

2020 Site Management

Periodic Review Report -

Love Canal Site

NYSDEC Site No. 932020

Niagara Falls, New York

Glenn Springs Holdings, Inc.





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1. Introduction

Operation of the Love Canal Site (Site) was transferred from the New York State Department of Environmental Conservation (NYSDEC) to Occidental Chemical Corporation (OCC) in April 1995. Effective July 1, 1998, Site responsibility was assigned by OCC to Glenn Springs Holdings, Inc. (GSH), an affiliate of OCC. Since October 1, 2008, GHD Services, Inc. (GHD), formerly Conestoga-Rovers & Associates (CRA), has performed operation, maintenance, monitoring, and reporting activities for the Site under contract to and direct management of GSH.

This report is the twenty-sixth annual report prepared by or on behalf of OCC and covers operation, maintenance, and monitoring activities for 2020. The completed 2020 NYSDEC Institutional and Engineering Controls Certification Form is included as Appendix A.

2. Remedial Systems

Operation of remedial systems to prevent the off-Site migration of chemical contaminants from the Site began in October 1978 with the installation of a barrier drain along the east and west sides of the South Sector of the Canal. The barrier drain was later extended to completely encompass the entire area of disposed waste within the Central and North Sectors of the Canal. The barrier drain, designed to intercept the shallow overburden lateral groundwater flow, consists of a trench approximately 4 feet wide that varies in depth from approximately 12 to 25 feet deep depending on location at the Site. Installed within the trench is a perforated vitrified clay tile pipe. The pipe is 6-inch diameter in the Central and North Sectors and both 6-inch and 8-inch diameter in the South Sector. The pipe is centered in a minimum of 2 feet of uniformly sized gravel, which is overlain with coarse sand extending to the existing ground surface present at the time of construction. Thirty-two lateral trenches, approximately 12 to 19 feet deep, filled with a minimum of 2 feet of gravel and overlain with sand similar to the barrier drain, were dug perpendicular to the barrier drain in the direction of the Canal. The majority of these laterals extend into the disposed waste. The barrier drain is graded from two highpoints, one in the southeast corner and the other in the northeast corner, toward a series of manholes which drain to four pump chambers (PC-1A/PC-2A in the North/Central Sector and PC-1/PC-2 in the South Sector) where the leachate is collected. The collected leachate is pumped from the four pump chambers to two other pump chambers connected to underground holding tanks (PC-3A in the North/Central Sector and PC-3 in the South Sector) where it is temporarily stored. From that point, the leachate is pumped to the on-Site Love Canal Treatment Facility (LCTF) where it is treated and discharged to the Niagara Falls Water Board (NFWB) sanitary sewer system under the Site's Significant Industrial User (SIU) Permit #44. The locations of the remedial system components are illustrated on the Site Plan presented as Figure 2.1.

The installation of a 22-acre clay cap over the entire former Canal area was completed in October 1980 following completion of the barrier drain collection system. The purpose of the cap is to reduce infiltration of precipitation. The thickness of the clay cap is a minimum of 3 feet. In 1985, a second (40-acre) cap was installed over the initial clay cap area. The newer cap consists of a 40-mil high density polyethylene (HDPE) liner covered by 18 inches of clean soil and vegetation.



In March 1999, the adjacent 102nd Street Landfill Site leachate collection system was connected to the Love Canal Site to facilitate the transfer of leachate from the 102nd Street landfill into Love Canal's pump chamber PC-3 for treatment at the LCTF.

2.1 Operations of the Barrier Drain and Collection System

2.1.1 Barrier Drain System

The barrier drain system continues to function as designed, with no major maintenance required during 2020. Semiannual inspections of the barrier drain components, including manholes and pump chambers, are required by the Site's NYSDEC-approved Operation and Maintenance (O&M) Manual (CRA, revised March 2015). Inspections of the barrier drain manholes were conducted on May 20 and October 19, 2020, and inspections of the barrier drain pump chambers were carried out on May 20 and November 10, 2020. The visual inspections showed that the manhole flumes were flowing freely and required no further maintenance. During both inspection events, limited buildup of sludge was noted at MH-6A, MH-6B, and MH-6C. However, the buildup was insufficient to warrant cleaning, as it did not impede flow through the manholes. The visual inspections were documented on the 2020 Semiannual Inspection Forms, which are presented in Appendix B. The manhole locations are presented on Figure 2.2.

2.1.2 Pumping System

The barrier drain system consists of two sectors, the Northern/Central and the Southern. Leachate from the Northern/Central Sector drains to pump chambers PC-1A and PC-2A where it is pumped to pump chamber PC-3A, while leachate from the Southern Sector is pumped from pump chambers PC-1 and PC-2 to the underground storage tank connected to pump chamber PC-3. From pump chambers PC-3 and PC-3A, the leachate is then transferred through a below ground metering chamber outside the LCTF on the southeast corner of the building and then into the LCTF for treatment. The pumping system is designed to operate continuously and was operational and functioned as designed throughout 2020.

2.1.3 102nd Street Landfill Forcemain

The leachate forcemain construction was completed in March 1999 and is used for the transfer of leachate from the 102nd Street Landfill to the LCTF. The forcemain begins at the northwest corner of the 102nd Street Landfill and extends northward beneath River Road, LaSalle Expressway, and Frontier Avenue to pump chamber PC-3 at the Site. During 2020, the leachate collection system at 102nd Street pumped 120,131 gallons of leachate to the LCTF.



Groundwater Treatment and Monitoring

3.1 Groundwater Treatment

3.1.1 Treatment System

The LCTF consists of clarification, bag filtration, and carbon treatment prior to discharge to the NFWB sanitary sewer system. A process schematic depicting the layout of the treatment system is presented as Figure 3.1.

Treated water from the Site is discharged essentially on a batch basis to the NFWB sanitary sewer system (i.e., when there is sufficient water in storage, the treatment system is operated); however, under seasonal high flow conditions, water is discharged continuously, The discharge is authorized under the Site's SIU Permit #44. The current permit is valid from January 10, 2020 to January 9, 2025. A copy of the NFWB permit is included as Appendix C.

3.1.2 Effluent Discharge

The LCTF discharged to the NFWB sanitary sewer system on 229 days in 2020.

Under high stormwater flow events, the NFWB periodically requires that the LCTF temporarily cease discharging to the sewer system. During an event of this type, the barrier drain pumping system will continue to operate and maintain a protective inward hydraulic gradient to capture leachate. The NFWB did not require the LCTF to temporarily cease discharging to the sewer system during 2020.

In 2020, the LCTF processed a total of 3,356,298 gallons of leachate. This total was comprised of 3,236,167 gallons of leachate from the Site and 120,131 gallons of leachate from the 102nd Street Landfill.

Table 3.1 shows the monthly total and average treated groundwater quantities from 2000 through 2020.

3.1.3 Effluent Sampling

Sampling of the effluent discharged to the NFWB sanitary sewer system occurred quarterly as required under the Site's SIU Discharge Permit #44. In accordance with the SIU permit, the quarterly monitoring periods for 2020 were as follows:

Quarter 1: December 1 – February 29 Quarter 2: March 1 – May 31

Quarter 3: June 1 – August 31 Quarter 4: September 1 – November 30

The quarterly effluent sampling for 2020 was performed on December 9, 2019 and March 16, June 8, and September 1, 2020. The sample results were submitted to the NFWB quarterly as required by the permit and to the NYSDEC. The results for each event were in compliance with the requirements of the Site's SIU permit.



3.1.4 Precipitation

In 2020, precipitation in the Niagara Falls region totaled 24.46 inches (Niagara Falls International Airport, National Climatic Data Center). Table 3.1 provides historical regional precipitation data from 2000 through 2020.

3.2 Groundwater Monitoring

Groundwater monitoring consists of both chemical monitoring to determine groundwater quality and hydraulic monitoring to demonstrate that the barrier drain is creating hydraulic containment. Monitoring and analytical protocols for the Site's groundwater monitoring program have been established and are set forth in the "Sampling Manual, Love Canal Site, Long-Term Groundwater Monitoring Program" (LTGMP), revised June 2013.

The monitoring results for 2020 are presented in the following sections.

3.2.1 Groundwater Quality

Chemical monitoring is performed annually by sampling select overburden and bedrock monitoring wells at the Site. On March 25, 2009, the NYSDEC communicated via email to GSH that the NYSDEC would no longer be providing an annual well sampling list for chemical monitoring and directed GSH to use the wells sampled in 2007 and 2008 for all future sampling events. Subsequent discussions between GSH and the NYSDEC regarding the well sampling list led to this decision being documented in an August 5, 2010 memo titled "Love Canal Annual Groundwater Sampling Schedule", presented in Appendix D.

It should be noted that Overburden Well 3151 is included on the Appendix D list; however, this well was noted in 2007 as "Well no longer available – destroyed" and could not be located. Therefore, this well has not been sampled since 2007. In addition, Overburden Well 10178A and Bedrock Wells MW-01 and MW-02 were added to the annual sampling program in 2011. The sampling frequency for Overburden Well 10178A became annual in 2016.

The 2020 annual groundwater chemical monitoring event was performed between June 24 and July 10, 2020. As part of the annual groundwater chemical monitoring in 2020, 33 monitoring wells were sampled, comprised of 13 overburden and 20 bedrock monitoring wells. In addition to the 12 overburden wells scheduled to be sampled in 2020 per the Appendix D list, the NYSDEC requested (via email on June 30, 2020) that a groundwater sample be collected from Overburden Well 10178B during the 2020 sampling event and analyzed for the same parameters as the other wells sampled. Well 10178B was requested to be sampled to provide additional data on groundwater quality in the overburden west of Well 10135.

As a component of the LTGMP, the NYSDEC has the option of collecting split samples during the annual sampling event and having those samples analyzed independently to verify data. However, based on a review of split sample data from 1995 to 2013, NYSDEC determined that there was no difference between the data sets and, therefore, split sampling was no longer required. No split samples were collected by the NYSDEC during the 2020 annual sampling event. The NYSDEC observed a portion of the groundwater sampling activities conducted on July 7, 2020.



Groundwater samples were submitted to ALS Environmental Group USA, Corp (ALS) located in Rochester, New York. ALS is a New York State Department of Health (NYSDOH) approved laboratory certified under the National Environmental Laboratory Approval Program (NELAP). The samples were analyzed for Site-specific volatile organic compounds (VOCs), semi-volatile organic compounds (SVOCs), pesticides, and polychlorinated biphenyls (PCBs). The raw data laboratory packages are presented in Appendix E. A GHD chemist performed the analytical Quality Assurance/Quality Control (QA/QC) review and data validation. The QA/QC report for this event is presented in Appendix F.

Figure 3.2 identifies the wells sampled and their locations. The Love Canal Annual Groundwater Sampling Schedule is presented in Appendix D. Table 3.2 provides a summary of the overburden wells that were sampled, the analytical data, and a summation of the number of compounds found at or above the detection limits in each well. Table 3.3 provides a summary of the bedrock wells that were sampled, the analytical data, and a summation of the number of compounds found at or above the detection limits in each well. Table 3.4 presents the number of discrete compounds detected in each individual well sampled, arranged by parameter group. This table also presents the total number of detections in the overburden and in the bedrock, as well as the total number of discrete compounds detected in the overburden in its entirety and the total number of discrete compounds detected in the bedrock in its entirety.

3.2.1.1 Overburden Monitoring Wells

The 2020 groundwater analytical results for the overburden monitoring wells (Table 3.2) are consistent with previous long-term monitoring analytical results. The analytical results were non-detect or were detected at low levels consistent with concentrations detected during previous monitoring events (with the exception of groundwater from Well 10135, which is installed in an area of known Site impacts).

Historically, Well 10135 has had the most detected compounds and the highest concentrations detected. This well is located in the southwestern portion of the Site and within the fenced boundaries of the Site. Although located outside the barrier drain, Well 10135 is within the influence of the barrier drain based on hydraulic monitoring conducted at Well 10135 and at adjacent nested piezometer string 1160, as demonstrated in Table 3.7 and on Figures 3.5 (1160 cross-section) and 3.9 (groundwater contours for overburden), and as further discussed in Section 3.2.2.

As stated in the LTGMP, Well 10135, located in an isolated area of known contamination, is sampled as an indicator well. A trend analysis of contaminants detected in Well 10135 for the years 1990 through 2019 was performed and presented in the 2019 PRR. Based on the demonstrated long-term stability of parameter concentrations in Well 10135 and its presence within the influence of the barrier drain, it was concluded that no further trend analysis for this well was warranted.

As indicated in Section 3.2.1, Well 10178B was sampled during the 2020 sampling event at the request of the NYSDEC to provide additional data on groundwater quality in the overburden west of Well 10135. No VOCs or PCBs were detected in Well 10178B. One SVOC (bis(2-ethylhexyl)phthalate) and four pesticides (alpha-BHC, delta-BHC, gamma-BHC, and methoxychlor) were detected at very low concentrations in this well. The reported concentrations of



delta-BHC and methoxychlor were estimated, as they were detected below the laboratory's reporting limits.

Overburden Well 10178A, which did not exhibit any pesticide detections above the laboratory's reporting limits, was measured to be approximately 10.94 feet deep during the 2020 sampling event. The four pesticides detected in Overburden Well 10178B ranged in concentration from 0.021 micrograms per liter (μ g/L) [estimated] (delta-BHC) to 0.069 μ g/L (alpha-BHC). This well was measured to be approximately 20.82 feet deep during the 2020 sampling event. The four pesticides detected in Well 10178B were detected in Overburden Well 10135 at concentrations ranging from 6.5 μ g/L (gamma-BHC) to 27 μ g/L (alpha-BHC). This well was measured to be approximately 29.50 feet deep during the 2020 sampling event. The four pesticides detected in Well 10178B were detected in Bedrock Well 10278 at concentrations ranging from 0.13 μ g/L (delta-BHC) to 0.48 μ g/L (gamma-BHC). This well was measured to be approximately 44.02 feet deep during the 2020 sampling event. All depths were relative to top of casing. Given that Well 10178B is shallower in depth than Well 10135 and that groundwater at Well 10135 flows eastwards into the barrier drain (see Section 3.2.2), it is unlikely that Well 10135 is the source of the pesticide detections in Well 10178B. Based on the very low concentrations of pesticides detected in Well 10178B, the detections are not a potential concern to groundwater quality.

3.2.1.2 Bedrock Monitoring Wells

The 2020 groundwater analytical results for the bedrock monitoring wells (Table 3.3) are consistent with previous long-term monitoring analytical results. Parameter concentrations were either non-detect or detected at low levels consistent with concentrations detected during previous monitoring events. Table 3.4 presents a summary of detections for each well sampled, arranged by parameter group.

3.2.1.3 Historical Compound Detections

Table 3.5 presents a summary of detected compounds of four long-term monitoring wells, consisting of three bedrock wells and one overburden well (Bedrock Wells 10210A, 10210B, and 10210C, and Overburden Well 10135) from 1990 to 2020. The data from these four wells are presented because they have the most consistent historical record of compound detections compared to the other overburden and bedrock wells. As discussed in Sections 3.2.1.1 and 3.2.1.2, the data from the additional Site wells not presented in Table 3.5 are typically non-detect or demonstrate sporadic low level detections and, therefore, do not present useful data in regards to a discussion of historical analytical trends at the Site. An evaluation of the 2020 sampling data for the four wells mentioned above shows that the compounds detected in 2020 are present at sporadic low-level concentrations or concentrations consistent with historical trends.

3.2.2 Hydraulic Containment

Hydraulic monitoring consists of water level measurements conducted quarterly from six nested piezometer strings (1140, 1150, 1160, 1170, 1180, and 1190) per the NYSDEC-approved long-term groundwater monitoring plan (LTGMP), as well as NYSDEC-requested water level measurements collected from four wells (7161, 9130, 9140, and 10135). Well 10135 was added to the hydraulic monitoring program during the first quarter of 2020, as per the United States Environmental Protection Agency's (USEPA's) recommendation in its Fourth Five-Year Review Report for the Site,



in order to demonstrate that this well is within the hydraulic influence of the barrier drain. In 2020, water levels were measured in March, June, September, and December. The water level data for the six nested piezometer strings are presented in Tables 3.6A to 3.6F. The wells on the tables are ordered from left to right on the table, beginning with the well furthest from the outside of the barrier drain to the well inside the area enclosed by the barrier drain. They are also ordered based on screen depth corresponding with the geologic stratum (e.g., fractured clay, soft, clay, and glacial till) as requested by NYSDEC in its letter dated October 21, 2016. Figures 3.3 to 3.8 show the overburden groundwater flow conditions for June 2020 at the six nested piezometers string locations based on geologic stratum consistent with the above.

Piezometer string groundwater elevation data from the remaining three quarters (March, September, and December 2020) demonstrate that the data from those monitoring periods are consistent with the June 2020 data.

In addition to the above-mentioned information, groundwater contour figures were prepared using the June 2020 water levels from the six nested piezometer strings, overburden well 10176C, seven bedrock monitoring wells, and four additional overburden wells (7161, 9130, 9140, and 10135) as requested by NYSDEC. Separate figures were prepared for the overburden and bedrock groundwater elevation contours. The June 2020 groundwater contour figures for the overburden and bedrock are presented as Figures 3.9 and 3.10, respectively.

The overburden groundwater contour figure (Figure 3.9) and Tables 3.6A to 3.6F illustrate that there is a minimum of 1.00 feet of inward gradient outside of the barrier drain at each of the six nested piezometer strings in the various geologic units. The term "feet of inward gradient" means the minimum difference in groundwater elevation between the wells on the outside of the barrier drain and the water level within the barrier drain, with the water level within the barrier drain representing the lowest water level elevation. This demonstrates that groundwater on the outside of the barrier drain (off-Site groundwater) is flowing toward and downward into the barrier drain. Based on the water level data from the six nested piezometer strings, an inward gradient can be inferred to exist around the collection drain system, demonstrating that the horizontal groundwater flow direction outside of the barrier drain is towards the barrier drain. A review of Figure 3.9 shows that groundwater flow on the inside of the barrier drain is also towards the barrier drain. Therefore, the barrier drain and lateral trenches are capturing leachate from the landfill area and a portion of groundwater outside the barrier drain, thereby preventing off-Site migration of chemicals and preventing off-Site groundwater from migrating into the landfill area.

Table 3.7 presents the 2020 quarterly groundwater elevations measured in Well 10135 and in piezometers 1161E, 1163A, and 1165A. The 1160 piezometer string is the closest piezometer string to Well 10135, and piezometers 1161E, 1163A, and 1165A are screened in the same stratigraphic unit as Well 10135 (soft clay). Piezometer 1165A is approximately the same horizontal distance to the barrier drain as Well 10135 is. As shown in Table 3.7, the quarterly groundwater elevations measured in Well 10135 were always lower than the groundwater elevations measured in 1165A in 2020. Therefore, there was an inward gradient extending from Piezometer 1165A through Well 10135 and into the barrier drain during all four quarters of 2020. Based on the inward gradient, chemical impacts identified at Well 10135 are being captured by the barrier drain (i.e., Well 10135 is hydraulically contained).



As demonstrated on Figure 3.10, groundwater in the bedrock was generally flowing from east to west across the Site during the June 2020 hydraulic monitoring event.

Monitoring will continue during 2021 as per the NYSDEC-approved LTGMP, and Well 10135 will continue to be monitored during the quarterly hydraulic monitoring events.

3.2.3 Colvin Boulevard Sewer System NAPL Presence

Overburden Monitoring Well MW-3 was installed on July 1, 2011, within the bedding material of a newly repaired sanitary sewer line on Colvin Boulevard (Figure 2.1). The purpose of this well was to monitor for the presence of residual non-aqueous phase liquid (NAPL) similar to that observed in the bedding material during construction/repair activities. Following well development, MW-3 was monitored for the presence of NAPL on a weekly basis starting on July 19, 2011 and ending on October 7, 2011. No NAPL or visible sheen was detected during these weekly monitoring events. Based on these results, it was concluded that the NAPL that had been observed sporadically during the sewer repair activities was likely limited in volume and mobility. As such, in the report "Colvin Boulevard Sewer Repair Supplemental Subsurface Investigation Report, Colvin Boulevard and 96th Street," dated October 2011, GSH recommended/requested that additional monitoring of MW-3 be continued on a quarterly basis. The NYSDEC approved this request in a letter dated June 5, 2012. The last weekly monitoring event was conducted on August 3, 2012.

Well MW-3 has been monitored on a quarterly basis for the presence of NAPL since November 5, 2012. To date, no NAPL or visible sheen has been detected in this well. As requested by the NYSDEC in an email dated August 10, 2018, the dates of all previous quarterly NAPL checks at MW-3 were included in the 2018 PRR. In 2020, the quarterly NAPL checks at MW-3 were conducted on March 2, June 2, September 2, and December 9. No NAPL or visible sheen was observed.

3.2.4 Well Maintenance

The 2020 well inspections identified the need for routine maintenance on several wells at the Site. Maintenance was conducted during the summer of 2020 and included the following:

Replacement of locks.

3.2.5 Summary of Treatment and Monitoring Results

The volume of effluent discharge from the LCTF decreased from 4,840,275 gallons in 2019 to 3,356,298 gallons in 2020, a number consistent with volumes from previous years with similar precipitation levels. Quarterly sampling and analysis results submitted to both the NFWB and NYSDEC indicated that all chemistry detected in the effluent samples for each event was either non-detect or present at very low levels within historical ranges and well below the Site's SIU Discharge Permit #44 discharge limits.

The inward hydraulic gradient observed at each of the six nested piezometer strings demonstrates that the barrier drain is effectively capturing leachate from the Site and preventing off-Site migration of chemicals. The analytical results from the monitoring wells sampled indicate that compounds were either not detected or were detected at low levels below or consistent with concentrations from



previous years (with the exception of groundwater from Well 10135, discussed in Section 3.2.1.1), further illustrating containment.

The presence of an overall inward hydraulic gradient towards the barrier drain and a review of groundwater quality for the groundwater monitoring wells demonstrate overall Site containment.

4. Activities

Summaries of normal activities and repairs performed in 2020 are presented below.

4.1 Process Activities

Process activities that occurred during the year included the following:

- Removal and disposal of hazardous waste
- Cleaning of all pump chambers
- Cleaning of all storage tanks
- · Cleaning of sludge from clarifier

4.2 Non-Process Activities

Non-process activities that occurred during the year included the following:

- Preventative maintenance
- Repair and maintenance of pump chambers and flow meters
- Repair and maintenance of fences and gates
- Landscape maintenance including grass cutting and tree and flower bed maintenance
- Upgrading computer system software
- · Heating and cooling system maintenance
- Repair of lights
- Regrading of two small, low areas in the surface of the landfill cap (Refer to Section 4.5)

4.3 Community Outreach

Community Outreach programs have included such activities as beautification of the area surrounding the Site and tours of the facility.

4.3.1 Beautification

The following beautification activities were conducted at Love Canal in 2020:

Maintenance and landscaping of the Site and surrounding areas



- Maintenance of flower beds and shrubs along Colvin Boulevard, 95th Street, and Frontier Avenue
- Cleanup of discarded debris along fence line

4.3.2 Tours

Tours of the facility have been given throughout the years to representatives of various environmental agencies (domestic and foreign) and educational groups. The tours include an informational orientation, accompanied with visual aids, followed by a guided tour of the treatment facility and landfill. No tours were given in 2020.

4.3.3 Communications

All required reports were prepared and submitted to various agencies throughout the year. Reports included the 2019 Annual Hazardous Waste Report to the NYSDEC, the 2019 Periodic Review Report (formerly titled the Annual Operations and Monitoring Report) to various agencies, quarterly SIU analytical reports to the NFWB and NYSDEC, and monthly SIU reports to the NFWB.

The Love Canal Annual Newsletter for 2019 was issued to surrounding citizens and agencies in March 2020. The report summarizes items such as the amount of groundwater treated on Site and then discharged to NFWB's sanitary sewer system, maintenance activities, and other non-operational activities for the year.

4.4 Waste Generation

Throughout 2020, both hazardous and nonhazardous waste was generated from various activities and disposed of off Site in accordance with applicable laws and regulations.

The tracking of hazardous waste is performed by regulated hazardous waste manifests. A summary of the Site's annual hazardous waste generation is reported to the NYSDEC in the Annual Hazardous Waste Report. The Annual Hazardous Waste Report summarizes the quantities, transporters, and disposal methods.

A total of 17,346 pounds of hazardous waste was generated from the activities listed below. The waste materials were sent off Site for disposal in accordance with applicable laws and regulations. Wastes generated in 2020 were disposed through incineration by Clean Harbors, LLC and Veolia ES Technical Solutions, LLC.

The hazardous waste disposed of in 2020 consisted of soil/debris and NAPL sludge as follows:

- Soil/Debris: 1,346 pounds (consisting of personal protective equipment [PPE], spent filter bags, and general debris)
- NAPL Sludge: 16,000 pounds (collected from LCTF process)

4.5 Routine Operations, Inspections, and Monitoring

A daily inspection of the system operations was performed for each day in 2020 in accordance with the O&M Manual for the Love Canal Site, dated March 2015. Inspection records are available upon request.



Monthly inspections, including fire extinguishers and carbon vapor phase vents for breakthrough were also completed in accordance with the O&M Manual. Inspection records are available upon request.

Semiannual inspection of the landfill cap surface (Appendix B) was conducted on May 21, 2020 and November 9, 2020. Two small, low areas in the surface of the landfill cap were observed southwest of PC-1 in 2019 (refer to 2019 PRR) and were noted again during the May 2020 inspection. As noted in a letter dated February 20, 2020 from GSH to the NYSDEC, GSH completed an engineering evaluation of the low areas and confirmed that GSH should fill the depressions with soil, as NYSDEC had suggested in its June 18, 2019 letter. These low areas were addressed from October 5 through October 13, 2020 by Niagara Grass Cutting. The two low areas were brought to grade with the surrounding cap area using approximately 200 yards of clean, imported topsoil. Depressions in the access road near PC-1 were regraded using approximately 45 tons of clean, imported 1-inch crusher run stone. A roller was used to compact the access road and cap area. The regraded cap area was then hand seeded and covered with straw. A photographic log of the cap and access road before, during, and after the regrading is included as Appendix I. The locations of the cap regrading are noted on Figure 2.1.

The NYSDEC conducted a Site inspection (landfill and treatment system) on July 7, 2020. No issues were identified other than the low areas in the cap, which were addressed in October 2020. The NYSDEC did not conduct a hazardous waste compliance inspection of the Site in 2020.

The NFWB performed an annual inspection of the LCTF and performed verification sampling of the effluent discharge on September 4, 2020. The inspection and the annual effluent verification sampling concluded that the Site is being maintained and operated in accordance with the Site's SIU discharge permit and other local, State, and Federal requirements. The completed NFWB 2020 Inspection Form is included as Appendix G.

The USEPA conducted its Five-Year Review inspection of the Site on November 6, 2018. Representatives of the USEPA, NYSDEC, Niagara County Department of Health, GSH, and GHD were in attendance. No issues were identified. The USEPA's Five-Year Review Report was published on April 11, 2019. In the report, the USEPA indicated that no issues were identified as part of the Five-Year Review, however several minor suggestions were incorporated in the 2019 PRR. The fifth Five-Year Review is scheduled for 2023.

