



# REVISED SEPTEMBER 2021 BIENNIAL GROUNDWATER MONITORING REPORT

Former Signore Inc. 55-57 Jefferson Street Ellicottville, New York 14731

January 26, 2022 File No. 21.0056491.81



### **PREPARED FOR:**

Iskalo Ellicottville Holdings LLC Williamsville, New York

### **GZA GeoEnvironmental of New York**

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### **VIA EMAIL**

January 26, 2022 File No. 21.0056491.81

Mr. David Chiazza Iskalo Ellicottville Holdings LLC Harbinger Square 5166 Main Street Williamsville, New York 14221 email: dchiazza@iskalo.com

Re: Revised September 2021 Biennial Groundwater Monitoring Report Former Signore, Inc. Facility 55-57 Jefferson Street Ellicottville, New York 14731

Mr. Chiazza:

GZA GeoEnvironmental of New York (GZA) is pleased to submit this biennial groundwater monitoring report to Iskalo Ellicottville Holdings LLC (Iskalo). This report summarizes the analytical results of the sampling event conducted in September 2021 at the above referenced Site. The biennial groundwater monitoring was performed as required by the New York State Department of Environmental Conservation (NYSDEC) as specified in the Record of Decision (ROD) dated January 1992. Based upon the work conducted and the rate of chlorinated VOC (volatile organic compound) reduction observed, NYSDEC approved modification of the ROD to allow for a reduced frequency of groundwater monitoring from annual to biennial in 2020.

This report provides the analytical results of the ROD-required monitoring (12 wells sampled). The analytical results of the groundwater sampling provide data for concentrations of VOCs present in the on-Site groundwater and inform the areal extent of these constituents. Both on-site and off-site monitoring wells have been sampled since 1994. Comparison of over 20 years of groundwater data confirms that concentrations of tetrachloroethene (PCE) and trichloroethene (TCE) and their breakdown products cis-1,2-dichloroethene (cis-DCE) and vinyl chloride (VC) continue to decline with less exceedances of NYSDEC Class GA groundwater standards observed.



Should you have any questions or require additional information following your review, please contact Thomas Bohlen at 716-844-7050.

Sincerely,

GZA GEOENVIRONMENTAL OF NEW YORK

homas Bohlen

Sat a. Konte

Thomas Bohlen, P.G. Project Manager

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cc: Megan Kuczka, NYSDEC

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for Karen Kinsella, Ph.D.

Senior Technical Specialist (retired)



APPENDIX D ANALYTICAL TEST RESULTS

January 26, 2022 Revised Former Signore Inc., Ellicottville, New York September Biennial Monitoring Well Sampling Report File No. 21.0056491.81

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### 1.0 INTRODUCTION

In accordance with our proposal dated July 12, 2021, GZA GeoEnvironmental of New York (GZA) collected groundwater samples at eight on-Site and four off-site monitoring wells associated with the Former Signore, Inc. facility located at 55-57 Jefferson Street, Ellicottville, New York (Site). The sampling was performed on September 15 through September 17, 2021. A Locus Plan and Site Plan are attached as Figure 1 and Figure 2, respectively.

### 1.1 BACKGROUND AND SITE HISTORY

The Former Signore Inc. Site is currently listed as a Class 4 Site on the New York State Department of Environmental Conservation (NYSDEC), Inactive Hazardous Waste Site (IHWS) registry (Site #905023). As part of the Record of Decision (ROD) dated January 1992 issued by NYSDEC, 12 monitoring wells were to be sampled on a semi-annual basis. The former owner, Signore Inc., ceased sampling of these wells in October 2006.

In December 2007, GZA completed a Phase II ESA at the Site as part of due diligence services for Iskalo. During the Phase II activities, VOC contamination was identified in on-Site soil and groundwater. Three areas of concern (AOC) were identified where VOC concentrations in soil were greater than the NYSDEC Unrestricted Soil Cleanup Objectives (6 NYCRR Part 375¹ criteria).

Iskalo Ellicottville Holdings LLC (Iskalo) took ownership of the property in February 2008 and conducted the required sampling activities since April 2009. The Site was accepted into the NYSDEC Brownfield Cleanup Program (BCP) as Site # C905034 in January 2011.

In late 2011, AOC-1 and the majority of AOC-2 were addressed under a NYSDEC-approved IRM work plan. VOC-impacted soils and underground storage tanks were removed for proper off-Site disposal. A portion of AOC-2 and AOC-3 were located beneath the former building during the time of the first IRM, which has since been demolished.

A Supplemental Remedial Investigation (SRI) was completed between January 2012 and January 2013. The activities included the following.

- Off-site soil vapor intrusion assessment of nine homes;
- Completion of 10 on-site test pits;
- Completion of 21 on-site soil probes;
- Collection and analysis of four on-site surface soil samples;
- Collection and analysis of 21 soil samples from the 21 soil probes; and
- Collection of 19 groundwater samples from the 14 new microwells installed as part of the SRI, and 5
  existing wells.

<sup>&</sup>lt;sup>1</sup>6 New York Code Rules and Regulation (6 NYCRR) Part 375 Environmental Remediation Programs, effective December 14, 2006 (Part 375).





The magnitude and areal extent of groundwater contamination was further defined within the Signore BCP Site boundaries, during the SRI activities.

The remaining portion of AOC-2 and AOC-3 were addressed in summer 2013 as part of a 2<sup>nd</sup> IRM. Impacted soils at these locations were removed for off-site landfill disposal. A groundwater pilot test was also implemented as part of the 2<sup>nd</sup> IRM. The pilot test consisted of the injection of an electron donor compound (EDC) material that was mixed with water. A total of 2,500 pounds of EDC material was injected, 2,000 pounds in the vicinity of SP-3 and 500 pounds at SP-32.

The EDC material enhances the anaerobic breakdown of the "parent" chlorinated compounds present at the Site (TCE and PCE) via reductive dechlorination to the "daughter" breakdown products (cis-dichloroethene (cis-DCE) and vinyl chloride (VC)), which degrade under both anaerobic and aerobic conditions.

The groundwater pilot test work plan included two post-injection sampling events:

- 1) 1st event: not more than 3 months (Fall 2013) after the pilot test injections; and
- 2) 2<sup>nd</sup> event: 9 to 12 months (late Spring 2014) after the pilot test injections.

Groundwater samples were collected from six locations (EW-1.25, SP-32, SP-37, SP-38, SP-43, and SP-45) in conjunction with the October 2013 and June 2014 semi-annual groundwater sampling events. The results from the first to second pilot sampling events were as follows.

- There was a decrease in parent compound concentrations at three of the six sampling locations: EW-1.25, SP-37, and SP-43. Concentrations at SP-32 were relatively consistent and low (generally at estimated concentrations).
- There was an increase in parent compound concentrations at the other two locations: SP-38 and SP-45.
   The slight increase in concentration was not a concern at that time and may have been due to sample variability for this particular sampling event.
- Results of the groundwater pilot test supported the use of a similar but larger scale in-situ injection program for Site groundwater remediation. A Remedial Action Work Plan, detailing the groundwater remedial program, was prepared and submitted to NYSDEC. NYSDEC approved the RAWP and the groundwater remedial injections were initiated on July 6, 2015.

The July 2015 remedial injection material consisted of Organic Carbon Electron Donor Substrate (OCEDS). The program consisted of the injection of approximately 7,000 pounds of OC material into groundwater over an approximately 12,000 ft² oval-shaped area of the Site. The OC material was composed of food-, feed-, and agricultural- grade additives consisting of an aqueous solution of approximately 53% lactose, 40% inactive brewer's yeast or yeast extract, 4% sodium bicarbonate, and 3% trace nutrients (inorganic nitrate, phosphate, potassium, and vitamin B12) by mass. The following materials were mixed into an injectable slurry and injected into the subsurface groundwater, per each five injection locations (10 injections):





Lactose: 264 pounds
 MicroBlend® Yeast Extract: 20 pounds
 Sensient® Yeast Extract: 143 pounds
 Sodium bicarbonate: 21 pounds
 Miracle-Gro®: 14 pounds

The additive slurry was injected in a grid pattern encompassing 70 injection locations. The 70 injection locations were spaced approximately 20 feet apart. One hundred pounds of OC material and 70 gallons of water were injected at each location. The slurry was injected in two intervals below the groundwater table at each location, for a total of 50 pounds of OC material and 35 gallons of water per interval. The deep injection was completed first at each location, at approximately 10 feet below the groundwater table. Groundwater levels were measured on-site in the morning prior to the start of injections and were utilized to determine injection depths across the Site as groundwater levels varied across the area of injections. The deep injection occurred at ~18-20 feet below ground surface (ft. bgs). The direct push soil probe rod was then brought up approximately five feet, and the shallow injection was completed.

Post-injection groundwater sampling events were conducted in August and October 2015, June and October 2016, July 2017, June 2018, and in June 2019 to assess the efficacy of the OCEDS injections in promoting continued natural attenuation of chlorinated VOCs (cVOCs) at the Site. The efficacy of the remedy is being managed and reported under the NYSDEC BCP.

Reductive dechlorination is the biologically- or chemically- mediated replacement of chlorine (as chloride) on a chlorinated organic compound with elemental hydrogen, in the presence of a suitable electron donor. This causes transformation of the cVOC to a less chlorinated product. An electron donor is a substance capable of supplying electrons during oxidation-reduction reactions. In biological reductive dechlorination, microorganisms obtain energy by transferring electrons from electron donors to electron acceptors. Electron donors are chemicallyreduced materials such as the OCEDS. Electron acceptors include oxygen, nitrate, ferric iron, sulfate, and cVOCs. Biological reductive dechlorination of cVOCs typically occurs sequentially from PCE to TCE, TCE to DCE, DCE to VC, VC to ethene, ethene to ethane, and ethane to carbon dioxide and water. Suitability for continued reductive dechlorination can be assessed by measuring groundwater biogeochemical parameters, including dissolved oxygen (DO), oxidation-reduction potential (ORP), reduced iron and manganese, methane, total organic carbon (TOC), nitrate, and sulfate, as well as PCE and TCE degradation products DCE, VC, ethene, ethane, and chloride. In the first few months following injection of an organic carbon additive, groundwater concentrations of PCE and TCE can increase, as their solubility is improved by additive fermentation products. The increased solubility makes the PCE and TCE more available to cVOC-degrading microorganisms, and is typically followed by decreasing PCE and TCE accompanied by an increase in degradation products DCE, VC, ethene, and ethane as bioremediation proceeds.

The analytical results of the groundwater sampling provide data for documentation of concentrations of cVOCs present in the on-Site groundwater. Groundwater cVOC concentrations measured at 74 months post-OCEDS injection (September 2021) follow trends typical for this stage of enhanced reductive dechlorination. As cVOC concentrations decline, biodegradation typically slows down due to less contact between cVOCs and dechlorination bacteria. Also, as PCE and TCE concentrations approach class GA criteria (i.e., PCE and TCE concentrations become a few micrograms per liter ( $\mu$ g/L)), concentrations of their degradation products DCE and





VC are likely to be below laboratory detection limits. At this time, over six years after the Organic Carbon Electron Donor Substrate (OCEDS) injections, with the exception of EW-1.25R located downgradient of the OCEDS injection area, groundwater biogeochemical conditions at the Site are generally less conducive to reductive dechlorination. Daughter product DCE; however, is above laboratory detection limits in monitoring wells that still contain PCE and/or TCE above the class GA groundwater criteria of 5  $\mu$ g/L. With the exception of EW-1.25R, located downgradient of the OCEDC injection area, TOC concentrations in groundwater are low. This is expected, as the OCEDS additive, by design, provides organic carbon for indigenous bacteria to consume while reducing electron acceptors that compete with cVOCs. Biomass generated by bacterial growth cycles provides a sustainable source of organic carbon, helping to maintain reductive dechlorination at the soil-porewater interface as the injected OCEDS is consumed. In GZA's opinion, groundwater concentrations of cVOCs will continue to decline over time. Monitoring will continue to document the dechlorination process.

### 2.0 PURPOSE AND SCOPE OF WORK

Groundwater samples were collected from the 12 monitoring wells to assess current conditions and provide an opinion regarding volatile organic compound (VOC) concentrations. The following was completed:

- Coordinated with Alpha Analytical located in Westborough, Massachusetts prior to commencement of field activities to obtain the analytical sample containers.
- Collected groundwater samples from each of the 12 monitoring wells for chemical analysis of VOCs via EPA Method 8260 Target Compound List (TCL).
- Prepared this report, which summarizes the data collected during this sampling event and compares the data to NYSDEC Class GA groundwater standards and historical data.

This report presents GZA's field observations, results, and opinions and is subject to the limitations presented in Appendix A and modifications if subsequent information is developed by GZA or another party.

### 3.0 FIELD METHODS

This section describes the field activities of GZA's groundwater sampling event.

### 3.1 GROUND WATER SAMPLING PROCEDURES

### **Equipment Cleaning**

Prior to GZA's arrival on-Site, the sampling equipment (water level indicator, water quality meter and flow-through cell) were cleaned by rinsing with potable water, washing with a solution of laboratory detergent (Alconox®) and potable water, and rinsing with de-ionized water.

New, disposable polyethylene tubing (for placement down into the well and connecting to the water quality meter) and silicone tubing (for the peristaltic pump head) was used for groundwater sampling at each location.



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A variable speed peristaltic pump was used to purge groundwater from most monitoring wells. Groundwater remained within the polyethylene and silicone tubing and did not come in contact with the pump. Therefore, the tubing and pump did not require decontamination between sample locations.

### **Equipment Calibration**

A water quality meter and organic vapor meter (OVM) were used during groundwater monitoring. Prior to use each day, the calibration of the water quality meter and OVM were checked to verify that the equipment was in working order.

### Monitoring & Purging Methodologies

An OVM, equipped with photoionization detector (PID) and a 10.6 eV ultraviolet lamp, was used to screen for volatile organics in air at the top of the well riser immediately following the removal of each monitoring well riser cap. OVM readings were recorded on each respective monitoring well field sampling log. OVM readings were non-detect and no odors were noted at the top of each of the 12 monitoring wells sampled.

The purging and water quality measurements were completed using two different types of pumps depending on the depth to water surface measured at the well location.

Nine of the 12 monitoring wells had water surface depths less than 20 feet below top of well riser, (wells EW-1.25, EW-1.5, EW-2.5, EW-4.5, MW-1I, MW-2I, MW-4S, MW-5S, and MW-9I). These wells were sampled using a Geotech® Geopump II peristaltic pump. Wells IRM-1 and IRM-2I had water surface depths greater than 20 feet below top of well riser and were sampled using a Proactive® Monsoon down-hole centrifugal pump. The below grade portion of the Town Well was not accessible. This well is discussed in the next section.

Prior to initiation of each well purge event, a static water level was measured from the top of the monitoring well riser and recorded on the monitoring well sampling log. At each monitoring well location, (with the exception of Town Well) new polyethylene tubing was lowered into the monitoring well and positioned with the bottom of the tubing at the approximate vertical center of the well screen. Well construction information was taken from the monitoring well logs previously generated by others. Following the sampling efforts, GZA measured and documented the depth of each monitoring well, which were consistent with the information provided on the existing monitoring well logs.

The peristaltic pump/centrifugal pump was started and operated at a flow rate that minimized draw-down of the water column within the well. The first set of water quality readings were collected when the flow-through cell was full and water began to flow out. Once a constant head was established, the pumping flow rate was not altered. Sampling flow rates were kept consistent with purging/monitoring flow rates. Readings were recorded on well development forms in the field, once a constant head had been established. Readings were continuously recorded every five minutes, until water quality readings stabilized for three successive readings, which generally consisted of  $\pm$  0.1 for pH,  $\pm$ 3% for conductivity,  $\pm$  10 mV for oxidation reduction potential (ORP) and  $\pm$  10% for turbidity and dissolved oxygen (DO). Copies of the well purging forms are included in Appendix B.

Groundwater samples were collected for lab analysis once a constant head was established, the water quality readings had stabilized and/or at least one well volume was removed. The polyethylene tubing from the pump to





the water quality meter was disconnected and used to fill the appropriate groundwater sample containers provided by the laboratory. Groundwater collected for analysis did not enter the flow-through cell.

After the appropriate sample containers were filled, the pump was shut off and the tubing was removed from the monitoring well and pump-head. The tubing was then disposed as a solid waste. The flow-through cell and water quality meter were rinsed with de-ionized water prior to use at each well. Water generated during the purging/monitoring and equipment decontamination was placed on the ground in the vicinity of the monitoring well from which it was generated.

### Town Well

The Town Well sample was collected from a spigot within the pump house shed at the well location, as the subsurface portion of this well is not accessible. The spigot was turned on to allow approximately five gallons of water to discharge into a graduated 5-gallon bucket, which was emptied into a floor drain within the shed. The flow-through cell was filled with water directly from the spigot for water quality readings. The sample was collected from the spigot after approximately five gallons were purged and water quality readings were recorded.

### 3.2 GROUNDWATER DATA COLLECTION

GZA collected groundwater samples from the eight on-site monitoring wells (MW-2I, MW-5S, MW-9I, MW-1I, MW-4S, EW-1.25R, EW-1.5, and EW-2.5) and four off-site monitoring wells (EW-4.5, IRM-1, IRM-2I and the Town Well). In addition, a duplicate sample (from IRM-1), and matrix spike/matrix spike duplicate (MS/MSD) sample (from EW-4.5) were collected.

The following table shows the volume of water purged and the number of well volumes removed from the respective wells after a constant head was established. Constant head was not applicable at the Town Well location, as the well was not sampled using low-flow methodologies.

Monitoring Well ID	Volume Purged (gallons)	Well Volumes (#)
EW-1.25R	0.9	0.38
EW-1.5	0.6	0.09
EW-2.5	0.5	0.08
EW-4.5	0.7	0.13
MW-1I	1.0	0.19
MW-2I	1.2	0.22
MW-4S	0.6	0.52
MW-5S	0.6	0.44
MW-9I	0.8	0.13
IRM-1	12.0	2.81
IRM-2I	10.0	2.36
Town Well	5.0	NA

Prior to sampling, static groundwater level measurements were recorded from the top of riser at the 11 accessible monitoring wells (see table below). Monitoring well reference point elevation data were available from previous reports completed by others. Depth to groundwater was measured at each well prior to purging. The measured



groundwater elevations collected during the September 2021 sampling event are shown on Figure 2. Groundwater flow is generally in a south to southeasterly direction, consistent with previous monitoring events.

Monitoring Well	Top of Riser Elevation	Depth to	Groundwater
Location	(ft. AMSL)	Groundwater (ft.)	Elevation (ft. AMSL)
EW-1.25R	1534.04	10.36	1523.68
EW-1.5	1533.92	10.52	1523.40
EW-2.5	1533.92	13.06	1520.86
EW-4.5	1535.65	16.93	1518.72
MW-1I	1531.79	10.61	1524.18
MW-2I	1540.87	15.24	1525.63
MW-4S	1535.42	9.95	1525.47
MW-5S	1534.16	9.07	1525.09
MW-9I	1532.30	10.88	1521.42
IRM-1	1534.75	23.76	1510.99
IRM-2I	1535.99	23.98	1512.01

### 4.0 ANALYTICAL LABORATORY TESTING

Twelve groundwater samples, one duplicate sample (IRM-1), one matrix spike/matrix spike duplicate (EW-4.5), and one trip blank, were submitted for analytical testing. The samples were packed in an ice-filled cooler and, following typical chain-of-custody procedures, sent to Alpha Analytical in Westborough, Massachusetts. Table 1 presents a summary of the samples collected, dates of sample collection, and analyses completed.

### 5.0 ANALYTICAL TEST RESULTS

Discussion of the laboratory results for the groundwater samples is presented below. The laboratory report is provided in Appendix D and summarized on Table 2. Analytical data that were available from January 1989 to September 2021 (specifically trichloroethene (TCE) and tetrachloroethene (PCE)) are summarized on Table 3. These data are also provided graphically, per well location, in Appendix C.





The analytical test results for the groundwater samples were compared to NYSDEC Class GA standards presented in the Division of Water Technical and Operational Guidance Series (TOGS 1.1.1), dated October 1993, revised June 1998, errata January 1999 and amended April 2000.

The analytical data generated as part of the annual monitoring program (the 12 wells) have also been provided to NYSDEC electronically for their Environmental Information Management System (EIMS). The data was provided in a standardized electronic data deliverable (EDD) format that uses the database software application EQuIS<sup>™</sup> (EQuIS) from EarthSoft® Inc. The laboratory data and required information were imported into the EQuIS Data Processor (EDP) and submitted to NYSDEC.

### 5.1 ON SITE WELLS

• <u>EW-1.25R</u>: Three VOCs were detected above method detection limits, but below their respective NYSDEC Class GA (groundwater) criteria (Vinyl Chloride, 1,1-Dichloroethane, cis-1,2-Dichloroethene).

Since the groundwater sampling was reinitiated in 2009, there has been a general downward trend of total VOC concentrations detected at this location. EW-1.25 total VOC mass was similar in June 2016 and July 2017 (86 and 97  $\mu$ g/L, respectively). Additionally, there has been a decrease in VOC mass observed in June 2019 (11  $\mu$ g/L) and September 2021 (3.87  $\mu$ g/L) at EW-1.25R.

- <u>EW-1.5</u>: No VOC's were detected above method detection limits. Since 2009, there has been a general downward trend of total VOC concentrations detected at this location.
- MW-1I: Two VOCs were detected above method detection limits but below their respective Class GA Criteria (1,1-Dichloroethane, cis-1,2-Dichloroethene). Since 2009, there has been a downward trend of VOC concentrations detected at this location and the results have generally been at or below the Class GA criteria since 1996, with the exception of a slight increase in VOC concentrations since the OCEDC injections. As PCE/TCE degradation product cis-DCE is still being produced at this location, we anticipate that VOC concentrations will continue to decline in this well.
- <u>MW-21</u>: No VOC's were detected above method detection limits. Historically, the results have been either below method detection limits or the Class GA criteria since 1994.
- <u>MW-4S</u>: One VOC (PCE) was detected above method detection limits but below its respective Class GA
   Criteria. Historically, the results have been either below method detection limits and/or the Class GA
   criteria since 1998.
- <u>MW-5S</u>: Two VOCs were detected above method detection limits but below their respective Class GA Criteria (PCE and TCE). Since 2009, there has been a downward trend of VOC concentrations detected at this location.

MW-5S is approximately 30 feet southeast and downgradient of AOC-2 and AOC-3 which were addressed as part of IRM activities in 2011 and 2013.



 MW-9I: Two VOCs, (PCE, and TCE) were detected above method detection limits but below their respective NYSDEC Class GA criteria. VOCs have been below Class GA criteria and indicating a downward trend since sampling was reinitiated in 2009.

### 5.2 OFF SITE WELLS

- <u>EW-4.5:</u> Two VOC's (PCE and TCE) were detected above method detection limits but below their respective NYSDEC Class GA criteria. Since 2009, there has been a downward trend of VOC concentrations detected at this location.
- IRM-1: One VOC (TCE) was detected above method detection limits but below the Class GA criteria. Historically, TCE and PCE concentrations have been below the Class GA criteria since 1996.
- IRM-2I: One VOC (TCE) was detected above method detection limits but below the Class GA criteria. Historically, TCE and PCE concentrations have been below the Class GA criteria since 1996.
- <u>Town Well</u>: Three VOCs (Bromodichloromethane, Dibromochloromethane, and TCE) were detected above method detection limits but below their respective Class GA criteria. VOCs have been below Class GA criteria and indicating a downward trend since sampling was reinitiated in 2009.

### 6.0 SUMMARY

A summary of GZAs findings follows:

- Static groundwater level measurements indicate that groundwater flows toward the south/southeast, consistent with previous monitoring events.
- VOCs were not detected at concentrations above NYSDEC Class GA criteria in the groundwater samples collected from the on-site or off-site wells.
- A general downward trend in VOC concentrations since 2009 is noted in monitoring wells EW-1.25/EW-1.25R, EW-1.5, MW-5S and EW-4.5.
- In general, the concentrations of VOCs at monitoring wells EW-2.5, MW-1I, MW-2I, MW-4S, and MW-9I have predominantly been below NYSDEC Class GA criteria.
- Off-site monitoring well results for locations IRM-1, IRM-2 and the Town Well, have consistently been non-detect or at concentrations below Class GA criteria since 1994.

Groundwater monitoring has been conducted for over 20 years. The body of data collected since remedial injections indicate that reductive dechlorination is continuing to reduce the cVOC concentrations as intended, and that a slow and steady overall trend of cVOC reduction has been established. None of the 12 wells monitored have cVOCs at concentrations greater than the Class GA groundwater standards.



### September 2021 Analytical Testing Program Summary Former Signore Facility 55-57 Jefferson Street Ellicottville, New York

Location	Date Collected	Screened Interval (ft bgs)	VOCs EPA Method 8260-TCL
Groundwater Samples			
EW-1.25R	9/17/2021	15-25	X
EW-1.5	9/15/2021	40-50	Х
EW-2.5	9/15/2021	40-50	X
EW-4.5 (MS/MSD)	9/15/2021	40-50	X
MW-1I	9/15/2021	30-50	X
MW-2I	9/16/2021	29-49	X
MW-4S	9/15/2021	7-17	X
MW-5S	9/16/2021	7.5-17.5	X
MW-9I	9/15/2021	29.5-49.5	X
IRM-1	9/16/2021	40-50	X
IRM-2I	9/16/2021	40-50	Χ
TOWN WELL	9/16/2021	NA	X
GW Duplicate (IRM-1)	9/16/2021	40-50	Χ

### Notes:

- 1. ft bgs = feet below ground surface
- 2. VOCs = Volatile Organic Compounds; TCL = Target Compound List
- 3. EPA = Environmental Protection Agency
- 4. MS/MSD = Matrix Spike/Matrix Spike Duplicate

# September 2021 Groundwater Analytical Testing Results Summary Former Signore Facility 55-57 Jefferson Street

Ellicottville, New York

Parameter C										EW-1.2	5 / EW-1.2	25R								
i didiliotoi	Class GA Criteria	4/23/09	10/22/09	6/3/10	4/14/11	10/14/11	5/9/12	10/31/12	6/25/13	10/16/13	6/10/14	10/14/14	6/4/15	10/21/15	6/15/16	10/25/16	7/13/17	6/21/18	6/14/19	9/17/21
Volatile Organic Compounds - E	EPA Method 8260	TCL (ug/L)	)																	
Methylene chloride	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1	<	<
Acetone	50	<	<	<	<	<	<	<	<	<	<	<	<	3.8	2.3 J	<1.5	<1.5	<5.0	6.8	<
2-Butanone	50	<	<	<	<	4.2J	< 5	< 5	< 5	< 5	< 5	< 2	<2	<2	<2	<2	<2	<5.0	<	<
Bromodichloromethane	5	<	<	<	<	٧	< 1	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1	<	<
Dibromochloromethane	50	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1	<	<
Chloromethane	NV	<	<	<	<	<	< 1	< 1	0.77J	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1	0.88 J	<
Chloroform	7	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1		<
Benzene	1	<	<	<	<	<	<	<	<	<	<	<	<	<	<	<	<	<	0.18 J	<
Bromoform	50	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1	<	<
Carbon disulfide	NV	<	<	1.4	<	1.2	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	1.8 J	<1	<	<
Iodomethane	NV	<	<b>'</b>	<	<	٧	< 1	< 1	< 1	< 1	< 1	NT	NT	NT	NT	NT	NT	NT	NT	<
Vinyl Chloride	2	9.7	9.1	8.4	6.3	6	3.8	16	4.6	5	2.4	4.7	2.6	3.3	3.2	6.6	<1	<1	<	0.17 J
1,1-Dichloroethene	5*	<	0.88	0.85	.86J	<b>'</b>	< 1	1.4	< 1	< 1	< 1	0.34 J	0.25 J	0.36 J	0.24 J	0.48 J	0.39 J	<1	<	<
1,1-Dichloroethane	5	8.6	8.7	6.0	6.1	6.7	4.8	5.9	4.1	4.1	2.9	3.8	3	4.2	2.9	3.9	3.0	<1	1.1 J	1.2 J
trans-1, 2-Dichloroethene	5	<	0.92	0.66	.91J	.81J	< 1	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	0.79 J	<1	<1	<	<
cis-1,2-Dichloroethene	5	60	69	39	45	44	32	98	31	32	23	32	29	44	28	98	57	<1	2.1 J	2.5
1,1,1-Trichloroethane	5	1.5	0.82	0.65	.78J	.64J	< 1	2	< 1	< 1	< 1	0.80 J	<1	<1	<1	0.70 J	<1	<1	<	<
Trichloroethene	5	88	90	73	56	90	59	1.7	51	59	41	54	47	58	47	0.27 J	35	<1	<	<
Tetrachloroethene	5	7.5	5.6	5.6	4.2	8.3	5.9	< 1	3.3	3.8	3.6	5.0	3.1	1.8	3.1	<1	0.73	<1	<	<
Naphthalene	10	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	NT	NT	NT	NT	NT	NT	<1	<	<
Total VOCs		175.3	185.0	135.6	120.15	161.85	105.50	125.00	94.77	103.90	72.90	100.64	84.95	115.46	86.74	110.74	97.92		11.06	3.87
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		Т	1	ı	1		1			N	/W-4S	1		1				1	1	ı
Parameter C	Class GA Criteria	4/23/09	10/22/09	6/2/10	4/14/11	10/13/11	5/10/12	10/31/12	6/25/13	10/15/13	6/6/14	10/15/14	6/3/15	10/21/15	6/15/16	10/25/16	7/12/17	6/20/18	6/11/19	9/15/21
Volatile Organic Compounds - E	EPA Method 8260	TCL (ug/L)	)																	
Methylene chloride	5	<																		
Acetone	50	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	< 1	<	<
	50	<	< <	< <	< <	< <	< 1	< 1	< 1	< 1 <	< 1	< 1	<1	<1 2.3 J	<1 <	<1 <	<1 <	< 1 < 5	< 3.0 J	< <
2-Butanone	50	,																		
Bromodichloromethane		<	<	<	<	<	<	<	<	<	<	<	<	2.3 J	<	<	<	< 5	3.0 J	<
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Bromodichloromethane	50 5	< < <	< <	< < <	< < <	< <	< 5 < 1	< 5 < 1	< 5 < 1	< 5 < 1	< 5 < 1	< 2 < 1	<2 <1	2.3 J <2 <1	<2 <1	< <2 <1	< <2 <1	< 5 < 5 < 1	3.0 J <	< < <
Bromodichloromethane Dibromochloromethane Chloromethane Chloroform	50 5 50 NV 7	< < < < < < < < < < < < < < < < < < <	< < < < < < < < < < < < < < < < < < <	< < < < < < < < < < < < < < < < < < <	<	< < < < < < < < < < < < < < < < < < <	< 5 < 1 < 1 < 1 < 1 < 1	< 5 < 1 < 1 < 1 < 1 < 1	< 5 < 1 < 1 < 1 < 1 < 1	< 5 < 1 < 1 < 1 < 1	< 5 < 1 < 1 < 1 < 1 < 1	< 2 < 1 < 1 < 1 < 1 < 1	<2 <1 <1 <1 <1 <1	2.3 J <2 <1 <1 <1 <1	<pre>&lt;   &lt;2   &lt;1   &lt;1   &lt;1   &lt;1   &lt;1   &lt;1   &lt;</pre>	< 2 <1 <1 <1 <1 <1 <1	<	< 5 < 5 < 1 < 1 < 1 < 1	3.0 J < < < 1.2 J	< < < < < < < < < < < < < < < < < < <
Bromodichloromethane Dibromochloromethane Chloromethane Chloroform Bromoform	50 5 50 NV 7 50	< < < < < < < < < < < < < < < < < < <	< < < <	<	< < < < < < < < < < < < < < < < < < <	< < < <	< 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1	< 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1	< 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1	< 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1	< 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1	<pre></pre>	<pre>&lt;   &lt;2   &lt;1   &lt;1   &lt;1   &lt;1   &lt;1   &lt;1   &lt;</pre>	2.3 J <2 <1 <1 <1 <1 <1	<pre>&lt; c2 &lt;1 </pre>	<pre></pre>	<	< 5 < 5 < 1 < 1 < 1 < 1 < 1	3.0 J < < < 1.2 J	<
Bromodichloromethane Dibromochloromethane Chloromethane Chloroform Bromoform Carbon Disulfide	50 5 50 NV 7 50 NV	< < < < < < < < < < < < < < < < < < <	< < < < < < < < < < < < < < < < < < <	< < < < < < < < < < < < < < < < < < <	<td>&lt;</td> <td>&lt; 5 &lt; 1 &lt; 1</td> <td>&lt; 5 &lt; 1 &lt; 1</td> <td>&lt; 5 &lt; 1 &lt; 1</td> <td>&lt; 5 &lt;1 &lt;1</td> <td>&lt; 5 &lt; 1 &lt; 1</td> <td><pre></pre></td> <td>&lt; &lt;2 &lt;1 &lt;</td> <td>2.3 J &lt;2 &lt;1 &lt;1</td> <td><pre> &lt;      &lt;2      &lt;1      &lt;</pre></td> <td><pre></pre></td> <td>&lt;</td> <td>&lt;5 &lt;5 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td> <td>3.0 J &lt; &lt; &lt; 1.2 J &lt; &lt;</td> <td></td>	<	< 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1	< 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1	< 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1	< 5 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	< 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1	<pre></pre>	< <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <	2.3 J <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<pre> &lt;      &lt;2      &lt;1      &lt;</pre>	<pre></pre>	<	<5 <5 <1 <1 <1 <1 <1 <1	3.0 J < < < 1.2 J < <	
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Bromodichloromethane Dibromochloromethane Chloromethane Chloroform Bromoform Carbon Disulfide Iodomethane Vinyl Chloride 1,1-Dichloroethene 1,1-Dichloroethane trans-1, 2-Dichloroethene cis-1,2-Dichloroethene 1,1,1-Trichloroethane Trichloroethene Trichloroethene Tetrachloroethene	50 5 50 NV 7 50 NV NV 2 5* 5 5 5 5 5	<pre></pre>	<pre></pre>	<pre></pre>	<pre></pre>	<td><pre></pre></td> <td><pre></pre></td> <td><pre></pre></td> <td><pre></pre></td> <td><pre></pre></td> <td><pre></pre></td> <td><pre> &lt;</pre></td> <td>2.3 J</td> <td><pre></pre></td> <td>&lt;</td> <td><pre></pre></td> <td>&lt;5 &lt;5 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td> <td>3.0 J</td> <td><pre></pre></td>	<pre></pre>	<pre></pre>	<pre></pre>	<pre></pre>	<pre></pre>	<pre></pre>	<pre> &lt;</pre>	2.3 J	<pre></pre>	<	<pre></pre>	<5 <5 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	3.0 J	<pre></pre>
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### Notes

- 1. Compounds detected in one or more samples are presented on this table.
- 2. Analytical testing completed by Alpha Analytical.
- 3. NYSDEC Class GA criteria obtained from Division of Water Technical and Operational Guidance Series (TOGS 1.1.1), dated October 1993, revised June 1998, January 1999 errata sheet,
- and April 2000 addendum. \* Guidance value (not a standard) for 1,1-Dichloroethene = 0.07 ug/L as per the January 1999 update.
- 4. ug/L = part per billion (ppb).
- 5. < indicates compound was not detected; < 1 indicates compound was not detected above its respective reporting limit.
- 6. Shading indicates exceedance of Class GA Criteria.
- 7. NT = not tested.
- 8. NV = no value.
- 9. Results shown for IRM-1 for the September 2021 sampling event are the higher results from it or its respective duplicate.
- 10. Lab qualifiers: CH = continuing calibration outside of lab acceptance limits; results may be biased high. J = estimated concentration.
- L2 = analyte recovery in the control sample was below quality control limits; results may be biased low. Qualifiers for detected compounds only shown.

