

2023

PERIODIC REVIEW REPORT

**FOR
FORMER MOBIL SERVICE STATION 99-MST
979 MAIN STREET (1001 MAIN STREET)
NYSDEC SITE #C915260
CITY OF BUFFALO, ERIE COUNTY, NEW YORK**

Prepared by:



C&S ENGINEERS, INC.
141 ELM STREET
BUFFALO, NEW YORK 14203

Prepared on Behalf of:

**KALEIDA HEALTH
KALEIDA PROPERTIES, INC.
SEAVEST CORE BUFFALO CONVENTUS, LLC**

APRIL 2023

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ACRONYM LIST

C&S	C&S ENGINEERS, INC.
BGS	BELOW GROUND SURFACE
BCP	BROWNFIELD CLEANUP PROGRAM
BCA	BROWNFIELD CLEANUP AGREEMENT
BTEX	BENZENE, TOLUENE, ETHYLBENZENE AND XYLENE
DUSR	DATA USABILITY AND SUMMARY REPORT
LNAPL	LIGHT NON-AQUEOUS PHASE LIQUID
IRM	INTERIM REMEDIAL MEASURES
NYSDEC	NEW YORK STATE DEPARTMENT OF ENVIRONMENTAL CONSERVATION
PCOC	PRIMARY CONTAMINATE OF CONCERN
PID	PHOTO-IONIZATION DETECTOR
PPM	PARTS PER MILLION
RI	REMEDIAL INVESTIGATION
SCO	SOIL CLEANUP OBJECTIVES
SMP	SITE MANAGEMENT PLAN
SVOC	SEMI-VOLATILE ORGANIC COMPOUNDS
VOC	VOLATILE ORGANIC COMPOUNDS

EXECUTIVE SUMMARY

C&S Engineers, Inc. (C&S) has prepared this 2022 - 2023 Periodic Review Report for the former Mobil Service Station 99-MST - 979 Main Street (1001 Main Street) (hereinafter referred to as the Site) located at 1001 Main Street in Buffalo, New York.

The Site was remediated in accordance with Brownfield Cleanup Agreement (BCA) Index #C915260-03-12, Site #C915260, which was executed on June 15, 2012 and amended on:

- August 7, 2012 to modify the BCP boundary increasing the size of the Site;
- January 3, 2014 to add Conventus Partners, LLC as an owner; and
- November 7, 2014 to change SBL numbers for the Site.

BCA Volunteers included Kaleida Properties Inc., Kaleida Health and F.L.C 50 High Street Corporation. In December 2020, Conventus Partners, LLC and Kaleida Properties Inc. entered a ground lease for the property with Seavest Core Buffalo Conventus LLC to operate and maintain the property. A figure showing the Site location and boundaries is provided in **Figure 1** and **Figure 2**.

Remedial activities consisted of installing steel shoring around the property and removing contaminated soil and groundwater to approximately 26 – 40 feet below ground surface. After completion of the remedial work, some contamination remained in the subsurface at this Site. A Site Management Plan (SMP) was prepared on November 28, 2014 to manage remaining groundwater contamination at the Site until the Environmental Easement is extinguished in accordance with ECL Article 71, Title 36.

Petroleum contaminated groundwater is present within a discontinuous layer of coarse sand and gravel located between 32 and 35 feet below ground surface. This layer generally ranges from 6 inches to three feet thick, provides a preferential pathway for groundwater flow, and is confined within dense silt and fine sand present above and below the groundwater bearing zone.

During the remedial efforts, seven groundwater monitoring wells were installed prior to the installation of the two floors of underground parking. These monitoring wells were used to conduct in-situ injections by gravity feeding chemical oxidants into the groundwater bearing zone. A total of 2,480 pounds of chemical oxidant was used over three treatment events. Treatments occurred from December 2013 to June 2015. Groundwater samples following the in-situ injections show minor reductions in petroleum compounds.

In 2016, C&S conducted a limited groundwater extraction on the wells with the highest contaminant levels. Contaminated groundwater was pumped from the wells and treated with 200 pounds of activated carbon before discharging into the sanitary sewer. A total of 4,762.2 gallons of contaminated groundwater was removed. Groundwater samples collected in December 2015, January 2016 and March 2016 showed a slight reduction in petroleum compound concentrations.

The current ISCO treatment method is smaller pressurized injections around each target location on a periodic schedule. A total of six temporary PVC injection points were installed around BCP-MW-6 and BCP-MW-5. Each treatment injects a total of 800 pounds (130 pounds per injection point) of chemical oxidant. Groundwater monitoring is conducted biannually.

All institutional and engineering controls are in compliance with the SMP. To address the continued elevated concentrations of petroleum compounds in the groundwater, C&S recommends the completion of additional treatment methods, including the implementation of a slow release chemical oxidation method.

The Institutional and Engineering Controls Certification form is provided in **Appendix B**.

1 SITE OVERVIEW

1.1 Site Description

The Site is located in the City of Buffalo County of Erie County, New York and is identified below on the Erie County Tax Map.

SBL: 100.79 – 1- 1.1

Street Number: 1001 Main Street, Buffalo

(formerly 979 Main Street)

Land Owner:

SBL: 100.79 – 1- 1.1/1

Street Number: 1001 Main Street, Buffalo

(formerly 979 Main Street)

Land Owner: Seavest Core Buffalo Conventus, LLC

SBL: 100.79 – 1- 1.1/2

Street Number: 1001 Main Street, Buffalo
(formerly 979 Main Street)

1st – 6th Floor Building Owner: Seavest Core Buffalo Conventus, LLC

SBL: 100.79 – 1- 1.1/3

Street Number: 1001 Main Street, Buffalo
(formerly 979 Main Street)

7th Floor Building Owner: Fort Schuyler Management Corp. (an entity New York State).

The Site is an approximately 1.72-acre area bounded by Goodrich Street to the north, High Street to the south, parking lot to the east, and Main Street to the west (see **Figure 1 and 2**).

1.2 Geology and Hydrogeology

The Conventus Medical Office Building currently occupies the Site. During remedial activities, steel shoring was installed to a depth of 40 to 50 feet below grade around the Site. Across the majority of the Site, soils were excavated to 26 feet below ground surface (bgs). Two floors of underground parking were constructed underneath the Conventus building.

The Site geology begins at 26 feet bgs. Subsurface soils consist of dry to moist fine sand and silt formation extends to nearly 70 feet bgs. Below this massive sand and silt formation is a discontinuous coarse sand and gravel layer that grades to a sand, gravel; and clay till formation. Underlying the overburden is a grey cherty limestone formation at approximately 90 feet bgs.

The principal groundwater bearing zone beneath the Site is located within the coarse sand and gravel layer between 32 and 35 feet bgs. This layer is of variable thickness (generally 6 inches to three feet) but is horizontally discontinuous. The layer is located within the central and northeastern portions of the Site, but does not extend completely to the southern, northwestern or southeastern areas of the Site

and is confined by the dense fine sands and silt above and below the groundwater bearing zone.

1.3 Nature and Extent of Contamination

During the Interim Remedial Measure (IRM), grossly contaminated soil and groundwater were removed from the Site. In total, 67,458 tons of soils were sent for disposal or treatment due to gasoline contamination. The remaining contamination left on-site consists of petroleum impacted groundwater. Groundwater sampling that occurred prior to the IRM confirmed that the Primary Contaminants of Concern (PCOCs) are limited to petroleum hydrocarbons.

Groundwater flows within the coarse sand/gravel groundwater bearing zone to the northeast. Groundwater recharge from the surface has been eliminated due to the concrete floor of the parking garage, which effectively covers 100% of the Site recharge area. Additionally, below grade migration has been effectively stopped by the presence of deep sheet piling that cuts off the groundwater bearing zone from the remaining off-site formation around the majority of the Site. The lack of a vertical recharge from the surface and the horizontal containment in the groundwater bearing zone was designed to contain the remaining groundwater on-site and reduces the future contaminant loading into the surrounding off-site formation. However, a small gap in the sheet piling along the southwestern corner may provide a route for off-site contamination to impact the Site's groundwater.

1.4 Site History

Contamination is related to the historic use of the property as a gas station and originally was sourced from leaking underground storage tanks located above the "Deep Excavation Area" (see **Figure 3**).

For over 40 years, the light non-aqueous phase liquid (LNAPL) filtered downward from the base of the tank to a depth of approximately 40 feet bgs. LNAPL intercepted the groundwater at approximately 32 feet bgs. The water table is present within a semi-confined coarse sand and gravel lens. This lens varies in thickness (1/2 to 3 feet) and extends to the northeast, confined laterally to the east and west. Because of low carbon in the fine sand silt and gravel formations, breakdown of benzene, toluene, ethylbenzene and xylene (BTEX) compounds was slow. This resulted in high volatile organic compounds (VOC) soil gas in the unsaturated zone below the release area and the continual loading of BTEX into the groundwater from the LNAPL. Soil Contamination (exceeding Residential Use SCOs), below the LNAPL layer was noted to extend to a depth of 35 to 40 feet bgs. This area has been identified as the Source Area for groundwater contamination.

Dissolved BTEX, once entering the groundwater bearing zone was transported via localized, preferential groundwater flow to the northeast corner of the Site (Following the location of the coarse sand/gravel lens).

To redevelop the property into a medical office building, the Applicants (BCP F.L.C. 50 High Street, Corporation, Kaleida Health, Kaleida Properties, Inc. and Conventus Partners, LLC) acting as Brownfield Cleanup Program (BCP) Volunteers, submitted a BCP Application for the Site on November 28, 2011. The Applicants and the New York State Department of Environmental Conservation (NYSDEC) signed the Brownfield Cleanup Agreement (BCA) on June 15, 2012.

The NYSDEC approved IRM was implemented on January 2013. The following is a summary of the IRM performed at the Site:

1. Excavation of soil/fill exceeding restricted residential SCOs to 26 feet bgs;
2. Excavation of soil from the Source Area to 40 feet bgs;
3. Removal of LNAPL and contaminated groundwater;
4. Backfilling with clean fill and construction of concrete floor;
5. Backfilling the Source Area with flowable fill; and
6. Execution and recording of an Environmental Easement to restrict land use and prevent future exposure to any contamination remaining at the Site.

The removal of soils in the Source Area (“Area of Deep Excavation” in **Figure 3**) also included the removal of the groundwater bearing zone. During soil removal, 1,997 tons of groundwater and LNAPL was removed from the excavation and properly disposed off-site. The groundwater bearing zone within the Source Area was replaced with flowable fill, sealing this area off from the adjacent groundwater bearing zone beneath the Site.

Remedial activities were completed at the Site in October 2013. Implementation of the IRM, including source removal, was effective in removing any remaining free product grossly contaminated soils and the groundwater containing the highest dissolved BTEX. However, residual groundwater contamination remains on-site.

Following mass excavation activities, seven new wells were installed on-site.

Table 1-1: Post-Remediation Wells

Well ID	Diameter
BCP-MW-1	2”

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BCP-MW-2	8"
BCP-MW-3	8"
BCP-MW-4	2"
BCP-MW-5	2"
BCP-MW-6	8"
BCP-MW-7	2"

Note that one well (BCP-MW-2) was installed adjacent to the flowable fill within the Source Area. This well did not produce water. A second well, BCP-MW-6, was installed along the western side of the deep excavation, along the tiered excavation area and did intercept the portion of the groundwater bearing zone remaining along the shoring. This well did produce water for sampling. All other wells were installed through native materials and the gravel water bearing layer. All wells were installed to an approximate depth of 43 feet below surrounding grade (approximately 16 feet below basement floor elevation).

The monitoring well locations were located in areas of previously identified groundwater contamination and to the south of the plume to confirm that contamination had not moved off-site to the south.

BCP-MW-2 was installed adjacent to the Source Area that was backfilled with flowable fill. Since its installation, this well has been dry. NYSDEC requested the well be modified to evaluate if groundwater underneath the flowable fill mass contains residual contamination. On October 7, 2015 Nature's Way Environmental installed a 1-inch PVC well through the existing BCP-MW-2 to a final depth of 50 feet bgs. The modified well has remained dry. This provides additional evidence that groundwater and petroleum contamination are limited to the coarse sand and gravel layer 32 to 35 feet bgs.

1.4.1 Pump and Treat Pilot Studies

Limited groundwater extraction was conducted on-site to determine if contaminant levels could be reduced using pump and treat techniques. The first pilot study was implemented from December 2015 to March 2016. A second pilot study was implemented from November 2016 to January 2017.

Groundwater was removed using submersible electric pumps. The total volume of groundwater removed was 4,762.2 gallons during the first event and 6,161 gallons during the second event.

Groundwater sampling was conducted before and after the groundwater withdrawal events. Based on sampling results, the Pump-and-Treat Pilot studies did not appear to have a significant effect on the groundwater contaminant concentrations.

1.4.2 In-situ Injections

The remedial method selected for the Site was in-situ chemical oxidation (ISCO) using RegenOX manufactured by Regenesis. RegenOX is sodium percarbonate formulated to degrade petroleum hydrocarbons through direct oxidation and through the generation of free radical compounds which will also oxidize contaminants. RegenOx produces minimal heat and pressure and is non-corrosive, making it a relatively safe chemical oxidant that is compatible for use in direct contact with underground infrastructure such as utilities, tanks, piping, and communication lines. This was an important characteristic when selecting the ISCO product due to the close proximity of the monitoring wells to the earth retention sheeting for the Conventus Building.

The amount of RegenOX used was calculated based on Site specific data and professional experience of C&S and Regenesis. RegenOX was mixed with tap water in 55 gallon drums at a concentration of 100 pounds of RegenOX with 110 gallons of water for each location.

In-situ treatment consisted of gravity-feeding a chemical oxidizer mixed with water directly into monitoring wells, BCP-MW-3, BCP-MW-4, BCP-MW-5, and BCP-MW-6,. Groundwater samples were collected approximately three months after treatment. The first ISCO treatment was conducted on December 12, 2013.

Evaluation of the gravity fed treatments determined this method was not effective at reducing groundwater contaminants. A work plan for increasing the amount of treatment solution using pressure injections was developed. Borings were advanced in the lower floor of underground parking to apply in-situ treatments under pressure directly into the contaminated sand and gravel lens.

The ISCO solution was directly injected into the soil in 12 borings in the sub-basement. Three borings were advanced adjacent to each monitoring wells listed below:

- BCP-MW-3
- BCP-MW-5
- BCP-MW-4
- BCP-MW-6

Each injection boring had to be carefully located to avoid hitting utilities located underneath the floor, with the intent of being within 10 to 15 feet of each monitoring well. Each injection boring was advanced into the coarse sand and gravel layer, approximately 15 feet below the concrete floor.

The ISCO solution was pumped from the mixing station to a truck mounted geo-probe and into the subsurface. The mix of RegenOX and water was injected under pressure in each boring, and the 12 injection borings received approximately 100 pounds of RegenOx. Additionally, 100 pounds of ISCO material was gravity fed directly into each monitoring well. A total of 1,600 pounds of RegenOx was used for each treatment event. For two treatments, a total of 3,200 pounds of RegenOX was used. These large treatment events resulted in mixed results; some locations showed an increase in contaminant concentrations, likely due to additional petroleum desorption, other locations indicated a significant decrease of petroleum contaminants.

The starting in 2018 ISCO treatment method changed to smaller pressurized injections around each target location on a quarterly schedule. A total of six temporary PVC injection points were installed around BCP-MW-6 and BCP-MW-5. Each quarterly treatment injects a total of 800 pounds (130 pounds per injection point) of chemical oxidant. Groundwater monitoring is conducted biannually.

The following provides all ISCO injections that have occurred to date:

December 12, 2013
July 28, 2014
June 24, 2015
May 24 – 25, 2017
September 13 - 15, 2017
May 31, 2018
July 6, 2018
October 1, 2018
January 8 – 11, 2019
June 11 - 12, 2019
October 16 – 17, 2019
November 12 – 13, 2019
March 31 – April 1, 2020
June 23 – 24, 2020
September 21 – 24, 2020
January 5 – 8, 2021
June 1 – 4, 2021

August 13 – 14, 2021

October 13 – 14, 2021

2 REMEDY PERFORMANCE, EFFECTIVENESS AND PROTECTIVENESS

The table below presents a comparison of total VOC and BTEX concentrations from each monitoring well and the percent change from pre-treatment and post-treatment groundwater monitoring.

Table 2-1: VOC Concentration Change

<i>Monitoring Well</i>	<i>Percent Change Post Injections April 2022 to August 2022</i>	<i>Percent Change Post Injections August 2022 to December 2022</i>	<i>Percent Change Post Remediation Maximum to December 2022</i>
BCP MW-1	-100	-100	-100
BCP MW-3	+2018	-98	-100
BCP MW-4	+4	+12	-29
BCP MW-5	+3003	-61	-68
BCP MW-6	-100	-100	-100
BCP MW-7	-100	-100	-100

Note: Negative value indicates decrease in concentration and positives value indicates increase in concentration
 BCP-MW-2 was dry. No samples were collected.

Table 2-2: BTEX Concentration Change

<i>Monitoring Well</i>	<i>Percent Change Post Injections April 2022 to August 2022</i>	<i>Percent Change Post Injections August 2022 to December 2022</i>	<i>Percent Change Post Remediation Maximum to December 2022</i>
BCP MW-1	-100	-100	-100
BCP MW-3	+2162	-97	-100
BCP MW-4	+22	+20	-67
BCP MW-5	+4950	-67	-81
BCP MW-6	-100	-100	-100
BCP MW-7	-100	-100	-100

Note: Negative value indicates decrease in concentration and positives value indicates increase in concentration
 BCP-MW-2 was dry. No samples were collected.

Table 2-1 and **Table 2-2** shows several consistent decreases in contaminant concentrations from April 2022 to December 2022. The tables indicate a rebound in BCP-MW-3 and BCP-MW-5 and to a lesser extent BCP-MW-4. It should be noted that BCP-MW-5 has consistently had the highest VOC concentrations over the past few years. These increases have been observed during the in-situ injection program

multiple times. The injection product has chemicals that actively desorb VOCs from the soil and still maybe having an effect over time. This temporary increases the dissolved VOC concentrations in the groundwater while other chemicals degrade the VOCs. In addition, injection product has a significant delayed reduction on contaminant concentrations is due to nearly zero groundwater movement within the footprint of the Conventus Building. Permanent steel shoring has cut off the groundwater underneath the building resulting in very slow movement of treatment solution to contaminated areas.

Graph 1 shows total BTEX concentrations over time. **Figure 3** shows the historic BTEX concentrations from each well.

Despite the increase in VOC concentrations in three out of the seven monitoring wells the remedy is demonstrating effectiveness and protectiveness for the following reasons:

- The residual contamination is over 30 feet below ground surface or seven to ten feet below the surface of the lowest parking floor. The general public would not be exposed to contaminated water. Most utilities in the area are much shallower than the contamination; therefore, the risk to construction workers is low.
- Based on previous investigations conducted by C&S (including downgradient sampling conducted 2016), the residual contamination does not extend offsite. The residual contamination is confined to the footprint of the building due to the steel shoring cutting off groundwater flow.
- Exposure to any potential soil vapor is significantly mitigated to the occupants of the building and surrounding area due to two floors of underground parking designed with ventilation systems, vapor barrier on the lower floor and a passive sub-slab depressurization system.

Considering that the residual contamination is isolated both vertically and horizontally on the Site, the remedy is protective of the public and environment.

3 IC/EC PLAN COMPLIANCE REPORT

3.1 IC/EC Requirements and Compliance

As stated in the 2014 Decision Document, the remedial action objectives (RAO) selected for this Site are:

3.1.1 Groundwater

RAOs for Public Health Protection

- Prevent ingestion of groundwater with contaminant levels exceeding drinking water standards.

- Prevent contact with, or inhalation of volatiles, from contaminated groundwater.

RAOs for Environmental Protection

- Restore ground water aquifer to pre-disposal/pre-release conditions, to the extent practicable.
- Prevent the discharge of contaminants to surface water.

3.1.2 Soil

RAOs for Public Health Protection

- Prevent ingestion/direct contact with contaminated soil.
- Prevent inhalation of or exposure from contaminants volatilizing from contaminants in soil.

RAOs for Environmental Protection

- Prevent migration of contaminants that would result in groundwater or surface water contamination.

3.1.3 Soil Vapor

RAOs for Public Health Protection

- Mitigate impacts to public health resulting from existing, or the potential for, soil vapor intrusion directly into buildings at a site.

3.1.4 Institutional Controls

The institutional controls for this Site are:

- The Site may only be used for restricted residential use provided that the long-term Engineering and Institutional Controls included in this SMP are employed;
- The Site may not be used for a higher level of use, unrestricted or residential use, without additional remediation and amendment of the Environmental Easement, as approved by the NYSDEC;
- All future activities on the Site that will disturb remaining contaminated material must be conducted in accordance with this SMP;
- The use of the groundwater underlying the Site is prohibited by the City of Buffalo; and
- Vegetable gardens and farming on the Site are prohibited.

The Site has not changed owners and the land use of the Site has not change. All intuitional controls for this Site are in accordance with requirements of the Environmental Easement.

3.1.5 Engineering Controls

The engineering controls for this Site are:

- Groundwater treatment and monitoring using the seven wells installed in the sub-basement of the building

All engineering controls for this Site are in accordance with requirements of the Environmental Easement.

3.2 IC/EC Certification

As required, the Site Management Periodic Review Report Notice – Institutional and Engineering Controls Certificate Form has been completed and a copy is provided in **Appendix B**.

4 MONITORING PLAN COMPLIANCE REPORT

The SMP identified the need for continued monitoring of groundwater conditions at the Site, including the periodic measuring of water levels and collecting groundwater samples for VOC analysis.

The following monitoring wells are included in the groundwater monitoring plan:

- BCP-MW-1
- BCP-MW-2
- BCP-MW-3
- BCP-MW-4
- BCP-MW-5
- BCP-MW-6
- BCP-MW-7

All monitoring wells were sampled with the exception of BCP-MW-2, which has remained dry since its installation.

The groundwater monitoring activities included the collection of depth-to-water measurements at each monitoring well and the collection of groundwater samples for laboratory analysis. Groundwater sampling was conducted in accordance with the U.S. Environmental Protection Agency Low flow sample procedure. Groundwater sample occurred on the dates below:

September 20, 2013 March 22, 2016 November 30, 2018 December 14, 2022

March 19, 2014	June 3, 2016	July 30, 2019
May 22, 2014	October 25, 2016	December 4, 2019
March 11, 2015	December 8, 2016	May 13, 2020
June 17, 2015	January 20, 2017	November 25, 2020
August 3, 2015	May 17, 2017	May 14, 2021
October 7, 2015	July 5, 2017	December 14, 2021
December 14, 2015	November 2, 2017	April 7, 2022
January 27, 2016	August 18, 2018	August 25, 2022

Figure 3 shows the location of the groundwater wells in the sub-basement of the Conventus building.

Table 1A presents detected compounds over previous monitoring events. **Table 1B** provides sample results for this reporting period.

5 OPERATION AND MAINTENANCE PLAN COMPLIANCE

The only maintenance items are those associated with the monitoring wells. Minor maintenance to the well caps, PVC risers and road boxes is recommended for some of the monitoring wells. These issues do not interfere with groundwater monitoring or the integrity of the samples.

6 CONCLUSIONS

Based upon the remedial activities performed, the following conclusions have been formulated:

- All of the required work was completed and is reported herein.
- The remedial activities performed at the Site have prevented any adverse risk to human health and the environment.

7 RECOMMENDATIONS

In November of 2021, C&S and the NYSDEC discussed the possibility to reduce the in-situ injection activities. C&S recommends the following:

- Groundwater sampling will be conducted annually on the all monitoring wells in the sub-basement of the Site. All groundwater samples will be collected for VOCs and analyzed using EPA Method 8260.
- If VOC concentrations show a trend upward over three monitoring events, then the in-situ injection program will be re-started.

**Periodic Review Report
Former Mobil Service Station 99-MST
979 Main Street (1001 Main Street), BCP No. C915260**

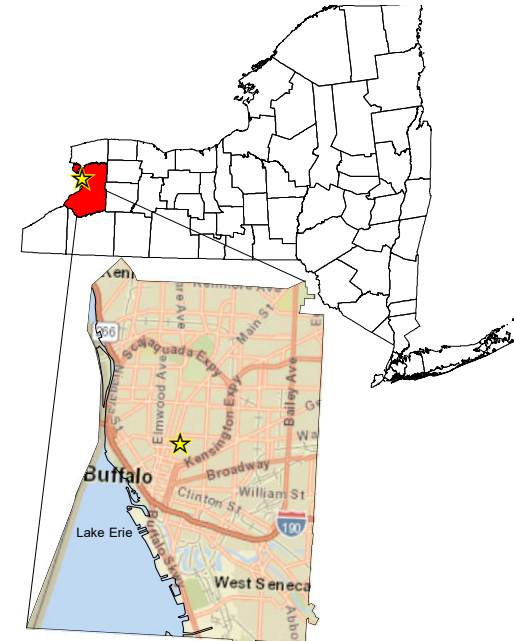
The NYSDEC denied changing to an annual sampling program. The injection program was temporarily halted unless a continued upward trend is identified.

Two monitoring wells show an upward trend. The next rounds of sampling will determine if the injection program will need to be restarted.

Semi-annual monitoring will continue for the Site until conditions change.

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FIGURES



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**FORMER MOBIL STATION 99-MST
979 MAIN ST (1001 MAIN ST)**

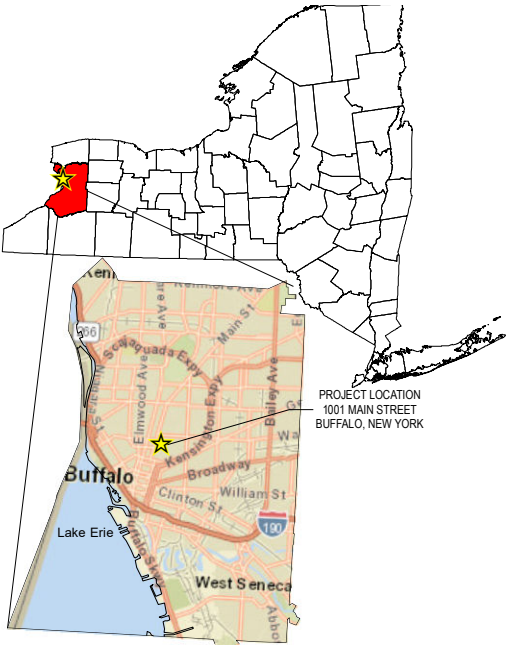
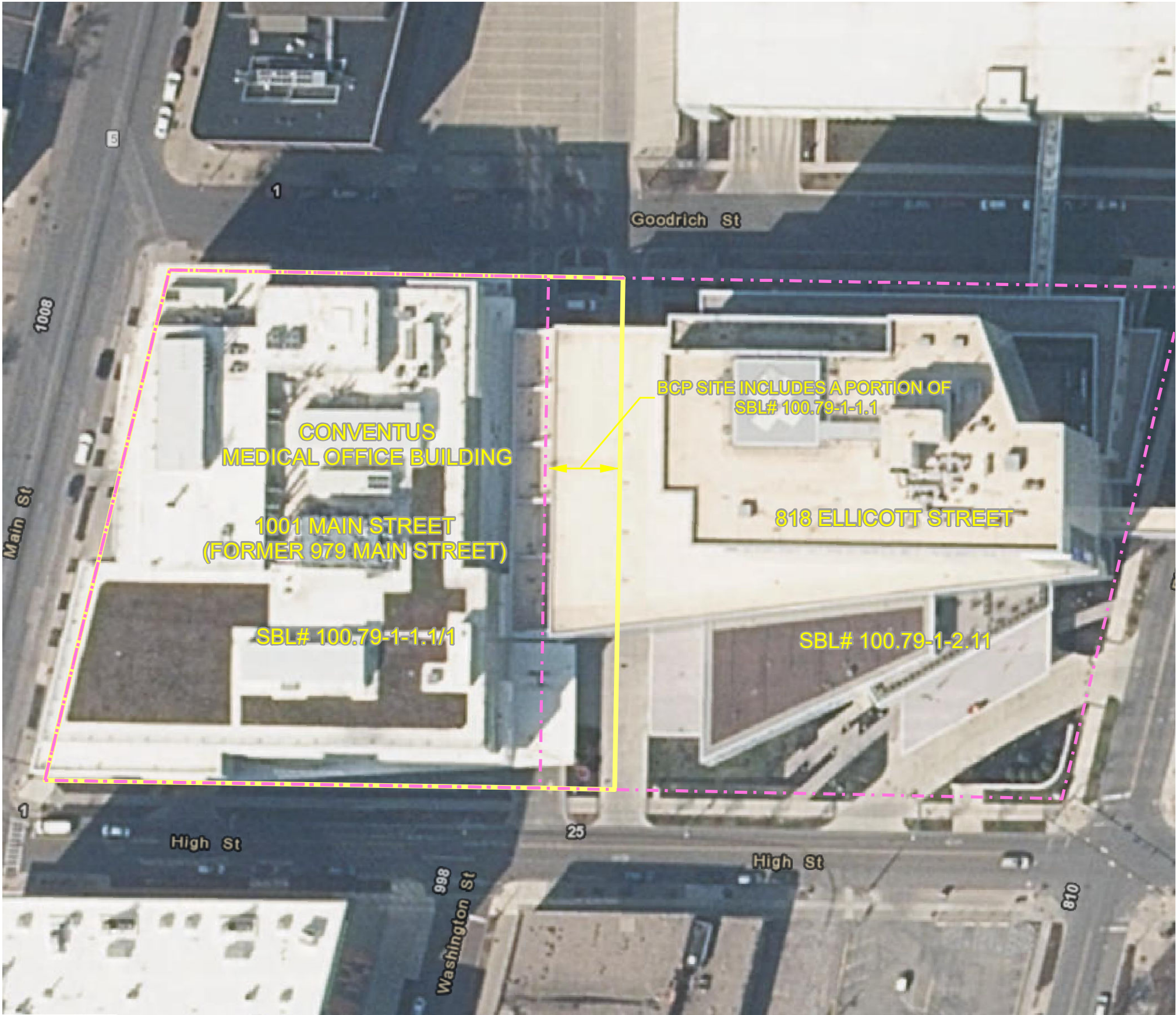
BUFFALO, NEW YORK

MARK	DATE	DESCRIPTION
REVISIONS		
PROJECT NO: U86.001.005		
DATE: APRIL 2023		
DRAWN BY: C. MARTIN		
DESIGNED BY: C. MARTIN		
CHECKED BY: D. RIKER		
NO ALTERATION PERMITTED HEREON EXCEPT AS PROVIDED UNDER SECTION 7209 SUBDIVISION 2 OF THE NEW YORK EDUCATION LAW		

SITE LOCATION

FIGURE 1

Path: F:\Project\U06 - Seavest Core Buffalo Conventus, LLC\PRR\2023\figure2_project boundaries.mxd



Legend

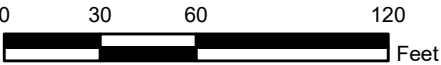
- Parcel Boundary
- Brownfield Cleanup Program Boundary

Property Note

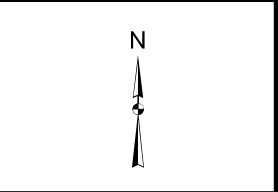
1) The BCP Project Area ("Site") includes the entire western parcel [1001 Main Street (formerly 979 Main Street)] and extends approximately 40 feet east onto the adjacent eastern parcel (818 Ellicott Street). Total acreage of the BCP Project Site is 1.72 acres.

Notes

- Groundwater elevation benchmark.
- Coordinate System: NAD 1983 StatePlane NY West FIPS 3103
Projection: Transverse Mercator
Datum: North American 1983
Units: Foot US



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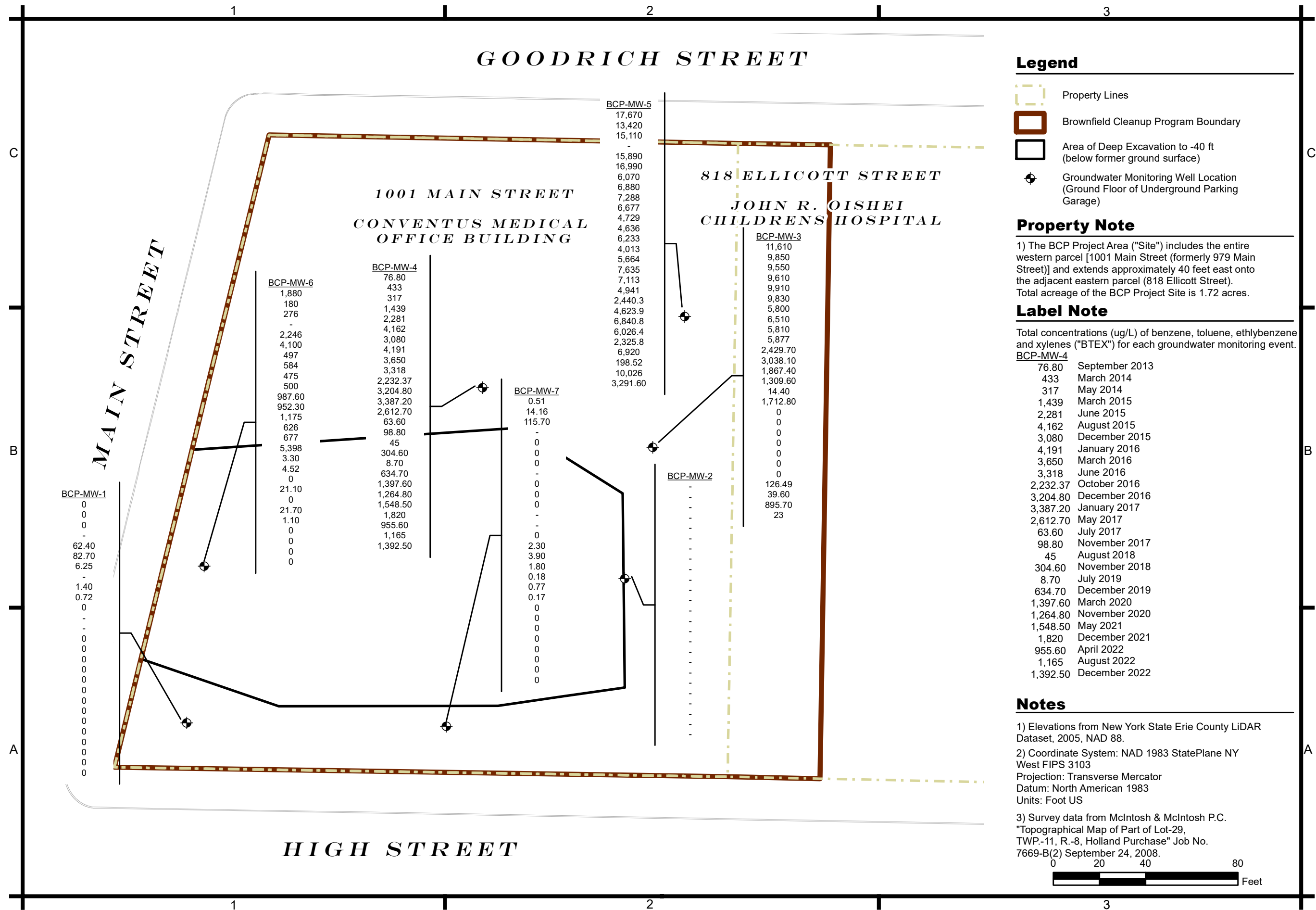
**FORMER MOBIL STATION 99-MST
979 MAIN ST (1001 MAIN ST)**

BUFFALO, NEW YORK

MARK	DATE	DESCRIPTION
REVISIONS		
PROJECT NO:	U06.001.005	
DATE:	APRIL 2023	
DRAWN BY:	C. MARTIN	
DESIGNED BY:	C. MARTIN	
CHECKED BY:		
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**PROJECT
BOUNDARIES**

FIGURE 2



**FORMER MOBIL STATION 99-MST
979 MAIN ST (1001 MAIN ST)**

MARK	DATE	DESCRIPTION
8661800065		
PROJECT NO: APRIL 2023		
DATE:		R. BACKERT
DRAWN BY:		C. MARTIN
DESIGNED BY:		D. RIKER
CHECKED BY:		
NO ALTERATION PERMITTED HEREON EXCEPT AS PROVIDED UNDER SECTION 7209 SUBDIVISION 2 OF THE NEW YORK EDUCATION LAW		

HISTORIC BTEX CONCENTRATIONS

FIGURE 3

TABLES

Table 1A Groundwater Analytical Results
Summary of Detected Compounds
Former Mobil Station 99-MST 979 Main Street (1001 Main Street) Brownfield Cleanup



NYSDEC Ambient Water Quality Standards & Guidance Values Volatile Organic Compound	Sample Name	MW-1	MW-1	MW-1	MW-1	MW-1	MW-1	MW-1	MW-1	MW-1	MW-1	MW-1	MW-1	MW-1	MW-1	MW-1	MW-1	MW-1	MW-1	MW-1	MW-1	MW-1	MW-1		
	Date Collected	9/20/2013	3/19/2014	5/22/2014	3/11/2015	6/17/2015	8/3/2014	12/15/2015	1/27/2016	3/22/2016	6/3/2016	10/25/2016	12/8/2016	1/20/2017	5/17/2017	7/5/2017	11/2/2017	8/16/2018	11/29/2018	7/30/2019	12/12/2019	3/31/2020	11/25/2020	5/14/2021	12/14/2021
	Matrix	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG
	Unit	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Groundwater																									
1,2-DICHLOROBENZENE	3	ND	ND	ND		ND	ND	ND		ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,2-DICHLOROETHANE	0.6	ND	ND	ND		ND	ND	ND		ND	ND	ND		ND	ND	ND	ND	ND	.15 J	ND	ND	ND	ND	ND	ND
1,2-DICHLOROPROPANE	1	ND	ND	ND		ND	ND	ND		ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,3-DICHLOROBENZENE	3	ND	ND	ND		ND	ND	ND		ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
2-HEXANONE	50	ND	ND	ND		ND	ND	3.5		ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
ACETONE	50	ND	ND	ND		ND	ND	ND		ND	ND	ND		ND	5.1	ND	ND	1.8J	2.4 J	1.7	ND	ND	1.8	J	ND
BENZENE	1	ND	ND	ND		35	39	5.7		1.4	0.72	ND		ND	ND	0.33	ND	ND	ND	ND	ND	ND	ND	ND	ND
BROMODICHLOROMETHANE	5	ND	ND	ND			ND			ND	ND	ND			ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
DIBROMOCHLOROMETHANE	50	ND	ND	ND			ND	ND		ND	ND	ND			ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
DICHLORODIFLUOROMETHANE	5	ND	ND	ND			ND	ND		ND	ND	ND			ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
ETHYLBENZENE	5	ND	ND	ND		2	1.5	ND		ND	ND	ND			ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
ISOPROPYLBENZENE (CUMENE)	5	ND	ND	ND		1.3	ND	ND		ND	ND	ND			ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
METHYL ETHYL KETONE (2-BUTANONE)	50	ND	ND	ND		ND	45	ND		ND	ND	ND			ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
METHYLENE CHLORIDE	5	ND	ND	ND		ND	ND	ND		ND	ND	ND			ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
TOLUENE	5	ND	ND	ND		19	38	0.55		ND	ND	ND			ND	ND	1.1	ND	ND	ND	ND	ND	ND	ND	ND
TRICHLOROETHYLENE (TCE)	5	ND	ND	ND		ND	ND	ND		ND	ND	ND			ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,1,2-TRICHLOROETHANE	1	ND	ND	ND		ND	ND	ND		0.33 J	ND	ND			ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
XYLENES, TOTAL	5	ND	ND	ND		6.4	4.2	ND		ND	ND	ND			ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
NAPHTHALENE	10	ND	ND	ND		ND	ND	ND		0.33 J	ND	ND			ND	ND	ND	ND	4.3	ND	ND	ND	ND	1.4	J
No Standard																									
CARBON DISULFIDE		ND	ND	0.94		ND	ND	ND		ND	ND	ND			ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
CYCLOHEXANE		ND	ND	ND		35	59	61		51	72	ND			ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
METHYL ISOBUTYL KETONE		ND	ND	ND		ND	13	ND		ND	ND	ND			ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
4-METHYL-2-PENTANONE																									
METHYLCYCLOHEXANE		ND	ND	0.47		3.2	17	15		11	ND	ND			ND	ND	ND	1.5	.88J	ND	ND	ND	ND	ND	ND
Total VOCs	0	0	1.41	0	101.90	216.70	85.75	0	63.40	72.72	0			0	5.10	1.40	1.50	6.98	2.55	1.70	0	0	3.20	0	
Total BTEX	0	0	0	-	62.40	82.70	6.25	0	1.40	0.72	0			0	0	0	0	0	0	0	0	0	0	0	
Non-Standard VOC List																									
1,3,5-TRIMETHYLBENZENE	5														ND	ND		ND		ND		ND	ND	ND	ND
1,2,4,5-TETRAMETHYLBENZENE	5														ND	ND		ND		ND		ND	ND	ND	ND
1,2,4-TRIMETHYLBENZENE	5														ND	ND		ND		ND		1.6	J	ND	ND
SEC-BUTYLBENZENE	5														ND	ND		ND		ND		ND	ND	ND	ND
N-PROPYLBENZENE	5														ND	ND		ND		ND		ND	ND	ND	ND
N-BUTYLBENZENE	5														ND	ND		ND		ND		ND	ND	ND	ND
P-ISOPROPYLTOLUENE															ND	ND		ND		ND		ND	ND	ND	ND
1,4-DIETHYLBENZENE															ND	ND		ND		ND		ND	ND	ND	ND

Notes:

Not Sampled

1) Blank space = analyte concentration not reported

2) BCP MW-2 was dry and not sampled

3) For the March 11, 2015 monitoring event well MW-1, MW-5, MW-6 and MW-7 were dry or not enough water was inside the well for a representative sample.