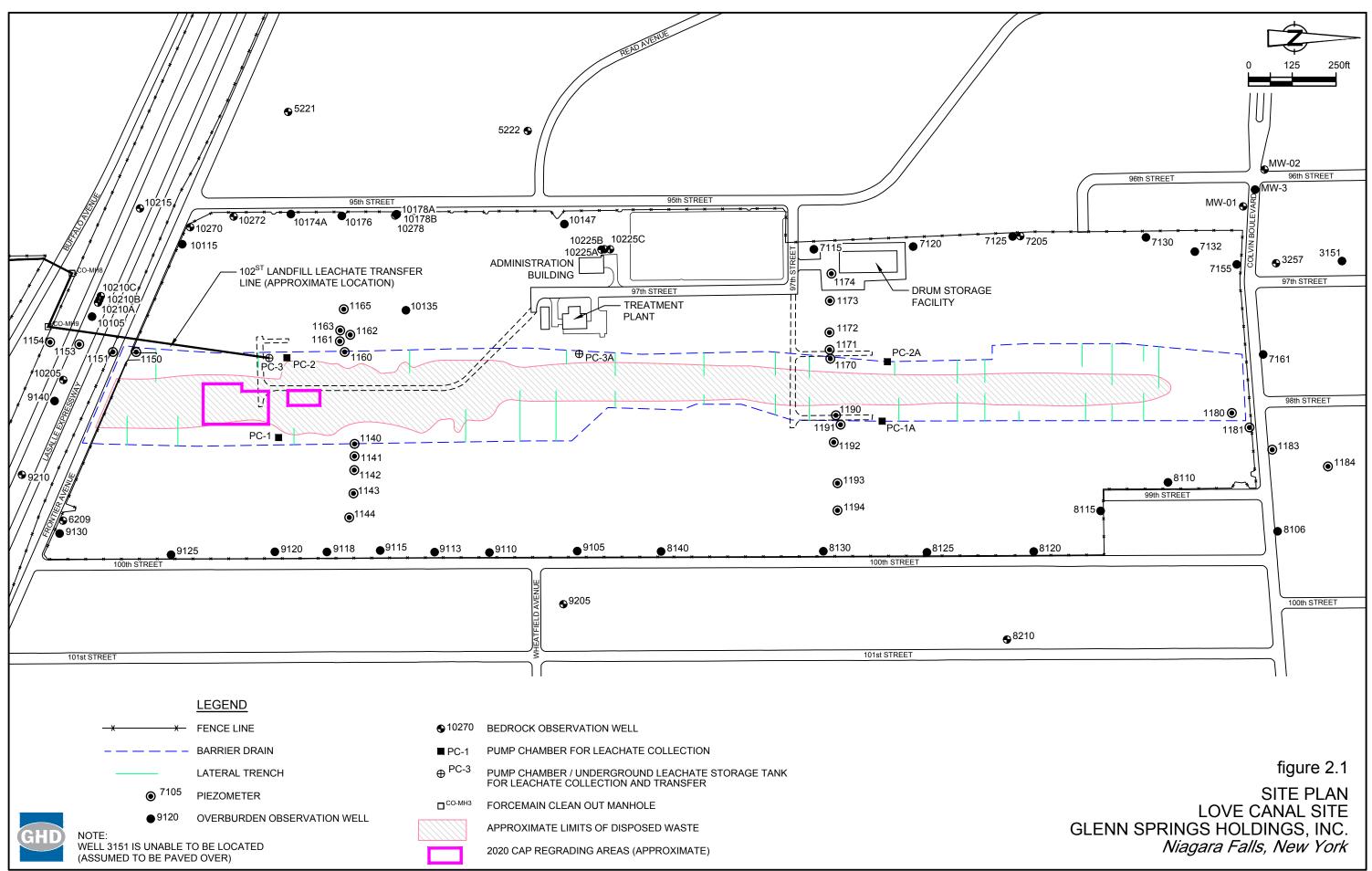
The backflow preventer system on the potable water supply lines was inspected and tested by CamTech Plumbing and Mechanical Inc. (CamTech) on February 25, 2020. CamTech is licensed and certified by the NFWB to perform the backflow preventer system inspections. All five backflow prevention devices were found to be operational with no maintenance required. A copy of the 2020 Test and Maintenance of Backflow Prevention Device Report for each device is presented in Appendix H.

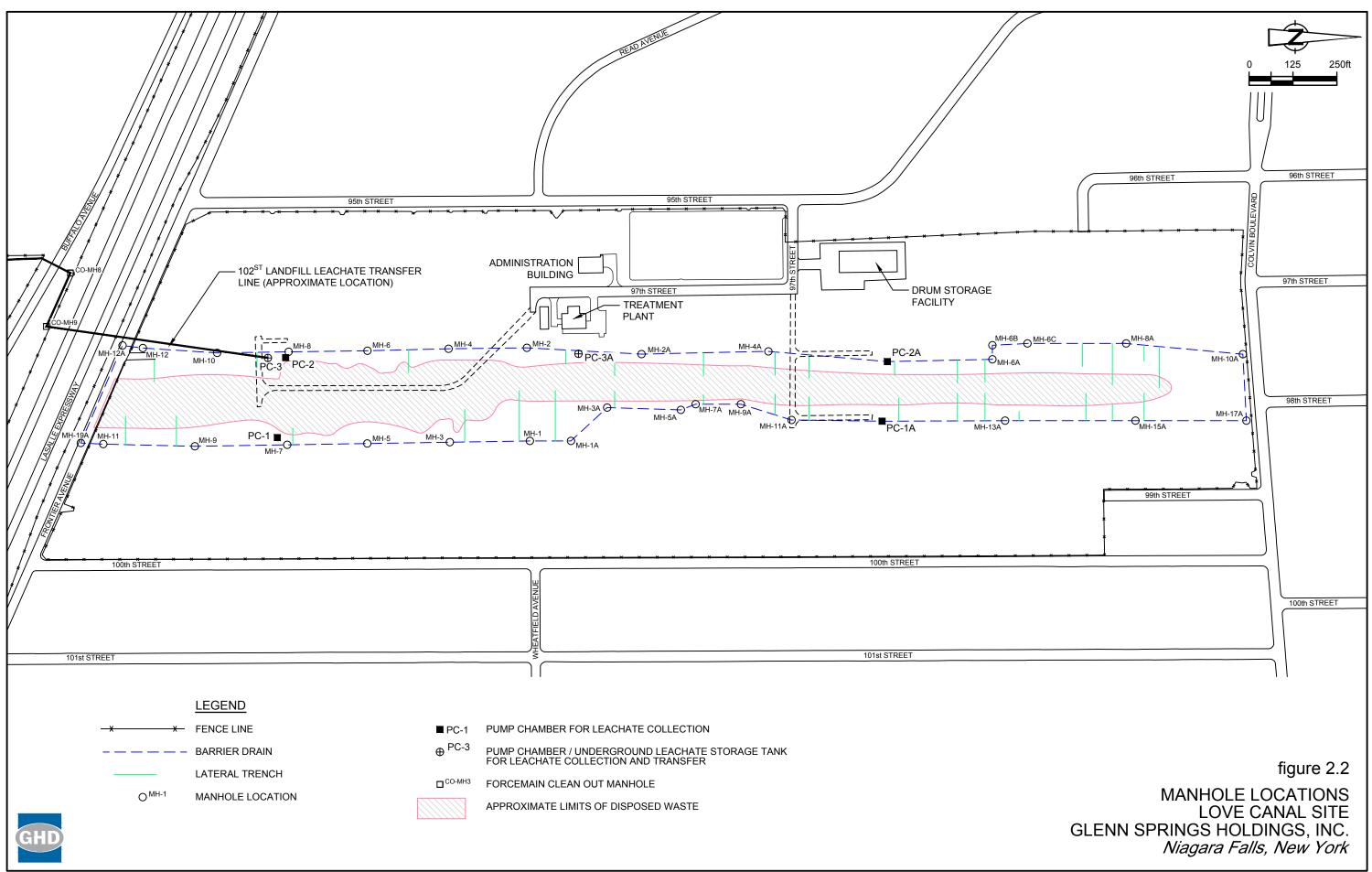
The annual fire system inspection was conducted on June 9, 2020. No issues were identified.

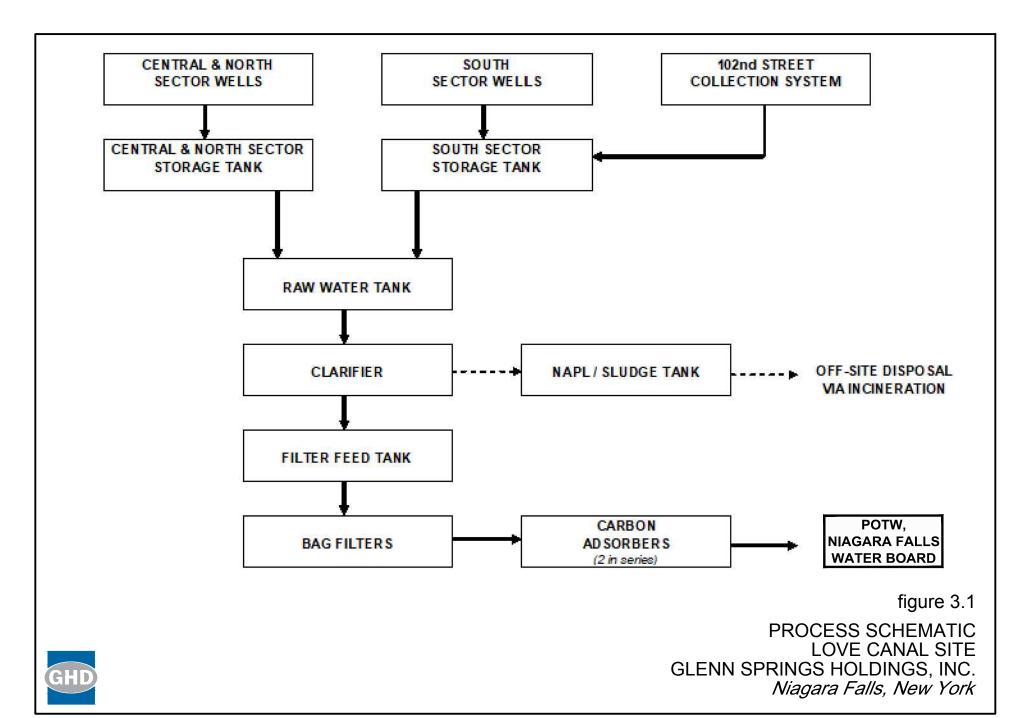


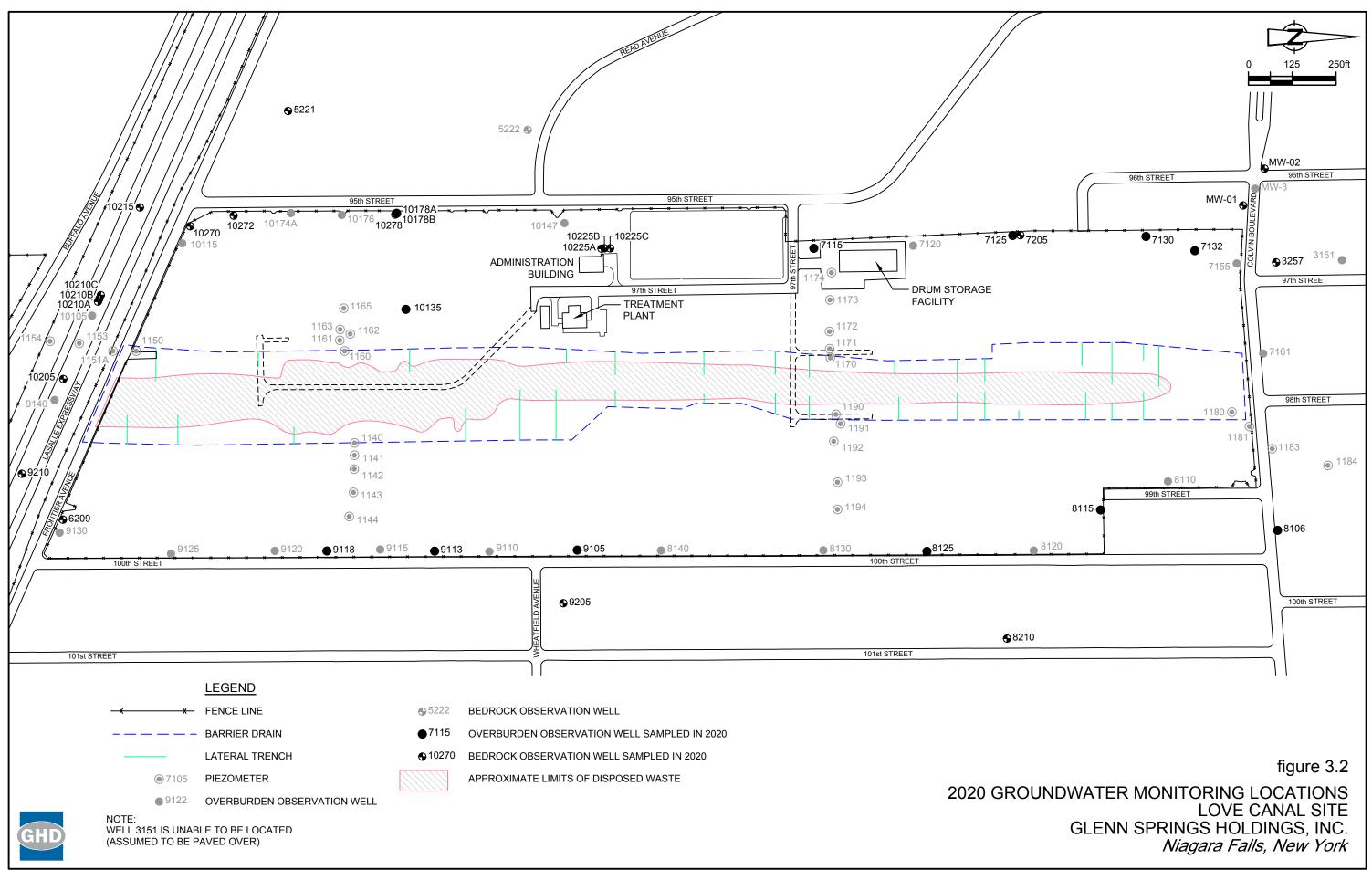
5. Conclusion

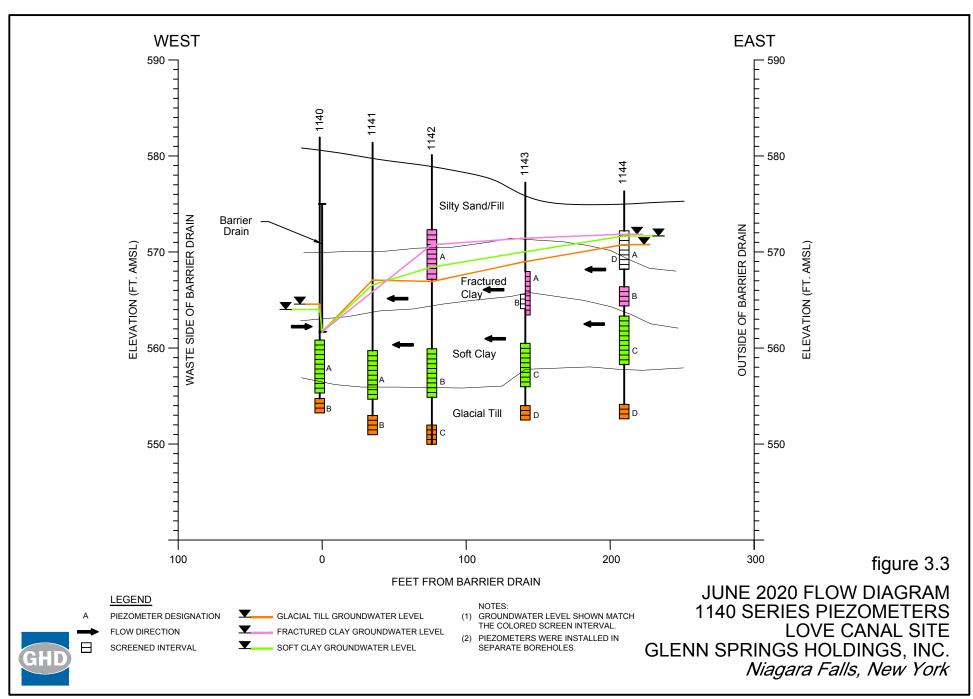
The 2020 monitoring results show that there has been no significant change in chemical concentration conditions and that the barrier drain system is successfully capturing leachate from the Site and preventing off-Site migration of contamination. The barrier drain continues to create an inward hydraulic gradient and capture leachate from the Site, preventing off-Site migration of chemicals, as evidenced by the groundwater gradients depicted on Figures 3.3 to 3.9 and analytical data from observation wells around the perimeter of the Site. The collection system is functioning as designed based on groundwater monitoring results and third-party inspections by the NYSDEC. The treatment system is functioning as designed based on inspections and sampling by the NFWB and sampling by GSH. Effluent quality is compliant with the Site's SIU discharge permit. There were 3,356,298 gallons of leachate collected, treated, and discharged from the Site, of which 3,236,167 gallons of leachate were collected from the Site, and 120,131 gallons were collected from the 102nd Street Site and pumped to the LCTF for treatment. Monitoring results continue to confirm that the remediation and containment system (i.e., the leachate collection and treatment system) is functioning as designed.