# September 2021 Groundwater Analytical Testing Results Summary Former Signore Facility

55-57 Jefferson Street Ellicottville, New York

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Parameter	Class GA Criteria	4/23/09	10/22/09	6/2/10	4/14/11	10/14/11	5/9/12	10/31/12	6/25/13	10/16/13	6/9/14	10/14/14	6/2/15	10/21/15	6/14/16	10/25/16	7/11/17	6/19/18	6/13/19	9/15/21
Volatile Organic Compound	ls - EPA Method 8260				<u> </u>				<u> </u>		<u> </u>	<u> </u>		<u> </u>	<u> </u>			<u> </u>		
Methylene chloride	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
Acetone	50	<	<	<	<	<	<	<	<	<	<	<	<	1.5 J	< 1.5	< 1.5	< 1.5	<5.0	3.0 J	<
2-Butanone	50	<	<	<	<	<	< 5	< 5	< 5	< 5	< 5	< 2	< 2	< 2	< 2	< 2	< 2	<5.0	<	<
Bromodichloromethane	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
Dibromochloromethane	50	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
Chloromethane	NV	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
Chloroform	7	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
Benzene	1	<	<	<	<	<	<	<	<	<	<	<	<	<	<	<	<	<	<	<
Bromoform	50	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
Carbon disulfide	NV	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1	<1	<	<
Iodomethane	NV	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	NT	NT	NT	NT	NT	NT	NT	<	<
Vinvl Chloride	2	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
1.1-Dichloroethene	5*	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
1,1-Dichloroethane	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
trans-1, 2-Dichloroethene	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
cis-1,2-Dichloroethene	5	2.1	4.6	2.2	3.3	1.7	2.1	2.9	1.3	< 1	1.6	2.7	2.0 J	2.1 J	1.6 J	1.2 J	1.3 J	<1	<	<
1.1.1-Trichloroethane	5	4.1	2.7	1.9	2.6	1.3	1.7	< 1	1.2	< 1	< 1	1.4 J	1.2 J	1.2 J	<1	0.90 J	1.2 J	<1	<	<
Trichloroethene	5	18	20	14	19	9.5	13.0	9.0	8.4	3.9	10	13	13	11	6.4	10	10	<1	<	<
Tetrachloroethene	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	0.22 J	0.20 J	0.22 J	<1	0.24 J	0.23 J	<1	<	<
Naphthalene	10	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	NT	NT	NT	NT	NT	NT	<1	<	<
Total VOCs		24.2	27.3	18.1	24.9	12.5	16.8	11.9	10.9	3.9	11.6	17.32	16.30	16.02	8.00	12.34	12.73		3.00	
	•		•																	l.
Parameter										M	IW-5S									
Farameter	Class GA Criteria	4/23/09	10/22/09	6/3/10	4/14/11	10/13/11	5/9/12	10/31/12	6/25/13	10/15/13	6/6/14	10/14/14	6/2/15	10/22/15	6/15/16	10/24/16	7/12/17	6/20/18	6/11/19	9/16/21
Volatile Organic Compound			10/22/09	6/3/10	4/14/11	10/13/11	5/9/12	10/31/12	6/25/13			10/14/14	6/2/15	10/22/15	6/15/16	10/24/16	7/12/17	6/20/18	6/11/19	9/16/21
			10/22/09	6/3/10	4/14/11	10/13/11	5/9/12	10/31/12	6/25/13			10/14/14	6/2/15	10/22/15	6/15/16	10/24/16	7/12/17	6/20/18	6/11/19	9/16/21
Volatile Organic Compound Methylene chloride	ls - EPA Method 8260									10/15/13	6/6/14									
Volatile Organic Compound	ls - EPA Method 8260	<	<	<	<	<	< 1	< 1	< 1	10/15/13	6/6/14	< 1	< 1	< 1	< 1	< 1	<1	< 1	<	<
Volatile Organic Compound Methylene chloride Acetone	1s - EPA Method 8260 5 50	< <	< <	< <	< <	< <	<1 <	< 1	< 1	10/15/13	6/6/14 < 1 <	< 1	< 1	< 1 4 J	< 1 3.4 J	< 1 <1.5	<1 <	< 1 < 5	< 1.6 J	< <
Volatile Organic Compound Methylene chloride Acetone 2-Butanone	5 5 50 50	< < <	< < <	< < <	< <	< < <	< 1 < < 5	< 1 < < 5	< 1 < < 5	10/15/13	6/6/14 <1 < 5	< 1 < 2	< 1 < < 2	< 1 4 J < 2	< 1 3.4 J < 2	<1 <1.5 <2	<1 < <2	< 1 < 5 < 5	< 1.6 J	< <
Volatile Organic Compound Methylene chloride Acetone 2-Butanone Bromodichloromethane	5 5 50 50 50	< < < < < < < < < < < < < < < < < < <	< < < < < < < < < < < < < < < < < < <	< < < < < < < < < < < < < < < < < < <	< < <	< < < < < < < < < < < < < < < < < < <	< 1 < < 5 < 1	< 1 < < 5 < 1	<1 < <5 <1	10/15/13 < 1 < 5 < 1	6/6/14 < 1 < 5 < 1	<1 < <2 <1	<1 < <2 <1	< 1 4 J < 2 < 1	< 1 3.4 J < 2 < 1	<1 <1.5 <2 <1	<1	< 1 < 5 < 5 < 1	< 1.6 J < <	< < <
Volatile Organic Compound Methylene chloride Acetone 2-Butanone Bromodichloromethane Dibromochloromethane	5 50 50 50 50	< < < < < < < < < < < < < < < < < < <	< < < < < < < < < < < < < < < < < < <	< < < < < < < < < < < < < < < < < < <	< < < < < < < < < < < < < < < < < < <	< < < < < < < < < < < < < < < < < < <	<1 <1 <5 <1 <1 <1	<1 < <5 <1 <1	<1 < 1 < 5 < 1 < 1 < 1	10/15/13  < 1   < 5   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   <	<pre>6/6/14 &lt; 1</pre>	<1 < 2 < 1 < 1 < 1 < 1	<1 < <2 <1 <1	< 1 4 J < 2 < 1 < 1	<1 3.4 J <2 <1 <1	<1 <1.5 <2 <1 <1	<1 <1 <2 <1 <1 <1	<1 <5 <5 <1 <1	< 1.6 J < < < < < < < < < < < < < < < < < <	< < < < < < < < < < < < < < < < < < <
Volatile Organic Compound Methylene chloride Acetone 2-Butanone Bromodichloromethane Dibromochloromethane Chloromethane	5 50 50 50 NV	<	<	<	< < < < < < < < < < < < < < < < < < <	<	< 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1	< 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1	< 1 < 5 < 1 < 1 0.99J	10/15/13  <1 <1 <5 <1 <1 <1 <1	<pre>6/6/14  &lt;1   &lt; 5   &lt; 1   &lt; 1</pre>	<1 <1 <2 <1 <1 <1 <1	<1 < 2 < 1 < 1 < 1 < 1 < 1 < 1	< 1 4 J < 2 < 1 < 1	<1 3.4 J <2 <1 <1	<1 <1.5 <2 <1 <1	<1 <1 <2 <1 <1 <1 <1 <1 <1	< 1 < 5 < 5 < 1 < 1		<
Volatile Organic Compound Methylene chloride Acetone 2-Butanone Bromodichloromethane Dibromochloromethane Chloromethane Chloroform	5 50 50 NV 7	<	<td>&lt;</td> <td>&lt;</td> <td>&lt;</td> <td>&lt;1 &lt;1 &lt;</td> <td>&lt;1 &lt;1 &lt;</td> <td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt; 1 &lt; 1 0.99J &lt; 1</td> <td>10/15/13  &lt;1 &lt;1 &lt;5 &lt;1 &lt;1</td> <td><pre>6/6/14  &lt;1 &lt;5 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 </pre></td> <td>&lt;1 &lt;1 &lt;2 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td> <td>&lt;1 &lt; 2 &lt; 1 &lt; 1 &lt; 1 &lt; 1 &lt; 1 &lt; 1 &lt; 1</td> <td>&lt; 1 4 J &lt; 2 &lt; 1 &lt; 1 &lt; 1</td> <td>&lt;1 3.4 J &lt;2 &lt;1 &lt;1 1.2 J &lt;1</td> <td>&lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1 &lt;1 &lt;1</td> <td>&lt;1 &lt; &lt;</td> <td>&lt;1 &lt;5 &lt;5 &lt;1 &lt;1 &lt;1 &lt;1</td> <td><pre></pre></td> <td>&lt;</td>	<	<	<	<1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <	<1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <	<1 < 1 < 5 < 1 < 1 < 1 0.99J < 1	10/15/13  <1 <1 <5 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<pre>6/6/14  &lt;1 &lt;5 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 </pre>	<1 <1 <2 <1 <1 <1 <1 <1 <1 <1	<1 < 2 < 1 < 1 < 1 < 1 < 1 < 1 < 1	< 1 4 J < 2 < 1 < 1 < 1	<1 3.4 J <2 <1 <1 1.2 J <1	<1 <1.5 <2 <1 <1 <1 <1	<1 < < < < < < < < < < < < < < < < < <	<1 <5 <5 <1 <1 <1 <1	<pre></pre>	<
Volatile Organic Compound Methylene chloride Acetone 2-Butanone Bromodichloromethane Dibromochloromethane Chloromethane Chloroform Bromoform	5 - EPA Method 8260 5 - 50 50 - 50 5 - 50 NV 7 - 50	<	<td><td>&lt;</td><td><td>&lt;1 &lt;1 &lt;&lt; 5 &lt;1 &lt;1</td><td>&lt;1 &lt;1 &lt;&lt; &lt;5 &lt;1 &lt;1</td><td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt; 1 &lt; 1 &lt; 0.99J &lt; 1 &lt; 1</td><td>10/15/13  &lt;1 &lt;1 &lt;5 &lt;1 &lt;1</td><td><pre>6/6/14  &lt;1     &lt; 5     &lt; 1     &lt;</pre></td><td>&lt;1 &lt;1 &lt;&lt; 2 &lt;1 &lt;1</td><td>&lt;1 &lt; 1 &lt; 2 &lt; 1 &lt;</td><td>&lt;1 4 J &lt; 2 &lt; 1 &lt; 1 &lt; 1 &lt; 1 &lt; 1</td><td>&lt;1 3.4 J &lt;2 &lt;1 &lt;1 1.2 J &lt;1</td><td>&lt;1 &lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td><td>&lt;1 &lt;1 &lt;2 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td><td>&lt;1 &lt;5 &lt;5 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td><td>&lt; 1.6 J &lt; &lt;</td><td>&lt;</td></td></td>	<td>&lt;</td> <td><td>&lt;1 &lt;1 &lt;&lt; 5 &lt;1 &lt;1</td><td>&lt;1 &lt;1 &lt;&lt; &lt;5 &lt;1 &lt;1</td><td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt; 1 &lt; 1 &lt; 0.99J &lt; 1 &lt; 1</td><td>10/15/13  &lt;1 &lt;1 &lt;5 &lt;1 &lt;1</td><td><pre>6/6/14  &lt;1     &lt; 5     &lt; 1     &lt;</pre></td><td>&lt;1 &lt;1 &lt;&lt; 2 &lt;1 &lt;1</td><td>&lt;1 &lt; 1 &lt; 2 &lt; 1 &lt;</td><td>&lt;1 4 J &lt; 2 &lt; 1 &lt; 1 &lt; 1 &lt; 1 &lt; 1</td><td>&lt;1 3.4 J &lt;2 &lt;1 &lt;1 1.2 J &lt;1</td><td>&lt;1 &lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td><td>&lt;1 &lt;1 &lt;2 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td><td>&lt;1 &lt;5 &lt;5 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td><td>&lt; 1.6 J &lt; &lt;</td><td>&lt;</td></td>	<	<td>&lt;1 &lt;1 &lt;&lt; 5 &lt;1 &lt;1</td> <td>&lt;1 &lt;1 &lt;&lt; &lt;5 &lt;1 &lt;1</td> <td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt; 1 &lt; 1 &lt; 0.99J &lt; 1 &lt; 1</td> <td>10/15/13  &lt;1 &lt;1 &lt;5 &lt;1 &lt;1</td> <td><pre>6/6/14  &lt;1     &lt; 5     &lt; 1     &lt;</pre></td> <td>&lt;1 &lt;1 &lt;&lt; 2 &lt;1 &lt;1</td> <td>&lt;1 &lt; 1 &lt; 2 &lt; 1 &lt;</td> <td>&lt;1 4 J &lt; 2 &lt; 1 &lt; 1 &lt; 1 &lt; 1 &lt; 1</td> <td>&lt;1 3.4 J &lt;2 &lt;1 &lt;1 1.2 J &lt;1</td> <td>&lt;1 &lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td> <td>&lt;1 &lt;1 &lt;2 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td> <td>&lt;1 &lt;5 &lt;5 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td> <td>&lt; 1.6 J &lt; &lt;</td> <td>&lt;</td>	<1 <1 << 5 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1 << <5 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 < 1 < 5 < 1 < 1 < 1 < 0.99J < 1 < 1	10/15/13  <1 <1 <5 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<pre>6/6/14  &lt;1     &lt; 5     &lt; 1     &lt;</pre>	<1 <1 << 2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 < 1 < 2 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 4 J < 2 < 1 < 1 < 1 < 1 < 1	<1 3.4 J <2 <1 <1 1.2 J <1	<1 <1 <1.5 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <5 <5 <1 <1 <1 <1 <1 <1	< 1.6 J < < < < < < < < < < < < < < < < < <	<
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Volatile Organic Compound Methylene chloride Acetone 2-Butanone Bromodichloromethane Dibromochloromethane Chloromethane Chloroform Bromoform Carbon Disulfide Iodomethane Vinyl Chloride	5 - EPA Method 8260 5 - 50 50 50 50 NV 7 - 50 NV NV NV 2	<td><td><td><td><td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td><td>&lt;1 &lt;1 &lt;</td><td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td><td>10/15/13  &lt;1 &lt;1 &lt;5 &lt;1 &lt;1</td><td><pre>6/6/14  &lt;1     &lt; 1     &lt; 5     &lt; 1     &lt;</pre></td><td>&lt;1 &lt;1 &lt; 2 &lt;1 &lt;1</td><td>&lt;1 &lt; 1 &lt; 2 &lt; 1 &lt;</td><td>&lt;1 4 J &lt; 2 &lt; 1 &lt; 1</td><td>&lt;1 3.4 J &lt;2 &lt;1 &lt;1 &lt;1 1.2 J &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td><td>&lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td><td>&lt;1 &lt;1 &lt;2 &lt;1 &lt;1</td><td>&lt;1 &lt;5 &lt;5 &lt;1 &lt;1</td><td><pre></pre></td><td></td></td></td></td></td>	<td><td><td><td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td><td>&lt;1 &lt;1 &lt;</td><td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td><td>10/15/13  &lt;1 &lt;1 &lt;5 &lt;1 &lt;1</td><td><pre>6/6/14  &lt;1     &lt; 1     &lt; 5     &lt; 1     &lt;</pre></td><td>&lt;1 &lt;1 &lt; 2 &lt;1 &lt;1</td><td>&lt;1 &lt; 1 &lt; 2 &lt; 1 &lt;</td><td>&lt;1 4 J &lt; 2 &lt; 1 &lt; 1</td><td>&lt;1 3.4 J &lt;2 &lt;1 &lt;1 &lt;1 1.2 J &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td><td>&lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td><td>&lt;1 &lt;1 &lt;2 &lt;1 &lt;1</td><td>&lt;1 &lt;5 &lt;5 &lt;1 &lt;1</td><td><pre></pre></td><td></td></td></td></td>	<td><td><td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td><td>&lt;1 &lt;1 &lt;</td><td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td><td>10/15/13  &lt;1 &lt;1 &lt;5 &lt;1 &lt;1</td><td><pre>6/6/14  &lt;1     &lt; 1     &lt; 5     &lt; 1     &lt;</pre></td><td>&lt;1 &lt;1 &lt; 2 &lt;1 &lt;1</td><td>&lt;1 &lt; 1 &lt; 2 &lt; 1 &lt;</td><td>&lt;1 4 J &lt; 2 &lt; 1 &lt; 1</td><td>&lt;1 3.4 J &lt;2 &lt;1 &lt;1 &lt;1 1.2 J &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td><td>&lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td><td>&lt;1 &lt;1 &lt;2 &lt;1 &lt;1</td><td>&lt;1 &lt;5 &lt;5 &lt;1 &lt;1</td><td><pre></pre></td><td></td></td></td>	<td><td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td><td>&lt;1 &lt;1 &lt;</td><td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td><td>10/15/13  &lt;1 &lt;1 &lt;5 &lt;1 &lt;1</td><td><pre>6/6/14  &lt;1     &lt; 1     &lt; 5     &lt; 1     &lt;</pre></td><td>&lt;1 &lt;1 &lt; 2 &lt;1 &lt;1</td><td>&lt;1 &lt; 1 &lt; 2 &lt; 1 &lt;</td><td>&lt;1 4 J &lt; 2 &lt; 1 &lt; 1</td><td>&lt;1 3.4 J &lt;2 &lt;1 &lt;1 &lt;1 1.2 J &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td><td>&lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td><td>&lt;1 &lt;1 &lt;2 &lt;1 &lt;1</td><td>&lt;1 &lt;5 &lt;5 &lt;1 &lt;1</td><td><pre></pre></td><td></td></td>	<td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td> <td>&lt;1 &lt;1 &lt;</td> <td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td> <td>10/15/13  &lt;1 &lt;1 &lt;5 &lt;1 &lt;1</td> <td><pre>6/6/14  &lt;1     &lt; 1     &lt; 5     &lt; 1     &lt;</pre></td> <td>&lt;1 &lt;1 &lt; 2 &lt;1 &lt;1</td> <td>&lt;1 &lt; 1 &lt; 2 &lt; 1 &lt;</td> <td>&lt;1 4 J &lt; 2 &lt; 1 &lt; 1</td> <td>&lt;1 3.4 J &lt;2 &lt;1 &lt;1 &lt;1 1.2 J &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td> <td>&lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td> <td>&lt;1 &lt;1 &lt;2 &lt;1 &lt;1</td> <td>&lt;1 &lt;5 &lt;5 &lt;1 &lt;1</td> <td><pre></pre></td> <td></td>	<1 < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <	<1 < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	10/15/13  <1 <1 <5 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<pre>6/6/14  &lt;1     &lt; 1     &lt; 5     &lt; 1     &lt;</pre>	<1 <1 < 2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 < 1 < 2 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 4 J < 2 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1	<1 3.4 J <2 <1 <1 <1 1.2 J <1 <1 <1 <1 <1	<1 <1.5 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <5 <5 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<pre></pre>	
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Volatile Organic Compound Methylene chloride Acetone 2-Butanone Bromodichloromethane Dibromochloromethane Chloroform Bromoform Carbon Disulfide Iodomethane Vinyl Chloride 1,1-Dichloroethane trans-1, 2-Dichloroethene	5 5 50 50 50 50 50 NV 7 50 NV 7 50 NV NV 2 5* 5	<td><td><td><td><td>&lt;1 &lt;1 &lt;5 &lt;1 &lt;1</td><td>&lt;1 &lt;1 &lt;5 &lt;1 &lt;1</td><td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td><td>10/15/13  &lt; 1 &lt; 5 &lt; 1 &lt; 2 &lt;</td><td>6/6/14  &lt; 1     &lt; 5     &lt; 1     &lt; 1</td><td>&lt;1 &lt;1 &lt;2 &lt;1 &lt;1</td><td>&lt;1 &lt; 1 &lt; 2 &lt; 1 &lt;</td><td>&lt;1 4 J</td><td>&lt;1 3.4 J &lt;2 &lt;1 &lt;1 &lt;1 2 J &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td><td>&lt;1 &lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td><td>&lt;1 &lt;1 &lt; 2 &lt;1 &lt;1</td><td>&lt;1 &lt;5 &lt;5 &lt;1 &lt;1</td><td><pre></pre></td><td></td></td></td></td></td>	<td><td><td><td>&lt;1 &lt;1 &lt;5 &lt;1 &lt;1</td><td>&lt;1 &lt;1 &lt;5 &lt;1 &lt;1</td><td>&lt;1 &lt; 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Volatile Organic Compound Methylene chloride Acetone 2-Butanone Bromodichloromethane Dibromochloromethane Chloroform Bromoform Carbon Disulfide Iodomethane Vinyl Chloride 1,1-Dichloroethene 1,1-Dichloroethane trans-1, 2-Dichloroethene cis-1,2-Dichloroethene 1,1,1-Trichloroethane	5 5 50 50 50 50 NV 7 50 NV NV NV NV 2 5* 5 5	< < < < < < < < < < <    3.4	<pre></pre>	<td><td><td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td><td>&lt;1 &lt;1 &lt; 5 &lt;1 &lt;1</td><td>&lt;1 &lt; 1 &lt; 5 &lt; 1 0.99J &lt; 1 &lt;</td><td>10/15/13  &lt;1 &lt;5 &lt;1 &lt;1</td><td>6/6/14  &lt;1 &lt;1</td><td>&lt;1 &lt;1 &lt;2 &lt;1 &lt;1</td><td>&lt;1 &lt; 1 &lt; 2 &lt; 1 &lt;</td><td>&lt;1 4 J</td><td>&lt;1 3.4 J &lt;2 &lt;1 &lt;1</td><td>&lt;1 &lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td><td>&lt;1 &lt;1 &lt; 2 &lt;1 &lt;1</td><td>&lt;1 &lt;5 &lt;5 &lt;1 &lt;1</td><td><pre></pre></td><td></td></td></td>	<td><td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td><td>&lt;1 &lt;1 &lt; 5 &lt;1 &lt;1</td><td>&lt;1 &lt; 1 &lt; 5 &lt; 1 0.99J &lt; 1 &lt;</td><td>10/15/13  &lt;1 &lt;5 &lt;1 &lt;1</td><td>6/6/14  &lt;1 &lt;1</td><td>&lt;1 &lt;1 &lt;2 &lt;1 &lt;1</td><td>&lt;1 &lt; 1 &lt; 2 &lt; 1 &lt;</td><td>&lt;1 4 J</td><td>&lt;1 3.4 J &lt;2 &lt;1 &lt;1</td><td>&lt;1 &lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td><td>&lt;1 &lt;1 &lt; 2 &lt;1 &lt;1</td><td>&lt;1 &lt;5 &lt;5 &lt;1 &lt;1</td><td><pre></pre></td><td></td></td>	<td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td> <td>&lt;1 &lt;1 &lt; 5 &lt;1 &lt;1</td> <td>&lt;1 &lt; 1 &lt; 5 &lt; 1 0.99J &lt; 1 &lt;</td> <td>10/15/13  &lt;1 &lt;5 &lt;1 &lt;1</td> <td>6/6/14  &lt;1 &lt;1</td> <td>&lt;1 &lt;1 &lt;2 &lt;1 &lt;1</td> <td>&lt;1 &lt; 1 &lt; 2 &lt; 1 &lt;</td> <td>&lt;1 4 J</td> <td>&lt;1 3.4 J &lt;2 &lt;1 &lt;1</td> <td>&lt;1 &lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td> <td>&lt;1 &lt;1 &lt; 2 &lt;1 &lt;1</td> <td>&lt;1 &lt;5 &lt;5 &lt;1 &lt;1</td> <td><pre></pre></td> <td></td>	<1 < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 <1 < 5 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 < 1 < 5 < 1 0.99J < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	10/15/13  <1 <5 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	6/6/14  <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 < 1 < 2 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 4 J	<1 3.4 J <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1 <1.5 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1 < 2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <5 <5 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<pre></pre>	
Volatile Organic Compound Methylene chloride Acetone 2-Butanone Bromodichloromethane Dibromochloromethane Chloroform Bromoform Carbon Disulfide Iodomethane Vinyl Chloride 1,1-Dichloroethene 1,1-Dichloroethene trans-1, 2-Dichloroethene cis-1,2-Dichloroethene 1,1,1-Trichloroethane Trichloroethene	5 50 50 50 85 - EPA Method 8260 50 50 80 80 80 80 80 80 80 80 80 80 80 80 80	<pre></pre>	<pre></pre>	<pre></pre>	<pre></pre>	<td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td> <td>&lt;1 &lt;1 &lt; 5 &lt;1 &lt;1</td> <td>&lt;1 &lt; 1 &lt; 5 &lt; 1 0.99J &lt; 1 &lt;</td> <td>10/15/13  &lt;1 &lt;1</td> <td>6/6/14  &lt;1 &lt;1</td> <td>&lt;1 &lt;1 &lt;2 &lt;1 &lt;1</td> <td>&lt;1 &lt; 1 &lt; 2 &lt; 1 &lt;</td> <td>&lt;1 4 J</td> <td>&lt;1 3.4 J &lt;2 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td> <td>&lt;1 &lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td> <td>&lt;1 &lt;1 &lt; 2 &lt;1 &lt;1</td> <td>&lt;1 &lt;5 &lt;5 &lt;1 &lt;1</td> <td><pre></pre></td> <td><pre></pre></td>	<1 < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 <1 < 5 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 < 1 < 5 < 1 0.99J < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	10/15/13  <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	6/6/14  <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 < 1 < 2 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 4 J	<1 3.4 J <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1 <1.5 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1 < 2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <5 <5 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<pre></pre>	<pre></pre>
Volatile Organic Compound Methylene chloride Acetone 2-Butanone Bromodichloromethane Dibromochloromethane Chloroform Bromoform Carbon Disulfide Iodomethane Vinyl Chloride 1,1-Dichloroethene 1,1-Dichloroethene trans-1, 2-Dichloroethene cis-1,2-Dichloroethene 1,1,1-Trichloroethane Trichloroethene Trichloroethene Tetrachloroethene Tetrachloroethene	5 50 50 50 85 50 85 50 80 80 80 80 80 80 80 80 80 80 80 80 80	<pre></pre>	<pre></pre>	<pre></pre>	<pre></pre>	<pre></pre>	<1 < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 <1 <5 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	10/15/13  <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	6/6/14  <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1 < 2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 4 J	<1 3.4 J <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1 <1.5 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1 < 2 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <5 <5 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<pre></pre>	<pre></pre>

### Notes

- 1. Compounds detected in one or more samples are presented on this table.
- 2. Analytical testing completed by Alpha Analytical.
- 3. NYSDEC Class GA criteria obtained from Division of Water Technical and Operational Guidance Series (TOGS 1.1.1), dated October 1993, revised June 1998, January 1999 errata sheet,

and April 2000 addendum. \* Guidance value (not a standard) for 1,1-Dichloroethene = 0.07 ug/L as per the January 1999 update.

- 4. ug/L = part per billion (ppb).
- 5. < indicates compound was not detected; < 1 indicates compound was not detected above its respective reporting limit.
- 6. Shading indicates exceedance of Class GA Criteria.
- 7. NT = not tested.
- 8. NV = no value.
- 9. Results shown for IRM-1 for the September 2021 sampling event are the higher results from it or its respective duplicate.
- 10. Lab qualifiers: CH = continuing calibration outside of lab acceptance limits; results may be biased high. J = estimated concentration.
- L2 = analyte recovery in the control sample was below quality control limits; results may be biased low. Qualifiers for detected compounds only shown.

# September 2021 Groundwater Analytical Testing Results Summary Former Signore Facility 55-57 Jefferson Street

Ellicottville, New York

											EW-2.5									
Parameter	Class GA Criteria	4/23/09	10/22/09	6/2/10	4/13/11	10/13/11	5/9/12	11/1/12	6/26/13	10/17/13	6/9/14	10/15/14	6/2/15	10/21/15	6/14/16	10/24/16	7/11/17	6/19/18	6/13/19	9/15/21
Volatile Organic Compound	s - EPA Method 8260																			
Methylene chloride	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
Acetone	50	<	<	<	<	<	<	<	<	<	<	<	<	2.4 J	1.7 J	<1.5	<1.5	<5.0	2.3 J	<
2-Butanone	50	<	<	<	<	<	< 5	< 5	< 5	< 5	< 5	< 2	< 2	< 2	< 2	< 2	< 2	<5.0	<	<
Bromodichloromethane	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
Dibromochloromethane	50	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
Chloromethane	NV	<	<	<	<	<	< 1	< 1	1.4	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
Chloroform	7	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
Benzene	1	<	<	<	<	<	<	<	<	<	<	<	<	<	<	<	<	<	<	<
Bromoform	50	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
Carbon disulfide	NV	<	<	<	0.94 J	<	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1	<1	<	<
Iodomethane	NV	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1	NT	<	<
Vinyl Chloride	2	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
1,1-Dichloroethene	5*	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
1,1-Dichloroethane	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
trans-1, 2-Dichloroethene	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
cis-1,2-Dichloroethene	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
1.1.1-Trichloroethane	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
Trichloroethene	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1	<1	<	<
Tetrachloroethene	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1	<1	<	<
Naphthalene	10	<	<	<	1.3	<	< 1	< 1	< 1	< 1	< 1	NT	NT	NT	NT	NT	NT	<1	<	<
Total VOCs					2.2				1.4					2.4	1.7				2.30	
					•						MW-9I				,				•	
Parameter	Class GA Criteria	4/23/09	10/22/09	6/2/10	4/14/11	10/13/11	5/9/12	11/1/12	6/25/13	10/15/13	6/9/14	10/15/14	6/3/15	10/22/15	6/14/16				6/13/19	0/45/04
<b>Volatile Organic Compound</b>	s - EPA Method 8260								0,20,10	10/13/13					0/14/10	10/24/16	7/11/17	6/20/18	0/13/19	9/15/21
Methylene chloride	_								0,20,10	10/13/13				10,22,10	0/14/10	10/24/16	7/11/17	6/20/18	0/13/19	9/15/21
Acetone	5	<	<	<	<	<	< 1	< 1	< 1	< 1	<1	< 1	< 1	< 1	< 1	< 1	7/11/17	6/20/18	<	9/15/21
	50		< <	< <	< <	< <	< 1	< 1			<1	<1 <	< 1							
2-Butanone	_	<			1				< 1	< 1				< 1	< 1	< 1	< 1	< 1	<	<
2-Butanone Bromodichloromethane	50	< <	<	<	<	<	<	<	< 1	< 1	<	<	<	< 1 2.7 J	< 1 1.6 J	< 1 <1.5	< 1 <1.5	< 1 < 5	< 1.9 J	< <
	50 50	< < <	< <	< <	< <	< <	< < 5	< < 5	< 1 < < 5	< 1 < < 5	< < 5	< <2	< <2	< 1 2.7 J < 2	< 1 1.6 J < 2	< 1 <1.5 < 2	< 1 <1.5 < 2	< 1 < 5 < 5	< 1.9 J	< <
Bromodichloromethane	50 50 5	< < <	< < <	< < <	< < <	< < <	< 5 < 1	< 5 < 1	<1 < <5 <1	< 1 < < 5 < 1	< 5 < 1	< 2 < 1	< 2 < 1	<1 2.7 J <2 <1	< 1 1.6 J < 2 < 1	<1 <1.5 <2 <1	<1 <1.5 <2 <1	< 1 < 5 < 5 < 1	< 1.9 J < < < <	< < < < < < < < < < < < < < < < < < <
Bromodichloromethane Dibromochloromethane	50 50 5 5	< < < <	< < < < < < < < < < < < < < < < < < <	< < <	< < < < < < < < < < < < < < < < < < <	< < <	< 5 < 1 < 1	< 5 < 1 < 1	<1 < <5 <1 <1	<1 < <5 <1 <1	< 5 < 1 < 1	< 2 < 1 < 1	< 2 < 1 < 1	<1 2.7 J <2 <1 <1	<1 1.6 J <2 <1 <1	<1 <1.5 <2 <1 <1	<1 <1.5 <2 <1 <1	<1 <5 <5 <1 <1	<pre></pre>	< < < < < < < < < < < < < < < < < < <
Bromodichloromethane Dibromochloromethane Chloromethane	50 50 5 5 NV	<	< < < < < < < < < < < < < < < < < < <	< < < < < < < < < < < < < < < < < < <	< < < < < < < < < < < < < < < < < < <	< < < < < < < < < < < < < < < < < < <	< 5 < 1 < 1 < 1	< 5 < 1 < 1 < 1	<1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1	<1 < <5 <1 <1 <1	< 5 < 1 < 1 < 1	< 2 < 1 < 1 < 1	< 2 < 1 < 1 < 1	<1 2.7 J <2 <1 <1	<1 1.6 J <2 <1 <1	<1 <1.5 <2 <1 <1	<1 <1.5 <2 <1 <1	< 1 < 5 < 5 < 1 < 1	< 1.9 J < < < < < < < < < < < < < < < < < <	< < < < < < < < < < < < < < < < < < <
Bromodichloromethane Dibromochloromethane Chloromethane Chloroform	50 50 5 5 50 NV 7	< < < < < < < < < < < < < < < < < < <	<	< < < < < < < < < < < < < < < < < < <	< < < < < < < < < < < < < < < < < < <	<	< 5 < 1 < 1 < 1 < 1 < 1	< 5 < 1 < 1 < 1 < 1 < 1	<1 < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 < <5 <1 <1 <1	< 5 < 1 < 1 < 1 < 1 < 1	< 2 < 1 < 1 < 1 < 1 < 1	< 2 <1 <1 <1 <1 <1	<1 2.7 J <2 <1 <1 <1	<1 1.6 J <2 <1 <1 <1	<1 <1.5 <2 <1 <1 <1	<1 <1.5 <2 <1 <1 <1	<1 <5 <5 <1 <1 <1 <1	<pre></pre>	<
Bromodichloromethane Dibromochloromethane Chloromethane Chloroform Bromoform	50 50 5 50 NV 7 50	<	<	<	<	< < < < < < < < < < < < < < < < < < <	< 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1	< 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1	<1 < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	< 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1	<pre> &lt; 2   &lt;1   &lt;1   &lt;1   &lt;1   &lt;1   &lt;1   &lt;1</pre>	< 2 <1 <1 <1 <1 <1 <1	<1 2.7 J <2 <1 <1 <1 <1	<1 1.6 J <2 <1 <1 <1 <1	<1 <1.5 <2 <1 <1 <1 <1	<1 <1.5 <2 <1 <1 <1 <1	<1 <5 <5 <1 <1 <1 <1 <1	< 1.9 J < < < < < < < < < < < < < < < < < <	<
Bromodichloromethane Dibromochloromethane Chloromethane Chloroform Bromoform Carbon Disulfide	50 50 5 50 NV 7 50 NV	<td>&lt;</td> <td>&lt;</td> <td>&lt;</td> <td>&lt;</td> <td>&lt; &lt;5 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td> <td>&lt;</td> <td>&lt;1 &lt;1 &lt;&lt; &lt;5 &lt;1 &lt;1</td> <td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td> <td>&lt; &lt;5 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td> <td><pre></pre></td> <td>&lt; 2 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td> <td>&lt;1 2.7 J &lt;2 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td> <td>&lt;1 1.6 J &lt;2 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td> <td>&lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td> <td>&lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td> <td>&lt;1 &lt;5 &lt;5 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td> <td><pre></pre></td> <td></td>	<	<	<	<	< <5 <1 <1 <1 <1 <1 <1 <1 <1 <1	<	<1 <1 << <5 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	< <5 <1 <1 <1 <1 <1 <1 <1 <1 <1	<pre></pre>	< 2 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 2.7 J <2 <1 <1 <1 <1 <1	<1 1.6 J <2 <1 <1 <1 <1 <1	<1 <1.5 <2 <1 <1 <1 <1 <1	<1 <1.5 <2 <1 <1 <1 <1 <1	<1 <5 <5 <1 <1 <1 <1 <1 <1	<pre></pre>	
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Bromodichloromethane Dibromochloromethane Chloromethane Chloroform Bromoform Carbon Disulfide Iodomethane Vinyl Chloride 1,1-Dichloroethene 1,1-Dichloroethane trans-1, 2-Dichloroethene cis-1,2-Dichloroethene	50 50 5 50 NV 7 50 NV NV NV 2 5* 5	<pre></pre>	<pre></pre>	<td><td><td><pre></pre></td><td><pre></pre></td><td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td><td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td><td><pre> &lt;</pre></td><td><pre></pre></td><td><pre></pre></td><td>&lt;1 2.7 J &lt; 2 &lt; 1 &lt; 1</td><td>&lt;1 1.6 J &lt;2 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td><td>&lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td><td>&lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td><td>&lt;1 &lt;5 &lt;5 &lt;1 &lt;1</td><td><pre></pre></td><td>&lt; <!--</td--></td></td></td>	<td><td><pre></pre></td><td><pre></pre></td><td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td><td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td><td><pre> &lt;</pre></td><td><pre></pre></td><td><pre></pre></td><td>&lt;1 2.7 J &lt; 2 &lt; 1 &lt; 1</td><td>&lt;1 1.6 J &lt;2 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td><td>&lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td><td>&lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td><td>&lt;1 &lt;5 &lt;5 &lt;1 &lt;1</td><td><pre></pre></td><td>&lt; <!--</td--></td></td>	<td><pre></pre></td> <td><pre></pre></td> <td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td> <td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td> <td><pre> &lt;</pre></td> <td><pre></pre></td> <td><pre></pre></td> <td>&lt;1 2.7 J &lt; 2 &lt; 1 &lt; 1</td> <td>&lt;1 1.6 J &lt;2 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td> <td>&lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td> <td>&lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td> <td>&lt;1 &lt;5 &lt;5 &lt;1 &lt;1</td> <td><pre></pre></td> <td>&lt; <!--</td--></td>	<pre></pre>	<pre></pre>	<1 < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<pre> &lt;</pre>	<pre></pre>	<pre></pre>	<1 2.7 J < 2 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1	<1 1.6 J <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1.5 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1.5 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <5 <5 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<pre></pre>	< </td
Bromodichloromethane Dibromochloromethane Chloromethane Chloroform Bromoform Carbon Disulfide Iodomethane Vinyl Chloride 1,1-Dichloroethene 1,1-Dichloroethane trans-1, 2-Dichloroethene cis-1,2-Dichloroethene 1,1,1-Trichloroethane	50 50 5 50 NV 7 50 NV NV 2 5* 5 5	<pre></pre>	<pre></pre>	<pre></pre>	<pre></pre>	<td><pre> &lt;     &lt;5     &lt;1     &lt;1</pre></td> <td><pre></pre></td> <td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td> <td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td> <td><pre> &lt;     &lt;5     &lt;1     &lt;1</pre></td> <td><pre></pre></td> <td><pre> &lt;</pre></td> <td>&lt;1 2.7 J &lt;2 &lt;1 &lt;1</td> <td>&lt;1 1.6 J &lt;2 &lt;1 &lt;1</td> <td>&lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td> <td>&lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td> <td>&lt;1 &lt;5 &lt;5 &lt;1 &lt;1</td> <td><pre></pre></td> <td>&lt; &lt; &lt;</td>	<pre> &lt;     &lt;5     &lt;1     &lt;1</pre>	<pre></pre>	<1 < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<pre> &lt;     &lt;5     &lt;1     &lt;1</pre>	<pre></pre>	<pre> &lt;</pre>	<1 2.7 J <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 1.6 J <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1.5 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1.5 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <5 <5 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<pre></pre>	< < <
Bromodichloromethane Dibromochloromethane Chloromethane Chloroform Bromoform Carbon Disulfide Iodomethane Vinyl Chloride 1,1-Dichloroethene 1,1-Dichloroethane trans-1, 2-Dichloroethene cis-1,2-Dichloroethene 1,1,1-Trichloroethane Trichloroethene	50 50 5 50 NV 7 50 NV NV 2 5* 5 5 5	<pre> &lt;</pre>	<pre></pre>	<pre></pre>	<pre></pre>	<td><pre> &lt;     &lt;5     &lt;1     &lt;1</pre></td> <td><pre></pre></td> <td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td> <td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td> <td><pre> &lt;     &lt;5     &lt;1     &lt;1</pre></td> <td><pre> &lt;</pre></td> <td><pre> &lt;</pre></td> <td>&lt;1 2.7 J &lt;2 &lt;1 &lt;1</td> <td>&lt;1 1.6 J &lt; 2 &lt; 1 &lt; 1</td> <td>&lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td> <td>&lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td> <td>&lt;1 &lt;5 &lt;5 &lt;1 &lt;1</td> <td><pre></pre></td> <td>&lt; &lt; &lt;</td>	<pre> &lt;     &lt;5     &lt;1     &lt;1</pre>	<pre></pre>	<1 < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<pre> &lt;     &lt;5     &lt;1     &lt;1</pre>	<pre> &lt;</pre>	<pre> &lt;</pre>	<1 2.7 J <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 1.6 J < 2 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1	<1 <1.5 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1.5 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <5 <5 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<pre></pre>	< < <

### Notes:

- 1. Compounds detected in one or more samples are presented on this table.
- 2. Analytical testing completed by Alpha Analytical.
- 3. NYSDEC Class GA criteria obtained from Division of Water Technical and Operational Guidance Series (TOGS 1.1.1), dated October 1993, revised June 1998, January 1999 errata sheet,

and April 2000 addendum. \* Guidance value (not a standard) for 1,1-Dichloroethene = 0.07 ug/L as per the January 1999 update.

- 4. ug/L = part per billion (ppb).
- 5. < indicates compound was not detected; < 1 indicates compound was not detected above its respective reporting limit.
- 6. Shading indicates exceedance of Class GA Criteria.
- 7. NT = not tested.
- 8. NV = no value.
- 9. Results shown for IRM-1 for the September 2021 sampling event are the higher results from it or its respective duplicate.
- 10. Lab qualifiers: CH = continuing calibration outside of lab acceptance limits; results may be biased high. J = estimated concentration.
- L2 = analyte recovery in the control sample was below quality control limits; results may be biased low. Qualifiers for detected compounds only shown.

### September 2021 Groundwater Analytical Testing Results Summary Former Signore Facility 55-57 Jefferson Street

Ellicottville, New York

											EW-4.5									
Parameter	Class GA Criteria	4/23/09	10/22/09	6/3/10	4/13/11	10/14/11	5/10/12	11/1/12	6/26/13	10/16/13	6/9/14	10/14/14	6/2/15	10/21/15	6/14/16	10/24/16	7/11/17	6/19/18	6/11/19	9/15/21
Volatile Organic Compounds	s - EPA Method 8260																			
Methylene chloride	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<	<
Acetone	50	<	<	<	<	<	<	<	<	<	<	<	<	<	4.1 J	<1.5	<1.5	< 5	3 J	<
2-Butanone	50	<	<	<	<	<	< 5	< 5	< 5	< 5	< 5	< 2	< 2	< 2	< 2	< 2	< 2	< 5	<	٧
Bromodichloromethane	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<	<
Dibromochloromethane	50	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<	<
Chloromethane	NV	<	<	<	<	<	< 1	< 1	2.5	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	0.73 J	<
Chloroform	7	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<	<
Benzene	1	<	<	<	<	<	<	<	<	<	<	<	<	<	<	<	<	<	<	<
Bromoform	50	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<	<
Carbon disulfide	NV	<	<	<	.63J	<	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1	<1	<	<
Iodomethane	NV	<	<	<	<	<	< 1	< 1	0.83J	< 1	< 1	NT	NT	NT	NT	NT	NT	NT	<	<
Vinyl Chloride	2	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<	<b>~</b>
1,1-Dichloroethene	5*	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<	<
1,1-Dichloroethane	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<	<
trans-1, 2-Dichloroethene	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<	<
cis-1,2-Dichloroethene	5	<	0.72	<	1.2	.51J	0.61J	< 1	0.76J	< 1	< 1	< 1	< 1	< 1	< 1	0.81 J	<1	<1	<	<
1,1,1-Trichloroethane	5	2.5	1.3	0.97	1.9	1.3	1.2	1.2	1.1	< 1	< 1	0.76 J	0.77 J	<1	<1	<1	<1	<1	<	<
Trichloroethene	5	8.0	7.9	5.5	10	6.9	7.6	7.0	6.8	5.8	5.0	5.4	5.4	3.9	4.6	4.6	1.6	1.1	5	2.4
Tetrachloroethene	5	2.0	1.7	1.1	2.5	1.5	1.5	1.6	1.6	1.4	1.7	1.5	1.7	1.2	1.3	1.6	0.76	< 1	1.50	0.78
Naphthalene	10	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	NT	NT	NT	NT	NT	NT	< 1	<	<
Total VOCs		12.5	11.6	7.6	16.2	10.2	10.9	9.8	13.6	7.2	6.7	7.66	7.86	5.10	10.00	7.01	2.36	1.10	10.23	3.18
_			1	ī	T	1	T	1	1		IRM-1	1		ı	1		1	1	ı	
Parameter	Class GA Criteria	4/23/09	10/22/09	6/3/10	4/13/11	10/14/11	5/10/12	11/1/12	6/26/13	40/40/40	0/0/4	40/44/44	0/0/45	40/04/45						
Volatile Organic Compounds					.,,			1 1/ 1/ 12	0/20/13	10/16/13	6/6/14	10/14/14	6/2/15	10/21/15	6/14/16	10/24/16	7/11/17	6/19/18	6/12/19	9/16/21
Mothylono oblazista	s - EPA Method 8260			3, 3, 1 3	1,, 10, 11			11/1/12	0/20/13	10/16/13	6/6/14	10/14/14	6/2/15	10/21/15	6/14/16	10/24/16	7/11/17	6/19/18	6/12/19	9/16/21
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Acetone 2-Butanone Bromodichloromethane Dibromochloromethane Chloroform Bromoform Carbon Disulfide Iodomethane Vinyl Chloride 1,1-Dichloroethene 1,1-Dichloroethene trans-1, 2-Dichloroethene 1,1,1-Trichloroethane Trichloroethene Trichloroethene	5 50 50 5 50 NV 7 50 NV NV 2 5* 5 5 5	<pre></pre>	<pre></pre>	<td><td><pre></pre></td><td><pre></pre></td><td>&lt;1 &lt;1 &lt; 1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td><td>&lt;1 &lt;1 &lt; 5 &lt;1 1.4 &lt;1 &lt;1</td><td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td><td>&lt;1 &lt;1 &lt; 1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td><td>&lt;1 &lt; 1 &lt; 2 &lt; 1 &lt;</td><td>&lt;1 &lt; 1 &lt; 2 &lt; 1 &lt;</td><td>&lt;1 &lt;1 &lt;</td><td>&lt;1 3.0 J &lt;2 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td><td>&lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td><td>&lt;1 &lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td><td>&lt;1 &lt;5 &lt;5 &lt;1 &lt;1</td><td><pre></pre></td><td></td></td>	<td><pre></pre></td> <td><pre></pre></td> <td>&lt;1 &lt;1 &lt; 1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td> <td>&lt;1 &lt;1 &lt; 5 &lt;1 1.4 &lt;1 &lt;1</td> <td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td> <td>&lt;1 &lt;1 &lt; 1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td> <td>&lt;1 &lt; 1 &lt; 2 &lt; 1 &lt;</td> <td>&lt;1 &lt; 1 &lt; 2 &lt; 1 &lt;</td> <td>&lt;1 &lt;1 &lt;</td> <td>&lt;1 3.0 J &lt;2 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td> <td>&lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td> <td>&lt;1 &lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td> <td>&lt;1 &lt;5 &lt;5 &lt;1 &lt;1</td> <td><pre></pre></td> <td></td>	<pre></pre>	<pre></pre>	<1 <1 < 1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1 < 5 <1 1.4 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 <1 < 1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 < 1 < 2 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 < 1 < 2 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <	<1 3.0 J <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1.5 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1 <1.5 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <5 <5 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<pre></pre>	

### Notes

- 1. Compounds detected in one or more samples are presented on this table.
- 2. Analytical testing completed by Alpha Analytical.
- 3. NYSDEC Class GA criteria obtained from Division of Water Technical and Operational Guidance Series (TOGS 1.1.1), dated October 1993, revised June 1998, January 1999 errata sheet,
- and April 2000 addendum. \* Guidance value (not a standard) for 1,1-Dichloroethene = 0.07 ug/L as per the January 1999 update.
- 4. ug/L = part per billion (ppb).
- 5. < indicates compound was not detected; < 1 indicates compound was not detected above its respective reporting limit.
- 6. Shading indicates exceedance of Class GA Criteria.
- 7. NT = not tested.
- 8. NV = no value.
- 9. Results shown for IRM-1 for the September 2021 sampling event are the higher results from it or its respective duplicate.
- 10. Lab qualifiers: CH = continuing calibration outside of lab acceptance limits; results may be biased high. J = estimated concentration.
- L2 = analyte recovery in the control sample was below quality control limits; results may be biased low. Qualifiers for detected compounds only shown.

# September 2021 Groundwater Analytical Testing Results Summary Former Signore Facility 55-57 Jefferson Street

Ellicottville, New York

											MW-1I									
Parameter	Class GA Criteria	4/23/09	10/22/09	6/2/10	4/14/11	10/14/11	5/9/12	10/5/12	6/25/13	10/15/13	6/9/14	10/15/14	6/2/15	10/22/15	6/14/16	10/25/16	7/11/17	6/20/18	6/13/19	9/15/21
<b>Volatile Organic Compound</b>	ls - EPA Method 8260		ı			<u> </u>							<u> </u>		<u> </u>	<u> </u>	<u> </u>	<u> </u>		
Methylene chloride	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
Acetone	50	<	<	<	<	<	<	<	<	<	<	<	<	<.1.5	<.1.5	<.1.5	1.9 J	<5.0	4.5 J	<
2-Butanone	50	<	<	<	<	<	< 5	< 5	< 5	< 5	< 5	< 2	< 2	< 2	< 2	< 2	< 2	<5.0	<	<
Bromodichloromethane	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	1.4	<
Dibromochloromethane	50	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	0.26 J	<
Chloromethane	NV	<	<	0.62	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	0.85 J	<
Chloroform	7	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
Benzene	1	<	<	<	<	<	<	<	<	<b>V</b>	٧	<	<	<	<	<	<	<	<	<
Bromoform	50	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
Carbon disulfide	NV	<	<	<	<	1.1	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1	<1	<	<
Iodomethane	NV	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	NT	NT	NT	NT	NT	NT	NT	<	<
Vinyl Chloride	2	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	0.53 J	<1	<1	<	<
1,1-Dichloroethene	5*	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
1,1-Dichloroethane	5	4.7	4.7	3.5	3.4	3.8	2.8	2.6	2.0	2.1	1.6	2.3 J	1.9 J	2.5	1.7 J	1.2 J	<1	1.1 L2	<	1.6 J
trans-1, 2-Dichloroethene	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
cis-1,2-Dichloroethene	5	4.2	5.7	2.2	2.5	2.2	1.2	3.1	2.9	1.8	< 1	1.8 J	0.87 J	0.80 J	1.6 J	7.1	<1	3.3	<	0.70 J
1,1,1-Trichloroethane	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	< 1	<1	<	<
Trichloroethene	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	2.8	2	<1	3	11	<1	15	<	<
Tetrachloroethene	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	2.4	1.3	<1	1.9	7.1	<1	11.6 CH	<	<
Naphthalene	10	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	NT	NT	NT	NT	NT	NT	<1	<	<
Total VOCs		8.9	10.4	6.3	5.9	7.1	4.0	5.7	4.9	3.9	1.6	9.0	6.1	3.3	8.2	26.9	1.9	31.0	7.01	2.30
											IDM OL									
Parameter	Class GA Criteria		1	ı			,				IRM-2I									
		4/23/09	10/22/09	6/3/10	4/13/11	10/14/11	5/10/12	11/1/12	6/26/13	10/16/13	6/6/14	10/14/14	6/2/15	10/21/15	6/14/16	10/24/16	7/11/17	6/19/18	6/12/19	9/16/21
Volatile Organic Compound	  s - FPA Method 8260		10/22/09	6/3/10	4/13/11	10/14/11	5/10/12	11/1/12	6/26/13		6/6/14	10/14/14	6/2/15	10/21/15	6/14/16	10/24/16	7/11/17	6/19/18	6/12/19	9/16/21
Volatile Organic Compound										10/16/13										
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Methylene chloride Acetone	5 50	< <	< <	< <	< <	< <	<1 <	< 1	< 1	10/16/13	< 1	< 1	<1 <	< 1 <1.5	< 1 2.9 J	< 1 <1.5	< 1 <1.5	< 1 < 5	< 2.7 J	< <
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### Notes:

- 1. Compounds detected in one or more samples are presented on this table.
- 2. Analytical testing completed by Alpha Analytical.
- 3. NYSDEC Class GA criteria obtained from Division of Water Technical and Operational Guidance Series (TOGS 1.1.1), dated October 1993, revised June 1998, January 1999 errata sheet,

and April 2000 addendum. \* Guidance value (not a standard) for 1,1-Dichloroethene = 0.07 ug/L as per the January 1999 update.

- 4. ug/L = part per billion (ppb).
- 5. < indicates compound was not detected; < 1 indicates compound was not detected above its respective reporting limit.
- 6. Shading indicates exceedance of Class GA Criteria.
- 7. NT = not tested.
- 8. NV = no value.
- 9. Results shown for IRM-1 for the September 2021 sampling event are the higher results from it or its respective duplicate.
- 10. Lab qualifiers: CH = continuing calibration outside of lab acceptance limits; results may be biased high. J = estimated concentration.

  L2 = analyte recovery in the control sample was below quality control limits; results may be biased low. Qualifiers for detected compounds only shown.