4) WG = groundwater

Table 1A Groundwater Analytical Results
Summary of Detected Compounds
Former Mobil Station 99-MST 979 Main Street (1001 Main Street) Brownfield Cleanup



NYSDEC Ambient Water Quality Standards & Guidance Values Volatile Organic Compound	Sample Name	MW-3	MW-3	MW-3	MW-3	MW-3	MW-3	MW-3	MW-3	MW-3	MW-3	MW-3	MW-3	MW-3	MW-3	MW-3	MW-3	MW-3	MW-3	MW-3	MW-3	MW-3	MW-3	MW-3	
	Date Collected	9/20/2013	3/19/2014	5/22/2014	3/11/2015	6/17/2015	8/3/2015	12/15/2015	1/27/2015	3/22/2016	6/3/2016	10/25/2016	12/8/2016	1/20/2017	5/17/2017	7/5/2017	11/2/2017	8/16/2018	11/29/2018	7/30/2019	12/12/2019	3/31/2020	11/25/2020	5/14/2021	12/14/2021
	Matrix	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG
	Unit	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
	Groundwater																								
1,2-DICHLOROBENZENE	3	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,2-DICHLOROETHANE	0.6	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,2-DICHLOROPROPANE	1	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,3-DICHLOROBENZENE	3	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
2-HEXANONE	50	ND	ND	ND	3	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	8	ND	ND	ND	ND	ND	ND	ND	ND	ND
ACETONE	50	ND	98	ND	17	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	166	ND	2.3	24.0	2.1 J	ND	ND	ND	3.8 J	6.2 J
BENZENE	1	6,600	4,500	4,700	3,700	4,300	4,100	2,100	2,200	1,900	3,100	1,390	635	363	451	3	364	ND	ND	ND	0.2J	ND	ND	ND	0.49 J
BROMODICHLOROMETHANE	5	ND	ND	ND														ND	ND	ND	ND	ND	ND	0.2 J	ND
DIBROMOCHLOROMETHANE	50	ND	ND	ND		ND		ND		ND		ND		ND		ND		ND	ND	ND	ND	ND	ND	ND	ND
DICHLORODIFLUOROMETHANE	5	ND	ND	ND		ND		ND		ND		ND		ND		ND		ND	ND	ND	ND	ND	ND	ND	ND
ETHYLBENZENE	5	1,200	1,600	1,500	1,600	1,500	1,700	1,400	1,600	1,600	610	194	899	517	197	2.4	384	ND	ND	ND	1.1 J	ND	ND	ND	42
ISOPROPYLBENZENE (CUMENE)	5	ND	37	ND	32	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		8.7	ND	ND	ND	ND	ND	ND	ND	ND
METHYL ETHYL KETONE (2-BUTANONE)	50	ND	71	ND	6.7	ND	ND	ND	ND	ND	ND	ND	ND	ND	201	51.4	51.4	ND	ND	ND	ND	ND	ND	ND	ND
METHYLENE CHLORIDE	5	ND	ND	ND	ND	ND	ND	ND	ND	ND	35	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
TOLUENE	5	110	150	150	110	110	130	100	110	110	67	39.4	74.5	38.4	22.6	1.6	34.8	ND	ND	ND	ND	ND	ND	ND	4.8
TRICHLOROETHYLENE (TCE)	5	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND
1,1,2-TRICHLOROETHANE	1	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
XYLENES, TOTAL	5	3,700	3,600	3,200	4200	4000	3900	2200	2600	2200	2100	806.3	1430	949	639	7.1	930.0	ND	ND	ND	1.3 J	ND	ND	ND	79.2 J
NAPHTHALENE	10	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	14	357	ND	ND	ND	ND	1.5 J	ND	18 J	
No Standard																									
CARBON DISULFIDE		ND	ND	ND	0.31	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
CYCLOHEXANE		120	320	270	390	330	210	100	93	110	170	ND	ND	ND	ND	ND	60.5	ND	ND	ND	ND	ND	3.4 J	1 J	16 J
METHYL ISOBUTYL KETONE		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
4-METHYL-2-PENTANONE																									ND
METHYLCYCLOHEXANE		ND	130	150	120	160	96	34	33	36 J	170	47.7	ND	ND	29.5	ND	33.4	ND	ND	ND	ND	ND	ND	ND	8.2 J
Total VOCs		11730.00	10506.00	9970.00	10179.01	10400.00	10136.00	5934.00	6636.00	5920.00	6252.00	2477.40	3038.10	1867.40	1540.10	253.80	2223.80	2.30	24.00	2.10	2.60	0	4.90	5	174.89
Total BTEX		11610.00	9850.00	9550.00	9610.00	9910.00	9830.00	5800.00	6510.00	5810.00	5877.00	2429.70	3038.10	1867.40	1309.60	14.40	1712.80	0	0	0	0	0	0	0	126.49
Non-Standard VOC List																									
1,3,5-TRIMETHYLBENZENE	5															ND	133	133	ND	ND	ND	ND	ND	ND	6.5
1,2,4,5-TETRAMETHYLBENZENE	5															ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,2,4-TRIMETHYLBENZENE	5															4.9	737	737	ND	ND	1.2 J	0.88J	ND	ND	68
SEC-BUTYLBENZENE	5															ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
N-PROPYLBENZENE	5															ND	ND	ND	ND	ND	ND	ND	ND	ND	2.3 J
N-BUTYLBENZENE	5															ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
P-ISOPROPYLTOLUENE																ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,4-DIETHYLBENZENE																ND	ND	ND	ND	ND	ND	ND	ND	ND	ND

Notes:

Not Sampled

1) Blank space = analyte concentration not reported
2) BCP MW-2 was dry and not sampled
3) For the March 11, 2015 monitoring event well MW- were dry or not enough water was inside the well for a
4) WG = groundwater

NYSDEC Ambient Water Quality Standards & Guidance Values Volatile Organic Compound	Sample Name	MW-4	MW-4	MW-4	MW-4	MW-4	MW-4	MW-4	MW-4	MW-4	MW-4	MW-4	MW-4	MW-4	MW-4	MW-4	MW-4	MW-4	MW-4	MW-4	MW-4	MW-4	MW-4	MW-4			
	Date Collected	9/20/2013	3/19/2014	5/22/2014	3/11/2015	6/17/2015	8/3/2015	12/15/2015	1/27/2016	3/22/2016	6/3/2016	10/25/2016	12/8/2016	1/20/2017	5/17/2017	7/5/2017	11/17/2017	8/16/2018	11/29/2018	7/30/2019	12/12/2019	3/31/2020	11/25/2020	5/14/2021	12/14/2021		
	Matrix	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG		
	Unit	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L		
	Groundwater																										
1,2-DICHLOROBENZENE	3	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
1,2-DICHLOROETHANE	0.6	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
1,2-DICHLOROPROPANE	1	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	1.0 J	ND	ND	ND	ND		
1,3-DICHLOROBENZENE	3	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
2-HEXANONE	50	ND	ND	ND	1.7	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
ACETONE	50	10	250	170	67	ND	210	ND	ND	ND	ND	ND	ND	ND	38.2	10	1.6	ND	ND	ND	ND	ND	ND	ND	ND		
BENZENE	1	42	29	15	26	24	242	ND	21	ND	21	9.57	12.8	10.2	10.8	1.3	97.0	45.0	36.0	6.7	6.4	7.6	7.8	8.5	5		
BROMODICHLOROMETHANE	5	ND	ND	ND																	ND	ND	ND	ND	ND		
DIBROMOCHLOROMETHANE	50	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
DICHLORODIFLUOROMETHANE	5	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
ETHYLBENZENE	5	4.7	34	32	560	1,000	680	1,100	1300	1,400	1400	1,000	1170	1,300	1220	28	1.8	ND	170	2.0 J	460	810	870	1100	1100		
ISOPROPYLBENZENE (CUMENE)	5	ND	ND	ND	9.8	15.0	26	ND	ND	ND	ND	19	30.3	28.7	ND	2.3	ND	ND	8.3	1.3 J	19	28	34	27	35		
METHYL ETHYL KETONE (2-BUTANONE)	50	ND	ND	ND	ND	8.50	ND	ND	ND	ND	ND	ND	ND	ND	ND	6.9	ND	ND	ND	ND	ND	ND	ND	ND	ND		
METHYLENE CHLORIDE	5	ND	ND	1 J	ND	ND	ND	ND	52	ND	42	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
TOLUENE	5	1.1	190	110	53	57	140	180	270	150	97	62.4	130	133	92.2	9.8	ND	ND	15	ND	11	46	29	22	J	64	
TRICHLOROETHYLENE (TCE)	5	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
1,1,2-TRICHLOROETHANE	1	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
XYLENES, TOTAL	5	29	180	160	800	1,200	3100	1,800	2600	2,100	1800	1,160	1892	1,944	1289.7	24.5	ND	ND	83.6	ND	157.3	534 J	358 J	418	J	651	J
NAPHTHALENE	10	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	1.9	ND	ND	36	ND	99	230	230	320	360	J	
No Standard																											
CARBON DISULFIDE		ND	ND	1.9 J	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
CYCLOHEXANE		8.2	11	7	170	170	110	160	220	250	340	189	259	276	235	276	5.5	ND	24	.41 J	60	100	140	160	190	J	
METHYL ISOBUTYL KETONE		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
4-METHYL-2-PENTANONE																											
METHYLCYCLOHEXANE		7.5	3.7	3.1	87	92	69	86	100	110	140	85.1	110	123	99.7	123	2.4	0.47	8.9	ND	8	22J	38 J	44	J	54	J
Total VOCs		102.50	697.70	497.10	1774.50	2566.50	4577.00	3326.00	4563.00	4010.00	3840.00	2525.47	3604.10	3814.90	2947.40	511.90	116.70	47.07	381.80	10.41	821.70	1777.60	1706.80	2099.50	2459.00		
Total BTEX		76.80	433.00	317.00	1439.00	2281.00	4162.00	3080.00	4191.00	3650.00	3318.00	2232.37	3204.80	3387.20	2612.70	63.60	98.80	45.00	304.60	8.70	634.70	1397.60	1264.80	1548.50	1820.00		
Non-Standard VOC List																											
1,3,5-TRIMETHYLBENZENE	5														2	ND	ND	1.4 J	ND	ND	7.0J	11 J	8.4	J	13	J	
1,2,4,5-TETRAMETHYLBENZENE	5														1.1	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
1,2,4-TRIMETHYLBENZENE	5														1.1	ND	ND	150	ND	470	1100	1300	1500	1500			
SEC-BUTYLBENZENE	5														ND	ND	ND	1.5 J	ND	2.9 J	ND	ND	ND	ND	ND	ND	
N-PROPYLBENZENE	5														2.3	ND	ND	37	ND	86	150	170	160	200			
N-BUTYLBENZENE	5														1.7	ND	ND	2.2 J	ND	4.1 J	10J	12 J	9.7	J	14	J	
P-ISOPROPYLTOLUENE															ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
1,4-DIETHYLBENZENE															ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	

Notes:

Not Sampled

1) Blank space = analyte concentration not reported
2) BCP MW-2 was dry and not sampled
3) For the March 11, 2015 monitoring event well MW- were dry or not enough water was inside the well for a
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Table 1A Groundwater Analytical Results
Summary of Detected Compounds
Former Mobil Station 99-MST 979 Main Street (1001 Main Street) Brownfield Cleanup



NYSDEC Ambient Water Quality Standards & Guidance Values Volatile Organic Compound	Sample Name	MW-5	MW-5	MW-5	MW-5	MW-5	MW-5	MW-5	MW-5	MW-5	MW-5	MW-5	MW-5	MW-5	MW-5	MW-5	MW-5	MW-5	MW-5	MW-5	MW-5	MW-5	MW-5	MW-5	
	Date Collected	9/20/2013	3/19/2014	5/22/2014	3/11/2015	6/17/2015	8/3/2015	12/15/2015	1/27/2016	3/22/2016	6/3/2016	10/25/2016	12/8/2016	1/20/2017	5/17/2017	7/5/2017	11/2/2017	8/16/2018	11/29/2018	7/30/2019	12/12/2019	3/31/2020	11/25/2020	5/14/2021	12/14/2021
	Matrix	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	
	Unit	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
	Groundwater																								
1,2-DICHLOROBENZENE	3	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,2-DICHLOROETHANE	0.6	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,2-DICHLOROPROPANE	1	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,3-DICHLOROBENZENE	3	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
2-HEXANONE	50	11	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	2.7 J	ND	ND	ND	ND	ND	ND
ACETONE	50	ND	520	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	15.3	ND	41	69 J	44	97 J	ND	43 J	45 J	ND
BENZENE	1	5,600	4,800	4,900		3,700	4,100	1,800	1,800	1,700	1,600	899	949	682	428	574	283	86	26	3.3	8.9 J	5.8J	3.4 J	5.8 J	ND
BROMODICHLOROMETHANE	5	ND	ND	ND																	ND	ND	ND	ND	ND
DIBROMOCHLOROMETHANE	50	ND	ND	ND		ND		ND		ND		ND		ND		ND		ND	ND	ND	ND	ND	ND	ND	ND
DICHLORODIFLUOROMETHANE	5	ND	ND	ND		ND		ND		ND		ND		ND		ND		ND	ND	ND	ND	ND	ND	ND	ND
ETHYLBENZENE	5	1,900	1,600	1,600		2,800	2,600	1,600	1,900	2,200	2,200	1,490	1,450	2,070	584	534	1,660	1,500	810	520 E	1200	1,700	1,700	770	1800
ISOPROPYLBENZENE (CUMENE)	5	28	29	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	13.6	ND	20	16 J	23	24 J	30J	33 J	13 J	33 J
METHYL ETHYL KETONE (2-BUTANONE)	50	10	350	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	5.1	ND	ND		ND	ND	ND	ND	ND	ND
METHYLENE CHLORIDE	5	ND	ND	ND		ND	ND	ND	ND	77	96	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND
TOLUENE	5	170	220	310		290	290	70	80	88	77	68.5	84.9	86.6	ND	36.2	82.0	66.0	39 J	38.0	42 J	49J	48 J	16.0 J	44 J
TRICHLOROETHYLENE (TCE)	5	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		.22 J	ND	ND	ND	ND	ND
1,1,2-TRICHLOROETHANE	1	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND
XYLENES, TOTAL	5	10,000	6,800	8,300		9,100	10,000	2,600	3,100	3,300	2,800	2,271.3	2,152.2	3,394.7	3,000.7	4,520.0	5,610.0	5,461.0	4,066.0	1879 E	3373	5,086.0	4,275	1,534.0	5,076
NAPHTHALENE	10	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	730	1,030	620	1,100		1100	940	820	430	1000 J
No Standard																									
CARBON DISULFIDE		ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	2.1	ND	ND		1.2 J	ND	ND	ND	ND	ND
CYCLOHEXANE		230	340	240		430	260	230	250	280	430	198	148	257	ND	257	238	150	130 J	140	220	250	240	130	330 J
METHYL ISOBUTYL KETONE		23	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		3.0 J	ND	ND	ND	ND	ND
4-METHYL-2-PENTANONE																									
METHYLCYCLOHEXANE		100	170	150		190	130	92	100	100	140	67.5	58.4	92.8	49	92.8	106	70	82 J	65	96	110J	110 J	60 J	120 J
Total VOCs		18072.00	14829.00	15500.00	0	16510.00	17380.00	6392.00	7230.00	7745.00	7343.00	4994.30	4842.50	6583.10	4061.70	6780.10	9009.00	8014.00	6338.00	2718.72	6160.90	8170.80	7.272.4	3003.80	8403.00
Total BTEX		17670.00	13420.00	15110.00	0	15890.00	16990.00	6070.00	6880.00	7288.00	6677.00	4728.80	4636.10	6233.30	4012.70	5664.20	7635.00	7113.00	4941.00	2440.30	4623.90	6840.80	6026.40	2325.80	6920.00
Non-Standard VOC List																									
1,3,5-TRIMETHYLBENZENE	5															823	ND	ND	630	ND	480	520	400	99	430
1,2,4,5-TETRAMETHYLBENZENE	5															135	ND	ND		ND	ND	ND	ND	ND	ND
1,2,4-TRIMETHYLBENZENE	5															2,280	2,490	2,400	2,300	ND	2200	2500	2500	1200	2500
SEC-BUTYLBENZENE	5															3.2	ND	ND		ND	ND	ND	ND	ND	ND
N-PROPYLBENZENE	5															34.8	ND	110	69	ND	110	140	150	64	170
N-BUTYLBENZENE	5															43.3	ND	ND		ND	4.1 J	ND	ND	ND	ND
P-ISOPROPYLTOLUENE																5.7	ND	ND		ND	ND	ND	ND	ND	ND
1,4-DIETHYLBENZENE																347	ND	ND		ND	ND	ND	ND	ND	ND

Notes:

Not Sampled

1) Blank space = analyte concentration not reported
2) BCP MW-2 was dry and not sampled
3) For the March 11, 2015 monitoring event well MW- were dry or not enough water was inside the well for a
4) WG = groundwater

Table 1A Groundwater Analytical Results
Summary of Detected Compounds
Former Mobil Station 99-MST 979 Main Street (1001 Main Street) Brownfield Cleanup



NYSDEC Ambient Water Quality Standards & Guidance Values Volatile Organic Compound	Sample Name	MW-6	MW-6	MW-6	MW-6	MW-6	MW-6	MW-6	MW-6	MW-6	MW-6	MW-6	MW-6	MW-6	MW-6	MW-6	MW-6	MW-6	MW-6	MW-6	MW-6	MW-6	MW-6	MW-6	
	Date Collected	9/20/2013	3/19/2014	5/22/2014	3/11/2015	6/17/2015	8/3/2015	12/14/2015	1/27/2016	3/22/2016	6/3/2016	10/25/2016	12/8/2016	1/20/2017	5/17/2017	7/5/2017	11/2/2017	8/16/2018	11/29/2018	7/30/2019	12/12/2019	3/31/2020	11/25/2020	5/14/2021	12/14/2021
	Matrix	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG
	Unit	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
	Groundwater																								
1,2-DICHLOROBENZENE	3	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,2-DICHLOROETHANE	0.6	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,2-DICHLOROPROPANE	1	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	.20 J	ND	ND	ND	ND
1,3-DICHLOROBENZENE	3	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
2-HEXANONE	50	ND	ND	ND		190	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
ACETONE	50	ND	ND	ND		480	340	ND	ND	ND	ND	ND	ND	ND	ND	102	ND	17	4.5 J	ND	6.4	1.6J	ND	ND	ND
BENZENE	1	190	33	16		470	890	250	230	200	120	302	168	200	113	131	774	ND	0.82	ND	4	ND	7.5	ND	ND
BROMODICHLOROMETHANE	5	ND	ND	ND															ND	ND	ND	ND	ND	ND	ND
DIBROMOCHLOROMETHANE	50	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
DICHLORODIFLUOROMETHANE	5	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
ETHYLBENZENE	5	130	20	31		36	210	22	44	67	50	163	169	173	175	85.5	154.0	3.3	1.7 J	ND	2.4 J	ND	2.7	ND	ND
ISOPROPYLBENZENE (CUMENE)	5	4.4	ND	1.9 J		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	2.5	ND	1.3	ND	ND	.90 J	ND	ND	ND	ND
METHYL ETHYL KETONE (2-BUTANONE)	50	ND	ND	ND		110	ND	ND	ND	ND	ND	ND	ND	ND	ND	19.6	ND	ND	ND	ND	ND	ND	ND	ND	ND
METHYLENE CHLORIDE	5	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
TOLUENE	5	810	42	79		1,000	1,900	85	120	78	120	130	255	351	147	22.5	2,970.0	ND	ND	ND	6.7	ND	9	ND	ND
TRICHLOROETHYLENE (TCE)	5	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,1,2-TRICHLOROETHANE	1	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
XYLENES, TOTAL	5	750	85	150		740	1,100	140	190	130	210	393	360	451	190.7	438	1,500	ND	2 J	ND	8	ND	10	1.1 J	ND
NAPHTHALENE	10	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	86.6	ND	1	.8 J	ND	4.8	ND	2.6	3.9	ND
No Standard																									
CARBON DISULFIDE		ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
CYCLOHEXANE		68	ND	130		270	41	62	110	110	91	81.5	ND	ND	ND	ND	84	7.4	3.7 J	.60 J	6.6 J	ND	7.2 J	ND	ND
METHYL ISOBUTYL KETONE		ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
4-METHYL-2-PENTANONE																									
METHYLCYCLOHEXANE		46	16	18		170	27	24	21	10	24	32.2	30.2	36.9	35.3	36.9	44	4.3	3.8 J	ND	4.5 J	ND	4.6 J	ND	ND
Total VOCs		1998.40	196.00	424.00	0	3466.00	4508.00	583.00	715.00	595.00	615.00	1101.30	982.50	1211.90	661.00	924.60	5526.00	34.50	17.32	0.60	44.50	1.60	43.60	5.00	0
Total BTEX		1880.00	180.00	276.00	0	2246.00	4100.00	497.00	584.00	475.00	500.00	987.60	952.30	1175.00	625.70	677.00	5398.00	3.30	4.52	0.00	21.10	0	21.70	1.10	0
Non-Standard VOC List																									
1,3,5-TRIMETHYLBENZENE	5															74.3	ND	ND	5.1	ND	1.4 J	ND	2.0 J	ND	ND
1,2,4,5-TETRAMETHYLBENZENE	5															14.3	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,2,4-TRIMETHYLBENZENE	5															134	ND	ND	ND	ND	2.2 J	ND	2.8	2 J	ND
SEC-BUTYLBENZENE	5																		ND	ND	0.88 J	ND	ND	ND	ND
N-PROPYLBENZENE	5															11.3	ND	4.7	1.7 J	ND	1.3 J	ND	1.2 J	ND	ND
N-BUTYLBENZENE	5															4.6	ND	0.72	ND	ND	4.1 J	ND	ND	ND	ND
P-ISOPROPYLTOLUENE																1.6	1.6	1.6	ND	ND	ND	ND	ND	ND	ND
1,4-DIETHYLBENZENE																32.9	32.9	32.9	ND	ND	ND	ND	ND	ND	ND

Notes:

Not Sampled

1) Blank space = analyte concentration not reported
2) BCP MW-2 was dry and not sampled
3) For the March 11, 2015 monitoring event well MW- were dry or not enough water was inside the well for a
4) WG = groundwater

Table 1A Groundwater Analytical Results
Summary of Detected Compounds
Former Mobil Station 99-MST 979 Main Street (1001 Main Street) Brownfield Cleanup



NYSDEC Ambient Water Quality Standards & Guidance Values	Sample Name	MW-7	MW-7	MW-7	MW-7	MW-7	MW-7	MW-7	MW-7	MW-7	MW-7	MW-7	MW-7	MW-7	MW-7	MW-7	MW-7	MW-7	MW-7	MW-7	MW-7	MW-7	MW-7	MW-7	
	Date Collected	9/20/2013	3/19/2014	5/22/2014	3/11/2015	6/17/2015	8/3/2015	12/15/2015	1/27/2016	3/22/2016	6/3/2016	10/25/2016	12/8/2016	1/20/2017	5/17/2017	7/5/2017	11/2/2017	8/16/2018	11/29/2018	7/30/2019	12/12/2019	3/31/2020	11/25/2020	5/14/2021	12/14/2021
	Matrix	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG	WG
	Unit	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
	Volatile Organic Compound	Groundwater																							
1,2-DICHLOROBENZENE	3	ND	ND	ND		ND	ND	ND		ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
1,2-DICHLOROETHANE	0.6	ND	ND	ND		ND	ND	ND		ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
1,2-DICHLOROPROPANE	1	ND	ND	ND		ND	ND	ND		ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
1,3-DICHLOROBENZENE	3	ND	ND	ND		ND	ND	ND		ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
2-HEXANONE	50	ND	ND	4.8		ND	ND	ND		ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
ACETONE	50	ND	3	ND		ND	ND	ND		ND	ND	ND		ND	ND	ND	1.5		4.2 J	ND	ND	ND	ND	ND	
BENZENE	1	0.51	8.8	14		ND	ND	ND		ND	ND	ND		ND	2.3	2.81	1.8	.18 J	.77	.17 J	ND	ND	ND	ND	
BROMODICHLOROMETHANE	5	ND	ND	ND		ND	ND	ND		ND	ND	ND		ND		ND				ND	ND	ND	ND	ND	
DIBROMOCHLOROMETHANE	50	ND	ND	ND		ND	ND	ND		ND	ND	ND		ND		ND			ND	ND	ND	ND	ND	ND	
DICHLORODIFLUOROMETHANE	5	ND	ND	ND		ND	ND	ND		ND	ND	ND		ND		ND			ND	ND	ND	ND	ND	ND	
ETHYLBENZENE	5	ND	ND	3		ND	ND	ND		ND	ND	ND		ND	ND	0	ND	ND	ND	ND	ND	ND	ND	ND	
ISOPROPYLBENZENE (CUMENE)	5	ND	ND	ND		ND	ND	ND		ND	ND	ND		ND	ND	0.45	ND	ND	ND	ND	ND	ND	ND	ND	
METHYL ETHYL KETONE (2-BUTANONE)	50	ND	ND	ND		ND	ND	ND		ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
METHYLENE CHLORIDE	5	ND	ND	ND		ND	ND	ND		ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
TOLUENE	5	ND	0.56	4.7		ND	ND	ND		ND	ND	ND		ND	ND	1.1	ND	ND	ND	ND	ND	ND	ND	ND	
TRICHLOROETHYLENE (TCE)	5	ND	ND	ND		ND	ND	ND		ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
1,1,2-TRICHLOROETHANE	1					ND	ND	ND		ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
XYLENES, TOTAL	5	0.96	4.8	94		ND	ND	ND		0.99 J	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
NAPHTHALENE	10																1.50	.86 J	ND	ND	ND	ND	1 J	ND	
No Standard																									
CARBON DISULFIDE		ND	ND	0.97		ND	ND	ND		ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
CYCLOHEXANE		ND	4.3	9.6		ND	ND	0.71		ND	ND	ND		ND	ND	0.99	0.66	ND	ND	ND	ND	ND	ND	ND	
METHYL ISOBUTYL KETONE		ND	ND	ND		ND	ND	ND		ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
4-METHYL-2-PENTANONE																									
METHYLCYCLOHEXANE		ND	1.7	5.1		0.18	ND	ND		ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
Total VOCs		1.47	23.16	136.17	0	0.18	0	0.71	0	0	0	0	0	0	0	2.30	5.35	3.66	1.04	4.97	0.17	0	0	1	0
Total BTEX		0.51	14.16	115.70	0	0	0	0.00	0	0	0	0	0	0	0	2.30	3.91	1.80	0.18	0.77	0.17	0	0	0	0
Non-Standard VOC List																									
1,3,5-TRIMETHYLBENZENE	5															ND	ND	3.2		3.2	ND	ND	ND	ND	ND
1,2,4,5-TETRAMETHYLBENZENE	5															ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,2,4-TRIMETHYLBENZENE	5															ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
SEC-BUTYLBENZENE	5															ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
N-PROPYLBENZENE	5															ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
N-BUTYLBENZENE	5															ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
P-ISOPROPYLTOLUENE																ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,4-DIETHYLBENZENE																ND	ND	ND	ND	ND	ND	ND	ND	ND	ND

Notes:

Not Sampled

1) Blank space = analyte concentration not reported
2) BCP MW-2 was dry and not sampled
3) For the March 11, 2015 monitoring event well MW- were dry or not enough water was inside the well for c
4) WG = groundwater

TABLE 1B

Groundwater Analytical Results
Summary of Detected Compounds
Former Mobil Station 99-MST 979 Main Street (1001 Main Street) Brownfield Cleanup



SAMPLE ID:		BCP-MW-1	BCP-MW-3	BCP-MW-4	BCP-MW-5	BCP-MW-6	BCP-MW-7	BCPMW-1	BCPMW-3	BCPMW-4	BCPMW-5	BCPMW-6	BCPMW-7	DUP	TRIP BLANK
COLLECTION DATE:		4/7/2022	4/7/2022	4/8/2022	4/7/2022	4/7/2022	4/7/2022	8/25/2022	8/25/2022	8/25/2022	8/25/2022	8/25/2022	8/25/2022	8/25/2022	8/25/2022
SAMPLE MATRIX:		WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER
NY-TOGS-GA		BCP-MW-1													
(ug/l)															
VOCs															
1,1,1-Trichloroethane	5	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,1,2,2-Tetrachloroethane	5	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,1,2-Trichloroethane	1	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,1-Dichloroethane	5	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,1-Dichloroethene	5	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,2,4-Trichlorobenzene	5	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,2,4-Trimethylbenzene	5	ND	15	1200	E 130	ND	ND	ND	560	1200	3100	ND	ND	ND	ND
1,2-Dibromo-3-chloropropane	0.04	ND	ND	ND	ND	ND	ND	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ
1,2-Dibromoethane	0.0006	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,2-Dichlorobenzene	3	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,2-Dichloroethane	0.6	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,2-Dichloropropane	1	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,3,5-Trimethylbenzene	5	ND	0.77 J	21	58	ND	ND	ND	47	ND	630	ND	ND	ND	ND
1,3-Dichlorobenzene	3	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,4-Dichlorobenzene	3	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
2-Butanone	50	ND	ND	ND	ND	ND	ND	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ
2-Hexanone	50	ND	ND	ND	ND	ND	ND	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ
4-Methyl-2-pentanone	NA	ND	2.2 J	ND	ND	ND	ND	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ
Acetone	50	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Benzene	1	ND	0.5	2.6	0.22 J	ND	ND	ND	1.7	5	ND	ND	ND	ND	ND
Bromodichloromethane	50	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Bromoform	50	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Bromomethane	5	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Carbon disulfide	60	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Carbon tetrachloride	5	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Chlorobenzene	5	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Chloroethane	5	ND	ND	ND	ND	ND	ND	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ
Chloroform	7	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Chloromethane	NA	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
cis-1,2-Dichloroethene	5	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
cis-1,3-Dichloropropene	0.4	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Cyclohexane	NA	ND	13	120	18	ND	ND	ND	40	130	240	ND	ND	ND	ND
Dibromochloromethane	50	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Dichlorodifluoromethane	5	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Ethylbenzene	5	ND	15	790	39	ND	ND	ND	380	1100	2100	ND	ND	ND	ND
Freon-113	5	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Isopropylbenzene	5	ND	ND	31	1.4 J	ND	ND	ND	5.2	24	29	ND	ND	ND	ND
Methyl Acetate	NA	ND	ND	ND	ND	ND	ND	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ
Methyl cyclohexane	NA	ND	6.2 J	43 J	11	ND	ND	ND	32	51	95	ND	ND	ND	ND
Methyl tert butyl ether	10	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Methylene chloride	5	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
n-Butylbenzene	5	ND	ND	10 J	0.77 J	ND	ND	ND	ND	9.4	ND	ND	ND	ND	ND
n-Propylbenzene	5	ND	1 J	150	3.6	ND	ND	ND	18	140	130	ND	ND	ND	ND
Naphthalene	10	ND	5.7	260	65	ND	ND	ND	170	200	840	ND	ND	ND	ND
o-Xylene	5	ND	1.9 J	18	8.4	ND	ND	ND	12	12	100	ND	ND	ND	ND
p-Isopropyltoluene	5	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
p/m-Xylene	5	ND	21	120	150	ND	ND	ND	480	14	7800	ND	ND	ND	ND

TABLE 1B

Groundwater Analytical Results
Summary of Detected Compounds
Former Mobil Station 99-MST 979 Main Street (1001 Main Street) Brownfield Cleanup



SAMPLE ID:		BCP-MW-1	BCP-MW-3	BCP-MW-4	BCP-MW-5	BCP-MW-6	BCP-MW-7	BCPMW-1	BCPMW-3	BCPMW-4	BCPMW-5	BCPMW-6	BCPMW-7	DUP	TRIP BLANK
COLLECTION DATE:		4/7/2022	4/7/2022	4/8/2022	4/7/2022	4/7/2022	4/7/2022	8/25/2022	8/25/2022	8/25/2022	8/25/2022	8/25/2022	8/25/2022	8/25/2022	8/25/2022
SAMPLE MATRIX:		WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER
NY-TOGS-GA (ug/l)														BCP-MW-1	
VOCs															
sec-Butylbenzene	5	ND	ND	4.6 J	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Styrene	5	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
tert-Butylbenzene	5	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Tetrachloroethene	5	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Toluene	5	ND	1.2 J	25	0.9 J	ND	ND	ND	22	34	26	ND	ND	ND	ND
trans-1,2-Dichloroethene	5	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
trans-1,3-Dichloropropene	0.4	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Trichloroethene	5	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Trichlorofluoromethane	5	ND	ND	ND	ND	ND	ND	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ
Vinyl chloride	2	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
VOC		0	83.47	2795.20	486.29	0	0	0	1767.90	2919.40	15090.00	0	0	0	0
BTEX		0	39.60	955.60	198.52	0	0	0	895.70	1165.00	10026.00	0	0	0	0

NY-TOGS-GA: New York TOGS 111 Groundwater Effluent Limitations criteria reflects all addendum to criteria through June 2004.

ND indicates analyte was not detected.

Blank space indicates analyte was not analyzed for in that sample.

*+ - LCS and/or LCSD is outside acceptance limits, high biased.

B - Compound was found in the blank and sample.

J - Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value.

TABLE 1B

Groundwater Analytical Results
Summary of Detected Compounds
Former Mobil Station 99-MST 979 Main Street (1001 Main Street) Brownfield Cleanup



		SAMPLE ID:	BCPMW-1	BCPMW-3	BCPMW-4	BCPMW-5	BCPMW-6	BCPMW-7	DUP	TRIP BLANK
		COLLECTION DATE:	12/14/2022	12/14/2022	12/14/2022	12/14/2022	12/14/2022	12/14/2022	12/14/2022	12/15/2022
		SAMPLE MATRIX:	WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER
		NY-TOGS-GA							BCP-MW-1	
		(ug/l)								
VOCs										
1,1,1-Trichloroethane	5	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,1,2,2-Tetrachloroethane	5	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,1,2-Trichloroethane	1	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND	ND UJ	ND
1,1-Dichloroethane	5	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,1-Dichloroethene	5	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,2,4-Trichlorobenzene	5	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,2,4-Trimethylbenzene	5	ND	13	1200	1400	ND	ND	ND	ND	ND
1,2-Dibromo-3-chloropropane	0.04	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND	ND UJ	ND
1,2-Dibromoethane	0.0006	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND	ND UJ	ND
1,2-Dichlorobenzene	3	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,2-Dichloroethane	0.6	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND	ND UJ	ND
1,2-Dichloropropane	1	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,3,5-Trimethylbenzene	5	ND	ND	ND	350	ND	ND	ND	ND	ND
1,3-Dichlorobenzene	3	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,4-Dichlorobenzene	3	ND	ND	ND	ND	ND	ND	ND	ND	ND
2-Butanone	50	ND	ND	ND	ND	ND	ND	ND	ND	ND
2-Hexanone	50	ND	ND	ND	ND	ND	ND	ND	ND	ND
4-Methyl-2-pentanone	NA	ND	ND	ND	ND	ND	ND	ND	ND	ND
Acetone	50	ND	ND	ND	ND	ND	ND	ND	ND	ND
Benzene	1	ND	ND	3.5	ND	ND	ND	ND	ND	ND
Bromodichloromethane	50	ND	ND	ND	ND	ND	ND	ND	ND	ND
Bromoform	50	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND	ND UJ	ND
Bromomethane	5	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND	ND UJ	ND UJ
Carbon disulfide	60	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ
Carbon tetrachloride	5	ND	ND	ND	ND	ND	ND	ND	ND	ND
Chlorobenzene	5	ND	ND	ND	ND	ND	ND	ND	ND	ND
Chloroethane	5	ND	ND	ND	ND	ND	ND	ND	ND	ND
Chloroform	7	ND	ND	ND	ND	ND	ND	ND	ND	ND
Chloromethane	NA	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND	ND UJ	ND
cis-1,2-Dichloroethene	5	ND	ND	ND	ND	ND	ND	ND	ND	ND
cis-1,3-Dichloropropene	0.4	ND	ND	ND	ND	ND	ND	ND	ND	ND
Cyclohexane	NA	ND	1.5	140	120	ND	ND	ND	ND	ND
Dibromochloromethane	50	ND	ND	ND	ND	ND	ND	ND	ND	ND
Dichlorodifluoromethane	5	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND	ND UJ	ND
Ethylbenzene	5	ND	11	1100	820	ND	ND	ND	ND	ND
Freon-113	5	ND	ND	ND	ND	ND	ND	ND	ND	ND
Isopropylbenzene	5	ND	ND	30	16	ND	ND	ND	ND	ND
Methyl Acetate	NA	ND	ND	ND	ND	ND	ND	ND	ND	ND
Methyl cyclohexane	NA	ND	0.65	43	54	ND	ND	ND	ND	ND
Methyl tert butyl ether	10	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND UJ	ND	ND UJ	ND
Methylene chloride	5	ND	ND	ND	ND	ND	ND	ND	ND	ND
n-Butylbenzene	5	ND	ND	11	ND	ND	ND	ND	ND	ND
n-Propylbenzene	5	ND	ND	160	71	ND	ND	ND	ND	ND
Naphthalene	10	ND	4.2	290	520	ND	ND	ND	ND	ND
o-Xylene	5	ND	ND	16	62	ND	ND	ND	ND	ND
p-Isopropyltoluene	5	ND	ND	ND	ND	ND	ND	ND	ND	ND
p/m-Xylene	5	ND	12	210	2400	ND	ND	ND	ND	ND

TABLE 1B

Groundwater Analytical Results
Summary of Detected Compounds
Former Mobil Station 99-MST 979 Main Street (1001 Main Street) Brownfield Cleanup



		SAMPLE ID:	BCPMW-1	BCPMW-3	BCPMW-4	BCPMW-5	BCPMW-6	BCPMW-7	DUP	TRIP BLANK
		COLLECTION DATE:	12/14/2022	12/14/2022	12/14/2022	12/14/2022	12/14/2022	12/14/2022	12/14/2022	12/15/2022
		SAMPLE MATRIX:	WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER
		NY-TOGS-GA							BCP-MW-1	
		(ug/l)								
VOCs										
sec-Butylbenzene	5		ND	ND	ND	ND	ND	ND	ND	ND
Styrene	5		ND	ND	ND	ND	ND	ND	ND	ND
tert-Butylbenzene	5		ND	ND	ND	ND	ND	ND	ND	ND
Tetrachloroethene	5		ND	ND	ND	ND	ND	ND	ND	ND
Toluene	5		ND	ND	63	9.6	ND	ND	ND	ND
trans-1,2-Dichloroethene	5		ND	ND	ND	ND	ND	ND	ND	ND
trans-1,3-Dichloropropene	0.4		ND JJ	ND JJ	ND JJ	ND JJ	ND JJ	ND	ND JJ	ND
Trichloroethene	5		ND	ND	ND	ND	ND	ND	ND	ND
Trichlorofluoromethane	5		ND	ND	ND	ND	ND	ND	ND	ND
Vinyl chloride	2		ND	ND	ND	ND	ND	ND	ND	ND
VOC			0	42.35	3266.50	5822.60	0	0	0	0
BTEX			0	23.00	1392.50	3291.60	0	0	0	0

NY-TOGS-GA: New York TOGS 111 Groundwater Effluent Limitat

ND indicates analyte was not detected.

Blank space indicates analyte was not analyzed for in that sampl

*+ - LCS and/or LCSD is outside acceptance limits, high biased.

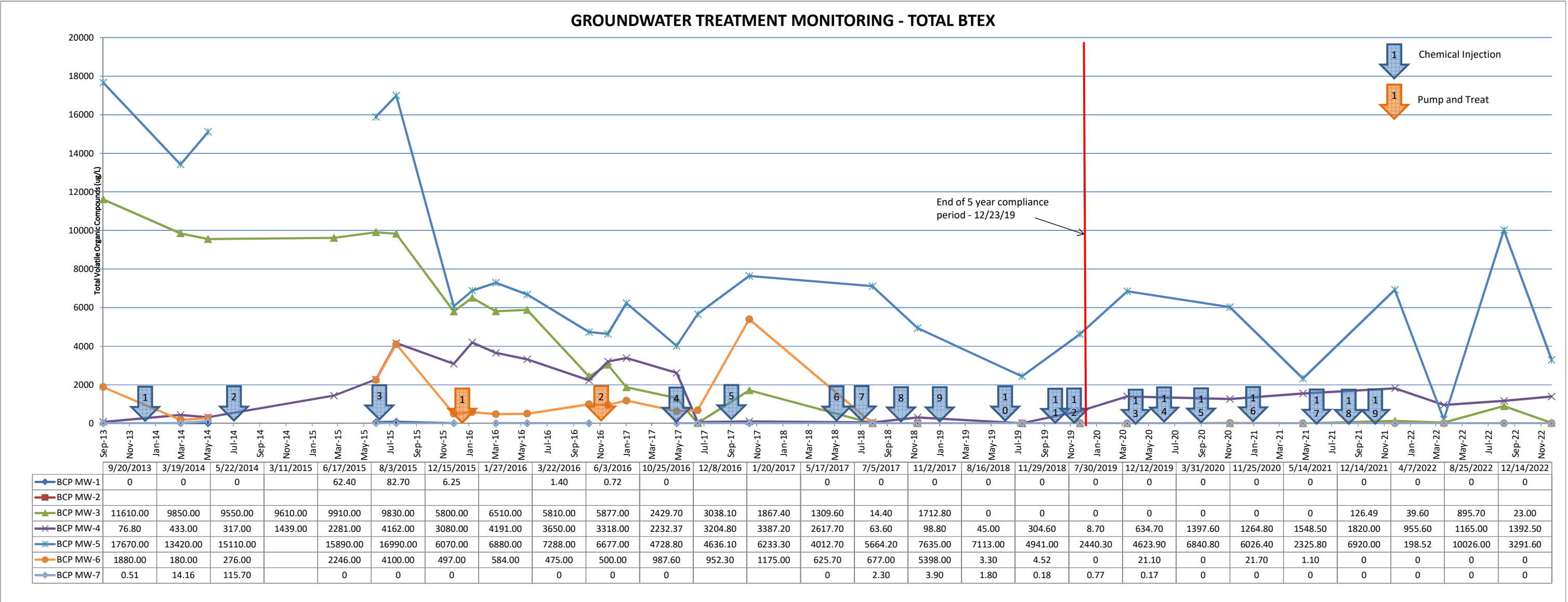
B - Compound was found in the blank and sample.

J - Result is less than the RL but greater than or equal to the MD

GRAPHS

GRAPH 1

Former Mobil Station 99-MST 979 Main Street (1001 Main Street)
Conventus Groundwater Remediation



APPENDICES

APPENDIX C

EXAMPLE (Minimum Requirements)

WELL PURGING-FIELD WATER QUALITY MEASUREMENTS FORM

Location (Site/Facility Name) <u>Conventus Parking Garage</u>		Depth to <u>6.72</u> / <u>14.90</u> of screen (below MP) top bottom	
Well Number <u>BEP-MW-1</u> Date <u>4-7-2022</u> <u>830</u>		Pump Intake at (ft. below MP) _____	
Field Personnel <u>TW & ES</u>		Purging Device; (pump type) <u>600 pump</u>	
Sampling Organization <u>Alpha</u>		Total Volume Purged <u>~ 3/4 gal</u>	
Identify MP <u>Top of Rider</u>			

Clock Time 24 HR	Water Depth below MP ft	Pump Dial ¹	Purge Rate ml/min	Cum. Volume Purged liters	Temp. °C	Spec. Cond. ² µS/cm mS/cm	pH	ORP ³ mv	DO mg/L	Turbidity NTU	Comments
835	6.72	1/4	~100ml		10.7	8.283	6.83	261.4	1.11	24.92	
840	6.73				10.7	8.289	6.94	258.0	1.04	26.73	
845	6.73				10.7	8.291	7.02	254.0	1.00	25.62	
850	6.72				10.7	8.291	7.05	249.3	0.99	23.59	
855	6.72				10.7	8.299	7.07	240.3	0.99	19.85	
900	6.73	↓	↓		10.7	8.308	7.08	232.6	0.97	20.38	
											well Sampled @
											900 on 4-7-2022

Stabilization Criteria

3%

3%

±0.1 ±10 mv

10%

10%

1. Pump dial setting (for example: hertz, cycles/min, etc).

2. µSiemens per cm (same as µmhos/cm) at 25°C.

3. Oxidation reduction potential (ORP)

EXAMPLE (Minimum Requirements)

WELL PURGING-FIELD WATER QUALITY MEASUREMENTS FORM

[illegible]

3%	3%	$\pm 0.1 \pm 10$ mv	10%	10%
3%	3%	$\pm 0.1 \pm 10$ mv	10%	10%

1. Pump dial setting (for example: hertz, cycles/min, etc).
2. μ Siemens per cm (same as μ mhos/cm) at 25°C.
3. Oxidation reduction potential (ORP)

APPENDIX C

EXAMPLE (Minimum Requirements)
WELL PURGING-FIELD WATER QUALITY MEASUREMENTS FORM

Location (Site/Facility Name) <u>Conventus Parking Garage</u>		Depth to <u>6.59</u> / <u>15.49</u> of screen (below MP) top bottom	
Well Number <u>BCP-MW-3</u> Date <u>4-7-2022</u> <u>10:55</u>		Pump Intake at (ft. below MP) _____	
Field Personnel <u>Tw/Es</u>		Purging Device; (pump type) <u>Geo Rmp</u>	
Sampling Organization <u>Alpha</u>		Total Volume Purged <u>~ 1 gal</u>	
Identify MP <u>Top of Riser</u>			

Clock Time 24 HR	Water Depth below MP ft	Pump Dial ¹	Purge Rate ml/min	Cum. Volume Purged liters	Temp. °C	Spec. Cond. ² $\mu\text{S}/\text{cm}$ <u>MS/cm</u>	pH	ORP ³ mv	DO mg/L	Turbidity NTU	Comments
1100	6.61	Y4	~100 mL		13.1	14.549	8.27	151.0	1.70	3.05	
1105	6.64				12.7	14.845	8.37	97.2	1.05	3.48	
1110	6.64				12.7	14.833	8.39	59.4	0.88	3.24	
1115	6.64				12.6	14.839	8.40	35.4	0.81	2.88	
1120	6.65				12.7	14.839	8.41	-4.9	0.75	2.84	
1125	6.65				12.7	14.839	8.42	-14.7	0.73	2.69	
1130	6.65				12.7	14.839	8.42	-19.6	0.73	2.65	
1135	6.65				12.7	14.840	8.42	-23.8	0.71	2.61	-well was sampled
											© 1135 on 4-7-22

Stabilization Criteria

- 3% 3% 3% 3% 3% 3% 3% 3% 3% 3% 3% 3%
1. Pump dial setting (for example: hertz, cycles/min, etc).
 2. $\mu\text{Siemens per cm}$ (same as $\mu\text{mhos/cm}$) at 25°C.
 3. Oxidation reduction potential (ORP)

APPENDIX C

EXAMPLE (Minimum Requirements)
WELL PURGING-FIELD WATER QUALITY MEASUREMENTS FORM

Location (Site/Facility Name) <u>Conventus Parking Garage</u>		Depth to <u>6.58'</u> / <u>14.85'</u> of screen (below MP) top bottom	
Well Number <u>BCP-MW-4</u> Date <u>4-8-2022</u> <u>935</u>		Pump Intake at (ft. below MP)	
Field Personnel <u>W & AF</u>		Purging Device; (pump type) <u>Geo pump</u>	
Sampling Organization <u>Alpha</u>		Total Volume Purged <u>~ 1 gal</u>	
Identify MP <u>Top of Riser</u>			

Clock Time 24 HR	Water Depth below MP ft	Pump Dial ¹	Purge Rate ml/min	Cum. Volume Purged liters	Temp. °C	Spec. Cond. ² $\mu S/cm$	pH	ORP ³ mv	DO mg/L	Turbidity NTU	Comments
940	6.83	1/4	~100 mL		12.5	4.282	8.01	-343.5	0.78	14.80	
945	6.72	↓	↓		12.5	3.843	8.00	-353.0	0.73	7.81	
950	6.72	↓	↓		12.4	3.587	7.92	-354.3	0.70	3.36	
955	6.70	↓	↓		12.3	3.490	7.86	-355.6	0.68	3.58	
1000	6.70	↓	↓		12.3	3.486	7.86	-358.0	0.68	3.30	
											well was sampled
											on 4-8-22

Stabilization Criteria

3% 3% ±0.1 ±10 mv 10% 10%

1. Pump dial setting (for example: hertz, cycles/min, etc).
2. μ Siemens per cm (same as μ mhos/cm) at 25°C.
3. Oxidation reduction potential (ORP)

APPENDIX C

EXAMPLE (Minimum Requirements)
WELL PURGING-FIELD WATER QUALITY MEASUREMENTS FORM

Location (Site/Facility Name) <u>Conventus Parking Garage</u>		Depth to <u>7.42</u> / <u>15.11</u> of screen (below MP) top bottom	
Well Number <u>BCP-MW-5</u> Date <u>4-7-22</u> <u>C 1140</u>		Pump Intake at (ft. below MP)	
Field Personnel <u>TV ES</u>		Purging Device; (pump type) <u>Geo Pump</u>	
Sampling Organization <u>Alpha</u>		Total Volume Purged <u>~1 gal</u>	
Identify MP <u>Top of Riser</u>			

Clock Time 24 HR	Water Depth below MP ft	Pump Dial ¹	Purge Rate ml/min	Cum. Volume Purged liters	Temp. °C	Spec. Cond. ² µS/cm mS/cm	pH	ORP ³ mv	DO mg/L	Turbidity NTU	Comments
1145	7.73	✓	~100 mL		12.0	16.010	10.17	32.4	0.87	9.97	
1150	7.79	✓	✓		11.9	15.409	10.19	21.8	0.81	8.55	
1155	7.80	✓	✓		11.9	14.923	10.19	14.1	0.78	7.88	
1200	7.81	✓	✓		11.9	14.776	10.20	9.0	0.76	7.60	
1205	7.81	✓	✓		11.9	14.580	10.20	5.0	0.75	7.28	
1210	7.81	✓	✓		11.9	14.524	10.20	4.2	0.75	7.31	
											well was sampled
											C 1210 on 4-7-22

Stabilization Criteria

1. Pump dial setting (for example: hertz, cycles/min, etc).
 2. µSiemens per cm (same as µmhos/cm) at 25°C.
 3. Oxidation reduction potential (ORP)

3% 3% ±0.1 ±10 mv 10% 10%

EXAMPLE (Minimum Requirements) WELL PURGING-FIELD WATER QUALITY MEASUREMENTS FORM

[illegible]

Stabilization Criteria

1. Pump dial setting (for example: hertz, cycles/min, etc).
2. μ Siemens per cm (same as μ mhos/cm) at 25°C.
3. Oxidation reduction potential (ORP)

APPENDIX C

EXAMPLE (Minimum Requirements)
WELL PURGING-FIELD WATER QUALITY MEASUREMENTS FORM

Location (Site/Facility Name) <u>Conventus Parking Garage</u>		Depth to <u>9.83</u> / <u>14.68</u> of screen (below MP) top bottom	
Well Number <u>BCP-44-7</u>		Pump Intake at (ft. below MP) _____	
Field Personnel <u>W & E S</u>		Purging Device; (pump type) <u>Geo Pump</u>	
Sampling Organization <u>Alpha</u>		Total Volume Purged <u>~ 1 gal</u>	
Identify MP <u>Top of Riser</u>			

Clock Time 24 HR	Water Depth below MP ft	Pump Dial ¹	Purge Rate ml/min	Cum. Volume Purged liters	Temp. °C	Spec. Cond. ² $\mu\text{S}/\text{cm}$ mS/cm	pH	ORP ³ mv	DO mg/L	Turbidity NTU	Comments
1005	9.83	1/4	~1000 mL		11.9	4.627	7.12	192.0	3.03	13.26	
1010	10.05	↓	↓		11.8	5.227	7.07	202.4	1.15	20.20	
1015	10.07	↓	↓		11.8	5.448	7.09	202.9	1.03	20.30	
1020	10.08	↓	↓		11.8	5.481	7.10	202.8	0.95	15.80	
1025	10.09	↓	↓		11.8	5.483	7.11	202.9	0.94	11.45	
1030	10.10	↓	↓		11.8	5.480	7.11	202.7	0.90	11.30	
1035	10.11	↓	↓		11.8	5.478	7.11	203.4	0.87	11.62	
											Well was Sampled
											at 1035 on 4-7-22

Stabilization Criteria

1. Pump dial setting (for example: hertz, cycles/min, etc).
 2. $\mu\text{Siemens}$ per cm (same as $\mu\text{mhos}/\text{cm}$) at 25°C.
 3. Oxidation reduction potential (ORP)
- 3% 3% $\pm 0.1 \pm 10 \text{ mv}$ 10% 10%



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Well Sampling Field Data Sheet

Well Casing Unit Volume

(gal/l.f.)

