04/11/2024 19:37	WG2264870
1,1,2,2-Tetrachloroethane	U		333	2500	2500	04/11/2024 19:37	WG2264870
1,1,2-Trichlorotrifluoroethane	U		450	2500	2500	04/11/2024 19:37	WG2264870
Tetrachloroethene	U		750	2500	2500	04/11/2024 19:37	WG2264870
Toluene	U		695	2500	2500	04/11/2024 19:37	WG2264870
1,2,3-Trichlorobenzene	U		575	2500	2500	04/11/2024 19:37	WG2264870
1,2,4-Trichlorobenzene	U		1200	2500	2500	04/11/2024 19:37	WG2264870
1,1,1-Trichloroethane	U		373	2500	2500	04/11/2024 19:37	WG2264870
1,1,2-Trichloroethane	U		395	2500	2500	04/11/2024 19:37	WG2264870
Trichloroethene	68500		475	2500	2500	04/11/2024 19:37	WG2264870
Trichlorofluoromethane	U		400	12500	2500	04/11/2024 19:37	WG2264870
1,2,3-Trichloropropane	U		592	6250	2500	04/11/2024 19:37	WG2264870
1,2,4-Trimethylbenzene	U		805	2500	2500	04/11/2024 19:37	WG2264870
1,2,3-Trimethylbenzene	U		260	2500	2500	04/11/2024 19:37	WG2264870
1,3,5-Trimethylbenzene	U		260	2500	2500	04/11/2024 19:37	WG2264870
Vinyl chloride	U		585	2500	2500	04/11/2024 19:37	WG2264870
Xylenes, Total	U		435	7500	2500	04/11/2024 19:37	WG2264870
(S) Toluene-d8	*110*			*80.0-120*		*04/11/2024 19:37*	WG2264870
(S) 4-Bromofluorobenzene	*96.3*			*77.0-126*		*04/11/2024 19:37*	WG2264870
(S) 1,2-Dichloroethane-d4	*85.4*			*70.0-130*		*04/11/2024 19:37*	WG2264870

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		11.3	50.0	1	04/11/2024 15:51	WG2264870
Acrolein	U		2.54	50.0	1	04/11/2024 15:51	WG2264870
Acrylonitrile	U		0.671	10.0	1	04/11/2024 15:51	WG2264870
Benzene	U		0.0941	1.00	1	04/11/2024 15:51	WG2264870
Bromobenzene	U		0.118	1.00	1	04/11/2024 15:51	WG2264870
Bromodichloromethane	U		0.136	1.00	1	04/11/2024 15:51	WG2264870
Bromoform	U		0.129	1.00	1	04/11/2024 15:51	WG2264870
Bromomethane	U		0.605	5.00	1	04/11/2024 15:51	WG2264870
1,3-Butadiene	U		0.299	2.00	1	04/11/2024 15:51	WG2264870
n-Butylbenzene	U		0.157	1.00	1	04/11/2024 15:51	WG2264870
sec-Butylbenzene	U		0.125	1.00	1	04/11/2024 15:51	WG2264870
tert-Butylbenzene	U		0.127	1.00	1	04/11/2024 15:51	WG2264870
Carbon tetrachloride	U		0.128	1.00	1	04/11/2024 15:51	WG2264870
Carbon disulfide	0.162	B1 E4	0.0962	1.00	1	04/11/2024 15:51	WG2264870
Chlorobenzene	U		0.116	1.00	1	04/11/2024 15:51	WG2264870
Chlorodibromomethane	U		0.140	1.00	1	04/11/2024 15:51	WG2264870
Chloroethane	U		0.192	5.00	1	04/11/2024 15:51	WG2264870
Chloroform	U		0.111	5.00	1	04/11/2024 15:51	WG2264870
Chloromethane	U		0.960	2.50	1	04/11/2024 15:51	WG2264870
Cyclohexane	U		0.188	1.00	1	04/11/2024 15:51	WG2264870
2-Chlorotoluene	U		0.106	1.00	1	04/11/2024 15:51	WG2264870
4-Chlorotoluene	U		0.114	1.00	1	04/11/2024 15:51	WG2264870
1,2-Dibromo-3-Chloropropane	U		0.276	5.00	1	04/11/2024 15:51	WG2264870
1,2-Dibromoethane	U		0.126	1.00	1	04/11/2024 15:51	WG2264870
Dibromomethane	U		0.122	1.00	1	04/11/2024 15:51	WG2264870
1,2-Dichlorobenzene	U		0.107	1.00	1	04/11/2024 15:51	WG2264870
1,3-Dichlorobenzene	U		0.110	1.00	1	04/11/2024 15:51	WG2264870
1,4-Dichlorobenzene	U		0.120	1.00	1	04/11/2024 15:51	WG2264870
Dichlorodifluoromethane	U		0.374	5.00	1	04/11/2024 15:51	WG2264870
1,1-Dichloroethane	U		0.100	1.00	1	04/11/2024 15:51	WG2264870
1,2-Dichloroethane	U		0.0819	1.00	1	04/11/2024 15:51	WG2264870
1,1-Dichloroethene	U		0.188	1.00	1	04/11/2024 15:51	WG2264870
cis-1,2-Dichloroethene	U		0.126	1.00	1	04/11/2024 15:51	WG2264870
trans-1,2-Dichloroethene	U		0.149	1.00	1	04/11/2024 15:51	WG2264870
1,2-Dichloropropane	U		0.149	1.00	1	04/11/2024 15:51	WG2264870
1,1-Dichloropropene	U		0.142	1.00	1	04/11/2024 15:51	WG2264870
1,3-Dichloropropane	U		0.110	1.00	1	04/11/2024 15:51	WG2264870
cis-1,3-Dichloropropene	U		0.111	1.00	1	04/11/2024 15:51	WG2264870
trans-1,3-Dichloropropene	U		0.118	1.00	1	04/11/2024 15:51	WG2264870
2,2-Dichloropropane	U		0.161	1.00	1	04/11/2024 15:51	WG2264870
Dicyclopentadiene	U		0.253	1.00	1	04/11/2024 15:51	WG2264870
Di-isopropyl ether	U		0.105	1.00	1	04/11/2024 15:51	WG2264870
Ethylbenzene	U		0.137	1.00	1	04/11/2024 15:51	WG2264870
4-Ethyltoluene	U		0.208	1.00	1	04/11/2024 15:51	WG2264870
Hexachloro-1,3-butadiene	U		0.337	1.00	1	04/11/2024 15:51	WG2264870
n-Hexane	U		0.749	10.0	1	04/11/2024 15:51	WG2264870
Isopropylbenzene	U		0.105	1.00	1	04/11/2024 15:51	WG2264870
p-Isopropyltoluene	U		0.120	1.00	1	04/11/2024 15:51	WG2264870
2-Butanone (MEK)	U		1.19	10.0	1	04/11/2024 15:51	WG2264870
Methyl Cyclohexane	U		0.660	1.00	1	04/11/2024 15:51	WG2264870
Methylene Chloride	U		0.430	5.00	1	04/11/2024 15:51	WG2264870
4-Methyl-2-pentanone (MIBK)	U		0.478	10.0	1	04/11/2024 15:51	WG2264870
Methyl tert-butyl ether	U		0.101	1.00	1	04/11/2024 15:51	WG2264870
Naphthalene	U		1.00	5.00	1	04/11/2024 15:51	WG2264870
Propene	U		0.936	2.50	1	04/11/2024 15:51	WG2264870
n-Propylbenzene	U		0.0993	1.00	1	04/11/2024 15:51	WG2264870

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

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Styrene	U		0.118	1.00	1	04/11/2024 15:51	WG2264870
1,1,1,2-Tetrachloroethane	U		0.147	1.00	1	04/11/2024 15:51	WG2264870
1,1,2,2-Tetrachloroethane	U		0.133	1.00	1	04/11/2024 15:51	WG2264870
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00	1	04/11/2024 15:51	WG2264870
Tetrachloroethene	U		0.300	1.00	1	04/11/2024 15:51	WG2264870
Toluene	U		0.278	1.00	1	04/11/2024 15:51	WG2264870
1,2,3-Trichlorobenzene	U		0.230	1.00	1	04/11/2024 15:51	WG2264870
1,2,4-Trichlorobenzene	U		0.481	1.00	1	04/11/2024 15:51	WG2264870
1,1,1-Trichloroethane	U		0.149	1.00	1	04/11/2024 15:51	WG2264870
1,1,2-Trichloroethane	U		0.158	1.00	1	04/11/2024 15:51	WG2264870
Trichloroethene	U		0.190	1.00	1	04/11/2024 15:51	WG2264870
Trichlorofluoromethane	U		0.160	5.00	1	04/11/2024 15:51	WG2264870
1,2,3-Trichloropropane	U		0.237	2.50	1	04/11/2024 15:51	WG2264870
1,2,4-Trimethylbenzene	U		0.322	1.00	1	04/11/2024 15:51	WG2264870
1,2,3-Trimethylbenzene	U		0.104	1.00	1	04/11/2024 15:51	WG2264870
1,3,5-Trimethylbenzene	U		0.104	1.00	1	04/11/2024 15:51	WG2264870
Vinyl chloride	U		0.234	1.00	1	04/11/2024 15:51	WG2264870
Xylenes, Total	U		0.174	3.00	1	04/11/2024 15:51	WG2264870
(S) Toluene-d8	*108*			*80.0-120*		*04/11/2024 15:51*	WG2264870
(S) 4-Bromofluorobenzene	*93.6*			*77.0-126*		*04/11/2024 15:51*	WG2264870
(S) 1,2-Dichloroethane-d4	*86.9*			*70.0-130*		*04/11/2024 15:51*	WG2264870

Method Blank (MB)

(MB) R4057423-1 04/13/24 20:11

Analyte	MB Result ug/l	MB Qualifier	MB MDL ug/l	MB RDL ug/l
Perchlorate	U		0.300	4.00

L1724110-01 Original Sample (OS) • Duplicate (DUP)

(OS) L1724110-01 04/13/24 22:25 • (DUP) R4057423-3 04/13/24 22:56

Analyte	Original Result ug/l	DUP Result ug/l	Dilution	DUP RPD %	DUP Qualifier	DUP RPD Limits %
Perchlorate	11100	10700	500	3.96		15

Laboratory Control Sample (LCS)

(LCS) R4057423-2 04/13/24 21:08

Analyte	Spike Amount ug/l	LCS Result ug/l	LCS Rec. %	Rec. Limits %	LCS Qualifier
Perchlorate	10.0	11.0	110	90.0-110	

L1724110-02 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L1724110-02 04/13/24 23:24 • (MS) R4057423-4 04/13/24 23:53 • (MSD) R4057423-5 04/14/24 00:21

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MSD Result ug/l	MS Rec. %	MSD Rec. %	Dilution	Rec. Limits %	MS Qualifier	MSD Qualifier	RPD %	RPD Limits %
Perchlorate	500000	817000	1330000	1270000	103	90.5	50000	80.0-120			4.63	15

Method Blank (MB)

(MB) R4056937-4 04/11/24 10:10

Analyte	MB Result ug/l	MB Qualifier	MB MDL ug/l	MB RDL ug/l
Acetone	U		11.3	50.0
Acrolein	U		2.54	50.0
Acrylonitrile	U		0.671	10.0
Benzene	U		0.0941	1.00
Bromobenzene	U		0.118	1.00
Bromodichloromethane	U		0.136	1.00
Bromoform	U		0.129	1.00
Bromomethane	U		0.605	5.00
1,3-Butadiene	U		0.299	2.00
n-Butylbenzene	U		0.157	1.00
sec-Butylbenzene	U		0.125	1.00
tert-Butylbenzene	U		0.127	1.00
Carbon tetrachloride	U		0.128	1.00
Carbon disulfide	0.224	E4	0.0962	1.00
Chlorobenzene	U		0.116	1.00
Chlorodibromomethane	U		0.140	1.00
Chloroethane	U		0.192	5.00
Chloroform	U		0.111	5.00
Chloromethane	U		0.960	2.50
Cyclohexane	U		0.188	1.00
2-Chlorotoluene	U		0.106	1.00
4-Chlorotoluene	U		0.114	1.00
1,2-Dibromo-3-Chloropropane	U		0.276	5.00
1,2-Dibromoethane	U		0.126	1.00
Dibromomethane	U		0.122	1.00
1,2-Dichlorobenzene	U		0.107	1.00
1,3-Dichlorobenzene	U		0.110	1.00
1,4-Dichlorobenzene	U		0.120	1.00
Dichlorodifluoromethane	U		0.374	5.00
1,1-Dichloroethane	U		0.100	1.00
1,2-Dichloroethane	U		0.0819	1.00
1,1-Dichloroethene	U		0.188	1.00
cis-1,2-Dichloroethene	U		0.126	1.00
trans-1,2-Dichloroethene	U		0.149	1.00
1,2-Dichloropropane	U		0.149	1.00
1,1-Dichloropropene	U		0.142	1.00
1,3-Dichloropropane	U		0.110	1.00
cis-1,3-Dichloropropene	U		0.111	1.00
trans-1,3-Dichloropropene	U		0.118	1.00
2,2-Dichloropropane	U		0.161	1.00

Method Blank (MB)

(MB) R4056937-4 04/11/24 10:10

Analyte	MB Result ug/l	MB Qualifier	MB MDL ug/l	MB RDL ug/l
Dicyclopentadiene	U		0.253	1.00
Di-isopropyl ether	U		0.105	1.00
Ethylbenzene	U		0.137	1.00
4-Ethyltoluene	U		0.208	1.00
Hexachloro-1,3-butadiene	U		0.337	1.00
n-Hexane	U		0.749	10.0
Isopropylbenzene	U		0.105	1.00
p-Isopropyltoluene	U		0.120	1.00
2-Butanone (MEK)	U		1.19	10.0
Methyl Cyclohexane	U		0.660	1.00
Methylene Chloride	U		0.430	5.00
4-Methyl-2-pentanone (MIBK)	U		0.478	10.0
Methyl tert-butyl ether	U		0.101	1.00
Naphthalene	U		1.00	5.00
Propene	U		0.936	2.50
n-Propylbenzene	U		0.0993	1.00
Styrene	U		0.118	1.00
1,1,1,2-Tetrachloroethane	U		0.147	1.00
1,1,2,2-Tetrachloroethane	U		0.133	1.00
1,1,2-Trichlorotrifluoroethane	U		0.180	1.00
Tetrachloroethene	U		0.300	1.00
Toluene	U		0.278	1.00
1,2,3-Trichlorobenzene	U		0.230	1.00
1,2,4-Trichlorobenzene	U		0.481	1.00
1,1,1-Trichloroethane	U		0.149	1.00
1,1,2-Trichloroethane	U		0.158	1.00
Trichloroethene	U		0.190	1.00
Trichlorofluoromethane	U		0.160	5.00
1,2,3-Trichloropropane	U		0.237	2.50
1,2,4-Trimethylbenzene	U		0.322	1.00
1,2,3-Trimethylbenzene	U		0.104	1.00
1,3,5-Trimethylbenzene	U		0.104	1.00
Vinyl chloride	U		0.234	1.00
Xylenes, Total	U		0.174	3.00
(S) Toluene-d8	*108*			*80.0-120*
(S) 4-Bromofluorobenzene	*93.3*			*77.0-126*
(S) 1,2-Dichloroethane-d4	*86.3*			*70.0-130*

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R4056937-1 04/11/24 08:48 • (LCSD) R4056937-2 04/11/24 09:08

Analyte	Spike Amount ug/l	LCS Result ug/l	LCSD Result ug/l	LCS Rec. %	LCSD Rec. %	Rec. Limits %	LCS Qualifier	LCSD Qualifier	RPD %	RPD Limits %
Acetone	25.0	23.5	21.8	94.0	87.2	19.0-160			7.51	27
Acrolein	25.0	15.9	15.8	63.6	63.2	30.0-160			0.631	26
Acrylonitrile	25.0	17.9	17.4	71.6	69.6	55.0-149			2.83	20
Benzene	5.00	4.63	4.60	92.6	92.0	70.0-123			0.650	20
Bromobenzene	5.00	4.71	4.79	94.2	95.8	73.0-121			1.68	20
Bromodichloromethane	5.00	4.68	4.73	93.6	94.6	75.0-120			1.06	20
Bromoform	5.00	4.35	4.32	87.0	86.4	68.0-132			0.692	20
Bromomethane	5.00	1.82	1.72	36.4	34.4	30.0-160			5.65	25
1,3-Butadiene	5.00	3.06	2.92	61.2	58.4	45.0-147			4.68	20
n-Butylbenzene	5.00	5.24	5.36	105	107	73.0-125			2.26	20
sec-Butylbenzene	5.00	5.02	4.92	100	98.4	75.0-125			2.01	20
tert-Butylbenzene	5.00	5.28	5.09	106	102	76.0-124			3.66	20
Carbon tetrachloride	5.00	5.29	4.90	106	98.0	68.0-126			7.65	20
Carbon disulfide	5.00	4.48	4.05	89.6	81.0	61.0-128			10.1	20
Chlorobenzene	5.00	6.04	6.01	121	120	80.0-121			0.498	20
Chlorodibromomethane	5.00	5.43	5.52	109	110	77.0-125			1.64	20
Chloroethane	5.00	4.45	4.05	89.0	81.0	47.0-150			9.41	20
Chloroform	5.00	4.85	4.49	97.0	89.8	73.0-120			7.71	20
Chloromethane	5.00	5.54	5.48	111	110	41.0-142			1.09	20
Cyclohexane	5.00	4.52	4.43	90.4	88.6	71.0-124			2.01	20
2-Chlorotoluene	5.00	5.17	5.08	103	102	76.0-123			1.76	20
4-Chlorotoluene	5.00	5.31	5.20	106	104	75.0-122			2.09	20
1,2-Dibromo-3-Chloropropane	5.00	3.91	4.53	78.2	90.6	58.0-134			14.7	20
1,2-Dibromoethane	5.00	5.54	5.21	111	104	80.0-122			6.14	20
Dibromomethane	5.00	4.89	4.86	97.8	97.2	80.0-120			0.615	20
1,2-Dichlorobenzene	5.00	5.25	5.36	105	107	79.0-121			2.07	20
1,3-Dichlorobenzene	5.00	5.52	5.42	110	108	79.0-120			1.83	20
1,4-Dichlorobenzene	5.00	5.27	5.09	105	102	79.0-120			3.47	20
Dichlorodifluoromethane	5.00	4.57	4.28	91.4	85.6	51.0-149			6.55	20
1,1-Dichloroethane	5.00	4.68	4.45	93.6	89.0	70.0-126			5.04	20
1,2-Dichloroethane	5.00	4.26	4.24	85.2	84.8	70.0-128			0.471	20
1,1-Dichloroethene	5.00	4.91	4.53	98.2	90.6	71.0-124			8.05	20
cis-1,2-Dichloroethene	5.00	5.41	5.03	108	101	73.0-120			7.28	20
trans-1,2-Dichloroethene	5.00	5.16	4.81	103	96.2	73.0-120			7.02	20
1,2-Dichloropropane	5.00	4.89	4.57	97.8	91.4	77.0-125			6.77	20
1,1-Dichloropropene	5.00	4.82	4.65	96.4	93.0	74.0-126			3.59	20
1,3-Dichloropropane	5.00	5.14	5.00	103	100	80.0-120			2.76	20
cis-1,3-Dichloropropene	5.00	4.77	4.45	95.4	89.0	80.0-123			6.94	20
trans-1,3-Dichloropropene	5.00	4.80	4.74	96.0	94.8	78.0-124			1.26	20
2,2-Dichloropropane	5.00	4.70	4.31	94.0	86.2	58.0-130			8.66	20