# September 2021 Groundwater Analytical Testing Results Summary Former Signore Facility 55-57 Jefferson Street

Ellicottville, New York

											MW-2I									
Parameter	Class GA Criteria	4/23/09	10/22/09	6/3/10	4/13/11	10/13/11	5/9/12	10/31/12	6/25/13	10/15/13	6/6/14	10/14/14	6/3/15	10/22/15	6/15/16	10/24/16	7/11/17	6/20/18	6/13/19	9/15/21
Volatile Organic Compound	ls - EPA Method 8260																			
Methylene chloride	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1	<	<
Acetone	50	<	<	<	<	<	<	<	<	<	<	<	<b>'</b>	<1.5	<1.5	<1.5	<1.5	<5.0	2.1 J	<
2-Butanone	50	<	<	<	<	<	< 5	< 5	< 5	< 5	< 5	< 2	<2	<2	<2	<2	<2	<5.0	<	<
Bromodichloromethane	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1	<	<
Dibromochloromethane	50	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1	<	<
Chloromethane	NV	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1	<	<
Chloroform	7	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1	<	<
Benzene	1	<	<	<	<	<	<	<	<	<	<	<	<	<	<	<	<	<	<	<
Bromoform	50	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1	<	<
Carbon disulfide	NV	<	<	12.0	0.90J	1.3	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1	<1	<	<
Iodomethane	NV	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	NT	NT	NT	NT	NT	NT	NT	<	<
Vinyl Chloride	2	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1	<	<
1,1-Dichloroethene	5*	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1	<	<
1,1-Dichloroethane	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1	<	<
trans-1, 2-Dichloroethene	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1	<	<
cis-1,2-Dichloroethene	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1	<	<
1,1,1-Trichloroethane	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1	<	<
Trichloroethene	5	<	<	<	<	<	0.83J	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1	<	<
Tetrachloroethene	5	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	<1	<1	<1	<1	<1	<1	<1	<	<
Naphthalene	10	<	<	<	<	<	< 1	< 1	< 1	< 1	< 1	NT	NT	NT	NT	NT	NT	<1	<	<
Total VOCs				12.0	0.9	1.3	0.83												2.10	
-	1									TO	A/NI \A/EI I									
Parameter	Class CA Critoria			I				I	l	TO\	WN WELI	L		I		ı	-	I		
Parameter	Class GA Criteria	4/23/09	10/22/09	6/2/10	4/13/11	10/14/11	5/10/12	11/1/12	6/26/13	TO\ 10/16/13	WN WELI 6/9/14	10/14/14	6/2/15	10/22/15	6/14/16	10/24/16	7/12/17	6/19/18	6/11/19	9/16/21
Parameter  Volatile Organic Compound			10/22/09	6/2/10	4/13/11	10/14/11	5/10/12	11/1/12	6/26/13				6/2/15	10/22/15	6/14/16	10/24/16	7/12/17	6/19/18	6/11/19	9/16/21
	ds - EPA Method 8260		10/22/09 NT	6/2/10	4/13/11	10/14/11	5/10/12	11/1/12	6/26/13				6/2/15	10/22/15	<1	10/24/16	7/12/17	6/19/18	6/11/19	9/16/21
Volatile Organic Compound Methylene chloride Acetone	ds - EPA Method 8260 5 50	,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,					<1 <	< 1	< 1	10/16/13	6/9/14	10/14/14	<1	<1 <1.5	<1 2.4 J	<1 <1.5	<1 <1.5	< 1 < 5	6/11/19 2.6 J	<
Volatile Organic Compound Methylene chloride Acetone 2-Butanone	5 5 50 50	<	NT	<	< < <	< < <	< 1 < < 5	< 1 < < 5	< 1 < < 5	10/16/13	6/9/14 < 1 < 5	10/14/14	<1 < <2	<1 <1.5 <2	<1 2.4 J <2	<1 <1.5 <2	<1	< 1	2.6 J	< <
Volatile Organic Compound Methylene chloride Acetone 2-Butanone Bromodichloromethane	5 5 50 50 50	< <	NT <	< <	< < < < < < .53J	< < < < < < 1.4	<1 < <5 0.67J	< 1	< 1 < < 5 < 1	10/16/13	6/9/14 < 1 < 5 < 1	10/14/14 < 1 < 2 < 1	<1 < <2 0.52	<1 <1.5 <2 0.27 J	<1 2.4 J <2 0.45 J	<1 <1.5 <2 0.53	<1 <1.5 <2 <1	< 1 < 5 < 5 < 1	2.6 J < 0.5	< < < 0.36 J
Volatile Organic Compound Methylene chloride Acetone 2-Butanone	5 5 50 50 50 50 50	< < <	NT	< < < < < < < < < < < < < < < < < < <	< < <	< < <	< 1 < < 5	< 1 < < 5 0.96J < 1	< 1 < < 5 < 1 < 1	10/16/13 < 1 < 5	6/9/14 < 1 < 5	10/14/14	<1 < <2	<1 <1.5 <2	<1 2.4 J <2	<1 <1.5 <2	<1 <1.5 <2 <1 <1	< 1 < 5 < 5	2.6 J	< <
Volatile Organic Compound Methylene chloride Acetone 2-Butanone Bromodichloromethane Dibromochloromethane Chloromethane	5 50 50 50 50 50 50 NV	< < < < < < < < < < < < < < < < < < <	NT   <   <   <   <   <   <   <   <   <	< < < < < < < < < < < < < < < < < < <	< < < < < < .53J	<ul><li>&lt;</li><li>&lt;</li><li>&lt;</li><li>1.4</li><li>1.7</li></ul>	<1 < 1 < 5 0.67J 1.2 < 1	< 1 < < 5 0.96J < 1 < 1	< 1 < 5 < 1 < 1 1.3	10/16/13  <1 <1 <5 <1 <1 <1 <1	6/9/14 <1 <5 <1 <1 <1	10/14/14 < 1 < 2 < 1 < 1 < 1	<1 < 2 0.52 0.99 <1	<1 <1.5 <2 0.27 J 0.54 <1	<1 2.4 J <2 0.45 J 3 <1	<1 <1.5 <2 0.53 0.97 <1	<1 <1.5 <2 <1 <1 <1 <1	< 1 < 5 < 5 < 1 1.3 < 1	2.6 J < 0.5	<ul><li>&lt;</li><li>&lt;</li><li>0.36 J</li><li>0.66</li><li>&lt;</li></ul>
Volatile Organic Compound Methylene chloride Acetone 2-Butanone Bromodichloromethane Dibromochloromethane Chloroform	5 50 50 50 50 50 NV 7	< < < < < < < < < < < < < < < < < < <	NT	<	<ul><li>&lt;</li><li>&lt;</li><li>&lt;</li><li>.53J</li><li>1.2</li><li>&lt;</li><li>&lt;</li></ul>	<ul><li></li><li></li><li></li><li>1.4</li><li>1.7</li><li></li><li>1.1</li></ul>	<1 < 1 < 5 0.67J 1.2 < 1 < 1	<1 < 1 < 5 0.96J < 1 < 1 0.82J	<1 <1 <5 <1 <1 <1 3 <1	10/16/13  <1 <1 <5 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	6/9/14  < 1     < 5     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1	10/14/14   < 1   < 2   < 1   < 1   < 1   < 1	<1 < 2 0.52 0.99 <1 <1	<1 <1.5 <2 0.27 J 0.54 <1 <1	<1 2.4 J <2 0.45 J 3 <1 <1	<1 <1.5 <2 0.53 0.97 <1 <1	<1 <1 <1.5 <2 <1 <1 <1 <1 <1 <1 <1	<1 <5 <5 <1 1.3 <1 <1	2.6 J < 0.5 0.73	<ul><li>&lt;</li><li>&lt;</li><li>0.36 J</li><li>0.66</li><li>&lt;</li><li>&lt;</li></ul>
Volatile Organic Compound Methylene chloride Acetone 2-Butanone Bromodichloromethane Dibromochloromethane Chloroform Bromoform	5 50 50 50 50 50 NV 7 50	< < < < < < < < < < < < < < < < < < <	NT	< < < < < < < < < < < < < < < < < < <	<ul><li>&lt;</li><li>&lt;</li><li>&lt;</li><li>.53J</li><li>1.2</li><li>&lt;</li></ul>	<ul><li>&lt;</li><li>&lt;</li><li>&lt;</li><li>1.4</li><li>1.7</li></ul>	<1 < 1 < 5 0.67J 1.2 < 1	<1 < 1 < 5 0.96J < 1 < 1 0.82J 1.6	<1 < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	10/16/13  <1 <1 <5 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	6/9/14  < 1     < 5     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1	10/14/14   <1   < 2   < 1   < 1   < 1   < 1   < 1	<1 < < 0.52	<1 <1.5 <2 0.27 J 0.54 <1 <1 <1	<1 2.4 J <2 0.45 J 3 <1	<1 <1.5 <2 0.53 0.97 <1	<1 <1.5 <2 <1 <1 <1 <1 <1	< 1 < 5 < 5 < 1 1.3 < 1	2.6 J < 0.5 0.73	<ul><li>&lt;</li><li>&lt;</li><li>0.36 J</li><li>0.66</li><li>&lt;</li></ul>
Volatile Organic Compound Methylene chloride Acetone 2-Butanone Bromodichloromethane Dibromochloromethane Chloromethane Chloroform Bromoform Carbon Disulfide	5 50 50 50 50 50 NV 7 50 NV	<	NT	<	<ul><li>&lt;</li><li>&lt;</li><li>&lt;</li><li>.53J</li><li>1.2</li><li>&lt;</li><li>&lt;</li></ul>	<ul><li></li><li></li><li></li><li>1.4</li><li>1.7</li><li></li><li>1.1</li></ul>	<1 < 1 < 5 0.67J 1.2 < 1 < 1 < 1 0.88J < 1	<1 < 1 < 5 0.96J < 1 < 1 0.82J 1.6 < 1	<1 < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	10/16/13  <1 <1 <5 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	6/9/14  < 1     < 5     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1	10/14/14   <1   <2   <1   <1   <1   <1   <1   <1	<1 < 2 0.52 0.99 <1 <1 1.2 J <1	<1 <1 <1.5 <2 0.27 J 0.54 <1 <1 <1 <1 <1 <1	<1 2.4 J <2 0.45 J 3 <1 <1 1.3 J	<1 <1.5 <2 0.53 0.97 <1 <1 1.3 J <1	<1 <1 <1.5 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <5 <5 <1 1.3 <1 <1 <1 <1	2.6 J < 0.5 0.73	<ul> <li></li> <li>0.36 J</li> <li>0.66</li> <li></li> <li></li> <li></li> <li></li> </ul>
Volatile Organic Compound Methylene chloride Acetone 2-Butanone Bromodichloromethane Dibromochloromethane Chloroform Bromoform Carbon Disulfide Iodomethane	5 50 50 NV 7 50 NV NV	<	NT	<	<ul><li>&lt;</li><li>&lt;</li><li>&lt;</li><li>.53J</li><li>1.2</li><li>&lt;</li><li>&lt;</li><li>1.7</li></ul>	<	<1 < 1 < 5 0.67J 1.2 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 < 1 < 5 0.96J < 1 < 1 0.82J 1.6 < 1 < 1 < 1	<1 <1 <5 <1 1.3 <1 <1 <1 <1 <1 <1 <1 <1 <1	10/16/13   <1   <   <5   <1   <1   <1   <1   <1   <1   <1	6/9/14  < 1   < 5   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1	10/14/14   <1   <2   <1   <1   <1   <1   <1   <1   <1   <1	<1 < 2 0.52 0.99 <1 <1 1.2 J <1 NT	<1 <1.5 <2 0.27 J 0.54 <1 <1 <1 <1 NT	<1 2.4 J <2 0.45 J 3 <1 <1 <1 1.3 J <1 NT	<1 <1.5 <2 0.53 0.97 <1 <1 1.3 J <1 NT	<1 <1.5 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <5 <5 <1 1.3 <1 <1 <1 <1 NT	2.6 J < 0.5 0.73 < <	<ul> <li></li> <li></li> <li>0.36 J</li> <li>0.66</li> <li></li> <li></li> <li></li> <li></li> </ul>
Volatile Organic Compound Methylene chloride Acetone 2-Butanone Bromodichloromethane Dibromochloromethane Chloroform Bromoform Carbon Disulfide Iodomethane Vinyl Chloride	5 50 50 NV 7 50 NV NV 2	<td>  NT  </td> <td>&lt;</td> <td><ul> <li>&lt;</li> <li>&lt;</li> <li>.53J</li> <li>1.2</li> <li>&lt;</li> <li>&lt;</li> <li>1.7</li> </ul></td> <td><ul> <li>&lt;</li> <li>&lt;</li> <li>1.4</li> <li>1.7</li> <li>&lt;</li> <li>1.1</li> <li>1.4</li> <li>&lt;</li> </ul></td> <td>&lt;1 &lt; 1 &lt; 5 0.67J 1.2 &lt; 1 &lt;</td> <td>&lt;1 &lt; 1 &lt; 5 0.96J &lt; 1 &lt; 1 0.82J 1.6 &lt; 1 &lt; 1 &lt; 1 &lt; 1 &lt; 1 &lt; 1 &lt; 1</td> <td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td> <td>10/16/13  &lt;1 &lt;1 &lt;5 &lt;1 &lt;1</td> <td>6/9/14  &lt; 1   &lt; 5   &lt; 1   &lt; 1</td> <td>  10/14/14   &lt; 1</td> <td>&lt;1 &lt; 2 0.52 0.99 &lt;1 &lt;1 1.2 J &lt;1 NT &lt;1</td> <td>&lt;1 &lt;1.5 &lt;2 0.27 J 0.54 &lt;1 &lt;1 &lt;1 &lt;1 NT &lt;1</td> <td>&lt;1 2.4 J &lt;2 0.45 J 3 &lt;1 &lt;1 1.3 J &lt;1 NT &lt;1</td> <td>&lt;1 &lt;1.5 &lt;2 0.53 0.97 &lt;1 &lt;1 &lt;1 NT &lt;1</td> <td>&lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 &lt;1 NT &lt;1</td> <td>&lt;1 &lt;5 &lt;5 &lt;1 1.3 &lt;1 &lt;1 &lt;1 &lt;1</td> <td>2.6 J &lt; 0.5 0.73 &lt; &lt; &lt;</td> <td><ul> <li></li> <li>0.36 J</li> <li>0.66</li> <li></li> <li></li> <li></li> <li></li> </ul></td>	NT	<	<ul> <li>&lt;</li> <li>&lt;</li> <li>.53J</li> <li>1.2</li> <li>&lt;</li> <li>&lt;</li> <li>1.7</li> </ul>	<ul> <li>&lt;</li> <li>&lt;</li> <li>1.4</li> <li>1.7</li> <li>&lt;</li> <li>1.1</li> <li>1.4</li> <li>&lt;</li> </ul>	<1 < 1 < 5 0.67J 1.2 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 < 1 < 5 0.96J < 1 < 1 0.82J 1.6 < 1 < 1 < 1 < 1 < 1 < 1 < 1	<1 < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	10/16/13  <1 <1 <5 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	6/9/14  < 1   < 5   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1	10/14/14   < 1	<1 < 2 0.52 0.99 <1 <1 1.2 J <1 NT <1	<1 <1.5 <2 0.27 J 0.54 <1 <1 <1 <1 NT <1	<1 2.4 J <2 0.45 J 3 <1 <1 1.3 J <1 NT <1	<1 <1.5 <2 0.53 0.97 <1 <1 <1 NT <1	<1 <1.5 <2 <1 <1 <1 <1 <1 <1 <1 NT <1	<1 <5 <5 <1 1.3 <1 <1 <1 <1	2.6 J < 0.5 0.73 < < < < < < < < < < < < < < < < < < <	<ul> <li></li> <li>0.36 J</li> <li>0.66</li> <li></li> <li></li> <li></li> <li></li> </ul>
Volatile Organic Compound Methylene chloride Acetone 2-Butanone Bromodichloromethane Dibromochloromethane Chloroform Bromoform Carbon Disulfide Iodomethane Vinyl Chloride 1,1-Dichloroethene	5 50 50 50 NV 7 50 NV NV NV NV 2 5*	<td>  NT  </td> <td><pre></pre></td> <td><ul> <li></li> <li></li> <li>.53J</li> <li>1.2</li> <li></li> <li></li> <li></li> <li></li> <li></li> <li></li> </ul></td> <td><ul> <li>&lt;</li> <li>&lt;</li> <li>1.4</li> <li>1.7</li> <li></li> <li>1.1</li> <li>1.4</li> <li>&lt;</li> <li>&lt;</li> <li>&lt;</li> </ul></td> <td>&lt;1 &lt; 1 &lt; 5 0.67J 1.2 &lt; 1 &lt;</td> <td>&lt;1 &lt; 1 &lt; 5 0.96J &lt; 1 &lt; 1 0.82J 1.6 &lt; 1 &lt;</td> <td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td> <td>  10/16/13   &lt;1   &lt;   &lt;5   &lt;1   &lt;1   &lt;1   &lt;1   &lt;1   &lt;1   &lt;1   &lt;1</td> <td>6/9/14  &lt; 1   &lt; 5   &lt; 1   &lt; 1</td> <td>  10/14/14   &lt; 1</td> <td>&lt;1 &lt; 2 0.52 0.99 &lt;1 &lt;1 1.2 J &lt;1 NT &lt;1 &lt;1 &lt;1</td> <td>&lt;1 &lt;1.5 &lt;2 0.27 J 0.54 &lt;1 &lt;1 &lt;1 NT &lt;1 &lt;1 &lt;1</td> <td>&lt;1 2.4 J &lt;2 0.45 J 3 &lt;1 &lt;1 &lt;1 1.3 J &lt;1 NT &lt;1 &lt;1 &lt;1</td> <td>&lt;1 &lt;1.5 &lt;2 0.53 0.97 &lt;1 &lt;1 &lt;1 NT &lt;1 &lt;1 &lt;1</td> <td>&lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td> <td>&lt;1 &lt;5 &lt;5 &lt;1 1.3 &lt;1 &lt;1 &lt;1 NT &lt;1 &lt;1 &lt;1</td> <td>2.6 J &lt; 0.5 0.73 &lt; &lt; &lt;</td> <td><ul> <li></li> <li></li> <li>0.36 J</li> <li>0.66</li> <li></li> <li></li> <li></li> <li></li> <li></li> </ul></td>	NT	<pre></pre>	<ul> <li></li> <li></li> <li>.53J</li> <li>1.2</li> <li></li> <li></li> <li></li> <li></li> <li></li> <li></li> </ul>	<ul> <li>&lt;</li> <li>&lt;</li> <li>1.4</li> <li>1.7</li> <li></li> <li>1.1</li> <li>1.4</li> <li>&lt;</li> <li>&lt;</li> <li>&lt;</li> </ul>	<1 < 1 < 5 0.67J 1.2 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 < 1 < 5 0.96J < 1 < 1 0.82J 1.6 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	10/16/13   <1   <   <5   <1   <1   <1   <1   <1   <1   <1   <1	6/9/14  < 1   < 5   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1   < 1	10/14/14   < 1	<1 < 2 0.52 0.99 <1 <1 1.2 J <1 NT <1 <1 <1	<1 <1.5 <2 0.27 J 0.54 <1 <1 <1 NT <1 <1 <1	<1 2.4 J <2 0.45 J 3 <1 <1 <1 1.3 J <1 NT <1 <1 <1	<1 <1.5 <2 0.53 0.97 <1 <1 <1 NT <1 <1 <1	<1 <1.5 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <5 <5 <1 1.3 <1 <1 <1 NT <1 <1 <1	2.6 J < 0.5 0.73 < < < < < < < < < < < < < < < < < < <	<ul> <li></li> <li></li> <li>0.36 J</li> <li>0.66</li> <li></li> <li></li> <li></li> <li></li> <li></li> </ul>
Volatile Organic Compound Methylene chloride Acetone 2-Butanone Bromodichloromethane Dibromochloromethane Chloroform Bromoform Carbon Disulfide Iodomethane Vinyl Chloride 1,1-Dichloroethane 1,1-Dichloroethane	5 50 50 50 NV 7 50 NV NV NV NV 2 5* 5	<td>  NT  </td> <td><pre></pre></td> <td><ul> <li></li> <li></li> <li>.53J</li> <li>1.2</li> <li></li> <li></li> <li></li> <li></li> <li></li> </ul></td> <td><ul> <li>&lt;</li> <li>&lt;</li> <li>1.4</li> <li>1.7</li> <li></li> <li>1.1</li> <li>1.4</li> <li>&lt;</li> <li>&lt;</li> <li>&lt;</li> <li>&lt;</li> </ul></td> <td>&lt;1 &lt; 1 &lt; 5 0.67J 1.2 &lt; 1 &lt;</td> <td>&lt;1 &lt; 1 &lt; 5 0.96J &lt; 1 &lt; 1 0.82J 1.6 &lt; 1 &lt;</td> <td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td> <td>  10/16/13   &lt;1   &lt;   &lt;5   &lt;1   &lt;1   &lt;1   &lt;1   &lt;1   &lt;1   &lt;1   &lt;1</td> <td>6/9/14  &lt; 1     &lt; 5     &lt; 1     &lt; 1</td> <td>  10/14/14   &lt; 1</td> <td>&lt;1 &lt; 2 0.52 0.99 &lt;1 &lt;1 1.2 J &lt;1 NT &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td> <td>&lt;1 &lt;1.5 &lt;2 0.27 J 0.54 &lt;1 &lt;1</td> <td>&lt;1 2.4 J &lt;2 0.45 J 3 &lt;1 &lt;1 &lt;1 Style="background-color: lightblue;"&gt;3 &lt;1 &lt;1</td> <td>&lt;1 &lt;1.5 &lt;2 0.53 0.97 &lt;1 &lt;1 1.3 J &lt;1 NT &lt;1 &lt;1 &lt;1 &lt;1 &lt;1</td> <td>&lt;1 &lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td> <td>&lt;1 &lt;5 &lt;5 &lt;1 1.3 &lt;1 &lt;1</td> <td>2.6 J &lt; 0.5 0.73 &lt; &lt; &lt;</td> <td><ul> <li></li> <li>0.36 J</li> <li>0.66</li> <li></li> <li></li> <li></li> <li></li> <li></li> <li></li> </ul></td>	NT	<pre></pre>	<ul> <li></li> <li></li> <li>.53J</li> <li>1.2</li> <li></li> <li></li> <li></li> <li></li> <li></li> </ul>	<ul> <li>&lt;</li> <li>&lt;</li> <li>1.4</li> <li>1.7</li> <li></li> <li>1.1</li> <li>1.4</li> <li>&lt;</li> <li>&lt;</li> <li>&lt;</li> <li>&lt;</li> </ul>	<1 < 1 < 5 0.67J 1.2 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 < 1 < 5 0.96J < 1 < 1 0.82J 1.6 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	10/16/13   <1   <   <5   <1   <1   <1   <1   <1   <1   <1   <1	6/9/14  < 1     < 5     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1	10/14/14   < 1	<1 < 2 0.52 0.99 <1 <1 1.2 J <1 NT <1 <1 <1 <1 <1	<1 <1.5 <2 0.27 J 0.54 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 2.4 J <2 0.45 J 3 <1 <1 <1 Style="background-color: lightblue;">3 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1.5 <2 0.53 0.97 <1 <1 1.3 J <1 NT <1 <1 <1 <1 <1	<1 <1 <1.5 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <5 <5 <1 1.3 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	2.6 J < 0.5 0.73 < < < < < < < < < < < < < < < < < < <	<ul> <li></li> <li>0.36 J</li> <li>0.66</li> <li></li> <li></li> <li></li> <li></li> <li></li> <li></li> </ul>
Volatile Organic Compound Methylene chloride Acetone 2-Butanone Bromodichloromethane Dibromochloromethane Chloroform Bromoform Carbon Disulfide Iodomethane Vinyl Chloride 1,1-Dichloroethane trans-1, 2-Dichloroethene	5 5 50 50 50 50 50 NV 7 50 NV 7 50 NV NV 2 5* 5	<td>  NT  </td> <td><pre></pre></td> <td><ul> <li></li> <li></li> <li>.53J</li> <li>1.2</li> <li></li> <li></li> <li></li> <li></li> <li></li> <li></li> </ul></td> <td><ul> <li>&lt;</li> <li>&lt;</li> <li>1.4</li> <li>1.7</li> <li></li> <li>1.1</li> <li>1.4</li> <li>&lt;</li> <li>&lt;</li> <li>&lt;</li> </ul></td> <td>&lt;1 &lt; 1 &lt; 5 0.67J 1.2 &lt; 1 &lt;</td> <td>&lt;1 &lt; 1 &lt; 5 0.96J &lt; 1 &lt; 1 0.82J 1.6 &lt; 1 &lt;</td> <td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td> <td>10/16/13  &lt;1 &lt;1</td> <td>6/9/14  &lt; 1 &lt; 5 &lt; 1 &lt; 1</td> <td>10/14/14  &lt; 1</td> <td>&lt;1 &lt; 2 0.52 0.99 &lt; 1 &lt; 1 1.2 J &lt; 1 NT &lt; 1 &lt; 1 &lt; 1 &lt; 1 &lt; 1 &lt; 1</td> <td>&lt;1 &lt;1.5 &lt;2 0.27 J 0.54 &lt;1 &lt;1</td> <td>&lt;1 2.4 J &lt;2 0.45 J 3 &lt;1 1.3 J &lt;1 NT &lt;1 &lt;1</td> <td>&lt;1 &lt;1.5 &lt;2 0.53 0.97 &lt;1 &lt;1</td> <td>&lt;1 &lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td> <td>&lt;1 &lt;5 &lt;5 &lt;1 1.3 &lt;1 &lt;1</td> <td>2.6 J &lt; 0.5 0.73 &lt; &lt; &lt;</td> <td><ul> <li></li> <li></li> <li>0.36 J</li> <li>0.66</li> <li></li> <li></li> <li></li> <li></li> <li></li> </ul></td>	NT	<pre></pre>	<ul> <li></li> <li></li> <li>.53J</li> <li>1.2</li> <li></li> <li></li> <li></li> <li></li> <li></li> <li></li> </ul>	<ul> <li>&lt;</li> <li>&lt;</li> <li>1.4</li> <li>1.7</li> <li></li> <li>1.1</li> <li>1.4</li> <li>&lt;</li> <li>&lt;</li> <li>&lt;</li> </ul>	<1 < 1 < 5 0.67J 1.2 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 < 1 < 5 0.96J < 1 < 1 0.82J 1.6 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	10/16/13  <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	6/9/14  < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1	10/14/14  < 1	<1 < 2 0.52 0.99 < 1 < 1 1.2 J < 1 NT < 1 < 1 < 1 < 1 < 1 < 1	<1 <1.5 <2 0.27 J 0.54 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 2.4 J <2 0.45 J 3 <1 1.3 J <1 NT <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1.5 <2 0.53 0.97 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1 <1.5 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <5 <5 <1 1.3 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	2.6 J < 0.5 0.73 < < < < < < < < < < < < < < < < < < <	<ul> <li></li> <li></li> <li>0.36 J</li> <li>0.66</li> <li></li> <li></li> <li></li> <li></li> <li></li> </ul>
Volatile Organic Compound Methylene chloride Acetone 2-Butanone Bromodichloromethane Dibromochloromethane Chloroform Bromoform Carbon Disulfide Iodomethane Vinyl Chloride 1,1-Dichloroethene 1,1-Dichloroethane trans-1, 2-Dichloroethene cis-1,2-Dichloroethene	5 50 50 50 50 NV 7 50 NV NV 2 5* 5 5 5 5 5 5 5 5 5 5 5 5 5 5 5 5 5	<td>  NT   &lt;   &lt;   &lt;   &lt;   &lt;   &lt;   &lt;   &lt;   &lt;  </td> <td><pre></pre></td> <td><ul> <li></li> <li></li> <li>.53J</li> <li>1.2</li> <li></li> <li></li></ul></td> <td><pre></pre></td> <td>&lt;1 &lt; 1 &lt; 5 0.67J 1.2 &lt; 1 &lt;</td> <td>&lt;1 &lt; 1 &lt; 5 0.96J &lt; 1 &lt; 1 0.82J 1.6 &lt; 1 &lt;</td> <td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td> <td>10/16/13  &lt;1 &lt;1</td> <td>6/9/14  &lt; 1 &lt; 5 &lt; 1 &lt; 1</td> <td>10/14/14  &lt; 1</td> <td>&lt;1 &lt; 2 0.52 0.99 &lt;1 &lt;1 1.2 J &lt;1 NT &lt;1 &lt;1</td> <td>&lt;1 &lt;1.5 &lt;2 0.27 J 0.54 &lt;1 &lt;1</td> <td>&lt;1 2.4 J</td> <td>&lt;1 &lt;1.5 &lt;2 0.53 0.97 &lt;1 &lt;1 1.3 J &lt;1 NT &lt;1 &lt;1</td> <td>&lt;1 &lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td> <td>&lt;1 &lt;5 &lt;5 &lt;1 1.3 &lt;1 &lt;1</td> <td>2.6 J &lt; 0.5 0.73 &lt; &lt; &lt;</td> <td><ul> <li></li> <li>0.36 J</li> <li>0.66</li> <li></li> <li></li></ul></td>	NT   <   <   <   <   <   <   <   <   <	<pre></pre>	<ul> <li></li> <li></li> <li>.53J</li> <li>1.2</li> <li></li> <li></li></ul>	<pre></pre>	<1 < 1 < 5 0.67J 1.2 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 < 1 < 5 0.96J < 1 < 1 0.82J 1.6 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	10/16/13  <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	6/9/14  < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1	10/14/14  < 1	<1 < 2 0.52 0.99 <1 <1 1.2 J <1 NT <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1.5 <2 0.27 J 0.54 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 2.4 J	<1 <1.5 <2 0.53 0.97 <1 <1 1.3 J <1 NT <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1 <1.5 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <5 <5 <1 1.3 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	2.6 J < 0.5 0.73 < < < < < < < < < < < < < < < < < < <	<ul> <li></li> <li>0.36 J</li> <li>0.66</li> <li></li> <li></li></ul>
Volatile Organic Compound Methylene chloride Acetone 2-Butanone Bromodichloromethane Dibromochloromethane Chloroform Bromoform Carbon Disulfide Iodomethane Vinyl Chloride 1,1-Dichloroethene 1,1-Dichloroethane trans-1, 2-Dichloroethene cis-1,2-Dichloroethene 1,1,1-Trichloroethane	5 50 50 50 85 - EPA Method 8260 50 50 80 80 80 80 80 80 80 80 80 80 80 80 80	<td>  NT  </td> <td><pre></pre></td> <td><ul> <li></li> <li></li> <li></li> <li>.53J</li> <li>1.2</li> <li></li> <li></li></ul></td> <td><pre></pre></td> <td>&lt;1 &lt; 1 &lt; 5</td> <td>&lt;1 &lt; 1 &lt; 5 0.96J &lt; 1 &lt; 1 0.82J 1.6 &lt; 1 &lt;</td> <td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td> <td>10/16/13  &lt;1 &lt;1</td> <td>6/9/14  &lt; 1 &lt; 5 &lt; 1 &lt; 1</td> <td>10/14/14  &lt; 1</td> <td>&lt;1 &lt; 2 0.52 0.99 &lt;1 &lt;1 1.2 J &lt;1 NT &lt;1 &lt;1</td> <td>&lt;1 &lt;1.5 &lt;2 0.27 J 0.54 &lt;1 &lt;1</td> <td>&lt;1 2.4 J</td> <td>&lt;1 &lt;1.5 &lt;2 0.53 0.97 &lt;1 1.3 J &lt;1 NT &lt;1 &lt;1</td> <td>&lt;1 &lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td> <td>&lt;1 &lt;5 &lt;5 &lt;1 1.3 &lt;1 &lt;1</td> <td>2.6 J &lt; 0.5 0.73 &lt; &lt; &lt;</td> <td><ul> <li></li> <li>0.36 J</li> <li>0.66</li> <li></li> <li></li></ul></td>	NT	<pre></pre>	<ul> <li></li> <li></li> <li></li> <li>.53J</li> <li>1.2</li> <li></li> <li></li></ul>	<pre></pre>	<1 < 1 < 5	<1 < 1 < 5 0.96J < 1 < 1 0.82J 1.6 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	10/16/13  <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	6/9/14  < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1	10/14/14  < 1	<1 < 2 0.52 0.99 <1 <1 1.2 J <1 NT <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1.5 <2 0.27 J 0.54 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 2.4 J	<1 <1.5 <2 0.53 0.97 <1 1.3 J <1 NT <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1 <1.5 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <5 <5 <1 1.3 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	2.6 J < 0.5 0.73 < < < < < < < < < < < < < < < < < < <	<ul> <li></li> <li>0.36 J</li> <li>0.66</li> <li></li> <li></li></ul>
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1 &lt; 5 &lt; 1 &lt; 1</td><td>10/14/14  &lt; 1     &lt; 2     &lt; 1     &lt; 1</td><td>&lt;1 &lt; 2 0.52 0.99 &lt; 1 &lt; 1 1.2 J &lt; 1 NT &lt; 1 &lt;</td><td>&lt;1 &lt;1.5 &lt;2 0.27 J 0.54 &lt;1 &lt;1</td><td>&lt;1 2.4 J</td><td>&lt;1 &lt;1.5 &lt;2 0.53 0.97 &lt;1 &lt;1 1.3 J &lt;1 &lt;1</td><td>&lt;1 &lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td><td>&lt;1 &lt;5 &lt;5 &lt;1 1.3 &lt;1 &lt;1</td><td>2.6 J &lt; 0.5 0.73 &lt; &lt; &lt;</td><td><ul> <li></li> <li>0.36 J</li> <li>0.66</li> <li></li> <li></li></ul></td></li<></ul></td>	NT	<pre></pre>	<ul> <li></li> <li></li> <li>.53J</li> <li>1.2</li> <li></li> <li></li></ul>	<ul> <li>&lt;</li> <li>&lt;</li> <li>1.4</li> <li>1.7</li> <li></li> <li>1.1</li> <li>1.4</li> <li>&lt;</li> <li></li> <li>&lt;</li> <li></li> <li>&lt;</li> <li>&lt;</li> <li>&lt;</li> <li>&lt;</li> <li>&lt;</li> <li></li> <li>&lt;</li> <li>&lt;</li> <li>&lt;</li> <li></li> <li< td=""><td>&lt;1 &lt; 1 &lt; 5</td><td>&lt;1 &lt; 1 &lt; 5 0.96J &lt; 1 &lt; 1 0.82J 1.6 &lt; 1 &lt; 1 &lt; 1 &lt; 1 &lt; 1 &lt; 1 &lt; 1 &lt; 1</td><td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td><td>10/16/13  &lt;1 &lt;1</td><td>6/9/14  &lt; 1 &lt; 5 &lt; 1 &lt; 1</td><td>10/14/14  &lt; 1     &lt; 2     &lt; 1     &lt; 1</td><td>&lt;1 &lt; 2 0.52 0.99 &lt; 1 &lt; 1 1.2 J &lt; 1 NT &lt; 1 &lt;</td><td>&lt;1 &lt;1.5 &lt;2 0.27 J 0.54 &lt;1 &lt;1</td><td>&lt;1 2.4 J</td><td>&lt;1 &lt;1.5 &lt;2 0.53 0.97 &lt;1 &lt;1 1.3 J &lt;1 &lt;1</td><td>&lt;1 &lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td><td>&lt;1 &lt;5 &lt;5 &lt;1 1.3 &lt;1 &lt;1</td><td>2.6 J &lt; 0.5 0.73 &lt; &lt; &lt;</td><td><ul> <li></li> <li>0.36 J</li> <li>0.66</li> <li></li> <li></li></ul></td></li<></ul>	<1 < 1 < 5	<1 < 1 < 5 0.96J < 1 < 1 0.82J 1.6 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1	<1 < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	10/16/13  <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	6/9/14  < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1	10/14/14  < 1     < 2     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1	<1 < 2 0.52 0.99 < 1 < 1 1.2 J < 1 NT < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 <1.5 <2 0.27 J 0.54 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 2.4 J	<1 <1.5 <2 0.53 0.97 <1 <1 1.3 J <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1 <1.5 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <5 <5 <1 1.3 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	2.6 J < 0.5 0.73 < < < < < < < < < < < < < < < < < < <	<ul> <li></li> <li>0.36 J</li> <li>0.66</li> <li></li> <li></li></ul>
Volatile Organic Compound Methylene chloride Acetone 2-Butanone Bromodichloromethane Dibromochloromethane Chloromethane Chloroform Bromoform Carbon Disulfide Iodomethane Vinyl Chloride 1,1-Dichloroethene 1,1-Dichloroethene trans-1, 2-Dichloroethene cis-1,2-Dichloroethene 1,1,1-Trichloroethane Trichloroethene Trichloroethene Tetrachloroethene Tetrachloroethene	5 50 50 50 85 - EPA Method 8260 50 50 80 80 80 80 80 80 80 80 80 80 80 80 80	<td>  NT  </td> <td><pre></pre></td> <td><ul> <li></li> <li></li> <li></li> <li>.53J</li> <li>1.2</li> <li></li> <li></li></ul></td> <td><pre></pre></td> <td>&lt;1 &lt; 1 &lt; 5</td> <td>&lt;1 &lt; 1 &lt; 5 0.96J &lt; 1 &lt; 1 0.82J 1.6 &lt; 1 &lt; 1</td> <td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td> <td>10/16/13  &lt;1 &lt;1</td> <td>6/9/14  &lt; 1     &lt; 5     &lt; 1     &lt; 1</td> <td>10/14/14  &lt; 1</td> <td>&lt;1 &lt; 2 0.52 0.99 &lt;1 &lt;1 1.2 J &lt;1 NT &lt;1 &lt;1</td> <td>&lt;1 &lt;1.5 &lt;2 0.27 J 0.54 &lt;1 &lt;1</td> <td>&lt;1 2.4 J</td> <td>&lt;1 &lt;1.5 &lt;2 0.53 0.97 &lt;1 &lt;1 1.3 J &lt;1 &lt;1</td> <td>&lt;1 &lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td> <td>&lt;1 &lt;5 &lt;5 &lt;5 &lt;1 1.3 &lt;1 &lt;1</td> <td>2.6 J &lt; 0.5 0.73 &lt; &lt; &lt;</td> <td><ul> <li></li> <li>0.36 J</li> <li>0.66</li> <li></li> <li></li></ul></td>	NT	<pre></pre>	<ul> <li></li> <li></li> <li></li> <li>.53J</li> <li>1.2</li> <li></li> <li></li></ul>	<pre></pre>	<1 < 1 < 5	<1 < 1 < 5 0.96J < 1 < 1 0.82J 1.6 < 1 < 1	<1 < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	10/16/13  <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	6/9/14  < 1     < 5     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1	10/14/14  < 1	<1 < 2 0.52 0.99 <1 <1 1.2 J <1 NT <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1.5 <2 0.27 J 0.54 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 2.4 J	<1 <1.5 <2 0.53 0.97 <1 <1 1.3 J <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1 <1.5 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <5 <5 <5 <1 1.3 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	2.6 J < 0.5 0.73 < < < < < < < < < < < < < < < < < < <	<ul> <li></li> <li>0.36 J</li> <li>0.66</li> <li></li> <li></li></ul>
Volatile Organic Compound Methylene chloride Acetone 2-Butanone Bromodichloromethane Dibromochloromethane Chloroform Bromoform Carbon Disulfide Iodomethane Vinyl Chloride 1,1-Dichloroethene 1,1-Dichloroethene trans-1, 2-Dichloroethene cis-1,2-Dichloroethene 1,1,1-Trichloroethane Trichloroethene	5 50 50 50 NV 7 50 NV NV 8 50 50 50 50 50 50 50 50 50 50 50 50 50	<td>  NT  </td> <td><pre></pre></td> <td><ul> <li></li> <li></li> <li>.53J</li> <li>1.2</li> <li></li> <li></li></ul></td> <td><ul> <li>&lt;</li> <li>&lt;</li> <li>1.4</li> <li>1.7</li> <li></li> <li>1.1</li> <li>1.4</li> <li>&lt;</li> <li></li> <li>&lt;</li> <li></li> <li>&lt;</li> <li>&lt;</li> <li>&lt;</li> <li>&lt;</li> <li>&lt;</li> <li></li> <li>&lt;</li> <li>&lt;</li> <li>&lt;</li> <li></li> <li< td=""><td>&lt;1 &lt; 1 &lt; 5</td><td>&lt;1 &lt; 1 &lt; 5 0.96J &lt; 1 &lt; 1 0.82J 1.6 &lt; 1 &lt; 1 &lt; 1 &lt; 1 &lt; 1 &lt; 1 &lt; 1 &lt; 1</td><td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td><td>10/16/13  &lt;1 &lt;1</td><td>6/9/14  &lt; 1 &lt; 5 &lt; 1 &lt; 1</td><td>10/14/14  &lt; 1     &lt; 2     &lt; 1     &lt; 1</td><td>&lt;1 &lt; 2 0.52 0.99 &lt; 1 &lt; 1 1.2 J &lt; 1 NT &lt; 1 &lt;</td><td>&lt;1 &lt;1.5 &lt;2 0.27 J 0.54 &lt;1 &lt;1</td><td>&lt;1 2.4 J</td><td>&lt;1 &lt;1.5 &lt;2 0.53 0.97 &lt;1 &lt;1 1.3 J &lt;1 &lt;1</td><td>&lt;1 &lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td><td>&lt;1 &lt;5 &lt;5 &lt;1 1.3 &lt;1 &lt;1</td><td>2.6 J &lt; 0.5 0.73 &lt; &lt; &lt;</td><td><ul> <li></li> <li>0.36 J</li> <li>0.66</li> <li></li> <li></li></ul></td></li<></ul></td>	NT	<pre></pre>	<ul> <li></li> <li></li> <li>.53J</li> <li>1.2</li> <li></li> <li></li></ul>	<ul> <li>&lt;</li> <li>&lt;</li> <li>1.4</li> <li>1.7</li> <li></li> <li>1.1</li> <li>1.4</li> <li>&lt;</li> <li></li> <li>&lt;</li> <li></li> <li>&lt;</li> <li>&lt;</li> <li>&lt;</li> <li>&lt;</li> <li>&lt;</li> <li></li> <li>&lt;</li> <li>&lt;</li> <li>&lt;</li> <li></li> <li< td=""><td>&lt;1 &lt; 1 &lt; 5</td><td>&lt;1 &lt; 1 &lt; 5 0.96J &lt; 1 &lt; 1 0.82J 1.6 &lt; 1 &lt; 1 &lt; 1 &lt; 1 &lt; 1 &lt; 1 &lt; 1 &lt; 1</td><td>&lt;1 &lt; 1 &lt; 5 &lt; 1 &lt;</td><td>10/16/13  &lt;1 &lt;1</td><td>6/9/14  &lt; 1 &lt; 5 &lt; 1 &lt; 1</td><td>10/14/14  &lt; 1     &lt; 2     &lt; 1     &lt; 1</td><td>&lt;1 &lt; 2 0.52 0.99 &lt; 1 &lt; 1 1.2 J &lt; 1 NT &lt; 1 &lt;</td><td>&lt;1 &lt;1.5 &lt;2 0.27 J 0.54 &lt;1 &lt;1</td><td>&lt;1 2.4 J</td><td>&lt;1 &lt;1.5 &lt;2 0.53 0.97 &lt;1 &lt;1 1.3 J &lt;1 &lt;1</td><td>&lt;1 &lt;1 &lt;1.5 &lt;2 &lt;1 &lt;1</td><td>&lt;1 &lt;5 &lt;5 &lt;1 1.3 &lt;1 &lt;1</td><td>2.6 J &lt; 0.5 0.73 &lt; &lt; &lt;</td><td><ul> <li></li> <li>0.36 J</li> <li>0.66</li> <li></li> <li></li></ul></td></li<></ul>	<1 < 1 < 5	<1 < 1 < 5 0.96J < 1 < 1 0.82J 1.6 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1	<1 < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	10/16/13  <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	6/9/14  < 1 < 5 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1	10/14/14  < 1     < 2     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1     < 1	<1 < 2 0.52 0.99 < 1 < 1 1.2 J < 1 NT < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 < 1 <	<1 <1.5 <2 0.27 J 0.54 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 2.4 J	<1 <1.5 <2 0.53 0.97 <1 <1 1.3 J <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <1 <1.5 <2 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	<1 <5 <5 <1 1.3 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1 <1	2.6 J < 0.5 0.73 < < < < < < < < < < < < < < < < < < <	<ul> <li></li> <li>0.36 J</li> <li>0.66</li> <li></li> <li></li></ul>

### Notes

- 1. Compounds detected in one or more samples are presented on this table.
- 2. Analytical testing completed by Alpha Analytical.
- 3. NYSDEC Class GA criteria obtained from Division of Water Technical and Operational Guidance Series (TOGS 1.1.1), dated October 1993, revised June 1998, January 1999 errata sheet,

and April 2000 addendum. \* Guidance value (not a standard) for 1,1-Dichloroethene = 0.07 ug/L as per the January 1999 update.

- 4. ug/L = part per billion (ppb).
- 5. < indicates compound was not detected; < 1 indicates compound was not detected above its respective reporting limit.
- 6. Shading indicates exceedance of Class GA Criteria.
- 7. NT = not tested.
- 8. NV = no value.
- 9. Results shown for IRM-1 for the September 2021 sampling event are the higher results from it or its respective duplicate.
- 10. Lab qualifiers: CH = continuing calibration outside of lab acceptance limits; results may be biased high. J = estimated concentration.
- L2 = analyte recovery in the control sample was below quality control limits; results may be biased low. Qualifiers for detected compounds only shown.