1 1/4" = 0.08 2" = 0.17 3" = 0.38
4" = 0.66 6" = 1.5 8" = 2.6

Client Name: _____

Site Name: CONVERSE

Project No.: _____

Field Staff: RICH BARNETT

WELL DATA

Date		8/25/22							
Well Number		BUP MW-1							
Diameter (inches)		2"							
Total Sounded Depth (feet)		15'							
Static Water Level (feet)		7.2'							
H ₂ O Column (feet)		7.8'							
Pump Intake (feet)									
Well Volume (gallons)									
Amount to Evacuate (gallons)		2 gal							
Amount Evacuated (gallons)		2 gal							

FIELD READINGS

Date	Stabilization	8/25/22							
Time	Criteria	9:40	9:45	9:50	9:55				
pH (Std. Units)	+/-0.1	8.01	7.30	7.20	7.19				
Conductivity (mS/cm)	3%	8.50	8.96	9.20	9.24				
Turbidity (NTU)	10%	0.00	0.00	0.00	0.00				
D.O. (mg/L)	10%	2.76	1.13	1.13	1.10				
Temperature (°C) (°F)	3%	18.37	15.78°C	14.89°C	14.78°C				
ORP ³ (mV)	+/-10 mv	107	111	108	43				
Appearance		C	C	C	C				
Free Product (Yes/No)		NONE	NONE	NONE	NONE				
Odor		NONE	NONE	NONE	NONE				
Comments	- DUP COLLECTED FROM BUP MW-1 - RID AT TIME SAMPLE COLLECTED - 0.1 PM								

C = Clear T = Turbid ST = Semi Turbid VT = Very Turbid



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Well Casing Unit Volume

(gal/l.f.)

1 1/4" = 0.08 2" = 0.17 3" = 0.38
4" = 0.66 6" = 1.5 8" = 2.6

Client Name:

Site Name:

Project No.:

Field Staff:

CONVENTUS

RICH BACKERO

WELL DATA

Date	8/28/22								
Well Number	11CPHW7								
Diameter (inches)	2"								
Total Sounded Depth (feet)	15'								
Static Water Level (feet)	10.0'								
H ₂ O Column (feet)	5'								
Pump Intake (feet)									
Well Volume (gallons)									
Amount to Evacuate (gallons)	1.5 gal								
Amount Evacuated (gallons)	1.5 gal								

FIELD READINGS

Date	8/28/22								
Time	10:15	10:20	10:25						
pH (Std. Units)	+/-0.1	7.90	7.18	7.07					
Conductivity (mS/cm)	3%	4.37	4.35	4.40					
Turbidity (NTU)	10%	0.00	0.00	0.00					
D.O. (mg/L)	10%	2.64	.97	.87					
Temperature (°C) (°F)	3%	16.81°C	16.41°C	16.18°C					
ORP ³ (mV)	+/-10 mv	127	87	25					
Appearance		C	C	C					
Free Product (Yes/No)		NONE	NONE	NONE					
Odor		NONE	NONE	NONE					
Comments	PED AT TIME SAMPLE WAS COLLECTED - 0.2 ppm								

C = Clear T = Turbid ST = Semi Turbid VT = Very Turbid



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Well Sampling Field Data Sheet

Well Casing Unit Volume

(gal/l.f.)

1 1/4" = 0.08 2" = 0.17 3" = 0.38
4" = 0.66 6" = 1.5 8" = 2.6

Client Name: _____

Site Name: CONVENTUS

Project No.: _____

Field Staff: RICK BACKERT

WELL DATA

Date	8/25/22							
Well Number	BCPMW4							
Diameter (inches)	2"							
Total Sounded Depth (feet)	14.8 ^{1/2}							
Static Water Level (feet)	7.2'							
H ₂ O Column (feet)	7.6'							
Pump Intake (feet)								
Well Volume (gallons)								
Amount to Evacuate (gallons)	29.6							
Amount Evacuated (gallons)	29.6							

FIELD READINGS

Date	8/25/22							
Time	10:45	10:50	10:55					
pH (Std. Units)	+/-0.1	7.54	7.67	7.77				
Conductivity (mS/cm)	3%	4.62	4.37	4.10				
Turbidity (NTU)	10%	0.00	0.00	0.00				
D.O. (mg/L)	10%	1.72	1.61	1.57				
Temperature (°C) (°F)	3%	17.49°C	17.16°C	16.93°C				
ORP ³ (mV)	+/-10 mv	-14	-174	-230				
Appearance		C	C	C				
Free Product (Yes/No)		NONE	NONE	NONE				
Odor		NONE	NONE	NONE				
Comments	RED AT THE TIME SAMPLE COLLECTED - 7.7 ppm							

C = Clear T = Turbid ST = Semi Turbid VT = Very Turbid



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Well Sampling Field Data Sheet

7.3

Well Casing Unit Volume

(gal/l.f.)

1 1/4" = 0.08 2" = 0.17 3" = 0.38
4" = 0.66 6" = 1.5 8" = 2.6

Client Name:

Site Name: CONVENTUS

Project No.:

Field Staff: RIEN BACKUS

WELL DATA

Date	8/25/22							
Well Number	BOPMW3							
Diameter (inches)	8"							
Total Sounded Depth (feet)	15'							
Static Water Level (feet)	7.3'							
H ₂ O Column (feet)	7.7'							
Pump Intake (feet)								
Well Volume (gallons)								
Amount to Evacuate (gallons)	2 gal							
Amount Evacuated (gallons)	2 gal							

FIELD READINGS

Date	8/25/22							
Time	11:20	11:25	11:30	11:35				
pH (Std. Units)	+/-0.1	8.06	7.98	7.93	7.93			
Conductivity (mS/cm)	3%	5.31	5.29	5.34	5.39			
Turbidity (NTU)	10%	0.00	0.00	0.00	0.00			
D.O. (mg/L)	10%	1.73	1.90	1.62	1.63			
Temperature (°C) (°F)	3%	18.11°C	17.98°C	17.21°C	16.81°C			
ORP ³ (mV)	+/-10 mv	-28	-153	-204	-213			
Appearance		C	C	C	C			
Free Product (Yes/No)		NONE	NONE	NONE	NONE			
Odor		NONE	NONE	NONE	NONE			
Comments	P.D. READING WHEN SAMPLE WAS COLLECTED - 15.7ppm							

C = Clear T = Turbid ST = Semi Turbid VT = Very Turbid



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Well Sampling Field Data Sheet

Well Casing Unit Volume

(gal/l.f.)

1 1/4" = 0.08 2" = 0.17 3" = 0.38
4" = 0.66 6" = 1.5 8" = 2.6

Client Name:

Site Name: CONVENTUS

Project No.:

Field Staff: RICH BACKERT

WELL DATA

Date	8/25/22								
Well Number	REP11010								
Diameter (inches)	8"								
Total Sounded Depth (feet)	15'								
Static Water Level (feet)	7.2'								
H ₂ O Column (feet)	7.8'								
Pump Intake (feet)									
Well Volume (gallons)									
Amount to Evacuate (gallons)	2.59 gal								
Amount Evacuated (gallons)	2.59 gal								

FIELD READINGS

Date	8/25/22								
Time	12:00	12:05	12:10	12:15					
pH (Std. Units)	+/-0.1	8.34	9.58	9.102	9.102				
Conductivity (mS/cm)	3%	11.3	11.5	11.10	11.7				
Turbidity (NTU)	10%	0.00	0.00	0.00	0.00				
D.O. (mg/L)	10%	3.93	1.710	1.65	1.63				
Temperature (°C) (°F)	3%	18.19°C	17.32°C	16.98°C	16.85°C				
ORP ³ (mV)	+/-10 mv	55	33	31	31				
Appearance		C	C	C	C				
Free Product (Yes/No)		NONE	NONE	NONE	NONE				
Odor		NONE	NONE	NONE	NONE				
Comments	PTS READING WHEN SAMPLE WAS COLLECTED - 0.2 ppm								

C = Clear T = Turbid ST = Semi Turbid VT = Very Turbid



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Well Sampling Field Data Sheet

Well Casing Unit Volume

(gal/l.f.)

1 1/4" = 0.08 2" = 0.17 3" = 0.38
4" = 0.66 6" = 1.5 8" = 2.6

Client Name: _____

Site Name: CONVENTUS

Project No.: _____

Field Staff: WILL BACKUS

WELL DATA

Date		12/14/22							
Well Number		BCEP MW-1							
Diameter (inches)		2"							
Total Sounded Depth (feet)		15'							
Static Water Level (feet)		7'							
H ₂ O Column (feet)		8'							
Pump Intake (feet)									
Well Volume (gallons)									
Amount to Evacuate (gallons)		2.8 gal							
Amount Evacuated (gallons)		2.8 gal							

FIELD READINGS

Date	Stabilization Criteria	12/14/22							
Time		10:45	10:50	10:55	11:00				
pH (Std. Units)	+/-0.1	7.10	7.12	7.13	7.12				
Conductivity (mS/cm)	3%	9.21	9.20	9.28	9.30				
Turbidity (NTU)	10%	0.00	0.00	0.00	0.00				
D.O. (mg/L)	10%	5.74	4.97	3.12	2.40				
Temperature (°C) (°F)	3%	11.97°C	12.02°C	12.10°C	12.10°C				
ORP ³ (mV)	+/-10 mv	24	23	19	18				
Appearance		C	C	C	C				
Free Product (Yes/No)		NONE	NONE	NONE	NONE				
Odor		NONE	NONE	NONE	NONE				
Comments	- PED READING OF WELL 15' @ 11:00 AM AT TIME OF SAMPLING 0.2 MPN - DUA COLLECTED FROM BCEP MW-1								

C = Clear T = Turbid ST = Semi Turbid VT = Very Turbid



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Well Sampling Field Data Sheet

Well Casing Unit Volume

(gal/l.f.)

1 1/4" = 0.08 2" = 0.17 3" = 0.38
4" = 0.66 6" = 1.5 8" = 2.6

Client Name:

Site Name:

Project No.:

Field Staff:

CONVENTUS

RICH BACKERT

WELL DATA

Date	12/14/22								
Well Number	BCPW 7								
Diameter (inches)	2"								
Total Sounded Depth (feet)	15'								
Static Water Level (feet)	10.0'								
H ₂ O Column (feet)	5'								
Pump Intake (feet)									
Well Volume (gallons)									
Amount to Evacuate (gallons)	2gal								
Amount Evacuated (gallons)	2gal								

FIELD READINGS

Date	12/14/22								
Time	11:25	11:30	11:35	11:40					
Stabilization Criteria									
pH (Std. Units)	+/-0.1	7.55	7.42	7.04	6.97				
Conductivity (mS/cm)	3%	4.56	4.50	4.50	4.52				
Turbidity (NTU)	10%	0.00	0.00	0.00	0.00				
D.O. (mg/L)	10%	2.84	2.34	2.16	1.98				
Temperature (°C) (°F)	3%	12.19°C	12.45°C	12.80°C	12.90°C				
ORP ³ (mV)	+/-10 mv	109	110	87	70				
Appearance		C	C	C	C				
Free Product (Yes/No)		NONE	NONE	NONE	NONE				
Odor		NONE	NONE	NONE	NONE				
Comments	-PEL READING FROM WELL HEAD AT TIME OF SAMPLING WAS 0.2PPM.								

C = Clear T = Turbid ST = Semi Turbid VT = Very Turbid



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Well Sampling Field Data Sheet

Well Casing Unit Volume

(gal/l.f.)

1 1/4" = 0.08 2" = 0.17 3" = 0.38
4" = 0.66 6" = 1.5 8" = 2.6

Client Name: _____

Site Name: CONVENTUS

Project No.: _____

Field Staff: ROCK BACKUS

WELL DATA

Date	12/14/22								
Well Number	Burns 4								
Diameter (inches)	2"								
Total Sounded Depth (feet)	14.8'								
Static Water Level (feet)	6.9'								
H ₂ O Column (feet)	7.9'								
Pump Intake (feet)									
Well Volume (gallons)									
Amount to Evacuate (gallons)	1.89 gal								
Amount Evacuated (gallons)	1.89 gal								

FIELD READINGS

Date	Stabilization	12/14/22							
Time	Criteria	12:10	12:18	12:20	12:26				
pH (Std. Units)	+/-0.1	7.48	7.46	7.51	7.57				
Conductivity (mS/cm)	3%	4.13	3.86	3.88	3.97				
Turbidity (NTU)	10%	0.00	0.00	0.00	0.00				
D.O. (mg/L)	10%	7.17	2.23	1.94	1.75				
Temperature (°C) (°F)	3%	12.34°C	13.41°C	13.63°C	13.68°C				
ORP ³ (mV)	+/-10 mv	-95	-181	-202	-209				
Appearance		C	C	C	C				
Free Product (Yes/No)		NONE	NONE	NONE	NONE				
Odor		YES	NONE	NONE	NONE				
Comments	RED READING IN WELL HEAD AT TIME OF SAMPLING WAS 0.3 ppm								

C = Clear T = Turbid ST = Semi Turbid VT = Very Turbid



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Well Sampling Field Data Sheet

Well Casing Unit Volume

(gal/l.f.)

1 1/4" = 0.08 2" = 0.17 3" = 0.38
4" = 0.66 6" = 1.5 8" = 2.6

Client Name: _____

Site Name: CONVENTUS

Project No.: _____

Field Staff: NINA BACKUS

WELL DATA

Date	12/14/22								
Well Number	BENUS3								
Diameter (inches)	8"								
Total Sounded Depth (feet)	15'								
Static Water Level (feet)	7'								
H ₂ O Column (feet)	8'								
Pump Intake (feet)									
Well Volume (gallons)									
Amount to Evacuate (gallons)	29.2								
Amount Evacuated (gallons)	29.2								

FIELD READINGS

Date	12/14/22								
Time	12:40	12:45	12:50	12:55					
pH (Std. Units)	+/-0.1	7.95	7.77	7.89	7.90				
Conductivity (mS/cm)	3%	5.73	5.91	5.97	5.88				
Turbidity (NTU)	10%	0.00	0.00	0.00	0.00				
D.O. (mg/L)	10%	2.95	2.45	1.51	1.49				
Temperature (°C) (°F)	3%	13.70°C	14.57°C	14.97°C	14.95°C				
ORP ³ (mV)	+/-10 mv	14	5	-12	-14				
Appearance		C	C	C	C				
Free Product (Yes/No)		NONE	NONE	NONE	NONE				
Odor		NONE	NONE	NONE	NONE				
Comments	P2D READING IN WELL HEAD AT TIME OF SAMPLING WAS 0.4 PPM								

C = Clear T = Turbid ST = Semi Turbid VT = Very Turbid



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Well Sampling Field Data Sheet

Well Casing Unit Volume

(gal/l.f.)

1 1/4" = 0.08 2" = 0.17 3" = 0.38
4" = 0.66 6" = 1.5 8" = 2.6

Client Name: _____

Site Name: CONVENTUS

Project No.: _____

Field Staff: RICH BACKERT

WELL DATA

Date		12/14/22							
Well Number		300421A							
Diameter (inches)		8"							
Total Sounded Depth (feet)		15'							
Static Water Level (feet)		7.1'							
H ₂ O Column (feet)		7.9'							
Pump Intake (feet)									
Well Volume (gallons)									
Amount to Evacuate (gallons)		2.59							
Amount Evacuated (gallons)		2.59							

FIELD READINGS

Date	Stabilization	12/14/22							
Time	Criteria	1:18	1:20	1:25	1:30				
pH (Std. Units)	+/-0.1	8.24	8.90	9.13	9.33				
Conductivity (mS/cm)	3%	11.1	11.1	11.2	11.2				
Turbidity (NTU)	10%	0.00	0.00	0.00	0.00				
D.O. (mg/L)	10%	7.37	6.37	5.85	5.50				
Temperature (°C) (°F)	3%	13.07°C	13.47°C	13.54°C	13.63°C				
ORP ³ (mV)	+/-10 mv	37	25	22	21				
Appearance		C	C	C	C				
Free Product (Yes/No)		NONE	NONE	NONE	NONE				
Odor		NONE	NONE	NONE	NONE				
Comments	PID DRIVING FROM WELL HEAD AT TIME OF SAMPLING 0.4 PPM								

C = Clear T = Turbid ST = Semi Turbid VT = Very Turbid



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Well Sampling Field Data Sheet

Well Casing Unit Volume

(gal/l.f.)

1 1/4" = 0.08 2" = 0.17 3" = 0.38
4" = 0.66 6" = 1.5 8" = 2.6

Client Name: _____

Site Name: CONVENTUS

Project No.: _____

Field Staff: MICHAEL RACKRAT

WELL DATA

Date	12/14/22								
Well Number	BCMHWS								
Diameter (inches)	2.5"								
Total Sounded Depth (feet)	15'								
Static Water Level (feet)	7.7'								
H ₂ O Column (feet)	7.8'								
Pump Intake (feet)									
Well Volume (gallons)									
Amount to Evacuate (gallons)	2.59 gal								
Amount Evacuated (gallons)	2.89 gal								

FIELD READINGS

Date	12/14/22								
Time	1:58	2:00	2:08	2:10					
pH (Std. Units)	+/-0.1	8.77	8.54	8.15	8.08				
Conductivity (mS/cm)	3%	116.0	15.3	9.15	8.165				
Turbidity (NTU)	10%	103.5	32.7	0.00	0.00				
D.O. (mg/L)	10%	1.80	1.30	4.75	5.79				
Temperature (°C) (°F)	3%	13.78°C	14.08°C	14.22°C	14.16°C				
ORP ³ (mV)	+/-10 mv	-86	-33	-117	-124				
Appearance		C	C	C	C				
Free Product (Yes/No)		NONE	NONE	NONE	NONE				
Odor		YES	YES	NONE	NONE				
Comments	PEA READING OF WELL HEAD AT TIME OF SAMPLE WAS 1.0 ppm								

C = Clear T = Turbid ST = Semi Turbid VT = Very Turbid



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Well Sampling Field Data Sheet

Well Casing Unit Volume		
(gal/l.f.)		
1 1/4" = 0.08	2" = 0.17	3" = 0.38
4" = 0.66	6" = 1.5	8" = 2.6

Client Name: _____
Site Name: CONVENTUS
Project No.: _____
Field Staff: RICH BACIKERT

WELL DATA

Date	8/28/22								
Well Number	RCR11W5								
Diameter (inches)	2"								
Total Sounded Depth (feet)	15'								
Static Water Level (feet)	8.2'								
H ₂ O Column (feet)	6.8'								
Pump Intake (feet)									
Well Volume (gallons)									
Amount to Evacuate (gallons)	324								
Amount Evacuated (gallons)	394								

FIELD READINGS

Date	8/28/22								
Time	12:40	12:45	12:50	12:55	1:00				
pH (Std. Units)	+/-0.1	9.18	8.55	8.50	8.48	8.46			
Conductivity (mS/cm)	3%	9.16	8.24	8.01	7.18	6.87			
Turbidity (NTU)	10%	0.00	0.00	0.00	0.00	0.00			
D.O. (mg/L)	10%	1.81	1.57	1.58	1.67	1.70			
Temperature (°C) (°F)	3%	17.25°C	17.34°C	17.04°C	16.71°C	16.96°C			
ORP ³ (mV)	+/-10 mv	116	-132	-150	-173	-176			
Appearance		C	C	C	C	C			
Free Product (Yes/No)		none	none	none	none	none			
Odor		none	none	none	at 10	petroleum			
Comments	PID READING AT TIME SAMPLE COLLECTED - 73.5 ppm								

C = Clear T = Turbid ST = Semi Turbid VT = Very Turbid

APPENDIX A

LABORATORY ANALYTICAL RESULTS



ANALYTICAL REPORT

Lab Number:	L2246299
Client:	C&S Companies 141 Elm Street Suite 100 Buffalo, NY 14203
ATTN:	Richard Backert
Phone:	(716) 955-3024
Project Name:	CONVENTUS
Project Number:	U86
Report Date:	09/09/22

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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).

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Project Name: CONVENTUS
Project Number: U86

Lab Number: L2246299
Report Date: 09/09/22

Alpha Sample ID	Client ID	Matrix	Sample Location	Collection Date/Time	Receive Date
L2246299-01	BCP MW-1	WATER	CONVENTUS/MAIN ST	08/25/22 09:55	08/25/22
L2246299-02	BCP MW-7	WATER	CONVENTUS/MAIN ST	08/25/22 10:25	08/25/22
L2246299-03	BCP MW-4	WATER	CONVENTUS/MAIN ST	08/25/22 10:55	08/25/22
L2246299-04	BCP MW-3	WATER	CONVENTUS/MAIN ST	08/25/22 11:35	08/25/22
L2246299-05	BCP MW-6	WATER	CONVENTUS/MAIN ST	08/25/22 12:15	08/25/22
L2246299-06	BCP MW-5	WATER	CONVENTUS/MAIN ST	08/25/22 13:00	08/25/22
L2246299-07	DUP	WATER	CONVENTUS/MAIN ST	08/25/22 09:55	08/25/22
L2246299-08	TRIP BLANK	WATER	CONVENTUS/MAIN ST	08/25/22 13:50	08/25/22

Project Name: CONVENTUS
Project Number: U86

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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: CONVENTUS
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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.

Volatile Organics

L2246299-05: The pH was greater than two; however, the sample was analyzed within the method required holding time.

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:

Tiffani Morrissey - Tiffani Morrissey

Title: Technical Director/Representative

Date: 09/09/22

ORGANICS

VOLATILES

Project Name: CONVENTUS

Lab Number: L2246299

Project Number: U86

Report Date: 09/09/22

SAMPLE RESULTS

Lab ID: L2246299-01
 Client ID: BCP MW-1
 Sample Location: CONVENTUS/MAIN ST

Date Collected: 08/25/22 09:55
 Date Received: 08/25/22
 Field Prep: Not Specified

Sample Depth:

Matrix: Water
 Analytical Method: 1,8260C
 Analytical Date: 09/01/22 20:47
 Analyst: MV

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: CONVENTUS

Lab Number: L2246299

Project Number: U86

Report Date: 09/09/22

SAMPLE RESULTS

Lab ID: L2246299-01

Date Collected: 08/25/22 09:55

Client ID: BCP MW-1

Date Received: 08/25/22

Sample Location: CONVENTUS/MAIN ST

Field Prep: 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
1,2-Dibromoethane	ND		ug/l	2.0	0.65	1
n-Butylbenzene	ND		ug/l	2.5	0.70	1
sec-Butylbenzene	ND		ug/l	2.5	0.70	1
tert-Butylbenzene	ND		ug/l	2.5	0.70	1
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70	1
Isopropylbenzene	ND		ug/l	2.5	0.70	1
p-Isopropyltoluene	ND		ug/l	2.5	0.70	1
Naphthalene	ND		ug/l	2.5	0.70	1
n-Propylbenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70	1
1,3,5-Trimethylbenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trimethylbenzene	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
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	99		70-130
Toluene-d8	103		70-130
4-Bromofluorobenzene	95		70-130
Dibromofluoromethane	104		70-130

Project Name: CONVENTUS

Lab Number: L2246299

Project Number: U86

Report Date: 09/09/22

SAMPLE RESULTS

Lab ID: L2246299-02
 Client ID: BCP MW-7
 Sample Location: CONVENTUS/MAIN ST

Date Collected: 08/25/22 10:25
 Date Received: 08/25/22
 Field Prep: Not Specified

Sample Depth:

Matrix: Water
 Analytical Method: 1,8260C
 Analytical Date: 09/01/22 21:10
 Analyst: MV

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: CONVENTUS

Lab Number: L2246299

Project Number: U86

Report Date: 09/09/22

SAMPLE RESULTS

Lab ID: L2246299-02
 Client ID: BCP MW-7
 Sample Location: CONVENTUS/MAIN ST

Date Collected: 08/25/22 10:25
 Date Received: 08/25/22
 Field Prep: 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
1,2-Dibromoethane	ND		ug/l	2.0	0.65	1
n-Butylbenzene	ND		ug/l	2.5	0.70	1
sec-Butylbenzene	ND		ug/l	2.5	0.70	1
tert-Butylbenzene	ND		ug/l	2.5	0.70	1
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70	1
Isopropylbenzene	ND		ug/l	2.5	0.70	1
p-Isopropyltoluene	ND		ug/l	2.5	0.70	1
Naphthalene	ND		ug/l	2.5	0.70	1
n-Propylbenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70	1
1,3,5-Trimethylbenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trimethylbenzene	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
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	94		70-130
Toluene-d8	104		70-130
4-Bromofluorobenzene	96		70-130
Dibromofluoromethane	103		70-130

Project Name: CONVENTUS

Lab Number: L2246299

Project Number: U86

Report Date: 09/09/22

SAMPLE RESULTS

Lab ID: L2246299-03 D
 Client ID: BCP MW-4
 Sample Location: CONVENTUS/MAIN ST

Date Collected: 08/25/22 10:55
 Date Received: 08/25/22
 Field Prep: Not Specified

Sample Depth:

Matrix: Water
 Analytical Method: 1,8260C
 Analytical Date: 09/02/22 03:00
 Analyst: MV

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westborough Lab						
Methylene chloride	ND		ug/l	25	7.0	10
1,1-Dichloroethane	ND		ug/l	25	7.0	10
Chloroform	ND		ug/l	25	7.0	10
Carbon tetrachloride	ND		ug/l	5.0	1.3	10
1,2-Dichloropropane	ND		ug/l	10	1.4	10
Dibromochloromethane	ND		ug/l	5.0	1.5	10
1,1,2-Trichloroethane	ND		ug/l	15	5.0	10
Tetrachloroethene	ND		ug/l	5.0	1.8	10
Chlorobenzene	ND		ug/l	25	7.0	10
Trichlorofluoromethane	ND		ug/l	25	7.0	10
1,2-Dichloroethane	ND		ug/l	5.0	1.3	10
1,1,1-Trichloroethane	ND		ug/l	25	7.0	10
Bromodichloromethane	ND		ug/l	5.0	1.9	10
trans-1,3-Dichloropropene	ND		ug/l	5.0	1.6	10
cis-1,3-Dichloropropene	ND		ug/l	5.0	1.4	10
Bromoform	ND		ug/l	20	6.5	10
1,1,2,2-Tetrachloroethane	ND		ug/l	5.0	1.7	10
Benzene	5.0		ug/l	5.0	1.6	10
Toluene	34		ug/l	25	7.0	10
Ethylbenzene	1100		ug/l	25	7.0	10
Chloromethane	ND		ug/l	25	7.0	10
Bromomethane	ND		ug/l	25	7.0	10
Vinyl chloride	ND		ug/l	10	0.71	10
Chloroethane	ND		ug/l	25	7.0	10
1,1-Dichloroethene	ND		ug/l	5.0	1.7	10
trans-1,2-Dichloroethene	ND		ug/l	25	7.0	10
Trichloroethene	ND		ug/l	5.0	1.8	10
1,2-Dichlorobenzene	ND		ug/l	25	7.0	10

Project Name: CONVENTUS

Lab Number: L2246299

Project Number: U86

Report Date: 09/09/22

SAMPLE RESULTS

Lab ID: L2246299-03 D
 Client ID: BCP MW-4
 Sample Location: CONVENTUS/MAIN ST

Date Collected: 08/25/22 10:55
 Date Received: 08/25/22
 Field Prep: 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	25	7.0	10
1,4-Dichlorobenzene	ND		ug/l	25	7.0	10
Methyl tert butyl ether	ND		ug/l	25	7.0	10
p/m-Xylene	14	J	ug/l	25	7.0	10
o-Xylene	12	J	ug/l	25	7.0	10
cis-1,2-Dichloroethene	ND		ug/l	25	7.0	10
Styrene	ND		ug/l	25	7.0	10
Dichlorodifluoromethane	ND		ug/l	50	10.	10
Acetone	ND		ug/l	50	15.	10
Carbon disulfide	ND		ug/l	50	10.	10
2-Butanone	ND		ug/l	50	19.	10
4-Methyl-2-pentanone	ND		ug/l	50	10.	10
2-Hexanone	ND		ug/l	50	10.	10
1,2-Dibromoethane	ND		ug/l	20	6.5	10
n-Butylbenzene	9.4	J	ug/l	25	7.0	10
sec-Butylbenzene	ND		ug/l	25	7.0	10
tert-Butylbenzene	ND		ug/l	25	7.0	10
1,2-Dibromo-3-chloropropane	ND		ug/l	25	7.0	10
Isopropylbenzene	24	J	ug/l	25	7.0	10
p-Isopropyltoluene	ND		ug/l	25	7.0	10
Naphthalene	200		ug/l	25	7.0	10
n-Propylbenzene	140		ug/l	25	7.0	10
1,2,4-Trichlorobenzene	ND		ug/l	25	7.0	10
1,3,5-Trimethylbenzene	ND		ug/l	25	7.0	10
1,2,4-Trimethylbenzene	1200		ug/l	25	7.0	10
Methyl Acetate	ND		ug/l	20	2.3	10
Cyclohexane	130		ug/l	100	2.7	10
Freon-113	ND		ug/l	25	7.0	10
Methyl cyclohexane	51	J	ug/l	100	4.0	10

Surrogate	% Recovery	Qualifier	Acceptance Criteria
1,2-Dichloroethane-d4	96		70-130
Toluene-d8	104		70-130
4-Bromofluorobenzene	91		70-130
Dibromofluoromethane	97		70-130

Project Name: CONVENTUS

Lab Number: L2246299

Project Number: U86

Report Date: 09/09/22

SAMPLE RESULTS

Lab ID: L2246299-04 D
 Client ID: BCP MW-3
 Sample Location: CONVENTUS/MAIN ST

Date Collected: 08/25/22 11:35
 Date Received: 08/25/22
 Field Prep: Not Specified

Sample Depth:

Matrix: Water
 Analytical Method: 1,8260C
 Analytical Date: 09/02/22 03:23
 Analyst: MV

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westborough Lab						
Methylene chloride	ND		ug/l	12	3.5	5
1,1-Dichloroethane	ND		ug/l	12	3.5	5
Chloroform	ND		ug/l	12	3.5	5
Carbon tetrachloride	ND		ug/l	2.5	0.67	5
1,2-Dichloropropane	ND		ug/l	5.0	0.68	5
Dibromochloromethane	ND		ug/l	2.5	0.74	5
1,1,2-Trichloroethane	ND		ug/l	7.5	2.5	5
Tetrachloroethene	ND		ug/l	2.5	0.90	5
Chlorobenzene	ND		ug/l	12	3.5	5
Trichlorofluoromethane	ND		ug/l	12	3.5	5
1,2-Dichloroethane	ND		ug/l	2.5	0.66	5
1,1,1-Trichloroethane	ND		ug/l	12	3.5	5
Bromodichloromethane	ND		ug/l	2.5	0.96	5
trans-1,3-Dichloropropene	ND		ug/l	2.5	0.82	5
cis-1,3-Dichloropropene	ND		ug/l	2.5	0.72	5
Bromoform	ND		ug/l	10	3.2	5
1,1,2,2-Tetrachloroethane	ND		ug/l	2.5	0.84	5
Benzene	1.7	J	ug/l	2.5	0.80	5
Toluene	22		ug/l	12	3.5	5
Ethylbenzene	380		ug/l	12	3.5	5
Chloromethane	ND		ug/l	12	3.5	5
Bromomethane	ND		ug/l	12	3.5	5
Vinyl chloride	ND		ug/l	5.0	0.36	5
Chloroethane	ND		ug/l	12	3.5	5
1,1-Dichloroethene	ND		ug/l	2.5	0.84	5
trans-1,2-Dichloroethene	ND		ug/l	12	3.5	5
Trichloroethene	ND		ug/l	2.5	0.88	5
1,2-Dichlorobenzene	ND		ug/l	12	3.5	5

Project Name: CONVENTUS

Lab Number: L2246299

Project Number: U86

Report Date: 09/09/22

SAMPLE RESULTS

Lab ID: L2246299-04 D
 Client ID: BCP MW-3
 Sample Location: CONVENTUS/MAIN ST

Date Collected: 08/25/22 11:35
 Date Received: 08/25/22
 Field Prep: 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	12	3.5	5
1,4-Dichlorobenzene	ND		ug/l	12	3.5	5
Methyl tert butyl ether	ND		ug/l	12	3.5	5
p/m-Xylene	480		ug/l	12	3.5	5
o-Xylene	12		ug/l	12	3.5	5
cis-1,2-Dichloroethene	ND		ug/l	12	3.5	5
Styrene	ND		ug/l	12	3.5	5
Dichlorodifluoromethane	ND		ug/l	25	5.0	5
Acetone	ND		ug/l	25	7.3	5
Carbon disulfide	ND		ug/l	25	5.0	5
2-Butanone	ND		ug/l	25	9.7	5
4-Methyl-2-pentanone	ND		ug/l	25	5.0	5
2-Hexanone	ND		ug/l	25	5.0	5
1,2-Dibromoethane	ND		ug/l	10	3.2	5
n-Butylbenzene	ND		ug/l	12	3.5	5
sec-Butylbenzene	ND		ug/l	12	3.5	5
tert-Butylbenzene	ND		ug/l	12	3.5	5
1,2-Dibromo-3-chloropropane	ND		ug/l	12	3.5	5
Isopropylbenzene	5.2	J	ug/l	12	3.5	5
p-Isopropyltoluene	ND		ug/l	12	3.5	5
Naphthalene	170		ug/l	12	3.5	5
n-Propylbenzene	18		ug/l	12	3.5	5
1,2,4-Trichlorobenzene	ND		ug/l	12	3.5	5
1,3,5-Trimethylbenzene	47		ug/l	12	3.5	5
1,2,4-Trimethylbenzene	560		ug/l	12	3.5	5
Methyl Acetate	ND		ug/l	10	1.2	5
Cyclohexane	40	J	ug/l	50	1.4	5
Freon-113	ND		ug/l	12	3.5	5
Methyl cyclohexane	32	J	ug/l	50	2.0	5

Surrogate	% Recovery	Qualifier	Acceptance Criteria
1,2-Dichloroethane-d4	94		70-130
Toluene-d8	106		70-130
4-Bromofluorobenzene	88		70-130
Dibromofluoromethane	100		70-130

Project Name: CONVENTUS

Lab Number: L2246299

Project Number: U86

Report Date: 09/09/22

SAMPLE RESULTS

Lab ID: L2246299-05
 Client ID: BCP MW-6
 Sample Location: CONVENTUS/MAIN ST

Date Collected: 08/25/22 12:15
 Date Received: 08/25/22
 Field Prep: Not Specified

Sample Depth:

Matrix: Water
 Analytical Method: 1,8260C
 Analytical Date: 09/01/22 21:33
 Analyst: MV

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: CONVENTUS

Lab Number: L2246299

Project Number: U86

Report Date: 09/09/22

SAMPLE RESULTS

Lab ID: L2246299-05
 Client ID: BCP MW-6
 Sample Location: CONVENTUS/MAIN ST

Date Collected: 08/25/22 12:15
 Date Received: 08/25/22
 Field Prep: 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
1,2-Dibromoethane	ND		ug/l	2.0	0.65	1
n-Butylbenzene	ND		ug/l	2.5	0.70	1
sec-Butylbenzene	ND		ug/l	2.5	0.70	1
tert-Butylbenzene	ND		ug/l	2.5	0.70	1
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70	1
Isopropylbenzene	ND		ug/l	2.5	0.70	1
p-Isopropyltoluene	ND		ug/l	2.5	0.70	1
Naphthalene	ND		ug/l	2.5	0.70	1
n-Propylbenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70	1
1,3,5-Trimethylbenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trimethylbenzene	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
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	98		70-130
Toluene-d8	103		70-130
4-Bromofluorobenzene	92		70-130
Dibromofluoromethane	104		70-130

Project Name: CONVENTUS

Lab Number: L2246299

Project Number: U86

Report Date: 09/09/22

SAMPLE RESULTS

Lab ID: L2246299-06 D
 Client ID: BCP MW-5
 Sample Location: CONVENTUS/MAIN ST

Date Collected: 08/25/22 13:00
 Date Received: 08/25/22
 Field Prep: Not Specified

Sample Depth:

Matrix: Water
 Analytical Method: 1,8260C
 Analytical Date: 09/02/22 03:46
 Analyst: MV

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westborough Lab						
Methylene chloride	ND		ug/l	50	14.	20
1,1-Dichloroethane	ND		ug/l	50	14.	20
Chloroform	ND		ug/l	50	14.	20
Carbon tetrachloride	ND		ug/l	10	2.7	20
1,2-Dichloropropane	ND		ug/l	20	2.7	20
Dibromochloromethane	ND		ug/l	10	3.0	20
1,1,2-Trichloroethane	ND		ug/l	30	10.	20
Tetrachloroethene	ND		ug/l	10	3.6	20
Chlorobenzene	ND		ug/l	50	14.	20
Trichlorofluoromethane	ND		ug/l	50	14.	20
1,2-Dichloroethane	ND		ug/l	10	2.6	20
1,1,1-Trichloroethane	ND		ug/l	50	14.	20
Bromodichloromethane	ND		ug/l	10	3.8	20
trans-1,3-Dichloropropene	ND		ug/l	10	3.3	20
cis-1,3-Dichloropropene	ND		ug/l	10	2.9	20
Bromoform	ND		ug/l	40	13.	20
1,1,2,2-Tetrachloroethane	ND		ug/l	10	3.3	20
Benzene	ND		ug/l	10	3.2	20
Toluene	26	J	ug/l	50	14.	20
Ethylbenzene	2100		ug/l	50	14.	20
Chloromethane	ND		ug/l	50	14.	20
Bromomethane	ND		ug/l	50	14.	20
Vinyl chloride	ND		ug/l	20	1.4	20
Chloroethane	ND		ug/l	50	14.	20
1,1-Dichloroethene	ND		ug/l	10	3.4	20
trans-1,2-Dichloroethene	ND		ug/l	50	14.	20
Trichloroethene	ND		ug/l	10	3.5	20
1,2-Dichlorobenzene	ND		ug/l	50	14.	20

Project Name: CONVENTUS

Lab Number: L2246299

Project Number: U86

Report Date: 09/09/22

SAMPLE RESULTS

Lab ID: L2246299-06 D
 Client ID: BCP MW-5
 Sample Location: CONVENTUS/MAIN ST

Date Collected: 08/25/22 13:00
 Date Received: 08/25/22
 Field Prep: 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	50	14.	20
1,4-Dichlorobenzene	ND		ug/l	50	14.	20
Methyl tert butyl ether	ND		ug/l	50	14.	20
p/m-Xylene	7800		ug/l	50	14.	20
o-Xylene	100		ug/l	50	14.	20
cis-1,2-Dichloroethene	ND		ug/l	50	14.	20
Styrene	ND		ug/l	50	14.	20
Dichlorodifluoromethane	ND		ug/l	100	20.	20
Acetone	ND		ug/l	100	29.	20
Carbon disulfide	ND		ug/l	100	20.	20
2-Butanone	ND		ug/l	100	39.	20
4-Methyl-2-pentanone	ND		ug/l	100	20.	20
2-Hexanone	ND		ug/l	100	20.	20
1,2-Dibromoethane	ND		ug/l	40	13.	20
n-Butylbenzene	ND		ug/l	50	14.	20
sec-Butylbenzene	ND		ug/l	50	14.	20
tert-Butylbenzene	ND		ug/l	50	14.	20
1,2-Dibromo-3-chloropropane	ND		ug/l	50	14.	20
Isopropylbenzene	29	J	ug/l	50	14.	20
p-Isopropyltoluene	ND		ug/l	50	14.	20
Naphthalene	840		ug/l	50	14.	20
n-Propylbenzene	130		ug/l	50	14.	20
1,2,4-Trichlorobenzene	ND		ug/l	50	14.	20
1,3,5-Trimethylbenzene	630		ug/l	50	14.	20
1,2,4-Trimethylbenzene	3100		ug/l	50	14.	20
Methyl Acetate	ND		ug/l	40	4.7	20
Cyclohexane	240		ug/l	200	5.4	20
Freon-113	ND		ug/l	50	14.	20
Methyl cyclohexane	95	J	ug/l	200	7.9	20

Surrogate	% Recovery	Qualifier	Acceptance Criteria
1,2-Dichloroethane-d4	93		70-130
Toluene-d8	103		70-130
4-Bromofluorobenzene	89		70-130
Dibromofluoromethane	96		70-130

Project Name: CONVENTUS

Lab Number: L2246299

Project Number: U86

Report Date: 09/09/22

SAMPLE RESULTS

Lab ID: L2246299-07
 Client ID: DUP
 Sample Location: CONVENTUS/MAIN ST

Date Collected: 08/25/22 09:55
 Date Received: 08/25/22
 Field Prep: Not Specified

Sample Depth:

Matrix: Water
 Analytical Method: 1,8260C
 Analytical Date: 09/01/22 21:57
 Analyst: MV

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: CONVENTUS

Lab Number: L2246299

Project Number: U86

Report Date: 09/09/22

SAMPLE RESULTS

Lab ID: L2246299-07

Date Collected: 08/25/22 09:55

Client ID: DUP

Date Received: 08/25/22

Sample Location: CONVENTUS/MAIN ST

Field Prep: 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
1,2-Dibromoethane	ND		ug/l	2.0	0.65	1
n-Butylbenzene	ND		ug/l	2.5	0.70	1
sec-Butylbenzene	ND		ug/l	2.5	0.70	1
tert-Butylbenzene	ND		ug/l	2.5	0.70	1
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70	1
Isopropylbenzene	ND		ug/l	2.5	0.70	1
p-Isopropyltoluene	ND		ug/l	2.5	0.70	1
Naphthalene	ND		ug/l	2.5	0.70	1
n-Propylbenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70	1
1,3,5-Trimethylbenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trimethylbenzene	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
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	99		70-130
Toluene-d8	102		70-130
4-Bromofluorobenzene	95		70-130
Dibromofluoromethane	104		70-130

Project Name: CONVENTUS

Lab Number: L2246299

Project Number: U86

Report Date: 09/09/22

SAMPLE RESULTS

Lab ID: L2246299-08
 Client ID: TRIP BLANK
 Sample Location: CONVENTUS/MAIN ST

Date Collected: 08/25/22 13:50
 Date Received: 08/25/22
 Field Prep: Not Specified

Sample Depth:

Matrix: Water
 Analytical Method: 1,8260C
 Analytical Date: 09/01/22 20:23
 Analyst: MV

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: CONVENTUS

Lab Number: L2246299

Project Number: U86

Report Date: 09/09/22

SAMPLE RESULTS

Lab ID: L2246299-08
 Client ID: TRIP BLANK
 Sample Location: CONVENTUS/MAIN ST

Date Collected: 08/25/22 13:50
 Date Received: 08/25/22
 Field Prep: 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
1,2-Dibromoethane	ND		ug/l	2.0	0.65	1
n-Butylbenzene	ND		ug/l	2.5	0.70	1
sec-Butylbenzene	ND		ug/l	2.5	0.70	1
tert-Butylbenzene	ND		ug/l	2.5	0.70	1
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70	1
Isopropylbenzene	ND		ug/l	2.5	0.70	1
p-Isopropyltoluene	ND		ug/l	2.5	0.70	1
Naphthalene	ND		ug/l	2.5	0.70	1
n-Propylbenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70	1
1,3,5-Trimethylbenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trimethylbenzene	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
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	93		70-130
Toluene-d8	103		70-130
4-Bromofluorobenzene	96		70-130
Dibromofluoromethane	102		70-130

Project Name: CONVENTUS

Lab Number: L2246299

Project Number: U86

Report Date: 09/09/22

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260C
 Analytical Date: 09/01/22 20:00
 Analyst: AJK

Parameter	Result	Qualifier	Units	RL	MDL
Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-08 Batch: WG1682814-5					
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
1,1,1,2-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

Project Name: CONVENTUS

Lab Number: L2246299

Project Number: U86

Report Date: 09/09/22

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260C
 Analytical Date: 09/01/22 20:00
 Analyst: AJK

Parameter	Result	Qualifier	Units	RL	MDL
Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-08 Batch: WG1682814-5					
1,4-Dichlorobenzene	ND		ug/l	2.5	0.70
Methyl tert butyl ether	ND		ug/l	2.5	0.70
p/m-Xylene	ND		ug/l	2.5	0.70
o-Xylene	ND		ug/l	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/l	5.0	1.0
Acetone	ND		ug/l	5.0	1.5
Carbon disulfide	ND		ug/l	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
1,2-Dibromoethane	ND		ug/l	2.0	0.65
n-Butylbenzene	ND		ug/l	2.5	0.70
sec-Butylbenzene	ND		ug/l	2.5	0.70
tert-Butylbenzene	ND		ug/l	2.5	0.70
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70
Isopropylbenzene	ND		ug/l	2.5	0.70
p-Isopropyltoluene	ND		ug/l	2.5	0.70
Naphthalene	ND		ug/l	2.5	0.70
n-Propylbenzene	ND		ug/l	2.5	0.70
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70
1,3,5-Trimethylbenzene	ND		ug/l	2.5	0.70
1,2,4-Trimethylbenzene	ND		ug/l	2.5	0.70
Methyl Acetate	ND		ug/l	2.0	0.23
Cyclohexane	ND		ug/l	10	0.27
Freon-113	ND		ug/l	2.5	0.70
Methyl cyclohexane	ND		ug/l	10	0.40

Project Name: CONVENTUS**Project Number:** U86**Lab Number:** L2246299**Report Date:** 09/09/22**Method Blank Analysis**
Batch Quality Control

Analytical Method: 1,8260C
Analytical Date: 09/01/22 20:00
Analyst: AJK

Parameter	Result	Qualifier	Units	RL	MDL
Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-08 Batch: WG1682814-5					

Surrogate	%Recovery	Qualifier	Acceptance Criteria
1,2-Dichloroethane-d4	94		70-130
Toluene-d8	103		70-130
4-Bromofluorobenzene	94		70-130
Dibromofluoromethane	102		70-130

Lab Control Sample Analysis Batch Quality Control

Project Name: CONVENTUS

Project Number: U86

Lab Number: L2246299

Report Date: 09/09/22

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-08 Batch: WG1682814-3 WG1682814-4								
Methylene chloride	97		97		70-130	0		20
1,1-Dichloroethane	98		94		70-130	4		20
Chloroform	97		92		70-130	5		20
Carbon tetrachloride	96		90		63-132	6		20
1,2-Dichloropropane	96		94		70-130	2		20
Dibromochloromethane	93		92		63-130	1		20
1,1,2-Trichloroethane	95		94		70-130	1		20
Tetrachloroethene	100		96		70-130	4		20
Chlorobenzene	98		95		75-130	3		20
Trichlorofluoromethane	130		120		62-150	8		20
1,2-Dichloroethane	91		94		70-130	3		20
1,1,1-Trichloroethane	97		92		67-130	5		20
Bromodichloromethane	92		92		67-130	0		20
trans-1,3-Dichloropropene	90		89		70-130	1		20
cis-1,3-Dichloropropene	91		89		70-130	2		20
Bromoform	82		80		54-136	2		20
1,1,2,2-Tetrachloroethane	86		86		67-130	0		20
Benzene	99		94		70-130	5		20
Toluene	100		96		70-130	4		20
Ethylbenzene	98		92		70-130	6		20
Chloromethane	87		85		64-130	2		20
Bromomethane	120		110		39-139	9		20
Vinyl chloride	100		99		55-140	1		20

Lab Control Sample Analysis Batch Quality Control

Project Name: CONVENTUS

Project Number: U86

Lab Number: L2246299

Report Date: 09/09/22

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-08 Batch: WG1682814-3 WG1682814-4								
Chloroethane	140	Q	130		55-138	7		20
1,1-Dichloroethene	100		93		61-145	7		20
trans-1,2-Dichloroethene	100		92		70-130	8		20
Trichloroethene	94		86		70-130	9		20
1,2-Dichlorobenzene	94		94		70-130	0		20
1,3-Dichlorobenzene	96		93		70-130	3		20
1,4-Dichlorobenzene	95		92		70-130	3		20
Methyl tert butyl ether	82		80		63-130	2		20
p/m-Xylene	95		95		70-130	0		20
o-Xylene	95		90		70-130	5		20
cis-1,2-Dichloroethene	98		95		70-130	3		20
Styrene	95		90		70-130	5		20
Dichlorodifluoromethane	100		93		36-147	7		20
Acetone	80		82		58-148	2		20
Carbon disulfide	100		94		51-130	6		20
2-Butanone	79		84		63-138	6		20
4-Methyl-2-pentanone	78		83		59-130	6		20
2-Hexanone	69		67		57-130	3		20
1,2-Dibromoethane	94		97		70-130	3		20
n-Butylbenzene	99		96		53-136	3		20
sec-Butylbenzene	100		95		70-130	5		20
tert-Butylbenzene	97		92		70-130	5		20
1,2-Dibromo-3-chloropropane	77		78		41-144	1		20

Lab Control Sample Analysis

Batch Quality Control

Project Name: CONVENTUS

Project Number: U86

Lab Number: L2246299

Report Date: 09/09/22

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-08 Batch: WG1682814-3 WG1682814-4								
Isopropylbenzene	98		92		70-130	6		20
p-Isopropyltoluene	97		92		70-130	5		20
Naphthalene	80		84		70-130	5		20
n-Propylbenzene	99		94		69-130	5		20
1,2,4-Trichlorobenzene	92		91		70-130	1		20
1,3,5-Trimethylbenzene	94		91		64-130	3		20
1,2,4-Trimethylbenzene	94		91		70-130	3		20
Methyl Acetate	77		84		70-130	9		20
Cyclohexane	96		91		70-130	5		20
Freon-113	110		97		70-130	13		20
Methyl cyclohexane	97		92		70-130	5		20

Surrogate	LCS %Recovery	Qual	LCSD %Recovery	Qual	Acceptance Criteria
1,2-Dichloroethane-d4	101		99		70-130
Toluene-d8	106		105		70-130
4-Bromofluorobenzene	94		91		70-130
Dibromofluoromethane	104		102		70-130

Project Name: CONVENTUS**Lab Number:** L2246299**Project Number:** U86**Report Date:** 09/09/22**Sample Receipt and Container Information**

Were project specific reporting limits specified?