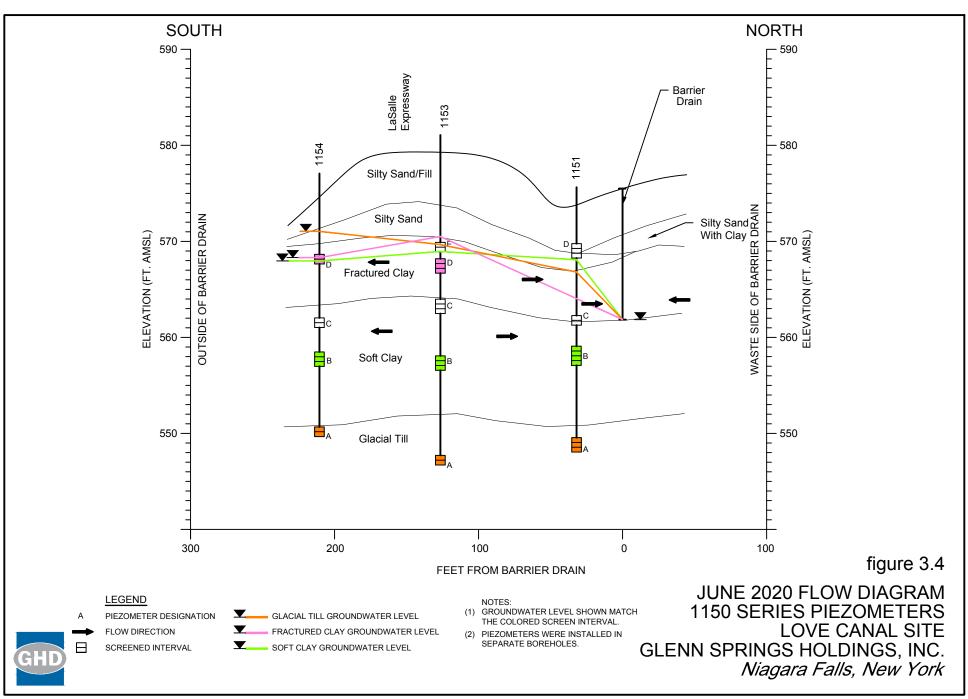


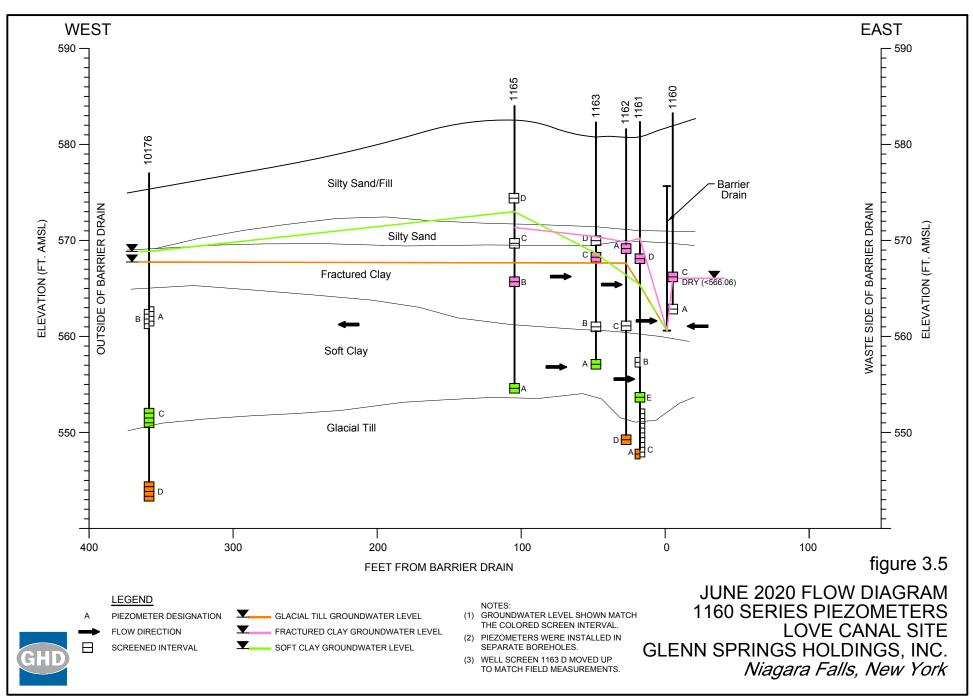


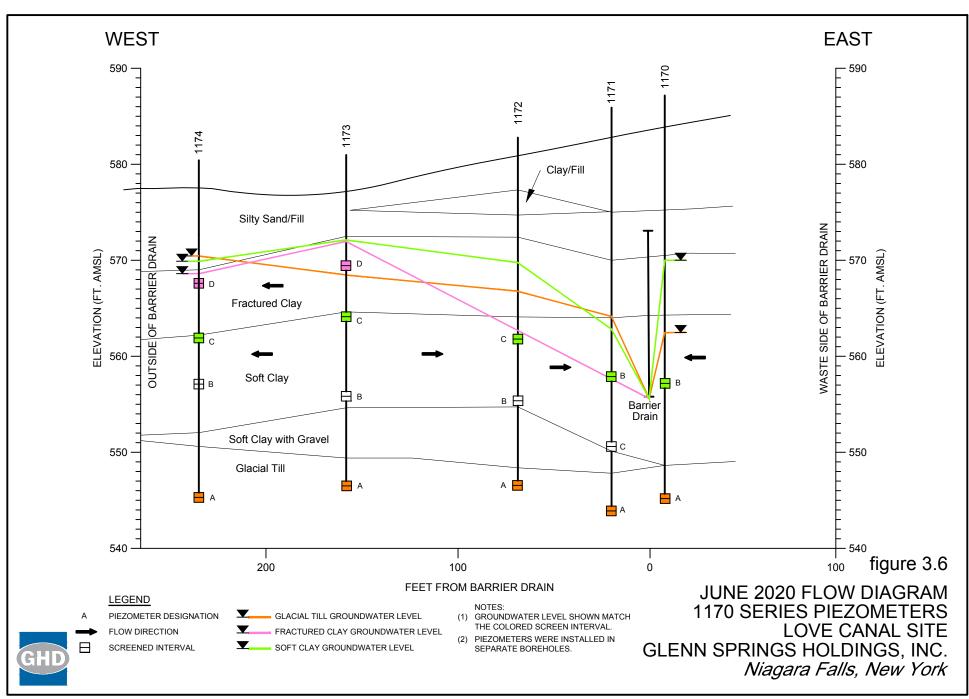


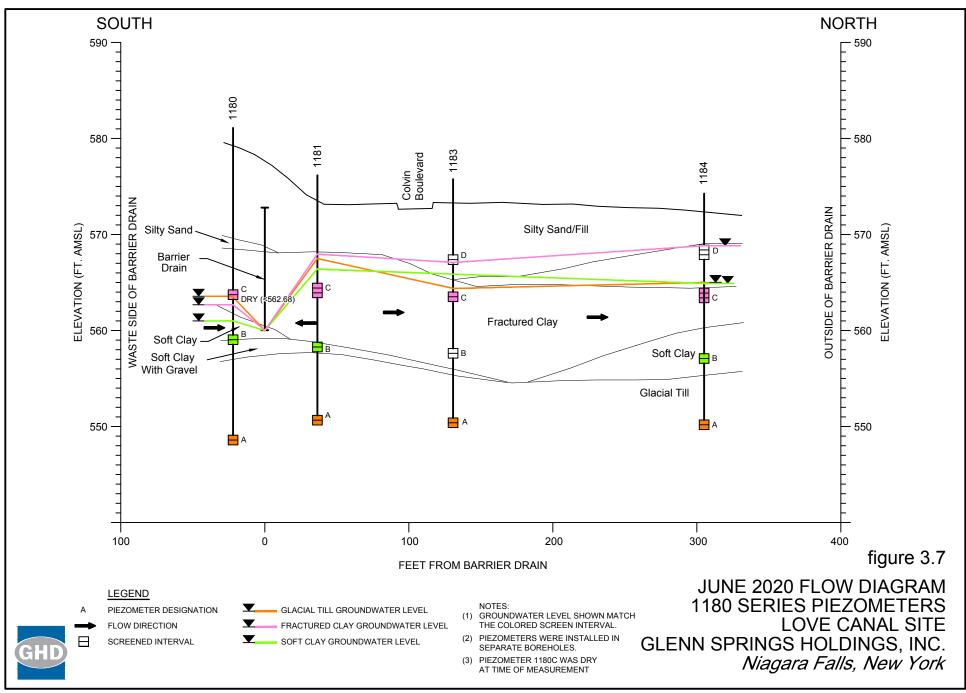


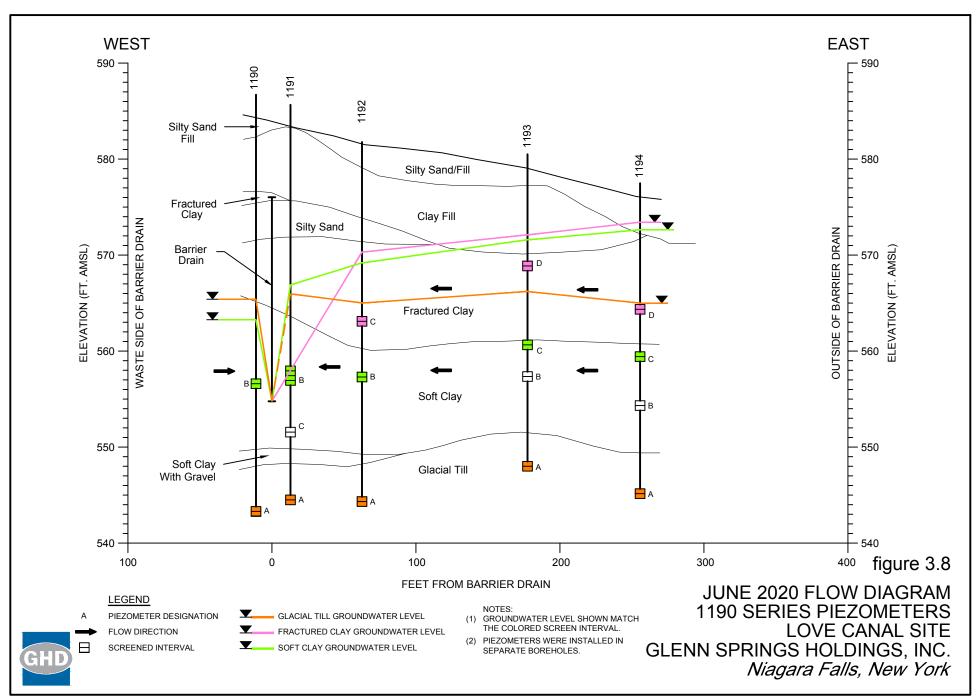


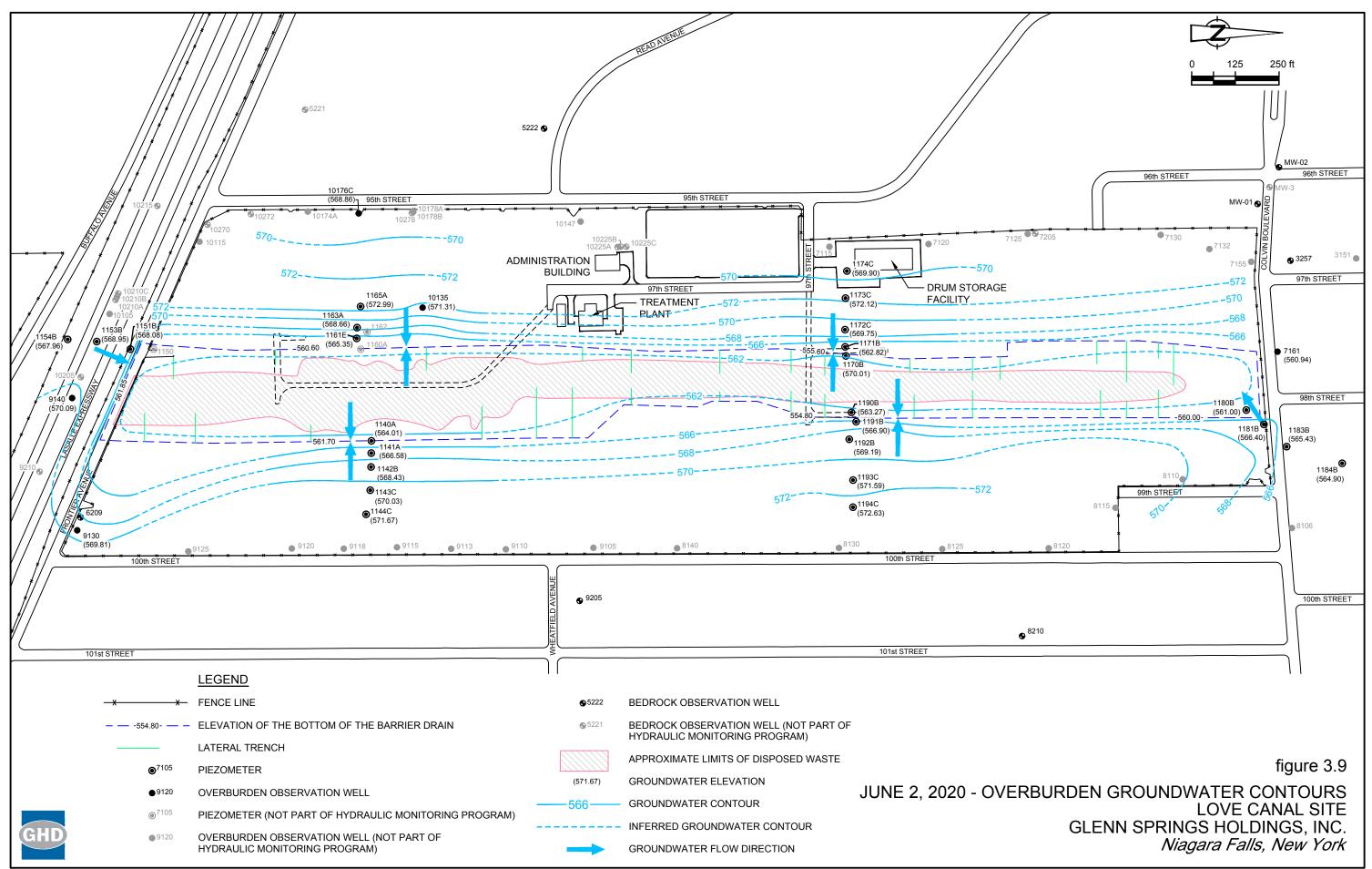


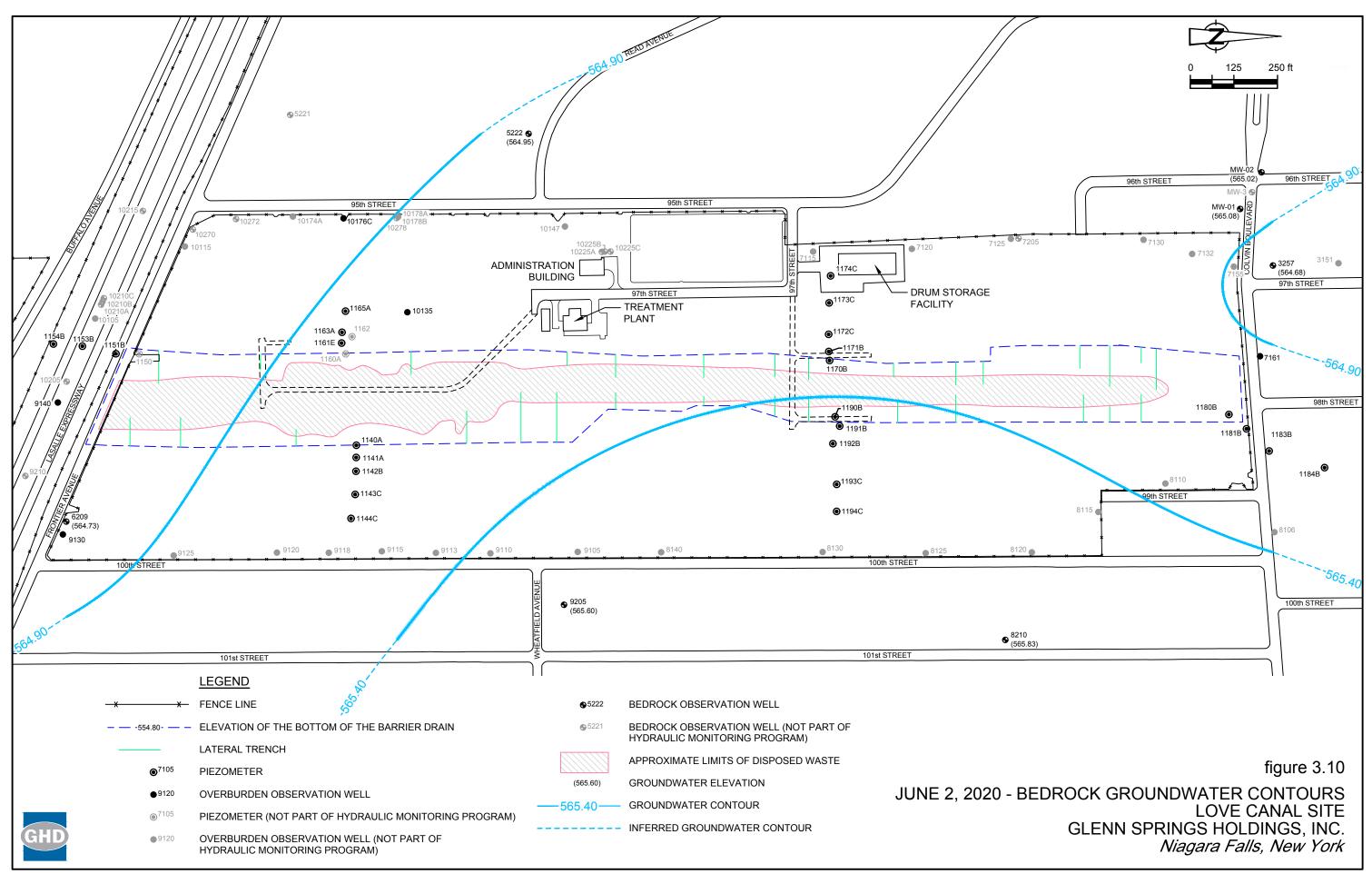








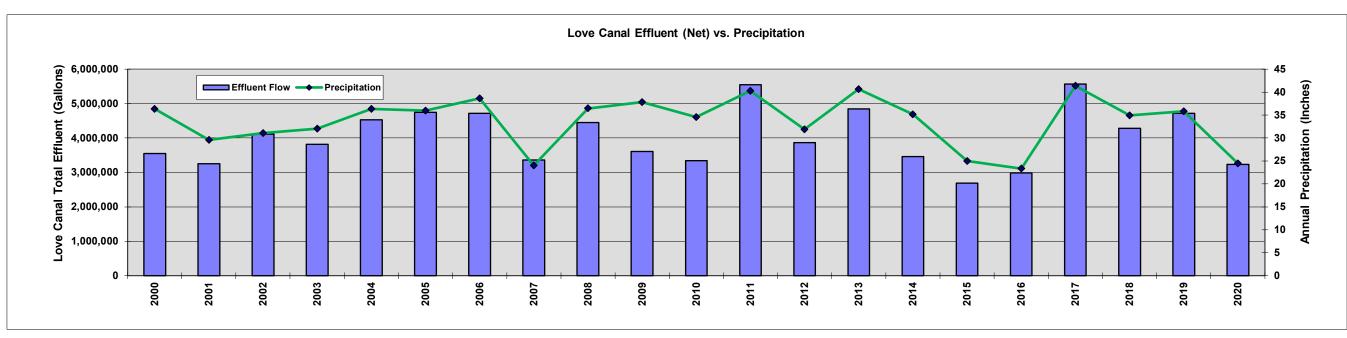




Monthly Volumes of Groundwater Treated Love Canal Long-Term Monitoring Program Niagara Falls, New York

		2000	2001	2002	2003	2004	2005	2006	2007	2008	2009	2010	2011	2012	2013	2014	2015	2016	2017	2018	2019	2020
January	Gross (1)	495,800	396,900	488,900	419,400	309,200	841,400	855,900	993,400	674,000	523,500	534,400	346,900	571,900	600,400	519,614	363,043	385,636	563,854	499,356	566,916	775,829
	Net ⁽²⁾	280,364	282,480	422,682	374,123	260,171	796,518	817,305	970,918	649,777	495,713	471,805	322,994	546,816	575,767	499,889	346,565	370,676	548,797	489,418	558,881	765,891
	Days ⁽³⁾	21	20	21	14	10	17	16	20	18	16	17	18	15	18	20	14	18	19	20	26	29
February	Gross	480,400	560,000	663,700	266,300	330,000	440,200	437,300	216,600	570,000	506,700	314,300	375,800	656,700	495,900	291,292	68,244	634,159	371,608	692,324	486,978	499,075
	Net Days	368,492 21	468,863 19	608,116 20	231,049 13	291,082	401,137	405,124 o	174,776 7	539,772 16	485,869 13	276,643 10	349,712 19	634,167 16	478,434 19	277,226 15	55,548 16	619,942	357,557 13	684,088 24	480,242 26	490,839 28
March	Gross	505,500	616,400	364,900	721,500	1,038,400	698,900	436,800	582,500	570,500	606,900	550,100	1,003,700	384,500	488,000	388,937	658,775	544,972	641,911	492,694	356,288	544,126
	Net	290,501	493,476	316,696	667,337	986,332	667,105	402,047	560,237	550,518	582,109	526,021	978,000	363,378	467,083	375,154	642,149	529,757	629,687	483,107	345,457	534,539
	Days	23	21	21	17	21	13	13	16	12	18	17	21	16	20	17	20	21	20	24	16	27
April	Gross	675,600 547,926	352,300 262,946	689,700 629,683	432,800 380,745	800,400 767,982	805,300 769,514	184,800 155,028	447,200 420,133	602,000 574,359	414,900 377,080	498,200 466,778	676,400 652,656	334,400 316,188	533,800 516,478	786,808 768,257	575,949 561,287	531,147 517,498	1,053,394 1,041,670	751,811 742,237	632,012 621,937	359,309 349,735
	Net Davs	20	202,940	20	16	17	14	6	420,133	12	16	15	11	18	22	20	17	19	1,041,070	28	25	28
May	Gross	473,300	311,200	589,500	425,400	326,500	183,400	121,800	323,200	172,900	306,200	379,400	942,700	363,100	148,500	444,598	113,599	175,158	983,450	169,570	579,868	214,573
	Net	335,331	207,580	532,251	379,299	294,612	156,846	93,394	297,471	147,715	267,700	348,837	917,206	341,424	129,687	428,177	99,179	163,324	971,685	160,642	570,737	205,645
	Days	20	17	20	14	10	5	4	12	11	14	18	17	16	18	21	12	15	25	18	19	23
June	Gross Net	632,200 486,721	202,200 132,132	395,100 347,485	367,900 303,576	253,200 208,659	160,800 118,979	130,700 104,449	173,300 148,638	128,700 107,411	110,000 79,200	205,200 174,305	473,100 449,046	142,000 118,568	497,300 478,285	168,921 152,639	262,025 245,083	98,255 83,122	178,582 164,597	110,862 101,522	248,940 239,997	147,455 138,115
	Days	20	16	14	13	9	6	5	4	6	7 9,200	13	16	12	18	132,039	19	11	15	101,322	12	130,113
July	Gross	333,900	182,200	194,500	187,700	137,700	92,600	195,500	129,100	164,760	187,900	85,600	79,700	98,400	280,000	151,772	138,495	77,140	335,930	103,168	134,549	91,978
	Net	184,955	111,941	145,344	142,849	111,217	78,234	183,084	99,026	141,442	153,170	55,670	53,632	72,435	260,823	123,921	122,874	62,847	322,782	95,872	125,685	84,682
	Days	20	16	16	11	7	3	5	6	6	7	4	5	9	19	15	16	12	18	8	10	8
August	Gross Net	437,100 286,925	267,200 194,821	151,300 107,928	158,600 114,497	301,900 269,934	98,800 55,055	322,440 293,900	120,800 106,040	197,340 191,068	369,400 347,425	184,300 162,562	193,900 166,652	73,960 49,422	193,144 168,418	98,166 83,010	108,376 91,308	65,714 50,772	242,754 228,321	91,721 79,817	105,894 99,061	77,128 64,652
	Days	23	18	17	8	10	5	10	5	6	18	8	13	8	21	9	7	6	17	79,017	8	8
September	Gross	209,600	144,900	148,600	105,800	484,800	317,900	249,160	68,400	152,200	101,500	88,100	47,800	161,100	131,289	139,016	151,905	96,279	114,926	95,188	217,213	101,043
	Net	82,263	81,619	94,401	60,350	435,482	284,315	213,343	49,041	122,101	76,057	56,678	21,679	136,728	110,397	111,392	134,935	79,011	100,242	83,595	204,446	89,450
0-4-1	Days	20	16	12	7	12	8	7	4	9	7	2	6	17	23	13	12	11	8	8	11	10
October	Gross Net	264,300 134,248	438,500 348,153	154,600 108,226	211,000 211,000	135,700 94,476	486,300 445,560	919,200 892,734	173,000 141,650	296,100 274,068	199,200 129,035	120,200 88,537	417,500 389,696	318,400 291,391	503,036 480,233	121,075 94,680	146,842 123,794	124,508 104,726	286,862 270,291	132,231 119,798	333,060 319,915	92,146 79,713
	Days	20	18	13	9	4	10	18	8	13	8	5	14	19	20	18	11	11	12	11	19	15
November	Gross	250,900	250,400	360,800	356,800	211,400	524,600	691,800	90,100	449,700	210,100	263,400	350,100	526,900	538,978	107,729	146,722	101,686	686,365	635,923	450,485	100,055
	Net	132,728	194,481	306,258	310,650	186,999	494,443	658,765	77,506	414,149	152,302	233,159	322,735	504,290	520,087	85,439	127,799	85,421	669,770	625,250	433,375	89,382
Dagamban	Days	17	16	14	12	5	14	14	3	14	12	15	12	20	19	12	15	11	22	21	25	18
December	Gross Net	522,600 421,149	555,300 475,856	549,600 496,556	692,300 643,735	674,400 622,403	502,000 476,165	510,400 492,900	345,700 317,790	757,500 733,582	506,200 467,578	510,900 483,221	952,000 926,201	517,700 493,061	677,411 660,890	471,085 456,099	155,368 138,929	335,448 318,036	278,817 264,686	626,070 616,013	728,072 716,352	353,581 343,524
	Days	17	18	15	14	14	12	12	8	20	17	17	19	14	17	14	13	14	19	28	28	22
	-		-						-												-	
Total	Gross	5,281,200	4,277,500	4,751,200	4,345,500	5,003,600	5,152,200	5,055,800	3,663,300	4,735,700	4,042,500	3,734,100	5,859,600	4,149,060	5,087,758	3,689,013	2,889,343	3,170,102	5,738,453	4,400,918	4,840,275	3,356,298
	Net	3,551,603	3,254,348	4,115,626	3,819,210	4,529,349	4,743,871	4,712,073	3,363,226	4,445,962	3,613,238	3,344,216	5,550,209	3,867,868	4,846,582	3,455,883	2,689,450	2,985,132	5,570,085	4,281,359	4,716,085	3,236,167
Monthly	Days Gross	242 440,100	215 356,458	203 395,933	148 362,125	128 416,967	118 429,350	119 421,317	107 305,275	143 394,642	153 336,875	141 311,175	171 488,300	180 345,755	234 423,980	186 307,418	172 240,779	168 264,175	212 478,204	207 366,743	225 403,356	230 279,692
Average	Net	295,967	271,196	342,969	318,268	377,446	395,323	392,673	280,269	370,497	301,103	278,685	462,517	322,322	403,882	287,990	224,121	248,761	464,174	356,780	393,007	269,681
	Days	20	18	17	12	11	10	10	9	12	13	12	14	15	20	16	14	14	18	17	19	19
Precipitation In	ches (4)	36.33	29.56	31.05	32.03	36.33	35.99	38.66	24.02	36.45	37.85	34.54	40.26	31.87	40.66	35.12	24.99	23.33	41.36	34.93	35.81	24.46

- Gross: Total volume of leachate treated in gallons; treatment at LCTF includes leachate collected from 102nd Street Landfill Site.
 Net: Love Canal leachate treated in gallons; net is equal to the total (gross) leachate treated less leachate received from 102nd Street.
 Days: Number of days treatment facility discharged to the sanitary sewer.
 Precipitation data obtained from the National Climatic Data Center for Niagara Falls International Airport.



Page 1 of 2

2020 Analytical Results Summary - Overburden Love Canal Long-Term Monitoring Program Niagara Falls, New York

Sa	mple Location: Sample ID: Sample Date:	7115 WG-9954-062420-SG-001 6/24/2020	7125 WG-9954-062420-SG-002 6/24/2020	7130 WG-9954-062420-SG-003 6/24/2020	7132 WG-9954-062420-SG-004 6/24/2020	8106 WG-9954-062620-RM-017 6/26/2020	8115 WG-9954-062420-SG-005 6/24/2020	8125 WG-9954-062420-SG-006 6/24/2020	9105 WG-9954-062420-SG-007 6/24/2020	9113 WG-9954-062520-SG-008 6/25/2020	9118 WG-9954-062520-SG-009 6/25/2020	9118 WG-9954-062520-SG-010 6/25/2020	10135 WG-9954-070120-SG-027 7/1/2020	10135 WG-9954-070120-SG-028 7/1/2020	10178A WG-9954-070720-SG-031 7/7/2020	10178B WG-9954-071020-SG-036 7/10/2020
Parameters	Uni	ts										(Duplicate)		(Duplicate)		
Volatile Organic Compounds																
1,1,1-Trichloroethane 1,1,2,2-Tetrachloroethane	hâ [,] hâ [,]		0.20 U 0.20 U	0.20 U 6.1	0.20 U 7.9	0.20 U 0.20 U	0.20 U 0.20 U									
1,1,2-Trichloroethane	μg		0.20 U	12	14	0.20 U	0.20 U									
1,1-Dichloroethane 1,1-Dichloroethene	μg		0.20 U 0.20 U	0.20 U 0.63 J	0.27 J 0.78 J	0.20 U 0.20 U	0.20 U 0.20 U									
1,2-Dichloroethane	μg. μg.		0.20 U	0.76 J 0.20 U	0.20 U	0.20 U										
1,2-Dichloropropane	μg		0.20 U	0.20 U 12	0.20 U	0.20 U	0.20 U									
2-Butanone (Methyl ethyl ketone) (MEK) 2-Hexanone	μg.		0.78 U 0.20 U	0.20 U	6.8 J 2.1 J	0.78 U 0.20 U	0.78 U 0.20 U									
4-Methyl-2-pentanone (Methyl isobutyl ketone) (M	IBK) μg.	L 0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.55 J	0.20 U	0.20 U				
Acetone Benzene	μg. μg.		5.0 U 0.20 U	79 6500	71 6300	5.0 U 0.20 U	5.0 U 0.20 U									
Bromodichloromethane	μg	L 0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U
Bromoform Bromomethane (Methyl bromide)	ha ha		0.25 U 0.70 U	0.25 U 0.70 U												
Carbon disulfide	μg	L 0.42 U	0.42 U	0.42 U	0.83 J	0.42 U	0.48 J	0.75 J	0.42 U	0.42 U	0.42 U	0.42 U				
Carbon tetrachloride Chlorobenzene	ha ha		0.34 U 0.20 U	0.34 U 2500	0.34 U 2400	0.34 U 0.20 U	0.34 U 0.20 U									
Chloroethane	μg	L 0.23 U	0.23 U	0.23 U	0.23 U	0.23 U	0.23 U	0.23 U	0.23 U	0.23 U	0.23 U	0.23 U	0.23 U	0.23 U	0.23 U	0.23 U
Chloroform (Trichloromethane) Chloromethane (Methyl chloride)	μg. μg.		0.24 U 0.28 U	0.24 U 0.28 U	0.24 U 0.28 U	0.24 U 0.38 J	0.24 U 0.28 U	0.24 U 0.28 U	0.24 U 0.28 U	0.24 U 0.32 J	0.24 U 0.38 J	0.24 U 0.28 U	140 0.28 U	140 0.28 U	0.24 U 0.28 U	0.24 U 0.28 U
cis-1,2-Dichloroethene	μg	L 0.23 U	0.23 U	0.23 U	0.23 U	0.23 U	0.23 U	0.23 U	0.23 U	0.23 U	0.23 U	0.23 U	26	28	0.23 U	0.23 U
cis-1,3-Dichloropropene Dibromochloromethane	μg. μg.		0.20 U 0.20 U	0.20 U 0.20 U												
Ethylbenzene	μg	L 0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	12	12	0.20 U	0.20 U
Methylene chloride Styrene	μg. μg.		0.65 U 0.20 U	6.4 0.20 U	7.2 0.20 U	0.65 U 0.20 U	0.65 U 0.20 U									
Tetrachloroethene	μg	L 0.21 U	0.21 U	0.21 U	0.21 U	0.21 U	0.21 U	0.21 U	0.21 U	0.21 U	0.21 U	0.21 U	20	23	0.21 U	0.21 U
Toluene trans-1,2-Dichloroethene	μg. μg.		0.20 U 0.20 U	22000 29	21000 35	0.20 U 0.20 U	0.20 U 0.20 U									
trans-1,3-Dichloropropene	μg.		0.23 U	0.23 U	0.23 U	0.23 U	0.20 U	0.23 U	0.23 U	0.23 U	0.23 U	0.20 U	0.23 U	0.23 U	0.23 U	0.23 U
Trichloroethene	μg		0.20 U 1.1 U	0.20 U 1.1 U	0.20 U 1.1 U	0.20 U 1.1 UJ	0.20 U	0.20 U	120	130 1.1 U	0.20 U 1.1 U	0.20 U				
Vinyl acetate Vinyl chloride	hâ [,] hâ [,]		0.20 U	1.1 U 0.20 U	1.1 U 0.20 U	1.1 U 10	11	0.20 U	1.1 U 0.20 U							
Xylenes (total)	μg		0.23 U	58	58	0.23 U	0.23 U									
Discrete Compo	unds Detected:	0	0	0	1	1	0	0	0	2	2	1	17	19	0	0
Semi-volatile Organic Compounds			4.4.111	4.4.11	4.4.11	4.4.11	4411	4.411	4.4.11	4.4.11	4.4.11	4411	02	100	4.4.11	4.4.11
1,2,4-Trichlorobenzene 1,2-Dichlorobenzene	ha ha		1.1 UJ 1.1 UJ	1.1 U 1.1 U	93 25	100 34	1.1 U 1.1 U	1.1 U 1.1 U								
1,3-Dichlorobenzene	μg	L 0.92 U	0.92 UJ	0.92 U	3.6 J	4.1 J	0.92 U	0.92 U								
1,4-Dichlorobenzene 2,2'-Oxybis(1-chloropropane) (bis(2-Chloroisoprop	μg. pyl) ether) μg.		1.1 UJ 1.3 UJ	1.1 U 1.3 U	73 1.3 U	97 1.3 U	1.1 U 1.3 U	1.1 U 1.3 U								
2,4,5-Trichlorophenol	μg	L 0.99 U	0.99 U	0.99 U	0.99 U	0.99 U	0.99 U	0.99 U	0.99 U	0.99 U	0.99 U	0.99 U	37	26	0.99 U	0.99 U
2,4,6-Trichlorophenol 2,4-Dichlorophenol	ha ha		1.3 U 1.2 U	1.3 U 330	1.3 U 350	1.3 U 1.2 U	1.3 U 1.2 U									
2,4-Dimethylphenol	μg	/L 1.3 U	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U	1.3 UJ	8.9 J	6.9 J	1.3 U	1.3 U				
2,4-Dinitrophenol 2,4-Dinitrotoluene	hâ [,] hâ [,]		19 U 2.2 UJ	19 U 2.2 U	19 U 2.2 U											
2,6-Dinitrotoluene	μg	L 1.2 U	1.2 UJ	1.2 U	1.2 U											
2-Chloronaphthalene 2-Chlorophenol	hâ [,] hâ [,]		1.3 UJ 0.97 U	1.3 U 0.97 U	1.3 U 0.97 U	1.3 U 0.97 U	1.3 U 0.97 U	1.3 U 0.97 U	1.3 U 0.97 U	1.3 U 0.97 U	1.3 U 0.97 U	1.3 U 0.97 U	1.3 U 27	1.3 U 31	1.3 U 0.97 U	1.3 U 0.97 U
2-Methylnaphthalene	μg	/L 1.2 U	1.2 UJ	1.2 U	1.4 J	1.2 U	1.2 U	1.2 U								
2-Methylphenol 2-Nitroaniline	ha ha		0.91 U 1.3 UJ	0.91 U 1.3 U	32 1.3 U	32 1.3 U	0.91 U 1.3 U	0.91 U 1.3 U								
2-Nitrophenol	μg	/L 1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U
3&4-Methylphenol 3,3'-Dichlorobenzidine	μg. μg.		1.1 U 1.1 UJ	1.1 U 1.1 U	71 R	71 1.1 U	1.1 U 1.1 U	1.1 U 1.1 U								
3-Nitroaniline	μg	L 2.3 U	2.3 UJ	2.3 U	2.3 U											
4,6-Dinitro-2-methylphenol 4-Bromophenyl phenyl ether	hâ [,] hâ [,]		18 U 1.5 UJ	18 U 1.5 U	18 U 1.5 U											
4-Chloro-3-methylphenol	μg	/L 0.98 U	0.98 U	0.98 U	0.98 U	0.98 U	0.98 U	0.98 UJ	46	40	0.98 U	0.98 U				
4-Chloroaniline 4-Chlorophenyl phenyl ether	μg. μg.		0.91 UJ 1.4 UJ	0.91 U 1.4 U	0.91 U 1.4 U											
4-Nitroaniline	μg	L 2.5 U	2.5 UJ	2.5 U	2.5 U	2.5 U	2.5 U	2.5 UJ	2.5 U	2.5 U	2.5 U	2.5 U				
4-Nitrophenol Acenaphthene	μg.		5.8 U 1.3 UJ	5.8 U 1.3 U	5.8 U 1.3 U											
Acenaphthylene	μg	/L 1.3 U	1.3 UJ	1.3 U	1.3 U											
Anthracene Benzo(a)anthracene	hâ [,] hâ [,]		1.2 UJ 1.5 UJ	1.2 U 1.5 U	1.2 U 1.5 U											
Benzo(a)pyrene	μg	/L 1.1 U	1.1 UJ	1.1 U	1.1 U											
Benzo(a h i)nerviene	μg		1.1 UJ 0.91 UJ	1.1 U 0.91 U	1.1 U 0.91 U											
Benzo(g,h,i)perylene Benzo(k)fluoranthene	hâ hâ	/L 1.1 U	1.1 UJ	1.1 U	1.1 U											
Benzoic acid	μg	L 33 U	33 UJ	33 U	33 U	33 U	33 U	33 U 1.5 U	33 U	33 U	33 U	33 U	11000	10000 290	33 U 1.5 U	33 U
Benzyl alcohol bis(2-Chloroethoxy)methane	hâ [,] hâ [,]		1.5 U 1.8 UJ	1.5 U 1.8 U	290 J 1.8 U	1.8 U	1.5 U 1.8 U	1.5 U 1.8 U								
bis(2-Chloroethyl)ether	μg	L 1.2 U	1.2 UJ	1.2 U	1.2 U	1.2 UJ	1.2 U	20	24	1.2 U	1.2 U					
bis(2-Ethylhexyl)phthalate (DEHP) Butyl benzylphthalate (BBP)	hâ [,] hâ [,]		0.91 UJ 1.3 UJ	0.91 U 1.3 U	1.3 J 1.3 U											
Chrysene	μg	/L 1.1 U	1.1 UJ	1.1 U	1.1 U											
Dibenz(a,h)anthracene Dibenzofuran	μg. μg.		0.93 UJ 1.3 UJ	0.93 U 1.3 U	0.93 U 1.3 U											