# Historical Analytical Data Summary Former Signore Facility 55-57 Jefferson Street Ellicottville, New York

Well I.D.   Analyte   September 2021   6/14/2019   6/20/2018   7/11/2017   10/25/2016   6/15/2016   10/21/2015   6/22015   10/14/2014   6/6/2014	W 11.15		Sample Date	Sample Date	Sample Date	Sample Date	Sample Date	Sample Date	Sample Date	Sample Date	Sample Date	Sample Date
EW-1.25	Well I.D.	Analyte	·		•	•			•	•		·
EW-1.65 TCE	EW 4.05	PCE										
EW-1.5 TCE	EVV-1.25				<	35	0.27 J	47	58	47	54	41
EW-2.5	T\\\ 1 E	PCE	<	<	<	0.23 J	0.24 J	<	0.22 J	0.2 J	0.22 J	<
EW-2.5	C.1-VV⊒	TCE	<	<	<	10	10	6.4	11	13	13	10
EW-3.5   PCE   NT   NT   NT   NT   NT   NT   NT   N	EW 25	PCE	<	<	<	<	<	<	<	<	<	<
EW-3.5 TCE NT	EVV-2.5	TCE	<	<	<	<	<	<	<	<	<	<
EW-4.5   PCE   0.78   1.5   0.8 J   1.6   4.6   4.6   3.9   5.4   5.4   5   5   1.1   1.6   4.6   4.6   3.9   5.4   5.4   5   5   1.1   1.6   4.6   4.6   3.9   5.4   5.4   5   5   5   1.1   1.6   4.6   4.6   3.9   5.4   5.4   5   5   5   5   5   5   5   5   5	E\M_3 5	PCE	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT
EW-4.5 TCE	L VV-3.3		NT	NT	NT	NT	NT	NT	NT	NT	NT	NT
NW-1    CE	E\M_4 5		0.78	1.5	<	0.8 J	1.6	1.3		1.7	1.5	1.7
MW-2    TCE	L VV -4.5		2.4	5	1.1	1.6		4.6	3.9	5.4		5
MW-2	M/M/-11	PCE	<	<	11.6	<	7.1	1.9	<	1.3	2.4	<
MW-9I         TCE <th< td=""><td>1010 0 - 11</td><td></td><td>&lt;</td><td>&lt;</td><td>15</td><td>&lt;</td><td>11</td><td>3</td><td>&lt;</td><td>2</td><td>2.8</td><td>&lt;</td></th<>	1010 0 - 11		<	<	15	<	11	3	<	2	2.8	<
MW-9	M\\\/-21		<	<	<	<	<	<	<	<	<	<
MW-9I         TCE         2.2         3.2         2.4         2.4         1.5         3         2.7         3         2.3           MW-4S         PCE         0.18 J  <	10100-21		<	<	<				<	<	<	<
MW-4S         PCE         0.18 J	M/M/_QI			l— — — — — — — -		0.73	l— — — — — — — -	0.34J	11		0.82	
MW-4S   TCE	10100-31		2.2	2.2	3.2	2.4	2.4	1.5	3	2.7	3	2.3
MW-5S	M\\/-4S	PCE	0.18 J	<	<	<	<	0.18J	0.32	0.22 J	0.36 J	<
TCE	10100 40					<						
ICE	M\\/-5S		3.6	4.8		<	l	4.9			3.7	4.6
TCE   0.24 J   0.35 J   <   0.33 J   0.36 J   0.32 J   0.38 J   0.35 J   0.34 J   <   0.44 J	10100-30		1.7		4.3	· ·		2.7		0.75	4	2.7
TCE   0.24 J   0.35 J   <   0.33 J   0.36 J   0.32 J   0.38 J   0.35 J   0.34 J   <   1	IRM-1	PCE			<		l					<
TCE   0.44 J   0.56   < 0.58   0.59   0.59   0.63   0.6   0.6   <   MAIN SCHOOL   PCE   NT   NT   NT   NT   NT   NT   NT   N	II XIVI- I		0.24 J	0.35 J	<	0.33 J	0.36 J	0.32J	0.38 J	0.35 J	0.34 J	<
TCE   0.44 J   0.56   <   0.58   0.59   0.59   0.63   0.6   0.6   <   MAIN SCHOOL   PCE   NT   NT   NT   NT   NT   NT   NT   N	IRM-2I				<		l			<		<
WELL         TCE         NT         NT			0.44 J					0.59	0.63		0.6	
TOWN WELL PCE < < < 0.23 J < 0.24 J < <				l— — — — — — — -			l— — — — — — — -					
	WELL		NT	NT	NT	NT		NT		NT	NT	NT
TCE   0.32 J   0.32 J   0.37   0.5   0.45 J   0.48 J   0.45 J   <	TOWN WELL				<	<	0.23 J					<
	TOWN WELL	TCE	0.32 J	0.32 J	<	0.37	0.5	0.45J	0.44 J	0.48 J	0.45 J	<

### Notes:

- 1. NT = Not Tested.
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- 3. Values shown are in ug/L (part per billion (ppb)).
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- \* Historical data provided date of 10/21/2004 However, based on historical sampling trend, sample was likely from April 2005 (sample date unknown).

# Historical Analytical Data Summary Former Signore Facility 55-57 Jefferson Street Ellicottville, New York

)A/ . II . I D	A 1 4 .	Sample Date									
Well I.D.	Analyte	10/15/2013	6/25/2013	10/31/2012	5/10/2012	10/13/2011	4/13/2011	6/2/2010	10/22/2009	4/23/2009	10/2/2007
EW 4.25	PCE	3.8	3.3	2	<	0.64J	0.78J	0.65	0.82	1.5	<
EW-1.25	TCE	59	51	1.7	59	90	56	73	90	88	5.1
EW-1.5	PCE	<	<	<	1.7	1.3	2.6	1.9	2.7	4.1	NT
LVV-1.5	TCE	3.9	8.4	9	13	9.5	19	14	20	18	NT
EW-2.5	PCE	<<	<	<	<	0	0	0	0	0	NT
	TCE	<	<	<	<	0	0	0	0	0	NT
EW-3.5	PCE	NT									
	TCE	NT									
EW-4.5	PCE	1.4	1.6	1.2	1.2	1.3	1.9	0.97	1.3	2.5	NT
	TCE	5.8	6.8	7	7.6	6.9	10	5.5	7.9	8	NT
MW-1I	PCE	<	<	<	<	<	<	<	<	<	<
	TCE	<	<	<	<	<	<	<	<	<	<
MW-2I	PCE	<u> </u>	<	<	<	<	<<	<	<	<	NT
	TCE	<	<	<	0.83J	<	<	<	<	<	NT
MW-9I	PCE		<	1.3	0.89J	1.4	1.4	0.9	1.6	2.2	1.1
	TCE	3.4	2.4	3.1	2.7	3.7	3.6	2.9	4.5	4.6	3.8
MW-4S	PCE	<	<	<	<	<	<	<	<	<	<
	TCE	<	<	<	<	<	<	<	<	<	<
MW-5S	PCE	3.9	3.5	2.6	0.59J	2.9	.61J	1.7	3.1	3.4	4.8
	TCE	8.1	6.9	3.1	17	15	12	14	22	30	19
IRM-1	PCE	<	<	<	<	<	.54J	<	<	<	NT
	TCE	<	0.52J	<	<	.52J	.69J	<	<	<	NT
IRM-2I	PCE	<	<	<	<	< <	<	<	<	<	NT
	TCE	<	0.72J	0.60J	0.74J	.86J	.88J	0.85	0.89	<	NT
IAIN SCHOOL	PCE	NT	NT	NTNT	NT						
WELL	TCE	NT									
TOWN WELL	PCE	<	<	<b>/</b>	<	<	<	<	NT	<	NT
TOVVIN VVLLL	TCE	<	0.63J	0.58J	<b>~</b>	.55J	.69J	<	NT	<	NT

### Notes:

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# Historical Analytical Data Summary Former Signore Facility 55-57 Jefferson Street Ellicottville, New York

EW-1.25	Wall I D	A male da	Sample Date									
EW-1.5   TCE   45.1   66   27.9   66.9   31   53   22   110   40   30   EW-1.5   PCE   16.7   18   15.6   25.3   28   1   20   <   13   6   EW-2.5   PCE   <   <   <   <   <   <   <   <   <	weii i.D.	Analyte	10/30/2006	4/25/2006	11/14/2005	4/1/2005*	10/21/2004	4/29/2004	10/16/2003	4/11/2003	10/23/2002	4/12/2002
TCE	EM 4.05	PCE	1.5	<	1.6	2	2	6.3	3	11	4	3
EW-1.5	EVV-1.25	TCE	45.1	66	27.9	66.9	31	53	22	110	40	30
EW-2.5   PCE   <	E\M_1 5		3.1	3	2.9	5	6	<	5	<	4	2
EW-2.5	L VV-1.5		16.7	18	15.6	25.3	28	1	20	<	13	6
EW-3.5   PCE   <	FW-2 5		<	<	<	<	<	<	<<	<	<	<
EW-4.5			<		<		<		<		<	
C	EW-3.5		<b></b>		<		<				+	
NT   NT   NT   NT   NT   NT   NT   NT											1	
NW-11	EW-4.5	. – – – – –						<u> </u>			<	
MW-2    TCE			5.9		4.6		5			NT	<	NT
N	MW-1I		<	<	<	<	<				<	<
MW-2I         TCE         <         NT         <         28         1         <         1         NT           MW-9I         PCE         1.4         NT         1.5         NT         2         NT         3         NT         2         NT           MW-9I         PCE         3.5         NT         3.3         NT         2         NT         3         NT         2         NT           MW-4S         PCE         <			<		<		<		0.6		<	
MW-9	MW-2I		<		<		<		<	NT NT	<	<del> </del>
MW-9I         TCE         3.5         NT         3.3         NT         3         NT         5         NT         3         NT           MW-4S         PCE         <									·		1	
MW-4S	MW-9I	1						<u> </u>			<del></del>	!
MW-4S   TCE   <   NT   <   NT   <   NT   <   NT   <   NT	10100 01		3.5		3.3		3		5		3	
MW-5S	M\\/-4S		<		<		<		<		<	
TCE   30.5   NT	10100 40											
ICE   30.5   N1   1   N1   26   N1   29   N1   26   N1   IRM-1   PCE   <   NT   <   NT     0.5   NT   <   NT   <   NT     NT     NT     NT     IRM-2    PCE   <   <   <   <   <   <   <   <   <	M\\/_5\$	1	2.3	NT	4.1	NT	10	NT	10	NT	10	NT
IRM-1	10100-30	TCE	30.5	NT	1	NT	26	NT	29	NT	26	NT
ICE   <   NI   <   NI   0.7   NI   0.7   NI   1   NI   NI	IRM-1	. – – – – –	<	NT	<	NT	0.5	NT	<	NT	<	NT
TCE	II XIVI- I	TCE	<	NT	<	NT	0.7	NT	0.7	NT	1	NT
ICE	IRM-2I			<		<	0.5	<	<	<		
WELL         TCE         <         NT         <         NT         <         NT         <         NT          NT         <         NT          NT         <         NT         <         NT         <         NT         <         NT          NT         <         NT            <			1.2		0.71	1	1		<		NT	
TOWN WELL PCE	MAIN SCHOOL	1	<		<		<	<u> </u>	<		<	
	WELL		<	NT								
TCE   0.52   <   0.62   0.8   0.9   0.9   1   <   1   1	TOWN WELL			<	<	<	<	<	<	<	<	<
	TOVVIN VVELL	TCE	0.52	<	0.62	0.8	0.9	0.9	1	<	1	

### Notes:

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- \* Historical data provided date of 10/21/2004 However, based on historical sampling trend, sample was likely from April 2005 (sample date unknown

# Historical Analytical Data Summary Former Signore Facility 55-57 Jefferson Street Ellicottville, New York

)		Sample Date									
Well I.D.	Analyte	11/9/2001	4/27/2001	10/25/2000	5/30/2000	10/15/1999	4/27/1999	11/5/1998	4/16/1998	10/30/1997	4/8/1997
EVA 4.05	PCE	4	3	3	2	32	ND	8	6	5	5
EW-1.25	TCE	11	39	35	31	25	55	63	66	120	78
EW-1.5	PCE	<	3	3	2	3	2	1	4	2	3
EVV-1.5	TCE	<	8	7	4	6	6	6	8	8	10
EW-2.5	PCE	<	<	<	<	<	<	<	NT	<	<
L VV - Z. J	TCE	6	<	<	<	<	<	<	NT	<	<
EW-3.5	PCE	<	8	<	2	<	2	<	3	<	4
LW 5.5	TCE	2	NT	2	NT	2	NT	1	NT	2	NT
EW-4.5	PCE	<<	NT	<	NT	< <	NT	<	NT	<	NT
L VV 4.0	TCE	<	NT								
MW-1I	PCE	<	<	<	<	< <	NT	<	NT	<	<
	TCE	<	<	<	<	<	NT	<	1	<	3
MW-2I	PCE	<	NT	NT	NT	<	NT	<	NT	<	NT
21	TCE	<	NT								
MW-9I	PCE	<<	NT NT	2	NT	<	NT	<	NT	NT	NT
	TCE	2	NT	3	NT	<	NT	2	NT	3	NT
MW-4S	PCE	<	NT	<	NT	< <	NT	<	NT	25	NT
	TCE	<	NT	1	NT	<	NT	<	NT	<	NT
MW-5S	PCE	8	NT	12	NT	12	NT	19	NT	<	NT
10177 00	TCE	21	NT	30	NT	18	NT	36	NT	80	NT
IRM-1	PCE	<	NT	<	NT	< <	<	<	NT	<	NT
11 (14)	TCE	NT	NT	1	NT	<	<	<	NT	2	NT
IRM-2I	PCE	NT	<	<	<	<	<	<	NT NT	<	<<
	TCE	<	2	2	1	<	<	<	2	2	3
MAIN SCHOOL	PCE	NT	NT	NT	NT	< <	NT	<	NT	<	NT
WELL	TCE	NT	NT	NT	NT	<	NT	<	NT	<	NT
TOWN WELL	PCE	<	<	<	<	<	<	<	NT NT	<	<<
	TCE	<	2	2	1	<	<	<	2	2	3

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# Historical Analytical Data Summary Former Signore Facility 55-57 Jefferson Street Ellicottville, New York

Well I.D.	Analyta	Sample Date	Sample Date	Sample Date	Sample Date	Sample Date	Sample Date	Sample Date	Sample Date	Sample Date
well i.D.	Analyte	10/17/1996	4/16/1996	2/8/1996	10/13/1994	7/11/1994	4/26/1994	2/14/1994	11/1/1993	7/13/1993
EW-1.25	PCE	14	39	26	23	NT	NT	NT	7	NT
EVV-1.25	TCE	78	86	83	100	NT	NT	NT	6	NT
EW-1.5	PCE	7	9	10	16	<	3	2	<	<
LVV-1.5	TCE	10	10	9	10	1	2	2	<	<
EW-2.5	PCE	<	<	<	<	NT	NT	NTNT	<<	NT
L VV 2.5	TCE	<	2	<	<	NT	NT	NT	<	NT
EW-3.5	PCE	<	NT	<	<	NT	NT	NT	<	NT
L VV 0.0	TCE	1	NT	4	1	NT	NT	NT	<	NT
EW-4.5	PCE	<	NT	<	<	<	< <	<	<<	<
L V 4.0	TCE	1	NT	2	2	<	<	<	2	<
MW-1I	PCE	<	<	11	<	<	<	<	<	<
	TCE	4	2	12	17	2	5	5	1	6
MW-2I	PCE	<	NT	<	2	NT	NT	NT	NT	NT
	TCE	<	NT	<	NT	NT	NT	NT	NT	NT
MW-9I	PCE	NT	NT	4	33	NT	2	NT	3	NT
	TCE	2	NT	5	6	NT	4	NT	7	NT
MW-4S	PCE	30	NT	2	<	NT	NT	NT	2	NT
	TCE	71	NT	2	2	NT	NT	NT	4	NT
MW-5S	PCE	<	NT	50	NT	NT	NT	NT	<	NT
	TCE	3	NT	63	NT	NT	NT	NT	6	NT
IRM-1	PCE	<	NT	<u>4</u>	1	4	2	3	5	4
	TCE	2	NT	5	4	5	4	5	6	5
IRM-2I	PCE		2 2	2	22	2	2	22	3	2
	TCE	2	4	4	4	5	4	5	6	4
MAIN SCHOOL	PCE	<	<del>-</del>	1	· <del>'</del>	NT		NT		<u>N</u> T
WELL	TCE	<	1	2	1	NT	1.5	NT	2	NT
TOWN WELL	PCE		1 1	$\frac{2}{4}$	1 1	NT	2	NT		2
	TCE	Notes:	3	4	4	NT	4.6	NT	5	4

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- \* Historical data provided date of 10/21/2004 However, based on historical sampling trend, sample was likely from April 2005 (sam

# Historical Analytical Data Summary Former Signore Facility 55-57 Jefferson Street Ellicottville, New York

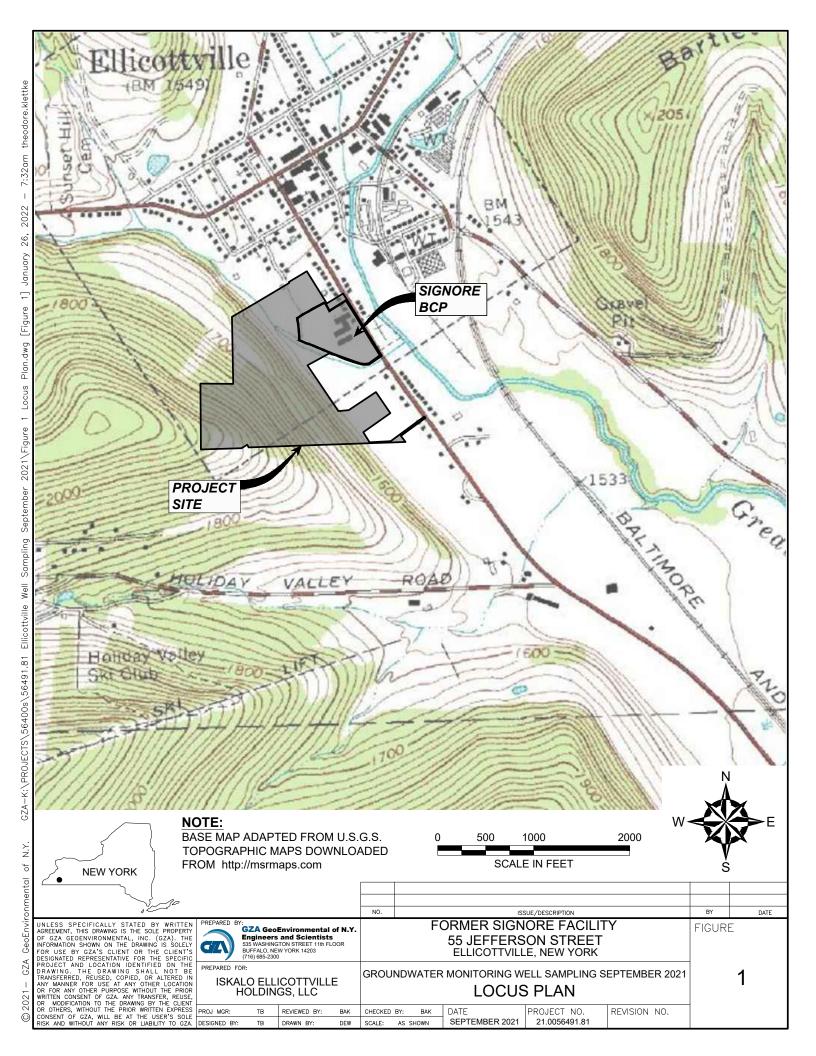
Wall I D	A   - 4	Sample Date	Sample Date	Sample Date	Sample Date	Sample Date	Sample Date	Sample Date	Sample Date	Sample Date
Well I.D.	Analyte	4/26/1993	1/26/1993	12/21/1992	1/7/1992	6/11/1991	6/28/1990	12/5/1990	6/25/1990	1/15/1989
EW-1.25	PCE	NT	NT	140	NT	NT	NT	NT	NT	NT
EVV-1.25	TCE	NT	NT	67	NT	NT	NT	NT	NT	NT
EW-1.5	PCE	9	<	14	NT	NT	NT	NT	NT	NT
LVV-1.5	TCE	7	<	7	NT	NT	NT	NT	NT	NT
EW-2.5	PCE	NT	NT	<	NT	NT	NT	NT	NT	NT
L V V - Z.J	TCE	NT	NT	2	NT	NT	NT	NT	NT	NT
EW-3.5	PCE	NT	NT	<	NT	NT	NT NT	NT	NT	NT
L VV 0.0	TCE	NT	NT	<	NT	NT	NT	NT	NT	NT
EW-4.5	PCE	<	2	4	NT	NT	NT	NTNT	NT	NT
EVV 4.0	TCE	2	<	8	NT	NT	NT	NT	NT	NT
MW-1I	PCE	11	11	2	NT	NT	6	NT	NT	19
	TCE	36	54	66	NT	NT	55	NT	NT	110
MW-2I	PCE	NT	NT	NT	NT	NT	NT	NT	NT	NT
	TCE	NT	NT	NT	NT	NT	NT	NT	NT	NT
MW-9I	PCE	<	NT	<	NT	NT	10	NT	NT	<
	TCE	3	NT	<	NT	NT	28	NT	NT	20
MW-4S	PCE	NT	NT	2	NT	NT_	10	NT	NT	15
	TCE	NT	NT	1	NT	NT	27	NT	NT	53
MW-5S	PCE	NT	NT NT	<	NT	NT	99	NT	NT	<
	TCE	NT	NT	74	NT	NT	100	NT	NT	150
IRM-1	PCE	33	4	5	NT	NT	4	NT	NT	NT
	TCE	5	6	5	NT	NT	6	NT	NT	NT
IRM-2I	PCE	2	3	3	NT	NT	5	NT	NT.	NT
	TCE	4	5 NT	4	NT	NT	9	NT	NT	NT
MAIN SCHOOL	PCE	<del>&lt;</del>	NT	0.6	1 1	0.8	NT			<u>NT</u>
WELL	TCE	) NT	NT	1.9	3	1.7	NT	2.2	2	NT
TOWN WELL	PCE	NT	<u>NT</u>	3.5	4	55	<u>N</u> T	5	NT	<u>N</u> T
	TCE	NT	NT	6.1	7	6.3	NT	8	NT	NT

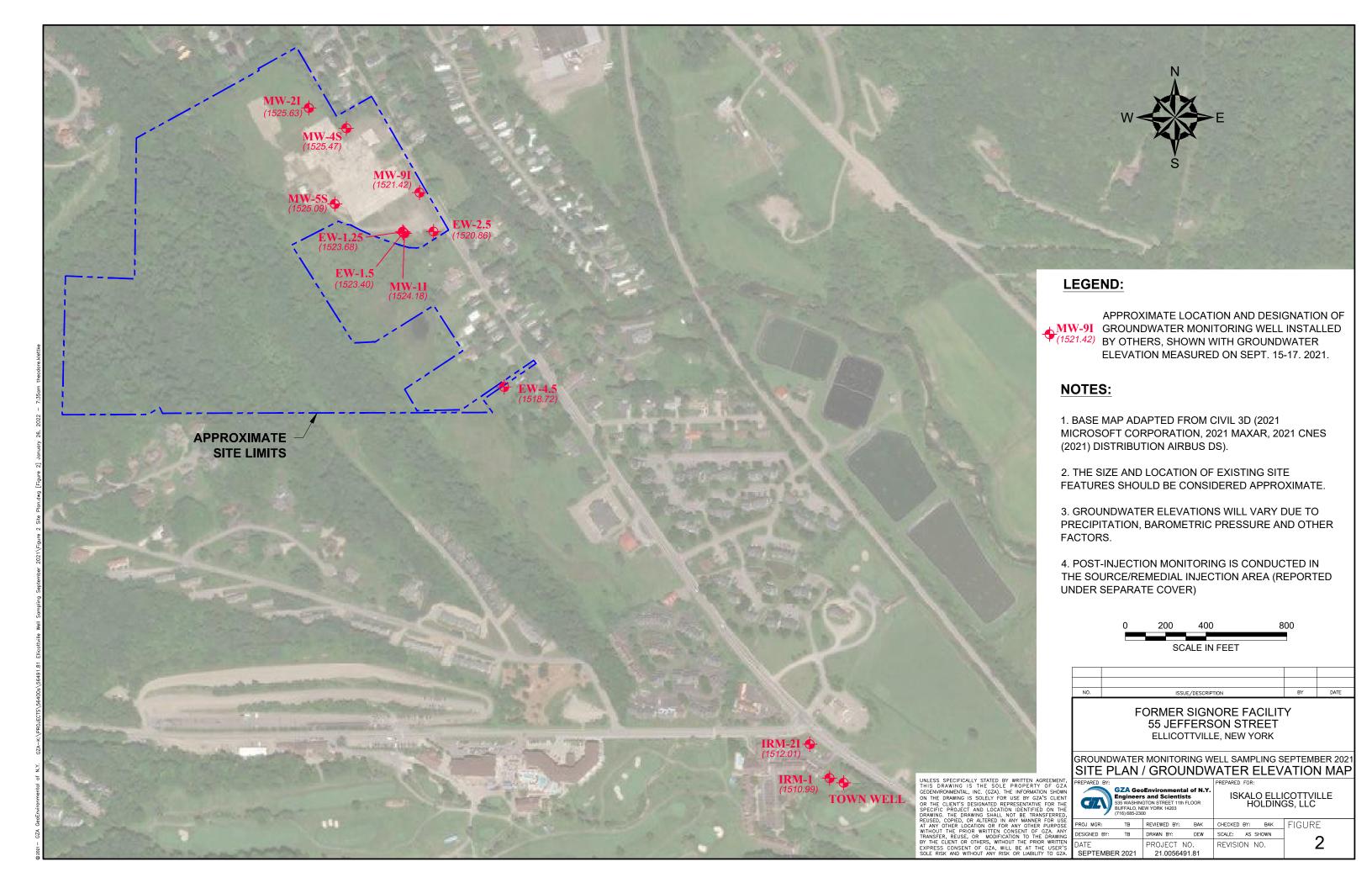
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- \* Historical data provided date of 10/21/2004 However, based on historical sampling trend, sample was likely from April 2005 (sam



**FIGURES** 







### **APPENDIX A**

**LIMITATIONS** 

# **GEOHYDROLOGICAL LIMITATIONS** 21.0056491.81



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### **USE OF REPORT**

1. GZA GeoEnvironmental, Inc. (GZA) prepared this report on behalf of, and for the exclusive use of our Client for the stated purpose(s) and location(s) identified in the Proposal for Services and/or Report. Use of this report, in whole or in part, at other locations, or for other purposes, may lead to inappropriate conclusions; and we do not accept any responsibility for the consequences of such use(s). Further, reliance by any party not expressly identified in the agreement, for any use, without our prior written permission, shall be at that party's sole risk, and without any liability to GZA.

### STANDARD OF CARE

- 2. GZA's findings and conclusions are based on the work conducted as part of the Scope of Services set forth in the Proposal for Services and/or Report and reflect our professional judgment. These findings and conclusions must be considered not as scientific or engineering certainties, but rather as our professional opinions concerning the limited data gathered during the course of our work. Conditions other than described in this report may be found at the subject location(s).
- 3. GZA's services were performed using the degree of skill and care ordinarily exercised by qualified professionals performing the same type of services, at the same time, under similar conditions, at the same or a similar property. No warranty, expressed or implied, is made. Specifically, GZA does not and cannot represent that the Site contains no hazardous material, oil, or other latent condition beyond that observed by GZA during its study. Additionally, GZA makes no warranty that any response action or recommended action will achieve all of its objectives or that the findings of this study will be upheld by a local, state or federal agency.
- 4. In conducting our work, GZA relied upon certain information made available by public agencies, Client and/or others. GZA did not attempt to independently verify the accuracy or completeness of that information. Inconsistencies in this information which we have noted, if any, are discussed in the Report.

### SUBSURFACE CONDITIONS

- 5. The generalized soil profile(s) provided in our Report are based on widely-spaced subsurface explorations and are intended only to convey trends in subsurface conditions. The boundaries between strata are approximate and idealized, and were based on our assessment of subsurface conditions. The composition of strata, and the transitions between strata, may be more variable and more complex than indicated. For more specific information on soil conditions at a specific location refer to the exploration logs. The nature and extent of variations between these explorations may not become evident until further exploration or construction. If variations or other latent conditions then become evident, it will be necessary to reevaluate the conclusions and recommendations of this report.
- 6. Water level readings have been made, as described in this Report, in and monitoring wells at the specified times and under the stated conditions. These data have been reviewed and interpretations have been made in this report. Fluctuations in the level of the groundwater however occur due to temporal or spatial variations in areal recharge rates, soil heterogeneities, the presence of subsurface utilities, and/or natural or artificially induced perturbations. The observed water table may be other than indicated in the Report.

### **COMPLIANCE WITH CODES AND REGULATIONS**

7. We used reasonable care in identifying and interpreting applicable codes and regulations necessary to execute our scope of work. These codes and regulations are subject to various, and possibly contradictory, interpretations. Interpretations and compliance with codes and regulations by other parties is beyond our control.



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### **SCREENING AND ANALYTICAL TESTING**

- 8. GZA collected environmental samples at the locations identified in the Report. These samples were analyzed for the specific parameters identified in the report. Additional constituents, for which analyses were not conducted, may be present in soil, groundwater, surface water, sediment and/or air. Future Site activities and uses may result in a requirement for additional testing.
- 9. Our interpretation of field screening and laboratory data is presented in the Report. Unless otherwise noted, we relied upon the laboratory's QA/QC program to validate these data.
- 10. Variations in the types and concentrations of contaminants observed at a given location or time may occur due to release mechanisms, disposal practices, changes in flow paths, and/or the influence of various physical, chemical, biological or radiological processes. Subsequently observed concentrations may be other than indicated in the Report.

### INTERPRETATION OF DATA

11. Our opinions are based on available information as described in the Report, and on our professional judgment. Additional observations made over time, and/or space, may not support the opinions provided in the Report.

### **ADDITIONAL INFORMATION**

12. In the event that the Client or others authorized to use this report obtain additional information on environmental or hazardous waste issues at the Site not contained in this report, such information shall be brought to GZA's attention forthwith. GZA will evaluate such information and, on the basis of this evaluation, may modify the conclusions stated in this report.

### **ADDITIONAL SERVICES**

13. GZA recommends that we be retained to provide services during any future investigations, design, implementation activities, construction, and/or property development/ redevelopment at the Site. This will allow us the opportunity to: i) observe conditions and compliance with our design concepts and opinions; ii) allow for changes in the event that conditions are other than anticipated; iii) provide modifications to our design; and iv) assess the consequences of changes in technologies and/or regulations.



### **APPENDIX B**

WELL DEVELOPMENT FORMS

File: 21.0056491.81

### FORMER SIGNORE, INC. FACILITY WELL DEVELOPMENT FORM 55-57 JEFFERSON STREET ELLICOTTVILLE, NEW YORK

						Historic Info	ormation	Agrange .			
Boring Log A	vailable (	es/no/attac	ched):								
Installation L			,								
						Summ	nary				
Monitoring Well: EW-1.25R Ground Surface Elevation: Riser/Screen Material: PVC											
	nstallation Date: 5/2019 Protective Casing Elevation: Top of Screen Depth: 15 ft.										
Installed By:											
				Elevation D							
Previous Fie	eld measur	ement Infor	mation Availa	ble (yes/ <b>no</b> /							
						s of Previous F					
Depth to	Water		pН	Specific (	Conductance	Tempera	ature	Tu	rbidity		Color
(ft)	)	(Standa	ard Units)	(uM	hos/cm)	( °C)	)	(1)	NTU)		
9.5	1	6	5.77		0.65	14.7	,	1	3.13		Clear
Notes:											
			Fie	eld Observa	tions				Parame	eters +/-	Sampling Information
Exterior Obs	ervations:	(+000)							рН	+/- 0.1	Sample ID: EW-1.25 R · 091721
									Conductivi		Sample Time: (220
Interior Obse	ervations	6000	G000								# of Sample Containers: 8
						Turbidity		Duplicate Sample ID:			
	_								ORP		Sample Analysis: VOCs 8260
Signs of Dar						~\· ·	Inin 14		DO	+/- 10%	MNA Parameters TOC, DISS-
Locked ((	yes/no)	Well Ca	p (ves/no)	Surfa	ace Seal Intact		PID Meas	urement:	0.0000	Odors: 🙏	one alved GAS (RSK175)
						Well Qual	ity Data		13.5		
Date	Time	Depth to	Cumulative	pH	Specific	Temperature	Turbidity	Color	Dissolved	Oxygen	Notes
Date	Time	Water	Volume	(Standard	'	'	(NTU)	Color		Reduction	Notes
		ft bgs	Purged	Units)	(uMhos/cm)	(°C)	(1010)		Oxygen	Potential	
9-17-21	100 11	10.38	o Purged	6.37	0.660	14.8	13.16	NONE	20.0	* 95.6	Depth of Water: /ø, 3 6
1	17.10	10.38	0.3	6.28	0.650	14,4	8.44	NONE	17.6	-95.7	Length of Water Column:
	1215	10.38	0.6	6128	0.645	14.2	9.18	None	17.9	-99.6	Depth of Well:
	1770	10.3%	0.9	6.28	0.641	14.1	9.32	NONE	17.1	-98.9	Sheen Observed: Y N
			,							1,-2,-3	DNAPL Observed: Y 9V
											Did Well Go Dry: Y 'N'
											Other:
		1									

File: 21.0056491.81

### FORMER SIGNORE, INC. FACILITY WELL DEVELOPMENT FORM 55-57 JEFFERSON STREET ELLICOTTVILLE, NEW YORK

	J.,			V 5 3 5 7		Historic Info	ormation						
Boring Log A	vailable ( <b>y</b>	es/no/attac	hed):										
Installation L	og Availab	le ( <b>yes</b> /no/a	attached)										
Summary													
Monitoring V	Vell:	EW-1.5			rface Elevation						inless Steel		
Installation E	nstallation Date: 7/90 Protective Casing Elevation: 1534.32 ft. Top of Screen Depth: 40 ft.												
nstalled By: Empire Soils Monitoring Point Elevation: 1533.92 ft. Bottom of Screen Depth: 50 ft.													
Elevation Datum:													
Previous Field measurement Information Available (yes/no/attached)													
Ranges of Previous Field Measurements													
Depth to	Water	F	рΗ	Specific (	Conductance	Tempera	ature	Tur	bidity		Color		
(ft)	)	(Standa	ard Units)	(uMl	hos/cm)	(°C)		(N	ITU)				
10.0	)9	7	.49	0	.575	14.1		6	5.51		Clear		
Notes:			•										
			Fie	ld Observa	tions				Parame	eters +/-	Sampling Information		
Exterior Obs	ervations:	60019							pН	+/- 0.1	Sample ID: EW-1.5 - 691521		
									Conductivi		Sample Time: 1140		
Interior Obse	ervations	Roadbox	Flooded								# of Sample Containers: 3		
									Turbidity		Duplicate Sample ID:		
									ORP		Sample Analysis: VOCs 8260		
Signs of Dar					19119-1-221				DO	+/- 10%			
Locked (	yés/no)	Well Ca	p (yes/no)	Surfa	ace Seal Intact		PID Meas	urement: ،	o. oppm	Odors: 🖊	SAD		
						Well Qual	ity Data						
Date	Time	Depth to	Cumulative	pН	Specific	Temperature	Turbidity	Color	Dissolved	Oxygen	Notes		
Date	I line	Water	Volume	(Standard	Conductance	(°C)	(NTU)	COIO	Oxygen	Reduction	140(65		
		ft bgs	Purged	Units)	(uMhos/cm)	( C)	(1410)		Oxygen	Potential			
9-15-21	1125	10.84	ruigeu O	11.31	0.929	15.3	8.80	NONE	33.4	-/5/19	Depth of Water: 70.5%		
1	1130	10.84	0.2	11,46	0,984	14.9	17,37	NONE	18.4	-150.6	Length of Water Column: 54.3%		
	1135	10.84	0:4	11.50	0.982	14.9	21,12	Mane	17.9	-145.2	Depth of Well: 64,90		
1	1140	10.84	0.6	11.52	0,984	14.8	20:26	110-01	17,4	-140,21	Sheen Observed: Y (N)		
Did Well Go Dry: Y (N)										Did Well Go Dry: Y (N)			
Other:													

						Historic Info	ormation				
Boring Log Available ( <b>yes</b> /no/attached):											
Installation L	og Availab	le ( <b>yes</b> /no/a	attached)								
						Summ	ary				
Monitoring V		EW-2.5			rface Elevation					al: Steel/Sta	inless Steel
Installation D		7/90			Casing Elevation				creen Depth		
Installed By:		Empire Soi	ls		Point Elevation	n: 1533.92 ft.		Bottom o	of Screen De	epth: 50 ft.	
				Elevation D						-	
Previous Fie	eid measur	ement Infor	mation Availa	ble (yes/ <b>no</b> /							
	107			0 10 4		s of Previous F					
Depth to			pH		Conductance	Tempera			rbidity		Color
(ft)	10		ard Units)		nos/cm)				NTU)		0.
	+2		7.54	0	.629	14			5.06		Clear
Notes:											
	7.7.5.0	S 7 70 L	Fie	eld Observa	fions		THE THE		Darame	eters +/-	Compling Information
Exterior Ohs	envations:	ha loc	h -> Plac			on well			pH		Sampling Information Sample ID: EW-2.5 -091521
Exterior Obs	servations.	100 100	re s Place	DO POED	1000	on well			Conductivit		Sample Time: 0858
Interior Obse	ervations										
		Turbidity +/- 10% Duplicate Sample ID:									
									ORP		Sample Analysis: VOCs 8260
Signs of Dar									DO	+/- 10%	
Locked (	yes(no)	Well Ca	p (yes/no)	Surfa	ace Seal Intact		PID Measu	urement:	0.0 ppm	Odors: 🙌	i a_tv
		Page 1			-1 1-5 V	Well Qual	ty Data		51115		
Date	Time	Depth to	Cumulative	pН	Specific	Temperature	Turbidity	Color	Dissolved	Owagon	Notes
Date	Tille	Water	Volume		Conductance	(°C)	(NTU)	Color	Oxygen	Oxygen Reduction	notes
		ft bgs	Purged	Units)	(uMhos/cm)	( 0)	(1410)		Oxygen	Potential	
9-15-21	0833	17,10	O O	9.08	0.211	14.3	43.92	None	40,0		Depth of Water: 13.06
	0838	13,10	0.10	8.98	0,210	13.9	30.45	None	17.3		Length of Water Column: 34,52
	0843	13,10	0.20	9,00	0.212	13,7	25.24	Noni	17,1		Depth of Well: 석가 5명
	0848	13.10	0.30	9:01	0.212	13.6	22.22	None	13.9		Sheen Observed: Y (N)
	0853	13.10	0.40	9.00	0.212	13.7	21,13	Nane	16.60		DNAPL Observed: Y N
	0858	13,10	0.50	9.00	0,212	13,7	20.58	None	110.5	-150.1	Did Well Go Dry: Y/N
											Other:
GZA GooEn		Logit Marris V									

	W. 3.					Historic Info	rmation				
Boring Log A	vailable (y	es/no/attac	hed):								
Installation L	og Availab	le (yes/no/a	attached)								
						Summ	ary				
Monitoring W	Vell:	EW-4.5		Ground Sur	face Elevation	: 1533.55		Riser/Scr	een Materia	al: Steel/Sta	inless Steel
Installation D		7/90		Protective 0	Casing Elevation	n: 1535.97 ft.			creen Depth		
Installed By:		Empire Soi	ls	Monitoring	Point Elevation	: 1535.65 ft.		Bottom o	f Screen De	pth: 50 ft.	
				Elevation D	atum:						
Previous Fie	ld measure	ement Inforr	mation Availa	ble (yes/ <b>no</b> /							
		V			Ranges	of Previous F	ield Measu	rements			
Depth to	Water	1	ЭΗ	Specific C	Conductance	Tempera	ature	Tur	bidity		Color
(ft)		(Standa	ard Units)	(uMl	nos/cm)	( °C)	)	(NTU)			
16.2			.99	0	.581	13.7		9	.12		Clear
Notes:											
10.00											
			Fie	eld Observa	tions				Parame	eters +/-	Sampling Information
Exterior Obs	ervations:	1000 1	Russed-C	0+					рН	+/- 0.1	Sample ID: EW-4.5 - 09157.)
									Conductivi		Sample Time: 13/5
nterior Observations Temperature +/- 10% # of Sample Containers: 9											
	Turbidity +/- 10%   Duplicate Sample ID:   M.S.   M.S.										
									ORP		Sample Analysis: VOCs 8260
Signs of Dar									DO	+/- 10%	
Locked (	yes/no)	Well Ca	p (yes/no)	Surfa	ace Seal Intact			urement:	0 000	Odors: 📈	OME
						Well Qual	ity Data				
_						<b>-</b> .		0.1	D:1		N
Date	Time	Depth to	Cumulative	pΗ	Specific	Temperature	Turbidity	Color	Dissolved	Oxygen	Notes
l .		Water	Volume	(Standard	Conductance	(°C)	(NTU)		Oxygen	Reduction	
		ft bgs	Purged	Units)	(uMhos/cm)	7	,		_	Potential	Double of Wilder
9-15-21	1240	17.02	0	9.00	0.786	14.0	105.20	NONE	25.6	-205.0	Depth of Water: 16,93 Length of Water Column: 30,22
	1305	17.02	0,5	7.03	0.692	13.8	40.82	None	30.8	-145,1	Depth of Well: 97, 15
	1310	17.02	0.6	7.01	0.693	14.1	34.27	Almine	29.1	-139.1	Sheen Observed: Y N
- 1	1315	17.07	0.7	1,00	0.613	. / 44 1	31161	JUDNIC	21,1	-137.8	DNAPL Observed: Y N
											Did Well Go Dry: Y N
		-								V	Other:
										9	

Historic Information											
Boring Log Available (yes/ <b>no</b> /attached):											
Installation Lo	og Availab	le ( <b>yes</b> /no/a	attached)								
						Summ					
Monitoring W		MW-1 I			face Elevation				een Materia		
Installation D		2/87			Casing Elevation				creen Depth		
Installed By:		Rochester I		_	Point Elevation	: 1531.79 ft.		Bottom of	f Screen De	pth: 50 ft.	
				Elevation D							
Previous Fiel	d measure	ement Inforr	nation Availat	ole (yes/ <b>no</b> /							
						of Previous Fi					
Depth to	Water		Н	Specific C	Conductance	Tempera			bidity		Color
(ft)		(Standa	ard Units)	(uMł	nos/cm)	(°C)			ITU)		
9.79	9	6	.87	(	).14	16.7		10	0.36		Clear
Notes:											
			Fie	eld Observa	tions					eters +/-	Sampling Information
Exterior Obse	ervations:	Good							рН	+/- 0.1	Sample ID: MW-1 I - 09/52/
									Conductivi		Sample Time: 1115
Interior Obse	ervations	Read box	Flooded								# of Sample Containers: 3
		Turbidity +/- 10% Duplicate Sample ID:									
	_								ORP DO	+/- 10mv +/- 10%	Sample Analysis: VOCs 8260
Signs of Dan				C	Caal lataat	(handha)	DID Mass			Odors: 🗸	
Locked (y	/es/no)	well Ca	p (yes/no)	Suna	ace Seal Intact	(yesmo) Well Quali	PID Meas	urement.	O. O ppm	Odors. 70	
						well Qual	ly Data				
Data	Time	Donth to	Cumulative	pН	Specific	Temperature	Turbidity	Color	Dissolved	Oxygen	Notes
Date	riine	Depth to Water	Volume	(Standard	Conductance	(°C)	(NTU)	COIO	Oxygen	Reduction	Notes
				Units)	(uMhos/cm)	( )	(1410)		Chygen	Potential	
9-15-21	1050	ft bgs	Purged	6.90	0 · 537	13.9	80.32	NONG	28.3	-61.4	Depth of Water: /o, 61
1-12-51	1100	10.71	0.4	6.10	0.533	12.5	58.14	None	17.8	-76,2	Length of Water Column: 39.01
	1105	10.71	0.6	6.60	0.535	12.7	42.61	NONE	17.7	-76.9	Depth of Well: 49,62
	1110	10.71	0.8	6.59	0.534	12.8	38.62	NUNE	17.5	-75.0	Sheen Observed: Y (N)
1	1115	10.71	1.0	6.59	01534	12.7	33.35	NONÉ	17.3	-75.2	DNAPL Observed: Y (N)
	- Info									,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,	Did Well Go Dry: Y/N
											Other:
		Ĺ								L	