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R4056937-1 04/11/24 08:48 • (LCSD) R4056937-2 04/11/24 09:08

Analyte	Spike Amount ug/l	LCS Result ug/l	LCSD Result ug/l	LCS Rec. %	LCSD Rec. %	Rec. Limits %	LCS Qualifier	LCSD Qualifier	RPD %	RPD Limits %
Dicyclopentadiene	5.00	4.42	4.27	88.4	85.4	74.0-126			3.45	20
Di-isopropyl ether	5.00	4.08	4.07	81.6	81.4	58.0-138			0.245	20
Ethylbenzene	5.00	5.61	5.45	112	109	79.0-123			2.89	20
4-Ethyltoluene	5.00	5.23	5.06	105	101	74.0-127			3.30	20
Hexachloro-1,3-butadiene	5.00	4.46	4.44	89.2	88.8	54.0-138			0.449	20
n-Hexane	5.00	4.40	3.87	88.0	77.4	57.0-133			12.8	20
Isopropylbenzene	5.00	5.38	5.29	108	106	76.0-127			1.69	20
p-Isopropyltoluene	5.00	5.26	5.05	105	101	76.0-125			4.07	20
2-Butanone (MEK)	25.0	20.8	19.7	83.2	78.8	44.0-160			5.43	20
Methyl Cyclohexane	5.00	4.26	4.16	85.2	83.2	68.0-126			2.38	20
Methylene Chloride	5.00	4.66	4.43	93.2	88.6	67.0-120			5.06	20
4-Methyl-2-pentanone (MIBK)	25.0	19.5	19.6	78.0	78.4	68.0-142			0.512	20
Methyl tert-butyl ether	5.00	4.30	4.27	86.0	85.4	68.0-125			0.700	20
Naphthalene	5.00	3.79	4.23	75.8	84.6	54.0-135			11.0	20
Propene	5.00	2.03	1.83	40.6	36.6	30.0-160			10.4	20
n-Propylbenzene	5.00	5.07	4.85	101	97.0	77.0-124			4.44	20
Styrene	5.00	5.20	5.38	104	108	73.0-130			3.40	20
1,1,1,2-Tetrachloroethane	5.00	5.79	5.62	116	112	75.0-125			2.98	20
1,1,2,2-Tetrachloroethane	5.00	4.69	4.72	93.8	94.4	65.0-130			0.638	20
1,1,2-Trichlorotrifluoroethane	5.00	4.70	4.58	94.0	91.6	69.0-132			2.59	20
Tetrachloroethene	5.00	5.92	5.78	118	116	72.0-132			2.39	20
Toluene	5.00	5.64	5.32	113	106	79.0-120			5.84	20
1,2,3-Trichlorobenzene	5.00	4.53	5.28	90.6	106	50.0-138			15.3	20
1,2,4-Trichlorobenzene	5.00	4.36	4.26	87.2	85.2	57.0-137			2.32	20
1,1,1-Trichloroethane	5.00	4.97	4.73	99.4	94.6	73.0-124			4.95	20
1,1,2-Trichloroethane	5.00	5.34	5.31	107	106	80.0-120			0.563	20
Trichloroethene	5.00	5.79	5.35	116	107	78.0-124			7.90	20
Trichlorofluoromethane	5.00	5.46	5.17	109	103	59.0-147			5.46	20
1,2,3-Trichloropropane	5.00	5.15	5.22	103	104	73.0-130			1.35	20
1,2,4-Trimethylbenzene	5.00	5.04	4.98	101	99.6	76.0-121			1.20	20
1,2,3-Trimethylbenzene	5.00	5.12	5.07	102	101	77.0-120			0.981	20
1,3,5-Trimethylbenzene	5.00	4.94	4.82	98.8	96.4	76.0-122			2.46	20
Vinyl chloride	5.00	4.52	4.09	90.4	81.8	67.0-131			9.99	20
Xylenes, Total	15.0	16.8	16.0	112	107	79.0-123			4.88	20
(S) Toluene-d8				*107*	*107*	*80.0-120*				
(S) 4-Bromofluorobenzene				*94.1*	*94.5*	*77.0-126*				
(S) 1,2-Dichloroethane-d4				*84.7*	*88.4*	*70.0-130*				

INTERNAL STANDARD SUMMARY

Instrument: VOCMS36 • File ID: 0411_02

04/11/24 08:48

Sample ID	File ID	8260-FLUOROBENZENE Response	8260-CHLOROBENZENE-D5 Response	8260-1,4-DICHLOROBENZENE-D4 Response
Standard	0411_02	526369	221954	196760
Upper Limit		1052738	443908	393520
Lower Limit		263185	110977	98380
LCS R4056937-1 WG2264870 1x	0411_02LCS	526369	221954	196760
LCSD R4056937-2 WG2264870 1x	0411_03	532817	222295	195924
BLANK R4056937-4 WG2264870 1x	0411_06A	518347	211264	182116
L1724110-03 WG2264870 1x	0411_20	499717	205677	175827
L1724110-01 WG2264870 20x	0411_30	503711	206161	175212
L1724110-02 WG2264870 2500x	0411_31	504283	204289	174603

GLOSSARY OF TERMS

Guide to Reading and Understanding Your Laboratory Report

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

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

Abbreviations and Definitions

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

Qualifier	Description
B1	Target analyte detected in method blank at or above the method reporting limit.
E4	Concentration estimated. Analyte was detected below laboratory minimum reporting level (MRL) but above MDL.

ACCREDITATIONS & LOCATIONS

Pace Analytical National 12065 Lebanon Rd Mount Juliet, TN 37122

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

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

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

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

Pinyon Environmental
3222 S. Vance St. Suite 200
Lakewood, CO 80227

Billing Information:
accounts Payable
ap@pinyon-env.com

Pres Chk

Report to: Andrew Parker

Email To: Parker@Pinyon-env.com

Project Description: Nammo TTU Groundwater

City/State Collected: Mesa, Az

Please Circle: PT MT CT ET

Phone: 303.785.7687

Client Project # 722152201.002

Lab Project #

Collected by (print): Crissa Wolf

Site/Facility ID #

P.O. #

Collected by (signature): Crissa Wolf

Immediately Packed on Ice N ___ Y X

Rush? (Lab MUST Be Notified)
___ Same Day X Five Day
___ Next Day ___ 5 Day (Rad Only)
___ Two Day ___ 10 Day (Rad Only)
___ Three Day

Quote #

Date Results Needed

Analysis / Container / Preservative

Sample ID	Comp/Grab	Matrix*	Depth	Date	Time	No. of Cntrs	VOC 8260 AZ/40ml amber HCl	Perchlorate 314.0/125ml	Remarks	Sample # (lab only)
TTU-15-75-20240409	Grab	GW	75	4-9-24	1046	4	X	X		-01
TTU-16-80-20240409	1	1	80	4-9-24	1026	4	X	X		-02
Trip Blank	—	—	—	—	—	1	X	X		-03

Chain of Custody Page 1 of 1

Pace Analytical®
National Center for Testing & Innovation

12065 Lebanon Rd
Mount Juliet, TN 37122
Phone: 615-758-5858
Phone: 800-767-5859
Fax: 615-758-5859

SDG # L1724110

D048

Acctnum: PINYONMAZ

Template:

Prelogin:

PM:

PB:

Shipped Via:

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

Remarks:

pH _____ Temp _____

Flow _____ Other _____

Samples returned via:
__ UPS __ FedEx __ Courier __ _____

Tracking # 724107050123

Sample Receipt Checklist

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

	Date:	Time:		
Relinquished by: (Signature) Crissa Wolf	4-9-24	1200	Received by: (Signature)	Trip Blank Received: (Yes) / No HCL / MeoH TBR: 1
Relinquished by: (Signature)	4-9-24	1800	Received by: (Signature) SWA	Temp: DPT °C 3.8+0.1=3.9 Bottles Received: 8
Relinquished by: (Signature)	Date:	Time:	Received for lab by: (Signature) Devanty	Date: 4/10/24 Time: 0915

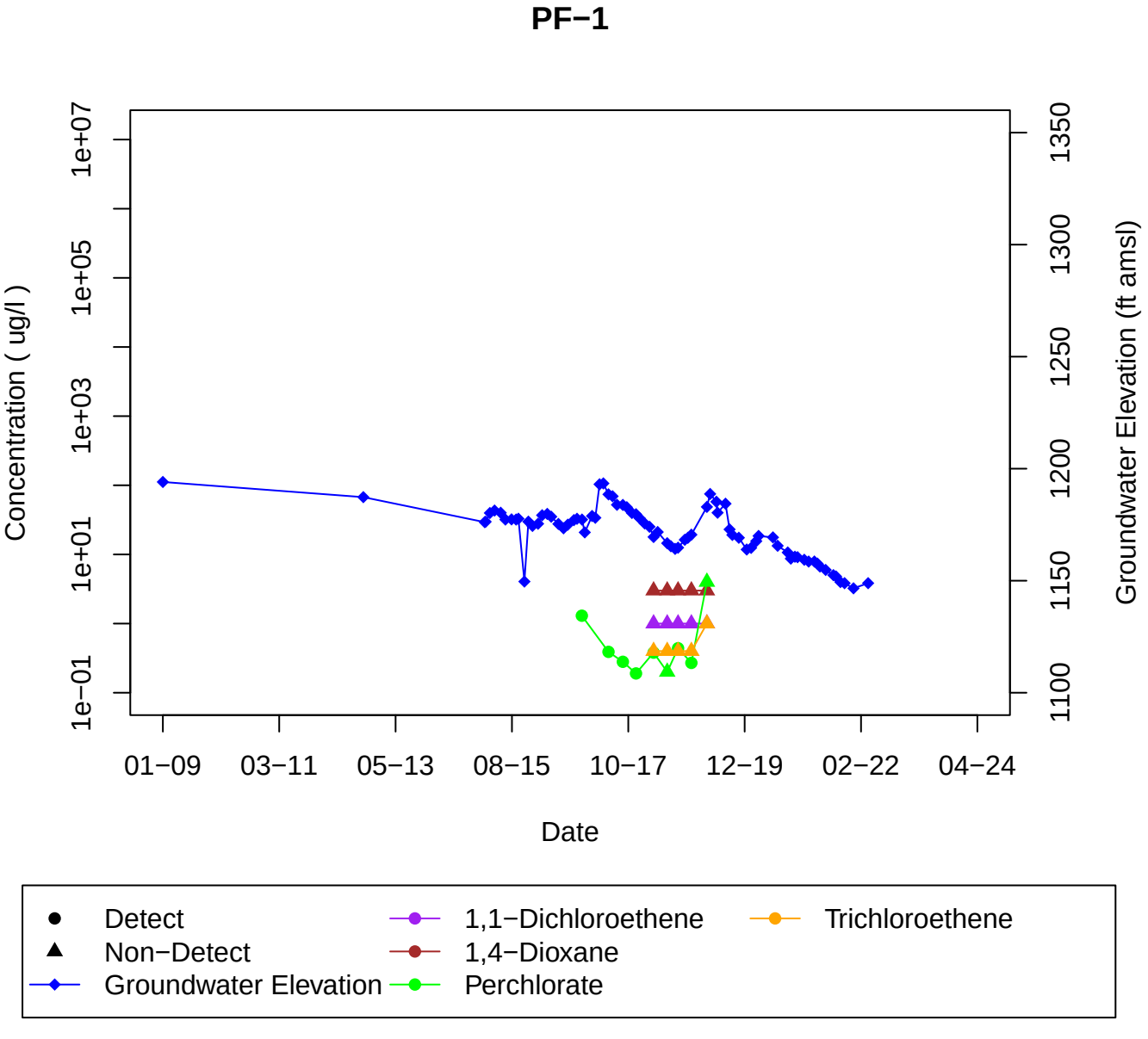
If preservation required by Login: Date/Time

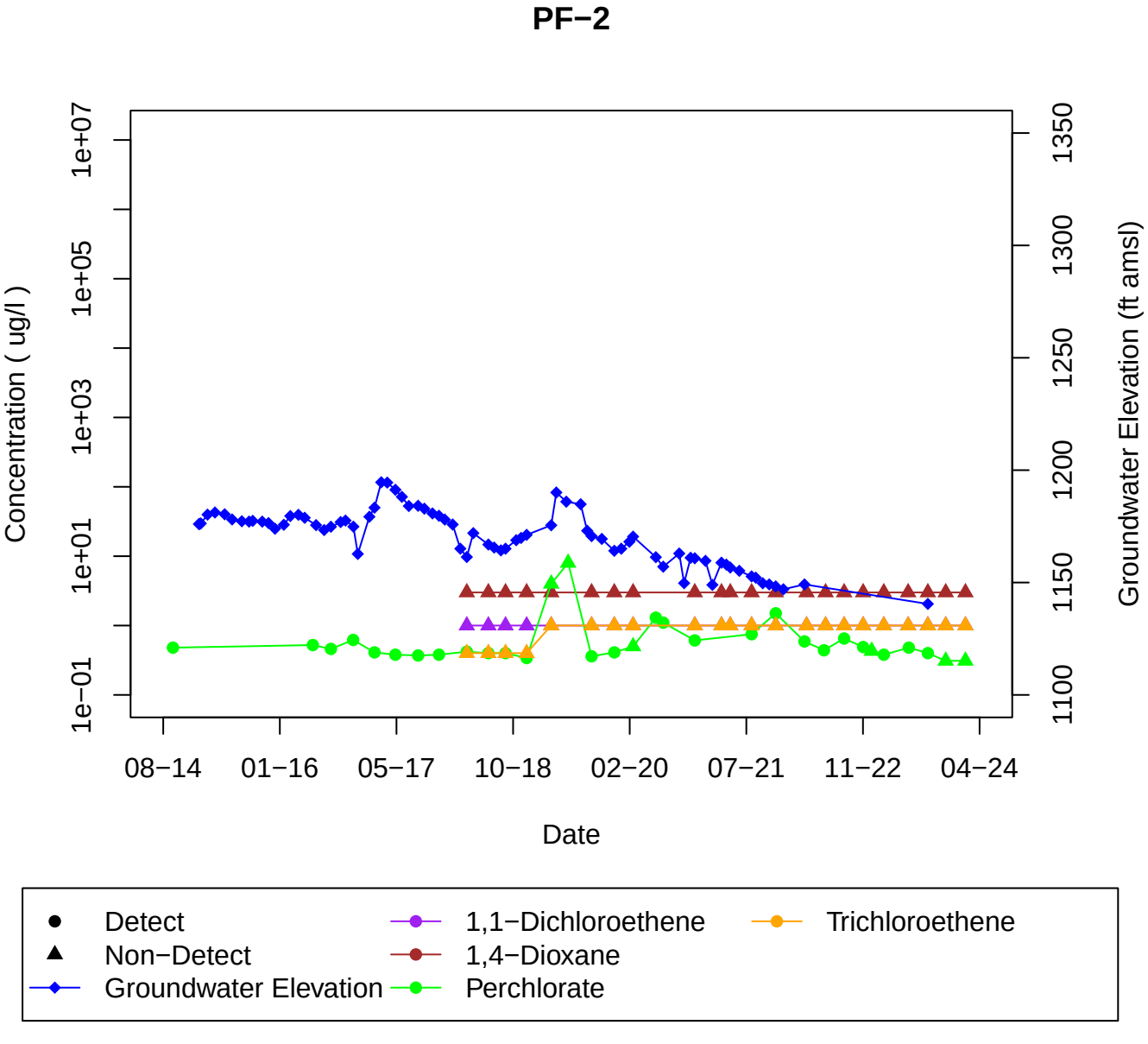
Hold:

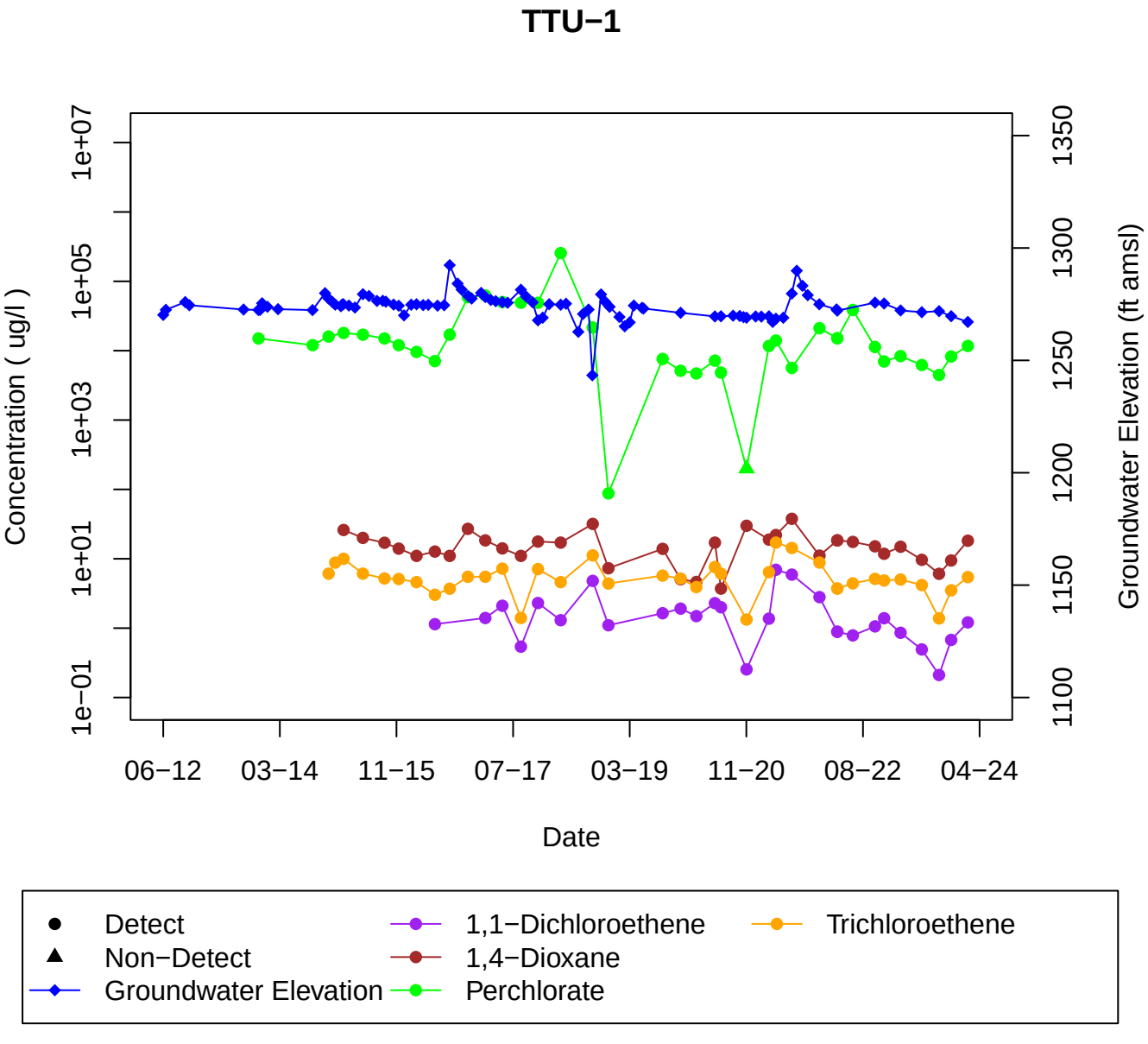
Condition: NCF / OK

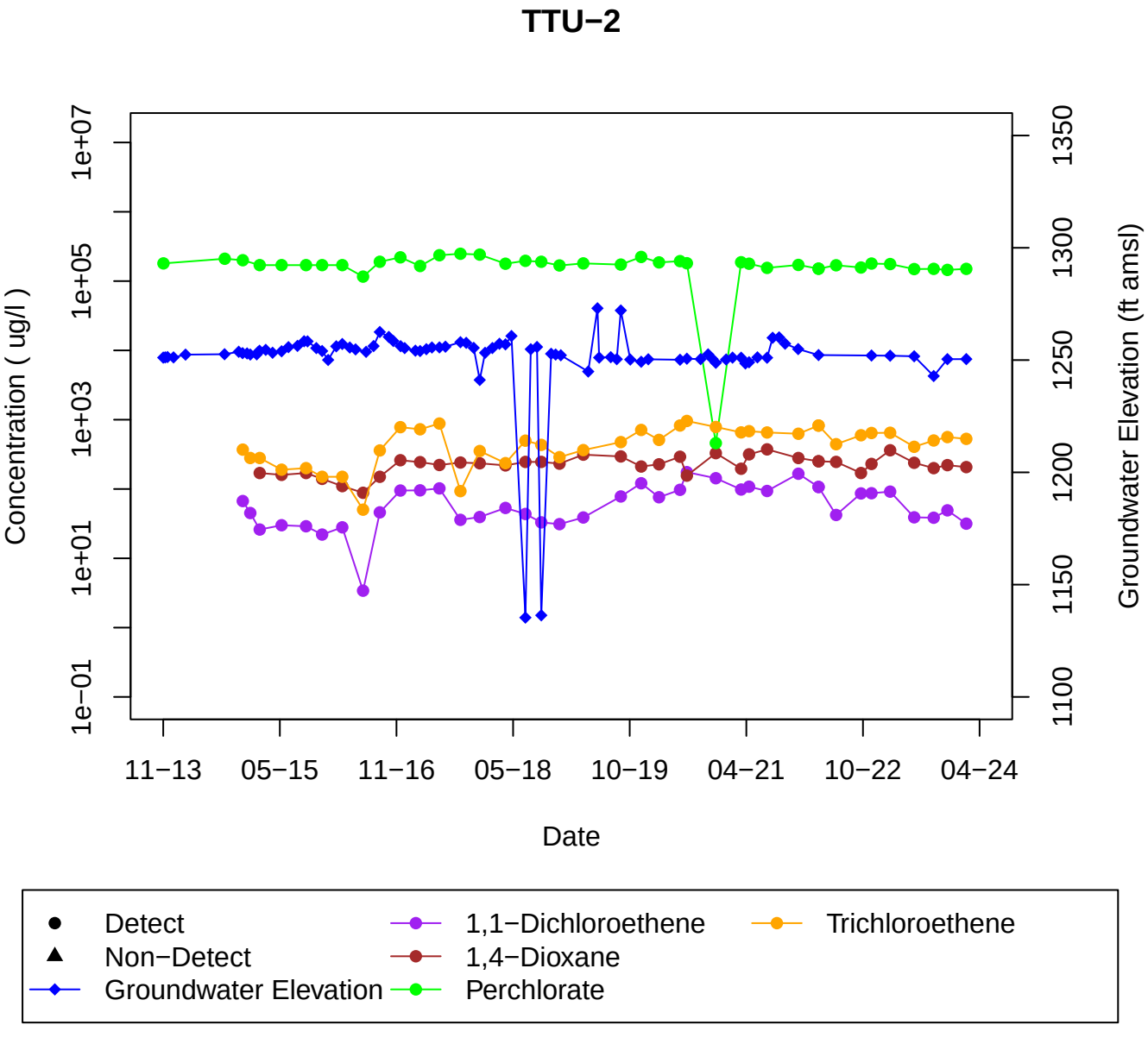
AJPAZ

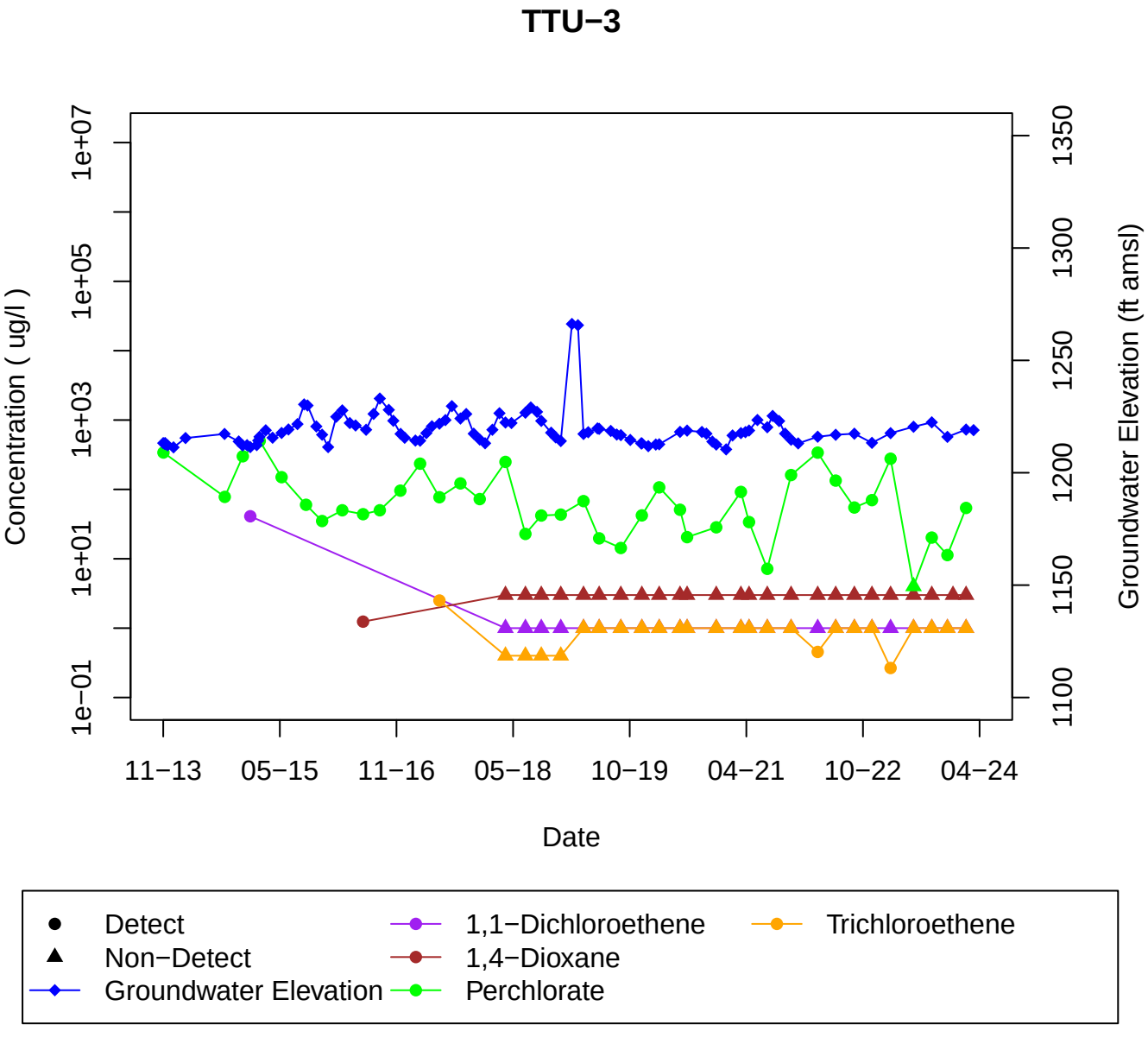
Attachment 3 – Concentration and Groundwater Elevation versus Time Plots

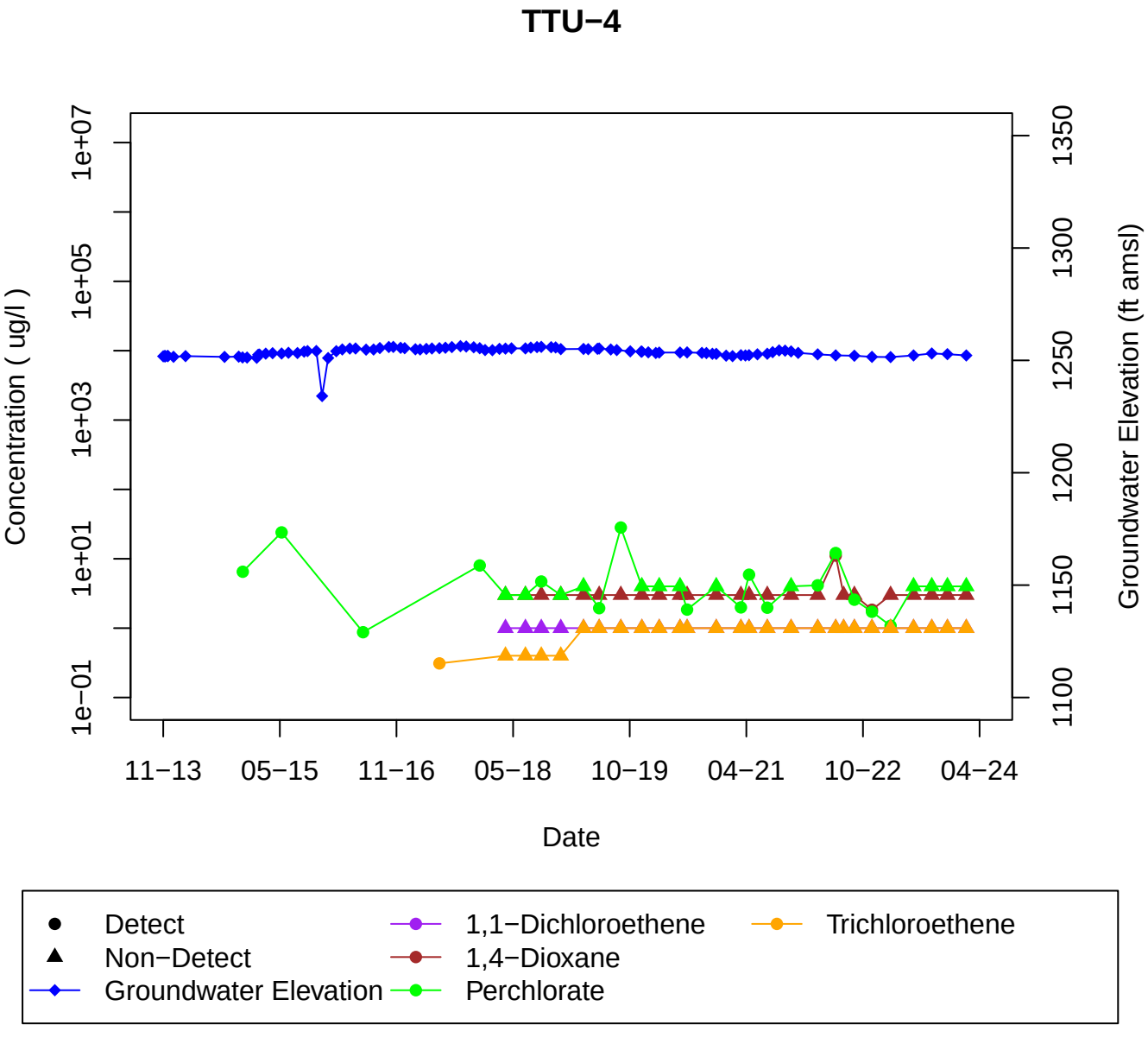

PF−1
Concentration (ug/l)
1e−01
1e+01
1e+03
1e+05
1e+07
Groundwater Elevation (ft amsl)
1100
1150
1200
1250
1300
1350
01−09
03−11
05−13
08−15
10−17
12−19
02−22
04−24
Date
Detect
Non−Detect
Groundwater Elevation
1,1−Dichloroethene
1,4−Dioxane
Perchlorate
Trichloroethene

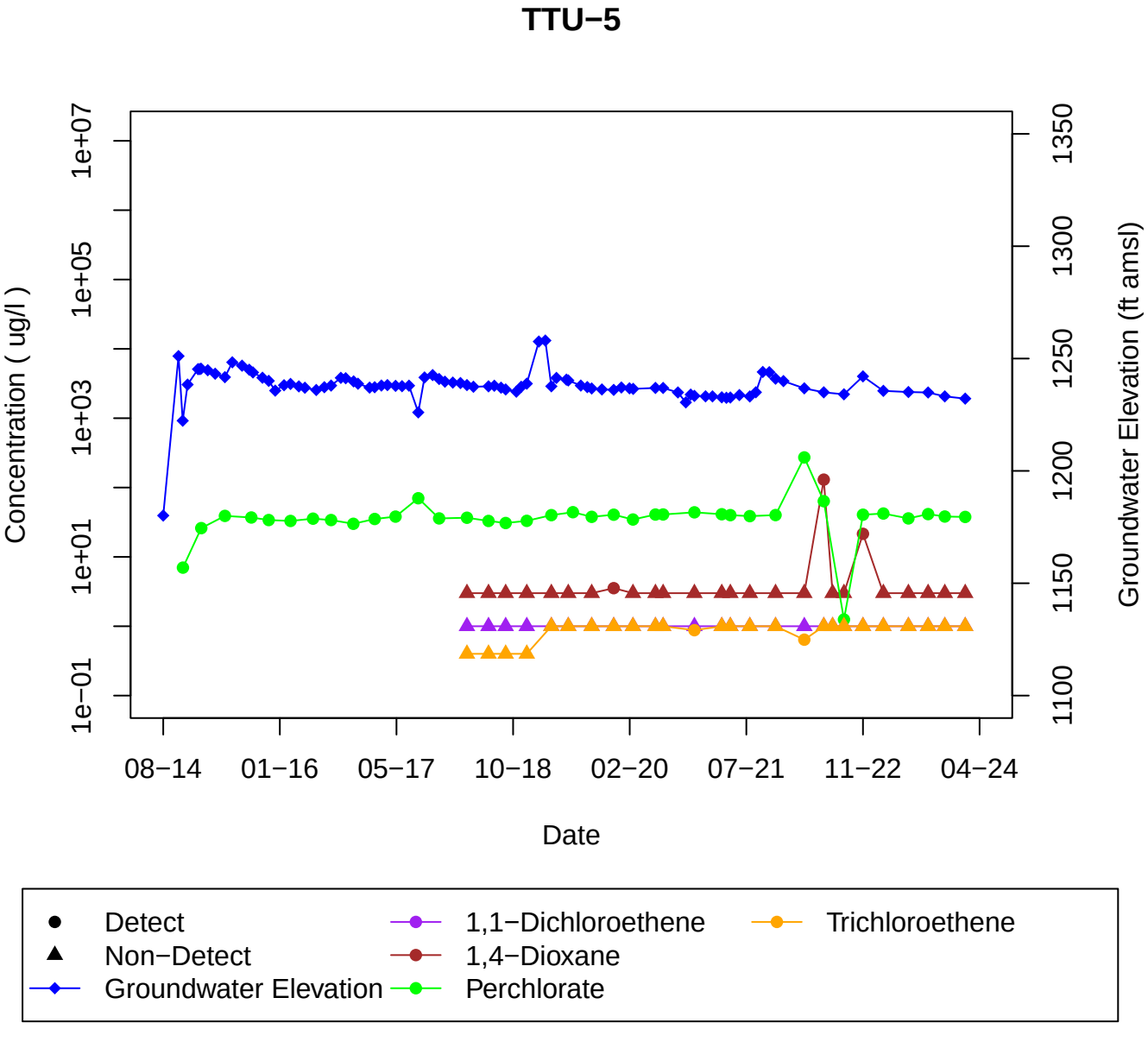

PF−2
Concentration (ug/l)
1e−01
1e+01
1e+03
1e+05
1e+07
Groundwater Elevation (ft amsl)
1100
1150
1200
1250
1300
1350
08−14
01−16
05−17
10−18
02−20
07−21
11−22
04−24
Date
Detect
Non−Detect
Groundwater Elevation
1,1−Dichloroethene
1,4−Dioxane
Perchlorate
Trichloroethene

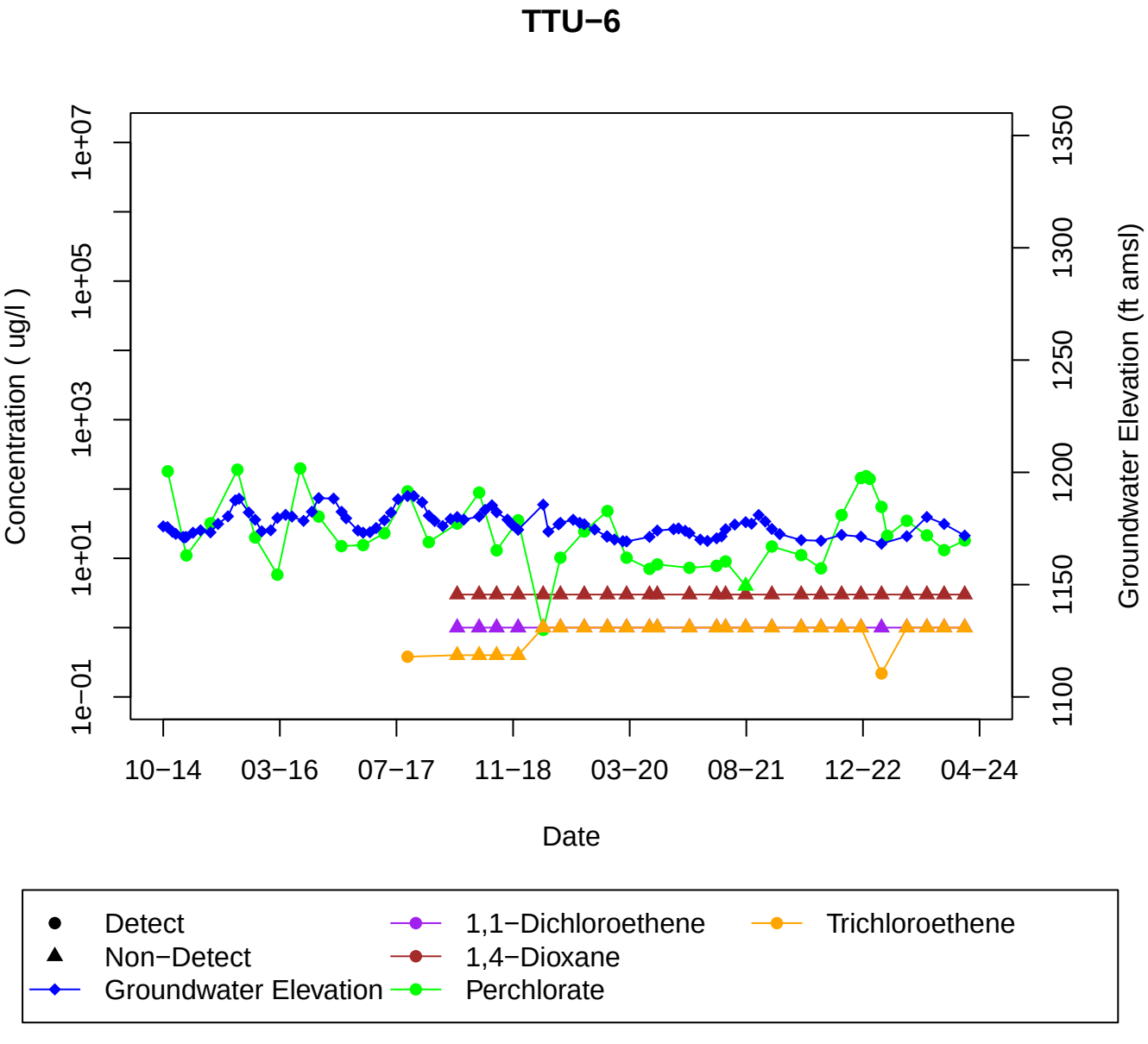

TTU−1
Concentration (ug/l)
1e−01
1e+01
1e+03
1e+05
1e+07
Groundwater Elevation (ft amsl)
1100
1150
1200
1250
1300
1350
06−12
03−14
11−15
07−17
03−19
11−20
08−22
04−24
Date
Detect
Non−Detect
Groundwater Elevation
1,1−Dichloroethene
1,4−Dioxane
Perchlorate
Trichloroethene