2020 Analytical Results Summary - Overburden Love Canal Long-Term Monitoring Program Niagara Falls, New York

	Sample Location:	7115	7125	7130	7132	8106	8115	8125	9105	9113	9118	9118	10135	10135	10178A	10178B
	Sample ID:	WG-9954-062420-SG-001	WG-9954-062420-SG-002		WG-9954-062420-SG-004	WG-9954-062620-RM-017	WG-9954-062420-SG-005	WG-9954-062420-SG-006	WG-9954-062420-SG-007	WG-9954-062520-SG-008	WG-9954-062520-SG-009	WG-9954-062520-SG-010	WG-9954-070120-SG-027	WG-9954-070120-SG-028	WG-9954-070720-SG-031	WG-9954-071020-SG-036
	Sample Date:	6/24/2020	6/24/2020	6/24/2020	6/24/2020	6/26/2020	6/24/2020	6/24/2020	6/24/2020	6/25/2020	6/25/2020	6/25/2020 (Duplicate)	7/1/2020	7/1/2020 (Duplicate)	7/7/2020	7/10/2020
Parameters	Unit	s										, ,		, ,		
Semi-volatile Organic Compounds (C																
Diethyl phthalate	μg/l		0.98 UJ	0.98 U	0.98 U	0.98 U	0.98 U	0.98 U	0.98 U	0.98 U	0.98 U	0.98 U	0.98 U	0.98 U	0.98 U	0.98 U
Dimethyl phthalate Di-n-butylphthalate (DBP)	μg/l		1.2 UJ 1.9 UJ	1.2 U 1.9 U	1.2 U 1.9 U	1.2 U 1.9 U	1.2 U 1.9 U	1.2 U 1.9 U	1.2 U 1.9 U	1.2 U 1.9 U	1.2 U 1.9 U	1.2 U 1.9 U	1.2 U 1.9 U	1.2 U 1.9 U	1.2 U 1.9 U	1.2 U 1.9 U
Di-n-octyl phthalate (DnOP)	րց/և րց/և		3.0 UJ	3.0 U	3.0 U	3.0 U	3.0 U	3.0 U	3.0 U	3.0 U	3.0 U	3.0 U	3.0 U	3.0 U	3.0 U	3.0 U
Fluoranthene	μg/l		1.4 UJ	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U
Fluorene	μg/l		1.2 UJ	1.2 U	1.2 U	1.2 U	1.2 U	1.2 U	1.2 U	1.2 U	1.2 U	1.2 U	1.2 U	1.2 U	1.2 U	1.2 U
Hexachlorobenzene	μg/l	_ 1.4 U	1.4 UJ	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U
Hexachlorobutadiene	μg/l		0.91 UJ	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U	1.0 J	0.91 U	0.91 U
Hexachlorocyclopentadiene	μg/l		2.0 UJ	2.0 U	2.0 U	2.0 U	2.0 U	2.0 U	2.0 U	2.0 U	2.0 U	2.0 U	2.0 U	2.0 U	2.0 U	2.0 U
Hexachloroethane Indeno(1,2,3-cd)pyrene	μg/l		0.96 UJ	0.96 U 1.6 U	0.96 U	0.96 U	0.96 U	0.96 U	0.96 U	0.96 U	0.96 U 1.6 U	0.96 U	0.96 U	0.96 U	0.96 U	0.96 U
Inderio(1,2,3-cd)pyrene Isophorone	րց/Լ րց/Լ		1.6 UJ 1.3 UJ	1.6 U	1.6 U 1.3 U	1.6 U 1.3 U	1.6 U 1.3 U	1.6 U 1.3 UJ	1.6 U R	1.6 U 1.3 U	1.6 U 1.3 U	1.6 U 1.3 U				
Naphthalene	μg/l		1.1 UJ	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U
Nitrobenzene	μg/l		1.4 UJ	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U
N-Nitrosodi-n-propylamine	μg/L		1.1 UJ	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U
N-Nitrosodiphenylamine	μg/l		2.4 UJ	2.4 U	2.4 U	2.4 U	2.4 U	2.4 U	2.4 U	2.4 U	2.4 U	2.4 U	2.4 U	2.4 U	2.4 U	2.4 U
Pentachlorophenol	μg/l		8.9 U	8.9 U	8.9 U	8.9 U	8.9 U	8.9 U	8.9 U	8.9 U	8.9 U	8.9 U	8.9 U	8.9 U	8.9 U	8.9 U
Phenal	μg/l		1.3 UJ	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U
Phenol Pyrene	μg/l		0.91 U 1.3 UJ	0.91 U 1.3 U	0.91 U 1.3 U	0.91 U 1.3 U	0.91 U 1.3 U	0.91 U 1.3 U	0.91 U 1.3 U	0.91 U 1.3 U	0.91 U 1.3 U	0.91 U 1.3 U	40 1.3 U	44 1.3 U	0.91 U 1.3 U	0.91 U 1.3 U
rytetie	μg/l	1.50	1.5 05	1.5 0	1.5 0	1.5 0	1.5 0	1.5 0	1.5 0	1.5 0	1.5 0	1.5 0	1.5 0	1.5 0	1.5 0	1.3 0
Discret	ete Compounds Detected:	1	0	0	0	0	0	0	0	0	0	0	16	16	0	1
Polychlorinated Biphenyls																
Aroclor-1016 (PCB-1016)	μg/l		0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U
Aroclor-1221 (PCB-1221)	μg/l		0.91 U	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1232 (PCB-1232) Aroclor-1242 (PCB-1242)	րց/Լ րց/Լ		0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U
Aroclor-1248 (PCB-1248)	μg/l		0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U
Aroclor-1254 (PCB-1254)	μg/l		0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U
Aroclor-1260 (PCB-1260)	μg/l	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U
Discret	ete Compounds Detected:	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
Pesticides																
4,4'-DDD	μg/l		0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.19 U	0.19 U	0.019 U	0.019 U
4,4'-DDE 4,4'-DDT	μg/l		0.019 U	0.019 U	0.019 U 0.019 U	0.019 U 0.019 U	0.019 U 0.019 U	0.019 U 0.019 U	0.019 U	0.019 U	0.019 U	0.019 U 0.019 U	0.19 U	0.19 U 0.19 U	0.019 U 0.019 U	0.019 U 0.019 U
Aldrin	րց/Լ րց/Լ		0.019 U 0.019 U	0.019 U 0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U 0.019 U	0.019 U 0.019 U	0.019 U 0.019 U	0.019 U	0.19 U 1.0 J	0.19 0	0.019 U	0.019 U
alpha-BHC	μg/l		0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	27	25	0.019 U	0.069
alpha-Chlordane	μg/l		0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.19 U	0.19 U	0.019 U	0.019 U
beta-BHC	μg/l		0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	6.8 J	6.9	0.019 U	0.019 U
delta-BHC	μg/l		0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	9.2	8.7	0.019 U	0.021 J
Dieldrin	μg/l		0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.19 U	0.19 U	0.019 U	0.019 U
Endosulfan I Endosulfan II	μg/l		0.019 U 0.019 U	0.019 U 0.019 U	0.019 U 0.019 U	0.019 U 0.019 U	0.019 U 0.019 U	0.019 U 0.019 U	0.019 U 0.019 U	0.019 U 0.019 U	0.019 U 0.019 U	0.019 U 0.019 U	0.19 U 0.19 U	0.19 U 0.19 U	0.019 U 0.019 U	0.019 U 0.019 U
Endosulfan sulfate	րց/Լ		0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.19 U	0.19 U	0.019 U	0.019 U
Endrin	μg/l		0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.19 U	0.19 U	0.019 U	0.019 U
Endrin ketone	μg/l		0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.19 U	0.19 U	0.019 U	0.019 U
gamma-BHC (lindane)	μg/l	_ 0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	6.5	6.2	0.019 U	0.058
gamma-Chlordane	μg/l		0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.19 U	0.19 U	0.019 U	0.019 U
Heptachlor	μg/l		0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.19 U	0.19 U	0.019 U	0.019 U
Heptachlor epoxide	μg/l		0.019 U 0.019 U	0.019 U 0.019 U	0.019 U 0.019 U	0.019 U 0.019 U	0.019 U 0.019 U	0.019 U 0.019 U	0.019 U 0.019 U	0.019 U 0.019 U	0.019 U	0.019 U 0.019 U	0.19 U 0.19 U	0.19 U 0.19 U	0.019 U 0.019 U	0.019 U 0.041 J
Methoxychlor Toxaphene	րց/Լ րց/Լ		0.019 U 0.46 U	0.019 U 0.46 U	0.019 U 0.46 U	0.019 U 0.46 U	0.019 U 0.46 U	0.019 U 0.46 U	0.019 U 0.46 U	0.019 U 0.46 U	0.019 U 0.46 U	0.019 U 0.46 U	0.19 U 4.6 U	0.19 U 4.6 U	0.019 U 0.46 U	0.041 J 0.46 U
		0.100	0	0.100	0	0.100	0.40 0	0	0	0	0.100	0		5	0	4
DISCRET	ete Compounds Detected:	U	U	U	U	U	U	U	U	U	U	U	ວ	ວ	U	4

Notes:
J - Estimated concentration
U - Not detected at the associated reporting limit
UJ - Not detected; associated reporting limit is estimated
R - Rejected

2020 Analytical Results Summary - Bedrock Love Canal Long-Term Monitoring Program Niagara Falls, New York

	Sample Location: Sample ID: Sample Date:	3257 WG-9954-062620-RM-018 6/26/2020	5221 WG-9954-070120-SG-025 7/1/2020	6209 WG-9954-062520-SG-011 6/25/2020	7205 WG-9954-062620-RM-014 6/26/2020	8210 WG-9954-062620-RM-015 6/26/2020	9205 WG-9954-062620-RM-016 6/26/2020	9210 WG-9954-062520-SG-012 6/25/2020	10205 WG-9954-062520-SG-013 6/25/2020	10210A WG-9954-070920-SG-033 7/9/2020	10210B WG-9954-063020-SG-019 6/30/2020	10210C WG-9954-063020-SG-020 6/30/2020	10210C WG-9954-063020-SG-021 6/30/2020
Parameters	Units	s											(Duplicate)
Volatile Organic Compounds 1,1,1-Trichloroethane 1,1,2,2-Tetrachloroethane	μg/L μg/.	0.20 U	0.20 U 0.20 U	0.20 U 0.20 U	0.20 U 0.20 U	0.20 U 0.20 U	0.20 U 0.20 U	0.20 U 0.20 U	0.20 U 0.20 U	0.20 U 0.20 U	0.20 U 0.20 U	0.20 U 0.20 U	0.20 U 0.20 U
1,1,2-Trichloroethane 1,1-Dichloroethane 1,1-Dichloroethene 1,2-Dichloroethane 1,2-Dichloropropane	µg/L µg/L µg/L µg/L µg/L	. 0.20 U . 0.20 U . 0.20 U . 0.20 U . 0.20 U	0.20 U 0.20 U 0.20 U 0.20 U 0.20 U										
2-Butanone (Methyl ethyl ketone) (MEK) 2-Hexanone 4-Methyl-2-pentanone (Methyl isobutyl ketone)	μg/L μg/L	. 0.78 U . 0.20 U . 0.20 U	0.78 U 0.20 U 0.20 U 5.0 U	0.78 U 0.20 U 0.20 U 5.0 U	0.78 U 0.20 U 2.4 J 5.0 U	0.78 U 0.20 U 0.20 U 5.0 U	0.78 U 0.20 U 0.20 U 5.0 U	0.78 U 0.20 U 0.20 U 5.0 U	0.78 U 0.20 U 0.20 U 5.0 U	0.78 U 0.20 U 0.20 U 5.0 U	0.78 U 0.20 U 0.20 U 5.0 U	0.78 U 0.20 U 0.20 U 5.0 U	0.78 U 0.20 U 0.20 U 5.0 U
Benzene Bromodichloromethane Bromoform Bromomethane (Methyl bromide)	. д µg/L µg/L µg/L	. 0.20 U	0.20 U 0.20 U 0.25 U 0.70 U	0.20 U 0.20 U 0.25 U 0.70 U	0.20 U 0.20 U 0.25 U 0.70 U	0.20 U 0.20 U 0.25 U 0.70 U	0.20 U 0.20 U 0.25 U 0.70 U	0.20 U 0.20 U 0.25 U 0.70 U	0.20 U 0.20 U 0.25 U 0.70 U	0.20 U 0.20 U 0.25 U 0.70 U	0.20 U 0.20 U 0.25 U 0.70 U	0.20 U 0.20 U 0.25 U 0.70 U	0.20 U 0.20 U 0.25 U 0.70 U
Carbon disulfide Carbon tetrachloride Chlorobenzene Chloroethane	µg/L µg/L µg/L	1.2 J 0.34 U 0.20 U 0.23 U	0.42 U 0.34 U 0.20 U 0.23 U	2.0 J 0.34 U 0.20 U 0.23 U	4.0 J 0.34 U 0.20 U 0.23 U	2.7 J 0.34 U 0.20 U 0.23 U	1.9 J 0.34 U 0.20 U 0.23 U	2.6 J 0.34 U 0.20 U 0.23 U	5.4 J 0.34 U 0.20 U 0.23 U	9.2 J 0.34 U 0.20 U 0.23 U	4.3 J 0.34 U 0.20 U 0.23 U	4.2 J 0.34 U 0.20 U 0.23 U	3.1 J 0.34 U 0.20 U 0.23 U
Chloroform (Trichloromethane) Chloromethane (Methyl chloride) cis-1,2-Dichloroethene cis-1,3-Dichloropropene Dibromochloromethane	µg/L µg/L µg/L µg/L	. 0.24 U . 0.28 U . 0.23 U . 0.20 U . 0.20 U	0.24 U 0.28 U 0.23 U 0.20 U 0.20 U	0.24 U 0.28 U 0.23 U 0.20 U 0.20 U	0.24 U 0.30 J 0.23 U 0.20 U 0.20 U	0.24 U 0.35 J 0.23 U 0.20 U 0.20 U	0.24 U 0.28 U 0.23 U 0.20 U 0.20 U	0.24 U 0.30 J 0.23 U 0.20 U 0.20 U	0.24 U 0.28 U 0.23 U 0.20 U 0.20 U				
Ethylbenzene Methylene chloride Styrene Tetrachloroethene	µg/L µg/L µg/L µg/L µg/L		0.20 U 0.20 U 0.65 U 0.20 U 0.21 U	0.20 U 0.65 U 0.20 U 0.21 U	0.20 U 0.20 U 0.65 U 0.20 U 0.21 U	0.20 U 0.65 U 0.20 U 0.21 U	0.20 U 0.20 U 0.65 U 0.20 U 0.21 U	0.20 U 0.20 U 0.65 U 0.20 U 0.21 U	0.20 U 0.65 U 0.20 U 0.21 U	0.20 U 0.20 U 0.65 U 0.20 U 0.21 U	0.20 U 0.20 U 0.65 U 0.20 U 0.21 U	0.20 U 0.65 U 0.20 U 0.21 U	0.20 U 0.65 U 0.20 U 0.21 U
Toluene trans-1,2-Dichloroethene trans-1,3-Dichloropropene Trichloroethene	µg/L µg/L µg/L µg/L		0.20 U 0.20 U 0.23 U 0.20 U										
Vinyl acetate Vinyl chloride Xylenes (total)	μg/L μg/L μg/L	1.1 U 0.20 U 0.23 U	1.1 U 0.20 U 0.23 U	1.1 U 0.20 U 0.23 U	1.1 UJ 0.20 U 0.23 U	1.1 UJ 0.20 U 0.23 U	1.1 U 0.20 U 0.23 U	1.1 U 0.20 U 0.23 U	1.1 U 0.20 U 0.23 U	1.1 U 0.20 U 0.23 U	1.1 U 0.20 U 0.23 U	1.1 U 0.20 U 0.23 U	1.1 U 0.20 U 0.23 U
Semi-volatile Organic Compounds	ne compoundo Detectou.	·	Ü	·	G	-	·	-	·	·	·	·	·
1,2,4-Trichlorobenzene 1,2-Dichlorobenzene 1,3-Dichlorobenzene 1,4-Dichlorobenzene	μg/L μg/L μg/L	. 1.1 U . 0.92 U . 1.1 U	1.1 U 1.1 U 0.92 U 1.1 U	1.1 U 1.1 U 0.92 U 1.1 U	1.1 U 1.1 U 0.92 U 1.1 U	1.1 U 1.1 U 0.92 U 1.1 U	1.1 U 1.1 U 0.92 U 1.1 U	1.1 U 1.1 U 0.92 U 1.1 U	1.1 U 1.1 U 0.92 U 1.1 U	1.1 U 1.1 U 0.92 U 1.1 U	1.1 U 1.1 U 0.92 U 1.1 U	1.1 U 1.1 U 0.92 U 1.1 U	1.1 U 1.1 U 0.92 U 1.1 U
2,2'-Oxybis(1-chloropropane) (bis(2-Chlor 2,4,5-Trichlorophenol 2,4,6-Trichlorophenol 2,4-Dichlorophenol 2,4-Dimethylphenol	roisopropyl) ether) µg/L µg/L µg/L µg/L µg/L	. 0.99 U	1.3 U 0.99 U 1.3 U 1.2 U 1.3 U	1.3 U 0.99 U 1.3 U 1.2 U 1.3 UJ	1.3 U 0.99 U 1.3 U 1.2 U 1.3 U	1.3 U 0.99 U 1.3 U 1.2 U 1.3 U	1.3 U 0.99 U 1.3 U 1.2 U 1.3 U	1.3 U 0.99 U 1.3 U 1.2 U 1.3 UJ	1.3 U 0.99 U 1.3 U 1.2 U 1.3 UJ	1.3 U 0.99 U 1.3 U 1.2 U 1.3 U			
2,4-Dinitrophenol2,4-Dinitrotoluene2,6-Dinitrotoluene2-Chloronaphthalene	µg/L µg/L µg/L		19 U 2.2 U 1.2 U 1.3 U										
2-Chlorophenol 2-Methylnaphthalene 2-Methylphenol 2-Nitroaniline 2-Nitrophenol	µg/L µg/L µg/L	. 0.91 U . 1.3 U	0.97 U 1.2 U 0.91 U 1.3 U 1.4 U	0.97 U 1.2 U 0.91 U 1.3 U 1.4 U	0.97 U 1.2 U 0.91 U 1.3 U 1.4 U	0.97 U 1.2 U 0.91 U 1.3 U 1.4 U	0.97 U 1.2 U 0.91 U 1.3 U 1.4 U	0.97 U 1.2 U 0.91 U 1.3 U 1.4 U	0.97 U 1.2 U 0.91 U 1.3 U 1.4 U	0.97 U 1.2 U 0.91 U 1.3 U 1.4 U	0.97 U 1.2 U 0.91 U 1.3 U 1.4 U	0.97 U 1.2 U 0.91 U 1.3 U 1.4 U	0.97 U 1.2 U 0.91 U 1.3 U 1.4 U
3&4-Methylphenol 3,3'-Dichlorobenzidine 3-Nitroaniline 4,6-Dinitro-2-methylphenol	µg/L µg/L µg/L µg/L µg/L	1.1 U 1.1 U 1.1 U 2.3 U 18 U	1.1 U 1.1 U 2.3 U 18 U	1.1 U 1.1 U 2.3 U 18 U	1.1 U 1.1 U 2.3 U 18 U	1.1 U 1.1 U 2.3 U 18 U	1.1 U 1.1 U 2.3 U 18 U	1.1 U 1.1 U 2.3 U 18 U	1.1 U 1.1 U 2.3 U 18 U	1.1 U 1.1 U 2.3 U 18 U	1.1 U 1.1 U 2.3 U 18 U	1.1 U 1.1 U 2.3 U 18 U	1.1 U 1.1 U 1.1 U 2.3 U 18 U
4-Bromophenyl phenyl ether 4-Chloro-3-methylphenol 4-Chloroaniline 4-Chlorophenyl phenyl ether	. дд/L µg/L µg/L µg/L	. 0.98 U . 0.91 U . 1.4 U	1.5 U 0.98 U 0.91 U 1.4 U	1.5 U 0.98 UJ 0.91 U 1.4 U	1.5 U 0.98 U 0.91 U 1.4 U	1.5 U 0.98 U 0.91 U 1.4 U	1.5 U 0.98 U 0.91 U 1.4 U	1.5 U 0.98 UJ 0.91 U 1.4 U	1.5 U 0.98 UJ 0.91 U 1.4 U	1.5 U 0.98 U 0.91 U 1.4 U	1.5 U 0.98 U 0.91 U 1.4 U	1.5 U 0.98 U 0.91 U 1.4 U	1.5 U 0.98 U 0.91 U 1.4 U
4-Nitroaniline 4-Nitrophenol Acenaphthene Acenaphthylene	μg/L μg/L μg/L	. 1.3 U . 1.3 U	2.5 U 5.8 U 1.3 U 1.3 U	2.5 UJ 5.8 U 1.3 U 1.3 U	2.5 U 5.8 U 1.3 U 1.3 U	2.5 U 5.8 U 1.3 U 1.3 U	2.5 U 5.8 U 1.3 U 1.3 U	2.5 UJ 5.8 U 1.3 U 1.3 U	2.5 UJ 5.8 U 1.3 U 1.3 U	2.5 U 5.8 U 1.3 U 1.3 U	2.5 U 5.8 U 1.3 U 1.3 U	2.5 U 5.8 U 1.3 U 1.3 U	2.5 U 5.8 U 1.3 U 1.3 U
Anthracene Benzo(a)anthracene Benzo(a)pyrene Benzo(b)fluoranthene Benzo(g,h,i)perylene	µg/L µg/L µg/L µg/L µg/L	. 1.2 U . 1.5 U . 1.1 U . 1.1 U . 0.91 U	1.2 U 1.5 U 1.1 U 1.1 U 0.91 U										
Benzo(k)fluoranthene Benzoic acid Benzyl alcohol bis(2-Chloroethoxy)methane	ру.с µg/L µg/L µg/L	. 1.1 U . 33 U . 1.5 U . 1.8 U	1.1 U 33 U 1.5 U 1.8 U	1.1 U 33 U 1.5 U 1.8 U	1.1 U 33 U 1.5 U 1.8 U	1.1 U 33 U 1.5 U 1.8 U	1.1 U 33 U 1.5 U 1.8 U	1.1 U 33 U 1.5 U 1.8 U	1.1 U 33 U 1.5 U 1.8 U	1.1 U 33 U 1.5 U 1.8 U	1.1 U 33 U 1.5 U 1.8 U	1.1 U 33 U 1.5 U 1.8 U	1.1 U 33 U 1.5 U 1.8 U
bis(2-Chloroethyl)ether bis(2-Ethylhexyl)phthalate (DEHP) Butyl benzylphthalate (BBP) Chrysene	µg/L µg/L µg/L µg/L	. 0.91 U . 1.3 U . 1.1 U	1.2 U 0.91 U 1.3 U 1.1 U	1.2 U 0.91 U 1.3 U 1.1 U	1.2 UJ 2.2 J 1.3 U 1.1 U	1.2 UJ 0.91 U 1.3 U 1.1 U	1.2 UJ 0.91 U 1.3 U 1.1 U	1.2 U 0.91 U 1.3 U 1.1 U	1.2 U 3.7 J 1.3 U 1.1 U	1.2 U 3.9 J 1.3 U 1.1 U	1.2 U 0.91 U 1.3 U 1.1 U	1.2 U 0.91 U 1.3 U 1.1 U	1.2 U 0.91 U 1.3 U 1.1 U
Dibenz(a,h)anthracene Dibenzofuran Diethyl phthalate Dimethyl phthalate Di-n-butylphthalate (DBP)	µg/L µg/L µg/L µg/L µg/L	. 1.3 U . 0.98 U . 1.2 U	0.93 U 1.3 U 0.98 U 1.2 U 1.9 U	0.93 U 1.3 U 0.98 U 1.2 U 1.9 U	0.93 U 1.3 U 0.98 U 1.2 U 1.9 U	0.93 U 1.3 U 0.98 U 1.2 U 1.9 U	0.93 U 1.3 U 0.98 U 1.2 U 1.9 U	0.93 U 1.3 U 0.98 U 1.2 U 1.9 U	0.93 U 1.3 U 0.98 U 1.2 U 1.9 U	0.93 U 1.3 U 0.98 U 1.2 U 2.3 J	0.93 U 1.3 U 0.98 U 1.2 U 1.9 U	0.93 U 1.3 U 0.98 U 1.2 U 1.9 U	0.93 U 1.3 U 0.98 U 1.2 U 1.9 U
Di-n-octyl phthalate (DnOP)	μg/L		3.0 U										

2020 Analytical Results Summary - Bedrock Love Canal Long-Term Monitoring Program Niagara Falls, New York

	Sample Location: Sample ID: Sample Date:	3257 WG-9954-062620-RM-018 6/26/2020	5221 WG-9954-070120-SG-025 7/1/2020	6209 WG-9954-062520-SG-011 6/25/2020	7205 WG-9954-062620-RM-014 6/26/2020	8210 WG-9954-062620-RM-015 6/26/2020	9205 WG-9954-062620-RM-016 6/26/2020	9210 WG-9954-062520-SG-012 6/25/2020	10205 WG-9954-062520-SG-013 6/25/2020	10210A WG-9954-070920-SG-033 7/9/2020	10210B WG-9954-063020-SG-019 6/30/2020	10210C WG-9954-063020-SG-020 6/30/2020	10210C WG-9954-063020-SG-021 6/30/2020 (Duplicate)
Parameters	Units												(Duplicato)
Semi-volatile Organic Compounds (Con	ntinued)												
Fluoranthene	μg/L	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U
Fluorene	μg/L	1.2 U	1.2 U	1.2 U	1.2 U	1.2 U	1.2 U	1.2 U	1.2 U	1.2 U	1.2 U	1.2 U	1.2 U
Hexachlorobenzene	μg/L	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U
Hexachlorobutadiene Hexachlorocyclopentadiene	μg/L	0.91 U	0.91 U 2.0 U	0.91 U 2.0 U	0.91 U	0.91 U	0.91 U	0.91 U 2.0 U	0.91 U	0.91 U 2.0 U	0.91 U 2.0 U	0.91 U	0.91 U
Hexachloroethane	μg/L μg/L	2.0 U 0.96 U	0.96 U	0.96 U	2.0 U 0.96 U	2.0 U 0.96 U	2.0 U 0.96 U	0.96 U	2.0 U 0.96 U	0.96 U	0.96 U	2.0 U 0.96 U	2.0 U 0.96 U
Indeno(1,2,3-cd)pyrene	μg/L	1.6 U	1.6 U	1.6 U	1.6 U	1.6 U	1.6 U	1.6 U	1.6 U	1.6 U	1.6 U	1.6 U	1.6 U
Isophorone	μg/L	1.3 U	1.3 U	1.3 UJ	1.3 U	1.3 U	1.3 U	1.3 UJ	1.3 UJ	1.3 U	1.3 U	1.3 U	1.3 U
Naphthalene	μg/L	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U
Nitrobenzene	μg/L	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U
N-Nitrosodi-n-propylamine	μg/L	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U
N-Nitrosodiphenylamine	μg/L	2.4 U	2.4 U	2.4 U	2.4 U	2.4 U	2.4 U	2.4 U	2.4 U	2.4 U	2.4 U	2.4 U	2.4 U
Pentachlorophenol	μg/L	8.9 U	8.9 U	8.9 U	8.9 U	8.9 U	8.9 U	8.9 U	8.9 U	8.9 U	8.9 U	8.9 U	8.9 U
Phenanthrene	μg/L	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U
Phenol	μg/L	0.91 U 1.3 U	0.91 U 1.3 U	0.91 U	0.91 U 1.3 U	0.91 U 1.3 U	0.91 U 1.3 U	0.91 U 1.3 U	0.91 U 1.3 U	0.91 U 1.3 U	0.91 U	0.91 U 1.3 U	0.91 U 1.3 U
Pyrene	μg/L	1.3 0	1.3 0	1.3 U	1.3 0	1.5 0	1.3 0	1.3 0	1.5 0	1.3 0	1.3 U	1.3 0	1.5 0
Discret	te Compounds Detected:	0	0	0	1	0	0	0	1	2	0	0	0
Polychlorinated Biphenyls													
Aroclor-1016 (PCB-1016)	μg/L	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U
Aroclor-1221 (PCB-1221)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1232 (PCB-1232)	μg/L	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U
Aroclor-1242 (PCB-1242)	μg/L	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U
Aroclor-1248 (PCB-1248) Aroclor-1254 (PCB-1254)	μg/L	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U
Aroclor-1264 (PCB-1264) Aroclor-1260 (PCB-1260)	μg/L μg/L	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U
Discret	te Compounds Detected:	0	0	0	0	0	0	0	0	0	0	0	0
		·	v	·	,	•	Č	,	v	· ·	v	· ·	·
Pesticides 4,4'-DDD	μg/L	R	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U
4,4'-DDE	μg/L	0.029 J	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U
4,4'-DDT	μg/L	R	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U
Aldrin	μg/L	R	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 J	0.019 U	0.019 U	0.019 U	0.019 U
alpha-BHC Pesticides (Continued)	μg/L	R	0.13	0.31	0.088	0.064	0.097	0.033 J	0.18	0.12	0.12	0.081	0.10
alpha-Chlordane	ug/l	R	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U
beta-BHC	μg/L μg/L	0.045 J	0.019 U	0.019 U	0.036 J	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U
delta-BHC	μg/L	R	0.029 J	0.13	0.32	0.042 J	0.10	0.068	0.28	0.088	0.076	0.068 J	0.12 J
Dieldrin	μg/L	R	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U
Endosulfan I	μg/L	0.037 J	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U
Endosulfan II	μg/L	R	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U
Endosulfan sulfate	μg/L	R -	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U
Endrin	μg/L	R	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U
Endrin ketone gamma-BHC (lindane)	μg/L	R 0.026 J	0.019 U 0.076	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U					
gamma-Chlordane	μg/L μg/l	0.026 J 0.020 NJ	0.076 0.019 U	0.25 0.019 U	0.11 0.019 U	0.060 0.019 U	0.096 0.019 U	0.037 J 0.028 J	0.21 0.019 U	0.15 0.019 U	0.10 0.019 U	0.089 0.019 U	0.11 0.019 U
Heptachlor	μg/L μg/L	0.020 NO R	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.028 J 0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U
Heptachlor epoxide	μg/L	R	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U
Methoxychlor	μg/L	R	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 UJ	0.019 U	0.019 U	0.019 U
Toxaphene	μg/L	R	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U
Discret	te Compounds Detected:	5	3	3	4	3	3	4	4	3	3	3	3