	- N. I.	5 1.0				Historic Info	rmation		E 70.00				
Boring Log Available (yes/no/attached):													
Installation Le	og Availab	le ( <b>yes</b> /no/a	attached)										
						Summ	ary						
Monitoring W		MW-2 I			rface Elevation				een Materia				
Installation D		2/87			Casing Elevation				creen Depth				
Installed By:		Rochester		~	Point Elevation	: 1540.87 ft.		Bottom o	f Screen De	pth: 49 ft.			
				Elevation D									
Previous Fiel	ld measure	ement Inforr	nation Availa	ble (yes/ <b>no</b> /		(5)							
- D (I :				0 10 1		of Previous Fi							
Depth to			)H	•	Conductance	Tempera		ı	bidity		Color		
(ft)			ard Units)		nos/cm)	(°C)			ITU)		01		
14.6	51	7	.44	0	.593	12		1	.91		Clear		
Notes:			_										
			Fi.	Id Observe	tiono				Doron	ntoro 1/	Carralina Information		
Exterior Obs	onuntional			eld Observa	tions				pH	eters +/-	Sampling Information		
Exterior Obs	ervations:	Well co	wer mis	Sing						+/- 0.1	Sample ID: MW-21 - 691421 Sample Time: 1315		
Interior Obse	ervations	6.00											
	51 <b>+ 44101113</b>	0.000	Turbidity +/- 10%   # or Sample Containers: 3										
									ORP		Sample Analysis: VOCs 8260		
Signs of Dan	nage/Tam <sub>l</sub>	pering:							DO	+/- 10%			
Locked ()	yes/no)	Well Ca	p (yes)no)	Surfa	ace Seal Intact	(yes/no)	PID Meas	urement:	0.0000	Odors: 🗸	NE		
						Well Quali	ty Data						
Б.	<b>T</b> .					_			_				
Date	Time	Depth to	Cumulative	pH	Specific	Temperature	Turbidity	Color	Dissolved	Oxygen	Notes		
		Water	Volume	(Standard		(°C)	(NTU)		Oxygen	Reduction			
9-16-21	1245	ft bgs	Purged	Units) フロス	(uMhos/cm)	tu =	7 07	A1 -	81.0	Potential	Donth of Water: 15 3/1		
1-14-21	1295	15.27	0.2	6.54	0.157	14.7	3.02	NONE		177.0	Depth of Water: 15,24 Length of Water Column: 31,42		
	1255	15.28	0.4	6.26	0.152	13.2	25, 32	NONE	63.2	201.6	Depth of Well: 46.66		
	1300	15.28	0.6	6.16	0.132	13.3	26,21	None	72,7	205,6	Sheen Observed: Y(N)		
	1305	15,28	0.8	6.14	01144	13.3	10.93	NONE	71.4	209.8	DNAPL Observed: YO		
	1310	15,28	1.0	6:18	0,148	13.1	11,12	None	68.4	205.9	Did Well Go Dry: YCN		
7	1315	15.20	1.2	6,20	6.148	13,/	10.36	NONE	66.7	210.7	Other:		
	L												

						Historic Info	ormation					
Boring Log Available ( <b>yes</b> /no/attached):												
Installation L	og Availab	le ( <b>yes</b> /no/a	attached)									
		= ::				Summ	ary					
Monitoring W	/ell:	MW-4 S		Ground Su	rface Elevation	:		Riser/Scr	een Materia	al: PVC		
Installation D	ate:	11/86		Protective (	Casing Elevation	n: 1535.47 ft.			creen Depth			
Installed By:		Rochester	Drilling Co.	Monitoring	Point Elevation	: 1535.42 ft.		Bottom o	f Screen De	pth: 17 ft.		
				Elevation D	atum:							
Previous Fie	ld measure	ement Inforr	mation Availal	ble (yes/ <b>no</b> /	attached)							
					Ranges	of Previous F	ield Measu	rements				
Depth to	Depth to Water pH Specific Conductance Temperature Turbidity Color											
(ft)		(Standa	ard Units)	· (uMl	nos/cm)	(°C)		(N	ITU)			
7.9		6.1 0.28 16.7 1.36 Clear									Clear	
Notes:												
1000			Fie	eld Observa	tions			E in an	Parame	eters +/-	Sampling Information	
Exterior Obs	ervations:	6000							pН	+/- 0.1	Sample ID: MW-4S - 091521	
									Conductivit	y +/- 3%	Sample Time: 1400	
Interior Obse	erior Observations Read box Freeded # of Sample Containers: 3											
	Turbidity +/- 10% Duplicate Sample ID:											
									ORP		Sample Analysis: VOCs 8260	
Signs of Dan									DO	+/- 10%		
Locked (	/e³/no)	Well Ca	p (yes/no)	Surfa	ace Seal Intact			urement:	سواؤه.ه	Odors: 🖊 🤋	Me.	
						Well Quali	ity Data					
Date	Time	Depth to	Cumulative	рН	Specific	Temperature	Turbidity	Color	Dissolved	Oxygen	Notes	
		Water	Volume	(Standard	Conductance	(°C)	(NTU)		Oxygen	Reduction		
		ft bgs	Purged	Units)	(uMhos/cm)					Potential		
9-15-21	1345	10,19	0	6.89	0.289	16.1	2.25	NaNE	66.3	90.3	Depth of Water: 9.95	
	1356	10.21	0.2	6.39	0.270	15.8	1.86	NENE	59.7	112,4	Length of Water Column: 6 . 2 %	
	1355	10,21	0.7	6.39	0.270	15.7	2.02	NONE	38,3	121,7	Depth of Well: /6,23	
V	1400	10.21	0.6	6,36	0:272	15.7	2,14	NONE	56.7	128.4	Sheen Observed: Y (N)	
											DNAPL Observed: Y W	
								2			Did Well Go Dry: Y Y	
											Other:	
		-										
									-			
			I									

	Historic Information											
Boring Log Available ( <b>yes</b> /no/attached):												
Installation L	og Availab	le (yes/no/a	attached)									
						Summ	ary					
Monitoring W	/ell:	MW-5 S		Ground Sur	face Elevation			Riser/Scr	een Materia	al: PVC		
Installation D		11/86			Casing Elevation				creen Depth			
Installed By:		Rochester			Point Elevation	: 1534.16 ft.		Bottom of	f Screen De	pth: 17.5 ft.		
				Elevation D	atum;							
Previous Fiel	ld measure	ement Inforr	mation Availal	ble (yes/ <b>no</b> /	attached)							
					Ranges	s of Previous Fi	eld Measu	rements				
Depth to	Water	F	рΗ	Specific C	Conductance	Tempera	iture	Tur	bidity		Color	
(ft)		(Standa	ard Units)	(uMł	nos/cm)	( °C)		(N	(NTU)			
8.96	6	7	.97	0	.595	13.9		3	.12		Clear	
Notes:												
	1111		Fie	eld Observa	tions				Parame	eters +/-	Sampling Information	
Exterior Obs	ervations:	6000							pН	+/- 0.1	Sample ID: MW-5S - 091621	
			- 1						Conductivit		Sample Time: 1210	
Interior Obse	ervations	Roadbox	Flooded								# of Sample Containers: 3	
	Turbidity +/- 10% Duplicate Sample ID:											
	ORP +/- 10mV Sample Analysis: VOCs 8260											
Signs of Dan									DO	+/- 10%		
Locked ()	(es/no)	Well Ca	p (yes)no)	Surfa	ace Seal Intact		PID Measu	urement:	0.079~	Odors: N	°NE	
						Well Quali	ty Data					
Date	Time	Donth to	Cumulative	рН	Specific	Temperature	Turbidity	Color	Dissolved	Oxygen	Notes	
Date	rine	Depth to Water	Volume	(Standard	Conductance		(NTU)	COIOI		Reduction	Notes	
1				Units)	(uMhos/cm)	(°C)	(1410)		Oxygen	Potential		
B. 17 . 2 1	11-6	ft bgs	Purged	6.66		16,9	4011	Al-al-			Depth of Water: 역, θ 7	
9-16-21	1155	9,27	0.2	6.34	0.305	16-8	3.86	NONE	53.2	164.1	Length of Water Column: 6.03	
	1205	9.40	0,4	6.34	0,369	16.9	4.21	News	52.3	184.4	Depth of Well: /5,10	
	1210	9,40	0.6	6.34	0,369	16.9	4.36	NONE	52.1	191.8	Sheen Observed: Y (N)	
``											DNAPL Observed: Y N>	
											Did Well Go Dry: Y (N)	
											Other:	
		L										

Historic Information											
Boring Log Available (yes/ <b>no</b> /attached):											
Installation Lo	og Availab	le ( <b>yes</b> /no/a	ittached)								
						Summa					
Monitoring W		MW-9 I			face Elevation				een Materia		
Installation D		1/87			Casing Elevation				creen Depth		
Installed By:		Rochester I			Point Elevation	: 1532.3 ft.		Bottom of	f Screen De	pth: 49.5 ft.	
				Elevation D							
Previous Fiel	ld measure	ement Inforr	mation Availal	ole (yes/no/							
						of Previous Fi					
Depth to			Н		Conductance	Tempera	- 1		bidity		Color
(ft)			ard Units)		nos/cm)	( °C)			ITU)		<u> </u>
10.2	3	7	.54	0	.618	13.5		6	.37		Clear
Notes:											
	1 1110		Fie	eld Observa	tions					eters +/-	Sampling Information
Exterior Obse	ervations:	600D							pH	+/- 0.1	Sample ID: MW-9 I - 0915Z1
	- 1								Conductivi		Sample Time: $950$
Interior Obse											
	9	Turbidity +/- 10% Duplicate Sample ID: ORP +/- 10mV Sample Analysis: VOCs 8260									
Signs of Dan	nago/Tami	ooring: Alex	A.F						DO	+/- 10111	Sample Analysis. VOCs 6200
Locked (s			p (yes/no)	Surfa	ace Seal Intact	(ves/no)	PID Measu			Odors. 🗚	2U22
LOCKED (S	(C3/110)	Well oa	p (yes/no)	Ourie	dec ocar irradi	Well Quali		ar official.		Odoro: 7	
						77011 QUOI	.,				
Date	Time	Depth to	Cumulative	pН	Specific	Temperature	Turbidity	Color	Dissolved	Oxygen	Notes
Dute		Water	Volume	(Standard	Conductance	(°C)	(NTU)		Oxygen	Reduction	
		ft bgs	Purged	Units)	(uMhos/cm)		, ,		, ,	Potential	
9-15-21	0930	10,9	0	7.49	0.633	16 x Z	24,94	NONE	45. 3	139.5	Depth of Water: 10,88
1	0935	10,9	0.2	7.33	0.633	141.1	39.52	None	32.6	132,6	Length of Water Column: 36.24
	0940	10.9	0.4	7.30	0.633	14.1	45,72	NONE	31.9	132.7	Depth of Well: 47, 12
	0945	10.9	0.6	7.27	0.633	14.2	43,49	None	31:4	133.7	Sheen Observed: Y (N)
J	0950	10.9	0.8	7.26	0.633	14,4	45.17	None	31.2	134.0	DNAPL Observed: Y N
											Did Well Go Dry: Y (N.)
			-								Other:

18318-	J 3 K					Historic Info	rmation				
Boring Log Available (yes/ <b>no</b> /attached):											
Installation L	og Availab	le (yes/no/a	attached)								
						Summa					
Monitoring W		IRM-1			face Elevation						inless Steel
Installation D		1990			Casing Elevation				creen Depth		
Installed By:		Empire Soi			Point Elevation	: 1534.75 ft,		Rottom o	f Screen De	ptn: 50 ft.	
	.,			Elevation D							
Previous Fiel	id measure	ement Inforr	mation Availat	ole (yes/ <b>no</b> /		of Decylone C	old Mogari	romonto			
D (1)	147			0:6:- 0		of Previous Fi			hidity I		Color
Depth to			oH		Conductance	Tempera			bidity		Coloi
(ft)			ard Units)		nos/cm)	(°C) 11.1			ITU)		Clear
22.3	12	/	.48	0	.652	11.1		8	). <i>I</i> Z		Cleal
Notes:											
			Eic	eld Observa	tions		-1000		Parame	eters +/-	Sampling Information
Exterior Obs	onvations	102 I-		d Observa	lions				Ha		Sample ID: IRM-1 - 0916 21
LYIGHOL ODS	ci valiui is.	NO 100							Conductivit		Sample Time: / o Z 5
Interior Obse	ior Observations Temperature +/- 10% # of Sample Containers: 6										
Turbidity +/- 10% Duplicate Sample ID: 6w Du										Duplicate Sample ID: ເພ DUP	
									ORP		Sample Analysis: VOCs 8260
Signs of Dan						,			DO	+/- 10%	
Locked (y	yes/no)	Well Ca	p (ýes/no)	Surfa	ace Seal Intact			urement:	O, Oppin	Odors: ^	10A1E
	17 17					Well Quali	ty Data				
Doto	Time	Donth to	Cumulative	pН	Specific	Temperature	Turbidity	Color	Dissolved	Oxygen	Notes
Date	Time	Depth to Water	Volume	(Standard	Conductance	(°C)	(NTU)	COIOI	Oxygen	Reduction	140(63
		ft bgs	Purged	Units)	(uMhos/cm)		(1410)		DAYGEIT	Potential	
9-16-21	0945	23.79	O	7,70	0.654	11,3	145.67	NONE	24.3	-84.6	Depth of Water: 23,76
1.0(1	1015	23.78	10	7126	0.666	/1,2	40.7	NONE	20.4	13.0	Length of Water Column: 21.24
	1020	23.78	1.1	7.26	0.666	11:2	3891	NONE	20.4	19.1	Depth of Well: 45.00
Ψ	1925	23158	. 12	7,26	0.666	[1,1	니기교내	None	20.4	15,5	Sheen Observed: Y (N)
											DNAPL Observed: Y (N)
											Did Well Go Dry: Y /N
											Other:
				-							
		100511 114									Page: 1 of 1

EE ST						Historic Info	rmation				
Boring Log Available (yes/ <b>no</b> /attached):											
Installation L	og Availab	le (yes/ <mark>no</mark> /a	attached)								
						Summa					
Monitoring W		IRM-2 I			face Elevation				70° C 70° 1 C 7	al: Steel/Sta	inless Steel
Installation D		1990			Casing Elevation				creen Depth		
Installed By:		Empire Soi			Point Elevation	: 1535.99 ft.		Bottom o	f Screen De	pth: 50 ft.	
				Elevation D							
Previous Fie	ld measure	ement Inforr	mation Availat	ole (yes/no/							
						of Previous Fi					
Depth to		·	oH		Conductance	Tempera	- 1		bidity		Color
(ft)			ard Units)		nos/cm)	(°C)					
22.6	2	7	.84	0	.547	10.9		2	2.8		
Notes:											
									D		Oline lefe-mation
	0.00			eld Observa	tions					eters +/- +/- 0.1	Sampling Information Sample ID: IRM-21 = 09162/
Exterior Obs	ervations:	No lo	ch						pH Conductivit		Sample Time: a9p 5
		V									# of Sample Containers: 3
Interior Obse	ervations	Temperature +/- 10% # of Sample Containers: 3  Turbidity +/- 10% Duplicate Sample ID:									
									ORP		Sample Analysis: VOCs 8260
Signs of Dar	nage/Tami	pering.							DO	+/- 10%	Cample / maryole. V C C C C C C C
Locked ()			p (yestno)	Surfa	ace Seal Intact	(ves/no)	PID Meas	urement:		Odors:	
						Well Quali					
Date	Time	Depth to Water	Cumulative Volume	pH (Standard	Specific Conductance	Temperature (°C)	Turbidity (NTU)	Color	Dissolved Oxygen	Oxygen Reduction	Notes
		ft bgs	Purged	Units)	(uMhos/cm)	( ),	(		,5	Potential	
9-16-21	0925	241.13	D.	10.63	0.335	11.10	28,15	NONE	30.5	48,3	Depth of Water: 23.98
T I	0855	24.58	8	7.33	0.566	11.0	18.54	None	23.7	-17.2	Length of Water Column: 17 (1)
	0900	24,58	9	7,33	0.566	11.0	10.75	Now	23.6	-16.5	Depth of Well: 44.09
$\rightarrow$	0905	24,58	10	7,32	0.566	11.1	9.76	NONE	23.6	-13.7	Sheen Observed: Y (N)
											DNAPL Observed: Y N
											Did Well Go Dry: Y (N)
											Other:
										L	

	-5.					Historic Info	rmation			100	
Boring Log Availa											
Installation Log A	vailabl	e (yes/no/a	attached)								
						Summa					
Monitoring Well:		Town Well			face Elevation						inless Steel
Installation Date:	- 3	1982			Casing Elevation				creen Depth		
Installed By:					Point Elevation	:		Bottom o	f Screen De	pth: 51.5 ft.	
				Elevation D							
Previous Field me	easure	ment Inforr	mation Availal	ole (yes/ <b>no</b> /							
					Ranges	of Previous Fi	eld Measu	rements			
Depth to Wate	er	ŗ	оН	Specific C	Conductance	Tempera	ture	Tur	bidity		Color
(ft)		(Standa	ard Units)	(uMł	nos/cm)	( °C)		(N	ITU)		
NA		7	.32	0	.566	16.2		2	.02		Clear
Notes:	lotes:										
		46.6	Fie	eld Observa	tions			350.00	Parame	eters +/-	Sampling Information
Exterior Observa	tions:	Good							pН	+/- 0.1	Sample ID: TOWN WELL -091621
	×-	- V							Conductivit		Sample Time: 0925
Interior Observations Good Temperature +/- 10% # of Sample Containers: 3											
Turbidity +/- 10% Duplicate Sample ID:											
	# <del>5</del>								ORP		Sample Analysis: VOCs 8260
Signs of Damage	e/Tamp								DO	+/- 10%	
Locked (yes/r	no)	Well Ca	p (yes/no)	Surfa	ace Seal Intact	(yes/no) NA		urement:	NA	Odors: N	ONE
		التنائل				Well Quali	ty Data			1 = W1-	
					V 19	_					N
Date Ti	ime	Depth to	Cumulative	рΗ	Specific	Temperature	Turbidity	Color	Dissolved	Oxygen	Notes
		Water	Volume	,		(°C)	(NTU)		Oxygen	Reduction	
		ft bgs	Purged	Units)	(uMhos/cm)					Potential	D # 61W .
9-16-21 09	25	NΑ	5	7.66	0.603	15.7	3,17	NONE	54.5	625.5	Depth of Water: A
											Length of Water Column:   A
											Depth of Well: NA Sheen Observed: Y (N)
											DNAPL Observed: Y
			ļ								Did Well Go Dry: Y /N
<b></b>											Other:
											Purged 5 gal then
											400h SIND SOUNDLY

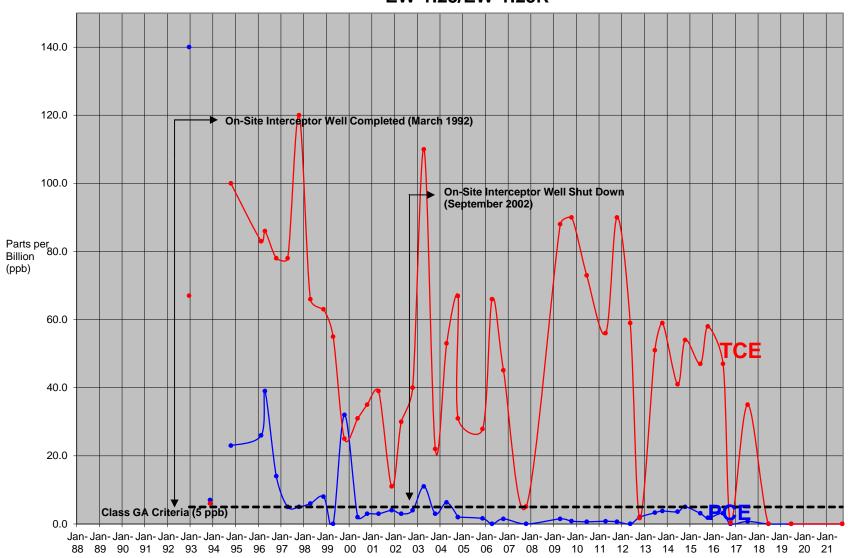


### **APPENDIX C**

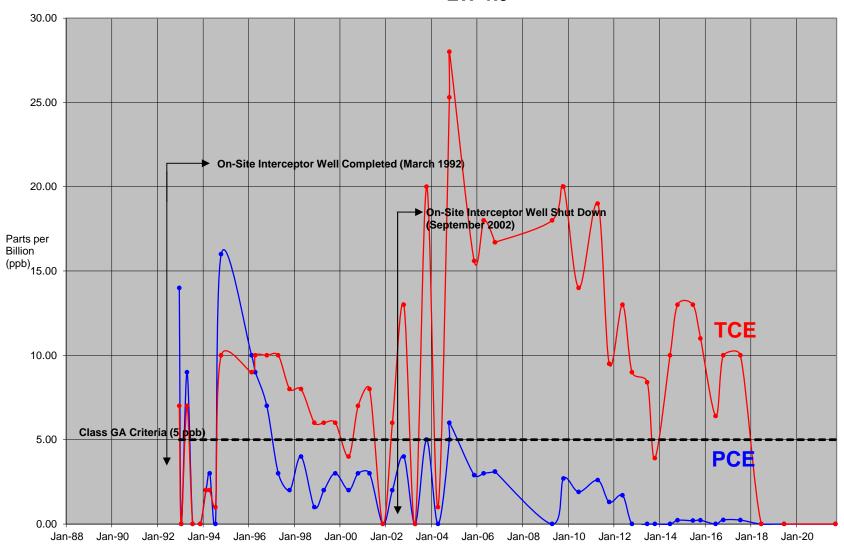
TCE AND PCE CONCENTRATION GRAPHS

### TCE and PCE Groundwater Concentrations

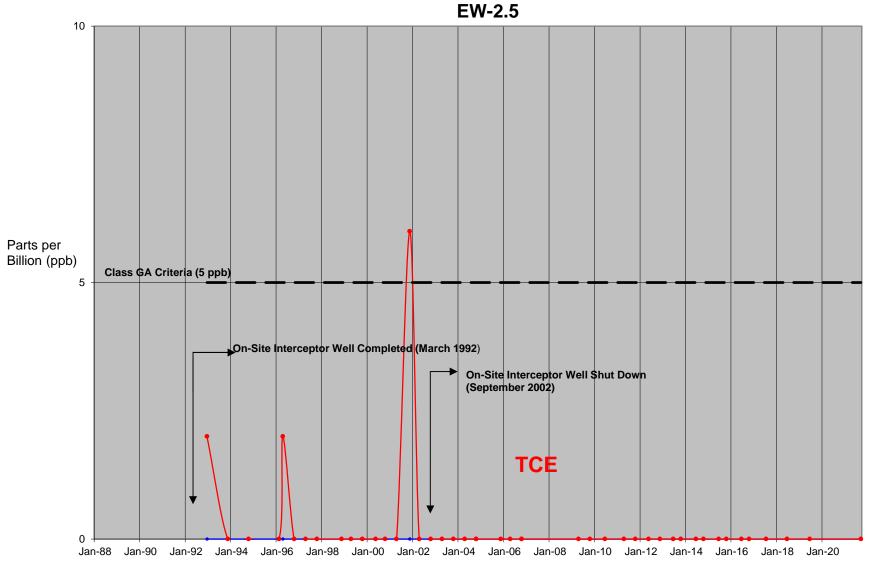
### EW-1.25/EW-1.25R



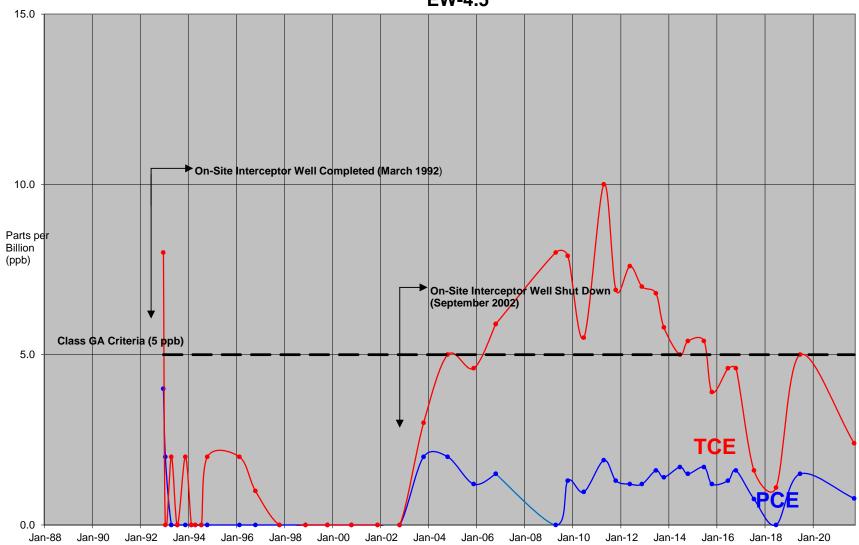
TCE and PCE Groundwater Concentrations **EW-1.5** 



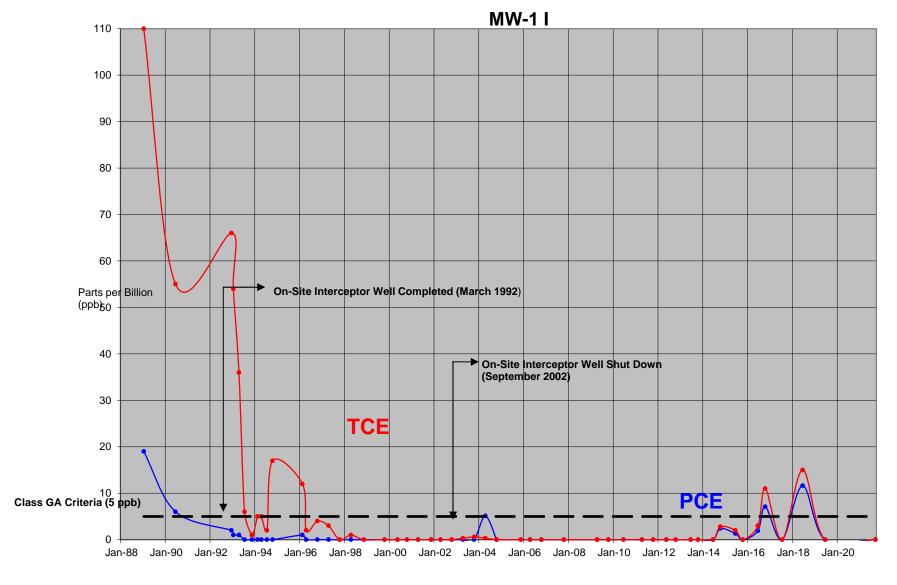
### TCE and PCE Groundwater Concentrations



TCE and PCE Groundwater Concentrations **EW-4.5** 

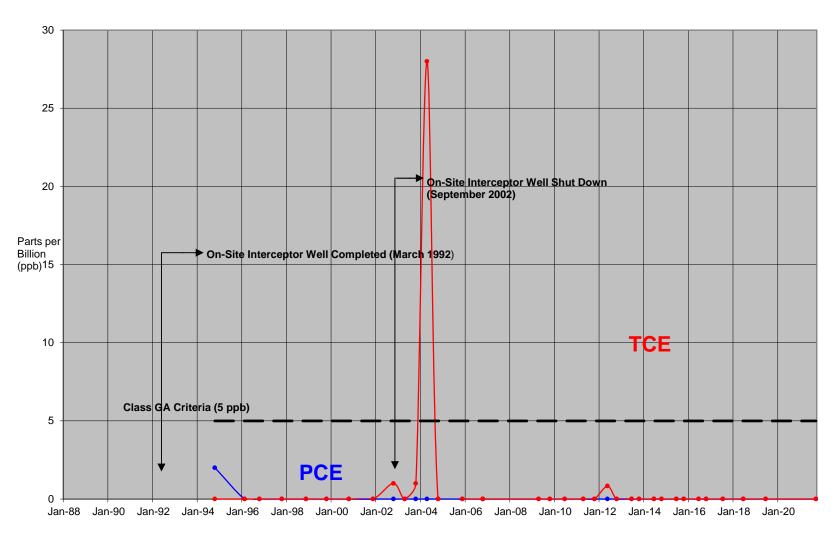


#### **TCE and PCE Groundwater Concentrations**

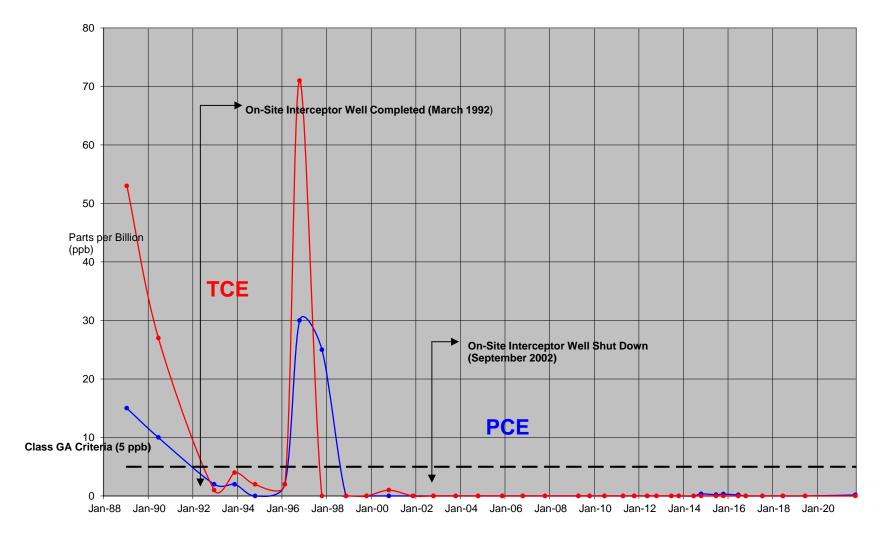


TCE and PCE Groundwater Concentrations

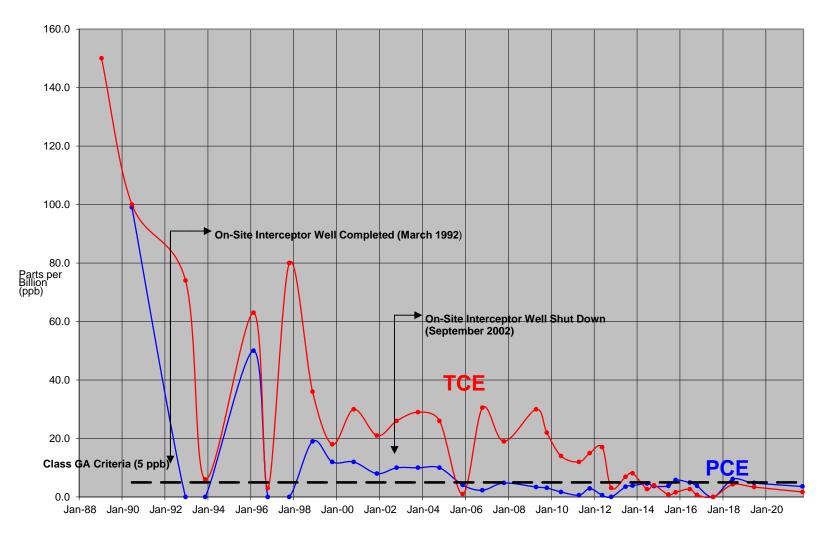
MW-2 I



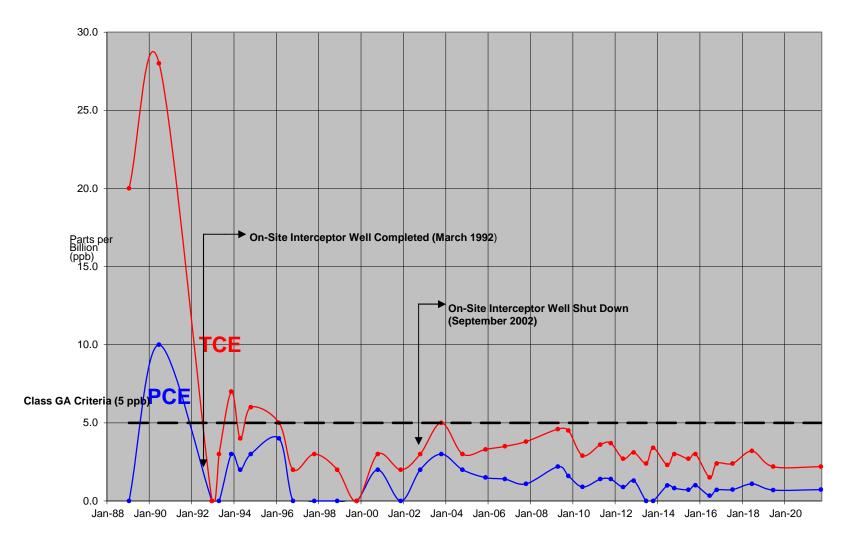
## TCE and PCE Groundwater Concentrations MW-4S



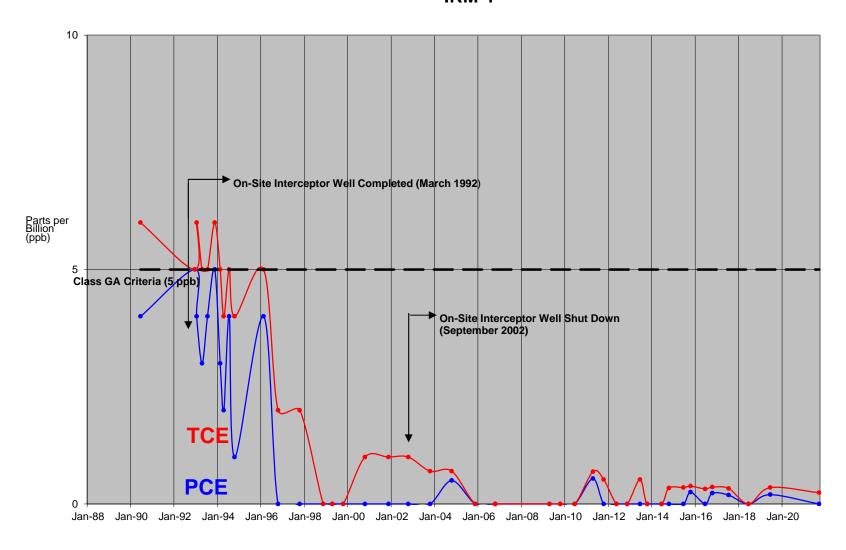
# TCE and PCE Groundwater Concentrations MW-5S



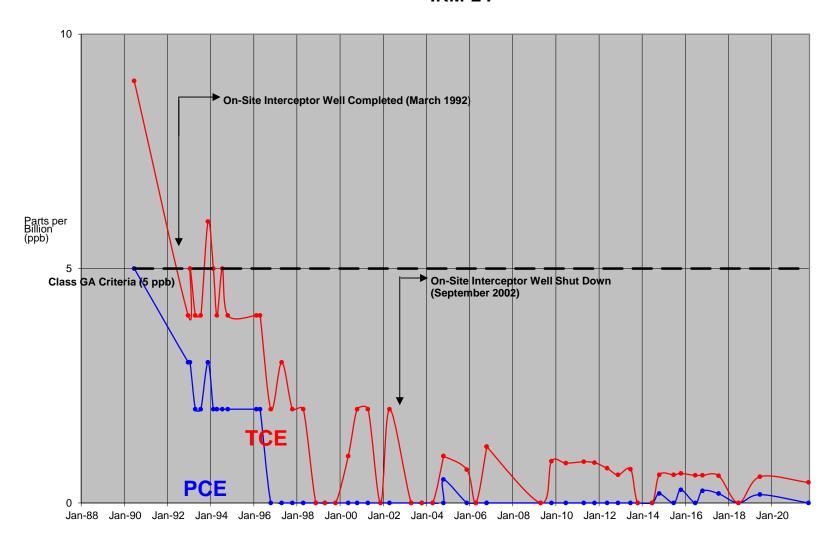
# TCE and PCE Groundwater Concentrations MW-9 I



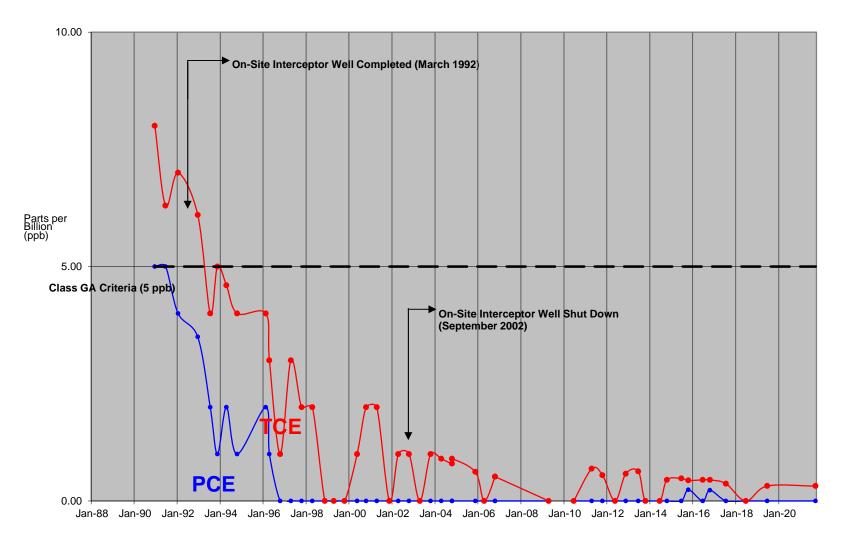
# TCE and PCE Groundwater Concentrations IRM-1



# TCE and PCE Groundwater Concentrations IRM-2 I



# TCE and PCE Groundwater Concentrations TOWN WELL





### **APPENDIX D**

**ANALYTICAL TEST RESULTS** 



#### ANALYTICAL REPORT

Lab Number: L2150447

Client: GZA GeoEnvironmental of New York

300 Pearl Street

Suite 700

Buffalo, NY 14202

ATTN: Thomas Bohlen Phone: (716) 844-7050

Project Name: SIGNORE POST INJECTION

Project Number: 21.0056367.66

Report Date: 09/24/21

The original project report/data package is held by Alpha Analytical. This report/data package is paginated and should be reproduced only in its entirety. Alpha Analytical holds no responsibility for results and/or data that are not consistent with the original.

Certifications & Approvals: MA (M-MA086), NH NELAP (2064), CT (PH-0574), IL (200077), ME (MA00086), MD (348), NJ (MA935), NY (11148), NC (25700/666), PA (68-03671), RI (LAO00065), TX (T104704476), VT (VT-0935), VA (460195), USDA (Permit #P330-17-00196).

Eight Walkup Drive, Westborough, MA 01581-1019 508-898-9220 (Fax) 508-898-9193 800-624-9220 - www.alphalab.com



Serial\_No:09242110:27

Project Name: SIGNORE POST INJECTION

**Project Number:** 21.0056367.66

Lab Number:

L2150447

Report Date:

09/24/21

Alpha Sample ID	Client ID	Matrix	Sample Location	Collection Date/Time	Receive Date
L2150447-01	EW-1.25R-091721	WATER	ELLICOTVILLE, NY	09/17/21 12:20	09/17/21



Project Name: SIGNORE POST INJECTION Lab Number: L2150447

Project Number: 21.0056367.66 Report Date: 09/24/21

#### **Case Narrative**

The samples were received in accordance with the Chain of Custody and no significant deviations were encountered during the preparation or analysis unless otherwise noted. Sample Receipt, Container Information, and the Chain of Custody are located at the back of the report.

Results contained within this report relate only to the samples submitted under this Alpha Lab Number and meet NELAP requirements for all NELAP accredited parameters unless otherwise noted in the following narrative. The data presented in this report is organized by parameter (i.e. VOC, SVOC, etc.). Sample specific Quality Control data (i.e. Surrogate Spike Recovery) is reported at the end of the target analyte list for each individual sample, followed by the Laboratory Batch Quality Control at the end of each parameter. Tentatively Identified Compounds (TICs), if requested, are reported for compounds identified to be present and are not part of the method/program Target Compound List, even if only a subset of the TCL are being reported. If a sample was re-analyzed or re-extracted due to a required quality control corrective action and if both sets of data are reported, the Laboratory ID of the re-analysis or re-extraction is designated with an "R" or "RE", respectively.

When multiple Batch Quality Control elements are reported (e.g. more than one LCS), the associated samples for each element are noted in the grey shaded header line of each data table. Any Laboratory Batch, Sample Specific % recovery or RPD value that is outside the listed Acceptance Criteria is bolded in the report. In reference to questions H (CAM) or 4 (RCP) when "NO" is checked, the performance criteria for CAM and RCP methods allow for some quality control failures to occur and still be within method compliance. In these instances, the specific failure is not narrated but noted in the associated QC Outlier Summary Report, located directly after the Case Narrative. QC information is also incorporated in the Data Usability Assessment table (Format 11) of our Data Merger tool, where it can be reviewed in conjunction with the sample result, associated regulatory criteria and any associated data usability implications.

Soil/sediments, solids and tissues are reported on a dry weight basis unless otherwise noted. Definitions of all data qualifiers and acronyms used in this report are provided in the Glossary located at the back of the report.

HOLD POLICY - For samples submitted on hold, Alpha's policy is to hold samples (with the exception of Air canisters) free of charge for 21 calendar days from the date the project is completed. After 21 calendar days, we will dispose of all samples submitted including those put on hold unless you have contacted your Alpha Project Manager and made arrangements for Alpha to continue to hold the samples. Air canisters will be disposed after 3 business days from the date the project is completed.

Please contact Project Management at 800-624-9220 with any questions.	



Serial\_No:09242110:27

Project Name:SIGNORE POST INJECTIONLab Number:L2150447Project Number:21.0056367.66Report Date:09/24/21

### **Case Narrative (continued)**

Report Submission

All non-detect (ND) or estimated concentrations (J-qualified) have been quantitated to the limit noted in the MDL column.