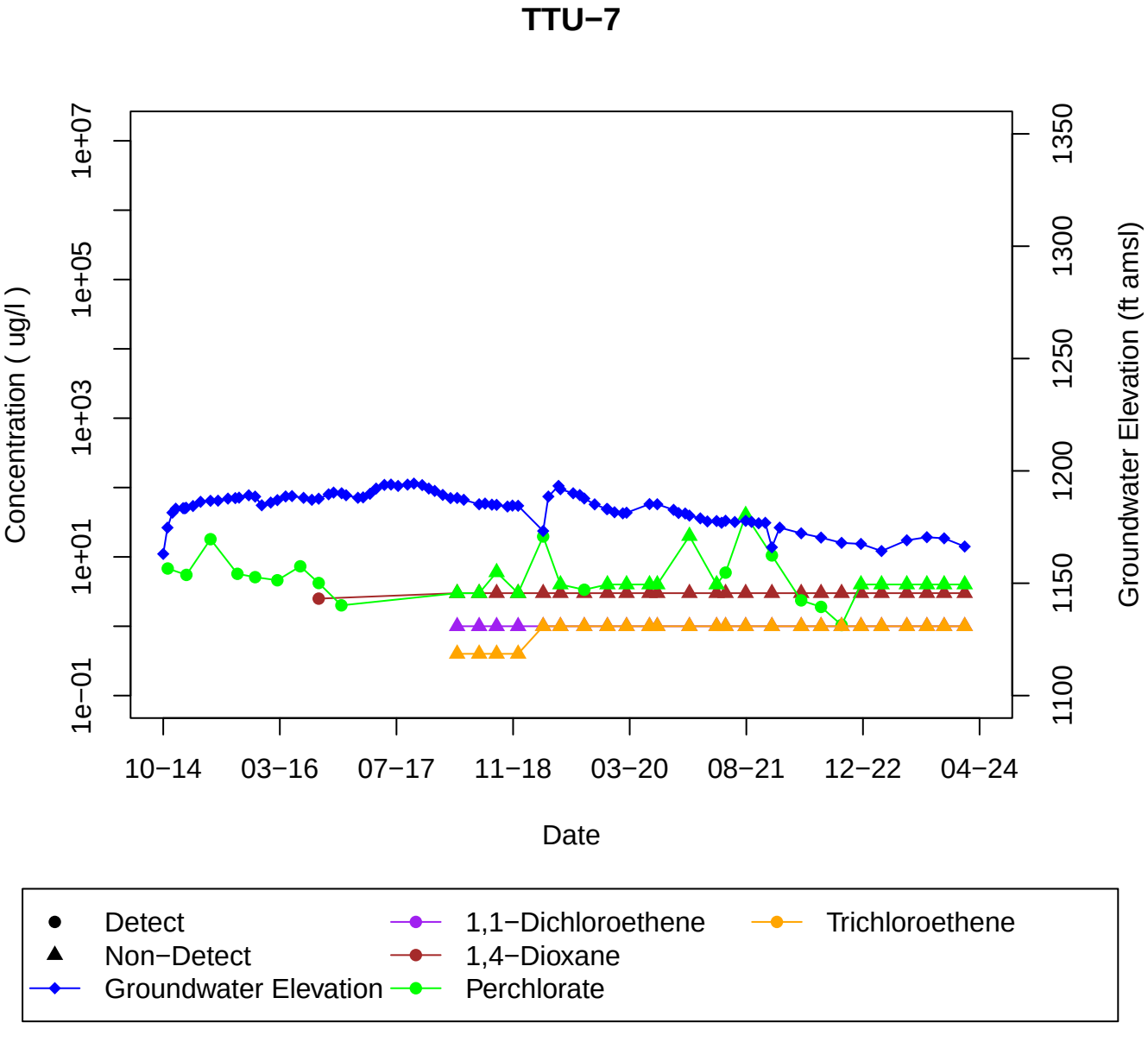

TTU−2
Concentration (ug/l)
1e−01
1e+01
1e+03
1e+05
1e+07
Groundwater Elevation (ft amsl)
1100
1150
1200
1250
1300
1350
11−13
05−15
11−16
05−18
10−19
04−21
10−22
04−24
Date
Detect
Non−Detect
Groundwater Elevation
1,1−Dichloroethene
1,4−Dioxane
Perchlorate
Trichloroethene

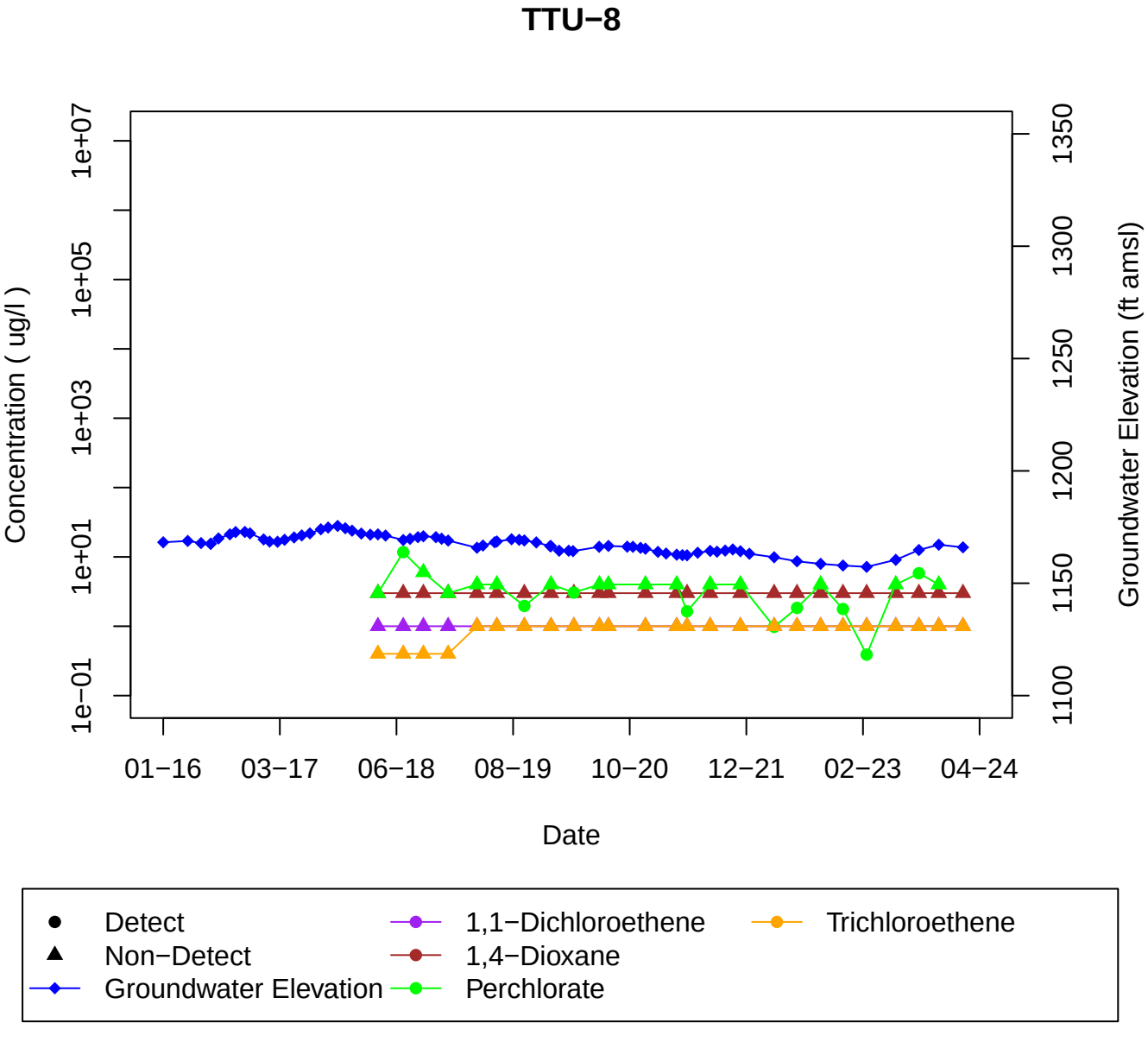

TTU−3
Concentration (ug/l)
1e−01
1e+01
1e+03
1e+05
1e+07
Groundwater Elevation (ft amsl)
1100
1150
1200
1250
1300
1350
11−13
05−15
11−16
05−18
10−19
04−21
10−22
04−24
Date
Detect
Non−Detect
Groundwater Elevation
1,1−Dichloroethene
1,4−Dioxane
Perchlorate
Trichloroethene

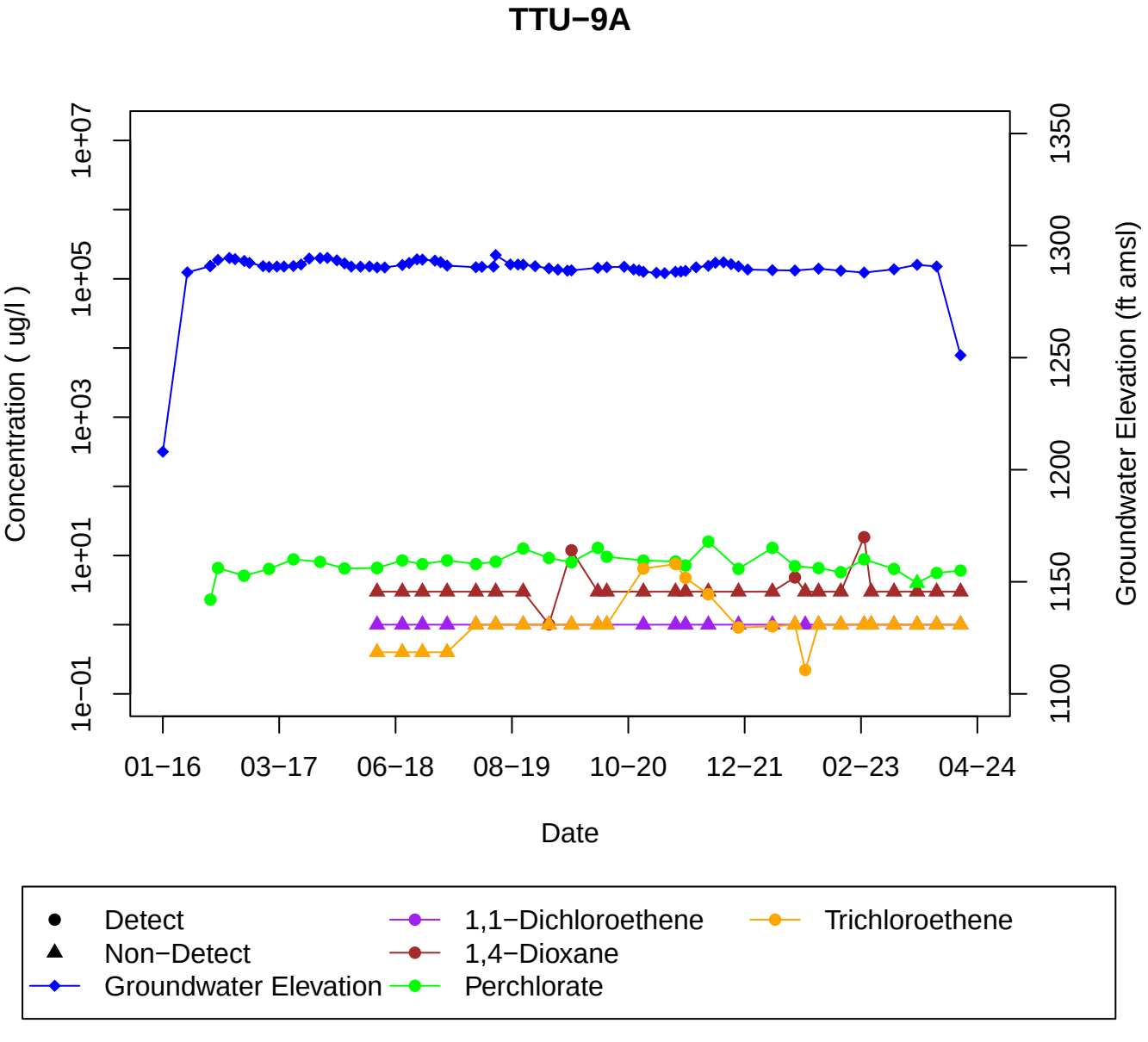

TTU−4
Concentration (ug/l)
1e−01
1e+01
1e+03
1e+05
1e+07
Groundwater Elevation (ft amsl)
1100
1150
1200
1250
1300
1350
11−13
05−15
11−16
05−18
10−19
04−21
10−22
04−24
Date
Detect
Non−Detect
Groundwater Elevation
1,1−Dichloroethene
1,4−Dioxane
Perchlorate
Trichloroethene

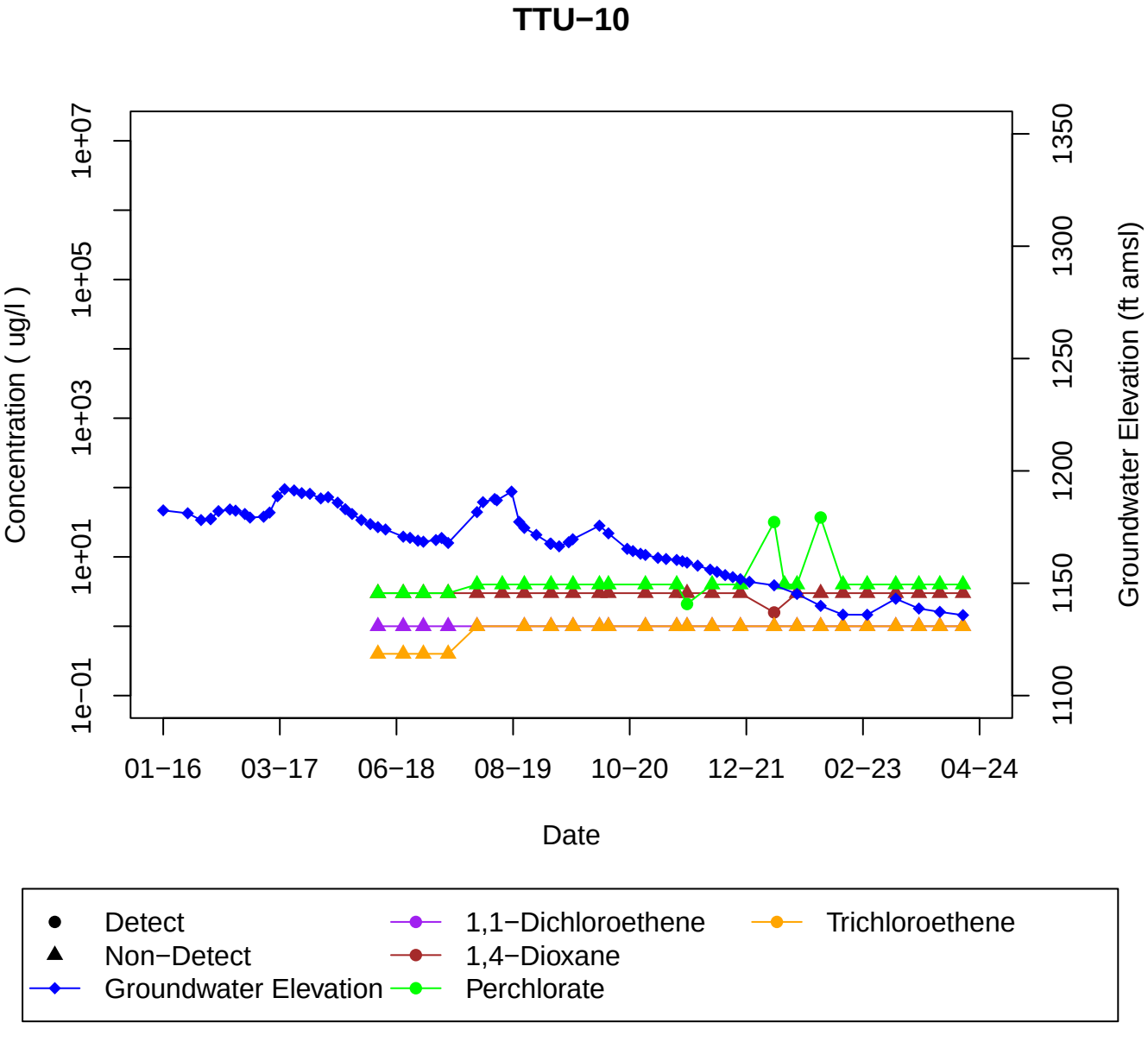

TTU−5
Concentration (ug/l)
1e−01
1e+01
1e+03
1e+05
1e+07
Groundwater Elevation (ft amsl)
1100
1150
1200
1250
1300
1350
08−14
01−16
05−17
10−18
02−20
07−21
11−22
04−24
Date
Detect
Non−Detect
Groundwater Elevation
1,1−Dichloroethene
1,4−Dioxane
Perchlorate
Trichloroethene