2020 Analytical Results Summary - Bedrock Love Canal Long-Term Monitoring Program Niagara Falls, New York

	ple Location: Sample ID: Sample Date:	10215 WG-9954-063020-SG-022 6/30/2020	10225A WG-9954-070920-SG-034 7/9/2020	10225B WG-9954-070920-SG-035 7/9/2020	10225C WG-9954-070120-SG-026 7/1/2020	10270 WG-9954-070720-SG-032 7/7/2020	10272 WG-9954-070720-SG-030 7/7/2020	10278 WG-9954-070620-SG-029 7/6/2020	MW-01 WG-9954-063020-RM-023 6/30/2020	MW-02 WG-9954-070120-SG-024 7/1/2020
Parameters	Units									
Volatile Organic Compounds 1,1,1-Trichloroethane 1,1,2,2-Tetrachloroethane 1,1,2-Trichloroethane 1,1-Dichloroethane 1,1-Dichloroethane 1,2-Dichloroethane 1,2-Dichloroethane 1,2-Dichloropropane 2-Butanone (Methyl ethyl ketone) (MEK) 2-Hexanone 4-Methyl-2-pentanone (Methyl isobutyl ketone) (MIBK) Acetone Benzene Bromodichloromethane Bromoform Bromomethane (Methyl bromide) Carbon disulfide Carbon tetrachloride Chlorobenzene Chloroethane Chloroform (Trichloromethane) Chloromethane (Methyl chloride) cis-1,2-Dichloroethene cis-1,3-Dichloropropene Dibromochloromethane Ethylbenzene	µg/L µg/L µg/L µg/L µg/L µg/L µg/L µg/L	0.20 U 0.78 U 0.20 U 0.20 U 0.20 U 0.20 U 0.20 U 0.25 U 0.70 U 3.5 J 0.34 U 0.20 U 0.23 U 0.24 U 0.28 U 0.23 U 0.20 U 0.20 U 0.20 U	1.0 U 3.9 U 1.0 U 1.0 U 25 U 1.0 U 1.3 U 3.5 U 160 1.7 U 1.0 U 1.2 U 1.2 U 1.4 U 1.2 U 1.0 U	1.0 U 3.9 U 1.0 U 1.0 U 25 U 1.0 U 1.3 U 3.5 U 25 J 1.7 U 1.0 U 1.2 U 1.2 U 1.4 U 1.2 U 1.0 U 1.0 U 1.0 U 1.0 U	0.20 U 0.25 U 0.70 U 3.3 J 0.34 U 0.89 J 0.23 U 0.24 U 0.28 U 3.3 J 0.20 U 0.20 U	0.20 U 0.25 U 0.70 U 3.7 J 0.34 U 0.20 U 0.23 U 0.24 U 0.28 U 0.23 U 0.20 U 0.20 U 0.20 U 0.20 U	0.20 U 0.25 U 0.70 U 3.4 J 0.34 U 0.20 U 0.23 U 0.24 U 0.28 U 0.20 U 0.20 U 0.20 U 0.20 U	0.20 U 0.21 U 0.22 U 0.25 U 0.70 U 6.4 J 0.34 U 0.20 U 0.23 U 0.24 U 0.28 U 0.20 U 0.20 U 0.20 U 0.20 U 0.20 U	0.20 U 0.25 U 0.70 U 2.9 J 0.34 U 0.20 U 0.23 U 0.24 U 0.28 U 0.20 U 0.20 U 0.20 U 0.20 U	0.20 U 0.25 U 0.70 U 0.42 U 0.34 U 0.20 U 0.23 U 0.24 U 0.28 U 0.23 U 0.20 U
Methylene chloride Styrene Tetrachloroethene Toluene trans-1,2-Dichloroethene trans-1,3-Dichloropropene Trichloroethene Vinyl acetate Vinyl chloride Xylenes (total)	µg/L µg/L µg/L µg/L µg/L µg/L µg/L µg/L	0.20 U 0.65 U 0.20 U 0.21 U 0.20 U 0.23 U 0.20 U 1.1 U 0.20 U 0.23 U	1.0 U 3.3 U 1.0 U 1.1 U 1.0 U 1.2 U 1.0 U 5.5 U 1.0 U 1.2 U	1.0 U 1.0 U 1.1 U 1.0 U 1.0 U 1.2 U 1.0 U 5.5 U 1.0 U 1.2 U	0.20 U 0.65 U 0.20 U 0.21 U 0.20 U 0.23 U 6.5 1.1 U 0.20 U 0.23 U	0.20 U 0.65 U 0.20 U 0.21 U 0.20 U 0.23 U 0.20 U 1.1 U 0.20 U 0.23 U	0.20 U 0.65 U 0.20 U 0.21 U 0.20 U 0.23 U 0.20 U 1.1 U 0.20 U 0.23 U	0.20 U 0.65 U 0.20 U 0.21 U 0.20 U 0.23 U 0.20 U 1.1 U 0.20 U 0.23 U	0.20 U 0.65 U 0.20 U 0.21 U 0.20 U 0.23 U 0.20 U 1.1 U 0.20 U 0.23 U	0.20 U 0.65 U 0.20 U 0.21 U 0.20 U 0.23 U 0.20 U 1.1 U 0.20 U 0.23 U
Discrete Compour	nds Detected:	1	1	1	4	1	1	1	1	0
Semi-volatile Organic Compounds 1,2,4-Trichlorobenzene 1,3-Dichlorobenzene 1,4-Dichlorobenzene 2,2'-Oxybis(1-chloropropane) (bis(2-Chloroisopropyl) e 2,4,5-Trichlorophenol 2,4,6-Trichlorophenol 2,4-Dirichlorophenol 2,4-Dinitrotoluene 2,6-Dinitrotoluene 2,6-Dinitrotoluene 2,6-Dinitrotoluene 2-Chloronaphthalene 2-Chlorophenol 2-Methylphenol 2-Mitroaniline 2-Nitroaniline 2-Nitrophenol 3,3'-Dichlorobenzidine 3-Nitroaniline 4,6-Dinitro-2-methylphenol 4-Bromophenyl phenyl ether 4-Chloro-3-methylphenol 4-Chloroaniline 4-Nitrophenol Acenaphthene Acenaphthylene Acenaphthylene Anthracene Benzo(a)anthracene Benzo(b)fluoranthene Benzo(b)fluoranthene Benzo(b)fluoranthene Benzo(c)acid Benzyl alcohol bis(2-Chloroethoxy)methane Dibenzofuran Diethyl phthalate Dimethyl phthalate (DBP) Di-n-octyl phthalate (DnOP)	µg/L µg/L µg/L µg/L µg/L µg/L µg/L µg/L	1.1 U 1.1 U 1.1 U 1.92 U 1.1 U 1.3 U 0.99 U 1.3 U 1.2 U 1.3 U 1.9 U 2.2 U 1.2 U 1.3 U 0.97 U 1.2 U 0.91 U 1.3 U 1.1 U 1.1 U 2.3 U 1.5 U 0.98 U 0.91 U 1.4 U 2.5 U 5.8 U 1.3 U 1.5 U 1.1 U	1.1 U 1.1 U 1.92 U 1.1 U 1.3 U 0.99 U 1.3 U 1.2 U 1.3 U 1.9 U 2.2 U 1.2 U 1.3 U 0.97 U 1.2 U 0.91 U 1.3 U 1.4 U 1.1 U 2.3 U 1.5 U 0.98 U 0.91 U 1.4 U 2.5 U 5.8 U 1.3 U 1.5 U 1.1 U 1.1 U 0.91 U 1.1 U 1.1 U 0.91 U 0.9	1.1 U 1.1 U 1.92 U 1.1 U 1.3 U 1.99 U 1.3 U 1.2 U 1.3 U 1.9 U 2.2 U 1.2 U 1.3 U 0.97 U 1.4 U 1.1 U 1.1 U 2.3 U 1.4 U 1.1 U 2.3 U 1.5 U 0.98 U 0.91 U 1.4 U 2.5 U 5.8 U 1.3 U 1.2 U 1.5 U 1.1 U 1.1 U 0.91 U 1.1 U 1.1 U 0.91	7.7 J 1.1 U 0.92 U 1.1 U 1.3 U 0.99 U 1.3 U 1.2 U 1.3 U 1.9 U 2.2 U 1.2 U 1.3 U 0.97 U 1.2 U 0.91 U 1.1 U 1.1 U 2.3 U 1.5 U 0.98 U 0.91 U 1.4 U 2.5 U 5.8 U 1.3 U 1.2 U 1.5 U 1.1 U 0.91 U 1.1 U 1.1 U 0.91 U 1.1 U 1.1 U 0.91 U 0.9	1.1 U 1.1 U 1.92 U 1.1 U 1.3 U 0.99 U 1.3 U 1.2 U 1.3 U 1.9 U 1.2 U 1.3 U 0.97 U 1.2 U 0.91 U 1.3 U 1.4 U 1.1 U 2.3 U 1.5 U 0.98 U 0.91 U 1.4 U 2.5 U 1.5 U 1.1 U	1.1 U 1.1 U 1.92 U 1.1 U 1.3 U 0.99 U 1.3 U 1.2 U 1.3 U 1.9 U 2.2 U 1.2 U 1.3 U 0.97 U 1.1 U 0.91 U 1.3 U 1.1 U 2.3 U 1.4 U 1.1 U 2.3 U 1.5 U 0.98 U 0.91 U 1.4 U 2.5 U 5.8 U 1.3 U 1.2 U 1.5 U 1.1 U 1.1 U 0.91 U 1.1 U 1.1 U 0.91 U	1.1 U 1.1 U 1.1 U 1.92 U 1.1 U 1.3 U 0.99 U 1.3 U 1.2 U 1.3 U 1.9 U 2.2 U 1.2 U 1.3 U 0.97 U 1.2 U 0.91 U 1.1 U 1.1 U 1.1 U 2.3 U 1.5 U 0.98 U 0.91 U 1.4 U 2.5 U 5.8 U 1.3 U 1.3 U 1.5 U 1.1 U 0.91 U 1.1 U 1.1 U 1.1 U 0.91 U 1.3 U 0.98 U 1.2 U 0.91 U 1.3 U 0.98 U 1.2 U 1.9 U 1.9 U 3.0 U	1.1 U 1.1 U 1.1 U 1.92 U 1.1 U 1.3 U 0.99 U 1.3 U 1.2 U 1.3 U 1.9 U 2.2 U 1.2 U 1.3 U 0.97 U 1.2 U 0.91 U 1.1 U 1.1 U 1.1 U 2.3 U 1.5 U 0.98 U 0.91 U 1.4 U 2.5 U 5.8 U 1.3 U 1.2 U 1.5 U 1.1 U 0.91 U 1.1 U 0.91 U 1.1 U 0.91 U 1.1 U 0.91 U 1.1 U 0.93 U 1.2 U 0.91 U 1.3 U 1.9 U 1.9 U 1.9 U 3.0 U	1.1 U 1.1 U 1.3 U 1.99 U 1.3 U 1.2 U 1.3 U 1.9 U 2.2 U 1.2 U 1.3 U 0.97 U 1.2 U 0.91 U 1.3 U 1.1 U 2.3 U 1.4 U 1.1 U 2.3 U 1.8 U 1.5 U 0.98 U 0.91 U 1.4 U 2.5 U 5.8 U 1.3 U 1.2 U 1.5 U 1.1 U 0.91 U 0.9

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2020 Analytical Results Summary - Bedrock Love Canal Long-Term Monitoring Program Niagara Falls, New York

	Sample Location: Sample ID: Sample Date:	10215 WG-9954-063020-SG-022 6/30/2020	10225A WG-9954-070920-SG-034 7/9/2020	10225B WG-9954-070920-SG-035 7/9/2020	10225C WG-9954-070120-SG-026 7/1/2020	10270 WG-9954-070720-SG-032 7/7/2020	10272 WG-9954-070720-SG-030 7/7/2020	10278 WG-9954-070620-SG-029 7/6/2020	MW-01 WG-9954-063020-RM-023 6/30/2020	MW-02 WG-9954-070120-SG-024 7/1/2020
Parameters	Units									
Semi-volatile Organic Compou	nds (Continued)									
Fluoranthene	μg/L	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U
Fluorene	μg/L	1.2 U	1.2 U	1.2 U	1.2 U	1.2 U	1.2 U	1.2 U	1.2 U	1.2 U
Hexachlorobenzene	μg/L	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U
Hexachlorobutadiene	μg/L	0.91 U 2.0 U	0.91 U 2.0 U	0.91 U 2.0 U	0.91 U 2.0 U	0.91 U 2.0 U	0.91 U R	0.91 U 2.0 U	0.91 U 2.0 U	0.91 U 2.0 U
Hexachlorocyclopentadiene Hexachloroethane	μg/L μg/L	0.96 U	0.96 U	0.96 U	0.96 U	0.96 U	0.96 U	0.96 U	0.96 U	0.96 U
Indeno(1,2,3-cd)pyrene	μg/L	1.6 U	1.6 U	1.6 U	1.6 U	1.6 U	1.6 U	1.6 U	1.6 U	1.6 U
Isophorone	μg/L	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U
Naphthalene	μg/L	1.1 U	1.2 J	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U
Nitrobenzene	μg/L	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U	1.4 U
N-Nitrosodi-n-propylamine	μg/L	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U	1.1 U
N-Nitrosodiphenylamine	μg/L	2.4 U	2.4 U	2.4 U	2.4 U	2.4 U	2.4 U	2.4 U	2.4 U	2.4 U
Pentachlorophenol	μg/L	8.9 U	8.9 U	8.9 U	8.9 U	8.9 U	8.9 U	8.9 U	8.9 U	8.9 U
Phenanthrene Phenol	μg/L	1.3 U 0.91 U	1.3 U 0.91 U	1.3 U 0.91 U	1.3 U 0.91 U	1.3 U 0.91 U	1.3 U 0.91 U	1.3 U 0.91 U	1.3 U 0.91 U	1.3 U 0.91 U
Pyrene	μg/L μg/L	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U	1.3 U
1 yrone	۳۹٬ ـ	1.0 0	1.0 0	1.5 5	1.0 0	1.0 0	1.5 6	1.5 C	1.0 0	1.0 0
	Discrete Compounds Detected:	0	1	0	1	0	0	0	0	0
Polychlorinated Biphenyls										
Aroclor-1016 (PCB-1016)	μg/L	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U
Aroclor-1221 (PCB-1221)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1232 (PCB-1232)	μg/L	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U
Aroclor-1242 (PCB-1242)	μg/L	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U
Aroclor-1248 (PCB-1248)	μg/L	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U
Aroclor-1254 (PCB-1254) Aroclor-1260 (PCB-1260)	μg/L μg/L	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U	0.46 U 0.46 U
A100101-1200 (1 OB-1200)										
	Discrete Compounds Detected:	0	0	0	0	0	0	0	0	0
Pesticides										
4,4'-DDD	μg/L	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U
4,4'-DDE	μg/L	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U
4,4'-DDT	μg/L	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U
Aldrin	μg/L	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U
alpha-BHC Pesticides (Continued)	μg/L	0.064	0.069	0.072	0.27	0.070	0.24	0.47	0.019 U	0.019 U
alpha-Chlordane	μg/L	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U
beta-BHC	μg/L	0.013 J	0.019 U	0.019 J	0.040 J	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U
delta-BHC	μg/L	0.29	0.065	0.19	0.17	0.067	0.084	0.13	0.019 U	0.019 U
Dieldrin	μg/L	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U
Endosulfan I	μg/L	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U
Endosulfan II	μg/L	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U
Endosulfan sulfate	μg/L	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U
Endrin Endrin ketone	μg/L	0.019 U 0.019 U	0.019 U 0.019 U	0.019 U 0.019 U	0.019 U 0.019 U	0.019 U 0.019 U	0.019 U 0.019 U	0.019 U 0.019 U	0.019 U 0.019 U	0.019 U 0.019 U
gamma-BHC (lindane)	μg/L μg/L	0.074	0.019 0	0.019 0	0.019 0	0.019 0	0.019 0	0.48	0.019 U	0.019 U
gamma-Chlordane	μg/L μg/L	0.019 U	0.017 0.019 U	0.019 U	0.21 0.019 U	0.019 U	0.20 0.019 U	0.48 0.019 U	0.019 U	0.019 U
Heptachlor	μg/L	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U
Heptachlor epoxide	μg/L	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U
Methoxychlor	μg/L	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U	0.019 U
Toxaphene	μg/L	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U
	Discrete Compounds Detected:	4	3	4	4	3	3	3	0	0

Notes:

J - Estimated concentration

U - Not detected at the associated reporting limit
UJ - Not detected; associated reporting limit is estimated

R - Rejected
NJ - Tentatively identified compound, estimated concentration

2020 Detection Summary -Love Canal Long-Term Monitoring Program Niagara Falls, New York

Overburden Wells			Number of Discret	te Compounds Det	ected
without well 10135)	Well Group	VOCs	SVOCs	PCBs	Pesticides
7115	_			_	
	В	0	1	0	0
7125	В	0	0	0	0
7130	A	0	0	0	0
7132	A	1	0	0	0
8106	Α	1	0	0	0
8115	В	0	0	0	0
8125	В	0	0	0	0
9105					
	В	0	0	0	0
9113	В	2	0	0	0
9118	В	2 (1)	0	0	0
10178A	В	0	0	0	0
10178B	***	0	1	0	4
Total Overburden Well					
Detections		6 (5)	2	0	4
Number of Discrete Compounds Detected		3	1	0	4
Well 10135	Well Group	VOCs	SVOCs	PCBs	Pesticides
		1000			
10135	A	17 (19)	16	0	5
Total Well 10135 Well		4		_	
Detections Number of Discrete		17 (19)	16	0	5
Compounds Detected		17 (19)	16	0	5
Bedrock Wells	Well Group	VOCs	SVOCs	PCBs	Pesticides
3257	Α	1	0	0	5*
5221	А	0	0	0	3
			0	0	3
6209	A	1			
7205	Α	3	1	0	4
8210	A	2	0	0	3
9205	Α	1	0	0	3
9210	Α	2	0	0	4
10205	А	1	1	0	4
10210A	Α	1	2	0	3
10210B	Α	1	0	0	3
10210C	Α	1	0	0	3
10215	Α	1	0	0	4
10225A	Α	1	1	0	3
10225B	А	1	0	0	4
10225C	Α	4	1	0	4
10270	A	1	0	0	3
10272	A	1	0	0	3
10278	Α	1	0	0	3
MW-01	X	1	0	0	0
MW-02	Х	0	0	0	0
Total Bedrock Well				_	<u></u>
Detections Number of Discrete		25	6	0	62
Compounds Detected		6	4	0	8
Notes:					
* - A portion of the data was	rejected during data valida	ation; rejected data i	s not included in tota	al. Rejected results w	vere non-detect.
A - Annual Well		•		,	
B - Biannual Well					
X - Additional annual well ad	ded to program in 2011				
*** - Well requested to be sa	mpled in 2020 by NYSDE	С			
() - Results for duplicate sar	nple, if different from pare	nt sample			
	· · ·				
PCBs - Polychlorinated Rinha	envis '				
PCBs - Polychlorinated Bipho SVOCs - Semi-Volatile Orga					

	Well Number: SampleDate:	10210A 7/24/1990	10210A 8/22/1991	10210A 8/26/1992	10210A 8/11/1993	10210A 5/25/1995	10210A 7/1/1996	10210A 7/10/1997	10210A 6/26/1998	10210A 6/23/1999	10210A 6/21/2000	10210A 5/18/2001	10210A 6/13/2002	10210A 5/27/2003	10210A 6/3/2004	10210A 6/28/2005	10210A 7/6/2006
Parameters																	
Volatile Organic Compounds (μg/L)																	
1,1,1-Trichloroethane																	
1,1,2,2-Tetrachloroethane					0.2												
1,1,2-Trichloroethane																	
1,1-Dichloroethane																	
1,1-Dichloroethene																	
1,2-Dichloroethane																	
1,2-Dichloroethene (total)																	
1,2-Dichloropropane					0.04												
2-Butanone (Methyl Ethyl Ketone)										2 J			3 J		4 J		
2-Hexanone					0.4					3 J							
Acetone		14C			13B				120 J			10 J					
Benzene					0.1												
Bromodichloromethane																	
Bromoform					0.03												
Bromomethane (Methyl bromide)																	
Carbon Disulfide					2	20	310					6 J			6 J	1.6 J	1 J
Chlorobenzene																	
Chloroform																	
Chlorotoluenes																	
cis-1,2-Dichloroethene																	
Dibromochloromethane					0.08												
Dichlorotoluene, total																	
Ethylbenzene					0.6												
m&p-Xylenes																	
Methylene Chloride					0.3												
o-Xylene																	
Styrene					0.1												
Tetrachloroethene					0.07												
Toluene					0.4					2 J						2.3 J	
trans-1,2-Dichloroethene																	
Trichloroethene					0.1												
Trichlorotoluene, total																	
Vinyl Acetate																	
Vinyl Chloride																	
Xylenes (total)					1												
Semi-volatile Organic Compounds (µ	a/L)																
1,2,4-Trichlorobenzene	. ,																
1,2-Dichlorobenzene																	
1,3-Dichlorobenzene																	
1,4-Dichlorobenzene																	
2,4,5-Trichlorophenol																	
_, .,o monorophonor																	

	Well Number: SampleDate:	10210A 7/24/1990	10210A 8/22/1991	10210A 8/26/1992	10210A 8/11/1993	10210A 5/25/1995	10210A 7/1/1996	10210A 7/10/1997	10210A 6/26/1998	10210A 6/23/1999	10210A 6/21/2000	10210A 5/18/2001	10210A 6/13/2002	10210A 5/27/2003	10210A 6/3/2004	10210A 6/28/2005	10210A 7/6/2006
Davamatava	Jan., p. 13 a. 13		0	0.20.1002		0.20.1000	.,		0.20.1000	0.20.1000	0.22000	0.10/2001	3.13.232	0.2.,2000	0.0.2001	0.20.200	
Parameters																	
Semi-volatile Organic Compounds	(μg/L) (Continued)																
2,4,6-Trichlorophenol																	
2,4-Dichlorophenol																	
2,4-Dimethylphenol																	
2-Chloronaphthalene																	
2-Chlorophenol																	
2-Methylnaphthalene																	
2-Methylphenol																	
3&4-Methylphenol																	
2-Nitrophenol																	
3,5-Dichlorotoluene																	
4-Chloro-3-methylphenol																	
4-Chlorophenyl phenyl ether																	
4-Methylphenol																	
Acetic acid																	
Anthracene																	
Benzo(a)pyrene																	
Benzo(b)fluoranthene																	
Benzo(g,h,i)perylene																	
Benzo(k)fluoranthene																	
Benzoic Acid								12 J						3 J	3 J	2.7 J	
Benzyl Alcohol																	
Bis(2-Chloroethyl)Ether																	
bis(2-Ethylhexyl)Phthalate			12	21	31	51									1 J	1.7 J	8 J
Butyl benzylphthalate (BBP)					3												
Camphor																	
Carbazole																	
Chlorobenzoic acid																	
Chrysene																	
Dibenz(a,h)anthracene																	
Diethyl phthalate																	
Dimethyl Phthalate		16															
Dimethyl tetrasulfide					22												
Di-n-butyl phthalate (DBP)			2		0.9												
Di-n-octyl phthalate (DnOP)		3B															
Fluoranthene																	
Hexachlorobenzene																	
Hexachloroethane																	
Indeno(1,2,3-cd)pyrene																	
Naphthalene																	
N-Nitrosodiphenylamine																	
Pentachlorophenol																	
Phenanthrene																	
Phenol										1 J				5 J	1 J	1.7 J	
Pyrene																	

	Well Number:	10210A	10210A	10210A	10210A	10210A	10210A	10210A	10210A	10210A	10210A	10210A	10210A	10210A	10210A	10210A	10210A
	SampleDate:	7/24/1990	8/22/1991	8/26/1992	8/11/1993	5/25/1995	7/1/1996	7/10/1997	6/26/1998	6/23/1999	6/21/2000	5/18/2001	6/13/2002	5/27/2003	6/3/2004	6/28/2005	7/6/2006
Parameters																	
Pesticides/PCBs (μg/L)																	
4,4'-DDD														0.013 J			
4,4'-DDE																	
Aldrin																	
Alpha-BHC										0.28							
Alpha-Chlordane																	
Aroclor-1260 (PCB-1260)																	
beta&gamma-BHC (sum of isomers)																	
Beta-BHC										0.035 J				0.020 J	0.011 J		
Delta-BHC					0.0061									0.062 J	0.043 J		
Dieldrin																	
Endosulfan I										0.046 J							
Endosulfan II																	
Endosulfan Sulfate																	
Endrin																	
Endrin ketone																	
Gamma-BHC (Lindane)										0.10 J				0.039 J			
Gamma-Chlordane																	
Heptachlor																	
Heptachlor epoxide																	
Methoxychlor																	

	Well Number: SampleDate:	10210A 7/26/2007	10210A 7/17/2008	10210A 7/15/2009	10210A 6/24/2010	10210A 7/19/2011	10210A 6/22/2012	10210A 6/13/2013	10210A 6/27/2014	10210A 6/26/2015	10210A 6/24/2016	10210A 7/10/2017	10210A 7/23/2018	10210A 6/14/2019	10210A 7/9/2020	10210B 7/24/1990	10210B 8/22/1991
Parameters																	
Volatile Organic Compounds (μg/L)																	
1,1,1-Trichloroethane																	
1,1,2,2-Tetrachloroethane																	
1,1,2-Trichloroethane																	
1,1-Dichloroethane																	
1,1-Dichloroethene																	
1,2-Dichloroethane																	
1,2-Dichloroethene (total)																	
1,2-Dichloropropane																	
2-Butanone (Methyl Ethyl Ketone)																	
2-Hexanone																	
Acetone					5.2 J						3.5 J	8.4 J		5.8 J			
Benzene																	
Bromodichloromethane																	
Bromoform																	
Bromomethane (Methyl bromide)											6.1						
Carbon Disulfide		8 J	24				2.7 J			2.9 J		2.0 J	340	51	9.2 J		
Chlorobenzene																	
Chloroform																	
Chlorotoluenes																	
cis-1,2-Dichloroethene																	
Dibromochloromethane																	
Dichlorotoluene, total																	
Ethylbenzene											0.28 J	0.35 J	0.22 J				
m&p-Xylenes																	
Methylene Chloride																	
o-Xylene																	
Styrene																	
Tetrachloroethene																	
Toluene																1.8	
trans-1,2-Dichloroethene																	
Trichloroethene					6.3												
Trichlorotoluene, total																	
Vinyl Acetate																	
Vinyl Chloride																	
Xylenes (total)													0.27 J	0.29 J			
Semi-volatile Organic Compounds (μ	g/L)																
1,2,4-Trichlorobenzene																	
1,2-Dichlorobenzene																	
1,3-Dichlorobenzene																	
1,4-Dichlorobenzene																	
2,4,5-Trichlorophenol																	

	Well Number: SampleDate:	10210A 7/26/2007	10210A 7/17/2008	10210A 7/15/2009	10210A 6/24/2010	10210A 7/19/2011	10210A 6/22/2012	10210A 6/13/2013	10210A 6/27/2014	10210A 6/26/2015	10210A 6/24/2016	10210A 7/10/2017	10210A 7/23/2018	10210A 6/14/2019	10210A 7/9/2020	10210B 7/24/1990	10210B 8/22/1991
Davamatava	-																
Parameters																	
Semi-volatile Organic Compounds (μg/L) (Continued)																
2,4,6-Trichlorophenol																	
2,4-Dichlorophenol																	
2,4-Dimethylphenol																	
2-Chloronaphthalene																	
2-Chlorophenol																	
2-Methylnaphthalene																	
2-Methylphenol																	
3&4-Methylphenol																	
2-Nitrophenol																	
3,5-Dichlorotoluene																	
4-Chloro-3-methylphenol																	
4-Chlorophenyl phenyl ether																	
4-Methylphenol																	
Acetic acid																	
Anthracene																	
Benzo(a)pyrene																	
Benzo(b)fluoranthene																	
Benzo(g,h,i)perylene																	
Benzo(k)fluoranthene																	
Benzoic Acid				5.8 J													
Benzyl Alcohol																	
Bis(2-Chloroethyl)Ether																	
bis(2-Ethylhexyl)Phthalate				2.5 J											3.9 J	7 B	13
Butyl benzylphthalate (BBP)																	
Camphor																	
Carbazole																	
Chlorobenzoic acid																	
Chrysene																	
Dibenz(a,h)anthracene																	
Diethyl phthalate																	
Dimethyl Phthalate																	
Dimethyl tetrasulfide																	
Di-n-butyl phthalate (DBP)															2.3 J	1	1
Di-n-octyl phthalate (DnOP)																	
Fluoranthene																	
Hexachlorobenzene																	
Hexachloroethane																	
Indeno(1,2,3-cd)pyrene																	
Naphthalene										0.41 J							
N-Nitrosodiphenylamine																	
Pentachlorophenol																1	
Phenanthrene																	
Phenol																3	3
Pyrene																	

	Well Number:	10210A	10210A	10210B	10210B												
	SampleDate:	7/26/2007	7/17/2008	7/15/2009	6/24/2010	7/19/2011	6/22/2012	6/13/2013	6/27/2014	6/26/2015	6/24/2016	7/10/2017	7/23/2018	6/14/2019	7/9/2020	7/24/1990	8/22/1991
Parameters																	
Pesticides/PCBs (μg/L)																	
4,4'-DDD																	
4,4'-DDE																	
Aldrin																	
Alpha-BHC					0.14 J						0.086			0.10	0.12		
Alpha-Chlordane										0.030							
Aroclor-1260 (PCB-1260)																	
beta&gamma-BHC (sum of isomers)																	
Beta-BHC			0.015 J		0.12 J												
Delta-BHC					0.12 J			0.067 J			0.067 J		0.032 J	0.045 J	0.088		
Dieldrin																	
Endosulfan I																	
Endosulfan II																	
Endosulfan Sulfate																	
Endrin																	
Endrin ketone																	
Gamma-BHC (Lindane)					0.12 J						0.083		0.021 J	0.080	0.150		
Gamma-Chlordane																	
Heptachlor																	
Heptachlor epoxide																	
Methoxychlor																	