YES

Cooler Information

Cooler	Custody Seal
A	Absent

Container Information

Container ID	Container Type	Cooler	Initial pH	Final pH	Temp deg C	Pres	Seal	Frozen Date/Time	Analysis(*)
L2246299-01A	Vial HCl preserved	A	NA		2.7	Y	Absent		NYTCL-8260-R2(14)
L2246299-01B	Vial HCl preserved	A	NA		2.7	Y	Absent		NYTCL-8260-R2(14)
L2246299-01C	Vial HCl preserved	A	NA		2.7	Y	Absent		NYTCL-8260-R2(14)
L2246299-02A	Vial HCl preserved	A	NA		2.7	Y	Absent		NYTCL-8260-R2(14)
L2246299-02B	Vial HCl preserved	A	NA		2.7	Y	Absent		NYTCL-8260-R2(14)
L2246299-02C	Vial HCl preserved	A	NA		2.7	Y	Absent		NYTCL-8260-R2(14)
L2246299-03A	Vial HCl preserved	A	NA		2.7	Y	Absent		NYTCL-8260-R2(14)
L2246299-03B	Vial HCl preserved	A	NA		2.7	Y	Absent		NYTCL-8260-R2(14)
L2246299-03C	Vial HCl preserved	A	NA		2.7	Y	Absent		NYTCL-8260-R2(14)
L2246299-04A	Vial HCl preserved	A	NA		2.7	Y	Absent		NYTCL-8260-R2(14)
L2246299-04B	Vial HCl preserved	A	NA		2.7	Y	Absent		NYTCL-8260-R2(14)
L2246299-04C	Vial HCl preserved	A	NA		2.7	Y	Absent		NYTCL-8260-R2(14)
L2246299-05A	Vial HCl preserved	A	NA		2.7	Y	Absent		NYTCL-8260-R2(14)
L2246299-05B	Vial HCl preserved	A	NA		2.7	Y	Absent		NYTCL-8260-R2(14)
L2246299-05C	Vial HCl preserved	A	NA		2.7	Y	Absent		NYTCL-8260-R2(14)
L2246299-06A	Vial HCl preserved	A	NA		2.7	Y	Absent		NYTCL-8260-R2(14)
L2246299-06B	Vial HCl preserved	A	NA		2.7	Y	Absent		NYTCL-8260-R2(14)
L2246299-06C	Vial HCl preserved	A	NA		2.7	Y	Absent		NYTCL-8260-R2(14)
L2246299-07A	Vial HCl preserved	A	NA		2.7	Y	Absent		NYTCL-8260-R2(14)
L2246299-07B	Vial HCl preserved	A	NA		2.7	Y	Absent		NYTCL-8260-R2(14)
L2246299-07C	Vial HCl preserved	A	NA		2.7	Y	Absent		NYTCL-8260-R2(14)
L2246299-08A	Vial HCl preserved	A	NA		2.7	Y	Absent		NYTCL-8260-R2(14)
L2246299-08B	Vial HCl preserved	A	NA		2.7	Y	Absent		NYTCL-8260-R2(14)

Project Name: CONVENTUS
Project Number: U86

Serial_No:09092213:30
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Container Information

Container ID	Container Type	Cooler	Initial pH	Final pH	Temp deg C	Pres	Seal	Frozen Date/Time	Analysis(*)
L2246299-08C	Vial HCl preserved	NA	NA			Y	Absent		NYTCL-8260-R2(14)

Project Name: CONVENTUS**Lab Number:** L2246299**Project Number:** U86**Report Date:** 09/09/22

GLOSSARY

Acronyms

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.)
EDL	- 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.
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.)
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 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 only.)
MDL	- 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.
MS	- 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

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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.

Chlordane: 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.)

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'.

Gasoline Range Organics (GRO): Gasoline Range Organics (GRO) results include all chromatographic peaks eluting from Methyl tert butyl ether through Naphthalene, with the exception of GRO analysis in support of State of Ohio programs, which includes all chromatographic peaks eluting from Hexane through Dodecane.

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.

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.
- B** - 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).
- C** - Co-elution: The target analyte co-elutes with a known lab standard (i.e. surrogate, internal standards, etc.) for co-extracted analyses.
- D** - 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.
- J** - 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

Report Format: DU Report with 'J' Qualifiers



Project Name: CONVENTUS
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Data Qualifiers

Identified Compounds (TICs).

- M** - 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.
- 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.
- V** - The surrogate associated with this target analyte has a recovery outside the QC acceptance limits. (Applicable to MassDEP DW Compliance samples only.)
- Z** - 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: CONVENTUS
Project Number: U86

Lab Number: L2246299
Report Date: 09/09/22

REFERENCES

- 1 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**

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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; 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 I, 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.

 NEW YORK CHAIN OF CUSTODY Westborough, MA 01581 8 Walkup Dr. TEL: 508-898-9220 FAX: 508-898-9193 Mansfield, MA 02048 320 Forbes Blvd TEL: 508-822-9300 FAX: 508-822-3288		Service Centers Mahwah, NJ 07430: 35 Whitney Rd, Suite 5 Albany, NY 12205: 14 Walker Way Tonawanda, NY 14150: 275 Cooper Ave, Suite 105		Page 1 of 1		Date Rec'd in Lab 8/26/22		ALPHA Job # 2246299																																																																																																																																																																																																	
		Project Information Project Name: CONVENTUS Project Location: CONVENTUS / MAIN ST. Project # U86 (Use Project name as Project #) ☐ Project Manager: RICH BAUKERT ALPHAQuote #: Turn-Around Time Standard ☒ Rush (only if pre approved) ☐ Due Date: # of Days:		Deliverables ☐ ASP-A ☒ ASP-B ☒ EQuIS (1 File) ☐ EQuIS (4 File) ☐ Other Regulatory Requirement ☐ NY TOGS ☒ NY Part 375 ☐ AWQ Standards ☐ NY CP-51 ☐ NY Restricted Use ☐ Other ☐ NY Unrestricted Use ☐ NYC Sewer Discharge		Billing Information ☐ Same as Client Info PO # Disposal Site Information Please identify below location of applicable disposal facilities. Disposal Facility: ☐ NJ ☒ NY ☐ Other:																																																																																																																																																																																																			
Client Information Client: C+S ENGINEERS Address: 141 ELM ST BUFFALO NY 14203 Phone: 716 796 3520 Fax: Email: Rbaukert@C805.com		These samples have been previously analyzed by Alpha ☐ Other project specific requirements/comments: Please specify Metals or TAL.		ANALYSIS 8200 VOL		Sample Filtration ☐ Done ☐ Lab to do Preservation ☐ Lab to do (Please Specify below) Sample Specific Comments		Total Bottles																																																																																																																																																																																																	
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DATA USABILITY SUMMARY REPORT (DUSR)

**Conventus
Buffalo, NY
Project # U86001005**

SDG: L2246299
7 Water Samples and 1 Trip Blank

Prepared for:

**C&S Companies
141 Elm Street, Suite 100
Buffalo, NY 14203
Attention: Cody Martin**

November 2022



Environmental Data Usability 10028 Deer Park Dr. Dansville, NY 14437 585-991-9156

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APPENDIX B	Laboratory QC Documentation
APPENDIX C	Validator Qualifications

Tables

Table 4-1	Data Validation Guidance Documents
Table 4-2	Quality Control Criteria for Validating Laboratory Analytical Data

Summaries of Validated Results

Table 6-1	VOCs
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REVIEWER'S NARRATIVE
C&S Companies SDG L2246299 JCC

The data associated with this Sample Delivery Groups (SDG) L2246299, analyzed by Alpha Analytical, Westborough, MA have been reviewed in accordance with assessment criteria provided by the New York State Department of Environmental Conservation following the review procedures provided in the USEPA Functional Guidelines for evaluating organic and inorganic data.

All analytical results reported by the laboratory are considered valid and acceptable except results that have been qualified as rejected, "R". Results qualified as estimated "J", or as non-detects, "U", are considered usable for the purpose of evaluating water and/or soil quality. However, these qualifiers indicate that the accuracy and/or precision of the analytical result is questionable. A summary of all data that have been qualified and the reasons for qualification are provided in the following data usability summary report (DUSR).

Two facts should be noted by all data users. First, the "R" qualifier means that the associated value is unusable. In other words, due to significant quality control (QC) problems, the analysis is invalid and provides no information as to whether the analyte is present or not. Values qualified with an "R" should not appear on the final data tables because they cannot be relied upon, even as the last resort. Second, no analyte concentration, even if it has passed all QC tests, is guaranteed to be accurate. Strict QC serves to increase confidence in data, but any value potentially contains error.

Reviewer's Signature: Michael K. Perry Date: 11/28/2022
Michael K. Perry
Chemist

1.0 EVENT SUMMARY

SITE: Conventus
Buffalo, NY
Project #: U86001005

SAMPLING DATES: August 25, 2022

SAMPLE TYPE: 7 water samples and 1 trip blank

LABORATORY: Alpha Analytical
Westborough, MA

SDG No.: SDGs L2246299

2.0 INTRODUCTION

This data usability summary report (DUSR) was prepared in accordance with guidance provided by the New York State Department of Environmental Conservation (NYSDEC). The DUSR is based on a review and evaluation of the laboratory analytical data package. Specifically, the NYSDEC guidance recommends review and evaluation of the following elements of the data package:

- Completeness of the data package as defined under the requirements of the NYSDEC Analytical Services Protocols (ASP) Category B or the United States Environmental Protection Agency (USEPA) Contract Laboratory Program (CLP) deliverables,
- Compliance with established analyte holding times,
- Adherence to quality control (QC) limits and specifications for blanks, instrument tuning and calibration, surrogate recoveries, spike recoveries, laboratory duplicate analyses, and other QC criteria,
- Adherence to established analytical protocols,
- Conformance of data summary sheets with raw analytical data, and
- Use of correct data qualifiers.

Data deficiencies, analytical protocol deviations, and quality control problems identified using the review criteria above and their effect on the analytical results are discussed in this report.

3.0 SAMPLE AND ANALYSIS SUMMARY

The data package consists of analytical results for 7 water samples and 1 trip blank collected on 8/25/22. These samples were analyzed for Volatile Organic Compounds (VOCs).

All laboratory analyses were submitted to Alpha Analytical, Westborough, MA and analyzed as SDG L2246299. The analytical results were provided in NYSDEC ASP Category B format, which includes all raw analytical data and laboratory QC data.

4.0 GUIDANCE DOCUMENTS AND DATA REVIEW CRITERIA

The guidance documents appropriate for reviewing laboratory quality control (QC) data and assigning data qualifiers (flags) to analytical results were selected from those listed in Table 4-1. The QC limits established in the documents applicable to this data review were used to assess the quality of the analytical results. In some cases, however, QC limits established internally by the laboratory were taken into account to determine data quality.

The QC criteria considered for assessing the usability of the reported analytical results provided for each analyte type (i.e. VOCs, SVOCs, metals, etc.) are listed in Table 4-2. These criteria may vary with the analytical method utilized by the laboratory. These criteria comply with the guidance recommended in Section 2.0 above.

5.0 DATA VALIDATION QUALIFIERS

The letter qualifiers (flags) used to define data usability are described briefly below. These letters are assigned by the data validator to analytical results having questionable accuracy and/or precision as determined by reviewing the laboratory QC data associated with the analytical results.

TABLE 4-1**Guidance Used For Validating Laboratory Analytical Data**

Analyte Group	Guidance	Date
Metals (ICP-AES)	USEPA SOP HW-3a, Rev. 1	September 2016
Metals (Hg & CN)	USEPA SOP HW-3c, Rev. 1	September 2016
Volatile Organic Compounds (by Methods 8260B & 8260C)	USEPA SOP HW-24, Rev. 4	September 2014
Semi-Volatile Organic Compounds (by Method 8270D)	USEPA SOP HW-22 Rev. 5	December 2010
Pesticides (by Method 8181B)	USEPA SOP HW-44, Rev. 1.1	December 2010
Chlorinated Herbicides (by Method 8151A)	USEPA SOP HW-17, Rev. 3.1	December 2010
Polychlorinated Biphenyls (PCBs)	USEPA SOP HW-37A, Rev. 0	June 2015
Volatile Organic Compounds (Air) (by Method TO-15)	USEPA SOP HW-31, Rev. 6	September 2016
Per- and PolyFluoroAlkyl Substances (PFAS)	* NYSDEC	January 2021
General Chemistry Parameters	per NYSDEC ASP	July 2005

* Sampling, Analysis, and Assessment of Per- and Polyfluoroalkyl Substances (PFAS) Under NYSDEC's Part 375 Remedial Programs, Appendix I

TABLE 4-2

**QUALITY CONTROL CRITERIA USED FOR VALIDATING
LABORATORY ANALYTICAL DATA**

VOCs	SVOCs	Pesticides/PCBs	Metals	Gen Chemistry	PFAS
Completeness of Pkg Sample Preservation Holding Time System Monitoring Compounds Lab Control Sample Matrix Spikes Blanks Instrument Tuning Internal Standards Initial Calibration Continuing Calibration Lab Qualifiers Field Duplicate	Completeness of Pkg Sample Preservation Holding Time Surrogate Recoveries Lab Control Sample Matrix Spikes Blanks Instrument Tuning Internal Standards Initial Calibration Continuing Calibration Lab Qualifiers Field Duplicate	Completeness of Pkg Sample Preservation Holding Time Surrogate Recoveries Matrix Spikes Blanks Instrument Calibration & Verification Comparison of duplicate GC column results Analyte ID Lab Qualifiers Field Duplicate	Completeness of Pkg Sample Preservation Holding Time Initial/Continuing Calibration CRDL Standards Blanks Interference Check Sample Spike Recoveries Lab Duplicate Lab Control Sample ICP Serial Dilutions Lab Qualifiers Field Duplicate	Completeness of Pkg Sample Preservation Holding Times Calibration Lab Control Samples Blanks Spike Recoveries Lab Duplicates	Completeness of Pkg Sample Preservation Holding Time Instr Performance Check Initial Calibration Continuing Calibration Blanks Surrogates Lab Fortified Blank Matrix Spikes Internal Standards

Method TO-15 (Air)

Completeness of Pkg
 Sample Preservation
 Holding Time
 Canister Certification
 Instrument Tuning
 Initial Calibration and
 Instrument Performance
 Daily Calibration
 Blanks
 Lab Control Sample
 Field Duplicate

The laboratory may also use various letters and symbols to flag analytical results generated when QC limits were exceeded. The meanings of these flags may differ from those used by the independent data validator. Those used by the laboratory are provided with the analytical results.

NOTE: The assignment of data qualifiers by the data reviewer (validator) to laboratory analytical results should not necessarily be interpreted by the data user as a measure of laboratory ability or proficiency. Rather, the qualifiers are intended to provide a measure of data accuracy and precision to the data user, which, for example, may provide a level of confidence in determining whether or not standards or cleanup objectives have been met.

- U** The analyte was analyzed for but was not detected at or above the sample quantitation limit.
- J** The analyte was positively identified; the associated numerical value is the *approximate* concentration of the analyte in the sample. (The magnitude of any $\pm$ value associated with the result is not determined by data validation).
- J+** The result is an estimated quantity and may be biased high.
- J-** The result is an estimated quantity and may be biased low.
- UJ** The analyte was analyzed for but not detected. The reported quantitation limit is approximate and may be inaccurate or imprecise.
- R** The sample result is rejected (i.e., is unusable) due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.
- NJ** The analysis indicates the presence of an analyte that has been “tentatively identified” and the associated numerical value represents its approximate concentration.

The validated analytical results are attached to this report. Validation qualifiers (flags) are indicated in red print. Data sheets having qualified data are signed and dated by the data reviewer.

6.0 RESULTS OF THE DATA REVIEW

The results of the data review are summarized in Table 6-1. The table lists the samples where QC criteria were found to exceed acceptable limits and the actions taken to qualify the associated analytical results.

7.0 TOTAL USABLE DATA

For SDG L2246299, eight samples were analyzed and results were reported for 464 analytes. Even though some results were flagged with a “J” as estimated, all results (100 %) are considered usable. See the summary table for the analyses that have been rejected and qualified and the associated QC reasons.

Table 6-1 VOCs

SAMPLES AFFECTED	ANALYTES	ACTION	QC VIOLATION	COMMENTS
BCP MW-1 BCP MW-7 BCP MW-4 BCP MW-3 BCP MW-6 BCP MW-5 DUP Trip Blank	Chloroethane Trichlorofluoromethane Methyl acetate 2-Butanone 4-Methyl-2-pentanone 2-Hexanone DBCP	J detects UJ non-detects	CCV > QC limit	Data are estimated

ACRONYMS

BSP	Blank Spike
CCAL	Continuing Calibration
CCB	Continuing Calibration Blank
CCV	Continuing Calibration Verification
CRDL	Contract Required Detection Limit
CRQL	Contract Required Quantitation Limit
%D	Percent Difference
ICAL	Initial Calibration
ICB	Initial Calibration Blank
IS	Internal Standard
LCS	Laboratory Control Sample
MS/MSD	Matrix Spike/Matrix Spike Duplicate
QA	Quality Assurance
QC	Quality Control
%R	Percent recovery
RPD	Relative Percent Difference
RRF	Relative Response Factor
%RSD	Percent Relative Standard Deviation
TAL	Target Analyte List (metals)
TCL	Target Compound List (organics)

Appendix A

Validated Analytical Results



www.alphalab.com



Alpha Analytical

Laboratory Code: 11148

SDG Number: L2246299

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.

Project Name: CONVENTUS
Project Number: U86

Lab Number: L2246299
Report Date: 09/09/22

Alpha Sample ID	Client ID	Matrix	Sample Location	Collection Date/Time	Receive Date
L2246299-01	BCP MW-1	WATER	CONVENTUS/MAIN ST	08/25/22 09:55	08/25/22
L2246299-02	BCP MW-7	WATER	CONVENTUS/MAIN ST	08/25/22 10:25	08/25/22
L2246299-03	BCP MW-4	WATER	CONVENTUS/MAIN ST	08/25/22 10:55	08/25/22
L2246299-04	BCP MW-3	WATER	CONVENTUS/MAIN ST	08/25/22 11:35	08/25/22
L2246299-05	BCP MW-6	WATER	CONVENTUS/MAIN ST	08/25/22 12:15	08/25/22
L2246299-06	BCP MW-5	WATER	CONVENTUS/MAIN ST	08/25/22 13:00	08/25/22
L2246299-07	DUP	WATER	CONVENTUS/MAIN ST	08/25/22 09:55	08/25/22
L2246299-08	TRIP BLANK	WATER	CONVENTUS/MAIN ST	08/25/22 13:50	08/25/22

Project Name: CONVENTUS
Project Number: U86

Lab Number: L2246299
Report Date: 09/09/22

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.

Volatile Organics

L2246299-05: The pH was greater than two; however, the sample was analyzed within the method required holding time.

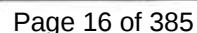
I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis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

Authorized Signature: *Jiffani Morrissey*

Report Date: 09/09/22

Title: Technical Director/Representative





GC/MS 8260

Analysis

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies
 Project Name : CONVENTUS
 Lab ID : L2246299-01
 Client ID : BCP MW-1
 Sample Location : CONVENTUS/MAIN ST
 Sample Matrix : WATER
 Analytical Method : 1,8260C
 Lab File ID : V05220901N07
 Sample Amount : 10 ml
 Level : LOW
 Extract Volume (MeOH) : N/A

Lab Number : L2246299
 Project Number : U86
 Date Collected : 08/25/22 09:55
 Date Received : 08/25/22
 Date Analyzed : 09/01/22 20:47
 Dilution Factor : 1
 Analyst : MV
 Instrument ID : VOA105
 GC Column : RTX-502.2
 %Solids : N/A
 Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
75-09-2	Methylene chloride	ND	2.5	0.70	U
75-34-3	1,1-Dichloroethane	ND	2.5	0.70	U
67-66-3	Chloroform	ND	2.5	0.70	U
56-23-5	Carbon tetrachloride	ND	0.50	0.13	U
78-87-5	1,2-Dichloropropane	ND	1.0	0.14	U
124-48-1	Dibromochloromethane	ND	0.50	0.15	U
79-00-5	1,1,2-Trichloroethane	ND	1.5	0.50	U
127-18-4	Tetrachloroethene	ND	0.50	0.18	U
108-90-7	Chlorobenzene	ND	2.5	0.70	U
75-69-4	Trichlorofluoromethane	ND	2.5	0.70	U
107-06-2	1,2-Dichloroethane	ND	0.50	0.13	U UJ
71-55-6	1,1,1-Trichloroethane	ND	2.5	0.70	U
75-27-4	Bromodichloromethane	ND	0.50	0.19	U
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.16	U
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.14	U
75-25-2	Bromoform	ND	2.0	0.65	U
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.17	U
71-43-2	Benzene	ND	0.50	0.16	U
108-88-3	Toluene	ND	2.5	0.70	U
100-41-4	Ethylbenzene	ND	2.5	0.70	U
74-87-3	Chloromethane	ND	2.5	0.70	U
74-83-9	Bromomethane	ND	2.5	0.70	U
75-01-4	Vinyl chloride	ND	1.0	0.07	U
75-00-3	Chloroethane	ND	2.5	0.70	U
75-35-4	1,1-Dichloroethene	ND	0.50	0.17	U UJ



Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies
 Project Name : CONVENTUS
 Lab ID : L2246299-01
 Client ID : BCP MW-1
 Sample Location : CONVENTUS/MAIN ST
 Sample Matrix : WATER
 Analytical Method : 1,8260C
 Lab File ID : V05220901N07
 Sample Amount : 10 ml
 Level : LOW
 Extract Volume (MeOH) : N/A

Lab Number : L2246299
 Project Number : U86
 Date Collected : 08/25/22 09:55
 Date Received : 08/25/22
 Date Analyzed : 09/01/22 20:47
 Dilution Factor : 1
 Analyst : MV
 Instrument ID : VOA105
 GC Column : RTX-502.2
 %Solids : N/A
 Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
156-60-5	trans-1,2-Dichloroethene	ND	2.5	0.70	U
79-01-6	Trichloroethene	ND	0.50	0.18	U
95-50-1	1,2-Dichlorobenzene	ND	2.5	0.70	U
541-73-1	1,3-Dichlorobenzene	ND	2.5	0.70	U
106-46-7	1,4-Dichlorobenzene	ND	2.5	0.70	U
1634-04-4	Methyl tert butyl ether	ND	2.5	0.70	U
179601-23-1	p/m-Xylene	ND	2.5	0.70	U
95-47-6	o-Xylene	ND	2.5	0.70	U
156-59-2	cis-1,2-Dichloroethene	ND	2.5	0.70	U
100-42-5	Styrene	ND	2.5	0.70	U
75-71-8	Dichlorodifluoromethane	ND	5.0	1.0	U
67-64-1	Acetone	ND	5.0	1.5	U
75-15-0	Carbon disulfide	ND	5.0	1.0	U
78-93-3	2-Butanone	ND	5.0	1.9	U
108-10-1	4-Methyl-2-pentanone	ND	5.0	1.0	U UJ
591-78-6	2-Hexanone	ND	5.0	1.0	U UJ
106-93-4	1,2-Dibromoethane	ND	2.0	0.65	U UJ
104-51-8	n-Butylbenzene	ND	2.5	0.70	U
135-98-8	sec-Butylbenzene	ND	2.5	0.70	U
98-06-6	tert-Butylbenzene	ND	2.5	0.70	U
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.5	0.70	U
98-82-8	Isopropylbenzene	ND	2.5	0.70	U UJ
99-87-6	p-Isopropyltoluene	ND	2.5	0.70	U
91-20-3	Naphthalene	ND	2.5	0.70	U
103-65-1	n-Propylbenzene	ND	2.5	0.70	U



Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies
 Project Name : CONVENTUS
 Lab ID : L2246299-01
 Client ID : BCP MW-1
 Sample Location : CONVENTUS/MAIN ST
 Sample Matrix : WATER
 Analytical Method : 1,8260C
 Lab File ID : V05220901N07
 Sample Amount : 10 ml
 Level : LOW
 Extract Volume (MeOH) : N/A

Lab Number : L2246299
 Project Number : U86
 Date Collected : 08/25/22 09:55
 Date Received : 08/25/22
 Date Analyzed : 09/01/22 20:47
 Dilution Factor : 1
 Analyst : MV
 Instrument ID : VOA105
 GC Column : RTX-502.2
 %Solids : N/A
 Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
120-82-1	1,2,4-Trichlorobenzene	ND	2.5	0.70	U
108-67-8	1,3,5-Trimethylbenzene	ND	2.5	0.70	U
95-63-6	1,2,4-Trimethylbenzene	ND	2.5	0.70	U
79-20-9	Methyl Acetate	ND	2.0	0.23	U
110-82-7	Cyclohexane	ND	10	0.27	U UJ
76-13-1	Freon-113	ND	2.5	0.70	U
108-87-2	Methyl cyclohexane	ND	10	0.40	U

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Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies
 Project Name : CONVENTUS
 Lab ID : L2246299-02
 Client ID : BCP MW-7
 Sample Location : CONVENTUS/MAIN ST
 Sample Matrix : WATER
 Analytical Method : 1,8260C
 Lab File ID : V05220901N08
 Sample Amount : 10 ml
 Level : LOW
 Extract Volume (MeOH) : N/A

Lab Number : L2246299
 Project Number : U86
 Date Collected : 08/25/22 10:25
 Date Received : 08/25/22
 Date Analyzed : 09/01/22 21:10
 Dilution Factor : 1
 Analyst : MV
 Instrument ID : VOA105
 GC Column : RTX-502.2
 %Solids : N/A
 Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
75-09-2	Methylene chloride	ND	2.5	0.70	U
75-34-3	1,1-Dichloroethane	ND	2.5	0.70	U
67-66-3	Chloroform	ND	2.5	0.70	U
56-23-5	Carbon tetrachloride	ND	0.50	0.13	U
78-87-5	1,2-Dichloropropane	ND	1.0	0.14	U
124-48-1	Dibromochloromethane	ND	0.50	0.15	U
79-00-5	1,1,2-Trichloroethane	ND	1.5	0.50	U
127-18-4	Tetrachloroethene	ND	0.50	0.18	U
108-90-7	Chlorobenzene	ND	2.5	0.70	U
75-69-4	Trichlorofluoromethane	ND	2.5	0.70	U
107-06-2	1,2-Dichloroethane	ND	0.50	0.13	U UJ
71-55-6	1,1,1-Trichloroethane	ND	2.5	0.70	U
75-27-4	Bromodichloromethane	ND	0.50	0.19	U
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.16	U
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.14	U
75-25-2	Bromoform	ND	2.0	0.65	U
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.17	U
71-43-2	Benzene	ND	0.50	0.16	U
108-88-3	Toluene	ND	2.5	0.70	U
100-41-4	Ethylbenzene	ND	2.5	0.70	U
74-87-3	Chloromethane	ND	2.5	0.70	U
74-83-9	Bromomethane	ND	2.5	0.70	U
75-01-4	Vinyl chloride	ND	1.0	0.07	U
75-00-3	Chloroethane	ND	2.5	0.70	U
75-35-4	1,1-Dichloroethene	ND	0.50	0.17	U UJ



Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies
 Project Name : CONVENTUS
 Lab ID : L2246299-02
 Client ID : BCP MW-7
 Sample Location : CONVENTUS/MAIN ST
 Sample Matrix : WATER
 Analytical Method : 1,8260C
 Lab File ID : V05220901N08
 Sample Amount : 10 ml
 Level : LOW
 Extract Volume (MeOH) : N/A

Lab Number : L2246299
 Project Number : U86
 Date Collected : 08/25/22 10:25
 Date Received : 08/25/22
 Date Analyzed : 09/01/22 21:10
 Dilution Factor : 1
 Analyst : MV
 Instrument ID : VOA105
 GC Column : RTX-502.2
 %Solids : N/A
 Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
156-60-5	trans-1,2-Dichloroethene	ND	2.5	0.70	U
79-01-6	Trichloroethene	ND	0.50	0.18	U
95-50-1	1,2-Dichlorobenzene	ND	2.5	0.70	U
541-73-1	1,3-Dichlorobenzene	ND	2.5	0.70	U
106-46-7	1,4-Dichlorobenzene	ND	2.5	0.70	U
1634-04-4	Methyl tert butyl ether	ND	2.5	0.70	U
179601-23-1	p/m-Xylene	ND	2.5	0.70	U
95-47-6	o-Xylene	ND	2.5	0.70	U
156-59-2	cis-1,2-Dichloroethene	ND	2.5	0.70	U
100-42-5	Styrene	ND	2.5	0.70	U
75-71-8	Dichlorodifluoromethane	ND	5.0	1.0	U
67-64-1	Acetone	ND	5.0	1.5	U
75-15-0	Carbon disulfide	ND	5.0	1.0	U
78-93-3	2-Butanone	ND	5.0	1.9	U
108-10-1	4-Methyl-2-pentanone	ND	5.0	1.0	U UJ
591-78-6	2-Hexanone	ND	5.0	1.0	U UJ
106-93-4	1,2-Dibromoethane	ND	2.0	0.65	U UJ
104-51-8	n-Butylbenzene	ND	2.5	0.70	U
135-98-8	sec-Butylbenzene	ND	2.5	0.70	U
98-06-6	tert-Butylbenzene	ND	2.5	0.70	U
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.5	0.70	U
98-82-8	Isopropylbenzene	ND	2.5	0.70	U UJ
99-87-6	p-Isopropyltoluene	ND	2.5	0.70	U
91-20-3	Naphthalene	ND	2.5	0.70	U
103-65-1	n-Propylbenzene	ND	2.5	0.70	U

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies
 Project Name : CONVENTUS
 Lab ID : L2246299-02
 Client ID : BCP MW-7
 Sample Location : CONVENTUS/MAIN ST
 Sample Matrix : WATER
 Analytical Method : 1,8260C
 Lab File ID : V05220901N08
 Sample Amount : 10 ml
 Level : LOW
 Extract Volume (MeOH) : N/A

Lab Number : L2246299
 Project Number : U86
 Date Collected : 08/25/22 10:25
 Date Received : 08/25/22
 Date Analyzed : 09/01/22 21:10
 Dilution Factor : 1
 Analyst : MV
 Instrument ID : VOA105
 GC Column : RTX-502.2
 %Solids : N/A
 Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
120-82-1	1,2,4-Trichlorobenzene	ND	2.5	0.70	U
108-67-8	1,3,5-Trimethylbenzene	ND	2.5	0.70	U
95-63-6	1,2,4-Trimethylbenzene	ND	2.5	0.70	U
79-20-9	Methyl Acetate	ND	2.0	0.23	U
110-82-7	Cyclohexane	ND	10	0.27	U UJ
76-13-1	Freon-113	ND	2.5	0.70	U
108-87-2	Methyl cyclohexane	ND	10	0.40	U

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Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies
 Project Name : CONVENTUS
 Lab ID : L2246299-03D
 Client ID : BCP MW-4
 Sample Location : CONVENTUS/MAIN ST
 Sample Matrix : WATER
 Analytical Method : 1,8260C
 Lab File ID : V05220901N23
 Sample Amount : 1 ml
 Level : LOW
 Extract Volume (MeOH) : N/A

Lab Number : L2246299
 Project Number : U86
 Date Collected : 08/25/22 10:55
 Date Received : 08/25/22
 Date Analyzed : 09/02/22 03:00
 Dilution Factor : 10
 Analyst : MV
 Instrument ID : VOA105
 GC Column : RTX-502.2
 %Solids : N/A
 Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
75-09-2	Methylene chloride	ND	25	7.0	U
75-34-3	1,1-Dichloroethane	ND	25	7.0	U
67-66-3	Chloroform	ND	25	7.0	U
56-23-5	Carbon tetrachloride	ND	5.0	1.3	U
78-87-5	1,2-Dichloropropane	ND	10	1.4	U
124-48-1	Dibromochloromethane	ND	5.0	1.5	U
79-00-5	1,1,2-Trichloroethane	ND	15	5.0	U
127-18-4	Tetrachloroethene	ND	5.0	1.8	U
108-90-7	Chlorobenzene	ND	25	7.0	U
75-69-4	Trichlorofluoromethane	ND	25	7.0	U
107-06-2	1,2-Dichloroethane	ND	5.0	1.3	U UJ
71-55-6	1,1,1-Trichloroethane	ND	25	7.0	U
75-27-4	Bromodichloromethane	ND	5.0	1.9	U
10061-02-6	trans-1,3-Dichloropropene	ND	5.0	1.6	U
10061-01-5	cis-1,3-Dichloropropene	ND	5.0	1.4	U
75-25-2	Bromoform	ND	20	6.5	U
79-34-5	1,1,2,2-Tetrachloroethane	ND	5.0	1.7	U
71-43-2	Benzene	5.0	5.0	1.6	
108-88-3	Toluene	34	25	7.0	
100-41-4	Ethylbenzene	1100	25	7.0	
74-87-3	Chloromethane	ND	25	7.0	U
74-83-9	Bromomethane	ND	25	7.0	U
75-01-4	Vinyl chloride	ND	10	0.71	U
75-00-3	Chloroethane	ND	25	7.0	U
75-35-4	1,1-Dichloroethene	ND	5.0	1.7	U UJ

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies
 Project Name : CONVENTUS
 Lab ID : L2246299-03D
 Client ID : BCP MW-4
 Sample Location : CONVENTUS/MAIN ST
 Sample Matrix : WATER
 Analytical Method : 1,8260C
 Lab File ID : V05220901N23
 Sample Amount : 1 ml
 Level : LOW
 Extract Volume (MeOH) : N/A

Lab Number : L2246299
 Project Number : U86
 Date Collected : 08/25/22 10:55
 Date Received : 08/25/22
 Date Analyzed : 09/02/22 03:00
 Dilution Factor : 10
 Analyst : MV
 Instrument ID : VOA105
 GC Column : RTX-502.2
 %Solids : N/A
 Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
156-60-5	trans-1,2-Dichloroethene	ND	25	7.0	U
79-01-6	Trichloroethene	ND	5.0	1.8	U
95-50-1	1,2-Dichlorobenzene	ND	25	7.0	U
541-73-1	1,3-Dichlorobenzene	ND	25	7.0	U
106-46-7	1,4-Dichlorobenzene	ND	25	7.0	U
1634-04-4	Methyl tert butyl ether	ND	25	7.0	U
179601-23-1	p/m-Xylene	14	25	7.0	J
95-47-6	o-Xylene	12	25	7.0	J
156-59-2	cis-1,2-Dichloroethene	ND	25	7.0	U
100-42-5	Styrene	ND	25	7.0	U
75-71-8	Dichlorodifluoromethane	ND	50	10.	U
67-64-1	Acetone	ND	50	15.	U
75-15-0	Carbon disulfide	ND	50	10.	U
78-93-3	2-Butanone	ND	50	19.	U
108-10-1	4-Methyl-2-pentanone	ND	50	10.	U UJ
591-78-6	2-Hexanone	ND	50	10.	U UJ
106-93-4	1,2-Dibromoethane	ND	20	6.5	U UJ
104-51-8	n-Butylbenzene	9.4	25	7.0	J
135-98-8	sec-Butylbenzene	ND	25	7.0	U
98-06-6	tert-Butylbenzene	ND	25	7.0	U
96-12-8	1,2-Dibromo-3-chloropropane	ND	25	7.0	U
98-82-8	Isopropylbenzene	24	25	7.0	J UJ
99-87-6	p-Isopropyltoluene	ND	25	7.0	U
91-20-3	Naphthalene	200	25	7.0	
103-65-1	n-Propylbenzene	140	25	7.0	

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies
 Project Name : CONVENTUS
 Lab ID : L2246299-03D
 Client ID : BCP MW-4
 Sample Location : CONVENTUS/MAIN ST
 Sample Matrix : WATER
 Analytical Method : 1,8260C
 Lab File ID : V05220901N23
 Sample Amount : 1 ml
 Level : LOW
 Extract Volume (MeOH) : N/A

Lab Number : L2246299
 Project Number : U86
 Date Collected : 08/25/22 10:55
 Date Received : 08/25/22
 Date Analyzed : 09/02/22 03:00
 Dilution Factor : 10
 Analyst : MV
 Instrument ID : VOA105
 GC Column : RTX-502.2
 %Solids : N/A
 Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
120-82-1	1,2,4-Trichlorobenzene	ND	25	7.0	U
108-67-8	1,3,5-Trimethylbenzene	ND	25	7.0	U
95-63-6	1,2,4-Trimethylbenzene	1200	25	7.0	
79-20-9	Methyl Acetate	ND	20	2.3	U
110-82-7	Cyclohexane	130	100	2.7	UJ
76-13-1	Freon-113	ND	25	7.0	U
108-87-2	Methyl cyclohexane	51	100	4.0	J

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Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies
 Project Name : CONVENTUS
 Lab ID : L2246299-04D
 Client ID : BCP MW-3
 Sample Location : CONVENTUS/MAIN ST
 Sample Matrix : WATER
 Analytical Method : 1,8260C
 Lab File ID : V05220901N24
 Sample Amount : 2 ml
 Level : LOW
 Extract Volume (MeOH) : N/A

Lab Number : L2246299
 Project Number : U86
 Date Collected : 08/25/22 11:35
 Date Received : 08/25/22
 Date Analyzed : 09/02/22 03:23
 Dilution Factor : 5
 Analyst : MV
 Instrument ID : VOA105
 GC Column : RTX-502.2
 %Solids : N/A
 Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
75-09-2	Methylene chloride	ND	12	3.5	U
75-34-3	1,1-Dichloroethane	ND	12	3.5	U
67-66-3	Chloroform	ND	12	3.5	U
56-23-5	Carbon tetrachloride	ND	2.5	0.67	U
78-87-5	1,2-Dichloropropane	ND	5.0	0.68	U
124-48-1	Dibromochloromethane	ND	2.5	0.74	U
79-00-5	1,1,2-Trichloroethane	ND	7.5	2.5	U
127-18-4	Tetrachloroethene	ND	2.5	0.90	U
108-90-7	Chlorobenzene	ND	12	3.5	U
75-69-4	Trichlorofluoromethane	ND	12	3.5	U
107-06-2	1,2-Dichloroethane	ND	2.5	0.66	U UJ
71-55-6	1,1,1-Trichloroethane	ND	12	3.5	U
75-27-4	Bromodichloromethane	ND	2.5	0.96	U
10061-02-6	trans-1,3-Dichloropropene	ND	2.5	0.82	U
10061-01-5	cis-1,3-Dichloropropene	ND	2.5	0.72	U
75-25-2	Bromoform	ND	10	3.2	U
79-34-5	1,1,2,2-Tetrachloroethane	ND	2.5	0.84	U
71-43-2	Benzene	1.7	2.5	0.80	J
108-88-3	Toluene	22	12	3.5	
100-41-4	Ethylbenzene	380	12	3.5	
74-87-3	Chloromethane	ND	12	3.5	U
74-83-9	Bromomethane	ND	12	3.5	U
75-01-4	Vinyl chloride	ND	5.0	0.36	U
75-00-3	Chloroethane	ND	12	3.5	U
75-35-4	1,1-Dichloroethene	ND	2.5	0.84	U UJ

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies
 Project Name : CONVENTUS
 Lab ID : L2246299-04D
 Client ID : BCP MW-3
 Sample Location : CONVENTUS/MAIN ST
 Sample Matrix : WATER
 Analytical Method : 1,8260C
 Lab File ID : V05220901N24
 Sample Amount : 2 ml
 Level : LOW
 Extract Volume (MeOH) : N/A

Lab Number : L2246299
 Project Number : U86
 Date Collected : 08/25/22 11:35
 Date Received : 08/25/22
 Date Analyzed : 09/02/22 03:23
 Dilution Factor : 5
 Analyst : MV
 Instrument ID : VOA105
 GC Column : RTX-502.2
 %Solids : N/A
 Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
156-60-5	trans-1,2-Dichloroethene	ND	12	3.5	U
79-01-6	Trichloroethene	ND	2.5	0.88	U
95-50-1	1,2-Dichlorobenzene	ND	12	3.5	U
541-73-1	1,3-Dichlorobenzene	ND	12	3.5	U
106-46-7	1,4-Dichlorobenzene	ND	12	3.5	U
1634-04-4	Methyl tert butyl ether	ND	12	3.5	U
179601-23-1	p/m-Xylene	480	12	3.5	
95-47-6	o-Xylene	12	12	3.5	
156-59-2	cis-1,2-Dichloroethene	ND	12	3.5	U
100-42-5	Styrene	ND	12	3.5	U
75-71-8	Dichlorodifluoromethane	ND	25	5.0	U
67-64-1	Acetone	ND	25	7.3	U
75-15-0	Carbon disulfide	ND	25	5.0	U
78-93-3	2-Butanone	ND	25	9.7	U
108-10-1	4-Methyl-2-pentanone	ND	25	5.0	U UJ
591-78-6	2-Hexanone	ND	25	5.0	U UJ
106-93-4	1,2-Dibromoethane	ND	10	3.2	U UJ
104-51-8	n-Butylbenzene	ND	12	3.5	U
135-98-8	sec-Butylbenzene	ND	12	3.5	U
98-06-6	tert-Butylbenzene	ND	12	3.5	U
96-12-8	1,2-Dibromo-3-chloropropane	ND	12	3.5	U
98-82-8	Isopropylbenzene	5.2	12	3.5	J UJ
99-87-6	p-Isopropyltoluene	ND	12	3.5	U
91-20-3	Naphthalene	170	12	3.5	
103-65-1	n-Propylbenzene	18	12	3.5	



Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies
 Project Name : CONVENTUS
 Lab ID : L2246299-04D
 Client ID : BCP MW-3
 Sample Location : CONVENTUS/MAIN ST
 Sample Matrix : WATER
 Analytical Method : 1,8260C
 Lab File ID : V05220901N24
 Sample Amount : 2 ml
 Level : LOW
 Extract Volume (MeOH) : N/A

Lab Number : L2246299
 Project Number : U86
 Date Collected : 08/25/22 11:35
 Date Received : 08/25/22
 Date Analyzed : 09/02/22 03:23
 Dilution Factor : 5
 Analyst : MV
 Instrument ID : VOA105
 GC Column : RTX-502.2
 %Solids : N/A
 Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
120-82-1	1,2,4-Trichlorobenzene	ND	12	3.5	U
108-67-8	1,3,5-Trimethylbenzene	47	12	3.5	
95-63-6	1,2,4-Trimethylbenzene	560	12	3.5	
79-20-9	Methyl Acetate	ND	10	1.2	U
110-82-7	Cyclohexane	40	50	1.4	J UJ
76-13-1	Freon-113	ND	12	3.5	U
108-87-2	Methyl cyclohexane	32	50	2.0	J

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Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies
 Project Name : CONVENTUS
 Lab ID : L2246299-05
 Client ID : BCP MW-6
 Sample Location : CONVENTUS/MAIN ST
 Sample Matrix : WATER
 Analytical Method : 1,8260C
 Lab File ID : V05220901N09
 Sample Amount : 10 ml
 Level : LOW
 Extract Volume (MeOH) : N/A

Lab Number : L2246299
 Project Number : U86
 Date Collected : 08/25/22 12:15
 Date Received : 08/25/22
 Date Analyzed : 09/01/22 21:33
 Dilution Factor : 1
 Analyst : MV
 Instrument ID : VOA105
 GC Column : RTX-502.2
 %Solids : N/A
 Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
75-09-2	Methylene chloride	ND	2.5	0.70	U
75-34-3	1,1-Dichloroethane	ND	2.5	0.70	U
67-66-3	Chloroform	ND	2.5	0.70	U
56-23-5	Carbon tetrachloride	ND	0.50	0.13	U
78-87-5	1,2-Dichloropropane	ND	1.0	0.14	U
124-48-1	Dibromochloromethane	ND	0.50	0.15	U
79-00-5	1,1,2-Trichloroethane	ND	1.5	0.50	U
127-18-4	Tetrachloroethene	ND	0.50	0.18	U
108-90-7	Chlorobenzene	ND	2.5	0.70	U
75-69-4	Trichlorofluoromethane	ND	2.5	0.70	U
107-06-2	1,2-Dichloroethane	ND	0.50	0.13	U UJ
71-55-6	1,1,1-Trichloroethane	ND	2.5	0.70	U
75-27-4	Bromodichloromethane	ND	0.50	0.19	U
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.16	U
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.14	U
75-25-2	Bromoform	ND	2.0	0.65	U
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.17	U
71-43-2	Benzene	ND	0.50	0.16	U
108-88-3	Toluene	ND	2.5	0.70	U
100-41-4	Ethylbenzene	ND	2.5	0.70	U
74-87-3	Chloromethane	ND	2.5	0.70	U
74-83-9	Bromomethane	ND	2.5	0.70	U
75-01-4	Vinyl chloride	ND	1.0	0.07	U
75-00-3	Chloroethane	ND	2.5	0.70	U
75-35-4	1,1-Dichloroethene	ND	0.50	0.17	U UJ