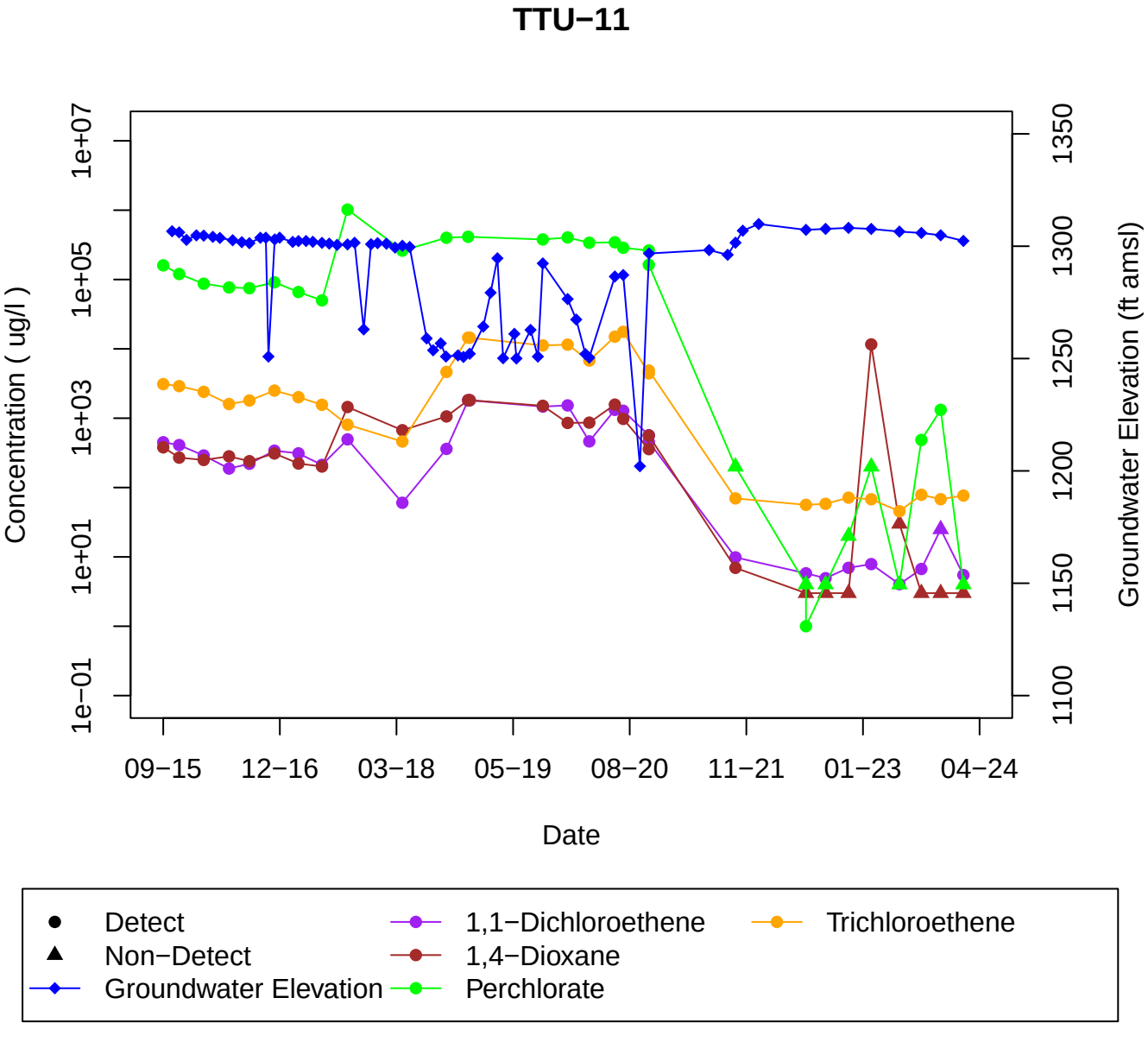

TTU−6
Concentration (ug/l)
1e−01
1e+01
1e+03
1e+05
1e+07
Groundwater Elevation (ft amsl)
1100
1150
1200
1250
1300
1350
10−14
03−16
07−17
11−18
03−20
08−21
12−22
04−24
Date
Detect
Non−Detect
Groundwater Elevation
1,1−Dichloroethene
1,4−Dioxane
Perchlorate
Trichloroethene

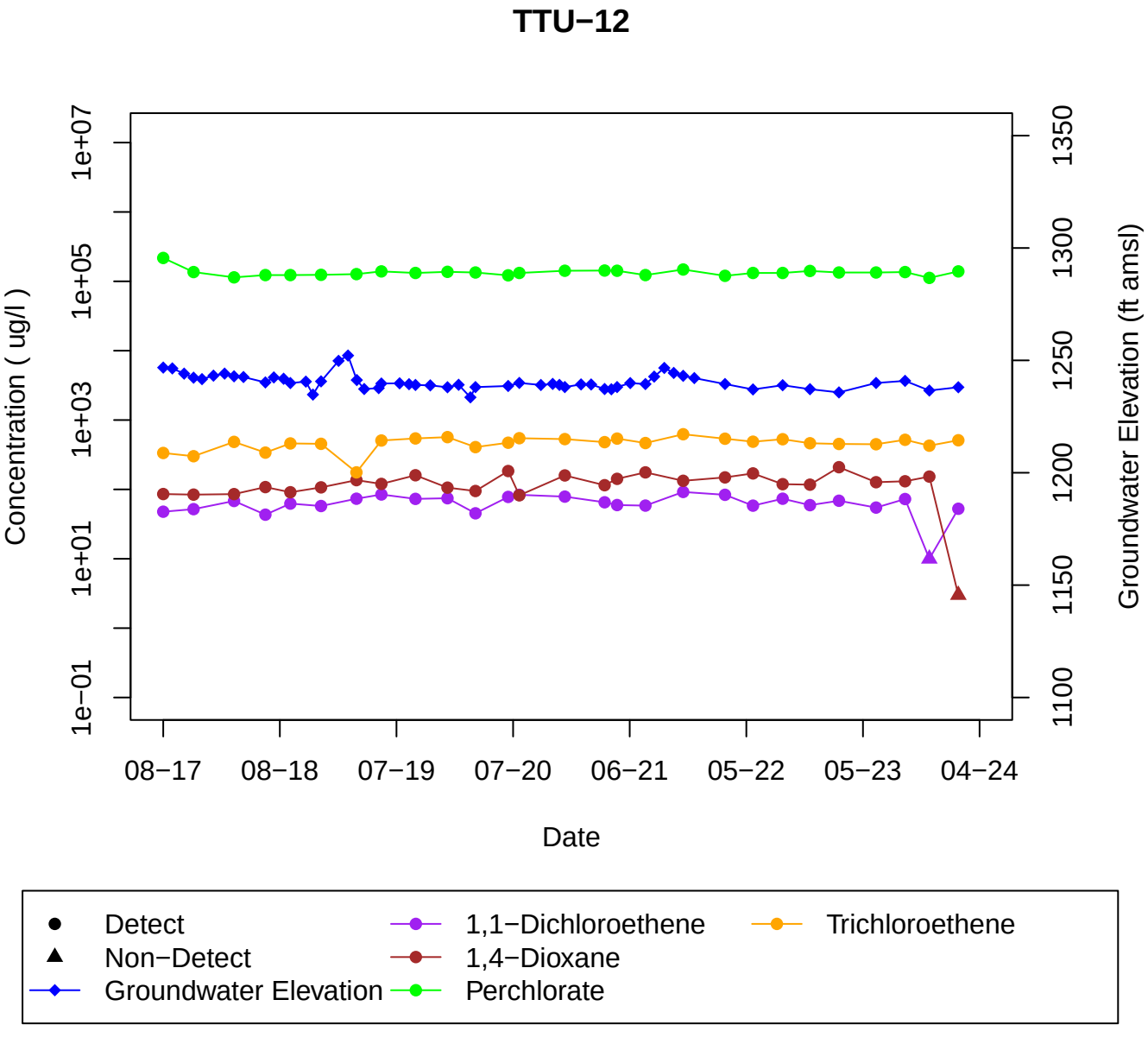

TTU−7
Concentration (ug/l)
1e−01
1e+01
1e+03
1e+05
1e+07
Groundwater Elevation (ft amsl)
1100
1150
1200
1250
1300
1350
10−14
03−16
07−17
11−18
03−20
08−21
12−22
04−24
Date
Detect
Non−Detect
Groundwater Elevation
1,1−Dichloroethene
1,4−Dioxane
Perchlorate
Trichloroethene

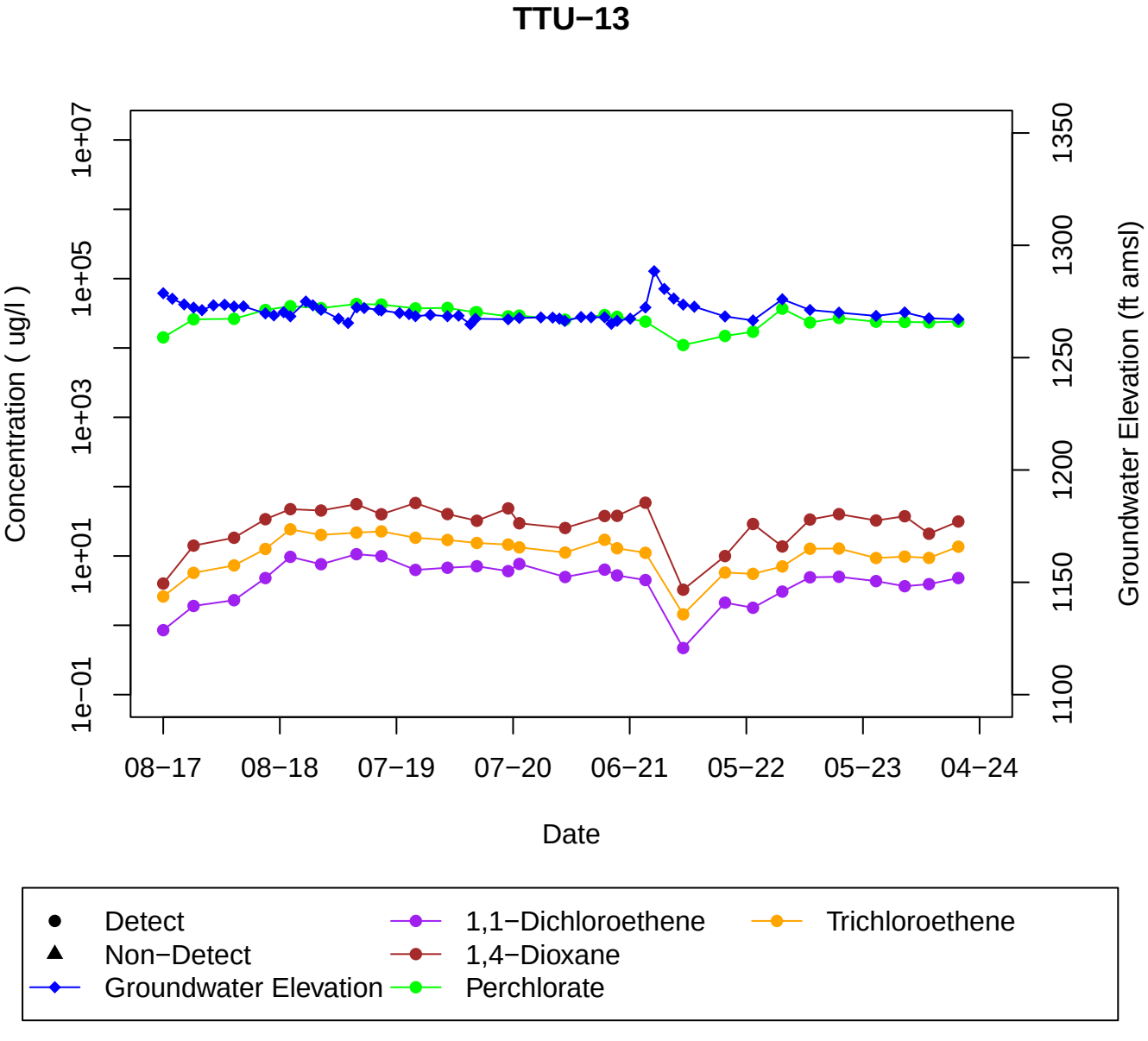

TTU−8
Concentration (ug/l)
1e−01
1e+01
1e+03
1e+05
1e+07
Groundwater Elevation (ft amsl)
1100
1150
1200
1250
1300
1350
01−16
03−17
06−18
08−19
10−20
12−21
02−23
04−24
Date
Detect
Non−Detect
Groundwater Elevation
1,1−Dichloroethene
1,4−Dioxane
Perchlorate
Trichloroethene

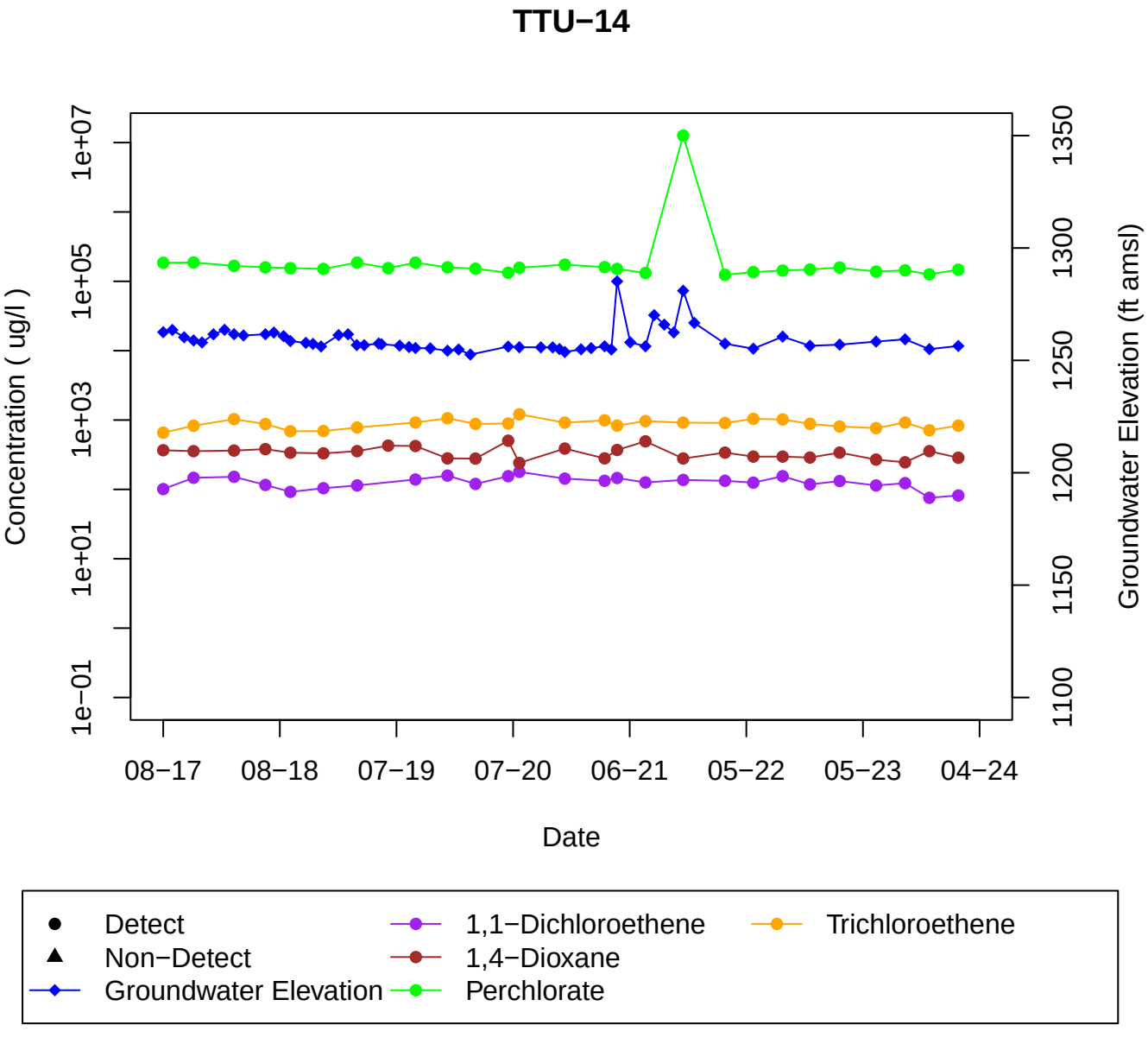

TTU−9A
Concentration (ug/l)
1e−01
1e+01
1e+03
1e+05
1e+07
Groundwater Elevation (ft amsl)
1100
1150
1200
1250
1300
1350
01−16
03−17
06−18
08−19
10−20
12−21
02−23
04−24
Date
Detect
Non−Detect
Groundwater Elevation
1,1−Dichloroethene
1,4−Dioxane
Perchlorate
Trichloroethene

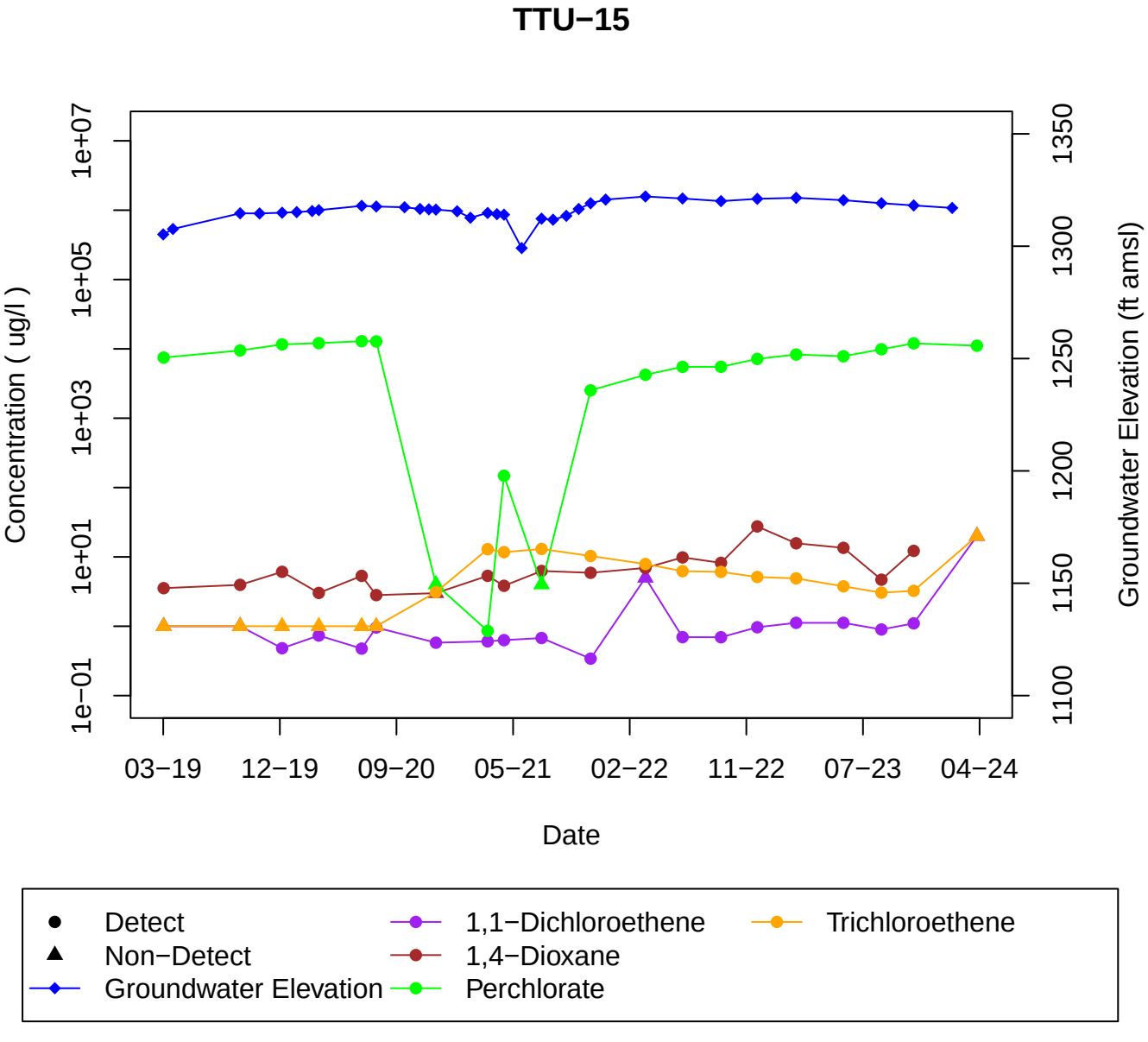

TTU−10
Concentration (ug/l)
1e−01
1e+01
1e+03
1e+05
1e+07
Groundwater Elevation (ft amsl)
1100
1150
1200
1250
1300
1350
01−16
03−17
06−18
08−19
10−20
12−21
02−23
04−24
Date
Detect
Non−Detect
Groundwater Elevation
1,1−Dichloroethene
1,4−Dioxane
Perchlorate
Trichloroethene