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	Well Number: SampleDate:	10210B 8/26/1992	10210B 8/11/1993	10210B 6/15/1994	10210B 6/1/1995	10210B 7/5/1996	10210B 7/1/1997	10210B 6/18/1998	10210B 6/24/1999	10210B 6/15/2000	10210B 5/17/2001	10210B 6/10/2002	10210B 5/23/2003	10210B 6/2/2004	10210B 6/24/2005	10210B 6/28/2006	10210B 7/26/2007
Parameters																	
Volatile Organic Compounds (μg/L)																	
1,1,1-Trichloroethane																	
1,1,2,2-Tetrachloroethane																	
1,1,2-Trichloroethane																	
1,1-Dichloroethane																	
1,1-Dichloroethene				0.06													
1,2-Dichloroethane																	
1,2-Dichloroethene (total)																	
1,2-Dichloropropane																	
2-Butanone (Methyl Ethyl Ketone)				4									23				
2-Hexanone																	
Acetone		31	6	12 B	23						12 J						
Benzene			0.3	0.3													
Bromodichloromethane																	
Bromoform																	
Bromomethane (Methyl bromide)			0.2														
Carbon Disulfide			2	0.4				8 J	2 J		14	3 J	2 J		1.4 J	1 J	6 J
Chlorobenzene				0.2										1 J			
Chloroform																	
Chlorotoluenes																	
cis-1,2-Dichloroethene																	
Dibromochloromethane																	
Dichlorotoluene, total				0.00													
Ethylbenzene			0.2	0.08													
m&p-Xylenes			0.4	0.0													
Methylene Chloride			0.4	0.2													
o-Xylene																	
Styrene			0.00											0.1			
Tetrachloroethene			0.06	0.5					2.1	4.1				9 J	441		
Toluene trans-1,2-Dichloroethene				0.5					2 J	1 J					1.1 J		
Trichloroethene			0.1	0.1													
Trichlorotoluene, total			0.1	0.1													
Vinyl Acetate	-																
Vinyl Chloride																	
Xylenes (total)			0.5	0.5													
Aylenes (total)			0.5	0.0													
Semi-volatile Organic Compounds (µ	g/L)																
1,2,4-Trichlorobenzene													3 J				
1,2-Dichlorobenzene																	
1,3-Dichlorobenzene																	
1,4-Dichlorobenzene																	
2,4,5-Trichlorophenol																	

Summary of Detected Compounds in Select Wells Love Canal Long-Term Monitoring Program Niagara Falls, New York

	Well Number: SampleDate:	10210B 8/26/1992	10210B 8/11/1993	10210B 6/15/1994	10210B 6/1/1995	10210B 7/5/1996	10210B 7/1/1997	10210B 6/18/1998	10210B 6/24/1999	10210B 6/15/2000	10210B 5/17/2001	10210B 6/10/2002	10210B 5/23/2003	10210B 6/2/2004	10210B 6/24/2005	10210B 6/28/2006	10210B 7/26/2007
Parameters		0.20.1002		G. 10. 100 1	S. II 1000			0.10.100	0. <u>=</u> 000	0.10.200	S	S. 10.2002	0.20.200	0/2/2001	0.22000	0.20.200	
Parameters																	
Semi-volatile Organic Compounds	s (μg/L) (Continued)																
2,4,6-Trichlorophenol																	
2,4-Dichlorophenol																	
2,4-Dimethylphenol																	
2-Chloronaphthalene																	
2-Chlorophenol																	
2-Methylnaphthalene				0.06													
2-Methylphenol																	
3&4-Methylphenol																	
2-Nitrophenol																	
3,5-Dichlorotoluene																	
4-Chloro-3-methylphenol																	
4-Chlorophenyl phenyl ether																	
4-Methylphenol				0.2													
Acetic acid																	
Anthracene																	
Benzo(a)pyrene				0.07													
Benzo(b)fluoranthene				0.08													
Benzo(g,h,i)perylene				0.1													
Benzo(k)fluoranthene				0.04													
Benzoic Acid																2 J	
Benzyl Alcohol																	
Bis(2-Chloroethyl)Ether																	
bis(2-Ethylhexyl)Phthalate			11	9			55	6 J						4 J	4.5 J	3 J	
Butyl benzylphthalate (BBP)				0.2													
Camphor																	
Carbazole				0.05													
Chlorobenzoic acid																	
Chrysene																	
Dibenz(a,h)anthracene				0.1													
Diethyl phthalate				0.3													
Dimethyl Phthalate																	
Dimethyl tetrasulfide																	
Di-n-butyl phthalate (DBP)				0.6							3 J						
Di-n-octyl phthalate (DnOP)				0.1													
Fluoranthene				0.04													
Hexachlorobenzene													1 J				
Hexachloroethane																	
Indeno(1,2,3-cd)pyrene				0.1													
Naphthalene																	
N-Nitrosodiphenylamine				0.2													
Pentachlorophenol				0.3													
Phenanthrene				0.07													
Phenol				2													
Pyrene				0.04													

0.35 J

Summary of Detected Compounds in Select Wells Love Canal Long-Term Monitoring Program Niagara Falls, New York

	Well Number: SampleDate:	10210B 8/26/1992	10210B 8/11/1993	10210B 6/15/1994	10210B 6/1/1995	10210B 7/5/1996	10210B 7/1/1997	10210B 6/18/1998	10210B 6/24/1999	10210B 6/15/2000	10210B 5/17/2001	10210B 6/10/2002	10210B 5/23/2003	10210B 6/2/2004	10210B 6/24/2005	10210B 6/28/2006	10210B 7/26/2007
Parameters	·																
Pesticides/PCBs (μg/L)																	
4,4'-DDD																	
4,4'-DDE														0.011 J			
Aldrin																.0089 J	
Alpha-BHC													19	2.4	0.37	.58	0.016 J
Alpha-Chlordane																	
Aroclor-1260 (PCB-1260)																	
beta&gamma-BHC (sum of isomers)																	
Beta-BHC													1.9	0.53	0.082 P	.082	
Delta-BHC													0.56 J	0.15		.047 J	
Dieldrin													0.13 J				
Endosulfan I													0.11 J				
Endosulfan II																	
Endosulfan Sulfate																	
Endrin																	
Endrin ketone																	
Gamma-BHC (Lindane)													2.1	0.39	0.046 J	.099	
Gamma-Chlordane													0.15 J				
Heptachlor																	

Heptachlor epoxide

Methoxychlor

Table 3.5 Page 10 of 24

	Well Number: SampleDate:	10210B 7/15/2009	10210B 7/15/2009	10210B 7/9/2009	10210B 6/15/2010	10210B 7/14/2011	10210B 6/18/2012	10210B 6/15/2013	10210B 6/26/2014	10210B 7/2/2015	10210B 6/28/2016	10210B 7/10/2017	10210B 7/19/2018	10210B 6/17/2019	10210B 6/30/2020	10210C 7/25/1990
Parameters																
Volatile Organic Compounds (μg/L)																
1,1,1-Trichloroethane																
1,1,2,2-Tetrachloroethane																
1,1,2-Trichloroethane																
1,1-Dichloroethane																
1,1-Dichloroethene																
1,2-Dichloroethane																
1,2-Dichloroethene (total)		 														
1,2-Dichloropropane		 														
2-Butanone (Methyl Ethyl Ketone)																
2-Hexanone																
Acetone												2.6 J				
Benzene																
Bromodichloromethane																
Bromoform																
Bromomethane (Methyl bromide)					40.1	4.0.1	0.4.1		0.0.1	7.0		0.4.1		4.0.1	401	
Carbon Disulfide					4.0 J	4.6 J	6.1 J		2.9 J	7.9		3.4 J	20	4.2 J	4.3 J	
Chloroform																-
Chloroform Chlorotoluenes																
cis-1,2-Dichloroethene																
Dibromochloromethane																
Dichlorotoluene, total																
Ethylbenzene																
m&p-Xylenes																
Methylene Chloride								1.2 J								
o-Xylene								0								
Styrene																
Tetrachloroethene																
Toluene																3
trans-1,2-Dichloroethene																
Trichloroethene																
Trichlorotoluene, total																
Vinyl Acetate																
Vinyl Chloride																
Xylenes (total)																
Semi-volatile Organic Compounds (µ	 ıg/L)															
1,2,4-Trichlorobenzene	- ,															
1,2-Dichlorobenzene	-															
1,3-Dichlorobenzene																_
1,4-Dichlorobenzene																
2,4,5-Trichlorophenol																

Summary of Detected Compounds in Select Wells Love Canal Long-Term Monitoring Program Niagara Falls, New York

	Well Number: SampleDate:	10210B 7/17/2008	10210B 7/15/2009	10210B 7/15/2009	10210B 7/9/2009	10210B 6/15/2010	10210B 7/14/2011	10210B 6/18/2012	10210B 6/15/2013	10210B 6/26/2014	10210B 7/2/2015	10210B 6/28/2016	10210B 7/10/2017	10210B 7/19/2018	10210B 6/17/2019	10210B 6/30/2020	10210C 7/25/1990
Parameters						0.10.2010		0.10.2012	0.10.20.10	0.20.20	.,	0.20.20.10			S <u>-</u>	0.00.2020	
raidificters																	
Semi-volatile Organic Compounds	(μg/L) (Continued)																
2,4,6-Trichlorophenol																	
2,4-Dichlorophenol																	
2,4-Dimethylphenol																	
2-Chloronaphthalene																	
2-Chlorophenol																	
2-Methylnaphthalene																	
2-Methylphenol																	
3&4-Methylphenol																	
2-Nitrophenol																	
3,5-Dichlorotoluene																	
4-Chloro-3-methylphenol																	
4-Chlorophenyl phenyl ether																	
4-Methylphenol																	
Acetic acid																	
Anthracene																	
Benzo(a)pyrene																	
Benzo(b)fluoranthene																	
Benzo(g,h,i)perylene																	
Benzo(k)fluoranthene																	
Benzoic Acid																	
Benzyl Alcohol																	
Bis(2-Chloroethyl)Ether																	
bis(2-Ethylhexyl)Phthalate																	7 B
Butyl benzylphthalate (BBP)																	1
Camphor																	
Carbazole																	
Chlorobenzoic acid																	
Chrysene																	
Dibenz(a,h)anthracene																	
Diethyl phthalate								1.8 J									1
Dimethyl Phthalate																	
Dimethyl tetrasulfide																	
Di-n-butyl phthalate (DBP)																	2
Di-n-octyl phthalate (DnOP)																	
Fluoranthene																	
Hexachlorobenzene																	
Hexachloroethane																	
Indeno(1,2,3-cd)pyrene																	
Naphthalene																	
N-Nitrosodiphenylamine																	
Pentachlorophenol																	
Phenanthrene																	
Phenol																	2

Pyrene

	Well Number: SampleDate:	10210B 7/17/2008	10210B 7/15/2009	10210B 7/15/2009	10210B 7/9/2009	10210B 6/15/2010	10210B 7/14/2011	10210B 6/18/2012	10210B 6/15/2013	10210B 6/26/2014	10210B 7/2/2015	10210B 6/28/2016	10210B 7/10/2017	10210B 7/19/2018	10210B 6/17/2019	10210B 6/30/2020	10210C 7/25/1990
Parameters																	
Pesticides/PCBs (μg/L)																	
4,4'-DDD																	
4,4'-DDE																	
Aldrin																	
Alpha-BHC			0.064	0.050	0.050 / 0.064		0.048 J					0.056 J			0.054	0.12	
Alpha-Chlordane																	
Aroclor-1260 (PCB-1260)																	
beta&gamma-BHC (sum of isomers)																	
Beta-BHC																	
Delta-BHC			0.032 J	0.028 J	0.028 J / 0.032 J	0.050 J	0.042 J					0.17 J				0.076	
Dieldrin																	
Endosulfan I																	
Endosulfan II																	
Endosulfan Sulfate																	
Endrin																	
Endrin ketone																	
Gamma-BHC (Lindane)			0.038 J	0.033 J	0.038 J / 0.033 J		0.061 J					0.084			0.055	0.10	
Gamma-Chlordane																	
Heptachlor							0.053 J										
Heptachlor epoxide																	
Methoxychlor																	

	Well Number: SampleDate:	10210C 8/22/1991	10210C 8/26/1992	10210C 8/11/1993	10210C 6/8/1994	10210C 6/1/1995	10210C 7/1/1996	10210C 7/1/1997	10210C 6/22/1998	10210C 6/24/1999	10210C 6/15/2000	10210C 5/17/2001	10210C 6/10/2002	10210C 5/23/2003	10210C 6/7/2004	10210C 6/23/2005	10210C 6/28/2006
Parameters																	
Volatile Organic Compounds (μg/L)																	
1,1,1-Trichloroethane																	
1,1,2,2-Tetrachloroethane																	
1,1,2-Trichloroethane																	
1,1-Dichloroethane																	
1,1-Dichloroethene																	
1,2-Dichloroethane																	
1,2-Dichloroethene (total)																	
1,2-Dichloropropane																	
2-Butanone (Methyl Ethyl Ketone)																	
2-Hexanone																	
Acetone			10 B		19 B					2100	8 J	9 J				1.9 J	
Benzene																	
Bromodichloromethane																	
Bromoform																	
Bromomethane (Methyl bromide)																	
Carbon Disulfide					0.6						3 J						
Chlorobenzene																	2 J
Chloroform																	
Chlorotoluenes																	
cis-1,2-Dichloroethene																	
Dibromochloromethane																	
Dichlorotoluene, total																	
Ethylbenzene																	
m&p-Xylenes					0.0												
Methylene Chloride					0.2												
o-Xylene																	
Styrene																	
Tetrachloroethene Toluene															29 / 23		6 J
trans-1,2-Dichloroethene															29 / 23		
Trichloroethene																	
Trichlorotoluene, total																	
Vinyl Acetate																	
Vinyl Chloride																	
Xylenes (total)																	
Aylones (total)																	
Semi-volatile Organic Compounds (բյ	g/L)																
1,2,4-Trichlorobenzene	- ·																6 J
1,2-Dichlorobenzene																	
1,3-Dichlorobenzene																	
1,4-Dichlorobenzene																	
2,4,5-Trichlorophenol																	

10210C

10210C

Summary of Detected Compounds in Select Wells Love Canal Long-Term Monitoring Program Niagara Falls, New York

10210C

10210C

10210C

10210C

10210C

10210C

10210C

10210C

	well number:	102106	10210C	10210C	10210C	10210C	10210C	10210C	10210C	102100	102100	10210C	10210C	10210C	10210C	10210C	10210C
	SampleDate:	8/22/1991	8/26/1992	8/11/1993	6/8/1994	6/1/1995	7/1/1996	7/1/1997	6/22/1998	6/24/1999	6/15/2000	5/17/2001	6/10/2002	5/23/2003	6/7/2004	6/23/2005	6/28/2006
Parameters																	
Semi-volatile Organic Compounds	(μg/L) (Continued)																
2,4,6-Trichlorophenol																	
2,4-Dichlorophenol																	
2,4-Dimethylphenol																	
2-Chloronaphthalene																	
2-Chlorophenol																	
2-Methylnaphthalene																	
2-Methylphenol					5												
3&4-Methylphenol																	
2-Nitrophenol																	
3,5-Dichlorotoluene																	
4-Chloro-3-methylphenol																	
4-Chlorophenyl phenyl ether																	
4-Methylphenol					6	29	110	62	0.6J								
Acetic acid				11													
Anthracene																	
Benzo(a)pyrene																	
Benzo(b)fluoranthene																	
Benzo(g,h,i)perylene																	
Benzo(k)fluoranthene																	
Benzoic Acid																	
Benzyl Alcohol																	
Bis(2-Chloroethyl)Ether																	
bis(2-Ethylhexyl)Phthalate		13			8										29 / 5 J		5 J
Butyl benzylphthalate (BBP)					0.4												
Camphor																	
Carbazole																	
Chlorobenzoic acid																	
Chrysene																	
Dibenz(a,h)anthracene																	
Diethyl phthalate					0.2												
Dimethyl Phthalate																	
Dimethyl tetrasulfide																	
Di-n-butyl phthalate (DBP)		3			0.5												
Di-n-octyl phthalate (DnOP)					0.04												
Fluoranthene																	
Hexachlorobenzene																	
Hexachloroethane		1															
Indeno(1,2,3-cd)pyrene																	
Naphthalene																	
N-Nitrosodiphenylamine																	
Pentachlorophenol																	
Phenanthrene					0.03												
Phenol		6				22		22									
Pyrene																	

Well Number:

10210C

10210C

10210C

10210C

10210C

10210C

	Well Number: SampleDate:	10210C 8/22/1991	10210C 8/26/1992	10210C 8/11/1993	10210C 6/8/1994	10210C 6/1/1995	10210C 7/1/1996	10210C 7/1/1997	10210C 6/22/1998	10210C 6/24/1999	10210C 6/15/2000	10210C 5/17/2001	10210C 6/10/2002	10210C 5/23/2003	10210C 6/7/2004	10210C 6/23/2005	10210C 6/28/2006
Parameters																	
Pesticides/PCBs (μg/L)																	
4,4'-DDD																	
4,4'-DDE																	
Aldrin																	0.061 J
Alpha-BHC																0.083	0.45 J
Alpha-Chlordane																	
Aroclor-1260 (PCB-1260)																	
beta&gamma-BHC (sum of isomers)																	
Beta-BHC																	0.048 J
Delta-BHC															0.019 J / 0.017 J		0.052 J
Dieldrin																	
Endosulfan I																	
Endosulfan II																	
Endosulfan Sulfate																	
Endrin																	0.14 J
Endrin ketone																	
Gamma-BHC (Lindane)																	0.11 J
Gamma-Chlordane																	0.018 J
Heptachlor																	
Heptachlor epoxide																	
Methoxychlor																	

	Well Number: SampleDate:	10210C 7/26/2007	10210C 7/16/2008	10210C 7/13/2009	10210C 6/15/2010	10210C 7/14/2011	10210C 6/22/2012	10210C 6/15/2013	10210C 6/26/2014	10210C 7/2/2015	10210C 6/28/2016	10210C 7/10/2017	10210C 7/19/2018	10210C 6/17/2019	10210C 6/30/2020	10135 9/13/1990	10135 8/29/1991
Parameters	·																
Volatile Organic Compounds (µg/L)																	
1,1,1-Trichloroethane																	
1,1,2,2-Tetrachloroethane																	
1,1,2-Trichloroethane																	
1,1-Dichloroethane																	
1,1-Dichloroethene																6	
1,2-Dichloroethane																	
1,2-Dichloroethene (total)																	_
1,2-Dichloropropane																	_
2-Butanone (Methyl Ethyl Ketone)																	
2-Hexanone																	
Acetone												4.3 J				50	
Benzene																6200	6700
Bromodichloromethane																	
Bromoform																	
Bromomethane (Methyl bromide)																	
Carbon Disulfide			2 J				U/1.4 J					0.77 J	6.3 J	2.6 J	4.2 J		
Chlorobenzene																2380	2400
Chloroform	-																
Chlorotoluenes																16600	16000
cis-1,2-Dichloroethene																	
Dibromochloromethane																	
Dichlorotoluene, total																14000	140
Ethylbenzene																12	10
m&p-Xylenes																	
Methylene Chloride																5	
o-Xylene																	
Styrene																	
Tetrachloroethene																50	
Toluene																22800	26000
trans-1,2-Dichloroethene																20	
Trichloroethene																260	450
Trichlorotoluene, total																40	
Vinyl Acetate																	
Vinyl Chloride																	
Xylenes (total)																50	30
Semi-volatile Organic Compounds (μ	ıg/L)																
1,2,4-Trichlorobenzene																80	290
1,2-Dichlorobenzene																50	16
1,3-Dichlorobenzene																	
1,4-Dichlorobenzene																120	47
2,4,5-Trichlorophenol																860	130

	Well Number: SampleDate:	10210C 7/26/2007	10210C 7/16/2008	10210C 7/13/2009	10210C 6/15/2010	10210C 7/14/2011	10210C 6/22/2012	10210C 6/15/2013	10210C 6/26/2014	10210C 7/2/2015	10210C 6/28/2016	10210C 7/10/2017	10210C 7/19/2018	10210C 6/17/2019	10210C 6/30/2020	10135 9/13/1990	10135 8/29/1991
Parameters	·																
Semi-volatile Organic Compounds	(µg/L) (Continued)																
2,4,6-Trichlorophenol																	120
2,4-Dichlorophenol																830	
2,4-Dimethylphenol																	
2-Chloronaphthalene																	
2-Chlorophenol																	
2-Methylnaphthalene																	
2-Methylphenol																	
3&4-Methylphenol																	
2-Nitrophenol																	
3,5-Dichlorotoluene																	
4-Chloro-3-methylphenol																	13
4-Chlorophenyl phenyl ether																	
4-Methylphenol																	10
Acetic acid																	
Anthracene																	
Benzo(a)pyrene																	
Benzo(b)fluoranthene																	
Benzo(g,h,i)perylene																	
Benzo(k)fluoranthene																	
Benzoic Acid																140000	580
Benzyl Alcohol																4200	1100
Bis(2-Chloroethyl)Ether																	
bis(2-Ethylhexyl)Phthalate																	
Butyl benzylphthalate (BBP)																	
Camphor																	
Carbazole																	
Chlorobenzoic acid																4000	
Chrysene																	
Dibenz(a,h)anthracene																	
Diethyl phthalate					4.4 J												
Dimethyl Phthalate					0.87 J												
Dimethyl tetrasulfide																	
Di-n-butyl phthalate (DBP)																	
Di-n-octyl phthalate (DnOP)																	
Fluoranthene																	
Hexachlorobenzene																	
Hexachloroethane																	
Indeno(1,2,3-cd)pyrene																	
Naphthalene																	
N-Nitrosodiphenylamine																	
Pentachlorophenol																	
Phenanthrene																	
Phenol														1.7 J			10
Pyrene																	

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	Well Number: SampleDate:	10210C 7/26/2007	10210C 7/16/2008	10210C 7/13/2009	10210C 6/15/2010	10210C 7/14/2011	10210C 6/22/2012	10210C 6/15/2013	10210C 6/26/2014	10210C 7/2/2015	10210C 6/28/2016	10210C 7/10/2017	10210C 7/19/2018	10210C 6/17/2019	10210C 6/30/2020	10135 9/13/1990	10135 8/29/1991
Parameters																	
Pesticides/PCBs (μg/L)																	
4,4'-DDD	-																
4,4'-DDE																	
Aldrin																	
Alpha-BHC											0.062			0.021 J	0.081		
Alpha-Chlordane																	
Aroclor-1260 (PCB-1260)																	
beta&gamma-BHC (sum of isomers)																	
Beta-BHC																	
Delta-BHC				0.048 J							0.14 J			0.028 J	0.068 J		
Dieldrin																	
Endosulfan I																	
Endosulfan II																	
Endosulfan Sulfate																	
Endrin										0.13							
Endrin ketone																	
Gamma-BHC (Lindane)											0.13			0.026 J	0.089		
Gamma-Chlordane																	
Heptachlor																	
Heptachlor epoxide																	
Methoxychlor																	

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Summary of Detected Compounds in Select Wells Love Canal Long-Term Monitoring Program Niagara Falls, New York

Table 3.5

	Well Number: SampleDate:	10135 8/26/1992	10135 8/19/1993	10135 6/22/1994	10135 6/1/1995	10135 6/27/1996	10135 7/7/1997	10135 6/17/1998	10135 6/16/1999	10135 6/22/2000	10135 5/11/2001	10135 6/12/2002	10135 5/19/2003	10135 5/28/2004	10135 6/17/2005	10135 6/28/2005	10135 6/26/2006
Parameters																	
Volatile Organic Compounds (μg/L)																	
1,1,1-Trichloroethane			8	14													
1,1,2,2-Tetrachloroethane			12	51		26		94 J	29 / 32	27 J / 26 J	120 J / 100 J	56	38				
1,1,2-Trichloroethane						14		29 J	15 / 12	16 J / 14 J	29 J / 34 J	27					
1,1-Dichloroethane																	
1,1-Dichloroethene			15	3					4 J / 3 J	4 J / 4 J	4 J / 4 J	4 J	3 J				
1,2-Dichloroethane																	
1,2-Dichloroethene (total)		700	840	650	670 JD	560						600 J / 560	490 J				
1,2-Dichloropropane																	
2-Butanone (Methyl Ethyl Ketone)			36							10 J	11 J / 12 J						
2-Hexanone																	
Acetone			270	100 B	100 J	60		110 J		28 J / 46 J		72	74				200 J
Benzene			5200	6000 E	4900 D	4800	5000 / 5600	5300 J	5700 / 5600	6900 J / 6400 E	0 8500 J / 7600	5900 / 6400	5500		5400	5700	6800
Bromodichloromethane																	_
Bromoform																	
Bromomethane (Methyl bromide)																	
Carbon Disulfide			4=00=	2222 =			0000 / N/D		2 J				4000			0.400	
Chlorobenzene		2600	1700 E	2900 E	2000 D	1500	2300 / ND	1900 J	1900 / 1800		J 3000 J / 2700 J	2200 / 2400	1900		2000	2100	2400
Chloroform			100	120	86 J	110		150 J	110 / 120	130 J / 100 J	160 J / 150 J	160	110				110 J
Chlorotoluenes	-																
cis-1,2-Dichloroethene																	630
Dibromochloromethane Diablaretalwana total																	
Dichlorotoluene, total Ethylbenzene			13	12				12	9 J / 10 J	12 J / 12 J	24 J / 22 J	15	10				
m&p-Xylenes			13	12				12	93/103	12 3 / 12 3	24 J / 22 J	39	29				
Methylene Chloride			41	8		11				24 J / 24 J		39	26			100 J	44 J
o-Xylene			71	<u> </u>						24 3 / 24 3		12	9 J			100 3	
Styrene			4									12	9.0				
Tetrachloroethene			8	32				40 J	13 / 12	16 J / 14 J	50 J / 61 J	38	18				
Toluene		2700	17000	21500	18000 D	14000	19000 / 17000	16000 J	16000 / 17000		24000 / 22000		15000		16000	18000	21000
trans-1,2-Dichloroethene		2100	17000	21000	10000 B	14000	100007 17000	58 J	67 / 70	67 J / 70 J	59 J / 60 J	20000 j 7 10000	10000		10000	10000	52 J
Trichloroethene			24	140	18 J	36		170 J	58 / 70	60 J / 72 J	180 J / 140 J	160 / 130 J	91				46 J
Trichlorotoluene, total	·									0007.120		1007 100 0	<u> </u>				
Vinyl Acetate	-	6800															
Vinyl Chloride	-			61	44 J	50		48 J	62 / 61	110 J / 85 J	66 J / 75 J	48	51				
Xylenes (total)			47	10 B	37 J	28		55 J	44 / 43	42 J / 44 J		-	-				
Semi-volatile Organic Compounds (μ	a/L)																
1,2,4-Trichlorobenzene	y' ∟)		74	87 B				78 J	45 J / 65 J	45 J / 36 J	65 J / 42 J		97 J		67 J		63
1,2-Dichlorobenzene			35	34				101	24 J / 30 J	18 J / 22 J	48 J		97 J 59 J		36 J		37
1,3-Dichlorobenzene			JÜ	4					24 J / 30 J	10 J / ZZ J	4 0 J		29.0		30 J		3 J
1,4-Dichlorobenzene		110	94	91					61 J / 74 J	59 J / 52 J	110 J / 69 J		160 J		100 J	110 J	100
2,4,5-Trichlorophenol		110	70	91 59				38 J	010/140	0.9 J	110 0 / 09 0		100 0		100 0	1100	8 J
z,4,0- moniorophenoi			70	<u>ეგ</u>				30 J		U.8 J							0 J

Summary of Detected Compounds in Select Wells Love Canal Long-Term Monitoring Program Niagara Falls, New York

Table 3.5

	SampleDate:	8/26/1992	8/19/1993	6/22/1994	6/1/1995	6/27/1996	7/7/1997	6/17/1998	6/16/1999	6/22/2000	5/11/2001	6/12/2002	5/19/2003	5/28/2004	6/17/2005	6/28/2005	6/26/2006
Parameters																	
i diameters	_																
Semi-volatile Organic Compounds (µ	g/L) (Continued)																
2,4,6-Trichlorophenol				8						1 J							
2,4-Dichlorophenol		1200B	420	610	150		2100 / 2100	2000	690 / 610	1400 J / 470 J	620 J / 1200 J	1800 J / 1500 J	1700		420	300 J	250
2,4-Dimethylphenol				9						2 J							
2-Chloronaphthalene					150						370 J / 550 J						
2-Chlorophenol				20				28 J	25 J								18
2-Methylnaphthalene																	
2-Methylphenol			51	46				55 J	42 J / 35 J	160 J	41 J		50 J		25 J		33
3&4-Methylphenol																	
2-Nitrophenol										1 J							
3,5-Dichlorotoluene			350														
4-Chloro-3-methylphenol				31					25 J / 33 J				41 J				15
4-Chlorophenyl phenyl ether				2													
4-Methylphenol			60	64				130 J	95 J / 120	99 J / 300 J	130 J / 86 J		210 J		49 J	98 J	110
Acetic acid																	
Anthracene				1													
Benzo(a)pyrene																	
Benzo(b)fluoranthene																	
Benzo(g,h,i)perylene																	
Benzo(k)fluoranthene																	
Benzoic Acid					6400 D	4000	27000 J / 30000 J	23000 J	4300 / 5000	4700 J / 19000 J	J 6200 J / 4400 J	31000 / 25000	26000		1400 J	4700 J	14000 J
Benzyl Alcohol					380		1900 / 1600	2700	680 / 540	14000 D / 3200 s	J 330 J / 630 J	2000 / 1700 J	640		23 J		48
Bis(2-Chloroethyl)Ether			23					24 J	26 J / 25 J						24 J		24
bis(2-Ethylhexyl)Phthalate			50	2						41 J / 24 J							53
Butyl benzylphthalate (BBP)																	
Camphor			130														
Carbazole																	
Chlorobenzoic acid																	
Chrysene				0.2													
Dibenz(a,h)anthracene				<u> </u>													
Diethyl phthalate				1													
Dimethyl Phthalate																	
Dimethyl tetrasulfide																	
Di-n-butyl phthalate (DBP)																	
Di-n-octyl phthalate (DnOP)																	
Fluoranthene				0.2													
Hexachlorobenzene				0.2													
Hexachloroethane																	
Indeno(1,2,3-cd)pyrene	-																
Naphthalene									1400 .L / 2000 .	J 4000 J / 1800 J	1400 / 1100				1800 J	4500 J	
N-Nitrosodiphenylamine									00 3 / 2000 0	. 1000 0 7 1000 0	1.0071100				.0000	10000	
Pentachlorophenol			52	4													
Phenanthrene				_													
Phenol			98	91	140				120 / 96 J		51 J		180 J			100 J	140
Pyrene				J1	170				1207 30 0		510		100 0			100 0	
i yione	-																