Sample Receipt

The project number was specified by the client.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Authorized Signature:

Title: Technical Director/Representative Date: 09/24/21

Custen Walker Cristin Walker

### **ORGANICS**



### **VOLATILES**



Serial\_No:09242110:27

L2150447

09/17/21 12:20

**Project Name:** SIGNORE POST INJECTION

**Project Number:** 21.0056367.66

**SAMPLE RESULTS** 

Report Date: 09/24/21

Lab Number:

Date Collected:

Lab ID: L2150447-01

Client ID: Date Received: 09/17/21 EW-1.25R-091721 Not Specified

Sample Location: Field Prep: ELLICOTVILLE, NY

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 09/22/21 15:18

Analyst: AJK

	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westboroug	jh Lab					
Methylene chloride	ND		ug/l	2.5	0.70	1
1,1-Dichloroethane	1.2	J	ug/l	2.5	0.70	1
Chloroform	ND		ug/l	2.5	0.70	1
Carbon tetrachloride	ND		ug/l	0.50	0.13	1
1,2-Dichloropropane	ND		ug/l	1.0	0.14	1
Dibromochloromethane	ND		ug/l	0.50	0.15	1
1,1,2-Trichloroethane	ND		ug/l	1.5	0.50	1
Tetrachloroethene	ND		ug/l	0.50	0.18	1
Chlorobenzene	ND		ug/l	2.5	0.70	1
Trichlorofluoromethane	ND		ug/l	2.5	0.70	1
1,2-Dichloroethane	ND		ug/l	0.50	0.13	1
1,1,1-Trichloroethane	ND		ug/l	2.5	0.70	1
Bromodichloromethane	ND		ug/l	0.50	0.19	1
trans-1,3-Dichloropropene	ND		ug/l	0.50	0.16	1
cis-1,3-Dichloropropene	ND		ug/l	0.50	0.14	1
Bromoform	ND		ug/l	2.0	0.65	1
1,1,2,2-Tetrachloroethane	ND		ug/l	0.50	0.17	1
Benzene	ND		ug/l	0.50	0.16	1
Toluene	ND		ug/l	2.5	0.70	1
Ethylbenzene	ND		ug/l	2.5	0.70	1
Chloromethane	ND		ug/l	2.5	0.70	1
Bromomethane	ND		ug/l	2.5	0.70	1
Vinyl chloride	0.17	J	ug/l	1.0	0.07	1
Chloroethane	ND		ug/l	2.5	0.70	1
1,1-Dichloroethene	ND		ug/l	0.50	0.17	1
trans-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1
Trichloroethene	ND		ug/l	0.50	0.18	1
1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1



Serial\_No:09242110:27

Project Name: SIGNORE POST INJECTION Lab Number: L2150447

**Project Number:** 21.0056367.66 **Report Date:** 09/24/21

SAMPLE RESULTS

Lab ID: L2150447-01 Date Collected: 09/17/21 12:20

Client ID: EW-1.25R-091721 Date Received: 09/17/21 Sample Location: ELLICOTVILLE, NY Field Prep: Not Specified

Sample Depth:

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor	
Volatile Organics by GC/MS - We	estborough Lab						
1,3-Dichlorobenzene	ND		ug/l	2.5	0.70	1	
1,4-Dichlorobenzene	ND		ug/l	2.5	0.70	1	
Methyl tert butyl ether	ND		ug/l	2.5	0.70	1	
p/m-Xylene	ND		ug/l	2.5	0.70	1	
o-Xylene	ND		ug/l	2.5	0.70	1	
cis-1,2-Dichloroethene	2.5		ug/l	2.5	0.70	1	
Styrene	ND		ug/l	2.5	0.70	1	
Dichlorodifluoromethane	ND		ug/l	5.0	1.0	1	
Acetone	ND		ug/l	5.0	1.5	1	
Carbon disulfide	ND		ug/l	5.0	1.0	1	
2-Butanone	ND		ug/l	5.0	1.9	1	
4-Methyl-2-pentanone	ND		ug/l	5.0	1.0	1	
2-Hexanone	ND		ug/l	5.0	1.0	1	
Bromochloromethane	ND		ug/l	2.5	0.70	1	
1,2-Dibromoethane	ND		ug/l	2.0	0.65	1	
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70	1	
Isopropylbenzene	ND		ug/l	2.5	0.70	1	
1,2,3-Trichlorobenzene	ND		ug/l	2.5	0.70	1	
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70	1	
Methyl Acetate	ND		ug/l	2.0	0.23	1	
Cyclohexane	ND		ug/l	10	0.27	1	
1,4-Dioxane	ND		ug/l	250	61.	1	
Freon-113	ND		ug/l	2.5	0.70	1	
Methyl cyclohexane	ND		ug/l	10	0.40	1	

Tentativel <sup>1</sup>	y Identified Compounds	3
I CHILALIVCI	y lacillilea Collipoullas	,

No Tentatively Identified Compounds ND ug/l 1

Surrogate	% Recovery	Acceptance Qualifier Criteria
1,2-Dichloroethane-d4	106	70-130
Toluene-d8	102	70-130
4-Bromofluorobenzene	107	70-130
Dibromofluoromethane	104	70-130



Project Name: SIGNORE POST INJECTION Lab Number: L2150447

**Project Number:** 21.0056367.66 **Report Date:** 09/24/21

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 09/22/21 09:33

Analyst: NLK

arameter	Result	Qualifier Units	RL	MDL
olatile Organics by GC/MS -	· Westborough Lab	for sample(s):	01 Batch:	WG1549680-5
Methylene chloride	ND	ug/l	2.5	0.70
1,1-Dichloroethane	ND	ug/l	2.5	0.70
Chloroform	ND	ug/l	2.5	0.70
Carbon tetrachloride	ND	ug/l	0.50	0.13
1,2-Dichloropropane	ND	ug/l	1.0	0.14
Dibromochloromethane	ND	ug/l	0.50	0.15
1,1,2-Trichloroethane	ND	ug/l	1.5	0.50
Tetrachloroethene	ND	ug/l	0.50	0.18
Chlorobenzene	ND	ug/l	2.5	0.70
Trichlorofluoromethane	ND	ug/l	2.5	0.70
1,2-Dichloroethane	ND	ug/l	0.50	0.13
1,1,1-Trichloroethane	ND	ug/l	2.5	0.70
Bromodichloromethane	ND	ug/l	0.50	0.19
trans-1,3-Dichloropropene	ND	ug/l	0.50	0.16
cis-1,3-Dichloropropene	ND	ug/l	0.50	0.14
Bromoform	ND	ug/l	2.0	0.65
1,1,2,2-Tetrachloroethane	ND	ug/l	0.50	0.17
Benzene	ND	ug/l	0.50	0.16
Toluene	ND	ug/l	2.5	0.70
Ethylbenzene	ND	ug/l	2.5	0.70
Chloromethane	ND	ug/l	2.5	0.70
Bromomethane	ND	ug/l	2.5	0.70
Vinyl chloride	ND	ug/l	1.0	0.07
Chloroethane	ND	ug/l	2.5	0.70
1,1-Dichloroethene	ND	ug/l	0.50	0.17
trans-1,2-Dichloroethene	ND	ug/l	2.5	0.70
Trichloroethene	ND	ug/l	0.50	0.18
1,2-Dichlorobenzene	ND	ug/l	2.5	0.70
1,3-Dichlorobenzene	ND	ug/l	2.5	0.70



L2150447

Project Name: SIGNORE POST INJECTION Lab Number:

**Project Number:** 21.0056367.66 **Report Date:** 09/24/21

Method Blank Analysis
Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 09/22/21 09:33

Analyst: NLK

Parameter	Result	Qualifier Unit	ts RL	MDL
olatile Organics by GC/MS - W	estborough Lab	for sample(s):	01 Batch:	WG1549680-5
1,4-Dichlorobenzene	ND	ug	/I 2.5	0.70
Methyl tert butyl ether	ND	ug	/I 2.5	0.70
p/m-Xylene	ND	ug	/I 2.5	0.70
o-Xylene	ND	ug	/I 2.5	0.70
cis-1,2-Dichloroethene	ND	ug	/I 2.5	0.70
Styrene	ND	ug	/I 2.5	0.70
Dichlorodifluoromethane	ND	ug	/I 5.0	1.0
Acetone	ND	ug	/I 5.0	1.5
Carbon disulfide	ND	ug	/I 5.0	1.0
2-Butanone	ND	ug	/I 5.0	1.9
4-Methyl-2-pentanone	ND	ug	/I 5.0	1.0
2-Hexanone	ND	ug	/I 5.0	1.0
Bromochloromethane	ND	ug	/I 2.5	0.70
1,2-Dibromoethane	ND	ug	/I 2.0	0.65
1,2-Dibromo-3-chloropropane	ND	ug	/I 2.5	0.70
Isopropylbenzene	ND	ug	/I 2.5	0.70
1,2,3-Trichlorobenzene	ND	ug	/I 2.5	0.70
1,2,4-Trichlorobenzene	ND	ug	/I 2.5	0.70
Methyl Acetate	ND	ug	/I 2.0	0.23
Cyclohexane	ND	ug	/I 10	0.27
1,4-Dioxane	ND	ug	/I 250	61.
Freon-113	ND	ug	/I 2.5	0.70
Methyl cyclohexane	ND	ug	/I 10	0.40

Tentatively Identified Compounds

No Tentatively Identified Compounds ND ug/l



Project Name: SIGNORE POST INJECTION Lab Number: L2150447

**Project Number:** 21.0056367.66 **Report Date:** 09/24/21

Method Blank Analysis
Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 09/22/21 09:33

Analyst: NLK

Parameter Result Qualifier Units RL MDL

Volatile Organics by GC/MS - Westborough Lab for sample(s): 01 Batch: WG1549680-5

		Acceptance
Surrogate	%Recovery Q	ualifier Criteria
1,2-Dichloroethane-d4	99	70-130
Toluene-d8	103	70-130
4-Bromofluorobenzene	109	70-130
Dibromofluoromethane	102	70-130



## Lab Control Sample Analysis Batch Quality Control

**Project Name:** SIGNORE POST INJECTION

**Project Number:** 21.0056367.66

Lab Number: L2150447

**Report Date:** 09/24/21

Methylene chloride	Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	RPD Qual Limits	
1,1-Dichloroethane         120         120         70-130         0         20           Chloroform         100         100         70-130         0         20           Carbon tetrachloride         96         92         63-132         4         20           1,2-Dichloropropane         120         110         70-130         9         20           Dibromochloromethane         91         88         63-130         3         20           1,1,2-Tichloroethane         100         100         70-130         0         20           Tetrachloroethane         98         95         70-130         3         20           Chlorobenzene         100         99         75-130         1         20           Trichlorofluoromethane         100         95         62-150         5         20           1,1,1-Trichloroethane         100         99         67-130         1         20           1,1,1-Trichloroethane         100         99         67-130         1         20           Bromodichloromethane         98         96         67-130         1         20           Itrans-1,3-Dichloropropene         100         98         70-130         0	Volatile Organics by GC/MS - Westborough L	ab Associated	sample(s): 01	Batch: WG1	549680-3	WG1549680-4			
Chloroform         100         100         70-130         0         20           Carbon tetrachloride         96         92         63-132         4         20           1,2-Dichloropropane         120         110         70-130         9         20           Dibromochloromethane         91         88         63-130         3         20           1,1,2-Trichloroethane         100         100         70-130         0         20           Tetrachloroethene         98         95         70-130         3         20           Chlorobenzene         100         99         75-130         1         20           Trichlorofluromethane         100         99         75-130         1         20           1,2-Dichloroethane         100         95         62-150         5         20           1,1-Trichloroethane         100         99         67-130         0         20           Irrans-1,3-Dichloropropene         100         99         67-130         2         20           Irrans-1,3-Dichloropropene         100         98         70-130         0         20           Bromoform         87         86         54-136         1	Methylene chloride	100		100		70-130	0	20	
Carbon tetrachloride         96         92         63-132         4         20           1,2-Dichloropropane         120         110         70-130         9         20           Dibromochloromethane         91         88         63-130         3         20           1,1,2-Trichloroethane         100         100         70-130         0         20           Tetrachloroethane         98         95         70-130         3         20           Chlorobenzene         100         99         75-130         1         20           Trichlorotuoromethane         100         95         62-150         5         20           1,2-Dichloroethane         100         100         70-130         0         20           1,1,1-Trichloroethane         100         99         67-130         1         20           Bromodichloromethane         98         96         67-130         2         20           trans-1,3-Dichloropropene         100         98         70-130         2         20           cis-1,3-Dichloropropene         100         100         70-130         0         20           Bromoform         87         86         54-136         1<	1,1-Dichloroethane	120		120		70-130	0	20	
1/2-Dichloropropane         120         110         70-130         9         20           Dibromochloromethane         91         88         63-130         3         20           1,1,2-Trichloroethane         100         100         70-130         0         20           Tetrachloroethane         98         95         70-130         3         20           Chlorobenzene         100         99         75-130         1         20           Trichlorofluoromethane         100         95         62-150         5         20           1,2-Dichloroethane         100         100         70-130         0         20           1,1,1-Trichloroethane         100         99         67-130         1         20           Bromodichloromethane         98         96         67-130         1         20           trans-1,3-Dichloropropene         100         98         70-130         2         20           trans-1,3-Dichloropropene         100         100         70-130         0         20           Bromoform         87         86         54-136         1         20           Benzene         100         100         67-130         0	Chloroform	100		100		70-130	0	20	
Dibromochloromethane         91         88         63-130         3         20           1,1,2-Trichloroethane         100         100         70-130         0         20           Tetrachloroethene         98         95         70-130         3         20           Chlorobenzene         100         99         75-130         1         20           Trichlorofluoromethane         100         95         62-150         5         20           1,2-Dichloroethane         100         100         70-130         0         20           1,1,1-Trichloroethane         100         99         67-130         1         20           Bromodichloromethane         98         96         67-130         1         20           trans-1,3-Dichloropropene         100         98         70-130         2         20           trans-1,3-Dichloropropene         100         100         70-130         0         20           is-1,3-Dichloropropene         100         100         70-130         0         20           Bromoform         87         86         54-136         1         20           1,1,2,2-Tetrachloroethane         100         100         70-130	Carbon tetrachloride	96		92		63-132	4	20	
1,1,2-Trichloroethane         100         100         70-130         0         20           Tetrachloroethene         98         95         70-130         3         20           Chlorobenzene         100         99         75-130         1         20           Trichlorofluoromethane         100         95         62-150         5         20           1,2-Dichloroethane         100         100         70-130         0         20           1,1,1-Trichloroethane         100         99         67-130         1         20           Bromodichloromethane         98         96         67-130         2         20           trans-1,3-Dichloropropene         100         98         70-130         2         20           cis-1,3-Dichloropropene         100         100         70-130         0         20           Bromoform         87         86         54-136         1         20           1,1,2,2-Tetrachloroethane         100         100         67-130         0         20           Benzene         100         110         70-130         10         20           Toluene         110         100         70-130         10	1,2-Dichloropropane	120		110		70-130	9	20	
Tetrachloroethene         98         95         70-130         3         20           Chlorobenzene         100         99         75-130         1         20           Trichlorofluoromethane         100         95         62-150         5         20           1,2-Dichloroethane         100         100         70-130         0         20           1,1,1-Trichloroethane         100         99         67-130         1         20           Bromodichloromethane         98         96         67-130         2         20           trans-1,3-Dichloropropene         100         98         70-130         2         20           cis-1,3-Dichloropropene         100         100         70-130         0         20           Bromoform         87         86         54-136         1         20           Benzene         100         100         67-130         0         20           Benzene         100         110         70-130         10         20           Ethylbenzene         110         100         70-130         10         20           Chloromethane         160         Q         150         Q         64-130 <t< td=""><td>Dibromochloromethane</td><td>91</td><td></td><td>88</td><td></td><td>63-130</td><td>3</td><td>20</td><td></td></t<>	Dibromochloromethane	91		88		63-130	3	20	
Chlorobenzene         100         99         75-130         1         20           Trichlorofluoromethane         100         95         62-150         5         20           1,2-Dichloroethane         100         100         70-130         0         20           1,1,1-Trichloroethane         100         99         67-130         1         20           Bromomethane         98         96         67-130         2         20           trans-1,3-Dichloropropene         100         98         70-130         2         20           cis-1,3-Dichloropropene         100         100         70-130         0         20           Bromoform         87         86         54-136         1         20           1,1,2,2-Tetrachloroethane         100         100         67-130         0         20           Benzene         100         110         70-130         10         20           Toluene         110         100         70-130         10         20           Ethylbenzene         110         100         70-130         10         20           Chloromethane         160         Q         150         Q         64-130	1,1,2-Trichloroethane	100		100		70-130	0	20	
Trichlorofluoromethane         100         95         62-150         5         20           1,2-Dichloroethane         100         100         70-130         0         20           1,1,1-Trichloroethane         100         99         67-130         1         20           Bromodichloromethane         98         96         67-130         2         20           trans-1,3-Dichloropropene         100         98         70-130         2         20           cis-1,3-Dichloropropene         100         100         70-130         0         20           Bromoform         87         86         54-136         1         20           1,1,2,2-Tetrachloroethane         100         100         67-130         0         20           Benzene         100         110         70-130         10         20           Toluene         110         100         70-130         10         20           Ethylbenzene         110         100         70-130         10         20           Chloromethane         84         81         39-139         4         20	Tetrachloroethene	98		95		70-130	3	20	
1,2-Dichloroethane       100       100       70-130       0       20         1,1,1-Trichloroethane       100       99       67-130       1       20         Bromodichloromethane       98       96       67-130       2       20         trans-1,3-Dichloropropene       100       98       70-130       2       20         cis-1,3-Dichloropropene       100       100       70-130       0       20         Bromoform       87       86       54-136       1       20         1,1,2,2-Tetrachloroethane       100       100       67-130       0       20         Benzene       100       110       70-130       10       20         Toluene       110       100       70-130       10       20         Ethylbenzene       110       100       70-130       10       20         Chloromethane       160       Q       150       Q       64-130       6       20         Bromomethane       84       81       39-139       4       20	Chlorobenzene	100		99		75-130	1	20	
1,1,1-Trichloroethane       100       99       67-130       1       20         Bromodichloromethane       98       96       67-130       2       20         trans-1,3-Dichloropropene       100       98       70-130       2       20         cis-1,3-Dichloropropene       100       100       70-130       0       20         Bromoform       87       86       54-136       1       20         1,1,2,2-Tetrachloroethane       100       100       67-130       0       20         Benzene       100       110       70-130       10       20         Toluene       110       100       70-130       10       20         Ethylbenzene       110       100       70-130       10       20         Chloromethane       160       Q       150       Q       64-130       6       20         Bromomethane       84       81       39-139       4       20	Trichlorofluoromethane	100		95		62-150	5	20	
Bromodichloromethane         98         96         67-130         2         20           trans-1,3-Dichloropropene         100         98         70-130         2         20           cis-1,3-Dichloropropene         100         100         70-130         0         20           Bromoform         87         86         54-136         1         20           1,1,2,2-Tetrachloroethane         100         100         67-130         0         20           Benzene         100         110         70-130         10         20           Toluene         110         100         70-130         10         20           Ethylbenzene         110         100         70-130         10         20           Chloromethane         160         Q         150         Q         64-130         6         20           Bromomethane         84         81         39-139         4         20	1,2-Dichloroethane	100		100		70-130	0	20	
trans-1,3-Dichloropropene         100         98         70-130         2         20           cis-1,3-Dichloropropene         100         100         70-130         0         20           Bromoform         87         86         54-136         1         20           1,1,2,2-Tetrachloroethane         100         100         67-130         0         20           Benzene         100         110         70-130         10         20           Toluene         110         100         70-130         10         20           Ethylbenzene         110         100         70-130         10         20           Chloromethane         160         Q         150         Q         64-130         6         20           Bromomethane         84         81         39-139         4         20	1,1,1-Trichloroethane	100		99		67-130	1	20	
cis-1,3-Dichloropropene         100         100         70-130         0         20           Bromoform         87         86         54-136         1         20           1,1,2,2-Tetrachloroethane         100         100         67-130         0         20           Benzene         100         110         70-130         10         20           Toluene         110         100         70-130         10         20           Ethylbenzene         110         100         70-130         10         20           Chloromethane         160         Q         150         Q         64-130         6         20           Bromomethane         84         81         39-139         4         20	Bromodichloromethane	98		96		67-130	2	20	
Bromoform       87       86       54-136       1       20         1,1,2,2-Tetrachloroethane       100       100       67-130       0       20         Benzene       100       110       70-130       10       20         Toluene       110       100       70-130       10       20         Ethylbenzene       110       100       70-130       10       20         Chloromethane       160       Q       150       Q       64-130       6       20         Bromomethane       84       81       39-139       4       20	trans-1,3-Dichloropropene	100		98		70-130	2	20	
1,1,2,2-Tetrachloroethane       100       100       67-130       0       20         Benzene       100       110       70-130       10       20         Toluene       110       100       70-130       10       20         Ethylbenzene       110       100       70-130       10       20         Chloromethane       160       Q       150       Q       64-130       6       20         Bromomethane       84       81       39-139       4       20	cis-1,3-Dichloropropene	100		100		70-130	0	20	
Benzene     100     110     70-130     10     20       Toluene     110     100     70-130     10     20       Ethylbenzene     110     100     70-130     10     20       Chloromethane     160     Q     150     Q     64-130     6     20       Bromomethane     84     81     39-139     4     20	Bromoform	87		86		54-136	1	20	
Toluene         110         100         70-130         10         20           Ethylbenzene         110         100         70-130         10         20           Chloromethane         160         Q         150         Q         64-130         6         20           Bromomethane         84         81         39-139         4         20	1,1,2,2-Tetrachloroethane	100		100		67-130	0	20	
Ethylbenzene         110         100         70-130         10         20           Chloromethane         160         Q         150         Q         64-130         6         20           Bromomethane         84         81         39-139         4         20	Benzene	100		110		70-130	10	20	
Chloromethane         160         Q         150         Q         64-130         6         20           Bromomethane         84         81         39-139         4         20	Toluene	110		100		70-130	10	20	
Bromomethane 84 81 39-139 4 20	Ethylbenzene	110		100		70-130	10	20	
	Chloromethane	160	Q	150	Q	64-130	6	20	
Vinyl chloride 100 89 55-140 12 20	Bromomethane	84		81		39-139	4	20	
	Vinyl chloride	100		89		55-140	12	20	



# Lab Control Sample Analysis Batch Quality Control

**Project Name:** SIGNORE POST INJECTION

**Project Number:** 21.0056367.66

Lab Number: L2150447

**Report Date:** 09/24/21

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/MS - Westborough L	_ab Associated	sample(s): 0	1 Batch: WG1	549680-3	WG1549680-4			
Chloroethane	66		65		55-138	2		20
1,1-Dichloroethene	110		100		61-145	10		20
trans-1,2-Dichloroethene	110		110		70-130	0		20
Trichloroethene	98		96		70-130	2		20
1,2-Dichlorobenzene	100		97		70-130	3		20
1,3-Dichlorobenzene	100		96		70-130	4		20
1,4-Dichlorobenzene	100		96		70-130	4		20
Methyl tert butyl ether	92		91		63-130	1		20
p/m-Xylene	100		100		70-130	0		20
o-Xylene	100		100		70-130	0		20
cis-1,2-Dichloroethene	110		100		70-130	10		20
Styrene	100		100		70-130	0		20
Dichlorodifluoromethane	100		100		36-147	0		20
Acetone	100		85		58-148	16		20
Carbon disulfide	110		110		51-130	0		20
2-Butanone	110		98		63-138	12		20
4-Methyl-2-pentanone	97		96		59-130	1		20
2-Hexanone	110		97		57-130	13		20
Bromochloromethane	96		96		70-130	0		20
1,2-Dibromoethane	93		90		70-130	3		20
1,2-Dibromo-3-chloropropane	91		80		41-144	13		20
Isopropylbenzene	100		100		70-130	0		20
1,2,3-Trichlorobenzene	110		100		70-130	10		20



# Lab Control Sample Analysis Batch Quality Control

**Project Name:** SIGNORE POST INJECTION

**Project Number:** 21.0056367.66

Lab Number: L2150447

**Report Date:** 09/24/21

Beremeter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits	
Parameter	76Necovery	Quai	7011CCOVERY	Quai	Lililis	KPU	Quai	LIIIIII	
Volatile Organics by GC/MS - Westborough	Lab Associated s	sample(s): 01	Batch: WG	1549680-3	WG1549680-4				
1,2,4-Trichlorobenzene	110		98		70-130	12		20	
Methyl Acetate	130		120		70-130	8		20	
Cyclohexane	130		130		70-130	0		20	
1,4-Dioxane	100		102		56-162	2		20	
Freon-113	110		100		70-130	10		20	
Methyl cyclohexane	99		94		70-130	5		20	

Surrogate	LCS %Recovery Qual	LCSD %Recovery Qual	Acceptance Criteria
1,2-Dichloroethane-d4	95	96	70-130
Toluene-d8	103	104	70-130
4-Bromofluorobenzene	109	112	70-130
Dibromofluoromethane	91	93	70-130

Lab Number: L2150447

Report Date: 09/24/21

Project Name: SIGNORE POST INJECTION

**Project Number:** 21.0056367.66

## Sample Receipt and Container Information

Were project specific reporting limits specified?

**Cooler Information** 

Cooler Custody Seal

B Absent

Container Info	ormation		Initial	Final	Temp			Frozen	
Container ID	Container Type	Cooler	рН	pН	deg C	Pres	Seal	Date/Time	Analysis(*)
L2150447-01A	Vial HCl preserved	В	NA		3.4	Υ	Absent		NYTCL-8260-R2(14)
L2150447-01B	Vial HCl preserved	В	NA		3.4	Υ	Absent		NYTCL-8260-R2(14)
L2150447-01C	Vial HCI preserved	В	NA		3.4	Υ	Absent		NYTCL-8260-R2(14)



Project Name:SIGNORE POST INJECTIONLab Number:L2150447Project Number:21.0056367.66Report Date:09/24/21

### **GLOSSARY**

#### **Acronyms**

**EDL** 

LOD

MDI

MS

DL - Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the limit of quantitation (LOQ). The DL includes any adjustments from dilutions, concentrations or moisture content, where applicable (DoD report formats only)

from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)

 Estimated Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The EDL includes any adjustments from dilutions, concentrations or moisture content, where applicable. The use of EDLs is specific to the analysis of PAHs using Solid-Phase Microextraction (SPME).

EMPC - Estimated Maximum Possible Concentration: The concentration that results from the signal present at the retention time of an analyte when the ions meet all of the identification criteria except the ion abundance ratio criteria. An EMPC is a worst-case

estimate of the concentration.

EPA - Environmental Protection Agency.

LCS - Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of

analytes or a material containing known and verified amounts of analytes.

LCSD - Laboratory Control Sample Duplicate: Refer to LCS.

LFB - Laboratory Fortified Blank: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.

 Limit of Detection: This value represents the level to which a target analyte can reliably be detected for a specific analyte in a specific matrix by a specific method. The LOD includes any adjustments from dilutions, concentrations or moisture content,

where applicable. (DoD report formats only.)

LOQ - Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats

Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats

- Method Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The MDL includes any

adjustments from dilutions, concentrations or moisture content, where applicable.

- Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for which an independent estimate of target analyte concentration is available. For Method 332.0, the spike recovery is calculated using the native concentration, including estimated values.

MSD - Matrix Spike Sample Duplicate: Refer to MS.

NA - Not Applicable.

NC - Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's

reporting unit.

NDPA/DPA - N-Nitrosodiphenylamine/Diphenylamine.

NI - Not Ignitable.

NP - Non-Plastic: Term is utilized for the analysis of Atterberg Limits in soil.

NR - No Results: Term is utilized when 'No Target Compounds Requested' is reported for the analysis of Volatile or Semivolatile

Organic TIC only requests.

RL - Reporting Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RL

includes any adjustments from dilutions, concentrations or moisture content, where applicable.

RPD - Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the

values; although the RPD value will be provided in the report.

SRM - Standard Reference Material: A reference sample of a known or certified value that is of the same or similar matrix as the

associated field samples.

STLP - Semi-dynamic Tank Leaching Procedure per EPA Method 1315.

TEF - Toxic Equivalency Factors: The values assigned to each dioxin and furan to evaluate their toxicity relative to 2,3,7,8-TCDD.

TEQ - Toxic Equivalent: The measure of a sample's toxicity derived by multiplying each dioxin and furan by its corresponding TEF

and then summing the resulting values.

TIC - Tentatively Identified Compound: A compound that has been identified to be present and is not part of the target compound list (TCL) for the method and/or program. All TICs are qualitatively identified and reported as estimated concentrations.

Report Format: DU Report with 'J' Qualifiers



Project Name:SIGNORE POST INJECTIONLab Number:L2150447Project Number:21.0056367.66Report Date:09/24/21

#### **Footnotes**

1 - The reference for this analyte should be considered modified since this analyte is absent from the target analyte list of the original method.

#### **Terms**

Analytical Method: Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum.

Difference: With respect to Total Oxidizable Precursor (TOP) Assay analysis, the difference is defined as the Post-Treatment value minus the Pre-Treatment value.

Final pH: As it pertains to Sample Receipt & Container Information section of the report, Final pH reflects pH of container determined after adjustment at the laboratory, if applicable. If no adjustment required, value reflects Initial pH.

Frozen Date/Time: With respect to Volatile Organics in soil, Frozen Date/Time reflects the date/time at which associated Reagent Water-preserved vials were initially frozen. Note: If frozen date/time is beyond 48 hours from sample collection, value will be reflected in 'bold'.

Initial pH: As it pertains to Sample Receipt & Container Information section of the report, Initial pH reflects pH of container determined upon receipt, if applicable.

PAH Total: With respect to Alkylated PAH analyses, the 'PAHs, Total' result is defined as the summation of results for all or a subset of the following compounds: Naphthalene, C1-C4 Naphthalenes, 2-Methylnaphthalene, 1-Methylnaphthalene, Biphenyl, Acenaphthylene, Acenaphthene, Fluorene, C1-C3 Fluorenes, Phenanthrene, C1-C4 Phenanthrenes/Anthracenes, Anthracene, Fluoranthene, Pyrene, C1-C4 Fluoranthenes/Pyrenes, Benza(a)anthracene, Chrysene, C1-C4 Chrysenes, Benzo(b)fluoranthene, Benzo(j)+(k)fluoranthene, Benzo(e)pyrene, Benzo(a)pyrene, Perylene, Indeno(1,2,3-cd)pyrene, Dibenz(ah)+(ac)anthracene, Benzo(g,h,i)perylene. If a 'Total' result is requested, the results of its individual components will also be reported.

PFAS Total: With respect to PFAS analyses, the 'PFAS, Total (5)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA and PFOS. In addition, the 'PFAS, Total (6)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA, PFDA and PFOS. For MassDEP DW compliance analysis only, the 'PFAS, Total (6)' result is defined as the summation of results at or above the RL. Note: If a 'Total' result is requested, the results of its individual components will also be reported.

The target compound Chlordane (CAS No. 57-74-9) is reported for GC ECD analyses. Per EPA, this compound "refers to a mixture of chlordane isomers, other chlorinated hydrocarbons and numerous other components." (Reference: USEPA Toxicological Review of Chlordane, In Support of Summary Information on the Integrated Risk Information System (IRIS), December 1997.)

Total: With respect to Organic analyses, a 'Total' result is defined as the summation of results for individual isomers or Aroclors. If a 'Total' result is requested, the results of its individual components will also be reported. This is applicable to 'Total' results for methods 8260, 8081 and 8082.

## Data Qualifiers

- A -Spectra identified as "Aldol Condensates" are byproducts of the extraction/concentration procedures when acetone is introduced in the process.
- The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For MCP-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For DOD-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank AND the analyte was detected above one-half the reporting limit (or above the reporting limit for common lab contaminants) in the associated method blank. For NJ-Air-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte above the reporting limit. For NJ-related projects (excluding Air), flag only applies to associated field samples that have detectable concentrations of the analyte, which was detected above the reporting limit in the associated method blank or above five times the reporting limit for common lab contaminants (Phthalates, Acetone, Methylene Chloride, 2-Butanone).
- Co-elution: The target analyte co-elutes with a known lab standard (i.e. surrogate, internal standards, etc.) for co-extracted analyses.
- Concentration of analyte was quantified from diluted analysis. Flag only applies to field samples that have detectable concentrations of the analyte.
- E Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.
- F The ratio of quantifier ion response to qualifier ion response falls outside of the laboratory criteria. Results are considered to be an estimated maximum concentration.
- G The concentration may be biased high due to matrix interferences (i.e, co-elution) with non-target compound(s). The result should be considered estimated.
- H The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.
- I The lower value for the two columns has been reported due to obvious interference.
- Estimated value. The Target analyte concentration is below the quantitation limit (RL), but above the Method Detection Limit (MDL) or Estimated Detection Limit (EDL) for SPME-related analyses. This represents an estimated concentration for Tentatively Identified Compounds (TICs).
- $\label{eq:main_equation} \textbf{M} \qquad \text{-Reporting Limit (RL) exceeds the MCP CAM Reporting Limit for this analyte.}$
- ND Not detected at the method detection limit (MDL) for the sample, or estimated detection limit (EDL) for SPME-related analyses.

Report Format: DU Report with 'J' Qualifiers



Project Name:SIGNORE POST INJECTIONLab Number:L2150447Project Number:21.0056367.66Report Date:09/24/21

#### **Data Qualifiers**

- NJ Presumptive evidence of compound. This represents an estimated concentration for Tentatively Identified Compounds (TICs), where the identification is based on a mass spectral library search.
- P The RPD between the results for the two columns exceeds the method-specified criteria.
- Q The quality control sample exceeds the associated acceptance criteria. For DOD-related projects, LCS and/or Continuing Calibration Standard exceedences are also qualified on all associated sample results. Note: This flag is not applicable for matrix spike recoveries when the sample concentration is greater than 4x the spike added or for batch duplicate RPD when the sample concentrations are less than 5x the RL. (Metals only.)
- **R** Analytical results are from sample re-analysis.
- RE Analytical results are from sample re-extraction.
- S Analytical results are from modified screening analysis.
- The surrogate associated with this target analyte has a recovery outside the QC acceptance limits. (Applicable to MassDEP DW Compliance samples only.)
- The batch matrix spike and/or duplicate associated with this target analyte has a recovery/RPD outside the QC acceptance limits.
   (Applicable to MassDEP DW Compliance samples only.)

Report Format: DU Report with 'J' Qualifiers



Project Name:SIGNORE POST INJECTIONLab Number:L2150447Project Number:21.0056367.66Report Date:09/24/21

## REFERENCES

Test Methods for Evaluating Solid Waste: Physical/Chemical Methods. EPA SW-846. Third Edition. Updates I - VI, 2018.

## LIMITATION OF LIABILITIES

Alpha Analytical performs services with reasonable care and diligence normal to the analytical testing laboratory industry. In the event of an error, the sole and exclusive responsibility of Alpha Analytical shall be to re-perform the work at it's own expense. In no event shall Alpha Analytical be held liable for any incidental, consequential or special damages, including but not limited to, damages in any way connected with the use of, interpretation of, information or analysis provided by Alpha Analytical.

We strongly urge our clients to comply with EPA protocol regarding sample volume, preservation, cooling, containers, sampling procedures, holding time and splitting of samples in the field.



Alpha Analytical, Inc.
Facility: Company-wide

Department: Quality Assurance

Title: Certificate/Approval Program Summary

ID No.:**17873** Revision 19

Published Date: 4/2/2021 1:14:23 PM

Page 1 of 1

## Certification Information

### The following analytes are not included in our Primary NELAP Scope of Accreditation:

### Westborough Facility

EPA 624/624.1: m/p-xylene, o-xylene, Naphthalene

EPA 625/625.1: alpha-Terpineol

**EPA 8260C/8260D:** NPW: 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene, Azobenzene; SCM: lodomethane (methyl iodide), 1,2,4,5-Tetramethylbenzene;

4-Ethyltoluene.

EPA 8270D/8270E: NPW: Dimethylnaphthalene,1,4-Diphenylhydrazine, alpha-Terpineol; SCM: Dimethylnaphthalene,1,4-Diphenylhydrazine.

SM4500: NPW: Amenable Cyanide; SCM: Total Phosphorus, TKN, NO2, NO3.

## **Mansfield Facility**

SM 2540D: TSS

EPA 8082A: NPW: PCB: 1, 5, 31, 87,101, 110, 141, 151, 153, 180, 183, 187.

EPA TO-15: Halothane, 2,4,4-Trimethyl-2-pentene, 2,4,4-Trimethyl-1-pentene, Thiophene, 2-Methylthiophene,

3-Methylthiophene, 2-Ethylthiophene, 1,2,3-Trimethylbenzene, Indan, Indene, 1,2,4,5-Tetramethylbenzene, Benzothiophene, 1-Methylnaphthalene.

Biological Tissue Matrix: EPA 3050B

## The following analytes are included in our Massachusetts DEP Scope of Accreditation

## Westborough Facility:

#### **Drinking Water**

EPA 300.0: Chloride, Nitrate-N, Fluoride, Sulfate; EPA 353.2: Nitrate-N, Nitrite-N; SM4500NO3-F: Nitrate-N, Nitrite-N; SM4500F-C, SM4500CN-CE,

EPA 180.1, SM2130B, SM4500CI-D, SM2320B, SM2540C, SM4500H-B, SM4500NO2-B

EPA 332: Perchlorate; EPA 524.2: THMs and VOCs; EPA 504.1: EDB, DBCP.

Microbiology: SM9215B; SM9223-P/A, SM9223B-Colilert-QT,SM9222D.

#### Non-Potable Water

SM4500H,B, EPA 120.1, SM2510B, SM2540C, SM2320B, SM4500CL-E, SM4500F-BC, SM4500NH3-BH: Ammonia-N and Kjeldahl-N, EPA 350.1: Ammonia-N, LACHAT 10-107-06-1-B: Ammonia-N, EPA 351.1, SM4500NO3-F, EPA 353.2: Nitrate-N, SM4500P-E, SM4500P-B, E, SM4500SO4-E, SM5220D, EPA 410.4, SM5210B, SM5310C, SM4500CL-D, EPA 1664, EPA 420.1, SM4500-CN-CE, SM2540D, EPA 300: Chloride, Sulfate, Nitrate. EPA 624.1: Volatile Halocarbons & Aromatics.

EPA 608.3: Chlordane, Toxaphene, Aldrin, alpha-BHC, beta-BHC, gamma-BHC, delta-BHC, Dieldrin, DDD, DDE, DDT, Endosulfan II, Endosulfan III, Endosulfan sulfate, Endrin, Endrin Aldehyde, Heptachlor, Heptachlor Epoxide, PCBs

EPA 625.1: SVOC (Acid/Base/Neutral Extractables), EPA 600/4-81-045: PCB-Oil.

Microbiology: SM9223B-Colilert-QT; Enterolert-QT, SM9221E, EPA 1600, EPA 1603, SM9222D.

## Mansfield Facility:

## **Drinking Water**

**EPA 200.7:** Al, Ba, Cd, Cr, Cu, Fe, Mn, Ni, Na, Ag, Ca, Zn. **EPA 200.8:** Al, Sb, As, Ba, Be, Cd, Cr, Cu, Pb, Mn, Ni, Se, Ag, TL, Zn. **EPA 245.1** Hg. **EPA 522, EPA 537.1.** 

#### Non-Potable Water

EPA 200.7: Al, Sb, As, Be, Cd, Ca, Cr, Co, Cu, Fe, Pb, Mg, Mn, Mo, Ni, K, Se, Ag, Na, Sr, TL, Ti, V, Zn.

EPA 200.8: Al, Sb, As, Be, Cd, Cr, Cu, Fe, Pb, Mn, Ni, K, Se, Ag, Na, TL, Zn.

EPA 245.1 Hg

SM2340B

For a complete listing of analytes and methods, please contact your Alpha Project Manager.

Document Type: Form Pre-Qualtrax Document ID: 08-113

Westborough, MA 01581 8 Walkup Dr. TEL: 508-898-9220 FAX: 508-898-9193	NEW YORK CHAIN OF CUSTODY Mansfield, MA 02048 320 Forbes Blvd TEL: 508-822-9300 FAX: 508-822-3288	Service Centers Mahwah, NJ 07430: 35 Whitne Albany, NY 12205: 14 Walker Tonawanda, NY 14150: 275 Co  Project Information  Project Name: Signo  Project Location: El	Way poper Ave, Suite 10 Post	- Injec	Page	-	Deliv	Date F in L erables ASP-A			×	ASP-			ALPHA Job # (2) 50 4 4 7  Billing Information  Same as Client Info
Client Information	THE PARTY OF	Project #		,,,,			Other NY					*	NEW CO.		
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Address: 300 Pearl	St. Suite 700	Project Manager: T. ALPHAQuote #:					THE REAL PROPERTY.	NY TO	38		=	NY Pa			Please identify below location of applicable disposal facilities.
Phone: 716 - 844	-7050	Turn-Around Time	SERVING.	ASS 2010	III SOURIS	CONTRACTOR OF	H				=		-51	- 1	Disposal Facility:
Fax:		Standar	Standard Due Date: sh (only if pre approved) # of Days:					NY Restricted Use Other  NY Unrestricted Use  NYC Sewer Discharge						NJ NY Other:	
These samples have be	C-						$\dashv$	Sample Filtration							
Other project specific  Please specify Metals		ents:					17CC								□ Done □ Lab to do Preservation □ Lab to do  (Please Specify below)
ALPHA Lab ID	Sa	mple ID	Colle	ection	Sample	Sampler's	26								
(Lab Use Only)	-	mpio ib	Date	Time	Matrix	Initials	8								Sample Specific Comments
50447-01	EW-1.25 R	-091721	9-17-21	1220	GW	PN	x								
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A = None B = HCI C = HNO <sub>3</sub> D = H <sub>2</sub> SO <sub>4</sub>	Container Code P = Plastic A = Amber Glass V = Vial G = Glass B = Bacteria Cup	Westboro: Certification N Mansfield: Certification N				ntainer Type									Please print clearly, legibly and completely. Samples can not be logged in and turnaround time clock will no
F = MeOH G = NaHSO <sub>4</sub> H = Na <sub>2</sub> S <sub>2</sub> O <sub>2</sub>	C = Cube O = Other E = Encore D = BOD Bottle	Relinquished	By:	Date/ 9-17-21 9/17/21		(M	Receiv	jed By:	Ad De		91	17/	Time 21 00°C	3/0	



## ANALYTICAL REPORT

Lab Number: L2150438

Client: GZA GeoEnvironmental of New York

300 Pearl Street

Suite 700

Buffalo, NY 14202

ATTN: Thomas Bohlen Phone: (716) 844-7050

Project Name: SIGNORE

Project Number: 21.0056491.81

Report Date: 09/24/21

The original project report/data package is held by Alpha Analytical. This report/data package is paginated and should be reproduced only in its entirety. Alpha Analytical holds no responsibility for results and/or data that are not consistent with the original.

Certifications & Approvals: MA (M-MA086), NH NELAP (2064), CT (PH-0574), IL (200077), ME (MA00086), MD (348), NJ (MA935), NY (11148), NC (25700/666), PA (68-03671), RI (LAO00065), TX (T104704476), VT (VT-0935), VA (460195), USDA (Permit #P330-17-00196).

Eight Walkup Drive, Westborough, MA 01581-1019 508-898-9220 (Fax) 508-898-9193 800-624-9220 - www.alphalab.com



Project Name: SIGNORE
Project Number: 21.0056491.81

**Lab Number:** L2150438 **Report Date:** 09/24/21

Alpha Sample ID	Client ID	Matrix	Sample Location	Collection Date/Time	Receive Date
L2150438-01	EW-2.5-091521	WATER	ELLICOTTVILLE, NY	09/15/21 08:58	09/17/21
L2150438-02	MW-9I-091521	WATER	ELLICOTTVILLE, NY	09/15/21 09:50	09/17/21
L2150438-03	MW-1I-091521	WATER	ELLICOTTVILLE, NY	09/15/21 11:15	09/17/21
L2150438-04	EW-1.5-091521	WATER	ELLICOTTVILLE, NY	09/15/21 11:40	09/17/21
L2150438-05	EW-4.5-091521	WATER	ELLICOTTVILLE, NY	09/15/21 13:15	09/17/21
L2150438-06	MW-4S-091521	WATER	ELLICOTTVILLE, NY	09/15/21 14:00	09/17/21
L2150438-07	IRM-2I-091621	WATER	ELLICOTTVILLE, NY	09/16/21 09:05	09/17/21
L2150438-08	TOWN WELL-091621	WATER	ELLICOTTVILLE, NY	09/16/21 09:25	09/17/21
L2150438-09	IRM-1-091621	WATER	ELLICOTTVILLE, NY	09/16/21 10:25	09/17/21
L2150438-10	MW-5S-091621	WATER	ELLICOTTVILLE, NY	09/16/21 12:10	09/17/21
L2150438-11	MM-2I-091621	WATER	ELLICOTTVILLE, NY	09/16/21 13:15	09/17/21
L2150438-12	GW-DUP-091621	WATER	ELLICOTTVILLE, NY	09/16/21 00:00	09/17/21
L2150438-13	TRIP BLANK	WATER	ELLICOTTVILLE, NY	09/16/21 00:00	09/17/21



Project Name:SIGNORELab Number:L2150438Project Number:21.0056491.81Report Date:09/24/21

## **Case Narrative**

The samples were received in accordance with the Chain of Custody and no significant deviations were encountered during the preparation or analysis unless otherwise noted. Sample Receipt, Container Information, and the Chain of Custody are located at the back of the report.

Results contained within this report relate only to the samples submitted under this Alpha Lab Number and meet NELAP requirements for all NELAP accredited parameters unless otherwise noted in the following narrative. The data presented in this report is organized by parameter (i.e. VOC, SVOC, etc.). Sample specific Quality Control data (i.e. Surrogate Spike Recovery) is reported at the end of the target analyte list for each individual sample, followed by the Laboratory Batch Quality Control at the end of each parameter. Tentatively Identified Compounds (TICs), if requested, are reported for compounds identified to be present and are not part of the method/program Target Compound List, even if only a subset of the TCL are being reported. If a sample was re-analyzed or re-extracted due to a required quality control corrective action and if both sets of data are reported, the Laboratory ID of the re-analysis or re-extraction is designated with an "R" or "RE", respectively.

When multiple Batch Quality Control elements are reported (e.g. more than one LCS), the associated samples for each element are noted in the grey shaded header line of each data table. Any Laboratory Batch, Sample Specific % recovery or RPD value that is outside the listed Acceptance Criteria is bolded in the report. In reference to questions H (CAM) or 4 (RCP) when "NO" is checked, the performance criteria for CAM and RCP methods allow for some quality control failures to occur and still be within method compliance. In these instances, the specific failure is not narrated but noted in the associated QC Outlier Summary Report, located directly after the Case Narrative. QC information is also incorporated in the Data Usability Assessment table (Format 11) of our Data Merger tool, where it can be reviewed in conjunction with the sample result, associated regulatory criteria and any associated data usability implications.

Soil/sediments, solids and tissues are reported on a dry weight basis unless otherwise noted. Definitions of all data qualifiers and acronyms used in this report are provided in the Glossary located at the back of the report.

HOLD POLICY - For samples submitted on hold, Alpha's policy is to hold samples (with the exception of Air canisters) free of charge for 21 calendar days from the date the project is completed. After 21 calendar days, we will dispose of all samples submitted including those put on hold unless you have contacted your Alpha Project Manager and made arrangements for Alpha to continue to hold the samples. Air canisters will be disposed after 3 business days from the date the project is completed.