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies
 Project Name : CONVENTUS
 Lab ID : L2246299-05
 Client ID : BCP MW-6
 Sample Location : CONVENTUS/MAIN ST
 Sample Matrix : WATER
 Analytical Method : 1,8260C
 Lab File ID : V05220901N09
 Sample Amount : 10 ml
 Level : LOW
 Extract Volume (MeOH) : N/A

Lab Number : L2246299
 Project Number : U86
 Date Collected : 08/25/22 12:15
 Date Received : 08/25/22
 Date Analyzed : 09/01/22 21:33
 Dilution Factor : 1
 Analyst : MV
 Instrument ID : VOA105
 GC Column : RTX-502.2
 %Solids : N/A
 Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
156-60-5	trans-1,2-Dichloroethene	ND	2.5	0.70	U
79-01-6	Trichloroethene	ND	0.50	0.18	U
95-50-1	1,2-Dichlorobenzene	ND	2.5	0.70	U
541-73-1	1,3-Dichlorobenzene	ND	2.5	0.70	U
106-46-7	1,4-Dichlorobenzene	ND	2.5	0.70	U
1634-04-4	Methyl tert butyl ether	ND	2.5	0.70	U
179601-23-1	p/m-Xylene	ND	2.5	0.70	U
95-47-6	o-Xylene	ND	2.5	0.70	U
156-59-2	cis-1,2-Dichloroethene	ND	2.5	0.70	U
100-42-5	Styrene	ND	2.5	0.70	U
75-71-8	Dichlorodifluoromethane	ND	5.0	1.0	U
67-64-1	Acetone	ND	5.0	1.5	U
75-15-0	Carbon disulfide	ND	5.0	1.0	U
78-93-3	2-Butanone	ND	5.0	1.9	U
108-10-1	4-Methyl-2-pentanone	ND	5.0	1.0	U UJ
591-78-6	2-Hexanone	ND	5.0	1.0	U UJ
106-93-4	1,2-Dibromoethane	ND	2.0	0.65	U UJ
104-51-8	n-Butylbenzene	ND	2.5	0.70	U
135-98-8	sec-Butylbenzene	ND	2.5	0.70	U
98-06-6	tert-Butylbenzene	ND	2.5	0.70	U
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.5	0.70	U
98-82-8	Isopropylbenzene	ND	2.5	0.70	U UJ
99-87-6	p-Isopropyltoluene	ND	2.5	0.70	U
91-20-3	Naphthalene	ND	2.5	0.70	U
103-65-1	n-Propylbenzene	ND	2.5	0.70	U

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies
 Project Name : CONVENTUS
 Lab ID : L2246299-05
 Client ID : BCP MW-6
 Sample Location : CONVENTUS/MAIN ST
 Sample Matrix : WATER
 Analytical Method : 1,8260C
 Lab File ID : V05220901N09
 Sample Amount : 10 ml
 Level : LOW
 Extract Volume (MeOH) : N/A

Lab Number : L2246299
 Project Number : U86
 Date Collected : 08/25/22 12:15
 Date Received : 08/25/22
 Date Analyzed : 09/01/22 21:33
 Dilution Factor : 1
 Analyst : MV
 Instrument ID : VOA105
 GC Column : RTX-502.2
 %Solids : N/A
 Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
120-82-1	1,2,4-Trichlorobenzene	ND	2.5	0.70	U
108-67-8	1,3,5-Trimethylbenzene	ND	2.5	0.70	U
95-63-6	1,2,4-Trimethylbenzene	ND	2.5	0.70	U
79-20-9	Methyl Acetate	ND	2.0	0.23	U
110-82-7	Cyclohexane	ND	10	0.27	U UJ
76-13-1	Freon-113	ND	2.5	0.70	U
108-87-2	Methyl cyclohexane	ND	10	0.40	U

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Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies
 Project Name : CONVENTUS
 Lab ID : L2246299-06D
 Client ID : BCP MW-5
 Sample Location : CONVENTUS/MAIN ST
 Sample Matrix : WATER
 Analytical Method : 1,8260C
 Lab File ID : V05220901N25
 Sample Amount : 0.5 ml
 Level : LOW
 Extract Volume (MeOH) : N/A

Lab Number : L2246299
 Project Number : U86
 Date Collected : 08/25/22 13:00
 Date Received : 08/25/22
 Date Analyzed : 09/02/22 03:46
 Dilution Factor : 20
 Analyst : MV
 Instrument ID : VOA105
 GC Column : RTX-502.2
 %Solids : N/A
 Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
75-09-2	Methylene chloride	ND	50	14.	U
75-34-3	1,1-Dichloroethane	ND	50	14.	U
67-66-3	Chloroform	ND	50	14.	U
56-23-5	Carbon tetrachloride	ND	10	2.7	U
78-87-5	1,2-Dichloropropane	ND	20	2.7	U
124-48-1	Dibromochloromethane	ND	10	3.0	U
79-00-5	1,1,2-Trichloroethane	ND	30	10.	U
127-18-4	Tetrachloroethene	ND	10	3.6	U
108-90-7	Chlorobenzene	ND	50	14.	U
75-69-4	Trichlorofluoromethane	ND	50	14.	U
107-06-2	1,2-Dichloroethane	ND	10	2.6	U UJ
71-55-6	1,1,1-Trichloroethane	ND	50	14.	U
75-27-4	Bromodichloromethane	ND	10	3.8	U
10061-02-6	trans-1,3-Dichloropropene	ND	10	3.3	U
10061-01-5	cis-1,3-Dichloropropene	ND	10	2.9	U
75-25-2	Bromoform	ND	40	13.	U
79-34-5	1,1,2,2-Tetrachloroethane	ND	10	3.3	U
71-43-2	Benzene	ND	10	3.2	U
108-88-3	Toluene	26	50	14.	J
100-41-4	Ethylbenzene	2100	50	14.	
74-87-3	Chloromethane	ND	50	14.	U
74-83-9	Bromomethane	ND	50	14.	U
75-01-4	Vinyl chloride	ND	20	1.4	U
75-00-3	Chloroethane	ND	50	14.	U
75-35-4	1,1-Dichloroethene	ND	10	3.4	U UJ

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies
 Project Name : CONVENTUS
 Lab ID : L2246299-06D
 Client ID : BCP MW-5
 Sample Location : CONVENTUS/MAIN ST
 Sample Matrix : WATER
 Analytical Method : 1,8260C
 Lab File ID : V05220901N25
 Sample Amount : 0.5 ml
 Level : LOW
 Extract Volume (MeOH) : N/A

Lab Number : L2246299
 Project Number : U86
 Date Collected : 08/25/22 13:00
 Date Received : 08/25/22
 Date Analyzed : 09/02/22 03:46
 Dilution Factor : 20
 Analyst : MV
 Instrument ID : VOA105
 GC Column : RTX-502.2
 %Solids : N/A
 Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
156-60-5	trans-1,2-Dichloroethene	ND	50	14.	U
79-01-6	Trichloroethene	ND	10	3.5	U
95-50-1	1,2-Dichlorobenzene	ND	50	14.	U
541-73-1	1,3-Dichlorobenzene	ND	50	14.	U
106-46-7	1,4-Dichlorobenzene	ND	50	14.	U
1634-04-4	Methyl tert butyl ether	ND	50	14.	U
179601-23-1	p/m-Xylene	7800	50	14.	
95-47-6	o-Xylene	100	50	14.	
156-59-2	cis-1,2-Dichloroethene	ND	50	14.	U
100-42-5	Styrene	ND	50	14.	U
75-71-8	Dichlorodifluoromethane	ND	100	20.	U
67-64-1	Acetone	ND	100	29.	U
75-15-0	Carbon disulfide	ND	100	20.	U
78-93-3	2-Butanone	ND	100	39.	U
108-10-1	4-Methyl-2-pentanone	ND	100	20.	U UJ
591-78-6	2-Hexanone	ND	100	20.	U UJ
106-93-4	1,2-Dibromoethane	ND	40	13.	U UJ
104-51-8	n-Butylbenzene	ND	50	14.	U
135-98-8	sec-Butylbenzene	ND	50	14.	U
98-06-6	tert-Butylbenzene	ND	50	14.	U
96-12-8	1,2-Dibromo-3-chloropropane	ND	50	14.	U
98-82-8	Isopropylbenzene	29	50	14.	J UJ
99-87-6	p-Isopropyltoluene	ND	50	14.	U
91-20-3	Naphthalene	840	50	14.	
103-65-1	n-Propylbenzene	130	50	14.	



Results Summary

Form 1

Volatile Organics by GC/MS

Client	: C&S Companies	Lab Number	: L2246299
Project Name	: CONVENTUS	Project Number	: U86
Lab ID	: L2246299-06D	Date Collected	: 08/25/22 13:00
Client ID	: BCP MW-5	Date Received	: 08/25/22
Sample Location	: CONVENTUS/MAIN ST	Date Analyzed	: 09/02/22 03:46
Sample Matrix	: WATER	Dilution Factor	: 20
Analytical Method	: 1,8260C	Analyst	: MV
Lab File ID	: V05220901N25	Instrument ID	: VOA105
Sample Amount	: 0.5 ml	GC Column	: RTX-502.2
Level	: LOW	%Solids	: N/A
Extract Volume (MeOH)	: N/A	Injection Volume	: N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
120-82-1	1,2,4-Trichlorobenzene	ND	50	14.	U
108-67-8	1,3,5-Trimethylbenzene	630	50	14.	
95-63-6	1,2,4-Trimethylbenzene	3100	50	14.	
79-20-9	Methyl Acetate	ND	40	4.7	U
110-82-7	Cyclohexane	240	200	5.4	UJ
76-13-1	Freon-113	ND	50	14.	U
108-87-2	Methyl cyclohexane	95	200	7.9	J

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Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies
 Project Name : CONVENTUS
 Lab ID : L2246299-07
 Client ID : DUP
 Sample Location : CONVENTUS/MAIN ST
 Sample Matrix : WATER
 Analytical Method : 1,8260C
 Lab File ID : V05220901N10
 Sample Amount : 10 ml
 Level : LOW
 Extract Volume (MeOH) : N/A

Lab Number : L2246299
 Project Number : U86
 Date Collected : 08/25/22 09:55
 Date Received : 08/25/22
 Date Analyzed : 09/01/22 21:57
 Dilution Factor : 1
 Analyst : MV
 Instrument ID : VOA105
 GC Column : RTX-502.2
 %Solids : N/A
 Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
75-09-2	Methylene chloride	ND	2.5	0.70	U
75-34-3	1,1-Dichloroethane	ND	2.5	0.70	U
67-66-3	Chloroform	ND	2.5	0.70	U
56-23-5	Carbon tetrachloride	ND	0.50	0.13	U
78-87-5	1,2-Dichloropropane	ND	1.0	0.14	U
124-48-1	Dibromochloromethane	ND	0.50	0.15	U
79-00-5	1,1,2-Trichloroethane	ND	1.5	0.50	U
127-18-4	Tetrachloroethene	ND	0.50	0.18	U
108-90-7	Chlorobenzene	ND	2.5	0.70	U
75-69-4	Trichlorofluoromethane	ND	2.5	0.70	U
107-06-2	1,2-Dichloroethane	ND	0.50	0.13	U UJ
71-55-6	1,1,1-Trichloroethane	ND	2.5	0.70	U
75-27-4	Bromodichloromethane	ND	0.50	0.19	U
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.16	U
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.14	U
75-25-2	Bromoform	ND	2.0	0.65	U
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.17	U
71-43-2	Benzene	ND	0.50	0.16	U
108-88-3	Toluene	ND	2.5	0.70	U
100-41-4	Ethylbenzene	ND	2.5	0.70	U
74-87-3	Chloromethane	ND	2.5	0.70	U
74-83-9	Bromomethane	ND	2.5	0.70	U
75-01-4	Vinyl chloride	ND	1.0	0.07	U
75-00-3	Chloroethane	ND	2.5	0.70	U
75-35-4	1,1-Dichloroethene	ND	0.50	0.17	U UJ



Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies
 Project Name : CONVENTUS
 Lab ID : L2246299-07
 Client ID : DUP
 Sample Location : CONVENTUS/MAIN ST
 Sample Matrix : WATER
 Analytical Method : 1,8260C
 Lab File ID : V05220901N10
 Sample Amount : 10 ml
 Level : LOW
 Extract Volume (MeOH) : N/A

Lab Number : L2246299
 Project Number : U86
 Date Collected : 08/25/22 09:55
 Date Received : 08/25/22
 Date Analyzed : 09/01/22 21:57
 Dilution Factor : 1
 Analyst : MV
 Instrument ID : VOA105
 GC Column : RTX-502.2
 %Solids : N/A
 Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
156-60-5	trans-1,2-Dichloroethene	ND	2.5	0.70	U
79-01-6	Trichloroethene	ND	0.50	0.18	U
95-50-1	1,2-Dichlorobenzene	ND	2.5	0.70	U
541-73-1	1,3-Dichlorobenzene	ND	2.5	0.70	U
106-46-7	1,4-Dichlorobenzene	ND	2.5	0.70	U
1634-04-4	Methyl tert butyl ether	ND	2.5	0.70	U
179601-23-1	p/m-Xylene	ND	2.5	0.70	U
95-47-6	o-Xylene	ND	2.5	0.70	U
156-59-2	cis-1,2-Dichloroethene	ND	2.5	0.70	U
100-42-5	Styrene	ND	2.5	0.70	U
75-71-8	Dichlorodifluoromethane	ND	5.0	1.0	U
67-64-1	Acetone	ND	5.0	1.5	U
75-15-0	Carbon disulfide	ND	5.0	1.0	U
78-93-3	2-Butanone	ND	5.0	1.9	U
108-10-1	4-Methyl-2-pentanone	ND	5.0	1.0	U UJ
591-78-6	2-Hexanone	ND	5.0	1.0	U UJ
106-93-4	1,2-Dibromoethane	ND	2.0	0.65	U UJ
104-51-8	n-Butylbenzene	ND	2.5	0.70	U
135-98-8	sec-Butylbenzene	ND	2.5	0.70	U
98-06-6	tert-Butylbenzene	ND	2.5	0.70	U
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.5	0.70	U
98-82-8	Isopropylbenzene	ND	2.5	0.70	U UJ
99-87-6	p-Isopropyltoluene	ND	2.5	0.70	U
91-20-3	Naphthalene	ND	2.5	0.70	U
103-65-1	n-Propylbenzene	ND	2.5	0.70	U

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies
 Project Name : CONVENTUS
 Lab ID : L2246299-07
 Client ID : DUP
 Sample Location : CONVENTUS/MAIN ST
 Sample Matrix : WATER
 Analytical Method : 1,8260C
 Lab File ID : V05220901N10
 Sample Amount : 10 ml
 Level : LOW
 Extract Volume (MeOH) : N/A

Lab Number : L2246299
 Project Number : U86
 Date Collected : 08/25/22 09:55
 Date Received : 08/25/22
 Date Analyzed : 09/01/22 21:57
 Dilution Factor : 1
 Analyst : MV
 Instrument ID : VOA105
 GC Column : RTX-502.2
 %Solids : N/A
 Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
120-82-1	1,2,4-Trichlorobenzene	ND	2.5	0.70	U
108-67-8	1,3,5-Trimethylbenzene	ND	2.5	0.70	U
95-63-6	1,2,4-Trimethylbenzene	ND	2.5	0.70	U
79-20-9	Methyl Acetate	ND	2.0	0.23	U
110-82-7	Cyclohexane	ND	10	0.27	U UJ
76-13-1	Freon-113	ND	2.5	0.70	U
108-87-2	Methyl cyclohexane	ND	10	0.40	U

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Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies
 Project Name : CONVENTUS
 Lab ID : L2246299-08
 Client ID : TRIP BLANK
 Sample Location : CONVENTUS/MAIN ST
 Sample Matrix : WATER
 Analytical Method : 1,8260C
 Lab File ID : V05220901N06
 Sample Amount : 10 ml
 Level : LOW
 Extract Volume (MeOH) : N/A

Lab Number : L2246299
 Project Number : U86
 Date Collected : 08/25/22 13:50
 Date Received : 08/25/22
 Date Analyzed : 09/01/22 20:23
 Dilution Factor : 1
 Analyst : MV
 Instrument ID : VOA105
 GC Column : RTX-502.2
 %Solids : N/A
 Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
75-09-2	Methylene chloride	ND	2.5	0.70	U
75-34-3	1,1-Dichloroethane	ND	2.5	0.70	U
67-66-3	Chloroform	ND	2.5	0.70	U
56-23-5	Carbon tetrachloride	ND	0.50	0.13	U
78-87-5	1,2-Dichloropropane	ND	1.0	0.14	U
124-48-1	Dibromochloromethane	ND	0.50	0.15	U
79-00-5	1,1,2-Trichloroethane	ND	1.5	0.50	U
127-18-4	Tetrachloroethene	ND	0.50	0.18	U
108-90-7	Chlorobenzene	ND	2.5	0.70	U
75-69-4	Trichlorofluoromethane	ND	2.5	0.70	U
107-06-2	1,2-Dichloroethane	ND	0.50	0.13	U UJ
71-55-6	1,1,1-Trichloroethane	ND	2.5	0.70	U
75-27-4	Bromodichloromethane	ND	0.50	0.19	U
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.16	U
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.14	U
75-25-2	Bromoform	ND	2.0	0.65	U
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.17	U
71-43-2	Benzene	ND	0.50	0.16	U
108-88-3	Toluene	ND	2.5	0.70	U
100-41-4	Ethylbenzene	ND	2.5	0.70	U
74-87-3	Chloromethane	ND	2.5	0.70	U
74-83-9	Bromomethane	ND	2.5	0.70	U
75-01-4	Vinyl chloride	ND	1.0	0.07	U
75-00-3	Chloroethane	ND	2.5	0.70	U
75-35-4	1,1-Dichloroethene	ND	0.50	0.17	U UJ

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies
 Project Name : CONVENTUS
 Lab ID : L2246299-08
 Client ID : TRIP BLANK
 Sample Location : CONVENTUS/MAIN ST
 Sample Matrix : WATER
 Analytical Method : 1,8260C
 Lab File ID : V05220901N06
 Sample Amount : 10 ml
 Level : LOW
 Extract Volume (MeOH) : N/A

Lab Number : L2246299
 Project Number : U86
 Date Collected : 08/25/22 13:50
 Date Received : 08/25/22
 Date Analyzed : 09/01/22 20:23
 Dilution Factor : 1
 Analyst : MV
 Instrument ID : VOA105
 GC Column : RTX-502.2
 %Solids : N/A
 Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
156-60-5	trans-1,2-Dichloroethene	ND	2.5	0.70	U
79-01-6	Trichloroethene	ND	0.50	0.18	U
95-50-1	1,2-Dichlorobenzene	ND	2.5	0.70	U
541-73-1	1,3-Dichlorobenzene	ND	2.5	0.70	U
106-46-7	1,4-Dichlorobenzene	ND	2.5	0.70	U
1634-04-4	Methyl tert butyl ether	ND	2.5	0.70	U
179601-23-1	p/m-Xylene	ND	2.5	0.70	U
95-47-6	o-Xylene	ND	2.5	0.70	U
156-59-2	cis-1,2-Dichloroethene	ND	2.5	0.70	U
100-42-5	Styrene	ND	2.5	0.70	U
75-71-8	Dichlorodifluoromethane	ND	5.0	1.0	U
67-64-1	Acetone	ND	5.0	1.5	U
75-15-0	Carbon disulfide	ND	5.0	1.0	U
78-93-3	2-Butanone	ND	5.0	1.9	U
108-10-1	4-Methyl-2-pentanone	ND	5.0	1.0	U UJ
591-78-6	2-Hexanone	ND	5.0	1.0	U UJ
106-93-4	1,2-Dibromoethane	ND	2.0	0.65	U UJ
104-51-8	n-Butylbenzene	ND	2.5	0.70	U
135-98-8	sec-Butylbenzene	ND	2.5	0.70	U
98-06-6	tert-Butylbenzene	ND	2.5	0.70	U
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.5	0.70	U
98-82-8	Isopropylbenzene	ND	2.5	0.70	U UJ
99-87-6	p-Isopropyltoluene	ND	2.5	0.70	U
91-20-3	Naphthalene	ND	2.5	0.70	U
103-65-1	n-Propylbenzene	ND	2.5	0.70	U

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies
 Project Name : CONVENTUS
 Lab ID : L2246299-08
 Client ID : TRIP BLANK
 Sample Location : CONVENTUS/MAIN ST
 Sample Matrix : WATER
 Analytical Method : 1,8260C
 Lab File ID : V05220901N06
 Sample Amount : 10 ml
 Level : LOW
 Extract Volume (MeOH) : N/A

Lab Number : L2246299
 Project Number : U86
 Date Collected : 08/25/22 13:50
 Date Received : 08/25/22
 Date Analyzed : 09/01/22 20:23
 Dilution Factor : 1
 Analyst : MV
 Instrument ID : VOA105
 GC Column : RTX-502.2
 %Solids : N/A
 Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
120-82-1	1,2,4-Trichlorobenzene	ND	2.5	0.70	U
108-67-8	1,3,5-Trimethylbenzene	ND	2.5	0.70	U
95-63-6	1,2,4-Trimethylbenzene	ND	2.5	0.70	U
79-20-9	Methyl Acetate	ND	2.0	0.23	U
110-82-7	Cyclohexane	ND	10	0.27	U UJ
76-13-1	Freon-113	ND	2.5	0.70	U
108-87-2	Methyl cyclohexane	ND	10	0.40	U

MKP 11/28/2022



Appendix B

Laboratory QC Documentation

Calibration Verification Summary

Form 7

Volatiles

Client : C&S Companies
 Project Name : CONVENTUS
 Instrument ID : VOA105
 Lab File ID : V05220901N02
 Sample No : WG1682814-2
 Channel :

Lab Number : L2246299
 Project Number : U86
 Calibration Date : 09/01/22 18:50
 Init. Calib. Date(s) : 08/19/22 08/19/22
 Init. Calib. Times : 14:23 17:52

Compound	Ave. RRF	RRF	Min RRF	%D	Max %D	Area%	Dev(min)
Fluorobenzene	1	1	-	0	20	67	0
Dichlorodifluoromethane	0.245	0.247	-	-0.8	20	63	0
Chloromethane	0.259	0.225	-	13.1	20	56	0
Vinyl chloride	0.28	0.296	-	-5.7	20	69	0
Bromomethane	0.283	0.333	-	-17.7	20	78	0
Chloroethane	0.214	0.299	-	-39.7*	20	88	0
Trichlorofluoromethane	0.558	0.71	-	-27.2*	20	80	0
Ethyl ether	0.12	0.144	-	-20	20	77	0
1,1-Dichloroethene	0.212	0.22	-	-3.8	20	68	0
Carbon disulfide	0.556	0.56	-	-0.7	20	67	0
Freon-113	0.235	0.252	-	-7.2	20	68	0
Iodomethane	0.292	0.12	-	58.9*	20	30	0
Acrolein	0.025	0.018	-	28*	20	49	0
Methylene chloride	0.229	0.221	-	3.5	20	65	0
Acetone	0.032	0.026	-	18.8	20	58	0
trans-1,2-Dichloroethene	0.238	0.241	-	-1.3	20	66	0
Methyl acetate	0.086	0.066	-	23.3*	20	55	0
Methyl tert-butyl ether	0.422	0.344	-	18.5	20	54	0
tert-Butyl alcohol	0.00767	0.00566*	-	26.2*	20	47	0
Diisopropyl ether	0.686	0.594	-	13.4	20	57	0
1,1-Dichloroethane	0.426	0.418	-	1.9	20	64	0
Halothane	0.189	0.187	-	1.1	20	66	0
Acrylonitrile	0.042	0.035	-	16.7	20	58	0
Ethyl tert-butyl ether	0.612	0.493	-	19.4	20	54	0
Vinyl acetate	0.428	0.316	-	26.2*	20	51	0
cis-1,2-Dichloroethene	0.263	0.258	-	1.9	20	64	0
2,2-Dichloropropane	0.373	0.334	-	10.5	20	59	0
Bromochloromethane	0.118	0.118	-	0	20	65	0
Cyclohexane	0.448	0.43	-	4	20	62	0
Chloroform	0.42	0.408	-	2.9	20	64	0
Ethyl acetate	0.131	0.101	-	22.9*	20	55	0
Carbon tetrachloride	0.359	0.345	-	3.9	20	65	0
Tetrahydrofuran	0.037	0.036	-	2.7	20	62	0
Dibromofluoromethane	0.26	0.27	-	-3.8	20	70	0
1,1,1-Trichloroethane	0.391	0.378	-	3.3	20	63	0
2-Butanone	0.051	0.04	-	21.6*	20	52	0
1,1-Dichloropropene	0.322	0.311	-	3.4	20	62	0
Benzene	0.935	0.923	-	1.3	20	65	0
tert-Amyl methyl ether	0.513	0.386	-	24.8*	20	51	0
1,2-Dichloroethane-d4	0.283	0.286	-	-1.1	20	65	0
1,2-Dichloroethane	0.291	0.266	-	8.6	20	60	0
Methyl cyclohexane	0.455	0.441	-	3.1	20	64	0
Trichloroethene	0.268	0.251	-	6.3	20	64	0

* Value outside of QC limits.



Calibration Verification Summary

Form 7

Volatiles

Client : C&S Companies
 Project Name : CONVENTUS
 Instrument ID : VOA105
 Lab File ID : V05220901N02
 Sample No : WG1682814-2
 Channel :

Lab Number : L2246299
 Project Number : U86
 Calibration Date : 09/01/22 18:50
 Init. Calib. Date(s) : 08/19/22 08/19/22
 Init. Calib. Times : 14:23 17:52

Compound	Ave. RRF	RRF	Min RRF	%D	Max %D	Area%	Dev(min)	
Dibromomethane	0.127	0.122	-	3.9	20	63	0	
1,2-Dichloropropane	0.233	0.224	-	3.9	20	62	0	
2-Chloroethyl vinyl ether	0.105	0.079	-	24.8*	20	48	0	
Bromodichloromethane	0.309	0.286*	-	7.4	20	61	0	NT
1,4-Dioxane	0.00087	0.00097*	-	-11.5	20	69	0	
cis-1,3-Dichloropropene	0.339	0.309	-	8.8	20	58	0	
Chlorobenzene-d5	1	1	-	0	20	66	0	
Toluene-d8	1.241	1.32	-	-6.4	20	69	0	
Toluene	0.769	0.769	-	0	20	65	0	
4-Methyl-2-pentanone	0.058	0.046	-	20.7*	20	49	0	
Tetrachloroethene	0.355	0.358	-	-0.8	20	65	0	
trans-1,3-Dichloropropene	0.355	0.319	-	10.1	20	54	0	
Ethyl methacrylate	0.252	0.198	-	21.4*	20	50	0	
1,1,2-Trichloroethane	0.18	0.171*	-	5	20	61	0	NT
Chlorodibromomethane	0.277	0.258	-	6.9	20	61	0	
1,3-Dichloropropane	0.394	0.379	-	3.8	20	60	0	
1,2-Dibromoethane	0.189	0.176*	-	6.9	20	58	0	
2-Hexanone	0.105	0.073	-	30.5*	20	46	0	
Chlorobenzene	0.883	0.869	-	1.6	20	65	0	
Ethylbenzene	1.55	1.513	-	2.4	20	63	0	
1,1,1,2-Tetrachloroethane	0.306	0.292	-	4.6	20	63	0	
p/m Xylene	0.634	0.618	-	2.5	20	64	0	
o Xylene	0.627	0.585	-	6.7	20	64	0	
Styrene	0.931	0.903	-	3	20	64	0	
1,4-Dichlorobenzene-d4	1	1	-	0	20	67	0	
Bromoform	0.294	0.24	-	18.4	20	60	0	
Isopropylbenzene	2.965	2.898	-	2.3	20	64	0	
4-Bromofluorobenzene	0.763	0.718	-	5.9	20	62	0	
Bromobenzene	0.676	0.636	-	5.9	20	64	0	
n-Propylbenzene	3.439	3.421	-	0.5	20	64	0	
1,4-Dichlorobutane	0.638	0.542	-	15	20	56	0	
1,1,2,2-Tetrachloroethane	0.463	0.397	-	14.3	20	59	0	
4-Ethyltoluene	2.905	2.835	-	2.4	20	65	0	
2-Chlorotoluene	1.888	1.828	-	3.2	20	65	0	
1,3,5-Trimethylbenzene	2.515	2.359	-	6.2	20	63	0	
1,2,3-Trichloropropane	0.359	0.289	-	19.5	20	54	0	
trans-1,4-Dichloro-2-buten	0.134	0.107	-	20.1*	20	50	0	
4-Chlorotoluene	2.008	1.866	-	7.1	20	62	0	NT
tert-Butylbenzene	2.176	2.121	-	2.5	20	65	0	
1,2,4-Trimethylbenzene	2.418	2.284	-	5.5	20	63	0	
sec-Butylbenzene	3.154	3.159	-	-0.2	20	66	0	
p-Isopropyltoluene	2.723	2.649	-	2.7	20	65	0	
1,3-Dichlorobenzene	1.346	1.291	-	4.1	20	65	0	

* Value outside of QC limits.



Calibration Verification Summary

Form 7

Volatiles

Client : C&S Companies
 Project Name : CONVENTUS
 Instrument ID : VOA105
 Lab File ID : V05220901N02
 Sample No : WG1682814-2
 Channel :

Lab Number : L2246299
 Project Number : U86
 Calibration Date : 09/01/22 18:50
 Init. Calib. Date(s) : 08/19/22 08/19/22
 Init. Calib. Times : 14:23 17:52

Compound	Ave. RRF	RRF	Min RRF	%D	Max %D	Area%	Dev(min)
1,4-Dichlorobenzene	1.375	1.304	-	5.2	20	66	0
p-Diethylbenzene	1.579	1.483	-	6.1	20	65	0
n-Butylbenzene	2.247	2.234	-	0.6	20	66	0
1,2-Dichlorobenzene	1.19	1.12	-	5.9	20	64	0
1,2,4,5-Tetramethylbenzene	2.102	1.86	-	11.5	20	62	0
1,2-Dibromo-3-chloropropan	0.061	0.047	-	23*	20	54	0
1,3,5-Trichlorobenzene	0.788	0.739	-	6.2	20	64	0
Hexachlorobutadiene	0.263	0.26	-	1.1	20	67	0
1,2,4-Trichlorobenzene	0.61	0.564	-	7.5	20	62	0
Naphthalene	1.194	0.959	-	19.7	20	54	0
1,2,3-Trichlorobenzene	0.484	0.424	-	12.4	20	61	0

* Value outside of QC limits.



Laboratory Control Sample Summary

Form 3

Volatiles

Client : C&S Companies
 Project Name : CONVENTUS
 Matrix : WATER
 LCS Sample ID : WG1682814-3 Analysis Date : 09/01/22 18:50 File ID : V05220901N02
 LCSD Sample ID : WG1682814-4 Analysis Date : 09/01/22 19:14 File ID : V05220901N03

Parameter	Laboratory Control Sample			Laboratory Control Duplicate			RPD	Recovery Limits	RPD Limit
	True (ug/l)	Found (ug/l)	%R	True (ug/l)	Found (ug/l)	%R			
Methylene chloride	10	9.7	97	10	9.7	97	0	70-130	20
1,1-Dichloroethane	10	9.8	98	10	9.4	94	4	70-130	20
Chloroform	10	9.7	97	10	9.2	92	5	70-130	20
Carbon tetrachloride	10	9.6	96	10	9.0	90	6	63-132	20
1,2-Dichloropropane	10	9.6	96	10	9.4	94	2	70-130	20
Dibromochloromethane	10	9.3	93	10	9.2	92	1	63-130	20
1,1,2-Trichloroethane	10	9.5	95	10	9.4	94	1	70-130	20
Tetrachloroethene	10	10	100	10	9.6	96	4	70-130	20
Chlorobenzene	10	9.8	98	10	9.5	95	3	75-130	20
Trichlorofluoromethane	10	13	130	10	12	120	8	62-150	20
1,2-Dichloroethane	10	9.1	91	10	9.4	94	3	70-130	20
1,1,1-Trichloroethane	10	9.7	97	10	9.2	92	5	67-130	20
Bromodichloromethane	10	9.2	92	10	9.2	92	0	67-130	20
trans-1,3-Dichloropropene	10	9.0	90	10	8.9	89	1	70-130	20
cis-1,3-Dichloropropene	10	9.1	91	10	8.9	89	2	70-130	20
Bromoform	10	8.2	82	10	8.0	80	2	54-136	20
1,1,2,2-Tetrachloroethane	10	8.6	86	10	8.6	86	0	67-130	20
Benzene	10	9.9	99	10	9.4	94	5	70-130	20
Toluene	10	10	100	10	9.6	96	4	70-130	20
Ethylbenzene	10	9.8	98	10	9.2	92	6	70-130	20
Chloromethane	10	8.7	87	10	8.5	85	2	64-130	20
Bromomethane	10	12	120	10	11	110	9	39-139	20
Vinyl chloride	10	10	100	10	9.9	99	1	55-140	20
Chloroethane	10	14	140 Q	10	13	130	7	55-138	20
1,1-Dichloroethene	10	10	100	10	9.3	93	7	61-145	20
trans-1,2-Dichloroethene	10	10	100	10	9.2	92	8	70-130	20



Appendix C

Validator Qualifications

KENNETH R. APPLIN

Geochemist/Data Validator

Ph.D., Geochemistry and Mineralogy, The Pennsylvania State University

M.S., Geochemistry and Mineralogy, The Pennsylvania State University

B.A., Geological Sciences, SUNY at Geneseo, NY

Dr. Applin has over 35 years of experience working with the geochemistry of natural waters. His prior experience includes working as an Assistant Professor of Geology at the University of Missouri-Columbia and as Chief Hydrogeologist and Geochemist with a leading engineering firm in Rochester, NY. In 1993, he established KR Applin and Associates, a small consulting business that focuses on the geochemistry of natural waters, especially as applied to problems involving the contamination of groundwater and surface water.

Dr. Applin is also an experienced analytical data validator and has provided data validation services since 1994 to a variety of clients performing brownfield cleanup projects, hazardous waste remediation, groundwater monitoring at solid waste facilities, and other projects requiring third-party data validation. Dr. Applin has several years of hands-on experience with the laboratory analysis of natural waters and has successfully completed the USEPA Region II certification courses for performing inorganic and organic analytical data validation.

MICHAEL K. PERRY

Chemist/Data Validator

B.S. Chemistry, Georgia State University, Atlanta, GA

A.A.S., Chemical Technology, Alfred State College, Alfred, NY

Mr. Perry has over 30 years of experience in the analytical laboratory business. During his early career, he spent several years as a laboratory analyst performing the analysis of soil, water, and air samples for inorganic and organic chemical parameters. During his last 20 years in the environmental laboratory business, he managed and directed two major analytical laboratories in Rochester, NY. His management responsibilities included oversight of the daily operations of the lab, staff training and supervision, the selection, purchase, and maintenance of analytical instruments, the introduction of new laboratory methods, analytical quality assurance and quality control, data acquisition and management, and other business-related activities.

Mr. Perry has an extensive working knowledge of the methods and procedures used for sampling and analyzing both inorganic and organic analytes in soil, water, and air. He is an accomplished laboratory chemist and is familiar with the analytical methods and procedures established under the USEPA Contract Laboratory Protocols (CLP), the NYSDEC Analytical Services Protocols (ASP), and the NYSDOH Environmental Laboratory Approval Program (ELAP).



ANALYTICAL REPORT

Lab Number:	L2270805
Client:	C&S Companies 141 Elm Street Suite 100 Buffalo, NY 14203
ATTN:	Richard Backert
Phone:	(716) 955-3024
Project Name:	CONVENTUS/MROW
Project Number:	1186
Report Date:	12/29/22

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: CONVENTUS/MROW
Project Number: 1186

Lab Number: L2270805
Report Date: 12/29/22

Alpha Sample ID	Client ID	Matrix	Sample Location	Collection Date/Time	Receive Date
L2270805-01	BCPMW-1	WATER	CONVENTUS 1 MAIN ST. BUFFALO NY	12/14/22 11:00	12/15/22
L2270805-02	DUP	WATER	CONVENTUS 1 MAIN ST. BUFFALO NY	12/14/22 11:00	12/15/22
L2270805-03	BCPMW-7	WATER	CONVENTUS 1 MAIN ST. BUFFALO NY	12/14/22 11:40	12/15/22
L2270805-04	BCPMW-4	WATER	CONVENTUS 1 MAIN ST. BUFFALO NY	12/14/22 12:25	12/15/22
L2270805-05	BCPMW-3	WATER	CONVENTUS 1 MAIN ST. BUFFALO NY	12/14/22 12:55	12/15/22
L2270805-06	BCPMW-6	WATER	CONVENTUS 1 MAIN ST. BUFFALO NY	12/14/22 13:30	12/15/22
L2270805-07	BCPMW-5	WATER	CONVENTUS 1 MAIN ST. BUFFALO NY	12/14/22 14:10	12/15/22
L2270805-08	TRIP BLANK	WATER	CONVENTUS 1 MAIN ST. BUFFALO NY	12/15/22 00:00	12/15/22
L2270805-09	MSMW-2	WATER	CONVENTUS 1 MAIN ST. BUFFALO NY	12/15/22 10:45	12/15/22

Project Name: CONVENTUS/MROW
Project Number: 1186

Lab Number: L2270805
Report Date: 12/29/22

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: CONVENTUS/MROW
Project Number: 1186

Lab Number: L2270805
Report Date: 12/29/22

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.

Volatile Organics

L2270805-07D: The pH was greater than two; however, the sample was analyzed within the method required holding time.

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: *Tiffani Morrissey* - Tiffani Morrissey

Title: Technical Director/Representative

Date: 12/29/22

ORGANICS

VOLATILES

Project Name: CONVENTUS/MROW**Lab Number:** L2270805**Project Number:** 1186**Report Date:** 12/29/22**SAMPLE RESULTS**

Lab ID: L2270805-01
 Client ID: BCPMW-1
 Sample Location: CONVENTUS 1 MAIN ST. BUFFALO NY

Date Collected: 12/14/22 11:00
 Date Received: 12/15/22
 Field Prep: Not Specified

Sample Depth:

Matrix: Water

Analytical Method: 1,8260D

Analytical Date: 12/21/22 14:10

Analyst: MJV

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: CONVENTUS/MROW**Lab Number:** L2270805**Project Number:** 1186**Report Date:** 12/29/22**SAMPLE RESULTS****Lab ID:** L2270805-01**Date Collected:** 12/14/22 11:00**Client ID:** BCPMW-1**Date Received:** 12/15/22**Sample Location:** CONVENTUS 1 MAIN ST. BUFFALO NY**Field Prep:** 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
1,2-Dibromoethane	ND		ug/l	2.0	0.65	1
n-Butylbenzene	ND		ug/l	2.5	0.70	1
sec-Butylbenzene	ND		ug/l	2.5	0.70	1
tert-Butylbenzene	ND		ug/l	2.5	0.70	1
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70	1
Isopropylbenzene	ND		ug/l	2.5	0.70	1
p-Isopropyltoluene	ND		ug/l	2.5	0.70	1
Naphthalene	ND		ug/l	2.5	0.70	1
n-Propylbenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70	1
1,3,5-Trimethylbenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trimethylbenzene	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
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	96		70-130
Toluene-d8	98		70-130
4-Bromofluorobenzene	102		70-130
Dibromofluoromethane	111		70-130

Project Name: CONVENTUS/MROW**Lab Number:** L2270805**Project Number:** 1186**Report Date:** 12/29/22**SAMPLE RESULTS**

Lab ID: L2270805-02

Date Collected: 12/14/22 11:00

Client ID: DUP

Date Received: 12/15/22

Sample Location: CONVENTUS 1 MAIN ST. BUFFALO NY

Field Prep: Not Specified

Sample Depth:

Matrix: Water

Analytical Method: 1,8260D

Analytical Date: 12/21/22 14:34

Analyst: MJV

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: CONVENTUS/MROW**Lab Number:** L2270805**Project Number:** 1186**Report Date:** 12/29/22**SAMPLE RESULTS****Lab ID:** L2270805-02**Date Collected:** 12/14/22 11:00**Client ID:** DUP**Date Received:** 12/15/22**Sample Location:** CONVENTUS 1 MAIN ST. BUFFALO NY**Field Prep:** 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
1,2-Dibromoethane	ND		ug/l	2.0	0.65	1
n-Butylbenzene	ND		ug/l	2.5	0.70	1
sec-Butylbenzene	ND		ug/l	2.5	0.70	1
tert-Butylbenzene	ND		ug/l	2.5	0.70	1
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70	1
Isopropylbenzene	ND		ug/l	2.5	0.70	1
p-Isopropyltoluene	ND		ug/l	2.5	0.70	1
Naphthalene	ND		ug/l	2.5	0.70	1
n-Propylbenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70	1
1,3,5-Trimethylbenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trimethylbenzene	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
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	89		70-130
Toluene-d8	98		70-130
4-Bromofluorobenzene	102		70-130
Dibromofluoromethane	103		70-130

Project Name: CONVENTUS/MROW**Lab Number:** L2270805**Project Number:** 1186**Report Date:** 12/29/22**SAMPLE RESULTS**

Lab ID: L2270805-03
 Client ID: BCPMW-7
 Sample Location: CONVENTUS 1 MAIN ST. BUFFALO NY

Date Collected: 12/14/22 11:40
 Date Received: 12/15/22
 Field Prep: Not Specified

Sample Depth:

Matrix: Water

Analytical Method: 1,8260D

Analytical Date: 12/22/22 22:51

Analyst: MJV

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: CONVENTUS/MROW**Lab Number:** L2270805**Project Number:** 1186**Report Date:** 12/29/22**SAMPLE RESULTS****Lab ID:** L2270805-03**Date Collected:** 12/14/22 11:40**Client ID:** BCPMW-7**Date Received:** 12/15/22**Sample Location:** CONVENTUS 1 MAIN ST. BUFFALO NY**Field Prep:** 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
1,2-Dibromoethane	ND		ug/l	2.0	0.65	1
n-Butylbenzene	ND		ug/l	2.5	0.70	1
sec-Butylbenzene	ND		ug/l	2.5	0.70	1
tert-Butylbenzene	ND		ug/l	2.5	0.70	1
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70	1
Isopropylbenzene	ND		ug/l	2.5	0.70	1
p-Isopropyltoluene	ND		ug/l	2.5	0.70	1
Naphthalene	ND		ug/l	2.5	0.70	1
n-Propylbenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70	1
1,3,5-Trimethylbenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trimethylbenzene	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
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	118		70-130
Toluene-d8	102		70-130
4-Bromofluorobenzene	113		70-130
Dibromofluoromethane	109		70-130

Project Name: CONVENTUS/MROW**Lab Number:** L2270805**Project Number:** 1186**Report Date:** 12/29/22**SAMPLE RESULTS**

Lab ID: L2270805-04 D
 Client ID: BCPMW-4
 Sample Location: CONVENTUS 1 MAIN ST. BUFFALO NY

Date Collected: 12/14/22 12:25
 Date Received: 12/15/22
 Field Prep: Not Specified

Sample Depth:

Matrix: Water

Analytical Method: 1,8260D

Analytical Date: 12/21/22 15:23

Analyst: MJV

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westborough Lab						
Methylene chloride	ND		ug/l	25	7.0	10
1,1-Dichloroethane	ND		ug/l	25	7.0	10
Chloroform	ND		ug/l	25	7.0	10
Carbon tetrachloride	ND		ug/l	5.0	1.3	10
1,2-Dichloropropane	ND		ug/l	10	1.4	10
Dibromochloromethane	ND		ug/l	5.0	1.5	10
1,1,2-Trichloroethane	ND		ug/l	15	5.0	10
Tetrachloroethene	ND		ug/l	5.0	1.8	10
Chlorobenzene	ND		ug/l	25	7.0	10
Trichlorofluoromethane	ND		ug/l	25	7.0	10
1,2-Dichloroethane	ND		ug/l	5.0	1.3	10
1,1,1-Trichloroethane	ND		ug/l	25	7.0	10
Bromodichloromethane	ND		ug/l	5.0	1.9	10
trans-1,3-Dichloropropene	ND		ug/l	5.0	1.6	10
cis-1,3-Dichloropropene	ND		ug/l	5.0	1.4	10
Bromoform	ND		ug/l	20	6.5	10
1,1,2,2-Tetrachloroethane	ND		ug/l	5.0	1.7	10
Benzene	3.5	J	ug/l	5.0	1.6	10
Toluene	63		ug/l	25	7.0	10
Ethylbenzene	1100		ug/l	25	7.0	10
Chloromethane	ND		ug/l	25	7.0	10
Bromomethane	ND		ug/l	25	7.0	10
Vinyl chloride	ND		ug/l	10	0.71	10
Chloroethane	ND		ug/l	25	7.0	10
1,1-Dichloroethene	ND		ug/l	5.0	1.7	10
trans-1,2-Dichloroethene	ND		ug/l	25	7.0	10
Trichloroethene	ND		ug/l	5.0	1.8	10
1,2-Dichlorobenzene	ND		ug/l	25	7.0	10

Project Name: CONVENTUS/MROW**Lab Number:** L2270805**Project Number:** 1186**Report Date:** 12/29/22**SAMPLE RESULTS**

Lab ID: L2270805-04 D

Date Collected: 12/14/22 12:25

Client ID: BCPMW-4

Date Received: 12/15/22

Sample Location: CONVENTUS 1 MAIN ST. BUFFALO NY

Field Prep: 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	25	7.0	10
1,4-Dichlorobenzene	ND		ug/l	25	7.0	10
Methyl tert butyl ether	ND		ug/l	25	7.0	10
p/m-Xylene	210		ug/l	25	7.0	10
o-Xylene	16	J	ug/l	25	7.0	10
cis-1,2-Dichloroethene	ND		ug/l	25	7.0	10
Styrene	ND		ug/l	25	7.0	10
Dichlorodifluoromethane	ND		ug/l	50	10.	10
Acetone	ND		ug/l	50	15.	10
Carbon disulfide	ND		ug/l	50	10.	10
2-Butanone	ND		ug/l	50	19.	10
4-Methyl-2-pentanone	ND		ug/l	50	10.	10
2-Hexanone	ND		ug/l	50	10.	10
1,2-Dibromoethane	ND		ug/l	20	6.5	10
n-Butylbenzene	11	J	ug/l	25	7.0	10
sec-Butylbenzene	ND		ug/l	25	7.0	10
tert-Butylbenzene	ND		ug/l	25	7.0	10
1,2-Dibromo-3-chloropropane	ND		ug/l	25	7.0	10
Isopropylbenzene	30		ug/l	25	7.0	10
p-Isopropyltoluene	ND		ug/l	25	7.0	10
Naphthalene	290		ug/l	25	7.0	10
n-Propylbenzene	160		ug/l	25	7.0	10
1,2,4-Trichlorobenzene	ND		ug/l	25	7.0	10
1,3,5-Trimethylbenzene	ND		ug/l	25	7.0	10
1,2,4-Trimethylbenzene	1200		ug/l	25	7.0	10
Methyl Acetate	ND		ug/l	20	2.3	10
Cyclohexane	140		ug/l	100	2.7	10
Freon-113	ND		ug/l	25	7.0	10
Methyl cyclohexane	43	J	ug/l	100	4.0	10