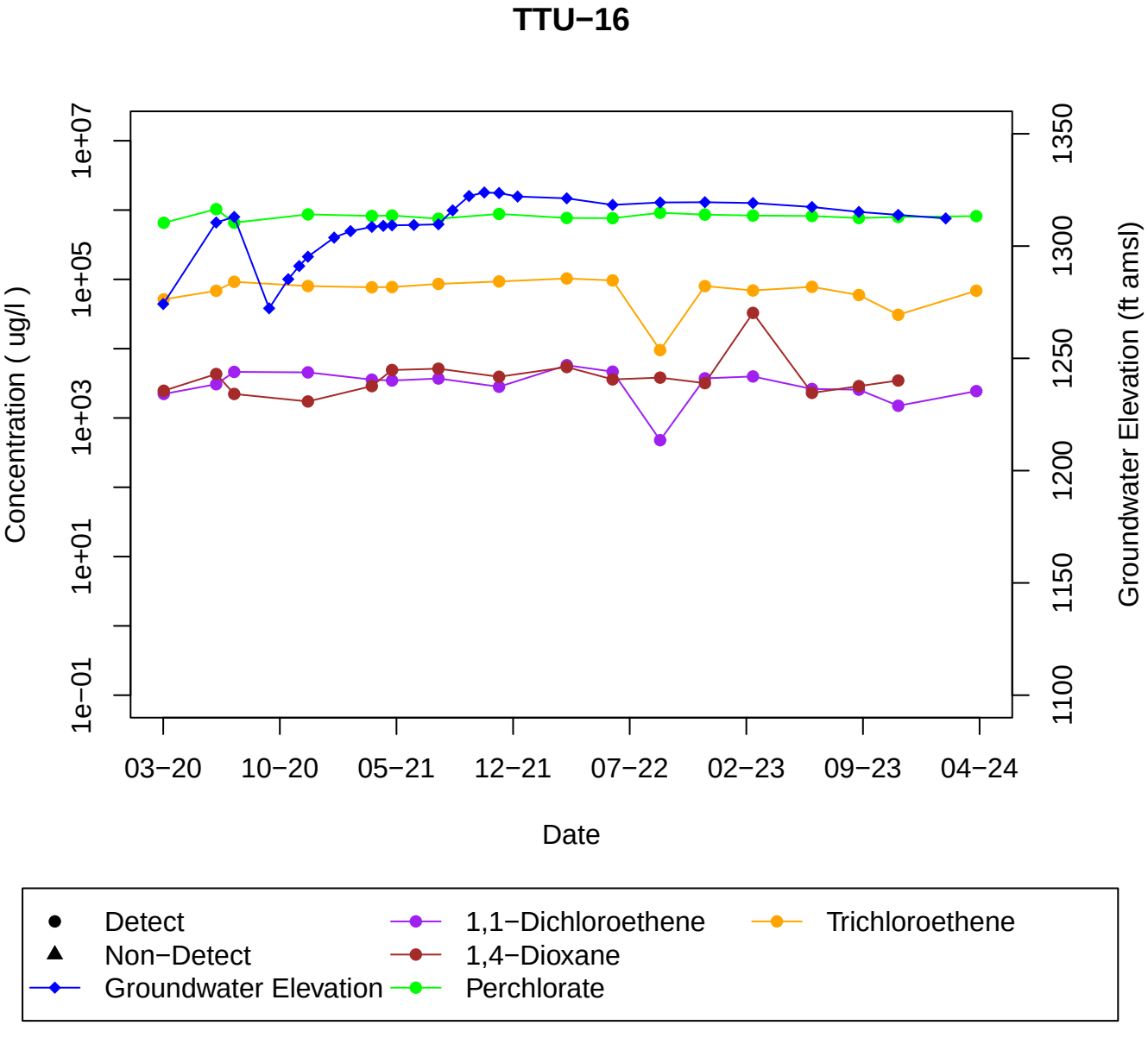

TTU−11
Concentration (ug/l)
1e−01
1e+01
1e+03
1e+05
1e+07
Groundwater Elevation (ft amsl)
1100
1150
1200
1250
1300
1350
09−15
12−16
03−18
05−19
08−20
11−21
01−23
04−24
Date
Detect
Non−Detect
Groundwater Elevation
1,1−Dichloroethene
1,4−Dioxane
Perchlorate
Trichloroethene

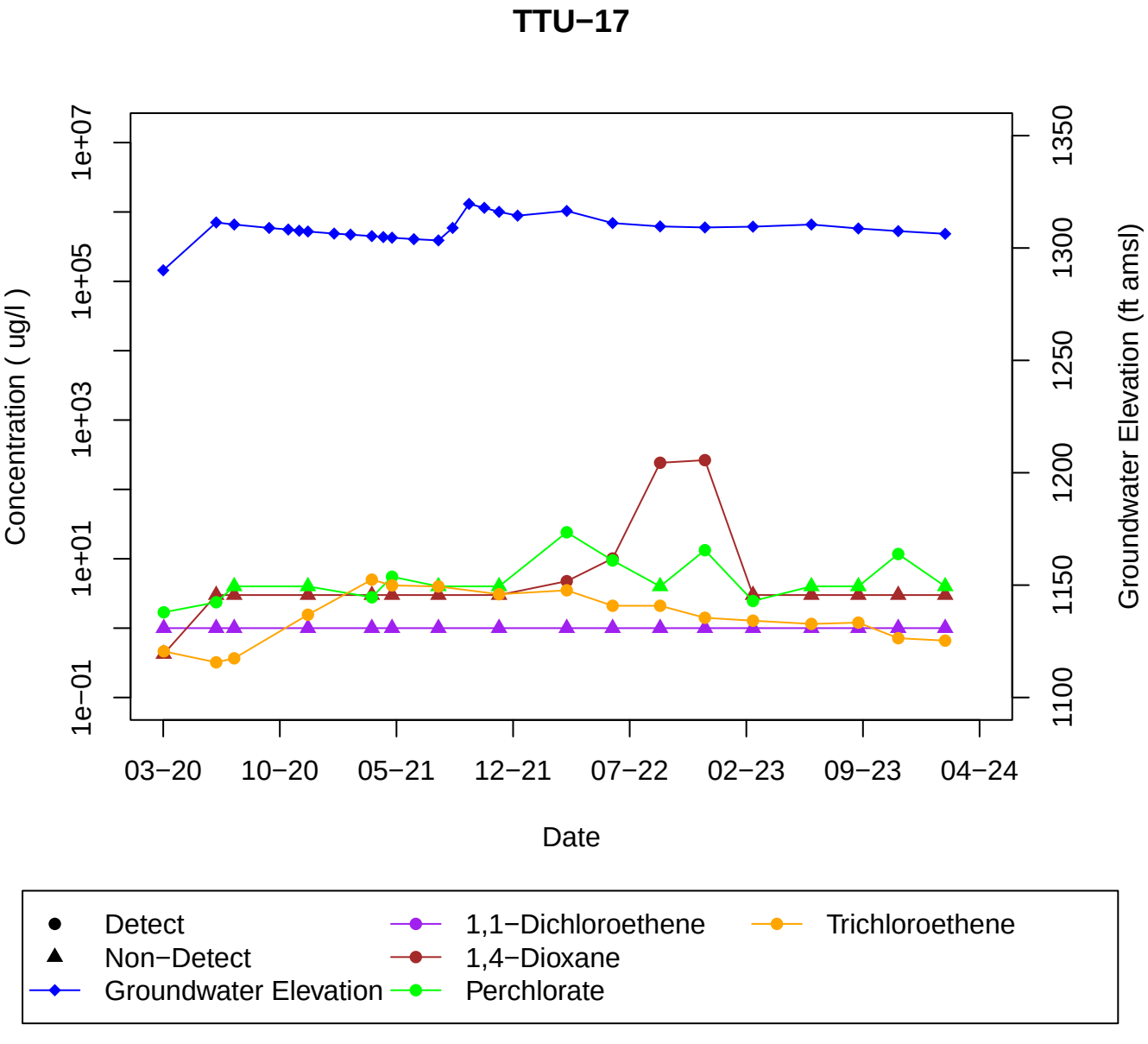

TTU−12
Concentration (ug/l)
1e−01
1e+01
1e+03
1e+05
1e+07
Groundwater Elevation (ft amsl)
1100
1150
1200
1250
1300
1350
08−17
08−18
07−19
07−20
06−21
05−22
05−23
04−24
Date
Detect
Non−Detect
Groundwater Elevation
1,1−Dichloroethene
1,4−Dioxane
Perchlorate
Trichloroethene

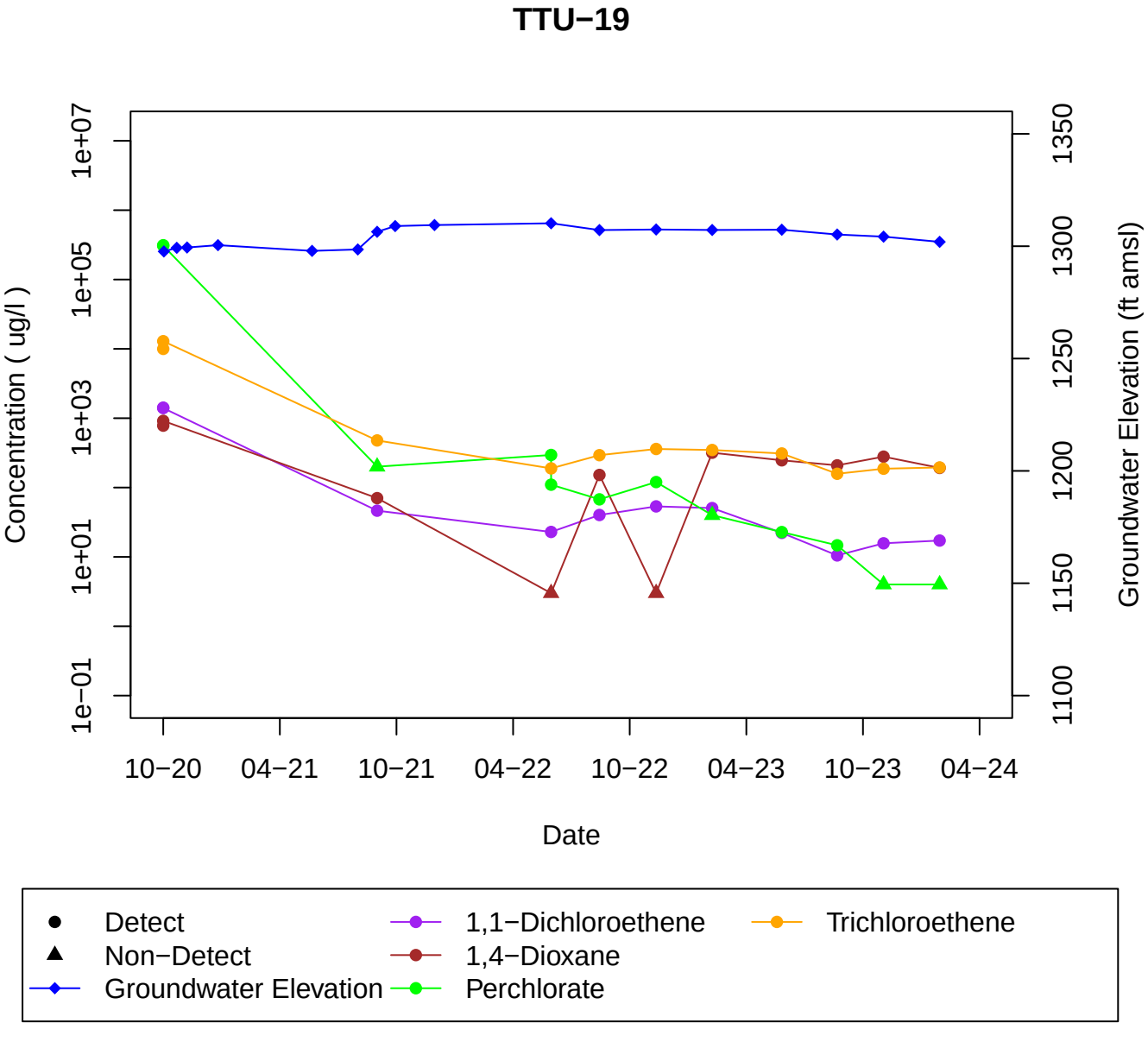

TTU−13
Concentration (ug/l)
1e−01
1e+01
1e+03
1e+05
1e+07
Groundwater Elevation (ft amsl)
1100
1150
1200
1250
1300
1350
08−17
08−18
07−19
07−20
06−21
05−22
05−23
04−24
Date
Detect
Non−Detect
Groundwater Elevation
1,1−Dichloroethene
1,4−Dioxane
Perchlorate
Trichloroethene

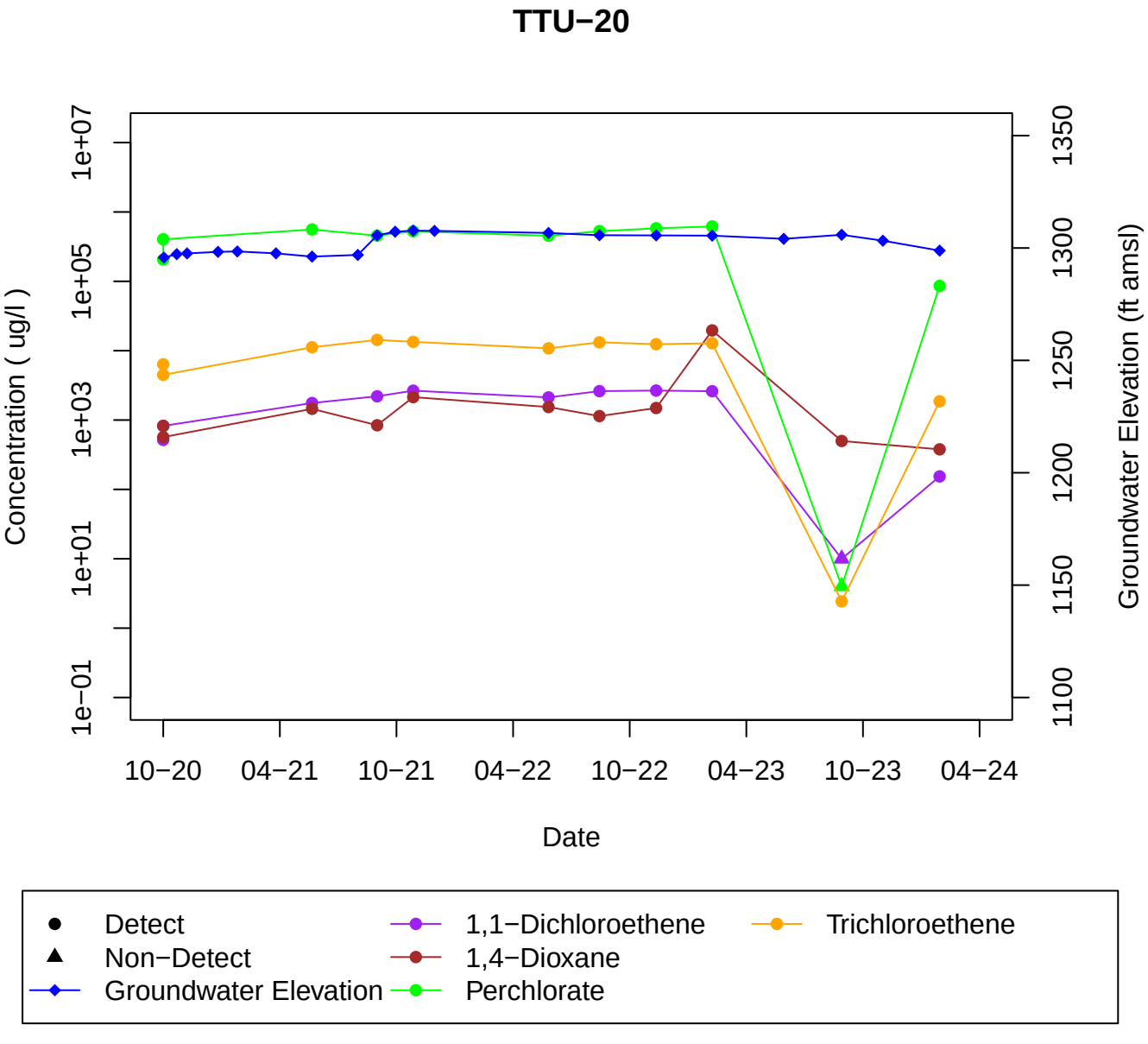

TTU−14
Concentration (ug/l)
1e−01
1e+01
1e+03
1e+05
1e+07
Groundwater Elevation (ft amsl)
1100
1150
1200
1250
1300
1350
08−17
08−18
07−19
07−20
06−21
05−22
05−23
04−24
Date
Detect
Non−Detect
Groundwater Elevation
1,1−Dichloroethene
1,4−Dioxane
Perchlorate
Trichloroethene