Please contact Project Management at 800-624-9220 with any questions.	



Project Name:SIGNORELab Number:L2150438Project Number:21.0056491.81Report Date:09/24/21

## **Case Narrative (continued)**

Report Submission

All non-detect (ND) or estimated concentrations (J-qualified) have been quantitated to the limit noted in the MDL column.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Authorized Signature:

Title: Technical Director/Representative Date: 09/24/21

Custin Walker Cristin Walker

# **ORGANICS**



## **VOLATILES**



09/15/21 08:58

**Project Name:** SIGNORE

**Project Number:** 21.0056491.81

**SAMPLE RESULTS** 

Lab Number: L2150438

Report Date: 09/24/21

Date Collected:

Lab ID: L2150438-01

Client ID: EW-2.5-091521 Sample Location: ELLICOTTVILLE, NY Date Received: 09/17/21 Field Prep: Not Specified

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 09/21/21 20:52

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westborough	Lab					
Methylene chloride	ND		ug/l	2.5	0.70	1
1,1-Dichloroethane	ND		ug/l	2.5	0.70	1
Chloroform	ND		ug/l	2.5	0.70	1
Carbon tetrachloride	ND		ug/l	0.50	0.13	1
1,2-Dichloropropane	ND		ug/l	1.0	0.14	1
Dibromochloromethane	ND		ug/l	0.50	0.15	1
1,1,2-Trichloroethane	ND		ug/l	1.5	0.50	1
Tetrachloroethene	ND		ug/l	0.50	0.18	1
Chlorobenzene	ND		ug/l	2.5	0.70	1
Trichlorofluoromethane	ND		ug/l	2.5	0.70	1
1,2-Dichloroethane	ND		ug/l	0.50	0.13	1
1,1,1-Trichloroethane	ND		ug/l	2.5	0.70	1
Bromodichloromethane	ND		ug/l	0.50	0.19	1
trans-1,3-Dichloropropene	ND		ug/l	0.50	0.16	1
cis-1,3-Dichloropropene	ND		ug/l	0.50	0.14	1
Bromoform	ND		ug/l	2.0	0.65	1
1,1,2,2-Tetrachloroethane	ND		ug/l	0.50	0.17	1
Benzene	ND		ug/l	0.50	0.16	1
Toluene	ND		ug/l	2.5	0.70	1
Ethylbenzene	ND		ug/l	2.5	0.70	1
Chloromethane	ND		ug/l	2.5	0.70	1
Bromomethane	ND		ug/l	2.5	0.70	1
Vinyl chloride	ND		ug/l	1.0	0.07	1
Chloroethane	ND		ug/l	2.5	0.70	1
1,1-Dichloroethene	ND		ug/l	0.50	0.17	1
trans-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1
Trichloroethene	ND		ug/l	0.50	0.18	1
1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1



**Project Name:** Lab Number: **SIGNORE** L2150438

**Project Number:** Report Date: 21.0056491.81 09/24/21

**SAMPLE RESULTS** 

Lab ID: L2150438-01 Date Collected: 09/15/21 08:58

Client ID: Date Received: 09/17/21 EW-2.5-091521 Sample Location: Field Prep: ELLICOTTVILLE, NY Not Specified

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Wes	stborough Lab					
1,3-Dichlorobenzene	ND		ug/l	2.5	0.70	1
1,4-Dichlorobenzene	ND		ug/l	2.5	0.70	1
Methyl tert butyl ether	ND		ug/l	2.5	0.70	1
p/m-Xylene	ND		ug/l	2.5	0.70	1
o-Xylene	ND		ug/l	2.5	0.70	1
cis-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1
Styrene	ND		ug/l	2.5	0.70	1
Dichlorodifluoromethane	ND		ug/l	5.0	1.0	1
Acetone	ND		ug/l	5.0	1.5	1
Carbon disulfide	ND		ug/l	5.0	1.0	1
2-Butanone	ND		ug/l	5.0	1.9	1
4-Methyl-2-pentanone	ND		ug/l	5.0	1.0	1
2-Hexanone	ND		ug/l	5.0	1.0	1
Bromochloromethane	ND		ug/l	2.5	0.70	1
1,2-Dibromoethane	ND		ug/l	2.0	0.65	1
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70	1
Isopropylbenzene	ND		ug/l	2.5	0.70	1
1,2,3-Trichlorobenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70	1
Methyl Acetate	ND		ug/l	2.0	0.23	1
Cyclohexane	ND		ug/l	10	0.27	1
1,4-Dioxane	ND		ug/l	250	61.	1
Freon-113	ND		ug/l	2.5	0.70	1
Methyl cyclohexane	ND		ug/l	10	0.40	1

Surrogate	% Recovery	Acceptance Qualifier Criteria	
1,2-Dichloroethane-d4	100	70-130	
Toluene-d8	98	70-130	
4-Bromofluorobenzene	105	70-130	
Dibromofluoromethane	96	70-130	



L2150438

Lab Number:

**Project Name: SIGNORE** 

**Project Number:** Report Date: 21.0056491.81 09/24/21

**SAMPLE RESULTS** 

Lab ID: L2150438-02 Date Collected: 09/15/21 09:50

Client ID: Date Received: 09/17/21 MW-9I-091521 Sample Location: Field Prep: Not Specified ELLICOTTVILLE, NY

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 09/21/21 21:13

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor	
Volatile Organics by GC/MS - Wes	stborough Lab						
Methylene chloride	ND		ug/l	2.5	0.70	1	
1,1-Dichloroethane	ND		ug/l	2.5	0.70	1	
Chloroform	ND		ug/l	2.5	0.70	1	
Carbon tetrachloride	ND		ug/l	0.50	0.13	1	
1,2-Dichloropropane	ND		ug/l	1.0	0.14	1	
Dibromochloromethane	ND		ug/l	0.50	0.15	1	
1,1,2-Trichloroethane	ND		ug/l	1.5	0.50	1	
Tetrachloroethene	0.73		ug/l	0.50	0.18	1	
Chlorobenzene	ND		ug/l	2.5	0.70	1	
Trichlorofluoromethane	ND		ug/l	2.5	0.70	1	
1,2-Dichloroethane	ND		ug/l	0.50	0.13	1	
1,1,1-Trichloroethane	ND		ug/l	2.5	0.70	1	
Bromodichloromethane	ND		ug/l	0.50	0.19	1	
trans-1,3-Dichloropropene	ND		ug/l	0.50	0.16	1	
cis-1,3-Dichloropropene	ND		ug/l	0.50	0.14	1	
Bromoform	ND		ug/l	2.0	0.65	1	
1,1,2,2-Tetrachloroethane	ND		ug/l	0.50	0.17	1	
Benzene	ND		ug/l	0.50	0.16	1	
Toluene	ND		ug/l	2.5	0.70	1	
Ethylbenzene	ND		ug/l	2.5	0.70	1	
Chloromethane	ND		ug/l	2.5	0.70	1	
Bromomethane	ND		ug/l	2.5	0.70	1	
Vinyl chloride	ND		ug/l	1.0	0.07	1	
Chloroethane	ND		ug/l	2.5	0.70	1	
1,1-Dichloroethene	ND		ug/l	0.50	0.17	1	
trans-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1	
Trichloroethene	2.2		ug/l	0.50	0.18	1	
1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1	



Project Name: SIGNORE Lab Number: L2150438

**Project Number:** 21.0056491.81 **Report Date:** 09/24/21

**SAMPLE RESULTS** 

Lab ID: L2150438-02 Date Collected: 09/15/21 09:50

Client ID: MW-9I-091521 Date Received: 09/17/21 Sample Location: ELLICOTTVILLE, NY Field Prep: Not Specified

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westboroug	gh Lab					
1,3-Dichlorobenzene	ND		ug/l	2.5	0.70	1
1,4-Dichlorobenzene	ND		ug/l	2.5	0.70	1
Methyl tert butyl ether	ND		ug/l	2.5	0.70	1
p/m-Xylene	ND		ug/l	2.5	0.70	1
o-Xylene	ND		ug/l	2.5	0.70	1
cis-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1
Styrene	ND		ug/l	2.5	0.70	1
Dichlorodifluoromethane	ND		ug/l	5.0	1.0	1
Acetone	ND		ug/l	5.0	1.5	1
Carbon disulfide	ND		ug/l	5.0	1.0	1
2-Butanone	ND		ug/l	5.0	1.9	1
4-Methyl-2-pentanone	ND		ug/l	5.0	1.0	1
2-Hexanone	ND		ug/l	5.0	1.0	1
Bromochloromethane	ND		ug/l	2.5	0.70	1
1,2-Dibromoethane	ND		ug/l	2.0	0.65	1
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70	1
Isopropylbenzene	ND		ug/l	2.5	0.70	1
1,2,3-Trichlorobenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70	1
Methyl Acetate	ND		ug/l	2.0	0.23	1
Cyclohexane	ND		ug/l	10	0.27	1
1,4-Dioxane	ND		ug/l	250	61.	1
Freon-113	ND		ug/l	2.5	0.70	1
Methyl cyclohexane	ND		ug/l	10	0.40	1

Surrogate	% Recovery	Acceptance Qualifier Criteria	
1,2-Dichloroethane-d4	98	70-130	
Toluene-d8	102	70-130	
4-Bromofluorobenzene	102	70-130	
Dibromofluoromethane	95	70-130	



09/15/21 11:15

**Project Name:** SIGNORE

**Project Number:** 21.0056491.81

**SAMPLE RESULTS** 

Lab Number: L2150438

Report Date: 09/24/21

Lab ID: L2150438-03

Client ID: MW-1I-091521 Sample Location: ELLICOTTVILLE, NY Date Received: 09/17/21 Field Prep: Not Specified

Date Collected:

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 09/21/21 21:33

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westborou	gh Lab					
Methylene chloride	ND		ug/l	2.5	0.70	1
1,1-Dichloroethane	1.6	J	ug/l	2.5	0.70	1
Chloroform	ND		ug/l	2.5	0.70	1
Carbon tetrachloride	ND		ug/l	0.50	0.13	1
1,2-Dichloropropane	ND		ug/l	1.0	0.14	1
Dibromochloromethane	ND		ug/l	0.50	0.15	1
1,1,2-Trichloroethane	ND		ug/l	1.5	0.50	1
Tetrachloroethene	ND		ug/l	0.50	0.18	1
Chlorobenzene	ND		ug/l	2.5	0.70	1
Trichlorofluoromethane	ND		ug/l	2.5	0.70	1
1,2-Dichloroethane	ND		ug/l	0.50	0.13	1
1,1,1-Trichloroethane	ND		ug/l	2.5	0.70	1
Bromodichloromethane	ND		ug/l	0.50	0.19	1
trans-1,3-Dichloropropene	ND		ug/l	0.50	0.16	1
cis-1,3-Dichloropropene	ND		ug/l	0.50	0.14	1
Bromoform	ND		ug/l	2.0	0.65	1
1,1,2,2-Tetrachloroethane	ND		ug/l	0.50	0.17	1
Benzene	ND		ug/l	0.50	0.16	1
Toluene	ND		ug/l	2.5	0.70	1
Ethylbenzene	ND		ug/l	2.5	0.70	1
Chloromethane	ND		ug/l	2.5	0.70	1
Bromomethane	ND		ug/l	2.5	0.70	1
Vinyl chloride	ND		ug/l	1.0	0.07	1
Chloroethane	ND		ug/l	2.5	0.70	1
1,1-Dichloroethene	ND		ug/l	0.50	0.17	1
trans-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1
Trichloroethene	ND		ug/l	0.50	0.18	1
1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1



**Project Name:** Lab Number: SIGNORE L2150438

**Project Number:** Report Date: 21.0056491.81 09/24/21

**SAMPLE RESULTS** 

Lab ID: Date Collected: 09/15/21 11:15 L2150438-03

Client ID: Date Received: 09/17/21 MW-1I-091521 Sample Location: Field Prep: ELLICOTTVILLE, NY Not Specified

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor				
Volatile Organics by GC/MS - Westborough Lab										
1,3-Dichlorobenzene	ND		ug/l	2.5	0.70	1				
1,4-Dichlorobenzene	ND		ug/l	2.5	0.70	1				
Methyl tert butyl ether	ND		ug/l	2.5	0.70	1				
p/m-Xylene	ND		ug/l	2.5	0.70	1				
o-Xylene	ND		ug/l	2.5	0.70	1				
cis-1,2-Dichloroethene	0.70	J	ug/l	2.5	0.70	1				
Styrene	ND		ug/l	2.5	0.70	1				
Dichlorodifluoromethane	ND		ug/l	5.0	1.0	1				
Acetone	ND		ug/l	5.0	1.5	1				
Carbon disulfide	ND		ug/l	5.0	1.0	1				
2-Butanone	ND		ug/l	5.0	1.9	1				
4-Methyl-2-pentanone	ND		ug/l	5.0	1.0	1				
2-Hexanone	ND		ug/l	5.0	1.0	1				
Bromochloromethane	ND		ug/l	2.5	0.70	1				
1,2-Dibromoethane	ND		ug/l	2.0	0.65	1				
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70	1				
Isopropylbenzene	ND		ug/l	2.5	0.70	1				
1,2,3-Trichlorobenzene	ND		ug/l	2.5	0.70	1				
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70	1				
Methyl Acetate	ND		ug/l	2.0	0.23	1				
Cyclohexane	ND		ug/l	10	0.27	1				
1,4-Dioxane	ND		ug/l	250	61.	1				
Freon-113	ND		ug/l	2.5	0.70	1				
Methyl cyclohexane	ND		ug/l	10	0.40	1				

Surrogate	% Recovery	Acceptance Qualifier Criteria	
1,2-Dichloroethane-d4	98	70-130	
Toluene-d8	101	70-130	
4-Bromofluorobenzene	99	70-130	
Dibromofluoromethane	95	70-130	



**Project Name:** SIGNORE

**Project Number:** 21.0056491.81

**SAMPLE RESULTS** 

Lab Number: L2150438

Report Date: 09/24/21

Lab ID: L2150438-04 Date Collected: 09/15/21 11:40

Client ID: Date Received: 09/17/21 EW-1.5-091521 Sample Location: Field Prep: Not Specified ELLICOTTVILLE, NY

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 09/21/21 21:54

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westb	orough Lab					
Methylene chloride	ND		ug/l	2.5	0.70	1
1,1-Dichloroethane	ND		ug/l	2.5	0.70	1
Chloroform	ND		ug/l	2.5	0.70	1
Carbon tetrachloride	ND		ug/l	0.50	0.13	1
1,2-Dichloropropane	ND		ug/l	1.0	0.14	1
Dibromochloromethane	ND		ug/l	0.50	0.15	1
1,1,2-Trichloroethane	ND		ug/l	1.5	0.50	1
Tetrachloroethene	ND		ug/l	0.50	0.18	1
Chlorobenzene	ND		ug/l	2.5	0.70	1
Trichlorofluoromethane	ND		ug/l	2.5	0.70	1
1,2-Dichloroethane	ND		ug/l	0.50	0.13	1
1,1,1-Trichloroethane	ND		ug/l	2.5	0.70	1
Bromodichloromethane	ND		ug/l	0.50	0.19	1
trans-1,3-Dichloropropene	ND		ug/l	0.50	0.16	1
cis-1,3-Dichloropropene	ND		ug/l	0.50	0.14	1
Bromoform	ND		ug/l	2.0	0.65	1
1,1,2,2-Tetrachloroethane	ND		ug/l	0.50	0.17	1
Benzene	ND		ug/l	0.50	0.16	1
Toluene	ND		ug/l	2.5	0.70	1
Ethylbenzene	ND		ug/l	2.5	0.70	1
Chloromethane	ND		ug/l	2.5	0.70	1
Bromomethane	ND		ug/l	2.5	0.70	1
Vinyl chloride	ND		ug/l	1.0	0.07	1
Chloroethane	ND		ug/l	2.5	0.70	1
1,1-Dichloroethene	ND		ug/l	0.50	0.17	1
trans-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1
Trichloroethene	ND		ug/l	0.50	0.18	1
1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1



Project Name: SIGNORE Lab Number: L2150438

**Project Number:** 21.0056491.81 **Report Date:** 09/24/21

**SAMPLE RESULTS** 

Lab ID: L2150438-04 Date Collected: 09/15/21 11:40

Client ID: EW-1.5-091521 Date Received: 09/17/21 Sample Location: ELLICOTTVILLE, NY Field Prep: Not Specified

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor				
Volatile Organics by GC/MS - Westborough Lab										
1,3-Dichlorobenzene	ND		ug/l	2.5	0.70	1				
1,4-Dichlorobenzene	ND		ug/l	2.5	0.70	1				
Methyl tert butyl ether	ND		ug/l	2.5	0.70	1				
p/m-Xylene	ND		ug/l	2.5	0.70	1				
o-Xylene	ND		ug/l	2.5	0.70	1				
cis-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1				
Styrene	ND		ug/l	2.5	0.70	1				
Dichlorodifluoromethane	ND		ug/l	5.0	1.0	1				
Acetone	ND		ug/l	5.0	1.5	1				
Carbon disulfide	ND		ug/l	5.0	1.0	1				
2-Butanone	ND		ug/l	5.0	1.9	1				
4-Methyl-2-pentanone	ND		ug/l	5.0	1.0	1				
2-Hexanone	ND		ug/l	5.0	1.0	1				
Bromochloromethane	ND		ug/l	2.5	0.70	1				
1,2-Dibromoethane	ND		ug/l	2.0	0.65	1				
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70	1				
Isopropylbenzene	ND		ug/l	2.5	0.70	1				
1,2,3-Trichlorobenzene	ND		ug/l	2.5	0.70	1				
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70	1				
Methyl Acetate	ND		ug/l	2.0	0.23	1				
Cyclohexane	ND		ug/l	10	0.27	1				
1,4-Dioxane	ND		ug/l	250	61.	1				
Freon-113	ND		ug/l	2.5	0.70	1				
Methyl cyclohexane	ND		ug/l	10	0.40	1				

Surrogate	% Recovery	Acceptance Qualifier Criteria	
1,2-Dichloroethane-d4	99	70-130	
Toluene-d8	100	70-130	
4-Bromofluorobenzene	102	70-130	
Dibromofluoromethane	94	70-130	



L2150438

**Project Name:** Lab Number: **SIGNORE** 

**Project Number:** Report Date: 21.0056491.81 09/24/21

**SAMPLE RESULTS** 

Lab ID: L2150438-05 Date Collected: 09/15/21 13:15

Client ID: Date Received: 09/17/21 EW-4.5-091521 Sample Location: Field Prep: Not Specified ELLICOTTVILLE, NY

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 09/22/21 00:58

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor	
Volatile Organics by GC/MS - Wes	stborough Lab						
Methylene chloride	ND		ug/l	2.5	0.70	1	
1,1-Dichloroethane	ND		ug/l	2.5	0.70	1	
Chloroform	ND		ug/l	2.5	0.70	1	
Carbon tetrachloride	ND		ug/l	0.50	0.13	1	
1,2-Dichloropropane	ND		ug/l	1.0	0.14	1	
Dibromochloromethane	ND		ug/l	0.50	0.15	1	
1,1,2-Trichloroethane	ND		ug/l	1.5	0.50	1	
Tetrachloroethene	0.78		ug/l	0.50	0.18	1	
Chlorobenzene	ND		ug/l	2.5	0.70	1	
Trichlorofluoromethane	ND		ug/l	2.5	0.70	1	
1,2-Dichloroethane	ND		ug/l	0.50	0.13	1	
1,1,1-Trichloroethane	ND		ug/l	2.5	0.70	1	
Bromodichloromethane	ND		ug/l	0.50	0.19	1	
trans-1,3-Dichloropropene	ND		ug/l	0.50	0.16	1	
cis-1,3-Dichloropropene	ND		ug/l	0.50	0.14	1	
Bromoform	ND		ug/l	2.0	0.65	1	
1,1,2,2-Tetrachloroethane	ND		ug/l	0.50	0.17	1	
Benzene	ND		ug/l	0.50	0.16	1	
Toluene	ND		ug/l	2.5	0.70	1	
Ethylbenzene	ND		ug/l	2.5	0.70	1	
Chloromethane	ND		ug/l	2.5	0.70	1	
Bromomethane	ND		ug/l	2.5	0.70	1	
Vinyl chloride	ND		ug/l	1.0	0.07	1	
Chloroethane	ND		ug/l	2.5	0.70	1	
1,1-Dichloroethene	ND		ug/l	0.50	0.17	1	
trans-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1	
Trichloroethene	2.4		ug/l	0.50	0.18	1	
1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1	



Project Name: SIGNORE Lab Number: L2150438

**Project Number:** 21.0056491.81 **Report Date:** 09/24/21

**SAMPLE RESULTS** 

Lab ID: L2150438-05 Date Collected: 09/15/21 13:15

Client ID: EW-4.5-091521 Date Received: 09/17/21 Sample Location: ELLICOTTVILLE, NY Field Prep: Not Specified

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor				
Volatile Organics by GC/MS - Westborough Lab										
1,3-Dichlorobenzene	ND		ug/l	2.5	0.70	1				
1,4-Dichlorobenzene	ND		ug/l	2.5	0.70	1				
Methyl tert butyl ether	ND		ug/l	2.5	0.70	1				
p/m-Xylene	ND		ug/l	2.5	0.70	1				
o-Xylene	ND		ug/l	2.5	0.70	1				
cis-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1				
Styrene	ND		ug/l	2.5	0.70	1				
Dichlorodifluoromethane	ND		ug/l	5.0	1.0	1				
Acetone	ND		ug/l	5.0	1.5	1				
Carbon disulfide	ND		ug/l	5.0	1.0	1				
2-Butanone	ND		ug/l	5.0	1.9	1				
4-Methyl-2-pentanone	ND		ug/l	5.0	1.0	1				
2-Hexanone	ND		ug/l	5.0	1.0	1				
Bromochloromethane	ND		ug/l	2.5	0.70	1				
1,2-Dibromoethane	ND		ug/l	2.0	0.65	1				
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70	1				
Isopropylbenzene	ND		ug/l	2.5	0.70	1				
1,2,3-Trichlorobenzene	ND		ug/l	2.5	0.70	1				
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70	1				
Methyl Acetate	ND		ug/l	2.0	0.23	1				
Cyclohexane	ND		ug/l	10	0.27	1				
1,4-Dioxane	ND		ug/l	250	61.	1				
Freon-113	ND		ug/l	2.5	0.70	1				
Methyl cyclohexane	ND		ug/l	10	0.40	1				

Surrogate	% Recovery	Qualifier	Acceptance Criteria	
1,2-Dichloroethane-d4	103		70-130	
Toluene-d8	101		70-130	
4-Bromofluorobenzene	99		70-130	
Dibromofluoromethane	97		70-130	



L2150438

Lab Number:

**Project Name: SIGNORE** 

Report Date:

**Project Number:** 21.0056491.81 09/24/21

**SAMPLE RESULTS** 

Lab ID: L2150438-06 Date Collected: 09/15/21 14:00

Client ID: Date Received: 09/17/21 MW-4S-091521 Sample Location: Field Prep: Not Specified ELLICOTTVILLE, NY

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 09/21/21 22:14

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westb	oorough Lab					
Methylene chloride	ND		ug/l	2.5	0.70	1
1,1-Dichloroethane	ND		ug/l	2.5	0.70	1
Chloroform	ND		ug/l	2.5	0.70	1
Carbon tetrachloride	ND		ug/l	0.50	0.13	1
1,2-Dichloropropane	ND		ug/l	1.0	0.14	1
Dibromochloromethane	ND		ug/l	0.50	0.15	1
1,1,2-Trichloroethane	ND		ug/l	1.5	0.50	1
Tetrachloroethene	0.18	J	ug/l	0.50	0.18	1
Chlorobenzene	ND		ug/l	2.5	0.70	1
Trichlorofluoromethane	ND		ug/l	2.5	0.70	1
1,2-Dichloroethane	ND		ug/l	0.50	0.13	1
1,1,1-Trichloroethane	ND		ug/l	2.5	0.70	1
Bromodichloromethane	ND		ug/l	0.50	0.19	1
trans-1,3-Dichloropropene	ND		ug/l	0.50	0.16	1
cis-1,3-Dichloropropene	ND		ug/l	0.50	0.14	1
Bromoform	ND		ug/l	2.0	0.65	1
1,1,2,2-Tetrachloroethane	ND		ug/l	0.50	0.17	1
Benzene	ND		ug/l	0.50	0.16	1
Toluene	ND		ug/l	2.5	0.70	1
Ethylbenzene	ND		ug/l	2.5	0.70	1
Chloromethane	ND		ug/l	2.5	0.70	1
Bromomethane	ND		ug/l	2.5	0.70	1
Vinyl chloride	ND		ug/l	1.0	0.07	1
Chloroethane	ND		ug/l	2.5	0.70	1
1,1-Dichloroethene	ND		ug/l	0.50	0.17	1
trans-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1
Trichloroethene	ND		ug/l	0.50	0.18	1
1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1



Project Name: SIGNORE Lab Number: L2150438

**Project Number:** 21.0056491.81 **Report Date:** 09/24/21

**SAMPLE RESULTS** 

Lab ID: L2150438-06 Date Collected: 09/15/21 14:00

Client ID: MW-4S-091521 Date Received: 09/17/21 Sample Location: ELLICOTTVILLE, NY Field Prep: Not Specified

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westboro	ugh Lab					
1,3-Dichlorobenzene	ND		ug/l	2.5	0.70	1
1,4-Dichlorobenzene	ND		ug/l	2.5	0.70	1
Methyl tert butyl ether	ND		ug/l	2.5	0.70	1
p/m-Xylene	ND		ug/l	2.5	0.70	1
o-Xylene	ND		ug/l	2.5	0.70	1
cis-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1
Styrene	ND		ug/l	2.5	0.70	1
Dichlorodifluoromethane	ND		ug/l	5.0	1.0	1
Acetone	ND		ug/l	5.0	1.5	1
Carbon disulfide	ND		ug/l	5.0	1.0	1
2-Butanone	ND		ug/l	5.0	1.9	1
4-Methyl-2-pentanone	ND		ug/l	5.0	1.0	1
2-Hexanone	ND		ug/l	5.0	1.0	1
Bromochloromethane	ND		ug/l	2.5	0.70	1
1,2-Dibromoethane	ND		ug/l	2.0	0.65	1
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70	1
Isopropylbenzene	ND		ug/l	2.5	0.70	1
1,2,3-Trichlorobenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70	1
Methyl Acetate	ND		ug/l	2.0	0.23	1
Cyclohexane	ND		ug/l	10	0.27	1
1,4-Dioxane	ND		ug/l	250	61.	1
Freon-113	ND		ug/l	2.5	0.70	1
Methyl cyclohexane	ND		ug/l	10	0.40	1

Surrogate	% Recovery	Acceptance Qualifier Criteria	
1,2-Dichloroethane-d4	100	70-130	
Toluene-d8	99	70-130	
4-Bromofluorobenzene	99	70-130	
Dibromofluoromethane	98	70-130	



09/16/21 09:05

Not Specified

09/17/21

**Project Name: SIGNORE** 

**Project Number:** 21.0056491.81

**SAMPLE RESULTS** 

Lab Number: L2150438

Date Collected:

Date Received:

Report Date: 09/24/21

Lab ID: L2150438-07

Client ID: IRM-2I-091621 Sample Location: ELLICOTTVILLE, NY

Field Prep:

Sample Depth:

Matrix: Water Analytical Method: 1,8260C

Analytical Date: 09/21/21 22:35

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor	
Volatile Organics by GC/MS - Westk	oorough Lab						
Methylene chloride	ND		ug/l	2.5	0.70	1	
1,1-Dichloroethane	ND		ug/l	2.5	0.70	1	
Chloroform	ND		ug/l	2.5	0.70	1	
Carbon tetrachloride	ND		ug/l	0.50	0.13	1	
1,2-Dichloropropane	ND		ug/l	1.0	0.14	1	
Dibromochloromethane	ND		ug/l	0.50	0.15	1	
1,1,2-Trichloroethane	ND		ug/l	1.5	0.50	1	
Tetrachloroethene	ND		ug/l	0.50	0.18	1	
Chlorobenzene	ND		ug/l	2.5	0.70	1	
Trichlorofluoromethane	ND		ug/l	2.5	0.70	1	
1,2-Dichloroethane	ND		ug/l	0.50	0.13	1	
1,1,1-Trichloroethane	ND		ug/l	2.5	0.70	1	
Bromodichloromethane	ND		ug/l	0.50	0.19	1	
trans-1,3-Dichloropropene	ND		ug/l	0.50	0.16	1	
cis-1,3-Dichloropropene	ND		ug/l	0.50	0.14	1	
Bromoform	ND		ug/l	2.0	0.65	1	
1,1,2,2-Tetrachloroethane	ND		ug/l	0.50	0.17	1	
Benzene	ND		ug/l	0.50	0.16	1	
Toluene	ND		ug/l	2.5	0.70	1	
Ethylbenzene	ND		ug/l	2.5	0.70	1	
Chloromethane	ND		ug/l	2.5	0.70	1	
Bromomethane	ND		ug/l	2.5	0.70	1	
Vinyl chloride	ND		ug/l	1.0	0.07	1	
Chloroethane	ND		ug/l	2.5	0.70	1	
1,1-Dichloroethene	ND		ug/l	0.50	0.17	1	
trans-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1	
Trichloroethene	0.44	J	ug/l	0.50	0.18	1	
1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1	



Project Name: SIGNORE Lab Number: L2150438

**Project Number:** 21.0056491.81 **Report Date:** 09/24/21

**SAMPLE RESULTS** 

Lab ID: Date Collected: 09/16/21 09:05

Client ID: IRM-2I-091621 Date Received: 09/17/21 Sample Location: ELLICOTTVILLE, NY Field Prep: Not Specified

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westb	orough Lab					
1,3-Dichlorobenzene	ND		ug/l	2.5	0.70	1
1,4-Dichlorobenzene	ND		ug/l	2.5	0.70	1
Methyl tert butyl ether	ND		ug/l	2.5	0.70	1
p/m-Xylene	ND		ug/l	2.5	0.70	1
o-Xylene	ND		ug/l	2.5	0.70	1
cis-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1
Styrene	ND		ug/l	2.5	0.70	1
Dichlorodifluoromethane	ND		ug/l	5.0	1.0	1
Acetone	ND		ug/l	5.0	1.5	1
Carbon disulfide	ND		ug/l	5.0	1.0	1
2-Butanone	ND		ug/l	5.0	1.9	1
4-Methyl-2-pentanone	ND		ug/l	5.0	1.0	1
2-Hexanone	ND		ug/l	5.0	1.0	1
Bromochloromethane	ND		ug/l	2.5	0.70	1
1,2-Dibromoethane	ND		ug/l	2.0	0.65	1
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70	1
Isopropylbenzene	ND		ug/l	2.5	0.70	1
1,2,3-Trichlorobenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70	1
Methyl Acetate	ND		ug/l	2.0	0.23	1
Cyclohexane	ND		ug/l	10	0.27	1
1,4-Dioxane	ND		ug/l	250	61.	1
Freon-113	ND		ug/l	2.5	0.70	1
Methyl cyclohexane	ND		ug/l	10	0.40	1

Surrogate	% Recovery	Acceptance Qualifier Criteria	
1,2-Dichloroethane-d4	93	70-130	
Toluene-d8	100	70-130	
4-Bromofluorobenzene	97	70-130	
Dibromofluoromethane	93	70-130	



L2150438

Project Name: SIGNORE Lab Number:

**Project Number:** 21.0056491.81 **Report Date:** 09/24/21

SAMPLE RESULTS

Lab ID: L2150438-08 Date Collected: 09/16/21 09:25

Client ID: TOWN WELL-091621 Date Received: 09/17/21 Sample Location: ELLICOTTVILLE, NY Field Prep: Not Specified

Sample Depth:

Matrix: Water
Analytical Method: 1,8260C
Analytical Date: 09/21/21 22:55

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Wes	tborough Lab					
Methylene chloride	ND		ug/l	2.5	0.70	1
1,1-Dichloroethane	ND		ug/l	2.5	0.70	1
Chloroform	ND		ug/l	2.5	0.70	1
Carbon tetrachloride	ND		ug/l	0.50	0.13	1
1,2-Dichloropropane	ND		ug/l	1.0	0.14	1
Dibromochloromethane	0.66		ug/l	0.50	0.15	1
1,1,2-Trichloroethane	ND		ug/l	1.5	0.50	1
Tetrachloroethene	ND		ug/l	0.50	0.18	1
Chlorobenzene	ND		ug/l	2.5	0.70	1
Trichlorofluoromethane	ND		ug/l	2.5	0.70	1
1,2-Dichloroethane	ND		ug/l	0.50	0.13	1
1,1,1-Trichloroethane	ND		ug/l	2.5	0.70	1
Bromodichloromethane	0.36	J	ug/l	0.50	0.19	1
trans-1,3-Dichloropropene	ND		ug/l	0.50	0.16	1
cis-1,3-Dichloropropene	ND		ug/l	0.50	0.14	1
Bromoform	ND		ug/l	2.0	0.65	1
1,1,2,2-Tetrachloroethane	ND		ug/l	0.50	0.17	1
Benzene	ND		ug/l	0.50	0.16	1
Toluene	ND		ug/l	2.5	0.70	1
Ethylbenzene	ND		ug/l	2.5	0.70	1
Chloromethane	ND		ug/l	2.5	0.70	1
Bromomethane	ND		ug/l	2.5	0.70	1
Vinyl chloride	ND		ug/l	1.0	0.07	1
Chloroethane	ND		ug/l	2.5	0.70	1
1,1-Dichloroethene	ND		ug/l	0.50	0.17	1
trans-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1
Trichloroethene	0.32	J	ug/l	0.50	0.18	1
1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1



MDL

**Dilution Factor** 

Project Name: SIGNORE Lab Number: L2150438

**Project Number:** 21.0056491.81 **Report Date:** 09/24/21

**SAMPLE RESULTS** 

Lab ID: L2150438-08 Date Collected: 09/16/21 09:25

Result

Client ID: TOWN WELL-091621 Date Received: 09/17/21 Sample Location: ELLICOTTVILLE, NY Field Prep: Not Specified

Qualifier

Units

RL

Sample Depth:

Parameter

i arameter	Nosun	Qualifici	Office			Dilation ractor	
Volatile Organics by GC/MS - Westb	orough Lab						
1,3-Dichlorobenzene	ND		ug/l	2.5	0.70	1	
1,4-Dichlorobenzene	ND		ug/l	2.5	0.70	1	
Methyl tert butyl ether	ND		ug/l	2.5	0.70	1	
p/m-Xylene	ND		ug/l	2.5	0.70	1	
o-Xylene	ND		ug/l	2.5	0.70	1	
cis-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1	
Styrene	ND		ug/l	2.5	0.70	1	
Dichlorodifluoromethane	ND		ug/l	5.0	1.0	1	
Acetone	ND		ug/l	5.0	1.5	1	
Carbon disulfide	ND		ug/l	5.0	1.0	1	
2-Butanone	ND		ug/l	5.0	1.9	1	
4-Methyl-2-pentanone	ND		ug/l	5.0	1.0	1	
2-Hexanone	ND		ug/l	5.0	1.0	1	
Bromochloromethane	ND		ug/l	2.5	0.70	1	
1,2-Dibromoethane	ND		ug/l	2.0	0.65	1	
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70	1	
Isopropylbenzene	ND		ug/l	2.5	0.70	1	
1,2,3-Trichlorobenzene	ND		ug/l	2.5	0.70	1	
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70	1	
Methyl Acetate	ND		ug/l	2.0	0.23	1	
Cyclohexane	ND		ug/l	10	0.27	1	
1,4-Dioxane	ND		ug/l	250	61.	1	
Freon-113	ND		ug/l	2.5	0.70	1	
Methyl cyclohexane	ND		ug/l	10	0.40	1	

Surrogate	% Recovery	Acceptance Qualifier Criteria	
1,2-Dichloroethane-d4	96	70-130	
Toluene-d8	102	70-130	
4-Bromofluorobenzene	101	70-130	
Dibromofluoromethane	94	70-130	



Project Name: SIGNORE Lab Number: L2150438

**Project Number:** 21.0056491.81 **Report Date:** 09/24/21

SAMPLE RESULTS

Lab ID: L2150438-09 Date Collected: 09/16/21 10:25

Client ID: IRM-1-091621 Date Received: 09/17/21 Sample Location: ELLICOTTVILLE, NY Field Prep: Not Specified

Sample Depth:

Matrix: Water
Analytical Method: 1,8260C
Analytical Date: 09/21/21 23:16

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westb	orough Lab					
Methylene chloride	ND		ug/l	2.5	0.70	1
1,1-Dichloroethane	ND		ug/l	2.5	0.70	1
Chloroform	ND		ug/l	2.5	0.70	1
Carbon tetrachloride	ND		ug/l	0.50	0.13	1
1,2-Dichloropropane	ND		ug/l	1.0	0.14	1
Dibromochloromethane	ND		ug/l	0.50	0.15	1
1,1,2-Trichloroethane	ND		ug/l	1.5	0.50	1
Tetrachloroethene	ND		ug/l	0.50	0.18	1
Chlorobenzene	ND		ug/l	2.5	0.70	1
Trichlorofluoromethane	ND		ug/l	2.5	0.70	1
1,2-Dichloroethane	ND		ug/l	0.50	0.13	1
1,1,1-Trichloroethane	ND		ug/l	2.5	0.70	1
Bromodichloromethane	ND		ug/l	0.50	0.19	1
trans-1,3-Dichloropropene	ND		ug/l	0.50	0.16	1
cis-1,3-Dichloropropene	ND		ug/l	0.50	0.14	1
Bromoform	ND		ug/l	2.0	0.65	1
1,1,2,2-Tetrachloroethane	ND		ug/l	0.50	0.17	1
Benzene	ND		ug/l	0.50	0.16	1
Toluene	ND		ug/l	2.5	0.70	1
Ethylbenzene	ND		ug/l	2.5	0.70	1
Chloromethane	ND		ug/l	2.5	0.70	1
Bromomethane	ND		ug/l	2.5	0.70	1
Vinyl chloride	ND		ug/l	1.0	0.07	1
Chloroethane	ND		ug/l	2.5	0.70	1
1,1-Dichloroethene	ND		ug/l	0.50	0.17	1
trans-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1
Trichloroethene	ND		ug/l	0.50	0.18	1
1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1



MDL

**Dilution Factor** 

Project Name: SIGNORE Lab Number: L2150438

**Project Number:** 21.0056491.81 **Report Date:** 09/24/21

**SAMPLE RESULTS** 

Lab ID: Date Collected: 09/16/21 10:25

Client ID: IRM-1-091621 Date Received: 09/17/21 Sample Location: ELLICOTTVILLE, NY Field Prep: Not Specified

Qualifier

Units

RL

Result

Sample Depth:

Parameter

i arameter	Nosun	Qualifici	Office			Dilation ractor	
Volatile Organics by GC/MS - Westb	orough Lab						
1,3-Dichlorobenzene	ND		ug/l	2.5	0.70	1	
1,4-Dichlorobenzene	ND		ug/l	2.5	0.70	1	
Methyl tert butyl ether	ND		ug/l	2.5	0.70	1	
p/m-Xylene	ND		ug/l	2.5	0.70	1	
o-Xylene	ND		ug/l	2.5	0.70	1	
cis-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1	
Styrene	ND		ug/l	2.5	0.70	1	
Dichlorodifluoromethane	ND		ug/l	5.0	1.0	1	
Acetone	ND		ug/l	5.0	1.5	1	
Carbon disulfide	ND		ug/l	5.0	1.0	1	
2-Butanone	ND		ug/l	5.0	1.9	1	
4-Methyl-2-pentanone	ND		ug/l	5.0	1.0	1	
2-Hexanone	ND		ug/l	5.0	1.0	1	
Bromochloromethane	ND		ug/l	2.5	0.70	1	
1,2-Dibromoethane	ND		ug/l	2.0	0.65	1	
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70	1	
Isopropylbenzene	ND		ug/l	2.5	0.70	1	
1,2,3-Trichlorobenzene	ND		ug/l	2.5	0.70	1	
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70	1	
Methyl Acetate	ND		ug/l	2.0	0.23	1	
Cyclohexane	ND		ug/l	10	0.27	1	
1,4-Dioxane	ND		ug/l	250	61.	1	
Freon-113	ND		ug/l	2.5	0.70	1	
Methyl cyclohexane	ND		ug/l	10	0.40	1	

Surrogate	% Recovery	Acceptance Qualifier Criteria	
1,2-Dichloroethane-d4	98	70-130	
Toluene-d8	101	70-130	
4-Bromofluorobenzene	99	70-130	
Dibromofluoromethane	95	70-130	



09/16/21 12:10

**Project Name: SIGNORE** 

**Project Number:** 21.0056491.81

**SAMPLE RESULTS** 

Lab Number: L2150438

Report Date: 09/24/21

Lab ID: L2150438-10

Client ID: MW-5S-091621 Sample Location: ELLICOTTVILLE, NY Date Received: 09/17/21 Field Prep: Not Specified

Date Collected:

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 09/21/21 23:36

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westl	oorough Lab					
Methylene chloride	ND		ug/l	2.5	0.70	1
1,1-Dichloroethane	ND		ug/l	2.5	0.70	1
Chloroform	ND		ug/l	2.5	0.70	1
Carbon tetrachloride	ND		ug/l	0.50	0.13	1
1,2-Dichloropropane	ND		ug/l	1.0	0.14	1
Dibromochloromethane	ND		ug/l	0.50	0.15	1
1,1,2-Trichloroethane	ND		ug/l	1.5	0.50	1
Tetrachloroethene	3.6		ug/l	0.50	0.18	1
Chlorobenzene	ND		ug/l	2.5	0.70	1
Trichlorofluoromethane	ND		ug/l	2.5	0.70	1
1,2-Dichloroethane	ND		ug/l	0.50	0.13	1
1,1,1-Trichloroethane	ND		ug/l	2.5	0.70	1
Bromodichloromethane	ND		ug/l	0.50	0.19	1
trans-1,3-Dichloropropene	ND		ug/l	0.50	0.16	1
cis-1,3-Dichloropropene	ND		ug/l	0.50	0.14	1
Bromoform	ND		ug/l	2.0	0.65	1
1,1,2,2-Tetrachloroethane	ND		ug/l	0.50	0.17	1
Benzene	ND		ug/l	0.50	0.16	1
Toluene	ND		ug/l	2.5	0.70	1
Ethylbenzene	ND		ug/l	2.5	0.70	1
Chloromethane	ND		ug/l	2.5	0.70	1
Bromomethane	ND		ug/l	2.5	0.70	1
Vinyl chloride	ND		ug/l	1.0	0.07	1
Chloroethane	ND		ug/l	2.5	0.70	1
1,1-Dichloroethene	ND		ug/l	0.50	0.17	1
trans-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1
Trichloroethene	1.7		ug/l	0.50	0.18	1
1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1



Project Name: SIGNORE Lab Number: L2150438

**Project Number:** 21.0056491.81 **Report Date:** 09/24/21

**SAMPLE RESULTS** 

Lab ID: Date Collected: 09/16/21 12:10

Client ID: MW-5S-091621 Date Received: 09/17/21 Sample Location: ELLICOTTVILLE, NY Field Prep: Not Specified

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westboroug	gh Lab					
1,3-Dichlorobenzene	ND		ug/l	2.5	0.70	1
1,4-Dichlorobenzene	ND		ug/l	2.5	0.70	1
Methyl tert butyl ether	ND		ug/l	2.5	0.70	1
p/m-Xylene	ND		ug/l	2.5	0.70	1
o-Xylene	ND		ug/l	2.5	0.70	1
cis-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1
Styrene	ND		ug/l	2.5	0.70	1
Dichlorodifluoromethane	ND		ug/l	5.0	1.0	1
Acetone	ND		ug/l	5.0	1.5	1
Carbon disulfide	ND		ug/l	5.0	1.0	1
2-Butanone	ND		ug/l	5.0	1.9	1
4-Methyl-2-pentanone	ND		ug/l	5.0	1.0	1
2-Hexanone	ND		ug/l	5.0	1.0	1
Bromochloromethane	ND		ug/l	2.5	0.70	1
1,2-Dibromoethane	ND		ug/l	2.0	0.65	1
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70	1
Isopropylbenzene	ND		ug/l	2.5	0.70	1
1,2,3-Trichlorobenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70	1
Methyl Acetate	ND		ug/l	2.0	0.23	1
Cyclohexane	ND		ug/l	10	0.27	1
1,4-Dioxane	ND		ug/l	250	61.	1
Freon-113	ND		ug/l	2.5	0.70	1
Methyl cyclohexane	ND		ug/l	10	0.40	1

Surrogate	% Recovery	Acceptance Qualifier Criteria
1,2-Dichloroethane-d4	99	70-130
Toluene-d8	102	70-130
4-Bromofluorobenzene	101	70-130
Dibromofluoromethane	93	70-130



L2150438

Lab Number:

Project Name: SIGNORE

Demant Date

**Project Number:** 21.0056491.81 **Report Date:** 09/24/21

**SAMPLE RESULTS** 

Lab ID: L2150438-11 Date Collected: 09/16/21 13:15

Client ID: MM-2I-091621 Date Received: 09/17/21 Sample Location: ELLICOTTVILLE, NY Field Prep: Not Specified

Sample Depth:

Matrix: Water
Analytical Method: 1,8260C
Analytical Date: 09/21/21 23:57

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor	
Volatile Organics by GC/MS - Wes	tborough Lab						
Methylene chloride	ND		ug/l	2.5	0.70	1	
1,1-Dichloroethane	ND		ug/l	2.5	0.70	1	
Chloroform	ND		ug/l	2.5	0.70	1	
Carbon tetrachloride	ND		ug/l	0.50	0.13	1	
1,2-Dichloropropane	ND		ug/l	1.0	0.14	1	
Dibromochloromethane	ND		ug/l	0.50	0.15	1	
1,1,2-Trichloroethane	ND		ug/l	1.5	0.50	1	
Tetrachloroethene	ND		ug/l	0.50	0.18	1	
Chlorobenzene	ND		ug/l	2.5	0.70	1	
Trichlorofluoromethane	ND		ug/l	2.5	0.70	1	
1,2-Dichloroethane	ND		ug/l	0.50	0.13	1	
1,1,1-Trichloroethane	ND		ug/l	2.5	0.70	1	
Bromodichloromethane	ND		ug/l	0.50	0.19	1	
trans-1,3-Dichloropropene	ND		ug/l	0.50	0.16	1	
cis-1,3-Dichloropropene	ND		ug/l	0.50	0.14	1	
Bromoform	ND		ug/l	2.0	0.65	1	
1,1,2,2-Tetrachloroethane	ND		ug/l	0.50	0.17	1	
Benzene	ND		ug/l	0.50	0.16	1	
Toluene	ND		ug/l	2.5	0.70	1	
Ethylbenzene	ND		ug/l	2.5	0.70	1	
Chloromethane	ND		ug/l	2.5	0.70	1	
Bromomethane	ND		ug/l	2.5	0.70	1	
Vinyl chloride	ND		ug/l	1.0	0.07	1	
Chloroethane	ND		ug/l	2.5	0.70	1	
1,1-Dichloroethene	ND		ug/l	0.50	0.17	1	
trans-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1	
Trichloroethene	ND		ug/l	0.50	0.18	1	
1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1	



**Project Name:** Lab Number: SIGNORE L2150438

**Project Number:** Report Date: 21.0056491.81 09/24/21

**SAMPLE RESULTS** 

Lab ID: Date Collected: 09/16/21 13:15 L2150438-11

Client ID: Date Received: 09/17/21 MM-2I-091621 Sample Location: Field Prep: ELLICOTTVILLE, NY Not Specified

Sample Depth:

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor				
Volatile Organics by GC/MS - Westborough Lab										
1,3-Dichlorobenzene	ND		ug/l	2.5	0.70	1				
1,4-Dichlorobenzene	ND		ug/l	2.5	0.70	1				
Methyl tert butyl ether	ND		ug/l	2.5	0.70	1				
p/m-Xylene	ND		ug/l	2.5	0.70	1				
o-Xylene	ND		ug/l	2.5	0.70	1				
cis-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1				
Styrene	ND		ug/l	2.5	0.70	1				
Dichlorodifluoromethane	ND		ug/l	5.0	1.0	1				
Acetone	ND		ug/l	5.0	1.5	1				
Carbon disulfide	ND		ug/l	5.0	1.0	1				
2-Butanone	ND		ug/l	5.0	1.9	1				
4-Methyl-2-pentanone	ND		ug/l	5.0	1.0	1				
2-Hexanone	ND		ug/l	5.0	1.0	1				
Bromochloromethane	ND		ug/l	2.5	0.70	1				
1,2-Dibromoethane	ND		ug/l	2.0	0.65	1				
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70	1				
Isopropylbenzene	ND		ug/l	2.5	0.70	1				
1,2,3-Trichlorobenzene	ND		ug/l	2.5	0.70	1				
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70	1				
Methyl Acetate	ND		ug/l	2.0	0.23	1				
Cyclohexane	ND		ug/l	10	0.27	1				
1,4-Dioxane	ND		ug/l	250	61.	1				
Freon-113	ND		ug/l	2.5	0.70	1				
Methyl cyclohexane	ND		ug/l	10	0.40	1				

Surrogate	% Recovery	Acceptance Qualifier Criteria	
1,2-Dichloroethane-d4	96	70-130	
Toluene-d8	103	70-130	
4-Bromofluorobenzene	99	70-130	
Dibromofluoromethane	95	70-130	



**Project Name:** SIGNORE

**Project Number:** 21.0056491.81

**SAMPLE RESULTS** 

Lab Number: L2150438

Report Date: 09/24/21

Lab ID: L2150438-12

Client ID: GW-DUP-091621

Sample Location: ELLICOTTVILLE, NY

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 09/22/21 00:17

Analyst: MKS Date Collected: 09/16/21 00:00 Date Received: 09/17/21 Field Prep: Not Specified

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - We	stborough Lab					
Methylene chloride	ND		ug/l	2.5	0.70	1
1,1-Dichloroethane	ND		ug/l	2.5	0.70	1
Chloroform	ND		ug/l	2.5	0.70	1
Carbon tetrachloride	ND		ug/l	0.50	0.13	1
1,2-Dichloropropane	ND		ug/l	1.0	0.14	1
Dibromochloromethane	ND		ug/l	0.50	0.15	1
1,1,2-Trichloroethane	ND		ug/l	1.5	0.50	1
Tetrachloroethene	ND		ug/l	0.50	0.18	1
Chlorobenzene	ND		ug/l	2.5	0.70	1
Trichlorofluoromethane	ND		ug/l	2.5	0.70	1
1,2-Dichloroethane	ND		ug/l	0.50	0.13	1
1,1,1-Trichloroethane	ND		ug/l	2.5	0.70	1
Bromodichloromethane	ND		ug/l	0.50	0.19	1
trans-1,3-Dichloropropene	ND		ug/l	0.50	0.16	1
cis-1,3-Dichloropropene	ND		ug/l	0.50	0.14	1
Bromoform	ND		ug/l	2.0	0.65	1
1,1,2,2-Tetrachloroethane	ND		ug/l	0.50	0.17	1
Benzene	ND		ug/l	0.50	0.16	1
Toluene	ND		ug/l	2.5	0.70	1
Ethylbenzene	ND		ug/l	2.5	0.70	1
Chloromethane	ND		ug/l	2.5	0.70	1
Bromomethane	ND		ug/l	2.5	0.70	1
Vinyl chloride	ND		ug/l	1.0	0.07	1
Chloroethane	ND		ug/l	2.5	0.70	1
1,1-Dichloroethene	ND		ug/l	0.50	0.17	1
trans-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1
Trichloroethene	0.24	J	ug/l	0.50	0.18	1
1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1



**Project Name:** Lab Number: **SIGNORE** L2150438

**Project Number:** Report Date: 21.0056491.81 09/24/21

**SAMPLE RESULTS** 

Lab ID: L2150438-12 Date Collected: 09/16/21 00:00

Client ID: Date Received: 09/17/21 GW-DUP-091621 Sample Location: Field Prep: ELLICOTTVILLE, NY Not Specified

Sample Depth:

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor				
Volatile Organics by GC/MS - Westborough Lab										
1,3-Dichlorobenzene	ND		ug/l	2.5	0.70	1				
1,4-Dichlorobenzene	ND		ug/l	2.5	0.70	1				
Methyl tert butyl ether	ND		ug/l	2.5	0.70	1				
p/m-Xylene	ND		ug/l	2.5	0.70	1				
o-Xylene	ND		ug/l	2.5	0.70	1				
cis-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1				
Styrene	ND		ug/l	2.5	0.70	1				
Dichlorodifluoromethane	ND		ug/l	5.0	1.0	1				
Acetone	ND		ug/l	5.0	1.5	1				
Carbon disulfide	ND		ug/l	5.0	1.0	1				
2-Butanone	ND		ug/l	5.0	1.9	1				
4-Methyl-2-pentanone	ND		ug/l	5.0	1.0	1				
2-Hexanone	ND		ug/l	5.0	1.0	1				
Bromochloromethane	ND		ug/l	2.5	0.70	1				
1,2-Dibromoethane	ND		ug/l	2.0	0.65	1				
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70	1				
Isopropylbenzene	ND		ug/l	2.5	0.70	1				
1,2,3-Trichlorobenzene	ND		ug/l	2.5	0.70	1				
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70	1				
Methyl Acetate	ND		ug/l	2.0	0.23	1				
Cyclohexane	ND		ug/l	10	0.27	1				
1,4-Dioxane	ND		ug/l	250	61.	1				
Freon-113	ND		ug/l	2.5	0.70	1				
Methyl cyclohexane	ND		ug/l	10	0.40	1				

Surrogate	% Recovery	Acceptance Qualifier Criteria	
1,2-Dichloroethane-d4	94	70-130	
Toluene-d8	99	70-130	
4-Bromofluorobenzene	99	70-130	
Dibromofluoromethane	96	70-130	



L2150438

09/24/21

**Project Name: SIGNORE** 

**Project Number:** 21.0056491.81

L2150438-13

TRIP BLANK

ELLICOTTVILLE, NY

**SAMPLE RESULTS** 

Date Collected: 09/16/21 00:00

Lab Number:

Report Date:

Date Received: 09/17/21 Field Prep: Not Specified

Sample Depth:

Sample Location:

Lab ID:

Client ID:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 09/22/21 00:38

Analyst: MKS

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westb	orough Lab					
Methylene chloride	ND		ug/l	2.5	0.70	1
1,1-Dichloroethane	ND		ug/l	2.5	0.70	1
Chloroform	ND		ug/l	2.5	0.70	1
Carbon tetrachloride	ND		ug/l	0.50	0.13	1
1,2-Dichloropropane	ND		ug/l	1.0	0.14	1
Dibromochloromethane	ND		ug/l	0.50	0.15	1
1,1,2-Trichloroethane	ND		ug/l	1.5	0.50	1
Tetrachloroethene	ND		ug/l	0.50	0.18	1
Chlorobenzene	ND		ug/l	2.5	0.70	1
Trichlorofluoromethane	ND		ug/l	2.5	0.70	1
1,2-Dichloroethane	ND		ug/l	0.50	0.13	1
1,1,1-Trichloroethane	ND		ug/l	2.5	0.70	1
Bromodichloromethane	ND		ug/l	0.50	0.19	1
trans-1,3-Dichloropropene	ND		ug/l	0.50	0.16	1
cis-1,3-Dichloropropene	ND		ug/l	0.50	0.14	1
Bromoform	ND		ug/l	2.0	0.65	1
1,1,2,2-Tetrachloroethane	ND		ug/l	0.50	0.17	1
Benzene	ND		ug/l	0.50	0.16	1
Toluene	ND		ug/l	2.5	0.70	1
Ethylbenzene	ND		ug/l	2.5	0.70	1
Chloromethane	ND		ug/l	2.5	0.70	1
Bromomethane	ND		ug/l	2.5	0.70	1
Vinyl chloride	ND		ug/l	1.0	0.07	1
Chloroethane	ND		ug/l	2.5	0.70	1
1,1-Dichloroethene	ND		ug/l	0.50	0.17	1
trans-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1
Trichloroethene	ND		ug/l	0.50	0.18	1
1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1



MDL

**Dilution Factor** 

Project Name: SIGNORE Lab Number: L2150438

**Project Number:** 21.0056491.81 **Report Date:** 09/24/21

**SAMPLE RESULTS** 

Lab ID: L2150438-13 Date Collected: 09/16/21 00:00

Client ID: TRIP BLANK Date Received: 09/17/21 Sample Location: ELLICOTTVILLE, NY Field Prep: Not Specified

Qualifier

Units

RL

Result

Sample Depth:

Parameter

i didilictoi					2	
Volatile Organics by GC/MS - Westbe	orough Lab					
1,3-Dichlorobenzene	ND	ug/l	2.5	0.70	1	
1,4-Dichlorobenzene	ND	ug/l	2.5	0.70	1	
Methyl tert butyl ether	ND	ug/l	2.5	0.70	1	
p/m-Xylene	ND	ug/l	2.5	0.70	1	
o-Xylene	ND	ug/l	2.5	0.70	1	
cis-1,2-Dichloroethene	ND	ug/l	2.5	0.70	1	
Styrene	ND	ug/l	2.5	0.70	1	
Dichlorodifluoromethane	ND	ug/l	5.0	1.0	1	
Acetone	ND	ug/l	5.0	1.5	1	
Carbon disulfide	ND	ug/l	5.0	1.0	1	
2-Butanone	ND	ug/l	5.0	1.9	1	
4-Methyl-2-pentanone	ND	ug/l	5.0	1.0	1	
2-Hexanone	ND	ug/l	5.0	1.0	1	
Bromochloromethane	ND	ug/l	2.5	0.70	1	
1,2-Dibromoethane	ND	ug/l	2.0	0.65	1	
1,2-Dibromo-3-chloropropane	ND	ug/l	2.5	0.70	1	
Isopropylbenzene	ND	ug/l	2.5	0.70	1	
1,2,3-Trichlorobenzene	ND	ug/l	2.5	0.70	1	
1,2,4-Trichlorobenzene	ND	ug/l	2.5	0.70	1	
Methyl Acetate	ND	ug/l	2.0	0.23	1	
Cyclohexane	ND	ug/l	10	0.27	1	
1,4-Dioxane	ND	ug/l	250	61.	1	
Freon-113	ND	ug/l	2.5	0.70	1	
Methyl cyclohexane	ND	ug/l	10	0.40	1	

Surrogate	% Recovery	Acceptance Qualifier Criteria
1,2-Dichloroethane-d4	99	70-130
Toluene-d8	102	70-130
4-Bromofluorobenzene	98	70-130
Dibromofluoromethane	95	70-130



**Project Name:** Lab Number: **SIGNORE** L2150438 **Project Number:** 21.0056491.81

Report Date: 09/24/21

### Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 09/21/21 18:08

Analyst: LAC

arameter	Result	Qualifier Units	RL RL	MDL
olatile Organics by GC/MS	- Westborough Lab	for sample(s):	01-13 Batch:	WG1549552-5
Methylene chloride	ND	ug/l	2.5	0.70
1,1-Dichloroethane	ND	ug/l	2.5	0.70
Chloroform	ND	ug/l	2.5	0.70
Carbon tetrachloride	ND	ug/l	0.50	0.13
1,2-Dichloropropane	ND	ug/l	1.0	0.14
Dibromochloromethane	ND	ug/l	0.50	0.15
1,1,2-Trichloroethane	ND	ug/l	1.5	0.50
Tetrachloroethene	ND	ug/l	0.50	0.18
Chlorobenzene	ND	ug/l	2.5	0.70
Trichlorofluoromethane	ND	ug/l	2.5	0.70
1,2-Dichloroethane	ND	ug/l	0.50	0.13
1,1,1-Trichloroethane	ND	ug/l	2.5	0.70
Bromodichloromethane	ND	ug/l	0.50	0.19
trans-1,3-Dichloropropene	ND	ug/l	0.50	0.16
cis-1,3-Dichloropropene	ND	ug/l	0.50	0.14
Bromoform	ND	ug/l	2.0	0.65
1,1,2,2-Tetrachloroethane	ND	ug/l	0.50	0.17
Benzene	ND	ug/l	0.50	0.16
Toluene	ND	ug/l	2.5	0.70
Ethylbenzene	ND	ug/l	2.5	0.70
Chloromethane	ND	ug/l	2.5	0.70
Bromomethane	ND	ug/l	2.5	0.70
Vinyl chloride	ND	ug/l	1.0	0.07
Chloroethane	ND	ug/l	2.5	0.70
1,1-Dichloroethene	ND	ug/l	0.50	0.17
trans-1,2-Dichloroethene	ND	ug/l	2.5	0.70
Trichloroethene	ND	ug/l	0.50	0.18
1,2-Dichlorobenzene	ND	ug/l	2.5	0.70
1,3-Dichlorobenzene	ND	ug/l	2.5	0.70



L2150438

Project Name: SIGNORE Lab Number:

**Project Number:** 21.0056491.81 **Report Date:** 09/24/21

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 09/21/21 18:08

Analyst: LAC

Parameter	Result	Qualifier Unit	S	RL	MDL	
Volatile Organics by GC/MS - We	stborough Lab	for sample(s):	01-13	Batch:	WG1549552-5	
1,4-Dichlorobenzene	ND	ug/	I	2.5	0.70	
Methyl tert butyl ether	ND	ug/	I	2.5	0.70	
p/m-Xylene	ND	ug/	l	2.5	0.70	
o-Xylene	ND	ug/		2.5	0.70	
cis-1,2-Dichloroethene	ND	ug/	l	2.5	0.70	
Styrene	ND	ug/	l	2.5	0.70	
Dichlorodifluoromethane	ND	ug/	I	5.0	1.0	
Acetone	ND	ug/	I	5.0	1.5	
Carbon disulfide	ND	ug/	I	5.0	1.0	
2-Butanone	ND	ug/	l	5.0	1.9	
4-Methyl-2-pentanone	ND	ug/	l	5.0	1.0	
2-Hexanone	ND	ug/	l	5.0	1.0	
Bromochloromethane	ND	ug/	I	2.5	0.70	
1,2-Dibromoethane	ND	ug/	I	2.0	0.65	
1,2-Dibromo-3-chloropropane	ND	ug/	I	2.5	0.70	
Isopropylbenzene	ND	ug/	l	2.5	0.70	
1,2,3-Trichlorobenzene	ND	ug/	I	2.5	0.70	
1,2,4-Trichlorobenzene	ND	ug/	l	2.5	0.70	
Methyl Acetate	ND	ug/	l	2.0	0.23	
Cyclohexane	ND	ug/		10	0.27	
1,4-Dioxane	ND	ug/		250	61.	
Freon-113	ND	ug/		2.5	0.70	
Methyl cyclohexane	ND	ug/	I	10	0.40	



Project Name: SIGNORE Lab Number: L2150438

**Project Number:** 21.0056491.81 **Report Date:** 09/24/21

Method Blank Analysis
Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 09/21/21 18:08

Analyst: LAC

Parameter Result Qualifier Units RL MDL

Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-13 Batch: WG1549552-5

	Acceptar					
Surrogate	%Recovery Quali	fier Criteria				
1,2-Dichloroethane-d4	98	70-130				
Toluene-d8	99	70-130				
4-Bromofluorobenzene	100	70-130				
Dibromofluoromethane	98	70-130				



# Lab Control Sample Analysis Batch Quality Control

Project Name: SIGNORE

**Project Number:** 

21.0056491.81

Lab Number: L2150438

**Report Date:** 09/24/21

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
/olatile Organics by GC/MS -	Westborough Lab Associated	sample(s):	01-13 Batch:	WG1549552-3	WG1549552-4			
Methylene chloride	100		100		70-130	0		20
1,1-Dichloroethane	110		110		70-130	0		20
Chloroform	100		100		70-130	0		20
Carbon tetrachloride	85		86		63-132	1		20
1,2-Dichloropropane	110		110		70-130	0		20
Dibromochloromethane	78		81		63-130	4		20
1,1,2-Trichloroethane	95		98		70-130	3		20
Tetrachloroethene	100		99		70-130	1		20
Chlorobenzene	100		100		75-130	0		20
Trichlorofluoromethane	100		100		62-150	0		20
1,2-Dichloroethane	94		96		70-130	2		20
1,1,1-Trichloroethane	94		95		67-130	1		20
Bromodichloromethane	86		86		67-130	0		20
trans-1,3-Dichloropropene	91		93		70-130	2		20
cis-1,3-Dichloropropene	93		93		70-130	0		20
Bromoform	70		75		54-136	7		20
1,1,2,2-Tetrachloroethane	94		99		67-130	5		20
Benzene	110		100		70-130	10		20
Toluene	100		100		70-130	0		20
Ethylbenzene	110		100		70-130	10		20
Chloromethane	110		100		64-130	10		20
Bromomethane	82		80		39-139	2		20
Vinyl chloride	130		120		55-140	8		20



### Lab Control Sample Analysis Batch Quality Control

Project Name: SIGNORE

**Project Number:** 21.0056491.81

Lab Number: L2150438

**Report Date:** 09/24/21

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits	
olatile Organics by GC/MS - Westb	oorough Lab Associated	sample(s):	01-13 Batch: V	VG1549552-3	WG1549552-4				
Chloroethane	130		130		55-138	0		20	
1,1-Dichloroethene	95		99		61-145	4		20	
trans-1,2-Dichloroethene	110		100		70-130	10		20	
Trichloroethene	100		100		70-130	0		20	
1,2-Dichlorobenzene	94		98		70-130	4		20	
1,3-Dichlorobenzene	100		100		70-130	0		20	
1,4-Dichlorobenzene	100		99		70-130	1		20	
Methyl tert butyl ether	88		94		63-130	7		20	
p/m-Xylene	105		100		70-130	5		20	
o-Xylene	105		105		70-130	0		20	
cis-1,2-Dichloroethene	97		100		70-130	3		20	
Styrene	105		105		70-130	0		20	
Dichlorodifluoromethane	110		110		36-147	0		20	
Acetone	130		170	Q	58-148	27	Q	20	
Carbon disulfide	110		100		51-130	10		20	
2-Butanone	96		120		63-138	22	Q	20	
4-Methyl-2-pentanone	84		91		59-130	8		20	
2-Hexanone	86		110		57-130	24	Q	20	
Bromochloromethane	97		100		70-130	3		20	
1,2-Dibromoethane	90		89		70-130	1		20	
1,2-Dibromo-3-chloropropane	70		77		41-144	10		20	
Isopropylbenzene	100		100		70-130	0		20	
1,2,3-Trichlorobenzene	85		92		70-130	8		20	



### Lab Control Sample Analysis Batch Quality Control

Project Name: SIGNORE

**Project Number:** 21.0056491.81

Lab Number:

L2150438

Report Date:

Parameter	LCS %Recovery	Qual	LCSD %Recov		%Recovery Limits	RPD	Qual	RPD Limits	
Volatile Organics by GC/MS - Westborough La	•		01-13 Bato	h: WG1549552-3	3 WG1549552-4				
1,2,4-Trichlorobenzene	92		93		70-130	1		20	
Methyl Acetate	83		87		70-130	5		20	
Cyclohexane	110		110		70-130	0		20	
1,4-Dioxane	82		94		56-162	14		20	
Freon-113	110		110		70-130	0		20	
Methyl cyclohexane	100		100		70-130	0		20	

Surrogate	LCS %Recovery Qual	LCSD %Recovery Qual	Acceptance Criteria
1,2-Dichloroethane-d4	97	100	70-130
Toluene-d8	103	102	70-130
4-Bromofluorobenzene	99	99	70-130
Dibromofluoromethane	96	97	70-130

# Matrix Spike Analysis Batch Quality Control

Project Name: SIGNORE

**Project Number:** 21.0056491.81

Lab Number:

L2150438

Report Date:

Parameter	Native Sample	MS Added	MS Found	MS %Recovery	<sup>,</sup> Qual	MSD Found	MSD %Recovery	Qual	Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/MS EW-4.5-091521	S - Westborough	Lab Asso	ciated sample	(s): 01-13 Q0	C Batch ID:	WG15495	552-6 WG1549	9552-7	QC Sample	e: L2150	438-05	Client ID:
Methylene chloride	ND	10	11	110		9.6	96		70-130	14		20
1,1-Dichloroethane	ND	10	12	120		11	110		70-130	9		20
Chloroform	ND	10	10	100		9.5	95		70-130	5		20
Carbon tetrachloride	ND	10	9.4	94		8.8	88		63-132	7		20
1,2-Dichloropropane	ND	10	12	120		10	100		70-130	18		20
Dibromochloromethane	ND	10	8.1	81		7.3	73		63-130	10		20
1,1,2-Trichloroethane	ND	10	9.9	99		9.0	90		70-130	10		20
Tetrachloroethene	0.78	10	12	112		10	92		70-130	18		20
Chlorobenzene	ND	10	10	100		9.5	95		75-130	5		20
Trichlorofluoromethane	ND	10	12	120		10	100		62-150	18		20
1,2-Dichloroethane	ND	10	9.8	98		8.9	89		70-130	10		20
1,1,1-Trichloroethane	ND	10	11	110		9.7	97		67-130	13		20
Bromodichloromethane	ND	10	9.1	91		7.9	79		67-130	14		20
trans-1,3-Dichloropropene	ND	10	9.1	91		8.2	82		70-130	10		20
cis-1,3-Dichloropropene	ND	10	9.2	92		8.2	82		70-130	11		20
Bromoform	ND	10	6.9	69		6.5	65		54-136	6		20
1,1,2,2-Tetrachloroethane	ND	10	9.4	94		9.3	93		67-130	1		20
Benzene	ND	10	11	110		10	100		70-130	10		20
Toluene	ND	10	11	110		10	100		70-130	10		20
Ethylbenzene	ND	10	11	110		10	100		70-130	10		20
Chloromethane	ND	10	12	120		11	110		64-130	9		20
Bromomethane	ND	10	11	110		9.8	98		39-139	12		20
Vinyl chloride	ND	10	14	140		13	130		55-140	7		20



# Matrix Spike Analysis Batch Quality Control

Project Name: SIGNORE

**Project Number:** 21.0056491.81

Lab Number:

L2150438

Report Date:

Parameter	Native Sample	MS Added	MS Found	MS %Recovery	Qual	MSD Found	MSD %Recovery	Qual	Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/MS - EW-4.5-091521	- Westborough	Lab Ass	ociated sample(	s): 01-13 QC	Batch ID: V	VG15495	552-6 WG1549	9552-7	QC Sample	: L2150	)438-05	Client ID:
Chloroethane	ND	10	14	140	Q	14	140	Q	55-138	0		20
1,1-Dichloroethene	ND	10	11	110		10	100		61-145	10		20
trans-1,2-Dichloroethene	ND	10	11	110		9.8	98		70-130	12		20
Trichloroethene	2.4	10	13	106		12	96		70-130	8		20
1,2-Dichlorobenzene	ND	10	9.8	98		9.3	93		70-130	5		20
1,3-Dichlorobenzene	ND	10	9.9	99		9.3	93		70-130	6		20
1,4-Dichlorobenzene	ND	10	10	100		9.4	94		70-130	6		20
Methyl tert butyl ether	ND	10	9.0	90		8.1	81		63-130	11		20
o/m-Xylene	ND	20	23	115		20	100		70-130	14		20
o-Xylene	ND	20	22	110		20	100		70-130	10		20
cis-1,2-Dichloroethene	ND	10	10	100		9.7	97		70-130	3		20
Styrene	ND	20	22	110		19	95		70-130	15		20
Dichlorodifluoromethane	ND	10	12	120		11	110		36-147	9		20
Acetone	ND	10	6.7	67		7.5	75		58-148	11		20
Carbon disulfide	ND	10	12	120		11	110		51-130	9		20
2-Butanone	ND	10	7.5	75		7.0	70		63-138	7		20
4-Methyl-2-pentanone	ND	10	8.9	89		8.0	80		59-130	11		20
2-Hexanone	ND	10	7.7	77		7.3	73		57-130	5		20
Bromochloromethane	ND	10	10	100		8.6	86		70-130	15		20
1,2-Dibromoethane	ND	10	9.2	92		8.4	84		70-130	9		20
1,2-Dibromo-3-chloropropane	ND	10	6.7	67		6.7	67		41-144	0		20
sopropylbenzene	ND	10	11	110		10	100		70-130	10		20
1,2,3-Trichlorobenzene	ND	10	8.8	88		8.4	84		70-130	5		20



# Matrix Spike Analysis Batch Quality Control

Project Name: SIGNORE

**Project Number:** 21.0056491.81

Lab Number:

L2150438

Report Date:

Parameter	Native Sample	MS Added	MS Found	MS %Recovery	/ Qual	MSD Found	MSD %Recovery	Qual	Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/MS - EW-4.5-091521	Westborough I	Lab Assoc	iated sample(s	s): 01-13 Q	C Batch ID:	WG15495	552-6 WG1549	9552-7	QC Sample	: L215	0438-05	Client ID:
1,2,4-Trichlorobenzene	ND	10	9.1	91		8.6	86		70-130	6		20
Methyl Acetate	ND	10	8.8	88		8.6	86		70-130	2		20
Cyclohexane	ND	10	13	130		11	110		70-130	17		20
1,4-Dioxane	ND	500	420	84		380	76		56-162	10		20
Freon-113	ND	10	12	120		12	120		70-130	0		20
Methyl cyclohexane	ND	10	12	120		11	110		70-130	9		20

	MS	MSD	Acceptance
Surrogate	% Recovery Qualifier	% Recovery Qualifier	Criteria
1,2-Dichloroethane-d4	96	99	70-130
4-Bromofluorobenzene	93	98	70-130
Dibromofluoromethane	98	98	70-130
Toluene-d8	101	103	70-130

Serial\_No:09242110:27 *Lab Number:* L2150438

Project Name: SIGNORE
Project Number: 21.0056491.81

Report Date: 09/24/21

### Sample Receipt and Container Information

Were project specific reporting limits specified?

YES

**Cooler Information** 

Cooler Custody Seal

A Absent

Container Information		rmation		Initial	Final	Temp			Frozen	
	Container ID	Container Type	Cooler	рН	pН		Pres	Seal	Date/Time	Analysis(*)
	L2150438-01A	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
	L2150438-01B	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
	L2150438-01C	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
	L2150438-02A	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
	L2150438-02B	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
	L2150438-02C	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
	L2150438-03A	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
	L2150438-03B	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
	L2150438-03C	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
	L2150438-04A	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
	L2150438-04B	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
	L2150438-04C	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
	L2150438-05A	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
	L2150438-05A1	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
	L2150438-05A2	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
	L2150438-05B	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
	L2150438-05B1	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
	L2150438-05B2	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
	L2150438-05C	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
	L2150438-05C1	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
	L2150438-05C2	Vial HCI preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
	L2150438-06A	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
	L2150438-06B	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)



**Lab Number:** L2150438

Report Date: 09/24/21

Project Name: SIGNORE **Project Number:** 21.0056491.81

Container Info	ormation		Initial	Final	Temp			Frozen	
Container ID	Container Type	Cooler	рН	pН	•	Pres	Seal	Date/Time	Analysis(*)
L2150438-06C	Vial HCI preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
L2150438-07A	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
L2150438-07B	Vial HCI preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
L2150438-07C	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
L2150438-08A	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
L2150438-08B	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
L2150438-08C	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
L2150438-09A	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
L2150438-09B	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
L2150438-09C	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
L2150438-10A	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
L2150438-10B	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
L2150438-10C	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
L2150438-11A	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
L2150438-11B	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
L2150438-11C	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
L2150438-12A	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
L2150438-12B	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
L2150438-12C	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
L2150438-13A	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)
L2150438-13B	Vial HCl preserved	Α	NA		5.4	Υ	Absent		NYTCL-8260-R2(14)



**Project Name:** Lab Number: **SIGNORE** L2150438 **Project Number:** 21.0056491.81 **Report Date:** 09/24/21

#### GLOSSARY

#### **Acronyms**

**EDL** 

**EPA** 

LOQ

MS

DL - Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the limit of quantitation (LOQ). The DL includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)

- Estimated Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The EDL includes any adjustments from dilutions, concentrations or moisture content, where applicable. The use of EDLs is specific to the analysis

of PAHs using Solid-Phase Microextraction (SPME).

**EMPC** - Estimated Maximum Possible Concentration: The concentration that results from the signal present at the retention time of an analyte when the ions meet all of the identification criteria except the ion abundance ratio criteria. An EMPC is a worst-case

estimate of the concentration.

LCS - Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of

analytes or a material containing known and verified amounts of analytes.

LCSD Laboratory Control Sample Duplicate: Refer to LCS.

Environmental Protection Agency.

LFB - Laboratory Fortified Blank: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.

LOD - Limit of Detection: This value represents the level to which a target analyte can reliably be detected for a specific analyte in a specific matrix by a specific method. The LOD includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)

- Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats

Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats

MDI - Method Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The MDL includes any adjustments from dilutions, concentrations or moisture content, where applicable.

> - Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for which an independent estimate of target analyte concentration is available. For Method 332.0, the spike recovery is calculated using the native concentration, including estimated values.

MSD - Matrix Spike Sample Duplicate: Refer to MS.

NA - Not Applicable.

NC - Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's

reporting unit.

NDPA/DPA - N-Nitrosodiphenylamine/Diphenylamine.

NI - Not Ignitable.

NP - Non-Plastic: Term is utilized for the analysis of Atterberg Limits in soil.

NR - No Results: Term is utilized when 'No Target Compounds Requested' is reported for the analysis of Volatile or Semivolatile

Organic TIC only requests.

RL - Reporting Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RL

includes any adjustments from dilutions, concentrations or moisture content, where applicable.

RPD - Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the

values; although the RPD value will be provided in the report.

SRM - Standard Reference Material: A reference sample of a known or certified value that is of the same or similar matrix as the

associated field samples.

STLP - Semi-dynamic Tank Leaching Procedure per EPA Method 1315.

TEF - Toxic Equivalency Factors: The values assigned to each dioxin and furan to evaluate their toxicity relative to 2,3,7,8-TCDD.

TEO - Toxic Equivalent: The measure of a sample's toxicity derived by multiplying each dioxin and furan by its corresponding TEF

and then summing the resulting values.

TIC - Tentatively Identified Compound: A compound that has been identified to be present and is not part of the target compound list (TCL) for the method and/or program. All TICs are qualitatively identified and reported as estimated concentrations.

Report Format: DU Report with 'J' Qualifiers



Project Name:SIGNORELab Number:L2150438Project Number:21.0056491.81Report Date:09/24/21

#### **Footnotes**

 The reference for this analyte should be considered modified since this analyte is absent from the target analyte list of the original method.

#### **Terms**

1

Analytical Method: Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum.

Difference: With respect to Total Oxidizable Precursor (TOP) Assay analysis, the difference is defined as the Post-Treatment value minus the Pre-Treatment value

Final pH: As it pertains to Sample Receipt & Container Information section of the report, Final pH reflects pH of container determined after adjustment at the laboratory, if applicable. If no adjustment required, value reflects Initial pH.

Frozen Date/Time: With respect to Volatile Organics in soil, Frozen Date/Time reflects the date/time at which associated Reagent Water-preserved vials were initially frozen. Note: If frozen date/time is beyond 48 hours from sample collection, value will be reflected in 'bold'.

Initial pH: As it pertains to Sample Receipt & Container Information section of the report, Initial pH reflects pH of container determined upon receipt, if applicable.

PAH Total: With respect to Alkylated PAH analyses, the 'PAHs, Total' result is defined as the summation of results for all or a subset of the following compounds: Naphthalene, C1-C4 Naphthalenes, 2-Methylnaphthalene, 1-Methylnaphthalene, Biphenyl, Acenaphthylene, Acenaphthene, Fluorene, C1-C3 Fluorenes, Phenanthrene, C1-C4 Phenanthrenes/Anthracenes, Anthracene, Fluoranthene, Pyrene, C1-C4 Fluoranthenes/Pyrenes, Benz(a)anthracene, Chrysene, C1-C4 Chrysenes, Benzo(b)fluoranthene, Benzo(j)+(k)fluoranthene, Benzo(e)pyrene, Benzo(a)pyrene, Perylene, Indeno(1,2,3-cd)pyrene, Dibenz(ah)+(ac)anthracene, Benzo(g,h,i)perylene. If a 'Total' result is requested, the results of its individual components will also be reported.

PFAS Total: With respect to PFAS analyses, the 'PFAS, Total (5)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA and PFOS. In addition, the 'PFAS, Total (6)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA, PFDA and PFOS. For MassDEP DW compliance analysis only, the 'PFAS, Total (6)' result is defined as the summation of results at or above the RL. Note: If a 'Total' result is requested, the results of its individual components will also be reported.

The target compound Chlordane (CAS No. 57-74-9) is reported for GC ECD analyses. Per EPA, this compound "refers to a mixture of chlordane isomers, other chlorinated hydrocarbons and numerous other components." (Reference: USEPA Toxicological Review of Chlordane, In Support of Summary Information on the Integrated Risk Information System (IRIS), December 1997.)

Total: With respect to Organic analyses, a 'Total' result is defined as the summation of results for individual isomers or Aroclors. If a 'Total' result is requested, the results of its individual components will also be reported. This is applicable to 'Total' results for methods 8260, 8081 and 8082.

#### Data Qualifiers

- A -Spectra identified as "Aldol Condensates" are byproducts of the extraction/concentration procedures when acetone is introduced in the process.
- The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For MCP-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentrations of the analyte at less than ten times (10x) the concentrations of the analyte was detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank AND the analyte was detected above one-half the reporting limit (or above the reporting limit for common lab contaminants) in the associated method blank. For NJ-Air-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte above the reporting limit. For NJ-related projects (excluding Air), flag only applies to associated field samples that have detectable concentrations of the analyte, which was detected above the reporting limit in the associated method blank or above five times the reporting limit for common lab contaminants (Phthalates, Acetone, Methylene Chloride, 2-Butanone).
- Co-elution: The target analyte co-elutes with a known lab standard (i.e. surrogate, internal standards, etc.) for co-extracted analyses.
- Concentration of analyte was quantified from diluted analysis. Flag only applies to field samples that have detectable concentrations of the analyte.
- E Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.
- F The ratio of quantifier ion response to qualifier ion response falls outside of the laboratory criteria. Results are considered to be an estimated maximum concentration.
- G The concentration may be biased high due to matrix interferences (i.e, co-elution) with non-target compound(s). The result should be considered estimated.
- H The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.
- I The lower value for the two columns has been reported due to obvious interference.
- Estimated value. The Target analyte concentration is below the quantitation limit (RL), but above the Method Detection Limit (MDL) or Estimated Detection Limit (EDL) for SPME-related analyses. This represents an estimated concentration for Tentatively Identified Compounds (TICs).
- $\label{eq:main_equation} \textbf{M} \qquad \text{-Reporting Limit (RL) exceeds the MCP CAM Reporting Limit for this analyte.}$
- ND Not detected at the method detection limit (MDL) for the sample, or estimated detection limit (EDL) for SPME-related analyses.

Report Format: DU Report with 'J' Qualifiers



Project Name:SIGNORELab Number:L2150438Project Number:21.0056491.81Report Date:09/24/21

#### **Data Qualifiers**

- NJ Presumptive evidence of compound. This represents an estimated concentration for Tentatively Identified Compounds (TICs), where the identification is based on a mass spectral library search.
- P The RPD between the results for the two columns exceeds the method-specified criteria.
- Q -The quality control sample exceeds the associated acceptance criteria. For DOD-related projects, LCS and/or Continuing Calibration Standard exceedences are also qualified on all associated sample results. Note: This flag is not applicable for matrix spike recoveries when the sample concentration is greater than 4x the spike added or for batch duplicate RPD when the sample concentrations are less than 5x the RL. (Metals only.)
- **R** Analytical results are from sample re-analysis.
- RE Analytical results are from sample re-extraction.
- S Analytical results are from modified screening analysis.
- The surrogate associated with this target analyte has a recovery outside the QC acceptance limits. (Applicable to MassDEP DW Compliance samples only.)
- The batch matrix spike and/or duplicate associated with this target analyte has a recovery/RPD outside the QC acceptance limits.
   (Applicable to MassDEP DW Compliance samples only.)

Report Format: DU Report with 'J' Qualifiers



Project Name:SIGNORELab Number:L2150438Project Number:21.0056491.81Report Date:09/24/21

#### REFERENCES

Test Methods for Evaluating Solid Waste: Physical/Chemical Methods. EPA SW-846. Third Edition. Updates I - VI, 2018.

### **LIMITATION OF LIABILITIES**

Alpha Analytical performs services with reasonable care and diligence normal to the analytical testing laboratory industry. In the event of an error, the sole and exclusive responsibility of Alpha Analytical shall be to re-perform the work at it's own expense. In no event shall Alpha Analytical be held liable for any incidental, consequential or special damages, including but not limited to, damages in any way connected with the use of, interpretation of, information or analysis provided by Alpha Analytical.

We strongly urge our clients to comply with EPA protocol regarding sample volume, preservation, cooling, containers, sampling procedures, holding time and splitting of samples in the field.



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Alpha Analytical, Inc. Facility: Company-wide

Department: Quality Assurance

Title: Certificate/Approval Program Summary

### Certification Information

The following analytes are not included in our Primary NELAP Scope of Accreditation:

Westborough Facility

EPA 624/624.1: m/p-xylene, o-xylene, Naphthalene

EPA 625/625.1: alpha-Terpineol

EPA 8260C/8260D: NPW: 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene, Azobenzene; SCM: Iodomethane (methyl iodide), 1,2,4,5-Tetramethylbenzene;

EPA 8270D/8270E: NPW: Dimethylnaphthalene,1,4-Diphenylhydrazine, alpha-Terpineol; SCM: Dimethylnaphthalene,1,4-Diphenylhydrazine.

SM4500: NPW: Amenable Cyanide; SCM: Total Phosphorus, TKN, NO2, NO3.

**Mansfield Facility** 

**SM 2540D:** TSS

EPA 8082A: NPW: PCB: 1, 5, 31, 87,101, 110, 141, 151, 153, 180, 183, 187.

EPA TO-15: Halothane, 2,4,4-Trimethyl-2-pentene, 2,4,4-Trimethyl-1-pentene, Thiophene, 2-Methylthiophene,

3-Methylthiophene, 2-Ethylthiophene, 1,2,3-Trimethylbenzene, Indan, Indene, 1,2,4,5-Tetramethylbenzene, Benzothiophene, 1-Methylnaphthalene.

Biological Tissue Matrix: EPA 3050B

The following analytes are included in our Massachusetts DEP Scope of Accreditation

Westborough Facility:

**Drinking Water** 

EPA 300.0: Chloride, Nitrate-N, Fluoride, Sulfate; EPA 353.2: Nitrate-N, Nitrite-N; SM4500NO3-F: Nitrate-N, Nitrite-N; SM4500F-C, SM4500CN-CE,

EPA 180.1, SM2130B, SM4500CI-D, SM2320B, SM2540C, SM4500H-B, SM4500NO2-B

EPA 332: Perchlorate; EPA 524.2: THMs and VOCs; EPA 504.1: EDB, DBCP.

Microbiology: SM9215B; SM9223-P/A, SM9223B-Colilert-QT,SM9222D.

Non-Potable Water

SM4500H,B, EPA 120.1, SM2510B, SM2540C, SM2320B, SM4500CL-E, SM4500F-BC, SM4500NH3-BH: Ammonia-N and Kjeldahl-N, EPA 350.1: Ammonia-N, LACHAT 10-107-06-1-B: Ammonia-N, EPA 351.1, SM4500NO3-F, EPA 353.2: Nitrate-N, SM4500P-E, SM4500P-B, E, SM4500SO4-E, SM5220D, EPA 410.4, SM5210B, SM5310C, SM4500CL-D, EPA 1664, EPA 420.1, SM4500-CN-CE, SM2540D, EPA 300: Chloride, Sulfate, Nitrate.

EPA 624.1: Volatile Halocarbons & Aromatics,

EPA 608.3: Chlordane, Toxaphene, Aldrin, alpha-BHC, beta-BHC, gamma-BHC, delta-BHC, Dieldrin, DDD, DDE, DDT, Endosulfan II, Endosulfan II,

Endosulfan sulfate, Endrin, Endrin Aldehyde, Heptachlor, Heptachlor Epoxide, PCBs

EPA 625.1: SVOC (Acid/Base/Neutral Extractables), EPA 600/4-81-045: PCB-Oil.

Microbiology: SM9223B-Colilert-QT; Enterolert-QT, SM9221E, EPA 1600, EPA 1603, SM9222D.

Mansfield Facility:

**Drinking Water** 

EPA 200.7: Al, Ba, Cd, Cr, Cu, Fe, Mn, Ni, Na, Ag, Ca, Zn. EPA 200.8: Al, Sb, As, Ba, Be, Cd, Cr, Cu, Pb, Mn, Ni, Se, Ag, TL, Zn. EPA 245.1 Hg.

EPA 522, EPA 537.1.

Non-Potable Water

EPA 200.7: Al, Sb, As, Be, Cd, Ca, Cr, Co, Cu, Fe, Pb, Mg, Mn, Mo, Ni, K, Se, Ag, Na, Sr, TL, Ti, V, Zn.

EPA 200.8: Al, Sb, As, Be, Cd, Cr, Cu, Fe, Pb, Mn, Ni, K, Se, Ag, Na, TL, Zn.

EPA 245.1 Hg

SM2340B

For a complete listing of analytes and methods, please contact your Alpha Project Manager.

Pre-Qualtrax Document ID: 08-113 Document Type: Form

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GZA GeoEnvironmental, Inc.