Surrogate	% Recovery	Qualifier	Acceptance Criteria
1,2-Dichloroethane-d4	94		70-130
Toluene-d8	102		70-130
4-Bromofluorobenzene	103		70-130
Dibromofluoromethane	92		70-130

Project Name: CONVENTUS/MROW**Lab Number:** L2270805**Project Number:** 1186**Report Date:** 12/29/22**SAMPLE RESULTS**

Lab ID: L2270805-05
 Client ID: BCPMW-3
 Sample Location: CONVENTUS 1 MAIN ST. BUFFALO NY

Date Collected: 12/14/22 12:55
 Date Received: 12/15/22
 Field Prep: Not Specified

Sample Depth:

Matrix: Water

Analytical Method: 1,8260D

Analytical Date: 12/21/22 15:48

Analyst: MJV

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	11		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: CONVENTUS/MROW**Lab Number:** L2270805**Project Number:** 1186**Report Date:** 12/29/22**SAMPLE RESULTS****Lab ID:** L2270805-05**Date Collected:** 12/14/22 12:55**Client ID:** BCPMW-3**Date Received:** 12/15/22**Sample Location:** CONVENTUS 1 MAIN ST. BUFFALO NY**Field Prep:** 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	12		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
1,2-Dibromoethane	ND		ug/l	2.0	0.65	1
n-Butylbenzene	ND		ug/l	2.5	0.70	1
sec-Butylbenzene	ND		ug/l	2.5	0.70	1
tert-Butylbenzene	ND		ug/l	2.5	0.70	1
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70	1
Isopropylbenzene	ND		ug/l	2.5	0.70	1
p-Isopropyltoluene	ND		ug/l	2.5	0.70	1
Naphthalene	4.2		ug/l	2.5	0.70	1
n-Propylbenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70	1
1,3,5-Trimethylbenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trimethylbenzene	13		ug/l	2.5	0.70	1
Methyl Acetate	ND		ug/l	2.0	0.23	1
Cyclohexane	1.5	J	ug/l	10	0.27	1
Freon-113	ND		ug/l	2.5	0.70	1
Methyl cyclohexane	0.65	J	ug/l	10	0.40	1

Surrogate	% Recovery	Qualifier	Acceptance Criteria
1,2-Dichloroethane-d4	99		70-130
Toluene-d8	99		70-130
4-Bromofluorobenzene	113		70-130
Dibromofluoromethane	111		70-130

Project Name: CONVENTUS/MROW**Lab Number:** L2270805**Project Number:** 1186**Report Date:** 12/29/22**SAMPLE RESULTS**

Lab ID: L2270805-06
 Client ID: BCPMW-6
 Sample Location: CONVENTUS 1 MAIN ST. BUFFALO NY

Date Collected: 12/14/22 13:30
 Date Received: 12/15/22
 Field Prep: Not Specified

Sample Depth:

Matrix: Water

Analytical Method: 1,8260D

Analytical Date: 12/21/22 16:12

Analyst: MJV

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: CONVENTUS/MROW**Lab Number:** L2270805**Project Number:** 1186**Report Date:** 12/29/22**SAMPLE RESULTS****Lab ID:** L2270805-06**Date Collected:** 12/14/22 13:30**Client ID:** BCPMW-6**Date Received:** 12/15/22**Sample Location:** CONVENTUS 1 MAIN ST. BUFFALO NY**Field Prep:** 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
1,2-Dibromoethane	ND		ug/l	2.0	0.65	1
n-Butylbenzene	ND		ug/l	2.5	0.70	1
sec-Butylbenzene	ND		ug/l	2.5	0.70	1
tert-Butylbenzene	ND		ug/l	2.5	0.70	1
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70	1
Isopropylbenzene	ND		ug/l	2.5	0.70	1
p-Isopropyltoluene	ND		ug/l	2.5	0.70	1
Naphthalene	ND		ug/l	2.5	0.70	1
n-Propylbenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70	1
1,3,5-Trimethylbenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trimethylbenzene	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
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	98		70-130
Toluene-d8	99		70-130
4-Bromofluorobenzene	100		70-130
Dibromofluoromethane	113		70-130

Project Name: CONVENTUS/MROW**Lab Number:** L2270805**Project Number:** 1186**Report Date:** 12/29/22**SAMPLE RESULTS**

Lab ID: L2270805-07 D
 Client ID: BCPMW-5
 Sample Location: CONVENTUS 1 MAIN ST. BUFFALO NY

Date Collected: 12/14/22 14:10
 Date Received: 12/15/22
 Field Prep: Not Specified

Sample Depth:

Matrix: Water

Analytical Method: 1,8260D

Analytical Date: 12/21/22 16:37

Analyst: MJV

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westborough Lab						
Methylene chloride	ND		ug/l	25	7.0	10
1,1-Dichloroethane	ND		ug/l	25	7.0	10
Chloroform	ND		ug/l	25	7.0	10
Carbon tetrachloride	ND		ug/l	5.0	1.3	10
1,2-Dichloropropane	ND		ug/l	10	1.4	10
Dibromochloromethane	ND		ug/l	5.0	1.5	10
1,1,2-Trichloroethane	ND		ug/l	15	5.0	10
Tetrachloroethene	ND		ug/l	5.0	1.8	10
Chlorobenzene	ND		ug/l	25	7.0	10
Trichlorofluoromethane	ND		ug/l	25	7.0	10
1,2-Dichloroethane	ND		ug/l	5.0	1.3	10
1,1,1-Trichloroethane	ND		ug/l	25	7.0	10
Bromodichloromethane	ND		ug/l	5.0	1.9	10
trans-1,3-Dichloropropene	ND		ug/l	5.0	1.6	10
cis-1,3-Dichloropropene	ND		ug/l	5.0	1.4	10
Bromoform	ND		ug/l	20	6.5	10
1,1,2,2-Tetrachloroethane	ND		ug/l	5.0	1.7	10
Benzene	ND		ug/l	5.0	1.6	10
Toluene	9.6	J	ug/l	25	7.0	10
Ethylbenzene	820		ug/l	25	7.0	10
Chloromethane	ND		ug/l	25	7.0	10
Bromomethane	ND		ug/l	25	7.0	10
Vinyl chloride	ND		ug/l	10	0.71	10
Chloroethane	ND		ug/l	25	7.0	10
1,1-Dichloroethene	ND		ug/l	5.0	1.7	10
trans-1,2-Dichloroethene	ND		ug/l	25	7.0	10
Trichloroethene	ND		ug/l	5.0	1.8	10
1,2-Dichlorobenzene	ND		ug/l	25	7.0	10

Project Name: CONVENTUS/MROW**Lab Number:** L2270805**Project Number:** 1186**Report Date:** 12/29/22**SAMPLE RESULTS**

Lab ID: L2270805-07 D
 Client ID: BCPMW-5
 Sample Location: CONVENTUS 1 MAIN ST. BUFFALO NY

Date Collected: 12/14/22 14:10
 Date Received: 12/15/22
 Field Prep: 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	25	7.0	10
1,4-Dichlorobenzene	ND		ug/l	25	7.0	10
Methyl tert butyl ether	ND		ug/l	25	7.0	10
p/m-Xylene	2400		ug/l	25	7.0	10
o-Xylene	62		ug/l	25	7.0	10
cis-1,2-Dichloroethene	ND		ug/l	25	7.0	10
Styrene	ND		ug/l	25	7.0	10
Dichlorodifluoromethane	ND		ug/l	50	10.	10
Acetone	ND		ug/l	50	15.	10
Carbon disulfide	ND		ug/l	50	10.	10
2-Butanone	ND		ug/l	50	19.	10
4-Methyl-2-pentanone	ND		ug/l	50	10.	10
2-Hexanone	ND		ug/l	50	10.	10
1,2-Dibromoethane	ND		ug/l	20	6.5	10
n-Butylbenzene	ND		ug/l	25	7.0	10
sec-Butylbenzene	ND		ug/l	25	7.0	10
tert-Butylbenzene	ND		ug/l	25	7.0	10
1,2-Dibromo-3-chloropropane	ND		ug/l	25	7.0	10
Isopropylbenzene	16	J	ug/l	25	7.0	10
p-Isopropyltoluene	ND		ug/l	25	7.0	10
Naphthalene	520		ug/l	25	7.0	10
n-Propylbenzene	71		ug/l	25	7.0	10
1,2,4-Trichlorobenzene	ND		ug/l	25	7.0	10
1,3,5-Trimethylbenzene	350		ug/l	25	7.0	10
1,2,4-Trimethylbenzene	1400		ug/l	25	7.0	10
Methyl Acetate	ND		ug/l	20	2.3	10
Cyclohexane	120		ug/l	100	2.7	10
Freon-113	ND		ug/l	25	7.0	10
Methyl cyclohexane	54	J	ug/l	100	4.0	10

Surrogate	% Recovery	Qualifier	Acceptance Criteria
1,2-Dichloroethane-d4	100		70-130
Toluene-d8	100		70-130
4-Bromofluorobenzene	103		70-130
Dibromofluoromethane	102		70-130

Project Name: CONVENTUS/MROW**Lab Number:** L2270805**Project Number:** 1186**Report Date:** 12/29/22**SAMPLE RESULTS**

Lab ID: L2270805-08
 Client ID: TRIP BLANK
 Sample Location: CONVENTUS 1 MAIN ST. BUFFALO NY

Date Collected: 12/15/22 00:00
 Date Received: 12/15/22
 Field Prep: Not Specified

Sample Depth:

Matrix: Water

Analytical Method: 1,8260D

Analytical Date: 12/23/22 14:38

Analyst: MJV

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: CONVENTUS/MROW**Lab Number:** L2270805**Project Number:** 1186**Report Date:** 12/29/22**SAMPLE RESULTS****Lab ID:** L2270805-08**Date Collected:** 12/15/22 00:00**Client ID:** TRIP BLANK**Date Received:** 12/15/22**Sample Location:** CONVENTUS 1 MAIN ST. BUFFALO NY**Field Prep:** 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
1,2-Dibromoethane	ND		ug/l	2.0	0.65	1
n-Butylbenzene	ND		ug/l	2.5	0.70	1
sec-Butylbenzene	ND		ug/l	2.5	0.70	1
tert-Butylbenzene	ND		ug/l	2.5	0.70	1
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70	1
Isopropylbenzene	ND		ug/l	2.5	0.70	1
p-Isopropyltoluene	ND		ug/l	2.5	0.70	1
Naphthalene	ND		ug/l	2.5	0.70	1
n-Propylbenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70	1
1,3,5-Trimethylbenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trimethylbenzene	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
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	102		70-130
Toluene-d8	104		70-130
4-Bromofluorobenzene	112		70-130
Dibromofluoromethane	99		70-130

Project Name: CONVENTUS/MROW**Lab Number:** L2270805**Project Number:** 1186**Report Date:** 12/29/22**SAMPLE RESULTS**

Lab ID: L2270805-09 D
 Client ID: MSMW-2
 Sample Location: CONVENTUS 1 MAIN ST. BUFFALO NY

Date Collected: 12/15/22 10:45
 Date Received: 12/15/22
 Field Prep: Not Specified

Sample Depth:

Matrix: Water
 Analytical Method: 1,8260D
 Analytical Date: 12/21/22 17:01
 Analyst: MJV

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westborough Lab						
Methylene chloride	ND		ug/l	25	7.0	10
1,1-Dichloroethane	ND		ug/l	25	7.0	10
Chloroform	ND		ug/l	25	7.0	10
Carbon tetrachloride	ND		ug/l	5.0	1.3	10
1,2-Dichloropropane	ND		ug/l	10	1.4	10
Dibromochloromethane	ND		ug/l	5.0	1.5	10
1,1,2-Trichloroethane	ND		ug/l	15	5.0	10
Tetrachloroethene	ND		ug/l	5.0	1.8	10
Chlorobenzene	ND		ug/l	25	7.0	10
Trichlorofluoromethane	ND		ug/l	25	7.0	10
1,2-Dichloroethane	ND		ug/l	5.0	1.3	10
1,1,1-Trichloroethane	ND		ug/l	25	7.0	10
Bromodichloromethane	ND		ug/l	5.0	1.9	10
trans-1,3-Dichloropropene	ND		ug/l	5.0	1.6	10
cis-1,3-Dichloropropene	ND		ug/l	5.0	1.4	10
Bromoform	ND		ug/l	20	6.5	10
1,1,2,2-Tetrachloroethane	ND		ug/l	5.0	1.7	10
Benzene	280		ug/l	5.0	1.6	10
Toluene	210		ug/l	25	7.0	10
Ethylbenzene	820		ug/l	25	7.0	10
Chloromethane	ND		ug/l	25	7.0	10
Bromomethane	ND		ug/l	25	7.0	10
Vinyl chloride	ND		ug/l	10	0.71	10
Chloroethane	ND		ug/l	25	7.0	10
1,1-Dichloroethene	ND		ug/l	5.0	1.7	10
trans-1,2-Dichloroethene	ND		ug/l	25	7.0	10
Trichloroethene	ND		ug/l	5.0	1.8	10
1,2-Dichlorobenzene	ND		ug/l	25	7.0	10

Project Name: CONVENTUS/MROW**Lab Number:** L2270805**Project Number:** 1186**Report Date:** 12/29/22**SAMPLE RESULTS**

Lab ID: L2270805-09 D

Date Collected: 12/15/22 10:45

Client ID: MSMW-2

Date Received: 12/15/22

Sample Location: CONVENTUS 1 MAIN ST. BUFFALO NY

Field Prep: 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	25	7.0	10
1,4-Dichlorobenzene	ND		ug/l	25	7.0	10
Methyl tert butyl ether	ND		ug/l	25	7.0	10
p/m-Xylene	1900		ug/l	25	7.0	10
o-Xylene	490		ug/l	25	7.0	10
cis-1,2-Dichloroethene	ND		ug/l	25	7.0	10
Styrene	ND		ug/l	25	7.0	10
Dichlorodifluoromethane	ND		ug/l	50	10.	10
Acetone	ND		ug/l	50	15.	10
Carbon disulfide	ND		ug/l	50	10.	10
2-Butanone	ND		ug/l	50	19.	10
4-Methyl-2-pentanone	ND		ug/l	50	10.	10
2-Hexanone	ND		ug/l	50	10.	10
1,2-Dibromoethane	ND		ug/l	20	6.5	10
n-Butylbenzene	12	J	ug/l	25	7.0	10
sec-Butylbenzene	ND		ug/l	25	7.0	10
tert-Butylbenzene	ND		ug/l	25	7.0	10
1,2-Dibromo-3-chloropropane	ND		ug/l	25	7.0	10
Isopropylbenzene	25		ug/l	25	7.0	10
p-Isopropyltoluene	ND		ug/l	25	7.0	10
Naphthalene	340		ug/l	25	7.0	10
n-Propylbenzene	94		ug/l	25	7.0	10
1,2,4-Trichlorobenzene	ND		ug/l	25	7.0	10
1,3,5-Trimethylbenzene	490		ug/l	25	7.0	10
1,2,4-Trimethylbenzene	1200		ug/l	25	7.0	10
Methyl Acetate	ND		ug/l	20	2.3	10
Cyclohexane	160		ug/l	100	2.7	10
Freon-113	ND		ug/l	25	7.0	10
Methyl cyclohexane	110		ug/l	100	4.0	10

Surrogate	% Recovery	Qualifier	Acceptance Criteria
1,2-Dichloroethane-d4	89		70-130
Toluene-d8	102		70-130
4-Bromofluorobenzene	100		70-130
Dibromofluoromethane	91		70-130

Project Name: CONVENTUS/MROW
Project Number: 1186

Lab Number: L2270805
Report Date: 12/29/22

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260D
 Analytical Date: 12/21/22 08:53
 Analyst: PID

Parameter	Result	Qualifier	Units	RL	MDL
Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-02,04-07,09 Batch: WG1726668-5					
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
1,1,1,2-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

Project Name: CONVENTUS/MROW
Project Number: 1186

Lab Number: L2270805
Report Date: 12/29/22

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260D
 Analytical Date: 12/21/22 08:53
 Analyst: PID

Parameter	Result	Qualifier	Units	RL	MDL
Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-02,04-07,09 Batch: WG1726668-5					
1,4-Dichlorobenzene	ND		ug/l	2.5	0.70
Methyl tert butyl ether	ND		ug/l	2.5	0.70
p/m-Xylene	ND		ug/l	2.5	0.70
o-Xylene	ND		ug/l	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/l	5.0	1.0
Acetone	ND		ug/l	5.0	1.5
Carbon disulfide	ND		ug/l	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
1,2-Dibromoethane	ND		ug/l	2.0	0.65
n-Butylbenzene	ND		ug/l	2.5	0.70
sec-Butylbenzene	ND		ug/l	2.5	0.70
tert-Butylbenzene	ND		ug/l	2.5	0.70
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70
Isopropylbenzene	ND		ug/l	2.5	0.70
p-Isopropyltoluene	ND		ug/l	2.5	0.70
Naphthalene	ND		ug/l	2.5	0.70
n-Propylbenzene	ND		ug/l	2.5	0.70
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70
1,3,5-Trimethylbenzene	ND		ug/l	2.5	0.70
1,2,4-Trimethylbenzene	ND		ug/l	2.5	0.70
Methyl Acetate	ND		ug/l	2.0	0.23
Cyclohexane	ND		ug/l	10	0.27
Freon-113	ND		ug/l	2.5	0.70
Methyl cyclohexane	ND		ug/l	10	0.40

Project Name: CONVENTUS/MROW
Project Number: 1186

Lab Number: L2270805
Report Date: 12/29/22

Method Blank Analysis
Batch Quality Control

Analytical Method: 1,8260D
Analytical Date: 12/21/22 08:53
Analyst: PID

Parameter	Result	Qualifier	Units	RL	MDL
Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-02,04-07,09 Batch: WG1726668-5					

Surrogate	%Recovery	Qualifier	Acceptance Criteria
1,2-Dichloroethane-d4	91		70-130
Toluene-d8	99		70-130
4-Bromofluorobenzene	103		70-130
Dibromofluoromethane	111		70-130

Project Name: CONVENTUS/MROW

Lab Number: L2270805

Project Number: 1186

Report Date: 12/29/22

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260D

Analytical Date: 12/22/22 17:51

Analyst: AJK

Parameter	Result	Qualifier	Units	RL	MDL
Volatile Organics by GC/MS - Westborough Lab for sample(s): 03 Batch: WG1726975-5					
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

Project Name: CONVENTUS/MROW
Project Number: 1186

Lab Number: L2270805
Report Date: 12/29/22

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260D
Analytical Date: 12/22/22 17:51
Analyst: AJK

Parameter	Result	Qualifier	Units	RL	MDL
Volatile Organics by GC/MS - Westborough Lab for sample(s): 03 Batch: WG1726975-5					
1,4-Dichlorobenzene	ND		ug/l	2.5	0.70
Methyl tert butyl ether	ND		ug/l	2.5	0.70
p/m-Xylene	ND		ug/l	2.5	0.70
o-Xylene	ND		ug/l	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/l	5.0	1.0
Acetone	ND		ug/l	5.0	1.5
Carbon disulfide	ND		ug/l	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
1,2-Dibromoethane	ND		ug/l	2.0	0.65
n-Butylbenzene	ND		ug/l	2.5	0.70
sec-Butylbenzene	ND		ug/l	2.5	0.70
tert-Butylbenzene	ND		ug/l	2.5	0.70
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70
Isopropylbenzene	ND		ug/l	2.5	0.70
p-Isopropyltoluene	ND		ug/l	2.5	0.70
Naphthalene	ND		ug/l	2.5	0.70
n-Propylbenzene	ND		ug/l	2.5	0.70
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70
1,3,5-Trimethylbenzene	ND		ug/l	2.5	0.70
1,2,4-Trimethylbenzene	ND		ug/l	2.5	0.70
Methyl Acetate	ND		ug/l	2.0	0.23
Cyclohexane	ND		ug/l	10	0.27
Freon-113	ND		ug/l	2.5	0.70
Methyl cyclohexane	ND		ug/l	10	0.40

Project Name: CONVENTUS/MROW
Project Number: 1186

Lab Number: L2270805
Report Date: 12/29/22

Method Blank Analysis
Batch Quality Control

Analytical Method: 1,8260D
 Analytical Date: 12/22/22 17:51
 Analyst: AJK

Parameter	Result	Qualifier	Units	RL	MDL
Volatile Organics by GC/MS - Westborough Lab for sample(s): 03 Batch: WG1726975-5					

Surrogate	%Recovery	Qualifier	Acceptance Criteria
1,2-Dichloroethane-d4	109		70-130
Toluene-d8	101		70-130
4-Bromofluorobenzene	110		70-130
Dibromofluoromethane	105		70-130

Project Name: CONVENTUS/MROW
Project Number: 1186

Lab Number: L2270805
Report Date: 12/29/22

Method Blank Analysis
Batch Quality Control

Analytical Method: 1,8260D
Analytical Date: 12/23/22 08:59
Analyst: PID

Parameter	Result	Qualifier	Units	RL	MDL
Volatile Organics by GC/MS - Westborough Lab for sample(s): 08 Batch: WG1727559-5					
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
1,1,1,2-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

Project Name: CONVENTUS/MROW
Project Number: 1186

Lab Number: L2270805
Report Date: 12/29/22

Method Blank Analysis
Batch Quality Control

Analytical Method: 1,8260D
 Analytical Date: 12/23/22 08:59
 Analyst: PID

Parameter	Result	Qualifier	Units	RL	MDL
Volatile Organics by GC/MS - Westborough Lab for sample(s): 08 Batch: WG1727559-5					
1,4-Dichlorobenzene	ND		ug/l	2.5	0.70
Methyl tert butyl ether	ND		ug/l	2.5	0.70
p/m-Xylene	ND		ug/l	2.5	0.70
o-Xylene	ND		ug/l	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/l	5.0	1.0
Acetone	ND		ug/l	5.0	1.5
Carbon disulfide	ND		ug/l	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
1,2-Dibromoethane	ND		ug/l	2.0	0.65
n-Butylbenzene	ND		ug/l	2.5	0.70
sec-Butylbenzene	ND		ug/l	2.5	0.70
tert-Butylbenzene	ND		ug/l	2.5	0.70
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70
Isopropylbenzene	ND		ug/l	2.5	0.70
p-Isopropyltoluene	ND		ug/l	2.5	0.70
Naphthalene	ND		ug/l	2.5	0.70
n-Propylbenzene	ND		ug/l	2.5	0.70
1,2,4-Trichlorobenzene	ND		ug/l	2.5	0.70
1,3,5-Trimethylbenzene	ND		ug/l	2.5	0.70
1,2,4-Trimethylbenzene	ND		ug/l	2.5	0.70
Methyl Acetate	ND		ug/l	2.0	0.23
Cyclohexane	ND		ug/l	10	0.27
Freon-113	ND		ug/l	2.5	0.70
Methyl cyclohexane	ND		ug/l	10	0.40

Project Name: CONVENTUS/MROW
Project Number: 1186

Lab Number: L2270805
Report Date: 12/29/22

Method Blank Analysis
Batch Quality Control

Analytical Method: 1,8260D
 Analytical Date: 12/23/22 08:59
 Analyst: PID

Parameter	Result	Qualifier	Units	RL	MDL
Volatile Organics by GC/MS - Westborough Lab for sample(s): 08 Batch: WG1727559-5					

Surrogate	%Recovery	Qualifier	Acceptance Criteria
1,2-Dichloroethane-d4	103		70-130
Toluene-d8	104		70-130
4-Bromofluorobenzene	113		70-130
Dibromofluoromethane	97		70-130

Lab Control Sample Analysis **Batch Quality Control**

Project Name: CONVENTUS/MROW

Project Number: 1186

Lab Number: L2270805

Report Date: 12/29/22

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-02,04-07,09 Batch: WG1726668-3 WG1726668-4								
Methylene chloride	100		110		70-130	10		20
1,1-Dichloroethane	100		100		70-130	0		20
Chloroform	93		86		70-130	8		20
Carbon tetrachloride	81		85		63-132	5		20
1,2-Dichloropropane	95		100		70-130	5		20
Dibromochloromethane	73		75		63-130	3		20
1,1,2-Trichloroethane	78		85		70-130	9		20
Tetrachloroethene	89		88		70-130	1		20
Chlorobenzene	92		93		75-130	1		20
Trichlorofluoromethane	97		97		62-150	0		20
1,2-Dichloroethane	79		82		70-130	4		20
1,1,1-Trichloroethane	86		87		67-130	1		20
Bromodichloromethane	82		84		67-130	2		20
trans-1,3-Dichloropropene	77		79		70-130	3		20
cis-1,3-Dichloropropene	84		91		70-130	8		20
Bromoform	72		80		54-136	11		20
1,1,2,2-Tetrachloroethane	91		98		67-130	7		20
Benzene	96		98		70-130	2		20
Toluene	90		91		70-130	1		20
Ethylbenzene	96		98		70-130	2		20
Chloromethane	110		110		64-130	0		20
Bromomethane	67		73		39-139	9		20
Vinyl chloride	110		120		55-140	9		20

Lab Control Sample Analysis

Batch Quality Control

Project Name: CONVENTUS/MROW

Project Number: 1186

Lab Number: L2270805

Report Date: 12/29/22

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-02,04-07,09 Batch: WG1726668-3 WG1726668-4								
Chloroethane	120		120		55-138	0		20
1,1-Dichloroethene	100		100		61-145	0		20
trans-1,2-Dichloroethene	99		99		70-130	0		20
Trichloroethene	85		91		70-130	7		20
1,2-Dichlorobenzene	96		98		70-130	2		20
1,3-Dichlorobenzene	99		100		70-130	1		20
1,4-Dichlorobenzene	95		99		70-130	4		20
Methyl tert butyl ether	76		83		63-130	9		20
p/m-Xylene	100		100		70-130	0		20
o-Xylene	100		100		70-130	0		20
cis-1,2-Dichloroethene	94		94		70-130	0		20
Styrene	95		100		70-130	5		20
Dichlorodifluoromethane	110		120		36-147	9		20
Acetone	100		92		58-148	8		20
Carbon disulfide	110		110		51-130	0		20
2-Butanone	82		89		63-138	8		20
4-Methyl-2-pentanone	87		91		59-130	4		20
2-Hexanone	89		110		57-130	21	Q	20
1,2-Dibromoethane	80		85		70-130	6		20
n-Butylbenzene	110		110		53-136	0		20
sec-Butylbenzene	110		110		70-130	0		20
tert-Butylbenzene	100		100		70-130	0		20
1,2-Dibromo-3-chloropropane	85		90		41-144	6		20

Lab Control Sample Analysis

Batch Quality Control

Project Name: CONVENTUS/MROW

Project Number: 1186

Lab Number: L2270805

Report Date: 12/29/22

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-02,04-07,09 Batch: WG1726668-3 WG1726668-4								
Isopropylbenzene	100		100		70-130	0		20
p-Isopropyltoluene	100		100		70-130	0		20
Naphthalene	89		96		70-130	8		20
n-Propylbenzene	110		110		69-130	0		20
1,2,4-Trichlorobenzene	83		89		70-130	7		20
1,3,5-Trimethylbenzene	98		100		64-130	2		20
1,2,4-Trimethylbenzene	95		99		70-130	4		20
Methyl Acetate	99		110		70-130	11		20
Cyclohexane	110		110		70-130	0		20
Freon-113	110		110		70-130	0		20
Methyl cyclohexane	98		99		70-130	1		20

Surrogate	LCS %Recovery	Qual	LCSD %Recovery	Qual	Acceptance Criteria
1,2-Dichloroethane-d4	90		92		70-130
Toluene-d8	102		101		70-130
4-Bromofluorobenzene	105		104		70-130
Dibromofluoromethane	99		100		70-130

Lab Control Sample Analysis

Batch Quality Control

Project Name: CONVENTUS/MROW

Lab Number: L2270805

Project Number: 1186

Report Date: 12/29/22

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 03 Batch: WG1726975-3 WG1726975-4								
Methylene chloride	100		96		70-130	4		20
1,1-Dichloroethane	100		95		70-130	5		20
Chloroform	99		94		70-130	5		20
Carbon tetrachloride	98		91		63-132	7		20
1,2-Dichloropropane	99		93		70-130	6		20
Dibromochloromethane	88		84		63-130	5		20
1,1,2-Trichloroethane	90		91		70-130	1		20
Tetrachloroethene	89		85		70-130	5		20
Chlorobenzene	92		89		75-130	3		20
Trichlorofluoromethane	110		98		62-150	12		20
1,2-Dichloroethane	100		97		70-130	3		20
1,1,1-Trichloroethane	100		93		67-130	7		20
Bromodichloromethane	97		92		67-130	5		20
trans-1,3-Dichloropropene	91		88		70-130	3		20
cis-1,3-Dichloropropene	94		90		70-130	4		20
Bromoform	77		76		54-136	1		20
1,1,2,2-Tetrachloroethane	84		88		67-130	5		20
Benzene	97		92		70-130	5		20
Toluene	92		90		70-130	2		20
Ethylbenzene	96		92		70-130	4		20
Chloromethane	110		97		64-130	13		20
Bromomethane	110		100		39-139	10		20
Vinyl chloride	110		96		55-140	14		20

Lab Control Sample Analysis **Batch Quality Control**

Project Name: CONVENTUS/MROW

Project Number: 1186

Lab Number: L2270805

Report Date: 12/29/22

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 03 Batch: WG1726975-3 WG1726975-4								
Chloroethane	100		91		55-138	9		20
1,1-Dichloroethene	100		93		61-145	7		20
trans-1,2-Dichloroethene	100		96		70-130	4		20
Trichloroethene	96		92		70-130	4		20
1,2-Dichlorobenzene	90		89		70-130	1		20
1,3-Dichlorobenzene	92		88		70-130	4		20
1,4-Dichlorobenzene	92		89		70-130	3		20
Methyl tert butyl ether	99		93		63-130	6		20
p/m-Xylene	95		90		70-130	5		20
o-Xylene	95		90		70-130	5		20
cis-1,2-Dichloroethene	100		92		70-130	8		20
Styrene	95		90		70-130	5		20
Dichlorodifluoromethane	110		100		36-147	10		20
Acetone	100		88		58-148	13		20
Carbon disulfide	110		97		51-130	13		20
2-Butanone	100		100		63-138	0		20
4-Methyl-2-pentanone	84		91		59-130	8		20
2-Hexanone	92		92		57-130	0		20
1,2-Dibromoethane	88		86		70-130	2		20
n-Butylbenzene	91		90		53-136	1		20
sec-Butylbenzene	94		91		70-130	3		20
tert-Butylbenzene	92		88		70-130	4		20
1,2-Dibromo-3-chloropropane	88		80		41-144	10		20

Lab Control Sample Analysis

Batch Quality Control

Project Name: CONVENTUS/MROW

Project Number: 1186

Lab Number: L2270805

Report Date: 12/29/22

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 03 Batch: WG1726975-3 WG1726975-4								
Isopropylbenzene	93		89		70-130	4		20
p-Isopropyltoluene	91		88		70-130	3		20
Naphthalene	85		85		70-130	0		20
n-Propylbenzene	95		90		69-130	5		20
1,2,4-Trichlorobenzene	84		83		70-130	1		20
1,3,5-Trimethylbenzene	89		86		64-130	3		20
1,2,4-Trimethylbenzene	88		86		70-130	2		20
Methyl Acetate	110		110		70-130	0		20
Cyclohexane	110		99		70-130	11		20
Freon-113	110		98		70-130	12		20
Methyl cyclohexane	96		89		70-130	8		20

Surrogate	LCS %Recovery	Qual	LCSD %Recovery	Qual	Acceptance Criteria
1,2-Dichloroethane-d4	109		105		70-130
Toluene-d8	99		101		70-130
4-Bromofluorobenzene	100		102		70-130
Dibromofluoromethane	104		101		70-130

Lab Control Sample Analysis

Batch Quality Control

Project Name: CONVENTUS/MROW

Project Number: 1186

Lab Number: L2270805

Report Date: 12/29/22

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 08 Batch: WG1727559-3 WG1727559-4								
Methylene chloride	98		99		70-130	1		20
1,1-Dichloroethane	110		100		70-130	10		20
Chloroform	100		100		70-130	0		20
Carbon tetrachloride	110		110		63-132	0		20
1,2-Dichloropropane	110		100		70-130	10		20
Dibromochloromethane	100		100		63-130	0		20
1,1,2-Trichloroethane	100		99		70-130	1		20
Tetrachloroethene	100		98		70-130	2		20
Chlorobenzene	98		97		75-130	1		20
Trichlorofluoromethane	110		110		62-150	0		20
1,2-Dichloroethane	110		100		70-130	10		20
1,1,1-Trichloroethane	110		110		67-130	0		20
Bromodichloromethane	110		100		67-130	10		20
trans-1,3-Dichloropropene	100		100		70-130	0		20
cis-1,3-Dichloropropene	100		100		70-130	0		20
Bromoform	100		100		54-136	0		20
1,1,2,2-Tetrachloroethane	100		100		67-130	0		20
Benzene	97		96		70-130	1		20
Toluene	100		100		70-130	0		20
Ethylbenzene	100		99		70-130	1		20
Chloromethane	96		94		64-130	2		20
Bromomethane	70		68		39-139	3		20
Vinyl chloride	110		100		55-140	10		20

Lab Control Sample Analysis **Batch Quality Control**

Project Name: CONVENTUS/MROW

Project Number: 1186

Lab Number: L2270805

Report Date: 12/29/22

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 08 Batch: WG1727559-3 WG1727559-4								
Chloroethane	110		110		55-138	0		20
1,1-Dichloroethene	100		100		61-145	0		20
trans-1,2-Dichloroethene	100		98		70-130	2		20
Trichloroethene	100		99		70-130	1		20
1,2-Dichlorobenzene	100		100		70-130	0		20
1,3-Dichlorobenzene	100		100		70-130	0		20
1,4-Dichlorobenzene	99		99		70-130	0		20
Methyl tert butyl ether	100		100		63-130	0		20
p/m-Xylene	100		100		70-130	0		20
o-Xylene	95		95		70-130	0		20
cis-1,2-Dichloroethene	100		100		70-130	0		20
Styrene	85		85		70-130	0		20
Dichlorodifluoromethane	110		100		36-147	10		20
Acetone	100		110		58-148	10		20
Carbon disulfide	100		99		51-130	1		20
2-Butanone	110		110		63-138	0		20
4-Methyl-2-pentanone	100		100		59-130	0		20
2-Hexanone	100		100		57-130	0		20
1,2-Dibromoethane	100		100		70-130	0		20
n-Butylbenzene	110		110		53-136	0		20
sec-Butylbenzene	110		100		70-130	10		20
tert-Butylbenzene	100		100		70-130	0		20
1,2-Dibromo-3-chloropropane	91		94		41-144	3		20

Lab Control Sample Analysis

Batch Quality Control

Project Name: CONVENTUS/MROW

Project Number: 1186

Lab Number: L2270805

Report Date: 12/29/22

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 08 Batch: WG1727559-3 WG1727559-4								
Isopropylbenzene	110		100		70-130	10		20
p-Isopropyltoluene	100		100		70-130	0		20
Naphthalene	100		110		70-130	10		20
n-Propylbenzene	100		100		69-130	0		20
1,2,4-Trichlorobenzene	100		100		70-130	0		20
1,3,5-Trimethylbenzene	100		100		64-130	0		20
1,2,4-Trimethylbenzene	100		100		70-130	0		20
Methyl Acetate	110		100		70-130	10		20
Cyclohexane	100		100		70-130	0		20
Freon-113	110		110		70-130	0		20
Methyl cyclohexane	100		99		70-130	1		20

Surrogate	LCS %Recovery	Qual	LCSD %Recovery	Qual	Acceptance Criteria
1,2-Dichloroethane-d4	101		100		70-130
Toluene-d8	101		100		70-130
4-Bromofluorobenzene	109		110		70-130
Dibromofluoromethane	99		100		70-130

Project Name: CONVENTUS/MROW**Lab Number:** L2270805**Project Number:** 1186**Report Date:** 12/29/22**Sample Receipt and Container Information**

Were project specific reporting limits specified?

YES

Cooler Information

Cooler	Custody Seal
A	Absent

Container Information

Container ID	Container Type	Cooler	Initial pH	Final pH	Temp deg C	Pres	Seal	Frozen Date/Time	Analysis(*)
L2270805-01A	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2270805-01B	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2270805-01C	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2270805-02A	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2270805-02B	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2270805-02C	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2270805-03A	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2270805-03B	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2270805-03C	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2270805-04A	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2270805-04B	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2270805-04C	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2270805-05A	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2270805-05B	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2270805-05C	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2270805-06A	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2270805-06B	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2270805-06C	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2270805-07A	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2270805-07B	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2270805-07C	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2270805-08A	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2270805-08B	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)

Project Name: CONVENTUS/MROW
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Container Information

Container ID	Container Type	Cooler	Initial pH	Final pH	Temp deg C	Pres	Seal	Frozen Date/Time	Analysis(*)
L2270805-09A	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2270805-09B	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2270805-09C	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)

Project Name: CONVENTUS/MROW
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GLOSSARY

Acronyms

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.)
EDL	- 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.
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.)
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 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 only.)
MDL	- 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.
MS	- 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.

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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.

Chlordane: 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.)

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'.

Gasoline Range Organics (GRO): Gasoline Range Organics (GRO) results include all chromatographic peaks eluting from Methyl tert butyl ether through Naphthalene, with the exception of GRO analysis in support of State of Ohio programs, which includes all chromatographic peaks eluting from Hexane through Dodecane.

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.

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.
- B** - 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).
- C** - Co-elution: The target analyte co-elutes with a known lab standard (i.e. surrogate, internal standards, etc.) for co-extracted analyses.
- D** - 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.
- J** - 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

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Data Qualifiers

Identified Compounds (TICs).

- M** - 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.
- 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.
- V** - The surrogate associated with this target analyte has a recovery outside the QC acceptance limits. (Applicable to MassDEP DW Compliance samples only.)
- Z** - 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: CONVENTUS/MROW
Project Number: 1186

Lab Number: L2270805
Report Date: 12/29/22

REFERENCES

- 1 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**

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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; 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 I, 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.

 NEW YORK CHAIN OF CUSTODY Westborough, MA 01581 8 Walkup Dr. TEL: 508-898-9220 FAX: 508-898-9193 Mansfield, MA 02048 320 Forbes Blvd TEL: 508-822-9300 FAX: 508-822-3288		Service Centers Mahwah, NJ 07430: 35 Whitney Rd, Suite 5 Albany, NY 12205: 14 Walker Way Tonawanda, NY 14150: 275 Cooper Ave, Suite 105		Page 1 of 1		Date Rec'd in Lab 12/16/22		ALPHA Job # L2270805	
		Project Information: Project Name: <u>CONVENTS / MROW</u> Project Location: <u>CONVENTS / MAIN ST.</u> Project # <u>U8L</u> (Use Project name as Project #) ☐		Deliverables ☐ ASP-A ☒ ASP-B ☒ EQuIS (1 File) ☐ EQuIS (4 File) ☐ Other		Billing Information ☐ Same as Client Info PO #			
Client Information Client: <u>COS ENGINEERS</u> Address: <u>141 ELDT ST.</u> <u>BUFFALO NY 14203</u> Phone: <u>716-796-3520</u> Fax: Email: <u>rbackett@cosos.com</u>		Project Manager: <u>ALPHA QUOTE #:</u> <u>ALPHA QUOTE #:</u> Turn-Around Time Standard ☒ Due Date: Rush (only if pre approved) ☐ # of Days:		Regulatory Requirement ☐ NY TOGS ☒ NY Part 375 ☐ AWQ Standards ☐ NY CP-51 ☐ NY Restricted Use ☐ Other ☐ NY Unrestricted Use ☐ NYC Sewer Discharge		Disposal Site Information Please identify below location of applicable disposal facilities. Disposal Facility: ☐ NJ ☒ NY ☐ Other:			
These samples have been previously analyzed by Alpha ☐ Other project specific requirements/comments:						ANALYSIS		Sample Filtration ☐ Done ☐ Lab to do Preservation ☐ Lab to do (Please Specify below)	
Please specify Metals or TAL.						VOCs 8200		Total Bottles	
ALPHA Lab ID (Lab Use Only)	Sample ID	Collection Date Time		Sample Matrix	Sampler's Initials				
70805-01	Bcp MW-1	12/14/22	11:00	GLW	RB3	x			
02	Dup	12/14/22	11:00	GLW	RB3	x			
03	Bcp MW-7	12/14/22	11:40	GLW	RB3	x			
04	Bcp MW-4	12/14/22	12:25	GLW	RB3	x			
05	Bcp MW-3	12/14/22	12:55	GLW	RB3	x			
06	Bcp MW-6	12/14/22	1:30	GLW	RB3	x			
07	Bcp MW-5	12/14/22	2:10	GLW	RB3	x			
08	TRIP BLANK	12/15/22	11:55	GLW	RB3	x			
09	MSHW-2	12/15/22	10:45	GLW	RB3	x			
Preservative Code:		Container Code		Westboro: Certification No: MA935		Container Type			
A = None		P = Plastic		Mansfield: Certification No: MA015		Preservative			
B = HCl		A = Amber Glass							
C = HNO ₃		V = Vial							
D = H ₂ SO ₄		G = Glass							
E = NaOH		B = Bacteria Cup							
F = MeOH		C = Cube							
G = NaHSO ₄		O = Other							
H = Na ₂ S ₂ O ₃		E = Encore							
K/E = Zn Ac/NaOH		D = BOD Bottle							
O = Other									
Form No: 01-25 HC (rev. 30-Sept-2013)		Relinquished By: <u>9 Ryt AAL</u>		Date/Time: <u>12/15/22 1320</u>		Received By: <u>9 Ryt AAL</u>		Date/Time: <u>12/15/22 1320</u>	

Please print clearly, legibly and completely. Samples can not be logged in and turnaround time clock will not start until any ambiguities are resolved. BY EXECUTING THIS COC, THE CLIENT HAS READ AND AGREES TO BE BOUND BY ALPHA'S TERMS & CONDITIONS. (See reverse side.)

DATA USABILITY SUMMARY REPORT (DUSR)

**Conventus/MROW
Buffalo, NY
Project # U86001005**

SDG: L2270805
8 Water Samples and 1 Trip Blank

Prepared for:

**C&S Companies
141 Elm Street, Suite 100
Buffalo, NY 14203
Attention: Cody Martin**

April 2023



Environmental Data Usability 10028 Deer Park Dr. Dansville, NY 14437 585-991-9156

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REVIEWER'S NARRATIVE

C&S Companies SDG L2270805 Conventus/MROW

The data associated with this Sample Delivery Groups (SDG) L2270805, analyzed by Alpha Analytical, Westborough, MA have been reviewed in accordance with assessment criteria provided by the New York State Department of Environmental Conservation following the review procedures provided in the USEPA Functional Guidelines for evaluating organic and inorganic data.

All analytical results reported by the laboratory are considered valid and acceptable except results that have been qualified as rejected, "R". Results qualified as estimated "J", or as non-detects, "U", are considered usable for the purpose of evaluating water and/or soil quality. However, these qualifiers indicate that the accuracy and/or precision of the analytical result is questionable. A summary of all data that have been qualified and the reasons for qualification are provided in the following data usability summary report (DUSR).

Two facts should be noted by all data users. First, the "R" qualifier means that the associated value is unusable. In other words, due to significant quality control (QC) problems, the analysis is invalid and provides no information as to whether the analyte is present or not. Values qualified with an "R" should not appear on the final data tables because they cannot be relied upon, even as the last resort. Second, no analyte concentration, even if it has passed all QC tests, is guaranteed to be accurate. Strict QC serves to increase confidence in data, but any value potentially contains error.

Reviewer's Signature: Michael K. Perry Date: 4/7/2023
Michael K. Perry
Chemist

1.0 EVENT SUMMARY

SITE: Conventus
Buffalo, NY
Project #: U86001005

SAMPLING DATES: December 14 and 15, 2022

SAMPLE TYPE: 8 water samples and 1 trip blank

LABORATORY: Alpha Analytical
Westborough, MA

SDG No.: SDG L2270805

2.0 INTRODUCTION

This data usability summary report (DUSR) was prepared in accordance with guidance provided by the New York State Department of Environmental Conservation (NYSDEC). The DUSR is based on a review and evaluation of the laboratory analytical data package. Specifically, the NYSDEC guidance recommends review and evaluation of the following elements of the data package:

- Completeness of the data package as defined under the requirements of the NYSDEC Analytical Services Protocols (ASP) Category B or the United States Environmental Protection Agency (USEPA) Contract Laboratory Program (CLP) deliverables,
- Compliance with established analyte holding times,
- Adherence to quality control (QC) limits and specifications for blanks, instrument tuning and calibration, surrogate recoveries, spike recoveries, laboratory duplicate analyses, and other QC criteria,
- Adherence to established analytical protocols,
- Conformance of data summary sheets with raw analytical data, and
- Use of correct data qualifiers.

Data deficiencies, analytical protocol deviations, and quality control problems identified using the review criteria above and their effect on the analytical results are discussed in this report.

3.0 SAMPLE AND ANALYSIS SUMMARY

The data package consists of analytical results for 8 water samples and 1 trip blank collected on 12/14/22 and 12/15/22. These samples were analyzed for Volatile Organic Compounds (VOCs).

All laboratory analyses were submitted to Alpha Analytical, Westborough, MA and analyzed as SDG L2270805. The analytical results were provided in NYSDEC ASP Category B format, which includes all raw analytical data and laboratory QC data.

4.0 GUIDANCE DOCUMENTS AND DATA REVIEW CRITERIA

The guidance documents appropriate for reviewing laboratory quality control (QC) data and assigning data qualifiers (flags) to analytical results were selected from those listed in Table 4-1. The QC limits established in the documents applicable to this data review were used to assess the quality of the analytical results. In some cases, however, QC limits established internally by the laboratory were taken into account to determine data quality.

The QC criteria considered for assessing the usability of the reported analytical results provided for each analyte type (i.e. VOCs, SVOCs, metals, etc.) are listed in Table 4-2. These criteria may vary with the analytical method utilized by the laboratory. These criteria comply with the guidance recommended in Section 2.0 above.

5.0 DATA VALIDATION QUALIFIERS

The letter qualifiers (flags) used to define data usability are described briefly below. These letters are assigned by the data validator to analytical results having questionable accuracy and/or precision as determined by reviewing the laboratory QC data associated with the analytical results.