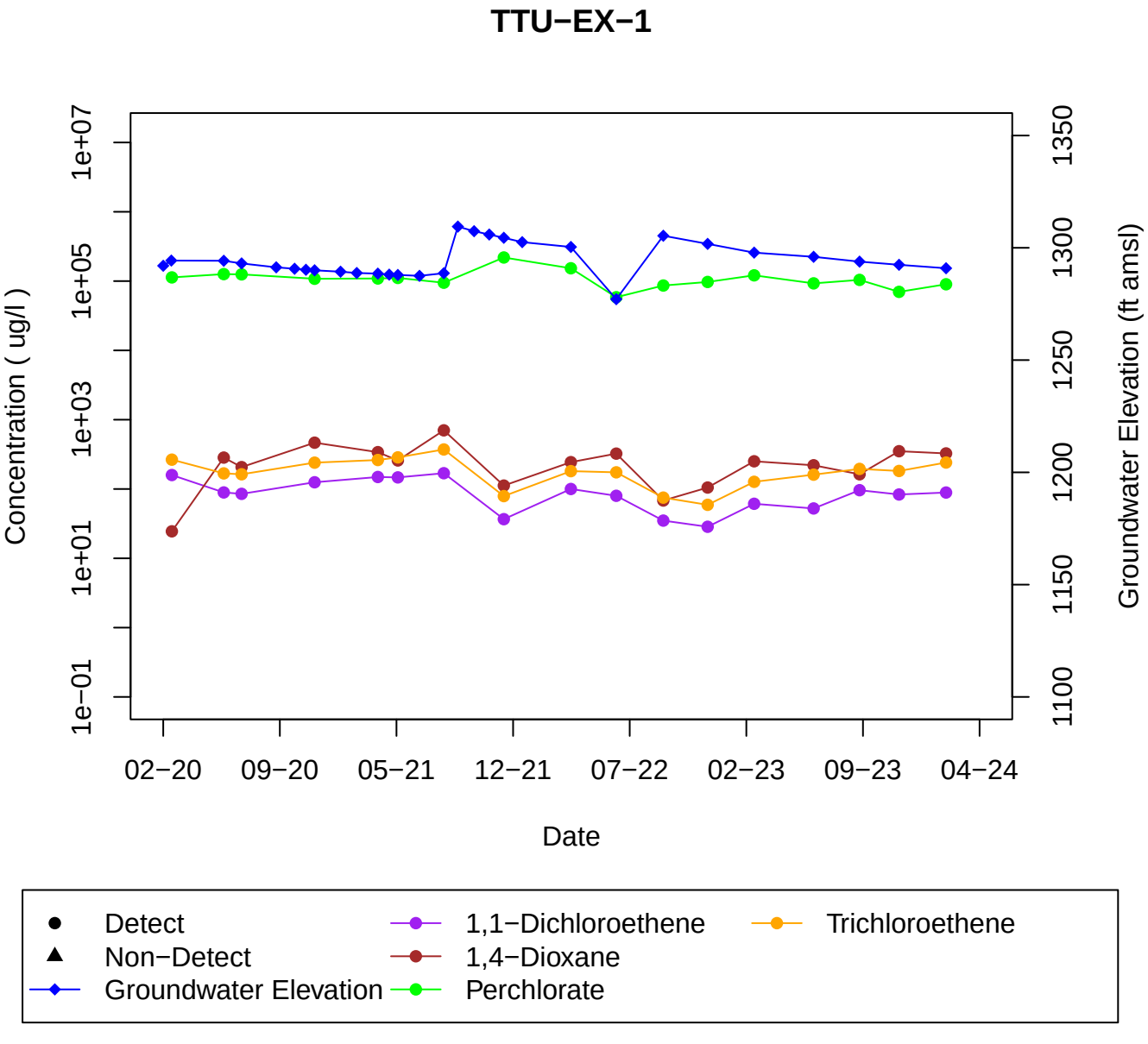

TTU−15
Concentration (ug/l)
1e−01
1e+01
1e+03
1e+05
1e+07
Groundwater Elevation (ft amsl)
1100
1150
1200
1250
1300
1350
03−19
12−19
09−20
05−21
02−22
11−22
07−23
04−24
Date
Detect
Non−Detect
Groundwater Elevation
1,1−Dichloroethene
1,4−Dioxane
Perchlorate
Trichloroethene

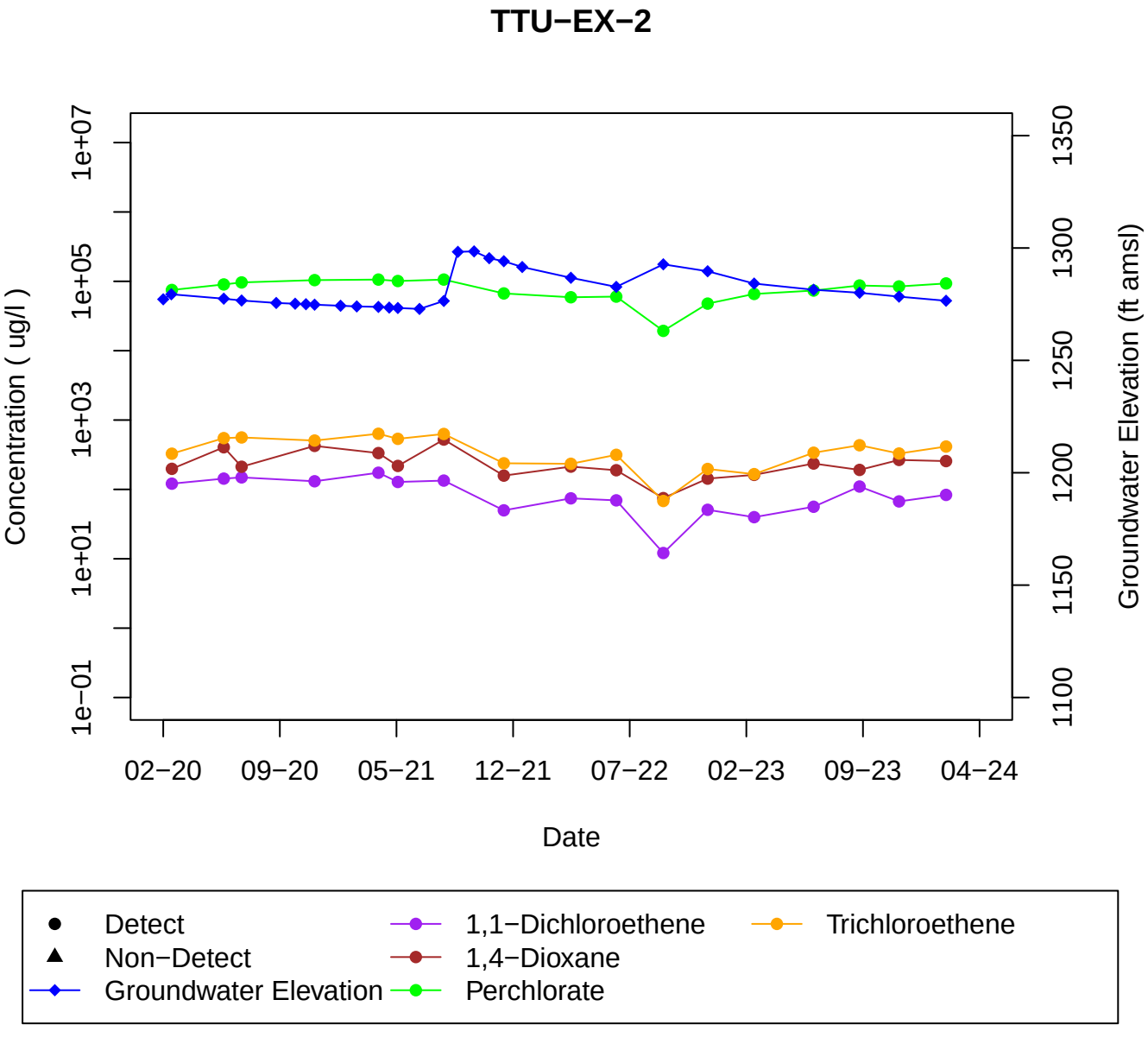

TTU−16
Concentration (ug/l)
1e−01
1e+01
1e+03
1e+05
1e+07
Groundwater Elevation (ft amsl)
1100
1150
1200
1250
1300
1350
03−20
10−20
05−21
12−21
07−22
02−23
09−23
04−24
Date
Detect
Non−Detect
Groundwater Elevation
1,1−Dichloroethene
1,4−Dioxane
Perchlorate
Trichloroethene

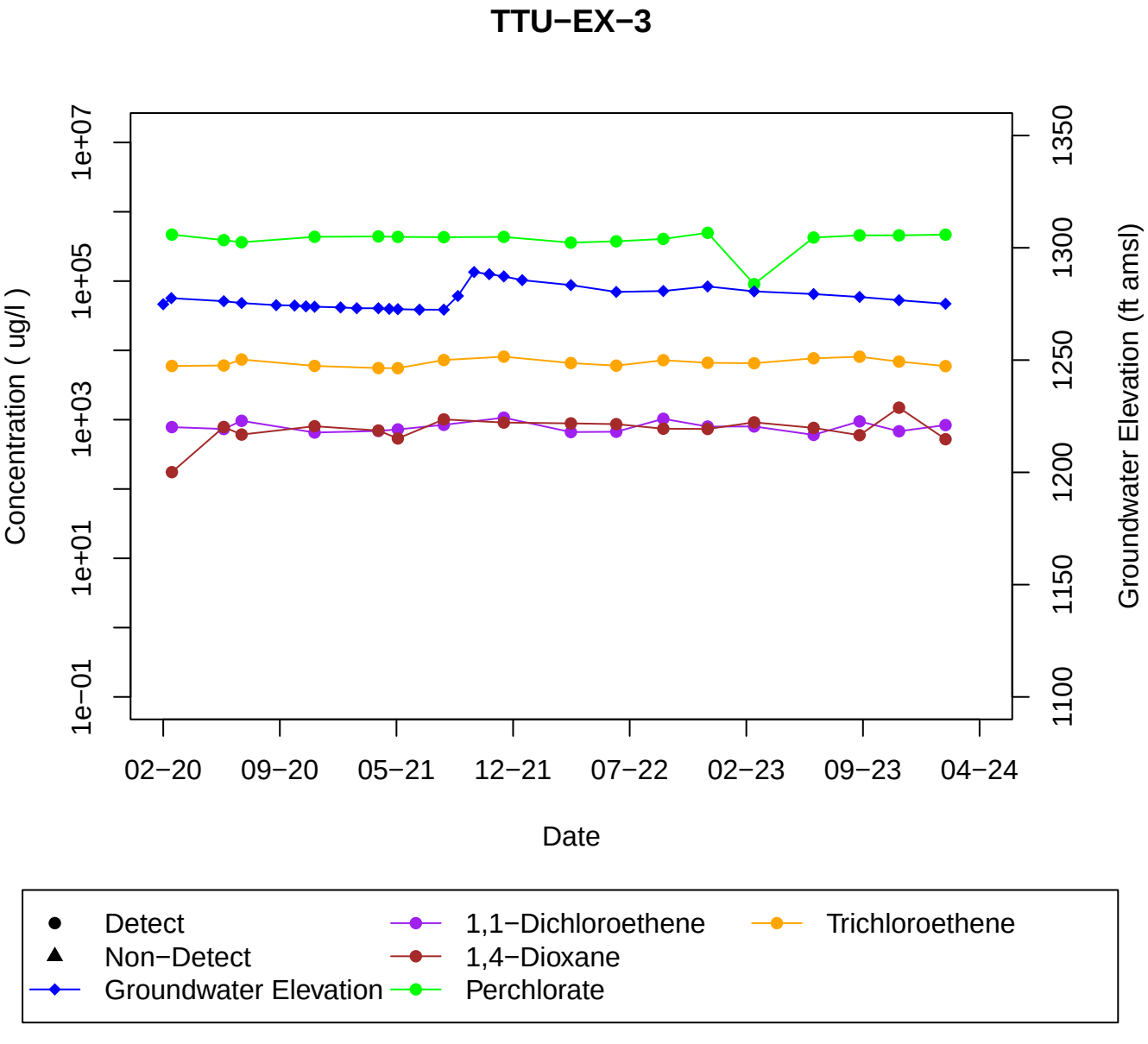

TTU−17
Concentration (ug/l)
1e−01
1e+01
1e+03
1e+05
1e+07
Groundwater Elevation (ft amsl)
1100
1150
1200
1250
1300
1350
03−20
10−20
05−21
12−21
07−22
02−23
09−23
04−24
Date
Detect
Non−Detect
Groundwater Elevation
1,1−Dichloroethene
1,4−Dioxane
Perchlorate
Trichloroethene

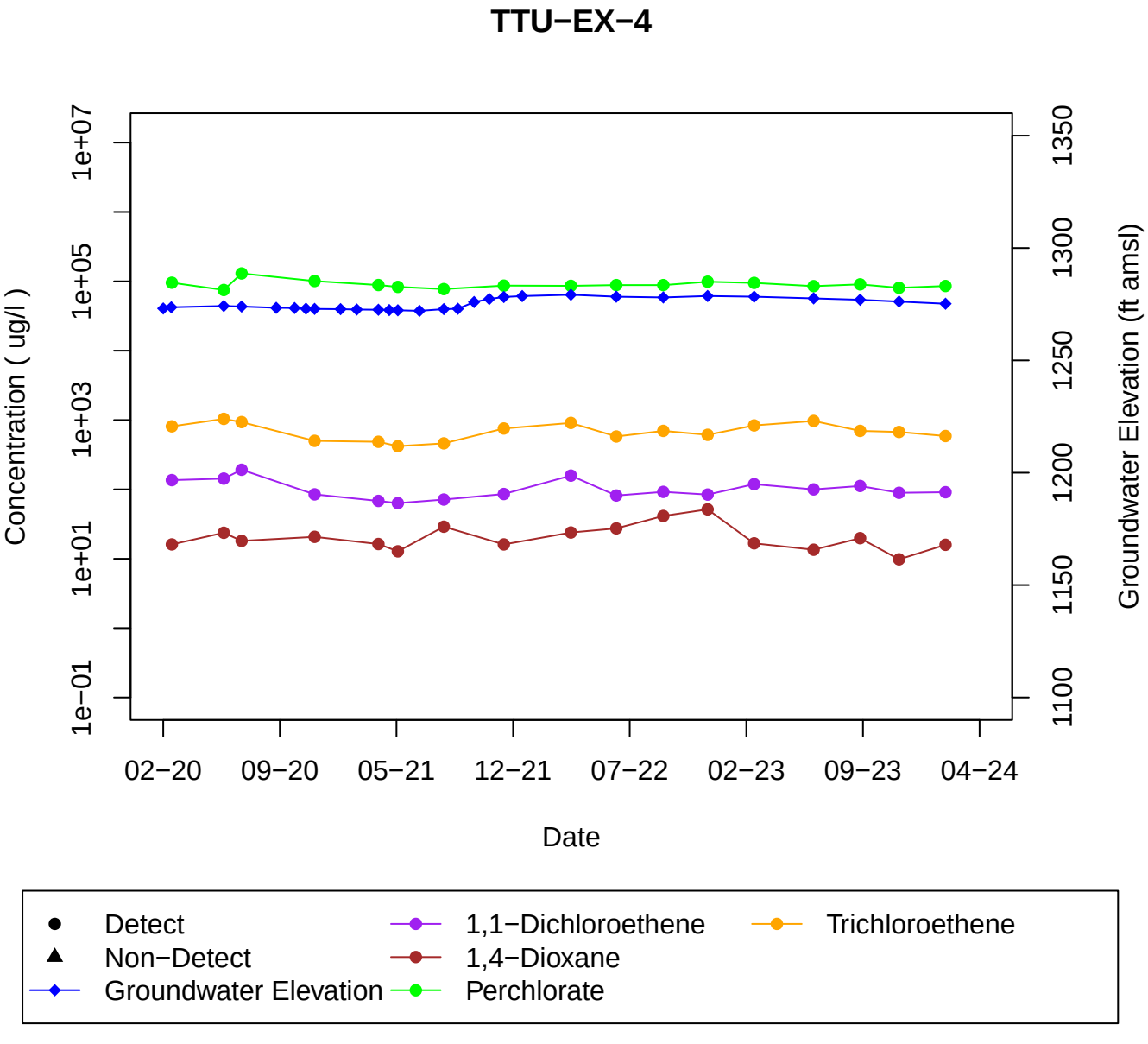

TTU−19
Concentration (ug/l)
1e−01
1e+01
1e+03
1e+05
1e+07
Groundwater Elevation (ft amsl)
1100
1150
1200
1250
1300
1350
10−20
04−21
10−21
04−22
10−22
04−23
10−23
04−24
Date
Detect
Non−Detect
Groundwater Elevation
1,1−Dichloroethene
1,4−Dioxane
Perchlorate
Trichloroethene

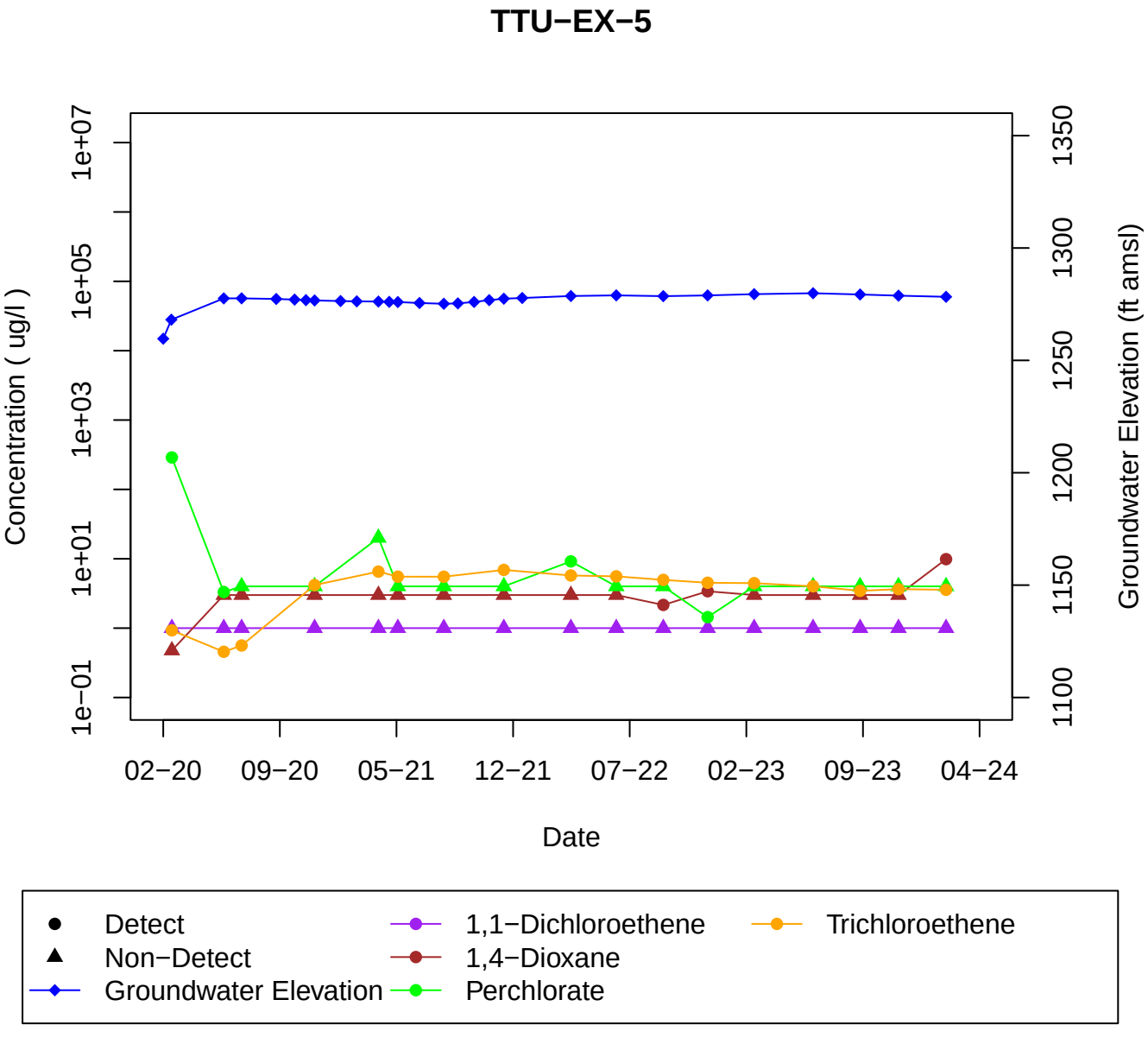

TTU−20
Concentration (ug/l)
1e−01
1e+01
1e+03
1e+05
1e+07
Groundwater Elevation (ft amsl)
1100
1150
1200
1250
1300
1350
10−20
04−21
10−21
04−22
10−22
04−23
10−23
04−24
Date
Detect
Non−Detect
Groundwater Elevation
1,1−Dichloroethene
1,4−Dioxane
Perchlorate
Trichloroethene


TTU−EX−1
Concentration (ug/l)
1e−01
1e+01
1e+03
1e+05
1e+07
Groundwater Elevation (ft amsl)
1100
1150
1200
1250
1300
1350
02−20
09−20
05−21
12−21
07−22
02−23
09−23
04−24
Date
Detect
Non−Detect
Groundwater Elevation
1,1−Dichloroethene
1,4−Dioxane
Perchlorate
Trichloroethene


TTU−EX−2
Concentration (ug/l)
1e−01
1e+01
1e+03
1e+05
1e+07
Groundwater Elevation (ft amsl)
1100
1150
1200
1250
1300
1350
02−20
09−20
05−21
12−21
07−22
02−23
09−23
04−24
Date
Detect
Non−Detect
Groundwater Elevation
1,1−Dichloroethene
1,4−Dioxane
Perchlorate
Trichloroethene


TTU−EX−3
Concentration (ug/l)
1e−01
1e+01
1e+03
1e+05
1e+07
Groundwater Elevation (ft amsl)
1100
1150
1200
1250
1300
1350
02−20
09−20
05−21
12−21
07−22
02−23
09−23
04−24
Date
Detect
Non−Detect
Groundwater Elevation
1,1−Dichloroethene
1,4−Dioxane
Perchlorate
Trichloroethene


TTU−EX−4
Concentration (ug/l)
1e−01
1e+01
1e+03
1e+05
1e+07
Groundwater Elevation (ft amsl)
1100
1150
1200
1250
1300
1350
02−20
09−20
05−21
12−21
07−22
02−23
09−23
04−24
Date
Detect
Non−Detect
Groundwater Elevation
1,1−Dichloroethene
1,4−Dioxane
Perchlorate
Trichloroethene


TTU−EX−5
Concentration (ug/l)
1e−01
1e+01
1e+03
1e+05
1e+07
Groundwater Elevation (ft amsl)
1100
1150
1200
1250
1300
1350
02−20
09−20
05−21
12−21
07−22
02−23
09−23
04−24
Date
Detect
Non−Detect
Groundwater Elevation
1,1−Dichloroethene
1,4−Dioxane
Perchlorate
Trichloroethene

Attachment 4 – Data Validation Documentation

Pinyon Environmental, Inc.