Well Number:

	Well Number: SampleDate:	10135 8/26/1992	10135 8/19/1993	10135 6/22/1994	10135 6/1/1995	10135 6/27/1996	10135 7/7/1997	10135 6/17/1998	10135 6/16/1999	10135 6/22/2000	10135 5/11/2001	10135 6/12/2002	10135 5/19/2003	10135 5/28/2004	10135 6/17/2005	10135 6/28/2005	10135 6/26/2006
	SampleDate.	0/20/1332	0/13/1333	0/22/1994	0/1/1995	0/2//1990	11111331	0/1//1330	0/10/1333	0/22/2000	3/11/2001	0/12/2002	3/13/2003	3/20/2004	0/11/2003	0/20/2003	0/20/2000
Parameters																	
Pesticides/PCBs (μg/L)																	
4,4'-DDD									0.21 / 0.20 J	0.13 J / 0.071 J					0.19 J	0.11 J	
4,4'-DDE				0.071													
Aldrin		0.53	0.24 P						0.21 J / 0.74 JN		1.5 JN / 0.95 JN	N 0.12 J / 0.12 J					
Alpha-BHC		84	42 C	24 CEP	28 D	29	39 / 39	59	40 / 37 J	50 / 50	43 J / 50 J	43 / 39	49		15	21 C	35
Alpha-Chlordane												0.031 J / 0.017 J					
Aroclor-1260 (PCB-1260)																	
beta&gamma-BHC (sum of isomers)			19.5	20.4													
Beta-BHC					10 D	11	8.1 / 8.6	12	12 / 11 J	15 / 16	16 J / 16 J	14 J / 13 J	15 J		3.4	5.6	7.1
Delta-BHC		15	9.8	7.5 CE	4.7	5.2	5.1 / ND	8.9	11 / 9.6 J	13 / 14	10 J / 12 J	9.0 J / 11 J	12		9.1	9.1	13
Dieldrin																	
Endosulfan I									0.34 J / 0.43 J		1.5 JN / 1.6 JN	ļ					
Endosulfan II										0.52 J / 0.69 J					0.15 J		
Endosulfan Sulfate			0.43 P						0.18 / 0.17 J	0.17 J			1.3 J				
Endrin				0.15 P													
Endrin ketone																	
Gamma-BHC (Lindane)		33				2.4 J	6.2 J / 5.1 J	6.5 J	5.5 / 4.1 J	6.4 / 8.0	7.3 J / 5.0 J	7.1 J / 6.1 J	7.1			0.32 J	4.8
Gamma-Chlordane										0.18 J / 0.16 J		0.29 J / 0.35 J					.33 J
Heptachlor									0.63 / 0.68 JN				0.61 J				
Heptachlor epoxide									0.043 J / 0.058 J	0.031 J / 0.029 J		0.016 J / 0.025 J	2.2 J		0.053		
Methoxychlor																	

	Well Number: SampleDate:	10135 7/18/2007	10135 7/23/2008	10135 6/25/2009	10135 6/16/2010	10135 7/13/2011	10135 6/12/2012	10135 6/18/2013	10135 6/13/2014	10135 6/25/2015	10135 7/5/2016	10135 7/7/2017	10135 7/10/2018	10135 6/19/2019	10135 7/1/2020
Parameters															
Volatile Organic Compounds (μg/L)															
1,1,1-Trichloroethane															
1,1,2,2-Tetrachloroethane		16 J		25 / 24								3.4 J		1.8 J / 1.8 J	6.1 / 7.9
1,1,2-Trichloroethane		15 J		9.1 J / 8.7 J								8.1		3.4 J / 3.8 J	12 / 14
1,1-Dichloroethane												0.32 J		0.22 J / 5.0 U	0.20 U / 0.27 J
1,1-Dichloroethene		2 J													0.63 J / 0.78 J
1,2-Dichloroethane															
1,2-Dichloroethene (total)															
1,2-Dichloropropane															
2-Butanone (Methyl Ethyl Ketone)				5.8 J / 6.1 J								5.1 J		5.7 J / 4.4 J	12 / 6.8 J
2-Hexanone															0.20 U / 2.1 J
Acetone		53 J		42 / 37	39							45		50 / 35	79 / 71
Benzene		7100	5300	7500 / 7600	3400	2200	5900	2500	6100	6200	7100	6300	6200	4900 / 5700	6500 / 6300
Bromodichloromethane													38 J		
Bromoform															
Bromomethane (Methyl bromide)															
Carbon Disulfide		2 J										0.31 J		0.31 J / 0.33 J	
Chlorobenzene		2100	1400	2900 J / 3000 J	1300	1100	2500	730	2300	2600	2700	2600	2400	2200 / 2400	2500 / 2400
Chloroform		140 J	99 J	96 / 97	160	67	130 J			180 J	400 J	100	230 J	59 / 60	140 / 140
Chlorotoluenes															
cis-1,2-Dichloroethene			79 J	79 / 76	110	38 J							39 J	8.1 / 8.4	26 / 28
Dibromochloromethane															
Dichlorotoluene, total															
Ethylbenzene		10 J		10 / 10	13							14		10 / 11	12 / 12
m&p-Xylenes															
Methylene Chloride		32 J		25 / 24	38	16 J				150 J		23		3.5 J / 3.8 J	6.4 / 7.2
o-Xylene															
Styrene															
Tetrachloroethene		13 J		14 / 14	19	9.5 J						17		5.0 J / 5.3	20 / 23
Toluene		23000	13000	24000 / 24000	11000	3100	14000	6100	20000	20000	23000	19000	21000	17000 / 19000	22000 / 21000
trans-1,2-Dichloroethene		50 J	32 J	30 / 30	48	17 J						31	34 J	27 / 27	29 / 35
Trichloroethene		89 J	27 J	91 / 89	140	52				110 J	78 J	73	71 J	26 / 28	120 / 130
Trichlorotoluene, total															
Vinyl Acetate															
Vinyl Chloride				27 / 17	31							15		3.9 J / 5.0 U	10 / 11
Xylenes (total)		37 J		44 / 53	51							58	35 J	47 / 48	58 / 58
Semi-volatile Organic Compounds (μց	g/L)														
1,2,4-Trichlorobenzene		47 J	28	110 / 110	78 J	76 J	74 J	69	64 J		22 J			28 J / 14 J	93 / 100
1,2-Dichlorobenzene		31 J	10 J	68 / 52	57 J	45 J		45	34 J	47 J	14 J				25 / 34
1,3-Dichlorobenzene		87 J		4.1 J / 5.5 J				5.2 J							3.6 J / 4.1 J
1,4-Dichlorobenzene		84 J	24	150 J / 100 J	150 J	130 J	110 J	130	94 J	140 J	36 J			45 J / 21 J	73 / 97
2,4,5-Trichlorophenol								10							37 / 26

Table 3.5

	Well Number: SampleDate:	10135 7/18/2007	10135 7/23/2008	10135 6/25/2009	10135 6/16/2010	10135 7/13/2011	10135 6/12/2012	10135 6/18/2013	10135 6/13/2014	10135 6/25/2015	10135 7/5/2016	10135 7/7/2017	10135 7/10/2018	10135 6/19/2019	10135 7/1/2020
Parameters															
Semi-volatile Organic Compounds	(μg/L) (Continued)														
2,4,6-Trichlorophenol			6 J	28 / 23				12							
2,4-Dichlorophenol		490	150	1200 / 1100	780	590	240	360	660	790	42 J		120 J	35 J / 94 U	330 / 350
2,4-Dimethylphenol															8.9 J / 6.9 J
2-Chloronaphthalene							150 J	210							
2-Chlorophenol			17 J	31 / 26				28			12 J				27 / 31
2-Methylnaphthalene															1.4 J / 1.2 U
2-Methylphenol		34 J	140	66 J / 50 J	42 J	30 J		29	23 J	44 J					32 / 32
3&4-Methylphenol												34	80 J	31 J / 94 U	71 / 71
2-Nitrophenol															
3,5-Dichlorotoluene															
4-Chloro-3-methylphenol			26	95 / 97	31 J			23							46 / 40
4-Chlorophenyl phenyl ether															
4-Methylphenol		120 J	110	170 J / 140 J	130 J	83 J		89	53 J	93 J					
Acetic acid															
Anthracene															
Benzo(a)pyrene															
Benzo(b)fluoranthene															
Benzo(g,h,i)perylene															
Benzo(k)fluoranthene															
Benzoic Acid		14000	7600 J	39000 J / 54000 J	9500	11000	8700	16000	14000	16000	1100		6000	1300 / 1100	11000 / 10000
Benzyl Alcohol		580	38	1200 / 1300	610	450	600 J	380 J	290	250	79 J		280 J	61 J / 43 J	290 J / 290
Bis(2-Chloroethyl)Ether		30 J	16 J	29 / 28	34 J	28 J		26	19 J	27 J					20 / 24
bis(2-Ethylhexyl)Phthalate				4.4 J / 4.2 J											
Butyl benzylphthalate (BBP)															
Camphor															
Carbazole															
Chlorobenzoic acid															
Chrysene															
Dibenz(a,h)anthracene															
Diethyl phthalate															
Dimethyl Phthalate															
Dimethyl tetrasulfide															
Di-n-butyl phthalate (DBP)															
Di-n-octyl phthalate (DnOP)															
Fluoranthene															
Hexachlorobenzene															
Hexachloroethane															
Indeno(1,2,3-cd)pyrene															
Naphthalene															
N-Nitrosodiphenylamine															
Pentachlorophenol															
Phenanthrene															
Phenol		130 J	96	140 J / 160 J	100	82	89 J	92	62	87 J	11 J				40 / 44
Pyrene					- -				<u> </u>		-				
. ,															

Table 3.5 Page 24 of 24

Summary of Detected Compounds in Select Wells Love Canal Long-Term Monitoring Program Niagara Falls, New York

	Well Number: SampleDate:	10135 7/18/2007	10135 7/23/2008	10135 6/25/2009	10135 6/16/2010	10135 7/13/2011	10135 6/12/2012	10135 6/18/2013	10135 6/13/2014	10135 6/25/2015	10135 7/5/2016	10135 7/7/2017	10135 7/10/2018	10135 6/19/2019	10135 7/1/2020
Parameters	·														
Pesticides/PCBs (µg/L)															
4,4'-DDD		0.081 J	0.13 J		0.048 J		0.036 J	0.089 J							
4,4'-DDE								0.053							
Aldrin		0.073	0.052 J	0.55 J / 0.55 J	0.063 J			0.16 J	0.060 J						1.0 J / 0.85
Alpha-BHC		12	17	27 J / 32 J	4.0	21	7.1 J	20	21 J	20	25		23	25 / 26	27 / 25
Alpha-Chlordane		0.011 J													
Aroclor-1260 (PCB-1260)				12 J / 11 J											
beta&gamma-BHC (sum of isomers)															
Beta-BHC		3.7	4.4	11 J / 9.1 J	4.1	7.1	3.1	5.9	5.3 J	4.7	6.7		4.3	6.7 / 7.5	6.8 J / 6.9
Delta-BHC		4.7	6.3	11 J / 12	0.28	7.3	1.6 J	5.2	4.8 J	7.4	11		7.2	7.6 / 6.3	9.2 / 8.7
Dieldrin															
Endosulfan I															
Endosulfan II				1.6 J / 2.3				0.053 J	0.12 J						
Endosulfan Sulfate		0.34	0.37 J	1.5 J						0.035 J					
Endrin		0.034 J		1.9 / 1.3 J						1.4					
Endrin ketone									0.067 J						
Gamma-BHC (Lindane)		2.1	2	7.4 J / 6.2 J	0.92	4.1	1.4 J	3.9	4.3 J	3.2	4.9		3.3	4.7 / 4.4	6.5 / 6.2
Gamma-Chlordane		0.017 J						0.065	0.064 J	1.1 J					
Heptachlor		0.092	0.19 J				0.71	0.15 J	0.23 J						
Heptachlor epoxide		0.29	0.13 J	1.6 J / 1.7 J	0.10 J		0.089 J	0.22 J	0.23 J						
Methoxychlor							0.036 J								

- D Sample result is from a dilution
- C Sample result was confirmed
- E Sample result was greater than the highest calibration level
- N Validator qualifier-presumptive certainty, usually used when there is a large difference in dual column results
- P Lab qualifier used when there are large differences in dual column results
- J Estimated concentration
- U Not detected at the associated reporting limit
- B Detected in the blank sample
- Blank Not detected
- PCBs Polychlorinated Biphenyls

Table 3.6A Page 1 of 1

1140 Series Piezometers Water Levels - 2020 Love Canal Long-Term Monitoring Program Niagara Falls, New York

Silty Sand/Fill Medium

	Well **					
	1144 A	Tile Drain				
Date	(ft. AMSL)	(ft. AMSL)				
03/02/20	572.73	561.70				
06/02/20	571.77	561.70				
09/02/20	559.52*	561.70				
12/09/20	568.37	561.70				

Fractured Clay Medium

	Well **							
	1144 B	1143 A	1142 A	Tile Drain				
Date	(ft. AMSL)	(ft. AMSL)	(ft. AMSL)	(ft. AMSL)				
03/02/20	572.75	571.80	570.93	561.70				
06/02/20	571.85	571.44	570.76	561.70				
09/02/20	569.41	569.64	569.70	561.70				
12/09/20	568.68	568.80	569.05	561.70				

Soft Clay Medium

			Well **			
1144 C	1143 B	1143 C	1142 B	1141 A	Tile Drain	1140 A
(ft. AMSL)	(ft. AMSL)	(ft. AMSL)	(ft. AMSL)	(ft. AMSL)	(ft. AMSL)	(ft. AMSL)
572.41	571.79	570.23	568.45	566.50	561.70	564.20
571.67	571.56	570.03	568.43	566.58	561.70	564.01
569.35	569.79	569.05	568.06	566.40	561.70	564.30
568.34	568.94	568.05	567.44	565.89	561.70	564.28
	(ft. AMSL) 572.41 571.67 569.35	(ft. AMSL) (ft. AMSL) 572.41 571.79 571.67 571.56 569.35 569.79	(ft. AMSL) (ft. AMSL) (ft. AMSL) 572.41 571.79 570.23 571.67 571.56 570.03 569.35 569.79 569.05	1144 C 1143 B 1143 C 1142 B (ft. AMSL) (ft. AMSL) (ft. AMSL) (ft. AMSL) 572.41 571.79 570.23 568.45 571.67 571.56 570.03 568.43 569.35 569.79 569.05 568.06	1144 C 1143 B 1143 C 1142 B 1141 A (ft. AMSL) (ft. AMSL) (ft. AMSL) (ft. AMSL) 572.41 571.79 570.23 568.45 566.50 571.67 571.56 570.03 568.43 566.58 569.35 569.79 569.05 568.06 566.40	1144 C 1143 B 1143 C 1142 B 1141 A Tile Drain (ft. AMSL) (ft. AMSL) (ft. AMSL) (ft. AMSL) (ft. AMSL) (ft. AMSL) 572.41 571.79 570.23 568.45 566.50 561.70 571.67 571.56 570.03 568.43 566.58 561.70 569.35 569.79 569.05 568.06 566.40 561.70

Glacial Till Medium

	Well **							
	1144 D	1143 D	1142 C	1141 B	Tile Drain	1140 B		
Date	(ft. AMSL)							
03/02/20	570.71	568.92	566.88	567.03	561.70	564.81		
06/02/20	570.77	569.00	566.93	567.08	561.70	564.57		
09/02/20	569.39	568.44	566.68	566.85	561.70	564.63		
12/09/20	567.89	567.52	566.24	566.33	561.70	564.58		

- * Groundwater elevation anomalous; suspected measurement error.
- ** Wells listed from left to right in order from most distant outside of tile drain, to tile drain, then inside of tile drain.
 - Indicates wells used in Figure 3.3 Piezometer Flow Diagram to generate Fractured Clay groundwater level
 - Indicates wells used in Figure 3.3 Piezometer Flow Diagram to generate Soft Clay groundwater level
 - Indicates wells used in Figure 3.3 Piezometer Flow Diagram to generate Glacial Till groundwater level
- ft. AMSL Feet above mean sea level.

Table 3.6B Page 1 of 1

1150 Series Piezometers Water Levels - 2020 Love Canal Long-Term Monitoring Program Niagara Falls, New York

Silty Sand/Fill Medium

	Well **				
	1151 D	Tile Drain			
Date	(ft. AMSL)	(ft. AMSL)			
03/02/20	571.52	561.85			
06/02/20	569.96	561.85			
09/02/20	DRY	561.85			
12/09/20	DRY	561.85			

Fractured Clay Medium

	Well **						
	1154 D	1153 E	1153 D	1151 C	Tile Drain		
Date	(ft. AMSL)						
03/02/20	567.98	569.68	571.49	569.31	561.85		
06/02/20	568.31	569.54	570.50	569.23	561.85		
09/02/20	567.92	568.94	569.86	567.57	561.85		
12/09/20	567.78	569.06	570.32	565.66	561.85		

Soft Clay Medium

	Well **								
	1154 B	1154 C	1153 B	1153 C	1151 B	Tile Drain			
Date	(ft. AMSL)								
03/02/20	568.26	568.52	570.06	577.27	568.01	561.85			
06/02/20	567.96	568.36	568.95	570.24	568.08	561.85			
09/02/20	567.66	567.91	569.11	570.13	567.34	561.85			
12/09/20	567.78	568.00	569.27	574.59	567.01	561.85			

Glacial Till Medium

	Well **							
	1154 A	1153 A	1151 A	Tile Drain				
Date	(ft. AMSL)	(ft. AMSL)	(ft. AMSL)	(ft. AMSL)				
03/02/20	571.03	573.62	567.45	561.85				
06/02/20	571.05	569.66	566.80	561.85				
09/02/20	570.89	568.72	567.34	561.85				
12/09/20	570.73	571.58	567.09	561.85				

Notes:

- Wells listed from left to right in order from most distant outside of tile drain, to tile drain, then inside of tile drain.

- Indicates wells used in Figure 3.4 Piezometer Flow Diagram to generate Fractured Clay groundwater level

- Indicates wells used in Figure 3.4 Piezometer Flow Diagram to generate Soft Clay groundwater level

- Indicates wells used in Figure 3.4 Piezometer Flow Diagram to generate Glacial Till groundwater level

ft. AMSL - Feet above mean sea level.

1160 Series Piezometers Water Levels - 2020 Love Canal Long-Term Monitoring Program Niagara Falls, New York

Silty Sand/Fill Medium

	Well **				
	1165 D	Tile Drain			
Date	(ft. AMSL)	(ft. AMSL)			
03/02/20	573.51	560.60			
06/02/20	572.46	560.60			
09/02/20	571.53	560.60			
12/09/20	571.11	560.60			

Silty Sand Medium

		Well **						
	1165 C	1163 D	Tile Drain					
Date	(ft. AMSL)	(ft. AMSL)	(ft. AMSL)					
03/02/20	572.40	DRY	560.60					
06/02/20	572.49	DRY	560.60					
09/02/20	572.27	DRY	560.60					
12/09/20	571.96	DRY	560.60					

Fractured Clay Medium

					Well **				
	1165 B	1163 B	1163 C	1162 A	1162 C	1161 D	Tile Drain	1160 A	1160 C
Date	(ft. AMSL)								
03/02/20	572.17	569.87	570.14	570.23	569.99	570.01	560.60	565.38	566.42
06/02/20	571.35	569.95	570.39	569.83	569.98	570.22	560.60	564.92	DRY
09/02/20	570.75	569.78	570.11	569.51	569.74	570.18	560.60	564.94	566.67
12/09/20	570.45	569.27	569.37	569.12	569.39	569.53	560.60	565.19	566.42

Soft Clay Medium

					Well **				
	10176 A	10176 B	10176 C	1165 A	1163 A	1161 B	1161 C	1161 E	Tile Drain
Date	(ft. AMSL)								
03/02/20	571.11	571.09	569.35	572.61	568.95	567.16	569.25	565.59	560.60
06/02/20	569.35	569.40	568.86	572.99	568.66	566.84	569.30	565.35	560.60
09/02/20	568.38	568.70	568.42	572.26	568.69	566.96	569.21	565.42	560.60
12/09/20	568.16	568.69	567.78	571.68	568.69	566.91	568.84	565.50	560.60

Glacial Till Medium

Well **					
10176 D	1162 D	1161 A	Tile Drain		
(ft. AMSL)	(ft. AMSL)	(ft. AMSL)	(ft. AMSL)		
568.15	567.94	565.63	560.60		
567.76	567.63	565.31	560.60		
567.59	567.53	565.24	560.60		
567.01	567.69	565.27	560.60		
	(ft. AMSL) 568.15 567.76 567.59	10176 D 1162 D (ft. AMSL) (ft. AMSL) 568.15 567.94 567.76 567.63 567.59 567.53	10176 D 1162 D 1161 A (ft. AMSL) (ft. AMSL) (ft. AMSL) 568.15 567.94 565.63 567.76 567.63 565.31 567.59 567.53 565.24		

Notes:

- Wells listed from left to right in order from most distant outside of tile drain, to tile drain, then inside of tile drain.

- Indicates wells used in Figure 3.5 Piezometer Flow Diagram to generate Fractured Clay groundwater level

- Indicates wells used in Figure 3.5 Piezometer Flow Diagram to generate Soft Clay groundwater level

- Indicates wells used in Figure 3.5 Piezometer Flow Diagram to generate Glacial Till groundwater level

ft. AMSL

- Feet above mean sea level.

Table 3.6D Page 1 of 1

1170 Series Piezometers Water Levels - 2020 Love Canal Long-Term Monitoring Program Niagara Falls, New York

	Well **					
	1174 D	1173 D	Tile Drain			
Date	(ft. AMSL)	(ft. AMSL)	(ft. AMSL)			
03/02/20	568.87	572.77	555.60			
06/02/20	568.61	571.96	555.60			
09/02/20	568.36	570.81	555.60			
12/09/20	568.78	570.50	555.60			

Soft Clay Medium

			Well **		
	1174 B	1174 C	1173 B	1173 C	1172 B
Date	(ft. AMSL)				
03/02/20	570.32	569.99	570.25	572.07	569.26
06/02/20	570.45	569.90	569.99	572.12	569.37
09/02/20	570.79	570.29	569.75	571.15	568.42
12/09/20	570.67	569.92	569.69	570.35	568.17
			_		
	1172 C	1171 B	1171 C	Tile Drain	1170 B
	(ft. AMSL)				
03/02/20	569.74	562.88	561.86	555.60	572.80
06/02/20	569.75	562.82	561.66	555.60	570.01
09/02/20	569.24	562.70	561.55	555.60	561.69
12/09/20	568.70	562.62	561.65	555.60	563.66

Glacial Till Medium

		Well **					
	1174 A	1173 A	1172 A	1171 A	Tile Drain	1170 A	
Date	(ft. AMSL)						
03/02/20	567.94	568.74	566.76	564.18	555.60	562.70	
06/02/20	570.45	568.47	566.78	564.14	555.60	562.48	
09/02/20	570.45	569.08	566.60	563.85	555.60	562.21	
12/09/20	570.55	568.97	566.48	563.61	555.60	562.35	

- ** Wells listed from left to right in order from most distant outside of tile drain, to tile drain, then inside of tile drain.
 - Indicates wells used in Figure 3.6 Piezometer Flow Diagram to generate Fractured Clay groundwater level
 - Indicates wells used in Figure 3.6 Piezometer Flow Diagram to generate Soft Clay groundwater level
 - Indicates wells used in Figure 3.6 Piezometer Flow Diagram to generate Glacial Till groundwater level
- ft. AMSL Feet above mean sea level.

1180 Series Piezometers Water Levels - 2020 Love Canal Long-Term Monitoring Program Niagara Falls, New York

Silty Sand Medium

	Well **				
	1183 D	Tile Drain			
Date	(ft. AMSL)	(ft. AMSL)			
03/02/20	566.78	560.00			
06/02/20	566.77	560.00			
09/02/20	DRY	560.00			
12/09/20	566.77	560.00			

Fractured Clay Medium

_			Well **		
	1184 C	1184 D	1183 C	1183 B	1181 C
Date	(ft. AMSL)				
03/02/20	568.08	568.40	567.18	565.71	569.71
06/02/20	568.82	567.92	567.09	565.43	567.96
09/02/20	562.70	DRY	565.57	565.01	566.85
12/09/20	DRY	DRY	565.40	567.18	568.75
	Tile Drain	1180 C			
	(ft. AMSL)	(ft. AMSL)			
03/02/20	560.00	DRY			
06/02/20	560.00	DRY			
09/02/20	560.00	DRY			
12/09/20	560.00	DRY			

Soft Clay Medium

	Well **					
	1184 B	1181 B	Tile Drain	1180 B		
Date	(ft. AMSL)	(ft. AMSL)	(ft. AMSL)	(ft. AMSL)		
03/02/20	565.21	567.17	560.00	561.15		
06/02/20	564.90	566.40	560.00	561.00		
09/02/20	563.67	565.78	560.00	561.06		
12/09/20	563.19	566.13	560.00	561.05		

Glacial Till Medium

_			Well **		
	1184 A	1183 A	1181 A	Tile Drain	1180 A
Date	(ft. AMSL)				
03/02/20	564.85	564.45	569.03	560.00	563.63
06/02/20	565.00	564.38	567.50	560.00	563.56
09/02/20	564.11	563.96	566.24	560.00	563.28
12/09/20	563.60	563.17	568.54	560.00	562.97

- * Wells listed from left to right in order from most distant outside of tile drain, to tile drain, then inside of tile drain.
 - Indicates wells used in Figure 3.7 Piezometer Flow Diagram to generate Fractured Clay groundwater level
 - Indicates wells used in Figure 3.7 Piezometer Flow Diagram to generate Soft Clay groundwater level
 - Indicates wells used in Figure 3.7 Piezometer Flow Diagram to generate Glacial Till groundwater level
- ft. AMSL Feet above mean sea level.

1190 Series Piezometers Water Levels - 2020 Love Canal Long-Term Monitoring Program Niagara Falls, New York

Fractured Clay Medium

	Well **					
	1194 D	1193 D	1192 C	Tile Drain		
Date	(ft. AMSL)	(ft. AMSL)	(ft. AMSL)	(ft. AMSL)		
03/02/20	574.54	572.17	569.97	554.80		
06/02/20	573.42	572.11	570.31	554.80		
09/02/20	569.41	571.11	570.04	554.80		
12/09/20	568.67	570.21	569.60	554.80		

Soft Clay Medium

	Well **						
	1194 B	1194 C	1193 B	1193 C	1192 B	1191 B	
Date	(ft. AMSL)						
03/02/20	570.74	573.42	569.46	571.28	568.87	566.01	
06/02/20	570.19	572.63	569.70	571.59	569.19	566.90	
09/02/20	568.13	569.21	568.94	570.94	568.86	566.38	
12/09/20	567.50	568.58	568.44	570.18	568.44	565.32	
	1191 C	Tile Drain	1190 B				
	(ft. AMSL)	(ft. AMSL)	(ft. AMSL)				
03/02/20	564.56	554.80	571.42				
06/02/20	564.44	554.80	563.27				
09/02/20	564.09	554.80	562.26				
12/09/20	563.96	554.80	562.74				

Glacial Till Medium

_	Well **					
	1194 A	1193 A	1192 A	1191 A	Tile Drain	1190 A
Date	(ft. AMSL)					
03/02/20	565.11	566.33	565.05	565.75	554.80	566.19
06/02/20	564.99	566.21	565.01	565.97	554.80	565.40
09/02/20	564.26	565.57	564.30	565.89	554.80	564.35
12/09/20	563.88	565.26	563.97	565.42	554.80	564.12

- Wells listed from left to right in order from most distant outside of tile drain, to tile drain, then inside of tile drain.
 - Indicates wells used in Figure 3.8 Piezometer Flow Diagram to generate Fractured Clay groundwater level
 - Indicates wells used in Figure 3.8 Piezometer Flow Diagram to generate Soft Clay groundwater level
 - Indicates wells used in Figure 3.8 Piezometer Flow Diagram to generate Glacial Till groundwater level
- ft. AMSL Feet above mean sea level.