TABLE 4-1**Guidance Used For Validating Laboratory Analytical Data**

Analyte Group	Guidance	Date
Metals (ICP-AES)	USEPA SOP HW-3a, Rev. 1	September 2016
Metals (Hg & CN)	USEPA SOP HW-3c, Rev. 1	September 2016
Volatile Organic Compounds (by Methods 8260B & 8260C)	USEPA SOP HW-24, Rev. 4	September 2014
Semi-Volatile Organic Compounds (by Method 8270D)	USEPA SOP HW-22 Rev. 5	December 2010
Pesticides (by Method 8181B)	USEPA SOP HW-44, Rev. 1.1	December 2010
Chlorinated Herbicides (by Method 8151A)	USEPA SOP HW-17, Rev. 3.1	December 2010
Polychlorinated Biphenyls (PCBs)	USEPA SOP HW-37A, Rev. 0	June 2015
Volatile Organic Compounds (Air) (by Method TO-15)	USEPA SOP HW-31, Rev. 6	September 2016
Per- and PolyFluoroAlkyl Substances (PFAS)	* NYSDEC	January 2021
General Chemistry Parameters	per NYSDEC ASP	July 2005

* Sampling, Analysis, and Assessment of Per- and Polyfluoroalkyl Substances (PFAS) Under NYSDEC's Part 375 Remedial Programs, Appendix I

TABLE 4-2

**QUALITY CONTROL CRITERIA USED FOR VALIDATING
LABORATORY ANALYTICAL DATA**

VOCs	SVOCs	Pesticides/PCBs	Metals	Gen Chemistry	PFAS
Completeness of Pkg Sample Preservation Holding Time System Monitoring Compounds Lab Control Sample Matrix Spikes Blanks Instrument Tuning Internal Standards Initial Calibration Continuing Calibration Lab Qualifiers Field Duplicate	Completeness of Pkg Sample Preservation Holding Time Surrogate Recoveries Lab Control Sample Matrix Spikes Blanks Instrument Tuning Internal Standards Initial Calibration Continuing Calibration Lab Qualifiers Field Duplicate	Completeness of Pkg Sample Preservation Holding Time Surrogate Recoveries Matrix Spikes Blanks Instrument Calibration & Verification Comparison of duplicate GC column results Analyte ID Lab Qualifiers Field Duplicate	Completeness of Pkg Sample Preservation Holding Time Initial/Continuing Calibration CRDL Standards Blanks Interference Check Sample Spike Recoveries Lab Duplicate Lab Control Sample ICP Serial Dilutions Lab Qualifiers Field Duplicate	Completeness of Pkg Sample Preservation Holding Times Calibration Lab Control Samples Blanks Spike Recoveries Lab Duplicates	Completeness of Pkg Sample Preservation Holding Time Instr Performance Check Initial Calibration Continuing Calibration Blanks Surrogates Lab Fortified Blank Matrix Spikes Internal Standards

Method TO-15 (Air)

Completeness of Pkg
 Sample Preservation
 Holding Time
 Canister Certification
 Instrument Tuning
 Initial Calibration and
 Instrument Performance
 Daily Calibration
 Blanks
 Lab Control Sample
 Field Duplicate

The laboratory may also use various letters and symbols to flag analytical results generated when QC limits were exceeded. The meanings of these flags may differ from those used by the independent data validator. Those used by the laboratory are provided with the analytical results.

NOTE: The assignment of data qualifiers by the data reviewer (validator) to laboratory analytical results should not necessarily be interpreted by the data user as a measure of laboratory ability or proficiency. Rather, the qualifiers are intended to provide a measure of data accuracy and precision to the data user, which, for example, may provide a level of confidence in determining whether or not standards or cleanup objectives have been met.

- U** The analyte was analyzed for but was not detected at or above the sample quantitation limit.
- J** The analyte was positively identified; the associated numerical value is the *approximate* concentration of the analyte in the sample. (The magnitude of any $\pm$ value associated with the result is not determined by data validation).
- J+** The result is an estimated quantity and may be biased high.
- J-** The result is an estimated quantity and may be biased low.
- UJ** The analyte was analyzed for but not detected. The reported quantitation limit is approximate and may be inaccurate or imprecise.
- R** The sample result is rejected (i.e., is unusable) due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.
- NJ** The analysis indicates the presence of an analyte that has been “tentatively identified” and the associated numerical value represents its approximate concentration.

The validated analytical results are attached to this report. Validation qualifiers (flags) are indicated in red print. Data sheets having qualified data are signed and dated by the data reviewer.

6.0 RESULTS OF THE DATA REVIEW

The results of the data review are summarized in Table 6-1. The table lists the samples where QC criteria were found to exceed acceptable limits and the actions taken to qualify the associated analytical results.

7.0 TOTAL USABLE DATA

For SDG L2270805, nine samples were analyzed and results were reported for 522 analytes. Even though some results were flagged with a “J” as estimated, all results (100 %) are considered usable. See the summary table for the analyses that have been rejected and qualified and the associated QC reasons.

Table 6-1 VOCs

SAMPLES AFFECTED	ANALYTES	ACTION	QC VIOLATION	COMMENTS
BCPMW-1 Dup BCPMW-4 BCPMW-3 BCPMW-6 BCPMW-5 MSMW-2	Bromomethane Chloromethane Carbon disulfide Dichlorodifluoromethane Methyl tert butyl ether 1,2-Dichloroethane trans-1,3-Dichloropropene 1,1,2-Trichloroethane Chlorodibromomethane 1,2-dibromoethane	J detects UJ non-detects	ICV and/or CCV > QC limit	Data are estimated
BCPMW-7	Carbon disulfide	J detects UJ non-detects	ICV > QC limit	Data are estimated
Trip blank	Bromomethane Carbon disulfide	J detects UJ non-detects	ICV and/or CCV > QC limit	Data are estimated

ACRONYMS

BSP	Blank Spike
CCAL	Continuing Calibration
CCB	Continuing Calibration Blank
CCV	Continuing Calibration Verification
CRDL	Contract Required Detection Limit
CRQL	Contract Required Quantitation Limit
%D	Percent Difference
ICAL	Initial Calibration
ICB	Initial Calibration Blank
IS	Internal Standard
LCS	Laboratory Control Sample
MS/MSD	Matrix Spike/Matrix Spike Duplicate
QA	Quality Assurance
QC	Quality Control
%R	Percent recovery
RPD	Relative Percent Difference
RRF	Relative Response Factor
%RSD	Percent Relative Standard Deviation
TAL	Target Analyte List (metals)
TCL	Target Compound List (organics)

Appendix A

Validated Analytical Results



www.alphalab.com



Alpha Analytical

Laboratory Code: 11148

SDG Number: L2270805

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.

Project Name: CONVENTUS/MROW
Project Number: 1186

Lab Number: L2270805
Report Date: 12/29/22

Alpha Sample ID	Client ID	Matrix	Sample Location	Collection Date/Time	Receive Date
L2270805-01	BCPMW-1	WATER	CONVENTUS 1 MAIN ST. BUFFALO NY	12/14/22 11:00	12/15/22
L2270805-02	DUP	WATER	CONVENTUS 1 MAIN ST. BUFFALO NY	12/14/22 11:00	12/15/22
L2270805-03	BCPMW-7	WATER	CONVENTUS 1 MAIN ST. BUFFALO NY	12/14/22 11:40	12/15/22
L2270805-04	BCPMW-4	WATER	CONVENTUS 1 MAIN ST. BUFFALO NY	12/14/22 12:25	12/15/22
L2270805-05	BCPMW-3	WATER	CONVENTUS 1 MAIN ST. BUFFALO NY	12/14/22 12:55	12/15/22
L2270805-06	BCPMW-6	WATER	CONVENTUS 1 MAIN ST. BUFFALO NY	12/14/22 13:30	12/15/22
L2270805-07	BCPMW-5	WATER	CONVENTUS 1 MAIN ST. BUFFALO NY	12/14/22 14:10	12/15/22
L2270805-08	TRIP BLANK	WATER	CONVENTUS 1 MAIN ST. BUFFALO NY	12/15/22 00:00	12/15/22
L2270805-09	MSMW-2	WATER	CONVENTUS 1 MAIN ST. BUFFALO NY	12/15/22 10:45	12/15/22

Project Name: CONVENTUS/MROW
Project Number: 1186

Lab Number: L2270805
Report Date: 12/29/22

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.

Volatile Organics

L2270805-07D: The pH was greater than two; however, the sample was analyzed within the method required holding time.

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: *Siffani Morrissey*

Report Date: 12/29/22

Title: Technical Director/Representative



GC/MS 8260

Analysis

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies	Lab Number : L2270805
Project Name : CONVENTUS/MROW	Project Number : 1186
Lab ID : L2270805-01	Date Collected : 12/14/22 11:00
Client ID : BCPMW-1	Date Received : 12/15/22
Sample Location : CONVENTUS 1 MAIN ST. BUFFALO NY	Date Analyzed : 12/21/22 14:10
Sample Matrix : WATER	Dilution Factor : 1
Analytical Method : 1,8260D	Analyst : MJV
Lab File ID : V22221221A18	Instrument ID : VOA122
Sample Amount : 10 ml	GC Column : RTX-502.2
Level : LOW	%Solids : N/A
Extract Volume (MeOH) : N/A	Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
75-09-2	Methylene chloride	ND	2.5	0.70	U
75-34-3	1,1-Dichloroethane	ND	2.5	0.70	U
67-66-3	Chloroform	ND	2.5	0.70	U
56-23-5	Carbon tetrachloride	ND	0.50	0.13	U
78-87-5	1,2-Dichloropropane	ND	1.0	0.14	U
124-48-1	Dibromochloromethane	ND	0.50	0.15	U
79-00-5	1,1,2-Trichloroethane	ND	1.5	0.50	U UJ
127-18-4	Tetrachloroethene	ND	0.50	0.18	U UJ
108-90-7	Chlorobenzene	ND	2.5	0.70	U
75-69-4	Trichlorofluoromethane	ND	2.5	0.70	U
107-06-2	1,2-Dichloroethane	ND	0.50	0.13	U
71-55-6	1,1,1-Trichloroethane	ND	2.5	0.70	U UJ
75-27-4	Bromodichloromethane	ND	0.50	0.19	U
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.16	U
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.14	U UJ
75-25-2	Bromoform	ND	2.0	0.65	U
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.17	U UJ
71-43-2	Benzene	ND	0.50	0.16	U
108-88-3	Toluene	ND	2.5	0.70	U
100-41-4	Ethylbenzene	ND	2.5	0.70	U
74-87-3	Chloromethane	ND	2.5	0.70	U
74-83-9	Bromomethane	ND	2.5	0.70	U UJ
75-01-4	Vinyl chloride	ND	1.0	0.07	U UJ
75-00-3	Chloroethane	ND	2.5	0.70	U
75-35-4	1,1-Dichloroethene	ND	0.50	0.17	U

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies	Lab Number : L2270805
Project Name : CONVENTUS/MROW	Project Number : 1186
Lab ID : L2270805-01	Date Collected : 12/14/22 11:00
Client ID : BCPMW-1	Date Received : 12/15/22
Sample Location : CONVENTUS 1 MAIN ST. BUFFALO NY	Date Analyzed : 12/21/22 14:10
Sample Matrix : WATER	Dilution Factor : 1
Analytical Method : 1,8260D	Analyst : MJV
Lab File ID : V22221221A18	Instrument ID : VOA122
Sample Amount : 10 ml	GC Column : RTX-502.2
Level : LOW	%Solids : N/A
Extract Volume (MeOH) : N/A	Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
156-60-5	trans-1,2-Dichloroethene	ND	2.5	0.70	U
79-01-6	Trichloroethene	ND	0.50	0.18	U
95-50-1	1,2-Dichlorobenzene	ND	2.5	0.70	U
541-73-1	1,3-Dichlorobenzene	ND	2.5	0.70	U
106-46-7	1,4-Dichlorobenzene	ND	2.5	0.70	U
1634-04-4	Methyl tert butyl ether	ND	2.5	0.70	U
179601-23-1	p/m-Xylene	ND	2.5	0.70	U UJ
95-47-6	o-Xylene	ND	2.5	0.70	U
156-59-2	cis-1,2-Dichloroethene	ND	2.5	0.70	U
100-42-5	Styrene	ND	2.5	0.70	U
75-71-8	Dichlorodifluoromethane	ND	5.0	1.0	U
67-64-1	Acetone	ND	5.0	1.5	U UJ
75-15-0	Carbon disulfide	ND	5.0	1.0	U
78-93-3	2-Butanone	ND	5.0	1.9	U UJ
108-10-1	4-Methyl-2-pentanone	ND	5.0	1.0	U
591-78-6	2-Hexanone	ND	5.0	1.0	U
106-93-4	1,2-Dibromoethane	ND	2.0	0.65	U
104-51-8	n-Butylbenzene	ND	2.5	0.70	U UJ
135-98-8	sec-Butylbenzene	ND	2.5	0.70	U
98-06-6	tert-Butylbenzene	ND	2.5	0.70	U
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.5	0.70	U
98-82-8	Isopropylbenzene	ND	2.5	0.70	U
99-87-6	p-Isopropyltoluene	ND	2.5	0.70	U
91-20-3	Naphthalene	ND	2.5	0.70	U
103-65-1	n-Propylbenzene	ND	2.5	0.70	U

Results Summary

Form 1

Volatile Organics by GC/MS

Client	: C&S Companies	Lab Number	: L2270805
Project Name	: CONVENTUS/MROW	Project Number	: 1186
Lab ID	: L2270805-01	Date Collected	: 12/14/22 11:00
Client ID	: BCPMW-1	Date Received	: 12/15/22
Sample Location	: CONVENTUS 1 MAIN ST. BUFFALO NY	Date Analyzed	: 12/21/22 14:10
Sample Matrix	: WATER	Dilution Factor	: 1
Analytical Method	: 1,8260D	Analyst	: MJV
Lab File ID	: V22221221A18	Instrument ID	: VOA122
Sample Amount	: 10 ml	GC Column	: RTX-502.2
Level	: LOW	%Solids	: N/A
Extract Volume (MeOH)	: N/A	Injection Volume	: N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
120-82-1	1,2,4-Trichlorobenzene	ND	2.5	0.70	U
108-67-8	1,3,5-Trimethylbenzene	ND	2.5	0.70	U
95-63-6	1,2,4-Trimethylbenzene	ND	2.5	0.70	U
79-20-9	Methyl Acetate	ND	2.0	0.23	U
110-82-7	Cyclohexane	ND	10	0.27	U
76-13-1	Freon-113	ND	2.5	0.70	U
108-87-2	Methyl cyclohexane	ND	10	0.40	U

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies	Lab Number : L2270805
Project Name : CONVENTUS/MROW	Project Number : 1186
Lab ID : L2270805-02	Date Collected : 12/14/22 11:00
Client ID : DUP	Date Received : 12/15/22
Sample Location : CONVENTUS 1 MAIN ST. BUFFALO NY	Date Analyzed : 12/21/22 14:34
Sample Matrix : WATER	Dilution Factor : 1
Analytical Method : 1,8260D	Analyst : MJV
Lab File ID : V22221221A19	Instrument ID : VOA122
Sample Amount : 10 ml	GC Column : RTX-502.2
Level : LOW	%Solids : N/A
Extract Volume (MeOH) : N/A	Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
75-09-2	Methylene chloride	ND	2.5	0.70	U
75-34-3	1,1-Dichloroethane	ND	2.5	0.70	U
67-66-3	Chloroform	ND	2.5	0.70	U
56-23-5	Carbon tetrachloride	ND	0.50	0.13	U
78-87-5	1,2-Dichloropropane	ND	1.0	0.14	U
124-48-1	Dibromochloromethane	ND	0.50	0.15	U
79-00-5	1,1,2-Trichloroethane	ND	1.5	0.50	U UJ
127-18-4	Tetrachloroethene	ND	0.50	0.18	U UJ
108-90-7	Chlorobenzene	ND	2.5	0.70	U
75-69-4	Trichlorofluoromethane	ND	2.5	0.70	U
107-06-2	1,2-Dichloroethane	ND	0.50	0.13	U
71-55-6	1,1,1-Trichloroethane	ND	2.5	0.70	U UJ
75-27-4	Bromodichloromethane	ND	0.50	0.19	U
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.16	U
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.14	U UJ
75-25-2	Bromoform	ND	2.0	0.65	U
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.17	U UJ
71-43-2	Benzene	ND	0.50	0.16	U
108-88-3	Toluene	ND	2.5	0.70	U
100-41-4	Ethylbenzene	ND	2.5	0.70	U
74-87-3	Chloromethane	ND	2.5	0.70	U
74-83-9	Bromomethane	ND	2.5	0.70	U UJ
75-01-4	Vinyl chloride	ND	1.0	0.07	U UJ
75-00-3	Chloroethane	ND	2.5	0.70	U
75-35-4	1,1-Dichloroethene	ND	0.50	0.17	U

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies	Lab Number : L2270805
Project Name : CONVENTUS/MROW	Project Number : 1186
Lab ID : L2270805-02	Date Collected : 12/14/22 11:00
Client ID : DUP	Date Received : 12/15/22
Sample Location : CONVENTUS 1 MAIN ST. BUFFALO NY	Date Analyzed : 12/21/22 14:34
Sample Matrix : WATER	Dilution Factor : 1
Analytical Method : 1,8260D	Analyst : MJV
Lab File ID : V22221221A19	Instrument ID : VOA122
Sample Amount : 10 ml	GC Column : RTX-502.2
Level : LOW	%Solids : N/A
Extract Volume (MeOH) : N/A	Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
156-60-5	trans-1,2-Dichloroethene	ND	2.5	0.70	U
79-01-6	Trichloroethene	ND	0.50	0.18	U
95-50-1	1,2-Dichlorobenzene	ND	2.5	0.70	U
541-73-1	1,3-Dichlorobenzene	ND	2.5	0.70	U
106-46-7	1,4-Dichlorobenzene	ND	2.5	0.70	U
1634-04-4	Methyl tert butyl ether	ND	2.5	0.70	U
179601-23-1	p/m-Xylene	ND	2.5	0.70	U UJ
95-47-6	o-Xylene	ND	2.5	0.70	U
156-59-2	cis-1,2-Dichloroethene	ND	2.5	0.70	U
100-42-5	Styrene	ND	2.5	0.70	U
75-71-8	Dichlorodifluoromethane	ND	5.0	1.0	U
67-64-1	Acetone	ND	5.0	1.5	U UJ
75-15-0	Carbon disulfide	ND	5.0	1.0	U
78-93-3	2-Butanone	ND	5.0	1.9	U UJ
108-10-1	4-Methyl-2-pentanone	ND	5.0	1.0	U
591-78-6	2-Hexanone	ND	5.0	1.0	U
106-93-4	1,2-Dibromoethane	ND	2.0	0.65	U
104-51-8	n-Butylbenzene	ND	2.5	0.70	U UJ
135-98-8	sec-Butylbenzene	ND	2.5	0.70	U
98-06-6	tert-Butylbenzene	ND	2.5	0.70	U
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.5	0.70	U
98-82-8	Isopropylbenzene	ND	2.5	0.70	U
99-87-6	p-Isopropyltoluene	ND	2.5	0.70	U
91-20-3	Naphthalene	ND	2.5	0.70	U
103-65-1	n-Propylbenzene	ND	2.5	0.70	U

Results Summary

Form 1

Volatile Organics by GC/MS

Client	: C&S Companies	Lab Number	: L2270805
Project Name	: CONVENTUS/MROW	Project Number	: 1186
Lab ID	: L2270805-02	Date Collected	: 12/14/22 11:00
Client ID	: DUP	Date Received	: 12/15/22
Sample Location	: CONVENTUS 1 MAIN ST. BUFFALO NY	Date Analyzed	: 12/21/22 14:34
Sample Matrix	: WATER	Dilution Factor	: 1
Analytical Method	: 1,8260D	Analyst	: MJV
Lab File ID	: V22221221A19	Instrument ID	: VOA122
Sample Amount	: 10 ml	GC Column	: RTX-502.2
Level	: LOW	%Solids	: N/A
Extract Volume (MeOH)	: N/A	Injection Volume	: N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
120-82-1	1,2,4-Trichlorobenzene	ND	2.5	0.70	U
108-67-8	1,3,5-Trimethylbenzene	ND	2.5	0.70	U
95-63-6	1,2,4-Trimethylbenzene	ND	2.5	0.70	U
79-20-9	Methyl Acetate	ND	2.0	0.23	U
110-82-7	Cyclohexane	ND	10	0.27	U
76-13-1	Freon-113	ND	2.5	0.70	U
108-87-2	Methyl cyclohexane	ND	10	0.40	U

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies	Lab Number : L2270805
Project Name : CONVENTUS/MROW	Project Number : 1186
Lab ID : L2270805-03	Date Collected : 12/14/22 11:40
Client ID : BCPMW-7	Date Received : 12/15/22
Sample Location : CONVENTUS 1 MAIN ST. BUFFALO NY	Date Analyzed : 12/22/22 22:51
Sample Matrix : WATER	Dilution Factor : 1
Analytical Method : 1,8260D	Analyst : MJV
Lab File ID : V30221222N20	Instrument ID : VOA130
Sample Amount : 10 ml	GC Column : RTX-502.2
Level : LOW	%Solids : N/A
Extract Volume (MeOH) : N/A	Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
75-09-2	Methylene chloride	ND	2.5	0.70	U
75-34-3	1,1-Dichloroethane	ND	2.5	0.70	U
67-66-3	Chloroform	ND	2.5	0.70	U
56-23-5	Carbon tetrachloride	ND	0.50	0.13	U
78-87-5	1,2-Dichloropropane	ND	1.0	0.14	U
124-48-1	Dibromochloromethane	ND	0.50	0.15	U
79-00-5	1,1,2-Trichloroethane	ND	1.5	0.50	U
127-18-4	Tetrachloroethene	ND	0.50	0.18	U
108-90-7	Chlorobenzene	ND	2.5	0.70	U
75-69-4	Trichlorofluoromethane	ND	2.5	0.70	U
107-06-2	1,2-Dichloroethane	ND	0.50	0.13	U
71-55-6	1,1,1-Trichloroethane	ND	2.5	0.70	U
75-27-4	Bromodichloromethane	ND	0.50	0.19	U
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.16	U
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.14	U
75-25-2	Bromoform	ND	2.0	0.65	U
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.17	U
71-43-2	Benzene	ND	0.50	0.16	U
108-88-3	Toluene	ND	2.5	0.70	U
100-41-4	Ethylbenzene	ND	2.5	0.70	U
74-87-3	Chloromethane	ND	2.5	0.70	U
74-83-9	Bromomethane	ND	2.5	0.70	U
75-01-4	Vinyl chloride	ND	1.0	0.07	U
75-00-3	Chloroethane	ND	2.5	0.70	U
75-35-4	1,1-Dichloroethene	ND	0.50	0.17	U



Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies	Lab Number : L2270805
Project Name : CONVENTUS/MROW	Project Number : 1186
Lab ID : L2270805-03	Date Collected : 12/14/22 11:40
Client ID : BCPMW-7	Date Received : 12/15/22
Sample Location : CONVENTUS 1 MAIN ST. BUFFALO NY	Date Analyzed : 12/22/22 22:51
Sample Matrix : WATER	Dilution Factor : 1
Analytical Method : 1,8260D	Analyst : MJV
Lab File ID : V30221222N20	Instrument ID : VOA130
Sample Amount : 10 ml	GC Column : RTX-502.2
Level : LOW	%Solids : N/A
Extract Volume (MeOH) : N/A	Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
156-60-5	trans-1,2-Dichloroethene	ND	2.5	0.70	U
79-01-6	Trichloroethene	ND	0.50	0.18	U
95-50-1	1,2-Dichlorobenzene	ND	2.5	0.70	U
541-73-1	1,3-Dichlorobenzene	ND	2.5	0.70	U
106-46-7	1,4-Dichlorobenzene	ND	2.5	0.70	U
1634-04-4	Methyl tert butyl ether	ND	2.5	0.70	U
179601-23-1	p/m-Xylene	ND	2.5	0.70	U
95-47-6	o-Xylene	ND	2.5	0.70	U
156-59-2	cis-1,2-Dichloroethene	ND	2.5	0.70	U
100-42-5	Styrene	ND	2.5	0.70	U
75-71-8	Dichlorodifluoromethane	ND	5.0	1.0	U
67-64-1	Acetone	ND	5.0	1.5	U
75-15-0	Carbon disulfide	ND	5.0	1.0	U
78-93-3	2-Butanone	ND	5.0	1.9	U UJ
108-10-1	4-Methyl-2-pentanone	ND	5.0	1.0	U
591-78-6	2-Hexanone	ND	5.0	1.0	U
106-93-4	1,2-Dibromoethane	ND	2.0	0.65	U
104-51-8	n-Butylbenzene	ND	2.5	0.70	U
135-98-8	sec-Butylbenzene	ND	2.5	0.70	U
98-06-6	tert-Butylbenzene	ND	2.5	0.70	U
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.5	0.70	U
98-82-8	Isopropylbenzene	ND	2.5	0.70	U
99-87-6	p-Isopropyltoluene	ND	2.5	0.70	U
91-20-3	Naphthalene	ND	2.5	0.70	U
103-65-1	n-Propylbenzene	ND	2.5	0.70	U

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies	Lab Number : L2270805
Project Name : CONVENTUS/MROW	Project Number : 1186
Lab ID : L2270805-03	Date Collected : 12/14/22 11:40
Client ID : BCPMW-7	Date Received : 12/15/22
Sample Location : CONVENTUS 1 MAIN ST. BUFFALO NY	Date Analyzed : 12/22/22 22:51
Sample Matrix : WATER	Dilution Factor : 1
Analytical Method : 1,8260D	Analyst : MJV
Lab File ID : V30221222N20	Instrument ID : VOA130
Sample Amount : 10 ml	GC Column : RTX-502.2
Level : LOW	%Solids : N/A
Extract Volume (MeOH) : N/A	Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
120-82-1	1,2,4-Trichlorobenzene	ND	2.5	0.70	U
108-67-8	1,3,5-Trimethylbenzene	ND	2.5	0.70	U
95-63-6	1,2,4-Trimethylbenzene	ND	2.5	0.70	U
79-20-9	Methyl Acetate	ND	2.0	0.23	U
110-82-7	Cyclohexane	ND	10	0.27	U
76-13-1	Freon-113	ND	2.5	0.70	U
108-87-2	Methyl cyclohexane	ND	10	0.40	U

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies
 Project Name : CONVENTUS/MROW
 Lab ID : L2270805-04D
 Client ID : BCPMW-4
 Sample Location : CONVENTUS 1 MAIN ST. BUFFALO NY
 Sample Matrix : WATER
 Analytical Method : 1,8260D
 Lab File ID : V22221221A21
 Sample Amount : 1 ml
 Level : LOW
 Extract Volume (MeOH) : N/A

Lab Number : L2270805
 Project Number : 1186
 Date Collected : 12/14/22 12:25
 Date Received : 12/15/22
 Date Analyzed : 12/21/22 15:23
 Dilution Factor : 10
 Analyst : MJV
 Instrument ID : VOA122
 GC Column : RTX-502.2
 %Solids : N/A
 Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
75-09-2	Methylene chloride	ND	25	7.0	U
75-34-3	1,1-Dichloroethane	ND	25	7.0	U
67-66-3	Chloroform	ND	25	7.0	U
56-23-5	Carbon tetrachloride	ND	5.0	1.3	U
78-87-5	1,2-Dichloropropane	ND	10	1.4	U
124-48-1	Dibromochloromethane	ND	5.0	1.5	U
79-00-5	1,1,2-Trichloroethane	ND	15	5.0	U UJ
127-18-4	Tetrachloroethene	ND	5.0	1.8	U UJ
108-90-7	Chlorobenzene	ND	25	7.0	U
75-69-4	Trichlorofluoromethane	ND	25	7.0	U
107-06-2	1,2-Dichloroethane	ND	5.0	1.3	U
71-55-6	1,1,1-Trichloroethane	ND	25	7.0	U UJ
75-27-4	Bromodichloromethane	ND	5.0	1.9	U
10061-02-6	trans-1,3-Dichloropropene	ND	5.0	1.6	U
10061-01-5	cis-1,3-Dichloropropene	ND	5.0	1.4	U UJ
75-25-2	Bromoform	ND	20	6.5	U
79-34-5	1,1,2,2-Tetrachloroethane	ND	5.0	1.7	U UJ
71-43-2	Benzene	3.5	5.0	1.6	J
108-88-3	Toluene	63	25	7.0	
100-41-4	Ethylbenzene	1100	25	7.0	
74-87-3	Chloromethane	ND	25	7.0	U
74-83-9	Bromomethane	ND	25	7.0	U UJ
75-01-4	Vinyl chloride	ND	10	0.71	U UJ
75-00-3	Chloroethane	ND	25	7.0	U
75-35-4	1,1-Dichloroethene	ND	5.0	1.7	U

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies
 Project Name : CONVENTUS/MROW
 Lab ID : L2270805-04D
 Client ID : BCPMW-4
 Sample Location : CONVENTUS 1 MAIN ST. BUFFALO NY
 Sample Matrix : WATER
 Analytical Method : 1,8260D
 Lab File ID : V22221221A21
 Sample Amount : 1 ml
 Level : LOW
 Extract Volume (MeOH) : N/A

Lab Number : L2270805
 Project Number : 1186
 Date Collected : 12/14/22 12:25
 Date Received : 12/15/22
 Date Analyzed : 12/21/22 15:23
 Dilution Factor : 10
 Analyst : MJV
 Instrument ID : VOA122
 GC Column : RTX-502.2
 %Solids : N/A
 Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
156-60-5	trans-1,2-Dichloroethene	ND	25	7.0	U
79-01-6	Trichloroethene	ND	5.0	1.8	U
95-50-1	1,2-Dichlorobenzene	ND	25	7.0	U
541-73-1	1,3-Dichlorobenzene	ND	25	7.0	U
106-46-7	1,4-Dichlorobenzene	ND	25	7.0	U
1634-04-4	Methyl tert butyl ether	ND	25	7.0	U
179601-23-1	p/m-Xylene	210	25	7.0	UJ
95-47-6	o-Xylene	16	25	7.0	J
156-59-2	cis-1,2-Dichloroethene	ND	25	7.0	U
100-42-5	Styrene	ND	25	7.0	U
75-71-8	Dichlorodifluoromethane	ND	50	10.	U
67-64-1	Acetone	ND	50	15.	UJ
75-15-0	Carbon disulfide	ND	50	10.	U
78-93-3	2-Butanone	ND	50	19.	UJ
108-10-1	4-Methyl-2-pentanone	ND	50	10.	U
591-78-6	2-Hexanone	ND	50	10.	U
106-93-4	1,2-Dibromoethane	ND	20	6.5	U
104-51-8	n-Butylbenzene	11	25	7.0	J UJ
135-98-8	sec-Butylbenzene	ND	25	7.0	U
98-06-6	tert-Butylbenzene	ND	25	7.0	U
96-12-8	1,2-Dibromo-3-chloropropane	ND	25	7.0	U
98-82-8	Isopropylbenzene	30	25	7.0	
99-87-6	p-Isopropyltoluene	ND	25	7.0	U
91-20-3	Naphthalene	290	25	7.0	
103-65-1	n-Propylbenzene	160	25	7.0	

Results Summary

Form 1

Volatile Organics by GC/MS

Client	: C&S Companies	Lab Number	: L2270805
Project Name	: CONVENTUS/MROW	Project Number	: 1186
Lab ID	: L2270805-04D	Date Collected	: 12/14/22 12:25
Client ID	: BCPMW-4	Date Received	: 12/15/22
Sample Location	: CONVENTUS 1 MAIN ST. BUFFALO NY	Date Analyzed	: 12/21/22 15:23
Sample Matrix	: WATER	Dilution Factor	: 10
Analytical Method	: 1,8260D	Analyst	: MJV
Lab File ID	: V22221221A21	Instrument ID	: VOA122
Sample Amount	: 1 ml	GC Column	: RTX-502.2
Level	: LOW	%Solids	: N/A
Extract Volume (MeOH)	: N/A	Injection Volume	: N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
120-82-1	1,2,4-Trichlorobenzene	ND	25	7.0	U
108-67-8	1,3,5-Trimethylbenzene	ND	25	7.0	U
95-63-6	1,2,4-Trimethylbenzene	1200	25	7.0	
79-20-9	Methyl Acetate	ND	20	2.3	U
110-82-7	Cyclohexane	140	100	2.7	
76-13-1	Freon-113	ND	25	7.0	U
108-87-2	Methyl cyclohexane	43	100	4.0	J

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies	Lab Number : L2270805
Project Name : CONVENTUS/MROW	Project Number : 1186
Lab ID : L2270805-05	Date Collected : 12/14/22 12:55
Client ID : BCPMW-3	Date Received : 12/15/22
Sample Location : CONVENTUS 1 MAIN ST. BUFFALO NY	Date Analyzed : 12/21/22 15:48
Sample Matrix : WATER	Dilution Factor : 1
Analytical Method : 1,8260D	Analyst : MJV
Lab File ID : V22221221A22	Instrument ID : VOA122
Sample Amount : 10 ml	GC Column : RTX-502.2
Level : LOW	%Solids : N/A
Extract Volume (MeOH) : N/A	Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
75-09-2	Methylene chloride	ND	2.5	0.70	U
75-34-3	1,1-Dichloroethane	ND	2.5	0.70	U
67-66-3	Chloroform	ND	2.5	0.70	U
56-23-5	Carbon tetrachloride	ND	0.50	0.13	U
78-87-5	1,2-Dichloropropane	ND	1.0	0.14	U
124-48-1	Dibromochloromethane	ND	0.50	0.15	U
79-00-5	1,1,2-Trichloroethane	ND	1.5	0.50	U UJ
127-18-4	Tetrachloroethene	ND	0.50	0.18	U UJ
108-90-7	Chlorobenzene	ND	2.5	0.70	U
75-69-4	Trichlorofluoromethane	ND	2.5	0.70	U
107-06-2	1,2-Dichloroethane	ND	0.50	0.13	U
71-55-6	1,1,1-Trichloroethane	ND	2.5	0.70	U UJ
75-27-4	Bromodichloromethane	ND	0.50	0.19	U
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.16	U
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.14	U UJ
75-25-2	Bromoform	ND	2.0	0.65	U
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.17	U UJ
71-43-2	Benzene	ND	0.50	0.16	U
108-88-3	Toluene	ND	2.5	0.70	U
100-41-4	Ethylbenzene	11	2.5	0.70	
74-87-3	Chloromethane	ND	2.5	0.70	U
74-83-9	Bromomethane	ND	2.5	0.70	U UJ
75-01-4	Vinyl chloride	ND	1.0	0.07	U UJ
75-00-3	Chloroethane	ND	2.5	0.70	U
75-35-4	1,1-Dichloroethene	ND	0.50	0.17	U

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies	Lab Number : L2270805
Project Name : CONVENTUS/MROW	Project Number : 1186
Lab ID : L2270805-05	Date Collected : 12/14/22 12:55
Client ID : BCPMW-3	Date Received : 12/15/22
Sample Location : CONVENTUS 1 MAIN ST. BUFFALO NY	Date Analyzed : 12/21/22 15:48
Sample Matrix : WATER	Dilution Factor : 1
Analytical Method : 1,8260D	Analyst : MJV
Lab File ID : V22221221A22	Instrument ID : VOA122
Sample Amount : 10 ml	GC Column : RTX-502.2
Level : LOW	%Solids : N/A
Extract Volume (MeOH) : N/A	Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
156-60-5	trans-1,2-Dichloroethene	ND	2.5	0.70	U
79-01-6	Trichloroethene	ND	0.50	0.18	U
95-50-1	1,2-Dichlorobenzene	ND	2.5	0.70	U
541-73-1	1,3-Dichlorobenzene	ND	2.5	0.70	U
106-46-7	1,4-Dichlorobenzene	ND	2.5	0.70	U
1634-04-4	Methyl tert butyl ether	ND	2.5	0.70	U
179601-23-1	p/m-Xylene	12	2.5	0.70	UJ
95-47-6	o-Xylene	ND	2.5	0.70	U
156-59-2	cis-1,2-Dichloroethene	ND	2.5	0.70	U
100-42-5	Styrene	ND	2.5	0.70	U
75-71-8	Dichlorodifluoromethane	ND	5.0	1.0	U
67-64-1	Acetone	ND	5.0	1.5	UJ
75-15-0	Carbon disulfide	ND	5.0	1.0	U
78-93-3	2-Butanone	ND	5.0	1.9	UJ
108-10-1	4-Methyl-2-pentanone	ND	5.0	1.0	U
591-78-6	2-Hexanone	ND	5.0	1.0	U
106-93-4	1,2-Dibromoethane	ND	2.0	0.65	U
104-51-8	n-Butylbenzene	ND	2.5	0.70	UJ
135-98-8	sec-Butylbenzene	ND	2.5	0.70	U
98-06-6	tert-Butylbenzene	ND	2.5	0.70	U
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.5	0.70	U
98-82-8	Isopropylbenzene	ND	2.5	0.70	U
99-87-6	p-Isopropyltoluene	ND	2.5	0.70	U
91-20-3	Naphthalene	4.2	2.5	0.70	
103-65-1	n-Propylbenzene	ND	2.5	0.70	U

Results Summary

Form 1

Volatile Organics by GC/MS

Client	: C&S Companies	Lab Number	: L2270805
Project Name	: CONVENTUS/MROW	Project Number	: 1186
Lab ID	: L2270805-05	Date Collected	: 12/14/22 12:55
Client ID	: BCPMW-3	Date Received	: 12/15/22
Sample Location	: CONVENTUS 1 MAIN ST. BUFFALO NY	Date Analyzed	: 12/21/22 15:48
Sample Matrix	: WATER	Dilution Factor	: 1
Analytical Method	: 1,8260D	Analyst	: MJV
Lab File ID	: V22221221A22	Instrument ID	: VOA122
Sample Amount	: 10 ml	GC Column	: RTX-502.2
Level	: LOW	%Solids	: N/A
Extract Volume (MeOH)	: N/A	Injection Volume	: N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
120-82-1	1,2,4-Trichlorobenzene	ND	2.5	0.70	U
108-67-8	1,3,5-Trimethylbenzene	ND	2.5	0.70	U
95-63-6	1,2,4-Trimethylbenzene	13	2.5	0.70	
79-20-9	Methyl Acetate	ND	2.0	0.23	U
110-82-7	Cyclohexane	1.5	10	0.27	J
76-13-1	Freon-113	ND	2.5	0.70	U
108-87-2	Methyl cyclohexane	0.65	10	0.40	J

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies
 Project Name : CONVENTUS/MROW
 Lab ID : L2270805-06
 Client ID : BCPMW-6
 Sample Location : CONVENTUS 1 MAIN ST. BUFFALO NY
 Sample Matrix : WATER
 Analytical Method : 1,8260D
 Lab File ID : V22221221A23
 Sample Amount : 10 ml
 Level : LOW
 Extract Volume (MeOH) : N/A

Lab Number : L2270805
 Project Number : 1186
 Date Collected : 12/14/22 13:30
 Date Received : 12/15/22
 Date Analyzed : 12/21/22 16:12
 Dilution Factor : 1
 Analyst : MJV
 Instrument ID : VOA122
 GC Column : RTX-502.2
 %Solids : N/A
 Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
75-09-2	Methylene chloride	ND	2.5	0.70	U
75-34-3	1,1-Dichloroethane	ND	2.5	0.70	U
67-66-3	Chloroform	ND	2.5	0.70	U
56-23-5	Carbon tetrachloride	ND	0.50	0.13	U
78-87-5	1,2-Dichloropropane	ND	1.0	0.14	U
124-48-1	Dibromochloromethane	ND	0.50	0.15	U
79-00-5	1,1,2-Trichloroethane	ND	1.5	0.50	U UJ
127-18-4	Tetrachloroethene	ND	0.50	0.18	U UJ
108-90-7	Chlorobenzene	ND	2.5	0.70	U
75-69-4	Trichlorofluoromethane	ND	2.5	0.70	U
107-06-2	1,2-Dichloroethane	ND	0.50	0.13	U
71-55-6	1,1,1-Trichloroethane	ND	2.5	0.70	U UJ
75-27-4	Bromodichloromethane	ND	0.50	0.19	U
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.16	U
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.14	U UJ
75-25-2	Bromoform	ND	2.0	0.65	U
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.17	U UJ
71-43-2	Benzene	ND	0.50	0.16	U
108-88-3	Toluene	ND	2.5	0.70	U
100-41-4	Ethylbenzene	ND	2.5	0.70	U
74-87-3	Chloromethane	ND	2.5	0.70	U
74-83-9	Bromomethane	ND	2.5	0.70	U UJ
75-01-4	Vinyl chloride	ND	1.0	0.07	U UJ
75-00-3	Chloroethane	ND	2.5	0.70	U
75-35-4	1,1-Dichloroethene	ND	0.50	0.17	U

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies	Lab Number : L2270805
Project Name : CONVENTUS/MROW	Project Number : 1186
Lab ID : L2270805-06	Date Collected : 12/14/22 13:30
Client ID : BCPMW-6	Date Received : 12/15/22
Sample Location : CONVENTUS 1 MAIN ST. BUFFALO NY	Date Analyzed : 12/21/22 16:12
Sample Matrix : WATER	Dilution Factor : 1
Analytical Method : 1,8260D	Analyst : MJV
Lab File ID : V22221221A23	Instrument ID : VOA122
Sample Amount : 10 ml	GC Column : RTX-502.2
Level : LOW	%Solids : N/A
Extract Volume (MeOH) : N/A	Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
156-60-5	trans-1,2-Dichloroethene	ND	2.5	0.70	U
79-01-6	Trichloroethene	ND	0.50	0.18	U
95-50-1	1,2-Dichlorobenzene	ND	2.5	0.70	U
541-73-1	1,3-Dichlorobenzene	ND	2.5	0.70	U
106-46-7	1,4-Dichlorobenzene	ND	2.5	0.70	U
1634-04-4	Methyl tert butyl ether	ND	2.5	0.70	U
179601-23-1	p/m-Xylene	ND	2.5	0.70	U UJ
95-47-6	o-Xylene	ND	2.5	0.70	U
156-59-2	cis-1,2-Dichloroethene	ND	2.5	0.70	U
100-42-5	Styrene	ND	2.5	0.70	U
75-71-8	Dichlorodifluoromethane	ND	5.0	1.0	U
67-64-1	Acetone	ND	5.0	1.5	U UJ
75-15-0	Carbon disulfide	ND	5.0	1.0	U
78-93-3	2-Butanone	ND	5.0	1.9	U UJ
108-10-1	4-Methyl-2-pentanone	ND	5.0	1.0	U
591-78-6	2-Hexanone	ND	5.0	1.0	U
106-93-4	1,2-Dibromoethane	ND	2.0	0.65	U
104-51-8	n-Butylbenzene	ND	2.5	0.70	U UJ
135-98-8	sec-Butylbenzene	ND	2.5	0.70	U
98-06-6	tert-Butylbenzene	ND	2.5	0.70	U
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.5	0.70	U
98-82-8	Isopropylbenzene	ND	2.5	0.70	U
99-87-6	p-Isopropyltoluene	ND	2.5	0.70	U
91-20-3	Naphthalene	ND	2.5	0.70	U
103-65-1	n-Propylbenzene	ND	2.5	0.70	U

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Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies	Lab Number : L2270805
Project Name : CONVENTUS/MROW	Project Number : 1186
Lab ID : L2270805-06	Date Collected : 12/14/22 13:30
Client ID : BCPMW-6	Date Received : 12/15/22
Sample Location : CONVENTUS 1 MAIN ST. BUFFALO NY	Date Analyzed : 12/21/22 16:12
Sample Matrix : WATER	Dilution Factor : 1
Analytical Method : 1,8260D	Analyst : MJV
Lab File ID : V22221221A23	Instrument ID : VOA122
Sample Amount : 10 ml	GC Column : RTX-502.2
Level : LOW	%Solids : N/A
Extract Volume (MeOH) : N/A	Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
120-82-1	1,2,4-Trichlorobenzene	ND	2.5	0.70	U
108-67-8	1,3,5-Trimethylbenzene	ND	2.5	0.70	U
95-63-6	1,2,4-Trimethylbenzene	ND	2.5	0.70	U
79-20-9	Methyl Acetate	ND	2.0	0.23	U
110-82-7	Cyclohexane	ND	10	0.27	U
76-13-1	Freon-113	ND	2.5	0.70	U
108-87-2	Methyl cyclohexane	ND	10	0.40	U

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies
 Project Name : CONVENTUS/MROW
 Lab ID : L2270805-07D
 Client ID : BCPMW-5
 Sample Location : CONVENTUS 1 MAIN ST. BUFFALO NY
 Sample Matrix : WATER
 Analytical Method : 1,8260D
 Lab File ID : V22221221A24
 Sample Amount : 1 ml
 Level : LOW
 Extract Volume (MeOH) : N/A

Lab Number : L2270805
 Project Number : 1186
 Date Collected : 12/14/22 14:10
 Date Received : 12/15/22
 Date Analyzed : 12/21/22 16:37
 Dilution Factor : 10
 Analyst : MJV
 Instrument ID : VOA122
 GC Column : RTX-502.2
 %Solids : N/A
 Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
75-09-2	Methylene chloride	ND	25	7.0	U
75-34-3	1,1-Dichloroethane	ND	25	7.0	U
67-66-3	Chloroform	ND	25	7.0	U
56-23-5	Carbon tetrachloride	ND	5.0	1.3	U
78-87-5	1,2-Dichloropropane	ND	10	1.4	U
124-48-1	Dibromochloromethane	ND	5.0	1.5	U
79-00-5	1,1,2-Trichloroethane	ND	15	5.0	U UJ
127-18-4	Tetrachloroethene	ND	5.0	1.8	U UJ
108-90-7	Chlorobenzene	ND	25	7.0	U
75-69-4	Trichlorofluoromethane	ND	25	7.0	U
107-06-2	1,2-Dichloroethane	ND	5.0	1.3	U
71-55-6	1,1,1-Trichloroethane	ND	25	7.0	U UJ
75-27-4	Bromodichloromethane	ND	5.0	1.9	U
10061-02-6	trans-1,3-Dichloropropene	ND	5.0	1.6	U
10061-01-5	cis-1,3-Dichloropropene	ND	5.0	1.4	U UJ
75-25-2	Bromoform	ND	20	6.5	U
79-34-5	1,1,2,2-Tetrachloroethane	ND	5.0	1.7	U UJ
71-43-2	Benzene	ND	5.0	1.6	U
108-88-3	Toluene	9.6	25	7.0	J
100-41-4	Ethylbenzene	820	25	7.0	
74-87-3	Chloromethane	ND	25	7.0	U
74-83-9	Bromomethane	ND	25	7.0	U UJ
75-01-4	Vinyl chloride	ND	10	0.71	U UJ
75-00-3	Chloroethane	ND	25	7.0	U
75-35-4	1,1-Dichloroethene	ND	5.0	1.7	U