Data Receipt Checklist

Project Name: Nammo TTU Date Collected: 2/14/24

Laboratory Report(s): J214412-1

Samples Submitted:

Number of Sample Type	Analysis: Perchlorate 314.0	Analysis:	Analysis:	Analysis:
Total	1			
Duplicate samples	0			
Trip blanks	0			
Field blanks	0			
Equipment blanks	0			
MS/MSD	0			
Other:	0			

How were samples relinquished to the lab? Drop-off ☑ Courier ☐ Shipped to Lab ☐

Date relinquished to lab: 2/15/24 Date received by lab: 2/15/24

Yes ☑ No ☐ Do names samples match the COC? PF-2-400-20240214

Yes ☑ No ☐ Do dates and times for samples match the COC? No time listed on COC

Yes ☐ No ☑ Did the lab assign date/time to samples (i.e. duplicate without date/time listed on COC)?
NA - No duplicate samples

Yes ☑ No ☐ Were samples preserved appropriately?

Preservatives: Ice

Yes ☑ No ☐ Where samples analyzed within holding time?

Holding Times: 28 days - analyzed 9 days post sampling

Yes ☑ No ☐ Where samples received within temperature?

Temperature Received: 5.1 C

Yes ☐ No ☑ Are tentatively identified compounds (TICs) reported?

Yes ☐ No ☑ Any analytes detect in the trip blank/field blank/equipment blank? NA,
no blanks submitted

Yes ☑ No ☐ Was an EDD received?

Yes ☑ No ☐ Is data validation required?

Data Checker: Daniel Jennings Date: 4/2/24

Pinyon Environmental, Inc.

Data Receipt Checklist

Sample Summary:

Original Sample Number	Original Sample Name	Original Sample Date	Duplicate Sample Number	Duplicate Sample Name	Duplicate Sample Date	Analysis

Data Checker: Daniel Jennings Date: 4/2/24

Pinyon

Data Validation Checklist

Project Name: Nammo TTU Date Collected: 2/14/24

Laboratory Report(s): J214412-1

Analysis Notes:

Yes☑ No☐ Was anything noted in the case narrative? Change to project name

E4, E8, T5 (Lab not licensed for this parameter)

Yes☐ No☐ Were surrogate recoveries within limits? NA

Yes☑ No☐ Were samples reported to MDLs?

Yes☐ No☑ Any elevated reporting limits for non-detects?

Yes☐ No☑ Any analytes detected in the method blank?

Yes☐ No☑ Are there qualifiers reported for laboratory QC samples?

Data Validator: Daniel Jennings Date: 4/2/24

Pinyon Environmental, Inc.

Data Validation Checklist

Laboratory Report(s): J214412-1

Laboratory QC Samples:

Batch Number	Analysis	Method Blank	MS/MSD	Lab Duplicate	LCS/LCSD	Other
Example1234	*8260B*	*1*	*2; 1 (MS only)*	*0*	*1*	*0*
-1	314.0	1	1/1	0	2/1	0

Data Validator: Daniel Jennings Date: 4/2/24 Page: 4/4

*Revised report received on 3/27/24

Pinyon Environmental, Inc.

Data Receipt Checklist

Project Name: Nammo TTU Groundwater Monitoring Date Collected: 2/12/24-2/13/24

Laboratory Report(s): L1705776

Samples Submitted:

Number of Sample Type	Analysis: 314.0 Mod	Analysis: 8260B	Analysis: 8260B-SIM	Analysis:
Total	22	22	22	
Duplicate samples	2	2	2	
Trip blanks	0	1	1	
Field blanks	0	0	0	
Equipment blanks	0	0	0	
MS/MSD	2	2	2	
Other:	0	0	0	

How were samples relinquished to the lab? Drop-off ☑ Courier ☐ Shipped to Lab ☐

Date relinquished to lab: 2/13/24 Date received by lab: 2/14/24

Yes☐ No☑ Do names samples match the COC? -03, -06, -07, -10, -11, -12, -15, -21, -22

Yes☑ No☐ Do dates and times for samples match the COC?

Yes☑ No☐ Did the lab assign date/time to samples (i.e. duplicate without date/time listed on COC)?

Dup-01 02/13/24 00:00; Dup-02 02/13/24 00:00; Trip Blank 02/13/24 00:00

Yes☑ No☐ Were samples preserved appropriately?

Preservatives:

Yes☑ No☐ Where samples analyzed within holding time?

Holding Times:

Yes☑ No☐ Where samples received within temperature? Based on Case Narrative

Temperature Received: Not listed on COC

Yes☐ No☑ Are tentatively identified compounds (TICs) reported?

Yes☐ No☑ Any analytes detect in the trip blank/field blank/equipment blank?

Yes☑ No☐ Was an EDD received?

Yes☑ No☐ Is data validation required?

Data Checker: Daniel Jennings Date: 3/16/24

Pinyon Environmental, Inc.

Data Receipt Checklist

Sample Summary:

Original Sample Number	Original Sample Name	Original Sample Date	Duplicate Sample Number	Duplicate Sample Name	Duplicate Sample Date	Analysis
-23	TTU-4-57	2/13/24	-23	Dup-01	2/13/24	314.0 Mod; 8260B; 8260B-SIM
-24	TTU-EX-5-80	2/13/24	-24	Dup-02	2/13/24	314.0 Mod; 8260B; 8260B-SIM

Data Checker: Daniel Jennings Date: 3/16/24

Data Validation Checklist

Project Name: Nammo TTU Groundwater Monitoring Date Collected: 2/12/24-2/13/24

Laboratory Report(s): L1705776

Analysis Notes:

Yes☑ No☐ Was anything noted in the case narrative? See report, insufficient sample volume to perform MS/MSD analyses per method GC requirements 8260B (12 samples), 8260B-SIM (4 samples)

Yes☐ No☑ Were surrogate recoveries within limits? -09, -22 qualified S10 Surrogate recovery was above laboratory and method acceptance limits.

Yes☑ No☐ Were samples reported to MDLs?

Yes☑ No☐ Any elevated reporting limits for non-detects? -02, -09, -10, -12, -13 (specific compounds), -14, -16, -17, -19, -20, -21, -22

Yes☐ No☑ Any analytes detected in the method blank?

Yes☑ No☐ Are there qualifiers reported for laboratory QC samples? E4, L1, L2, M1, M2, M3, R5, R8 S10

Data Validator: Daniel Jennings Date: 3/16/24

Pinyon Environmental, Inc.

Data Validation Checklist

Laboratory Report(s): L1705776

Laboratory QC Samples:

Batch Number	Analysis	Method Blank	MS/MSD	Lab Duplicate	LCS/LCSD	Other
Example1234	*8260B*	*1*	*2; 1 (MS only)*	*0*	*1*	*0*
WG2228552	314.0 Mod	2	2; 2	2	2	0
WG2228553	314.0 Mod	4	3; 2	2	1	0
WG2228368	8260B	1	2	0	1	0
WG2228641	8260B	1	1	0	1	0
WG2229520	8260B	1	0	0	1	0
WG2227341	8260B-SIM	1	1	0	1	0
WG2230157	8260B-SIM	1	2	0	1	0
WG2230690	8260B-SIM	1	0	0	1	0
WG2231600	8260B-SIM	1	0	0	1	0

Data Validator: Daniel Jennings Date: 3/16/24

*Report and COC description lists the project as WBO; however records are for TTU

Pinyon

Data Receipt Checklist

Project Name: Nammo TTU Groundwater Monitoring Date Collected: 2/14/24

Laboratory Report(s): L1706682

Samples Submitted:

Number of Sample Type	Analysis: 314.0 MOD	Analysis: 8260B	Analysis: 8260B-SIM	Analysis:
Total	2	3	3	
Duplicate samples	1	1	1	
Trip blanks	0	0	0	No analysis requested
Field blanks	0	0	0	
Equipment blanks	0	0	0	
MS/MSD	0	0	0	
Other:	0	0	0	

How were samples relinquished to the lab? Drop-off ☑ Courier ☐ Shipped to Lab ☐

Date relinquished to lab: 2/15/24 Date received by lab: 2/15/24

Yes☑ No☐ Do names samples match the COC?

Yes☑ No☐ Do dates and times for samples match the COC?

Yes☑ No☐ Did the lab assign date/time to samples (i.e. duplicate without date/time listed on COC)?
L1706682-04 assigned 02/14/24 00:00

Yes☑ No☐ Were samples preserved appropriately?

Preservatives: HCl

Yes☑ No☐ Where samples analyzed within holding time?

Holding Times: 314.0: 16 - 18days, 8260B: 7 - 8 days, 8260B-SIM: 8-11 days

Yes☑ No☐ Where samples received within temperature?

Temperature Received: 4.4 degrees Celsius

Yes☐ No☑ Are tentatively identified compounds (TICs) reported?

Yes☐ No☑ Any analytes detect in the trip blank/field blank/equipment blank?
Trip blank not analyzed

Yes☑ No☐ Was an EDD received?

Yes☑ No☐ Is data validation required?

Data Checker: Mary G. Weiss Date: 3/27/24

Pinyon Environmental, Inc.

Data Receipt Checklist

Sample Summary:

Original Sample Number	Original Sample Name	Original Sample Date	Duplicate Sample Number	Duplicate Sample Name	Duplicate Sample Date	Analysis
-03	TTU-10-172-20240214	2-14-24	-04	Dup-03	2-14-24	314.0 Mod, 8260B, 8260B-SIM

Data Checker: Mary G. Weiss Date: 3/27/24

Data Validation Checklist

Project Name: Nammo TTU Groundwater Monitoring Date Collected: 2/14/24

Laboratory Report(s): L1706682

Analysis Notes:

Yes☑ No☐ Was anything noted in the case narrative? -04 for 8260B: Insufficient sample volume to perform MS/MSD analyses per method QC requirements.

Yes☑ No☐ Were surrogate recoveries within limits?

Yes☑ No☐ Were samples reported to MDLs?

Yes☐ No☑ Any elevated reporting limits for non-detects?

Yes☑ No☐ Any analytes detected in the method blank? Estimated concentration of carbon disulfide detected in R4036492-4

Yes☑ No☐ Are there qualifiers reported for laboratory QC samples? B1, E4, L1, M1, M2, M3, R5, R7, R8

Data Validator: Mary G. Weiss Date: 3/27/24

Pinyon Environmental, Inc.

Data Validation Checklist

Laboratory Report(s): L1706682

Laboratory QC Samples:

Batch Number	Analysis	Method Blank	MS/MSD	Lab Duplicate	LCS/LCSD	Other
Example1234	*8260B*	*1*	*2; 1 (MS only)*	*0*	*1*	*0*
WG2228553	314.0	4	5 ; 2	2	1	0
WG2234696	314.0	2	3 ; 3	2	1	0
WG2230378	8260B	1	1 ; 1	0	1 ; 1	0
WG2230965	8260B	1	0 ; 0	0	1 ; 1	0
WG2231221	8260B-SIM	1	1 ; 1	0	1 ; 1	0
WG2231965	8260B-SIM	1	1 ; 1	0	1 ; 1	0

Data Validator: Mary G. Weiss Date: 3/27/24

Data Receipt Checklist

Project Name: Nammo TTU Groundwater Monitoring Date Collected: 3/18/24

Laboratory Report(s): L1716428

Samples Submitted:

Number of Sample Type	Analysis: 8260B-SIM	Analysis:	Analysis:	Analysis:
Total	1			
Duplicate samples	0			
Trip blanks	0			
Field blanks	0			
Equipment blanks	0			
MS/MSD	0			
Other:	0			

How were samples relinquished to the lab? Drop-off ☑ Courier ☐ Shipped to Lab ☐

Date relinquished to lab: 3/18/24 Date received by lab: 3/18/24

Yes ☑ No ☐ Do names samples match the COC?

Yes ☑ No ☐ Do dates and times for samples match the COC?

Yes ☐ No ☑ Did the lab assign date/time to samples (i.e. duplicate without date/time listed on COC)?

Yes ☑ No ☐ Were samples preserved appropriately?

Preservatives: HCl

Yes ☑ No ☐ Where samples analyzed within holding time?

Holding Times: 8260B-SIM: 2 days

Yes ☑ No ☐ Where samples received within temperature?

Temperature Received: 1.2 degrees celsius

Yes ☐ No ☑ Are tentatively identified compounds (TICs) reported?

Yes ☐ No ☑ Any analytes detect in the trip blank/field blank/equipment blank? NA

Yes ☑ No ☐ Was an EDD received?

Yes ☑ No ☐ Is data validation required?

Data Checker: Mary G. Weiss Date: 3/27/24

Pinyon Environmental, Inc.

Data Receipt Checklist

Sample Summary:

Original Sample Number	Original Sample Name	Original Sample Date	Duplicate Sample Number	Duplicate Sample Name	Duplicate Sample Date	Analysis
None						

Data Checker: Mary G. Weiss Date: 3/27/24

Data Validation Checklist

Project Name: Nammo TTU Groundwater Monitoring Date Collected: 3/18/24

Laboratory Report(s): L1716428

Analysis Notes:

Yes☑ No☐ Was anything noted in the case narrative? -01 for 8260B-SIM: Insufficient sample volume to perform MS/MSD analyses per method QC requirements.

Yes☑ No☐ Were surrogate recoveries within limits?

Yes☑ No☐ Were samples reported to MDLs?

Yes☐ No☑ Any elevated reporting limits for non-detects?

Yes☐ No☑ Any analytes detected in the method blank?

Yes☐ No☑ Are there qualifiers reported for laboratory QC samples?

Data Validator: Mary G. Weiss Date: 3/27/24

Pinyon Environmental, Inc.

Data Validation Checklist

Laboratory Report(s): L1716428

Laboratory QC Samples:

Batch Number	Analysis	Method Blank	MS/MSD	Lab Duplicate	LCS/LCSD	Other
Example1234	*8260B*	*1*	*2; 1 (MS only)*	*0*	*1*	*0*
WG2249950	8260B-SIM	1	0	0	1;1	0

Data Validator: Mary G. Weiss Date: 3/27/24

Data Receipt Checklist

Project Name: Nammo TTU Date Collected: 4/9/24

Laboratory Report(s): L1724110

Samples Submitted:

Number of Sample Type	Analysis: 314.0 Mod	Analysis: 8260B	Analysis:	Analysis:
Total	2	2		
Duplicate samples	0	0		
Trip blanks	1	1		
Field blanks	0	0		
Equipment blanks	0	0		
MS/MSD	0	0		
Other:	0	0		

How were samples relinquished to the lab? Drop-off ☑ Courier ☐ Shipped to Lab ☐

Date relinquished to lab: 4/9/24 Date received by lab: 4/9/24

Yes☑ No☐ Do names samples match the COC?

Yes☑ No☐ Do dates and times for samples match the COC?

Yes☑ No☐ Did the lab assign date/time to samples (i.e. duplicate without date/time listed on COC)?

L1724110-03 assigned 4/9/24 00:00 as date/time.

Yes☑ No☐ Were samples preserved appropriately?

Preservatives: HCl

Yes☑ No☐ Where samples analyzed within holding time?

Holding Times: 314.0 Mod: 4 days ; 8260B: 2 days

Yes☑ No☐ Where samples received within temperature?

Temperature Received: 3.9 degrees celsius

Yes☐ No☑ Are tentatively identified compounds (TICs) reported?

Yes☑ No☐ Any analytes detect in the trip blank/field blank/equipment blank?

Carbon disulfide detected at estimated concentration

Yes☑ No☐ Was an EDD received?

Yes☑ No☐ Is data validation required?

Data Checker: Mary G. Weiss Date: 4/24/24

Pinyon

Data Validation Checklist

Project Name: Nammo TTU Date Collected: 4/9/24

Laboratory Report(s): L1724110

Analysis Notes:

Yes☑ No☐ Was anything noted in the case narrative? For 8260B samples -01 to -03: Insufficient sample volume to perform MS/MSD analyses per method QC requirements.

Yes☑ No☐ Were surrogate recoveries within limits?

Yes☑ No☐ Were samples reported to MDLs?

Yes☑ No☐ Any elevated reporting limits for non-detects? samples -01 and -02 for 314.0 Mod and 8260B: Elevated detection limits.

Yes☑ No☐ Any analytes detected in the method blank? Carbon disulfide detected at estimated concentration in MB R4056937-4

Yes☑ No☐ Are there qualifiers reported for laboratory QC samples? B1 and E4

Data Validator: Mary G. Weiss Date: 4/24/24

Pinyon Environmental, Inc.

Data Validation Checklist

Laboratory Report(s): L1724110

Laboratory QC Samples:

Batch Number	Analysis	Method Blank	MS/MSD	Lab Duplicate	LCS/LCSD	Other
Example1234	*8260B*	*1*	*2; 1 (MS only)*	*0*	*1*	*0*
WG2264996	314.0 Mod	1	1;1	1	1	0
WG2264870	8260B	1	0;0	0	1;1	0

Data Validator: Mary G. Weiss Date: 4/24/24