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies	Lab Number : L2270805
Project Name : CONVENTUS/MROW	Project Number : 1186
Lab ID : L2270805-07D	Date Collected : 12/14/22 14:10
Client ID : BCPMW-5	Date Received : 12/15/22
Sample Location : CONVENTUS 1 MAIN ST. BUFFALO NY	Date Analyzed : 12/21/22 16:37
Sample Matrix : WATER	Dilution Factor : 10
Analytical Method : 1,8260D	Analyst : MJV
Lab File ID : V22221221A24	Instrument ID : VOA122
Sample Amount : 1 ml	GC Column : RTX-502.2
Level : LOW	%Solids : N/A
Extract Volume (MeOH) : N/A	Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
156-60-5	trans-1,2-Dichloroethene	ND	25	7.0	U
79-01-6	Trichloroethene	ND	5.0	1.8	U
95-50-1	1,2-Dichlorobenzene	ND	25	7.0	U
541-73-1	1,3-Dichlorobenzene	ND	25	7.0	U
106-46-7	1,4-Dichlorobenzene	ND	25	7.0	U
1634-04-4	Methyl tert butyl ether	ND	25	7.0	U
179601-23-1	p/m-Xylene	2400	25	7.0	UJ
95-47-6	o-Xylene	62	25	7.0	
156-59-2	cis-1,2-Dichloroethene	ND	25	7.0	U
100-42-5	Styrene	ND	25	7.0	U
75-71-8	Dichlorodifluoromethane	ND	50	10.	U
67-64-1	Acetone	ND	50	15.	UJ
75-15-0	Carbon disulfide	ND	50	10.	U
78-93-3	2-Butanone	ND	50	19.	UJ
108-10-1	4-Methyl-2-pentanone	ND	50	10.	U
591-78-6	2-Hexanone	ND	50	10.	U
106-93-4	1,2-Dibromoethane	ND	20	6.5	U
104-51-8	n-Butylbenzene	ND	25	7.0	UJ
135-98-8	sec-Butylbenzene	ND	25	7.0	U
98-06-6	tert-Butylbenzene	ND	25	7.0	U
96-12-8	1,2-Dibromo-3-chloropropane	ND	25	7.0	U
98-82-8	Isopropylbenzene	16	25	7.0	J
99-87-6	p-Isopropyltoluene	ND	25	7.0	U
91-20-3	Naphthalene	520	25	7.0	
103-65-1	n-Propylbenzene	71	25	7.0	

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies	Lab Number : L2270805
Project Name : CONVENTUS/MROW	Project Number : 1186
Lab ID : L2270805-07D	Date Collected : 12/14/22 14:10
Client ID : BCPMW-5	Date Received : 12/15/22
Sample Location : CONVENTUS 1 MAIN ST. BUFFALO NY	Date Analyzed : 12/21/22 16:37
Sample Matrix : WATER	Dilution Factor : 10
Analytical Method : 1,8260D	Analyst : MJV
Lab File ID : V22221221A24	Instrument ID : VOA122
Sample Amount : 1 ml	GC Column : RTX-502.2
Level : LOW	%Solids : N/A
Extract Volume (MeOH) : N/A	Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
120-82-1	1,2,4-Trichlorobenzene	ND	25	7.0	U
108-67-8	1,3,5-Trimethylbenzene	350	25	7.0	
95-63-6	1,2,4-Trimethylbenzene	1400	25	7.0	
79-20-9	Methyl Acetate	ND	20	2.3	U
110-82-7	Cyclohexane	120	100	2.7	
76-13-1	Freon-113	ND	25	7.0	U
108-87-2	Methyl cyclohexane	54	100	4.0	J

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies	Lab Number : L2270805
Project Name : CONVENTUS/MROW	Project Number : 1186
Lab ID : L2270805-08	Date Collected : 12/15/22 00:00
Client ID : TRIP BLANK	Date Received : 12/15/22
Sample Location : CONVENTUS 1 MAIN ST. BUFFALO NY	Date Analyzed : 12/23/22 14:38
Sample Matrix : WATER	Dilution Factor : 1
Analytical Method : 1,8260D	Analyst : MJV
Lab File ID : V01221223A18	Instrument ID : VOA101
Sample Amount : 10 ml	GC Column : RTX-502.2
Level : LOW	%Solids : N/A
Extract Volume (MeOH) : N/A	Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
75-09-2	Methylene chloride	ND	2.5	0.70	U
75-34-3	1,1-Dichloroethane	ND	2.5	0.70	U
67-66-3	Chloroform	ND	2.5	0.70	U
56-23-5	Carbon tetrachloride	ND	0.50	0.13	U
78-87-5	1,2-Dichloropropane	ND	1.0	0.14	U
124-48-1	Dibromochloromethane	ND	0.50	0.15	U
79-00-5	1,1,2-Trichloroethane	ND	1.5	0.50	U
127-18-4	Tetrachloroethene	ND	0.50	0.18	U
108-90-7	Chlorobenzene	ND	2.5	0.70	U
75-69-4	Trichlorofluoromethane	ND	2.5	0.70	U
107-06-2	1,2-Dichloroethane	ND	0.50	0.13	U
71-55-6	1,1,1-Trichloroethane	ND	2.5	0.70	U
75-27-4	Bromodichloromethane	ND	0.50	0.19	U
10061-02-6	trans-1,3-Dichloropropene	ND	0.50	0.16	U
10061-01-5	cis-1,3-Dichloropropene	ND	0.50	0.14	U
75-25-2	Bromoform	ND	2.0	0.65	U
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50	0.17	U
71-43-2	Benzene	ND	0.50	0.16	U
108-88-3	Toluene	ND	2.5	0.70	U
100-41-4	Ethylbenzene	ND	2.5	0.70	U
74-87-3	Chloromethane	ND	2.5	0.70	U
74-83-9	Bromomethane	ND	2.5	0.70	U
75-01-4	Vinyl chloride	ND	1.0	0.07	U UJ
75-00-3	Chloroethane	ND	2.5	0.70	U
75-35-4	1,1-Dichloroethene	ND	0.50	0.17	U

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies	Lab Number : L2270805
Project Name : CONVENTUS/MROW	Project Number : 1186
Lab ID : L2270805-08	Date Collected : 12/15/22 00:00
Client ID : TRIP BLANK	Date Received : 12/15/22
Sample Location : CONVENTUS 1 MAIN ST. BUFFALO NY	Date Analyzed : 12/23/22 14:38
Sample Matrix : WATER	Dilution Factor : 1
Analytical Method : 1,8260D	Analyst : MJV
Lab File ID : V01221223A18	Instrument ID : VOA101
Sample Amount : 10 ml	GC Column : RTX-502.2
Level : LOW	%Solids : N/A
Extract Volume (MeOH) : N/A	Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
156-60-5	trans-1,2-Dichloroethene	ND	2.5	0.70	U
79-01-6	Trichloroethene	ND	0.50	0.18	U
95-50-1	1,2-Dichlorobenzene	ND	2.5	0.70	U
541-73-1	1,3-Dichlorobenzene	ND	2.5	0.70	U
106-46-7	1,4-Dichlorobenzene	ND	2.5	0.70	U
1634-04-4	Methyl tert butyl ether	ND	2.5	0.70	U
179601-23-1	p/m-Xylene	ND	2.5	0.70	U
95-47-6	o-Xylene	ND	2.5	0.70	U
156-59-2	cis-1,2-Dichloroethene	ND	2.5	0.70	U
100-42-5	Styrene	ND	2.5	0.70	U
75-71-8	Dichlorodifluoromethane	ND	5.0	1.0	U
67-64-1	Acetone	ND	5.0	1.5	U
75-15-0	Carbon disulfide	ND	5.0	1.0	U
78-93-3	2-Butanone	ND	5.0	1.9	U UJ
108-10-1	4-Methyl-2-pentanone	ND	5.0	1.0	U
591-78-6	2-Hexanone	ND	5.0	1.0	U
106-93-4	1,2-Dibromoethane	ND	2.0	0.65	U
104-51-8	n-Butylbenzene	ND	2.5	0.70	U
135-98-8	sec-Butylbenzene	ND	2.5	0.70	U
98-06-6	tert-Butylbenzene	ND	2.5	0.70	U
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.5	0.70	U
98-82-8	Isopropylbenzene	ND	2.5	0.70	U
99-87-6	p-Isopropyltoluene	ND	2.5	0.70	U
91-20-3	Naphthalene	ND	2.5	0.70	U
103-65-1	n-Propylbenzene	ND	2.5	0.70	U

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies	Lab Number : L2270805
Project Name : CONVENTUS/MROW	Project Number : 1186
Lab ID : L2270805-08	Date Collected : 12/15/22 00:00
Client ID : TRIP BLANK	Date Received : 12/15/22
Sample Location : CONVENTUS 1 MAIN ST. BUFFALO NY	Date Analyzed : 12/23/22 14:38
Sample Matrix : WATER	Dilution Factor : 1
Analytical Method : 1,8260D	Analyst : MJV
Lab File ID : V01221223A18	Instrument ID : VOA101
Sample Amount : 10 ml	GC Column : RTX-502.2
Level : LOW	%Solids : N/A
Extract Volume (MeOH) : N/A	Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
120-82-1	1,2,4-Trichlorobenzene	ND	2.5	0.70	U
108-67-8	1,3,5-Trimethylbenzene	ND	2.5	0.70	U
95-63-6	1,2,4-Trimethylbenzene	ND	2.5	0.70	U
79-20-9	Methyl Acetate	ND	2.0	0.23	U
110-82-7	Cyclohexane	ND	10	0.27	U
76-13-1	Freon-113	ND	2.5	0.70	U
108-87-2	Methyl cyclohexane	ND	10	0.40	U

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies	Lab Number : L2270805
Project Name : CONVENTUS/MROW	Project Number : 1186
Lab ID : L2270805-09D	Date Collected : 12/15/22 10:45
Client ID : MSMW-2	Date Received : 12/15/22
Sample Location : CONVENTUS 1 MAIN ST. BUFFALO NY	Date Analyzed : 12/21/22 17:01
Sample Matrix : WATER	Dilution Factor : 10
Analytical Method : 1,8260D	Analyst : MJV
Lab File ID : V22221221A25	Instrument ID : VOA122
Sample Amount : 1 ml	GC Column : RTX-502.2
Level : LOW	%Solids : N/A
Extract Volume (MeOH) : N/A	Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
75-09-2	Methylene chloride	ND	25	7.0	U
75-34-3	1,1-Dichloroethane	ND	25	7.0	U
67-66-3	Chloroform	ND	25	7.0	U
56-23-5	Carbon tetrachloride	ND	5.0	1.3	U
78-87-5	1,2-Dichloropropane	ND	10	1.4	U
124-48-1	Dibromochloromethane	ND	5.0	1.5	U
79-00-5	1,1,2-Trichloroethane	ND	15	5.0	U UJ
127-18-4	Tetrachloroethene	ND	5.0	1.8	U UJ
108-90-7	Chlorobenzene	ND	25	7.0	U
75-69-4	Trichlorofluoromethane	ND	25	7.0	U
107-06-2	1,2-Dichloroethane	ND	5.0	1.3	U
71-55-6	1,1,1-Trichloroethane	ND	25	7.0	U UJ
75-27-4	Bromodichloromethane	ND	5.0	1.9	U
10061-02-6	trans-1,3-Dichloropropene	ND	5.0	1.6	U
10061-01-5	cis-1,3-Dichloropropene	ND	5.0	1.4	U UJ
75-25-2	Bromoform	ND	20	6.5	U
79-34-5	1,1,2,2-Tetrachloroethane	ND	5.0	1.7	U UJ
71-43-2	Benzene	280	5.0	1.6	
108-88-3	Toluene	210	25	7.0	
100-41-4	Ethylbenzene	820	25	7.0	
74-87-3	Chloromethane	ND	25	7.0	U
74-83-9	Bromomethane	ND	25	7.0	U UJ
75-01-4	Vinyl chloride	ND	10	0.71	U UJ
75-00-3	Chloroethane	ND	25	7.0	U
75-35-4	1,1-Dichloroethene	ND	5.0	1.7	U

Results Summary

Form 1

Volatile Organics by GC/MS

Client : C&S Companies	Lab Number : L2270805
Project Name : CONVENTUS/MROW	Project Number : 1186
Lab ID : L2270805-09D	Date Collected : 12/15/22 10:45
Client ID : MSMW-2	Date Received : 12/15/22
Sample Location : CONVENTUS 1 MAIN ST. BUFFALO NY	Date Analyzed : 12/21/22 17:01
Sample Matrix : WATER	Dilution Factor : 10
Analytical Method : 1,8260D	Analyst : MJV
Lab File ID : V22221221A25	Instrument ID : VOA122
Sample Amount : 1 ml	GC Column : RTX-502.2
Level : LOW	%Solids : N/A
Extract Volume (MeOH) : N/A	Injection Volume : N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
156-60-5	trans-1,2-Dichloroethene	ND	25	7.0	U
79-01-6	Trichloroethene	ND	5.0	1.8	U
95-50-1	1,2-Dichlorobenzene	ND	25	7.0	U
541-73-1	1,3-Dichlorobenzene	ND	25	7.0	U
106-46-7	1,4-Dichlorobenzene	ND	25	7.0	U
1634-04-4	Methyl tert butyl ether	ND	25	7.0	U
179601-23-1	p/m-Xylene	1900	25	7.0	UJ
95-47-6	o-Xylene	490	25	7.0	
156-59-2	cis-1,2-Dichloroethene	ND	25	7.0	U
100-42-5	Styrene	ND	25	7.0	U
75-71-8	Dichlorodifluoromethane	ND	50	10.	U
67-64-1	Acetone	ND	50	15.	UJ
75-15-0	Carbon disulfide	ND	50	10.	U
78-93-3	2-Butanone	ND	50	19.	UJ
108-10-1	4-Methyl-2-pentanone	ND	50	10.	U
591-78-6	2-Hexanone	ND	50	10.	U
106-93-4	1,2-Dibromoethane	ND	20	6.5	U
104-51-8	n-Butylbenzene	12	25	7.0	J UJ
135-98-8	sec-Butylbenzene	ND	25	7.0	U
98-06-6	tert-Butylbenzene	ND	25	7.0	U
96-12-8	1,2-Dibromo-3-chloropropane	ND	25	7.0	U
98-82-8	Isopropylbenzene	25	25	7.0	
99-87-6	p-Isopropyltoluene	ND	25	7.0	U
91-20-3	Naphthalene	340	25	7.0	
103-65-1	n-Propylbenzene	94	25	7.0	

Results Summary

Form 1

Volatile Organics by GC/MS

Client	: C&S Companies	Lab Number	: L2270805
Project Name	: CONVENTUS/MROW	Project Number	: 1186
Lab ID	: L2270805-09D	Date Collected	: 12/15/22 10:45
Client ID	: MSMW-2	Date Received	: 12/15/22
Sample Location	: CONVENTUS 1 MAIN ST. BUFFALO NY	Date Analyzed	: 12/21/22 17:01
Sample Matrix	: WATER	Dilution Factor	: 10
Analytical Method	: 1,8260D	Analyst	: MJV
Lab File ID	: V22221221A25	Instrument ID	: VOA122
Sample Amount	: 1 ml	GC Column	: RTX-502.2
Level	: LOW	%Solids	: N/A
Extract Volume (MeOH)	: N/A	Injection Volume	: N/A

CAS NO.	Parameter	ug/L			Qualifier
		Results	RL	MDL	
120-82-1	1,2,4-Trichlorobenzene	ND	25	7.0	U
108-67-8	1,3,5-Trimethylbenzene	490	25	7.0	
95-63-6	1,2,4-Trimethylbenzene	1200	25	7.0	
79-20-9	Methyl Acetate	ND	20	2.3	U
110-82-7	Cyclohexane	160	100	2.7	
76-13-1	Freon-113	ND	25	7.0	U
108-87-2	Methyl cyclohexane	110	100	4.0	

Appendix B

Laboratory QC Documentation

Evaluate Continuing Calibration Report

Data Path : I:\VOLATILES\VOA122\2022\221209AICAL\
 Data File : V22221209A18.D
 Acq On : 09 Dec 2022 09:33 pm
 Operator : VOA122:PID
 Sample : C8260STD10PPB
 Misc : WG1722111,ICAL (Sig #1); WG,ICAL (Sig #2)
 ALS Vial : 18 Sample Multiplier: 1

Quant Time: Dec 12 08:25:57 2022
 Quant Method : I:\VOLATILES\VOA122\2022\221209AICAL\V122_221209A_8260.m
 Quant Title : VOLATILES BY GC/MS
 QLast Update : Mon Dec 12 08:24:31 2022
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 20% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Fluorobenzene	1.000	1.000	0.0	99	0.00
2 TP	Dichlorodifluoromethane	0.137	0.197	-43.8#	124	0.00
3 TP	Chloromethane	0.223	0.300	-34.5#	119	0.00
4 TC	Vinyl chloride	0.162	0.197	-21.6#	107	0.00
5 TP	Bromomethane	* 10.000	12.528	-25.3#	111	0.00
6 TP	Chloroethane	0.100	0.121	-21.0#	104	0.00
7 TP	Trichlorofluoromethane	0.238	0.287	-20.6#	105	0.00
8 TP	Ethyl ether	0.048	0.063	-31.3#	121	0.00
10 TC	1,1-Dichloroethene	0.138	0.144	-4.3	94	0.00
11 TP	Carbon disulfide	0.264	0.404	-53.0#	137	0.00
12 TP	Freon-113	0.133	0.162	-21.8#	101	0.00
13 TP	Iodomethane	0.191	0.192	-0.5	88	0.00
14 TP	Acrolein	0.017	0.017	0.0	90	0.00
15 TP	Methylene chloride	0.150	0.169	-12.7	103	0.00
17 TP	Acetone	* 10.000	11.573	-15.7	103	0.00
18 TP	trans-1,2-Dichloroethene	0.151	0.160	-6.0	95	0.00
19 TP	Methyl acetate	* 10.000	11.882	-18.8	99	0.00
21 TP	Methyl tert-butyl ether	0.230	0.245	-6.5	103	0.00
22 TP	tert-Butyl alcohol	0.00810	0.00814#	-0.5	94	0.00
24 TP	Diisopropyl ether	0.420	0.403	4.0	95	0.00
25 TP	1,1-Dichloroethane	0.317	0.356	-12.3	99	0.00
26 TP	Halothane	0.101	0.115	-13.9	101	0.00
27 TP	Acrylonitrile	0.038	0.042	-10.5	100	0.00
28 TP	Ethyl tert-butyl ether	0.361	0.336	6.9	96	0.00
29 TP	Vinyl acetate	0.214	0.126	41.1#	64	0.00
30 TP	cis-1,2-Dichloroethene	0.122	0.124#	-1.6	94	0.00
31 TP	2,2-Dichloropropane	0.245	0.244	0.4	88	0.00
32 TP	Bromochloromethane	0.070	0.075#	-7.1	97	0.00
33 TP	Cyclohexane	0.327	0.352	-7.6	96	0.00
34 TC	Chloroform	0.299	0.324	-8.4	105	0.00
35 TP	Ethyl acetate	0.076	0.073	3.9	103	0.00
36 TP	Carbon tetrachloride	0.229	0.240	-4.8	92	0.00
37 TP	Tetrahydrofuran	* 10.000	10.118	-1.2	96	0.00
38 S	Dibromofluoromethane	0.313	0.342	-9.3	105	0.00
39 TP	1,1,1-Trichloroethane	0.257	0.276	-7.4	97	0.00
41 TP	2-Butanone	0.035	0.030	14.3	90	0.00
42 TP	1,1-Dichloropropene	0.187	0.191	-2.1	95	0.00
44 TP	Benzene	0.519	0.511	1.5	92	0.00
45 TP	tert-Amyl methyl ether	0.260	0.239	8.1	92	0.00

Evaluate Continuing Calibration Report

Data Path : I:\VOLATILES\VOA101\2022\221216A\
 Data File : V01221216A17.D
 Acq On : 16 Dec 2022 7:06 pm
 Operator : VOA101:PID
 Sample : C8260STD10PPB
 Misc : WG1724555,ICAL
 ALS Vial : 17 Sample Multiplier: 1

Quant Time: Dec 17 10:07:50 2022
 Quant Method : I:\VOLATILES\VOA101\2022\221216A\V101_221216A_8260.m
 Quant Title : VOLATILES BY GC/MS
 QLast Update : Sat Dec 17 10:04:09 2022
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 20% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Fluorobenzene	1.000	1.000	0.0	103	0.00
2 TP	Dichlorodifluoromethane	0.227	0.225	0.9	102	0.00
3 TP	Chloromethane	0.264	0.282	-6.8	111	0.00
4 TC	Vinyl chloride	0.369	0.378	-2.4	102	0.00
5 TP	Bromomethane	0.252	0.260	-3.2	125	0.00
6 TP	Chloroethane	0.301	0.321	-6.6	109	0.00
7 TP	Trichlorofluoromethane	0.640	0.690	-7.8	113	0.00
8 TP	Ethyl ether	0.152	0.201	-32.2#	146	0.00
10 TC	1,1-Dichloroethene	0.359	0.363	-1.1	107	0.00
11 TP	Carbon disulfide	0.659	0.922	-39.9#	150	0.00
12 TP	Freon-113	0.403	0.433	-7.4	110	0.00
13 TP	Iodomethane	0.335	0.261	22.1#	95	0.00
14 TP	Acrolein	0.023	0.022	4.3	105	0.00
15 TP	Methylene chloride	0.233	0.221	5.2	103	0.00
17 TP	Acetone	0.047	0.044	6.4	95	0.00
18 TP	trans-1,2-Dichloroethene	0.242	0.231	4.5	104	0.00
19 TP	Methyl acetate	0.109	0.113	-3.7	121	0.00
20 TP	Methyl tert-butyl ether	0.445	0.498	-11.9	128	0.00
21 TP	tert-Butyl alcohol	0.010	0.012	-20.0	147	0.00
22 TP	Diisopropyl ether	0.829	0.809	2.4	110	0.00
23 TP	1,1-Dichloroethane	0.451	0.444	1.6	105	0.00
24 TP	Halothane	0.186	0.191	-2.7	112	0.00
25 TP	Acrylonitrile	0.046	0.048	-4.3	120	0.00
26 TP	Ethyl tert-butyl ether	0.664	0.640	3.6	112	0.00
27 TP	Vinyl acetate	0.458	0.341	25.5#	100	0.00
28 TP	cis-1,2-Dichloroethene	0.269	0.248	7.8	101	0.00
29 TP	2,2-Dichloropropane	0.350	0.316	9.7	99	0.00
30 TP	Bromochloromethane	0.125	0.121	3.2	107	0.00
31 TP	Cyclohexane	0.466	0.424	9.0	102	0.00
32 TC	Chloroform	0.442	0.424	4.1	104	0.00
33 TP	Ethyl acetate	0.155	0.158	-1.9	121	0.00
34 TP	Carbon tetrachloride	0.329	0.315	4.3	111	0.00
35 TP	Tetrahydrofuran	0.044	0.045	-2.3	113	0.00
36 S	Dibromofluoromethane	0.279	0.277	0.7	104	0.00
37 TP	1,1,1-Trichloroethane	0.371	0.382	-3.0	112	0.00
39 TP	2-Butanone	0.062	0.058	6.5	111	0.00
40 TP	1,1-Dichloropropene	0.328	0.331	-0.9	108	0.00
41 TP	Benzene	1.016	0.900	11.4	101	0.00
42 TP	tert-Amyl methyl ether	0.540	0.525	2.8	112	0.00

Evaluate Continuing Calibration Report

Data Path : I:\VOLATILES\VOA130\2022\221221AICAL\
 Data File : V30221221A17.D
 Acq On : 21 Dec 2022 04:43 pm
 Operator : VOA130:PID
 Sample : C8260STD10PPB
 Misc : WG1726700,ICAL
 ALS Vial : 17 Sample Multiplier: 1

Quant Time: Dec 22 13:35:32 2022
 Quant Method : I:\VOLATILES\VOA130\2022\221221AICAL\VOA130_221221A_8260.m
 Quant Title : VOLATILES BY GC/MS
 QLast Update : Thu Dec 22 13:31:59 2022
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 20% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Fluorobenzene	1.000	1.000	0.0	96	0.00
2 TP	Dichlorodifluoromethane	0.233	0.264	-13.3	103	0.00
3 TP	Chloromethane	0.240	0.257	-7.1	104	0.00
4 TC	Vinyl chloride	0.248	0.271	-9.3	106	0.00
5 TP	Bromomethane	0.206	0.236	-14.6	112	0.00
6 TP	Chloroethane	0.178	0.194	-9.0	109	0.00
7 TP	Trichlorofluoromethane	0.424	0.478	-12.7	106	0.00
8 TP	Ethyl ether	0.089	0.116	-30.3#	119	0.00
10 TC	1,1-Dichloroethene	0.236	0.246	-4.2	103	0.00
11 TP	Carbon disulfide	0.411	0.569	-38.4#	133	0.00
12 TP	Freon-113	0.257	0.271	-5.4	99	0.00
13 TP	Iodomethane	0.258	0.219	15.1	99	0.00
14 TP	Acrolein	* 10.000	8.511	14.9	88	0.00
15 TP	Methylene chloride	0.258	0.268	-3.9	99	0.00
17 TP	Acetone	* 10.000	9.869	1.3	101	0.00
18 TP	trans-1,2-Dichloroethene	0.249	0.265	-6.4	99	0.00
19 TP	Methyl acetate	0.077	0.083	-7.8	104	0.00
20 TP	Methyl tert-butyl ether	0.437	0.476	-8.9	106	0.00
21 TP	tert-Butyl alcohol	0.00993	0.00926#	6.7	90	0.00
22 TP	Diisopropyl ether	0.639	0.584	8.6	96	0.00
23 TP	1,1-Dichloroethane	0.394	0.418	-6.1	102	0.00
24 TP	Halothane	0.171	0.184	-7.6	102	0.00
25 TP	Acrylonitrile	0.038	0.039	-2.6	103	0.00
26 TP	Ethyl tert-butyl ether	0.528	0.488	7.6	97	0.00
27 TP	Vinyl acetate	0.298	0.215	27.9#	81	0.00
28 TP	cis-1,2-Dichloroethene	0.263	0.267	-1.5	98	0.00
29 TP	2,2-Dichloropropane	0.355	0.345	2.8	95	0.00
30 TP	Bromochloromethane	0.120	0.123	-2.5	96	0.00
31 TP	Cyclohexane	0.376	0.378	-0.5	96	0.00
32 TC	Chloroform	0.438	0.465	-6.2	102	0.00
33 TP	Ethyl acetate	0.094	0.091	3.2	101	0.00
34 TP	Carbon tetrachloride	0.334	0.342	-2.4	99	0.00
35 TP	Tetrahydrofuran	0.027	0.027	0.0	95	0.00
36 S	Dibromofluoromethane	0.279	0.286	-2.5	98	0.00
37 TP	1,1,1-Trichloroethane	0.380	0.410	-7.9	102	0.00
39 TP	2-Butanone	0.042	0.036	14.3	84	0.00
40 TP	1,1-Dichloropropene	0.306	0.334	-9.2	103	0.00
41 TP	Benzene	0.922	0.942	-2.2	99	0.00
42 TP	tert-Amyl methyl ether	0.478	0.417	12.8	91	0.00

Calibration Verification Summary

Form 7

Volatiles

Client : C&S Companies
 Project Name : CONVENTUS/MROW
 Instrument ID : VOA122
 Lab File ID : V22221221A01
 Sample No : WG1726668-2
 Channel :

Lab Number : L2270805
 Project Number : 1186
 Calibration Date : 12/21/22 07:16
 Init. Calib. Date(s) : 12/09/22 12/09/22
 Init. Calib. Times : 15:23 19:30

Compound	Ave. RRF	RRF	Min RRF	%D	Max %D	Area%	Dev(min)
Fluorobenzene	1	1	-	0	20	165	0
Dichlorodifluoromethane	0.137	0.155	-	-13.1	20	164	0
Chloromethane	0.223	0.254	-	-13.9	20	168	0
Vinyl chloride	0.162	0.186	-	-14.8	20	168	0
Bromomethane	10	6.748	-	32.5*	20	102	0
Chloroethane	0.1	0.118	-	-18	20	169	0
Trichlorofluoromethane	0.238	0.232	-	2.5	20	141	0
Ethyl ether	0.048	0.045	-	6.3	20	144	0
1,1-Dichloroethene	0.138	0.145	-	-5.1	20	157	0
Carbon disulfide	0.264	0.293	-	-11	20	165	0
Freon-113	0.133	0.143	-	-7.5	20	148	0
Acrolein	0.017	0.019	-	-11.8	20	167	0
Methylene chloride	0.15	0.158	-	-5.3	20	160	0
Acetone	10	9.967	-	0.3	20	151	0
trans-1,2-Dichloroethene	0.151	0.15	-	0.7	20	148	0
Methyl acetate	10	9.896	-	1	20	142	0
Methyl tert-butyl ether	0.23	0.174	-	24.3*	20	122	0
tert-Butyl alcohol	0.0081	0.00664*	-	18	20	127	-0.01
Diisopropyl ether	0.42	0.387	-	7.9	20	151	0
1,1-Dichloroethane	0.317	0.33	-	-4.1	20	153	0
Halothane	0.101	0.097	-	4	20	143	0
Acrylonitrile	0.038	0.039	-	-2.6	20	153	0
Ethyl tert-butyl ether	0.361	0.287	-	20.5*	20	137	0
Vinyl acetate	0.214	0.188	-	12.1	20	160	0
cis-1,2-Dichloroethene	0.122	0.115*	-	5.7	20	144	0
2,2-Dichloropropane	0.245	0.239	-	2.4	20	143	0
Bromochloromethane	0.07	0.063*	-	10	20	136	0
Cyclohexane	0.327	0.348	-	-6.4	20	158	0
Chloroform	0.299	0.278	-	7	20	150	0
Ethyl acetate	0.076	0.061	-	19.7	20	144	0
Carbon tetrachloride	0.229	0.186	-	18.8	20	118	0
Tetrahydrofuran	10	9.009	-	9.9	20	146	0
Dibromofluoromethane	0.313	0.31	-	1	20	159	0
1,1,1-Trichloroethane	0.257	0.22	-	14.4	20	129	0
2-Butanone	0.035	0.029	-	17.1	20	140	0
1,1-Dichloropropene	0.187	0.169	-	9.6	20	141	0
Benzene	0.519	0.496	-	4.4	20	149	0
tert-Amyl methyl ether	0.26	0.201	-	22.7*	20	129	0
1,2-Dichloroethane-d4	0.361	0.326	-	9.7	20	141	0
1,2-Dichloroethane	0.212	0.168	-	20.8*	20	128	0
Methyl cyclohexane	0.231	0.226	-	2.2	20	153	0
Trichloroethene	0.141	0.119*	-	15.6	20	135	0
Dibromomethane	0.079	0.07	-	11.4	20	136	0

* Value outside of QC limits.



Calibration Verification Summary

Form 7

Volatiles

Client : C&S Companies
 Project Name : CONVENTUS/MROW
 Instrument ID : VOA122
 Lab File ID : V22221221A01
 Sample No : WG1726668-2
 Channel :

Lab Number : L2270805
 Project Number : 1186
 Calibration Date : 12/21/22 07:16
 Init. Calib. Date(s) : 12/09/22 12/09/22
 Init. Calib. Times : 15:23 19:30

Compound	Ave. RRF	RRF	Min RRF	%D	Max %D	Area%	Dev(min)
1,2-Dichloropropane	0.147	0.14	-	4.8	20	151	0
Bromodichloromethane	0.209	0.173*	-	17.2	20	133	0
1,4-Dioxane	0.00083	0.00078*	-	6	20	139	0
cis-1,3-Dichloropropene	0.212	0.179*	-	15.6	20	139	0
Chlorobenzene-d5	1	1	-	0	20	157	0
Toluene-d8	1.313	1.335	-	-1.7	20	159	0
Toluene	0.477	0.431	-	9.6	20	144	0
4-Methyl-2-pentanone	0.039	0.034	-	12.8	20	140	0
Tetrachloroethene	0.191	0.17	-	11	20	137	0
trans-1,3-Dichloropropene	0.246	0.19*	-	22.8*	20	130	0
Ethyl methacrylate	0.139	0.119	-	14.4	20	133	0
1,1,2-Trichloroethane	0.102	0.08*	-	21.6*	20	136	0
Chlorodibromomethane	0.157	0.115*	-	26.8*	20	121	0
1,3-Dichloropropane	0.215	0.182	-	15.3	20	142	0
1,2-Dibromoethane	0.117	0.093*	-	20.5*	20	133	0
2-Hexanone	0.065	0.058	-	10.8	20	136	0
Chlorobenzene	0.483	0.445	-	7.9	20	141	0
Ethylbenzene	0.901	0.868	-	3.7	20	143	0
1,1,1,2-Tetrachloroethane	0.157	0.119	-	24.2*	20	128	0
p/m Xylene	0.344	0.34	-	1.2	20	144	0
o Xylene	0.321	0.315	-	1.9	20	146	0
Styrene	0.524	0.507	-	3.2	20	143	0
1,4-Dichlorobenzene-d4	1	1	-	0	20	147	0
Bromoform	0.171	0.123	-	28.1*	20	114	0
Isopropylbenzene	1.675	1.726	-	-3	20	142	0
4-Bromofluorobenzene	0.983	1.031	-	-4.9	20	144	0
Bromobenzene	0.426	0.4	-	6.1	20	135	0
n-Propylbenzene	2.21	2.442	-	-10.5	20	148	0
1,4-Dichlorobutane	0.432	0.401	-	7.2	20	140	0
1,1,2,2-Tetrachloroethane	0.242	0.221	-	8.7	20	146	0
4-Ethyltoluene	1.652	1.709	-	-3.5	20	144	0
2-Chlorotoluene	1.196	1.239	-	-3.6	20	143	0
1,3,5-Trimethylbenzene	1.44	1.42	-	1.4	20	136	0
1,2,3-Trichloropropane	0.203	0.182	-	10.3	20	136	0
trans-1,4-Dichloro-2-buten	0.094	0.081	-	13.8	20	131	0
4-Chlorotoluene	1.266	1.294	-	-2.2	20	143	0
tert-Butylbenzene	1.185	1.183	-	0.2	20	138	0
1,2,4-Trimethylbenzene	1.462	1.387	-	5.1	20	132	0
sec-Butylbenzene	1.714	1.821	-	-6.2	20	145	0
p-Isopropyltoluene	1.464	1.514	-	-3.4	20	140	0
1,3-Dichlorobenzene	0.726	0.716	-	1.4	20	140	0
1,4-Dichlorobenzene	0.734	0.696	-	5.2	20	137	0
p-Diethylbenzene	0.857	0.856	-	0.1	20	137	0

* Value outside of QC limits.



Calibration Verification Summary

Form 7

Volatiles

Client : C&S Companies
 Project Name : CONVENTUS/MROW
 Instrument ID : VOA101
 Lab File ID : V01221223A02
 Sample No : WG1727559-2
 Channel :

Lab Number : L2270805
 Project Number : 1186
 Calibration Date : 12/23/22 07:41
 Init. Calib. Date(s) : 12/16/22 12/16/22
 Init. Calib. Times : 12:37 16:31

Compound	Ave. RRF	RRF	Min RRF	%D	Max %D	Area%	Dev(min)
Fluorobenzene	1	1	-	0	20	103	0
Dichlorodifluoromethane	0.227	0.241	-	-6.2	20	109	0
Chloromethane	0.264	0.255	-	3.4	20	99	0
Vinyl chloride	0.369	0.393	-	-6.5	20	105	0
Bromomethane	0.252	0.175	-	30.6	20	84	0
Chloroethane	0.301	0.334	-	-11	20	113	0
Trichlorofluoromethane	0.64	0.71	-	-10.9	20	116	0
Ethyl ether	0.152	0.151	-	0.7	20	109	0
1,1-Dichloroethene	0.359	0.38	-	-5.8	20	112	0
Carbon disulfide	0.659	0.664	-	-0.8	20	108	0
Freon-113	0.403	0.447	-	-10.9	20	113	0
Acrolein	0.023	0.025	-	-8.7	20	121	0
Methylene chloride	0.233	0.229	-	1.7	20	106	0
Acetone	0.047	0.05	-	-6.4	20	106	0
trans-1,2-Dichloroethene	0.242	0.247	-	-2.1	20	111	0
Methyl acetate	0.109	0.118	-	-8.3	20	125	0
Methyl tert-butyl ether	0.445	0.446	-	-0.2	20	114	0
tert-Butyl alcohol	0.01	0.011	-	-10	20	138	0
Diisopropyl ether	0.829	0.897	-	-8.2	20	121	0
1,1-Dichloroethane	0.451	0.479	-	-6.2	20	113	0
Halothane	0.186	0.194	-	-4.3	20	113	0
Acrylonitrile	0.046	0.047	-	-2.2	20	116	0
Ethyl tert-butyl ether	0.664	0.683	-	-2.9	20	119	0
Vinyl acetate	0.458	0.487	-	-6.3	20	143	0
cis-1,2-Dichloroethene	0.269	0.274	-	-1.9	20	111	0
2,2-Dichloropropane	0.35	0.392	-	-12	20	122	0
Bromochloromethane	0.125	0.123	-	1.6	20	108	0
Cyclohexane	0.466	0.491	-	-5.4	20	118	0
Chloroform	0.442	0.454	-	-2.7	20	111	0
Ethyl acetate	0.155	0.165	-	-6.5	20	126	0
Carbon tetrachloride	0.329	0.367	-	-11.6	20	128	0
Tetrahydrofuran	0.044	0.048	-	-9.1	20	119	0
Dibromofluoromethane	0.279	0.277	-	0.7	20	104	0
1,1,1-Trichloroethane	0.371	0.405	-	-9.2	20	118	0
2-Butanone	0.062	0.067	-	-8.1	20	126	0
1,1-Dichloropropene	0.328	0.363	-	-10.7	20	118	0
Benzene	1.016	0.989	-	2.7	20	111	0
tert-Amyl methyl ether	0.54	0.541	-	-0.2	20	115	0
1,2-Dichloroethane-d4	0.305	0.309	-	-1.3	20	105	0
1,2-Dichloroethane	0.304	0.323	-	-6.3	20	117	0
Methyl cyclohexane	0.433	0.442	-	-2.1	20	117	0
Trichloroethene	0.278	0.28	-	-0.7	20	111	0
Dibromomethane	0.135	0.138	-	-2.2	20	113	0

* Value outside of QC limits.



Appendix C

Validator Qualifications

KENNETH R. APPLIN

Geochemist/Data Validator

Ph.D., Geochemistry and Mineralogy, The Pennsylvania State University

M.S., Geochemistry and Mineralogy, The Pennsylvania State University

B.A., Geological Sciences, SUNY at Geneseo, NY

Dr. Applin has over 35 years of experience working with the geochemistry of natural waters. His prior experience includes working as an Assistant Professor of Geology at the University of Missouri-Columbia and as Chief Hydrogeologist and Geochemist with a leading engineering firm in Rochester, NY. In 1993, he established KR Applin and Associates, a small consulting business that focuses on the geochemistry of natural waters, especially as applied to problems involving the contamination of groundwater and surface water.

Dr. Applin is also an experienced analytical data validator and has provided data validation services since 1994 to a variety of clients performing brownfield cleanup projects, hazardous waste remediation, groundwater monitoring at solid waste facilities, and other projects requiring third-party data validation. Dr. Applin has several years of hands-on experience with the laboratory analysis of natural waters and has successfully completed the USEPA Region II certification courses for performing inorganic and organic analytical data validation.

MICHAEL K. PERRY

Chemist/Data Validator

B.S. Chemistry, Georgia State University, Atlanta, GA

A.A.S., Chemical Technology, Alfred State College, Alfred, NY

Mr. Perry has over 30 years of experience in the analytical laboratory business. During his early career, he spent several years as a laboratory analyst performing the analysis of soil, water, and air samples for inorganic and organic chemical parameters. During his last 20 years in the environmental laboratory business, he managed and directed two major analytical laboratories in Rochester, NY. His management responsibilities included oversight of the daily operations of the lab, staff training and supervision, the selection, purchase, and maintenance of analytical instruments, the introduction of new laboratory methods, analytical quality assurance and quality control, data acquisition and management, and other business-related activities.

Mr. Perry has an extensive working knowledge of the methods and procedures used for sampling and analyzing both inorganic and organic analytes in soil, water, and air. He is an accomplished laboratory chemist and is familiar with the analytical methods and procedures established under the USEPA Contract Laboratory Protocols (CLP), the NYSDEC Analytical Services Protocols (ASP), and the NYSDOH Environmental Laboratory Approval Program (ELAP).

APPENDIX B

INSTITUTIONAL AND ENGINEERING CONTROLS CERTIFICATION FORM



Enclosure 2
NEW YORK STATE DEPARTMENT OF ENVIRONMENTAL CONSERVATION
Site Management Periodic Review Report Notice
Institutional and Engineering Controls Certification Form



Site No. **C915260** **Site Details** **Box 1**

Site Name Former Mobil Service Station 99-MST

Site Address: 979 Main Street Zip Code: 14203
City/Town: Buffalo
County: Erie
Site Acreage: 1.725

Reporting Period: March 24, 2022 to March 24, 2023

- | | YES | NO |
|--|--------------------------|-------------------------------------|
| 1. Is the information above correct? | ☐ | ☒ |
| If NO, include handwritten above or on a separate sheet. | | |
| 2. Has some or all of the site property been sold, subdivided, merged, or undergone a tax map amendment during this Reporting Period? | ☐ | ☒ |
| 3. Has there been any change of use at the site during this Reporting Period (see 6NYCRR 375-1.11(d))? | ☐ | ☒ |
| 4. Have any federal, state, and/or local permits (e.g., building, discharge) been issued for or at the property during this Reporting Period? | ☐ | ☒ |
| If you answered YES to questions 2 thru 4, include documentation or evidence that documentation has been previously submitted with this certification form. | | |
| 5. Is the site currently undergoing development? | ☐ | ☒ |

- | | Box 2 | |
|---|-------------------------------------|--------------------------|
| | YES | NO |
| 6. Is the current site use consistent with the use(s) listed below?
Restricted-Residential, Commercial, and Industrial | ☒ | ☐ |
| 7. Are all ICs in place and functioning as designed? | ☒ | ☐ |

**IF THE ANSWER TO EITHER QUESTION 6 OR 7 IS NO, sign and date below and
DO NOT COMPLETE THE REST OF THIS FORM. Otherwise continue.**

A Corrective Measures Work Plan must be submitted along with this form to address these issues.

Signature of Owner, Remedial Party or Designated Representative

Date

Box 2A

YES NO

8. Has any new information revealed that assumptions made in the Qualitative Exposure Assessment regarding offsite contamination are no longer valid?

☐ YES ☒ NO

If you answered YES to question 8, include documentation or evidence that documentation has been previously submitted with this certification form.

9. Are the assumptions in the Qualitative Exposure Assessment still valid?
(The Qualitative Exposure Assessment must be certified every five years)

☐ YES ☒ NO

If you answered NO to question 9, the Periodic Review Report must include an updated Qualitative Exposure Assessment based on the new assumptions.

SITE NO. C915260

See Section 1.1

Box 3**Description of Institutional Controls**ParcelOwnerInstitutional Control**100.79-1-1.1**

Kaleida Properties, Inc.

Ground Water Use Restriction
Soil Management Plan
Landuse Restriction
Monitoring Plan
Site Management Plan
IC/EC Plan

1. Prohibition of use of groundwater.
2. Landuse Restriction for Restricted Residential, Commercial or Industrial use.
3. Soil Management or Excavation Work Plan for any future intrusive work.
4. Groundwater Monitoring Plan.

100.79-1-1.1/2

Seavest Core Buffalo Conventus, LLC

Ground Water Use Restriction
Soil Management Plan
Landuse Restriction
Monitoring Plan
Site Management Plan
IC/EC Plan

1. Prohibition of use of groundwater.
2. Landuse Restriction for Restricted Residential, Commercial or Industrial use.
3. Soil Management or Excavation Work Plan for any future intrusive work.
4. Groundwater Monitoring Plan.

100.79-1-2.11

Kaleida Health

Ground Water Use Restriction
Soil Management Plan
Landuse Restriction
Monitoring Plan
Site Management Plan
IC/EC Plan

1. Prohibition of use of groundwater.
2. Landuse Restriction for Restricted Residential, Commercial or Industrial use.
3. Soil Management or Excavation Work Plan for any future intrusive work.
4. Groundwater Monitoring Plan.

Box 4**Description of Engineering Controls**

Parcel

Engineering Control

100.79-1-1.1

Groundwater Treatment System

Groundwater will be treated in-situ by injections of oxygen release compounds (ORC) to degrade petroleum hydrocarbons to harmless compounds.

100.79-1-1.1/2

Groundwater Treatment System

Groundwater will be treated in-situ by injections of oxygen release compounds (ORC) to degrade petroleum hydrocarbons to harmless compounds.

100.79-1-2.11

Groundwater Treatment System

Groundwater will be treated in-situ by injections of oxygen release compounds (ORC) to degrade petroleum hydrocarbons to harmless compounds

Box 5

Periodic Review Report (PRR) Certification Statements

1. I certify by checking "YES" below that:

a) the Periodic Review report and all attachments were prepared under the direction of, and reviewed by, the party making the Engineering Control certification;

b) to the best of my knowledge and belief, the work and conclusions described in this certification are in accordance with the requirements of the site remedial program, and generally accepted engineering practices; and the information presented is accurate and complete.

YES NO

☒ ☐

2. For each Engineering control listed in Box 4, I certify by checking "YES" below that all of the following statements are true:

(a) The Engineering Control(s) employed at this site is unchanged since the date that the Control was put in-place, or was last approved by the Department;

(b) nothing has occurred that would impair the ability of such Control, to protect public health and the environment;

(c) access to the site will continue to be provided to the Department, to evaluate the remedy, including access to evaluate the continued maintenance of this Control;

(d) nothing has occurred that would constitute a violation or failure to comply with the Site Management Plan for this Control; and

(e) if a financial assurance mechanism is required by the oversight document for the site, the mechanism remains valid and sufficient for its intended purpose established in the document.

YES NO

☒ ☐

IF THE ANSWER TO QUESTION 2 IS NO, sign and date below and DO NOT COMPLETE THE REST OF THIS FORM. Otherwise continue.

A Corrective Measures Work Plan must be submitted along with this form to address these issues.

Signature of Owner, Remedial Party or Designated Representative

Date

**IC CERTIFICATIONS
SITE NO. C915260**

Box 6

SITE OWNER OR DESIGNATED REPRESENTATIVE SIGNATURE

I certify that all information and statements in Boxes 1,2, and 3 are true. I understand that a false statement made herein is punishable as a Class "A" misdemeanor, pursuant to Section 210.45 of the Penal Law.

I _____ at C&S Engineers, Inc. 141 Elm Street, Buffalo, NY 14203 _____,
print name print business address

Representative for Kaleida Health, Kaledia Properties, Inc.
am certifying as and Seavest Core Buffalo Conventus, LLC (Owner or Remedial Party)

for the Site named in the Site Details Section of this form.

Signature of Owner, Remedial Party, or Designated Representative
Rendering Certification

Date

EC CERTIFICATIONS

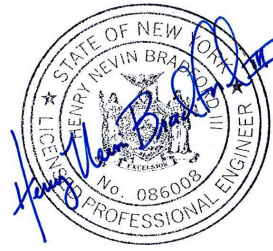
Box 7

Professional Engineer Signature

I certify that all information in Boxes 4 and 5 are true. I understand that a false statement made herein is punishable as a Class "A" misdemeanor, pursuant to Section 210.45 of the Penal Law.

I H. Nevin Bradford at C&S Engineers, Inc. 141 Elm Street, Buffalo, NY 14203,
print name print business address

Kaleida Health, Kaledia Properties, Inc.
am certifying as a Professional Engineer for the and Seavest Core Buffalo Conventus, LLC
(Owner or Remedial Party)



Signature of Professional Engineer, for the Owner or
Remedial Party, Rendering Certification

Stamp
(Required for PE)

April 21, 2023

Date