Table 3.7

Groundwater Elevations in the Vicinty of Well 10135 Love Canal Long-Term Monitoring Program Niagara Falls, New York

			Well Location	1	
	Well 1165A	Well 10135	Well 1163A	Well 1161E	Barrier Drain *
Date	(ft. AMSL)	(ft. AMSL)	(ft. AMSL)	(ft. AMSL)	(ft. AMSL)
March 2020	572.61	571.60	568.95	565.59	560.60
June 2020	572.99	571.31	568.66	565.35	560.60
Sept 2020	572.26	570.64	568.69	565.42	560.60
Dec 2020	571.68	569.14	568.69	565.50	560.60

Notes:

* Barrier Drain Elevation not measured; elevation taken from design data specific to 1160 piezometer location.



Appendix A Institutional and Engineering Controls Certification Form

NEW YORK STATE DEPARTMENT OF ENVIRONMENTAL CONSERVATION

Division of Environmental Remediation 625 Broadway, 11th Floor, Albany, NY 12233-7020 P: (518)402-9543 | F: (518)402-9547 www.dec.ny.gov

11/20/2020

Joseph Branch Project Coordinator OCC/Glenn Springs Holdings, Inc. 7601 Old Channel Trail Montaque, MI 49437

RE: Reminder Notice: Site Management Periodic Review Report and IC/EC Certification Submittal

Site Name: Love Canal

Site No.: 932020

Site Address: 805 97th Street

Dear Mr. Branch:

This letter serves as a reminder that sites in active Site Management (SM) require the submittal of a periodic progress report. This report, referred to as the Periodic Review Report (PRR), must document the implementation of and compliance with, site specific SM requirements. Section 6.3(b) of DER-10 Technical Guidance for Site Investigation and Remediation (available online at http://www.dec.nv.gov/regulations/67386.html) provides guidance regarding the information that must be included in the PRR. Further, if the site is comprised of multiple parcels, then you as the Certifying Party must arrange to submit one PRR for all parcels that comprise the site. The PRR must be received by the Department no later than January 30, 2021. Guidance on the content of a PRR is enclosed.

Site Management is defined in regulation (6 NYCRR 375-1.2(at)) and in Chapter 6 of DER-10. Depending on when the remedial program for your site was completed, SM may be governed by multiple documents (e.g., Operation, Maintenance, and Monitoring Plan; Soil Management Plan) or one comprehensive Site Management Plan.

A Site Management Plan (SMP) may contain one or all of the following elements, as applicable to the site: a plan to maintain institutional controls and/or engineering controls ("IC/EC Plan"); a plan for monitoring the performance and effectiveness of the selected remedy ("Monitoring Plan"); and/or a plan for the operation and maintenance of the selected remedy ("O&M Plan"). Additionally, the technical requirements for SM are stated in the decision document (e.g., Record of Decision) and, in some cases, the legal agreement directing the remediation of the site (e.g., order on consent, voluntary agreement, etc.).

When you submit the PRR (by the due date above), include the enclosed forms documenting that all SM requirements are being met. The Institutional Controls (ICs) portion of the form (Box 6) must be signed by you or your designated representative. The Engineering Controls (ECs) portion of the form (Box 7) must be signed by a Qualified Environmental Professional (QEP). If you cannot certify that all SM requirements are being met, you must submit a Corrective Measures Work Plan that identifies the actions to be taken to restore compliance. The work plan must include a schedule to be approved by the Department. The Periodic Review process will not be considered complete until all necessary corrective measures are completed and all required controls are certified. Instructions for completing the certifications are enclosed.



All site-related documents and data, including the PRR, are to be submitted in electronic format to the Department of Environmental Conservation. The Department will not approve the PRR unless all documents and data generated in support of that report have been submitted in accordance with the electronic submissions protocol. In addition, the certification forms are required to be submitted in both paper and electronic formats.

Information on the format of the data submissions can be found at: http://www.dec.ny.gov/regulations/2586.html

The signed certification forms should be sent to Brian Sadowski, Project Manager, at the following address:

New York State Department of Environmental Conservation 270 Michigan Avenue Buffalo, NY 14203-2915

Phone number: 716-851-7220 E-mail: brian.sadowski@dec.ny.gov

The contact information above is also provided so that you may notify the project manager about upcoming inspections, or for any other questions or concerns that may arise in regard to the site.

Enclosures

PRR General Guidance Certification Form Instructions Certification Forms

ec: w/enclosures Brian Sadowski, Project Manager Stanley Radon, Hazardous Waste Remediation Geologist, Region 9 John Pentilchuk, GHD Group (john.pentilchuk@ghd.com)

Enclosure 1 Certification

Instructions

I. Verification of Site Details (Box 1 and Box 2):

Answer the three questions in the Verification of Site Details Section. The Owner and/or Qualified Environmental Professional (QEP) may include handwritten changes and/or other supporting documentation, as necessary.

II. Certification of Institutional / Engineering Controls (Boxes 3, 4, and 5)

Review the listed IC/ECs, confirming that all existing controls are listed, and that all existing controls are still applicable. If there is a control that is no longer applicable the Owner / Remedial Party should petition the Department separately to request approval to remove the control.

In Box 5, complete certifications for all Plan components, as applicable, by checking the corresponding checkbox.

If you cannot certify "YES" for each Control listed in Box 3 & Box 4, sign and date the form in Box 5. Attach supporting documentation that explains why the **Certification** cannot be rendered, as well as a plan of proposed corrective measures, and an associated schedule for completing the corrective measures. Note that this **Certification** form must be submitted even if an IC or EC cannot be certified; however, the certification process will not be considered complete until corrective action is completed.

If the Department concurs with the explanation, the proposed corrective measures, and the proposed schedule, a letter authorizing the implementation of those corrective measures will be issued by the Department's Project Manager. Once the corrective measures are complete, a new Periodic Review Report (with IC/EC Certification) must be submitted within 45 days to the Department. If the Department has any questions or concerns regarding the PRR and/or completion of the IC/EC Certification, the Project Manager will contact you.

III. IC/EC Certification by Signature (Box 6 and Box 7):

If you certified "YES" for each Control, please complete and sign the IC/EC Certifications page as follows:

- Where the only control is an Institutional Control on the use of the property, the certification statement in Box 6 shall be completed and may be made by the property owner.
- Where the site has Institutional and Engineering Controls, the certification statement in Box 7 must be completed by a Professional Engineer or Qualified Environmental Professional, as noted on the form.



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



	Site D	Petails		Box 1
ite N	o. 932020			
ite Na	ame Love Canal			
ity/To	ddress: 805 97th Street own: Niagara Falls v: Niagara creage: 70.0	Zip Code: 14304		
eport	ing Period: January 1, 202	20 to December 31, 2020		
			YES	NO
1.	Is the information above	correct?	d	
	If NO, include handwritte	en above or on a separate sheet.		
2.	Has some or all of the si or undergone a tax map	ite property been sold, subdivided, merged, amendment during this Reporting Period?		Ø
3.	Has there been any cha Period (see 6NYCRR 37	nge of use at the site during this Reporting 75-1.11(d))?		Ø
4.	Have any federal, state, been issued for or at the	and/or local permits (e.g., building, discharge) e property during this Reporting Period?		
	If you answered YES to documentation has be	o questions 2 thru 4, include documentation or en previously submitted with this certification	evidence the	nat
5.	Is the site currently unde	ergoing development?	0	S
				Box 2
			YES	NO
6.	Is the current site use co	ensistent with the use(s) listed below?	Ø	
7.	Are all ICs/ECs in place	and functioning as designed?		
	IF THE ANS	WER TO EITHER QUESTION 6 OR 7 IS NO, sign a OT COMPLETE THE REST OF THIS FORM. Otherw	and date belo vise continue	ow and
	A Corrective Measures \	Nork Plan must be submitted along with this form	n to address	these issues.
		edial Party or Designated Representative		

SITE NO. 932020			
	n of Engineering and Institutional Controls	Boxes 3 and 4	
Parcel	Engineering Control	Institutional Control	
232 Parcels	Cover System Fencing/Access Control Groundwater Containment Leachate Collection Pump & Treat	Building Use Restriction Ground Water Use Restriction Landuse Restriction Monitoring Plan O&M Plan	
161.19-1-1			
161.57-1-1			
161.65-1-1			
161.73-1-1			
161.57-1-2			
161.65-1-2			
161.73-1-2			
161.57-1-3			
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161.73-1-3			
161.57-1-4			
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161.57-1-8			
61.65-1-8 61.73-1-8			
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TEACHER ASSET	

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	Periodic Review Report (PRR) Certification Statements Box	5	
1.	I certify by checking "YES" below that:		
	 a) the Periodic Review report and all attachments were prepared under the direction of, are reviewed by, the party making the certification; 	nd	
	 to the best of my knowledge and belief, the work and conclusions described in this certiare in accordance with the requirements of the site remedial program, and generally accept engineering practices; and the information presented is accurate and complete. 	fication ed YES	NO
2.	If this site has an IC/EC Plan (or equivalent as required in the Decision Document), for each Insor Engineering control listed in Boxes 3 and/or 4, I certify by checking "YES" below that all of the following statements are true:	titutional e	
(a) Co	the Institutional Control and/or Engineering Control(s) employed at this site is unchanged since ntrol was put in-place, or was last approved by the Department;	the date	that the
(b) the	nothing has occurred that would impair the ability of such Control, to protect public health and environment;		
(c) eva	access to the site will continue to be provided to the Department, to evaluate the remedy, includate the continued maintenance of this Control;	ling acces	ss to
(d) Co	nothing has occurred that would constitute a violation or failure to comply with the Site Manager ntrol; and	ment Plan	for this
(e) and	if a financial assurance mechanism is required by the oversight document for the site, the mech d sufficient for its intended purpose established in the document.	anism rer	mains vali
		YES	NO
		1	
	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 the	se issues	à.
	Signature of Owner, Remedial Party or Designated Representative Date		
	Date		

IC CERTIFICATIONS SITE NO. 932020

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.

JOSEPH BEANCH print name	at 7601 OLD CHANNEL TRAIL MONTAGUE, print business address	MI 49437
am certifying as OVINER	(Owner or Remedial Party)	
for the Site named in the Site Details Se	clion of this form.	
Signature of Owner or Remedial Party R	endering Certification 1/25/2021	

IC/EC CERTIFICATIONS

Box 7

Qualified Environmental Professional 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.

PILHARD J. SNYDER	at 2055 NEAGARA FALLS BLVD. NEAGARA print business address	FAHS, N 14301
am certifying as a Qualified Environment	ntal Professional for the Remedial PARTY (Owner or Remedial Party)	
	STATE OF MEN TORK	
0 111		
Signature of Qualified Environmental F the Owner or Remedial Party, Renderi	Professional, for Asiamp Date	127,202/

Enclosure 3 Periodic Review Report (PRR) General Guidance

- 1. Executive Summary: (1/2-page or less)
 - Provide a brief summary of site, nature and extent of contamination, and remedial history.
 - B. Effectiveness of the Remedial Program Provide overall conclusions regarding;
 - 1. progress made during the reporting period toward meeting the remedial objectives for the site
 - 2. the ultimate ability of the remedial program to achieve the remedial objectives for the site.
 - C. Compliance
 - Identify any areas of non-compliance regarding the major elements of the Site Management Plan (SMP, i.e., the Institutional/Engineering Control (IC/EC) Plan, the Monitoring Plan, and the Operation & Maintenance (O&M) Plan).
 - 2. Propose steps to be taken and a schedule to correct any areas of non-compliance.
 - Recommendations
 - 1. recommend whether any changes to the SMP are needed
 - 2. recommend any changes to the frequency for submittal of PRRs (increase, decrease)
 - 3. recommend whether the requirements for discontinuing site management have been met.
- II. Site Overview (one page or less)
 - A. Describe the site location, boundaries (figure), significant features, surrounding area, and the nature and extent of contamination prior to site remediation.
 - B. Describe the chronology of the main features of the remedial program for the site, the components of the selected remedy, cleanup goals, site closure criteria, and any significant changes to the selected remedy that have been made since remedy selection.
- III. Evaluate Remedy Performance, Effectiveness, and Protectiveness

Using tables, graphs, charts and bulleted text to the extent practicable, describe the effectiveness of the remedy in achieving the remedial goals for the site. Base findings, recommendations, and conclusions objective data. Evaluations and should be presented simply and concisely.

IV. IC/EC Plan Compliance Report (if applicable)

- A. IC/EC Requirements and Compliance
 - 1. Describe each control, its objective, and how performance of the control is evaluated.
 - 2. Summarize the status of each goal (whether it is fully in place and its effectiveness).
 - 3. Corrective Measures: describe steps proposed to address any deficiencies in ICECs.
 - 4. Conclusions and recommendations for changes.
- B. IC/EC Certification
 - 1. The certification must be complete (even if there are IC/EC deficiencies), and certified by the appropriate party as set forth in a Department-approved certification form(s).
- V. Monitoring Plan Compliance Report (if applicable)
 - A. Components of the Monitoring Plan (tabular presentations preferred) Describe the requirements of the monitoring plan by media (i.e., soil, groundwater, sediment, etc.) and by any remedial technologies being used at the site.
 - B. Summary of Monitoring Completed During Reporting Period Describe the monitoring tasks actually completed during this PRR reporting period. Tables and/or figures should be used to show all data.
 - C. Comparisons with Remedial Objectives Compare the results of all monitoring with the remedial objectives for the site. Include trend analyses where possible.
 - D. Monitoring Deficiencies Describe any ways in which monitoring did not fully comply with the monitoring plan.
 - E. Conclusions and Recommendations for Changes Provide overall conclusions regarding the monitoring completed and the resulting evaluations regarding remedial effectiveness.
- VI. Operation & Maintenance (O&M) Plan Compliance Report (if applicable)
 - A. Components of O&M Plan Describe the requirements of the O&M plan including required activities, frequencies, recordkeeping, etc.
 - B. Summary of O&M Completed During Reporting Period Describe the O&M tasks actually completed during this PRR reporting period.
 - C. Evaluation of Remedial Systems Based upon the results of the O&M activities completed, evaluated the ability of each component of the remedy subject to O&M requirements to perform as

designed/expected.

- D. O&M Deficiencies Identify any deficiencies in complying with the O&M plan during this PRR reporting period.
- E. Conclusions and Recommendations for Improvements Provide an overall conclusion regarding O&M for the site and identify any suggested improvements requiring changes in the O&M Plan.

VII. Overall PRR Conclusions and Recommendations

- A. Compliance with SMP For each component of the SMP (i.e., IC/EC, monitoring, O&M), summarize;
 - 1. whether all requirements of each plan were met during the reporting period

2. any requirements not met

3. proposed plans and a schedule for coming into full compliance.

B. Performance and Effectiveness of the Remedy - Based upon your evaluation of the components of the SMP, form conclusions about the performance of each component and the ability of the remedy to achieve the remedial objectives for the site.

C. Future PRR Submittals

- 1. Recommend, with supporting justification, whether the frequency of the submittal of PRRs should be changed (either increased or decreased).
- If the requirements for site closure have been achieved, contact the Departments Project Manager for the site to determine what, if any, additional documentation is needed to support a decision to discontinue site management.

VIII. Additional Guidance

Additional guidance regarding the preparation and submittal of an acceptable PRR can be obtained from the Departments Project Manager for the site.

Appendix B Semiannual Inspection Forms



Love Canal Semiannual Barrier System / Pump Chamber Inspections

Date:	5/20/2020		Weather:	65 °F	
Inspector:	Darrell Crockett		- -		
Check the	Following as Appr	opriate:			
Verification Inspection	ection of chamber n of level probe pe of pump chamber of pump chamber	rformance integrity			
	Wells	Satisfactory	Needs Maintenance		
	PC-1 PC-2 PC-3 PC-1A PC-2A PC-3A	Y Y Y Y Y			
Comments	:				

Signature: <u>Darrell Crockett</u>



Love Canal Semiannual Barrier System / Pump Chamber Inspections

Date:	11/10/2020		We	eather:	50 °F	
Inspector:	Darrell Crockett		-			
Check the F	Following as Appr	opriate:				
Verification Inspection	ection of chamber of level probe pe of pump chamber of pump chamber	erformance integrity				
	Wells	Satisfactory	Needs Maintenance			
	PC-1 PC-2 PC-3 PC-1A PC-2A PC-3A	Y Y Y Y Y				
Comments	:					
Signature:	Darrell Crod	<u>ckett</u>				

SEMIANNUAL LANDFILL CAP, SITE COVER, AND FENCE INSPECTION

Site:	Love Canal				
Date:	5/21/2020	Weather:	Sunny 50 °F		
Inspector:	Darrell Crocket	Darrell Crockett			
Inspection It	Applicable tem to Site	Inspect For			
. Landfill Cap	Y	- signs of erosion (cap, ditches, swales)	N		
		- exposure of the HDPE Liner	N		
		- areas of insufficient grass coverage	N		
		signs of dead/dying grass	N		
		- presence of washouts	N		
		- settlement causing ponding of water	Υ		
		- signs of slope instability	Υ		
		- signs of burrowing by animals	N		
		 presence of rooting trees (cap, ditches, swales) 	N		
		- signs of poor drainage in ditches/swales	N		
Site Cover	N	- signs of erosion (cover, ditches, swales)	Y/N		
(Asphalt, Gra	ass, Vegetation)	- areas of insufficient asphalt, grass, vegetation coverage	Y / N		
		 signs of dead/dying grass/vegetation 	Y / N		
		- presence of washouts	Y/N		
		 settlement causing ponding of water 	Y/N		
		- signs of slope instability	Y/N		
		- signs of burrowing by animals	Y/N		
		- presence of rooting trees (cover, ditches, swales)	Y/N		
		- signs of poor drainage in ditches/swales	Y/N		
Perimeter Fo	<u>ence</u> Y	- breaches in fence	N		
		- gates secure	Υ		
		- locks in place	Υ		
		- missing or illegible signage	N		
Comments/	Remarks (I	Note: If repair/maintenance is recommended, describe its location	on/extent below)		
		the road leading to PC1 and south of the road generally consistent with I	-		

SEMIANNUAL LANDFILL CAP, SITE COVER, AND FENCE INSPECTION

Sit	te:	Love Canal		
Da	ite:	11/11/2020	Weather:	Sunny 45 °F
Inspector:		Darrell Crockett		
lns	spection Item	Applicable to Site	Inspect For	
La	ndfill Cap	Υ	- signs of erosion (cap, ditches, swales)	N
	<u></u>	•	- exposure of the HDPE Liner	N
			- areas of insufficient grass coverage	N
			- signs of dead/dying grass	N
			- presence of washouts	N
			 settlement causing ponding of water 	N
			- signs of slope instability	N
			- signs of burrowing by animals	N
			- presence of rooting trees (cap, ditches, swales)	N
			- signs of poor drainage in ditches/swales	N
Sit	te Cover	N	- signs of erosion (cover, ditches, swales)	Y/N
(A:	sphalt, Grass, V	egetation)	- areas of insufficient asphalt, grass, vegetation coverage	Y / N
			 signs of dead/dying grass/vegetation 	Y / N
			- presence of washouts	Y / N
			- settlement causing ponding of water	Y / N
			- signs of slope instability	Y / N
			- signs of burrowing by animals	Y / N
			- presence of rooting trees (cover, ditches, swales)	Y / N
			- signs of poor drainage in ditches/swales	Y/N
<u>Pe</u>	erimeter Fence	Υ	- breaches in fence	N
			- gates secure	Υ
			- locks in place	Υ
			- missing or illegible signage	N
Co	omments/Rema	ırks (Not	e: If repair/maintenance is recommended, describe its location	on/extent below)
			ement and wet area on the road leading to PC1 and south of the road	observed
du	ring the first sem	ii-annual inspection.		
	mg the matach	aaaspeed.o		

Love Canal Semi-Annual Barrier Drain Manhole Inspection

							Date	5/20/2020
Sector	MH No.	Location	Water Y/N	Level Feet	Debris Y/N	Structure OK	Cleaning Y/N	Comments
North Colvin	MH-10A	NW	N	3"	N	Υ	N	
	MH-8A	NW	Υ	3"	N	Υ	N	
	MH-6C	NW	Υ	3"	Υ	Υ	N	Some built up sludge - no flow restriction
	MH-6B	NW	Υ	3"	Υ	Υ	N	Some built up sludge - no flow restriction
	MH-6A	NW	Υ	3"	Υ	Υ	N	Some built up sludge - no flow restriction
	PC-2A	NW	Υ	3'	N	Υ	N	
	MH-4A	NW	Υ	2"	N	Υ	N	
	MH-2A	NW	Υ	2"	N	Υ	N	
	MH-2	SW	Υ	3"	Υ	Υ	N	
	MH-4	SW	Υ	3"	Υ	Υ	N	
	MH-6	SW	Υ	3"	N	Υ	N	
	MH-8/PC2	SW	Υ	2.8'	N	Υ	N	
	MH-10	SW	Υ	2"	Υ	Υ	N	
	MH-12	SW	Υ	2"	N	Υ	N	
South Frontier	MH-14	SW	Υ	2"	N	Υ	N	
North Colvin	NH-17A	NE	N	3"	N	Υ	N	
	MH-15A	NE	Υ	3"	N	Υ	N	
	MH-13A	NE	Υ	3"	Υ	Υ	N	
	PC1A	NE	Υ	3.0'	Υ	Υ	N	
	MH-11A	NE	Υ	3"	N	Υ	N	
	MH-9A	NE	Υ	3"	N	Υ	N	
	MH-7A	NE	Υ	3"	Υ	Υ	N	
	MH-5A	NE	N	3"	Ν	Υ	N	
	MH-3A	NE	Υ	3"	N	Υ	N	
	MH-1A	NE	Υ	3"	N	Υ	N	
	MH-1	SE	N	3"	N	Υ	N	
	MH-3	SE	Υ	3"	Υ	Υ	N	
	MH-5	SE	N	2"	N	Υ	N	
	MH-7/PC1	SE	Υ	2.7'	N	Υ	N	
	MH-9	SE	Υ	2"	N	Υ	N	
	MH-11	SE	Υ	2"	Υ	Υ	N	
South Frontier	MH-13	SE	Υ	2"	N	Υ	N	

Signature: Darrell Crockett

Love Canal Semi-Annual Barrier Drain Manhole Inspection

Date 10/19/2020 МН Water **Debris** Structure Cleaning Level **Comments** Sector No. Location Y/N Y/N ОК **Feet** Y/N MH-10A NW0 North Colvin Ν Ν Υ Ν MH-8A NW Υ 0.4 Ν Υ Ν 1" MH-6C Υ Υ Υ Some built up sludge - no flow restriction NW Ν 1" MH-6B NWΥ Υ Υ Some built up sludge - no flow restriction Ν MH-6A 1" Some built up sludge - no flow restriction NW Υ Υ Υ Ν PC-2A Υ 4' Υ NW Ν Ν 1" MH-4A NW Υ Ν Υ Ν 2" MH-2A NW Υ Ν Υ Ν MH-2 3" Υ SW Υ Υ Ν MH-4 5" SW Υ Υ Υ Ν 3" MH-6 SW Υ Ν Υ Ν MH-8/PC2 3' SW Υ Ν Υ Ν 3" Υ MH-10 SW Υ Υ Ν MH-12 SW Υ 4" Ν Υ Ν South Frontier MH-14 SW Υ 1" Ν Υ Ν North Colvin NH-17A NE Ν 0 Ν Υ Ν MH-15A NE Υ Ν Υ Ν 2" MH-13A Υ Υ Υ NE Ν Υ 2' PC1A NE Υ Υ Ν 2" MH-11A NE Υ Ν Υ Ν MH-9A 1" Υ Υ NE Ν Ν MH-7A 2" NE Υ Υ Υ Ν 1" Υ MH-5A Ν Ν NE Ν MH-3A Υ 2" Υ NE Ν Ν MH-1A 2" Υ NE Υ Ν Ν MH-1 2" SE Ν Ν Υ Ν MH-3 SE Υ 2" Υ Υ Ν 2" Υ MH-5 SE Ν Ν Ν MH-7/PC1 SE Υ 3.5' Ν Υ Ν 1" SE MH-9 Υ Ν Υ Ν 3" MH-11 SE Υ Υ Υ Ν South Frontier MH-13 1" Υ SE Υ Ν Ν

Signature: Darrell Crockett

Appendix C Niagara Falls Water Board Wastewater Discharge Permit #44

NIAGARA FALLS WATER BOARD

SIGNIFICANT INDUSTRIAL USER WASTEWATER DISCHARGE

Permit No. 44

In accordance with all terms and conditions of the Niagara Falls Water Board Wastewater Regulations Part 1960 and also with all applicable provisions of Federal and State Law or regulation:

Permission Is Hereby Granted To:

GLENN SPRINGS HOLDINGS, INC. – LOVE CANAL LEACHATE TREATMENT FACILITY

Located at: 805 - 97th Street, Niagara Falls, NY 14304

Classified by SIC Number: 4952

For the contribution of wastewater into the Niagara Falls Water Board Publicly-Owned Treatment Works (POTW).

Effective this 10th day of January 2020 To Expire this 9th day of January 2025

Signed this 20th day of December 2019

For Pat Fama Executive Director of the Niagara Falls Water Board

DISCHARGE IDENTIFICATION

OUTFALL	DESCRIPTION	LOCATION	RECEIVING
#1	97 th Street Discharge	97 th Street	Carbon treated leachate from the Love Canal Leachate Treatment Facility and the 102 nd Street landfill

A. <u>Discharges to the Niagara Falls Water Board (NFWB) Sewer</u>

	TEWATER DISCHARGE PERMIT JIREMENTS FOR:	ACTION REQUIRED	REQUIRED DATE OF SUBMISSION
1.	Identification of all discharges to the NFWB Sewer System on a current plant sewer map certified by a New York State licensed professional engineer.	None	Submitted 12/06/2019
2.	Identification of each contributing waste stream to each discharge to the NFWB Sewer System clearly marked on, or referenced to, a current plant sewer map certified by a New York State licensed professional engineer.	None	Submitted 12/06/2019
3.	Elimination of all uncontaminated discharges to the NFWB Sewer System. All uncontaminated flows should be clearly identified on a current sewer map certified by a New York State licensed professional engineer.	N/A	
4.	Establishment of a control manhole that is continuously and immediately accessible for each discharge to the NFWB Sewer System.	None	Previously Established

B. <u>Wastewater Discharge Management Practices</u>

 Identification of a responsible person(s) (day to day and in emergencies). Updated as needed by NFWB personnel

C. <u>General Wastewater Discharge Permit Conditions</u>

- 1. Flow monitoring should be performed concurrently with any Wastewater Discharge Permit sampling and should be reported at the same time as analytical results. If it is not feasible to perform flow monitoring, an estimate of flow (method of estimated flow preapproved by the Niagara Falls Water Board) should be submitted with the analytical results.
- 2. All sampling for billing and pretreatment compliance purposes will be coordinated through the Niagara Falls Water Board Industrial Monitoring Coordinator.
- 3. All analysis must be performed by a State certified laboratory using analytical methods promulgated and consistent with 40 CFR 136 and amendments thereto. The permittee will request their contract laboratory to report both Practical Quantitation Limit (PQL) and Method Detection Limit (MDL). The PQL and MDL are defined in the NYSDEC Technical Guidance Series 1.3.7.

The permittee should report results that are less than the MDL or PQL on the NFWB Self Monitoring Report, as non-detect (ND), by placing a less than sign (<) followed by the analytical result. Every effort should be made to attain results down to the MDL. If this is not possible, then results less than PQL but greater than MDL must also be additionally flagged with the qualifier "J" on the Self Monitoring Report. For example, a result less than 5 PQL would be reported <5 (J). In either case the calculated load in lbs per day would be zero. Monitoring results which are lower than the PQL must be reported but will not be used to determine compliance with the permit limit.

- 4. An estimate of relative production levels for wastewater contributing processes at the time of any pretreatment compliance sampling will be submitted upon request of the Director of Niagara Falls Water Board Wastewater Facilities.
- 5. All samples will be handled in accordance with EPA approved methods. Chain of Custody records will be submitted with all sampling results.
- 6. All conditions, standards and numeric limitations of Niagara Falls Water Board Wastewater Regulations are hereby incorporated into this permit by reference. These conditions, standards and numeric limitations must be complied with. Failure to comply with any part of said regulations constitutes a violation and is subject to enforcement actions(s) described in Section 1960.9 of said Regulations, and in the Niagara Falls Water Board Pretreatment Administrative Procedure Number Five (5) "Enforcement Response Guide." Violators are subject to all applicable *Civil* and *Criminal* penalties. In the event of a violation, including slug discharges or spills, the Niagara Falls Water Board must be notified immediately by phone and confirmed by letter within five (5) working days. (C6. continued)

C. <u>General Wastewater Discharge Permit Conditions</u> (continued)

6.cont. Any person adjudicated of violating any provision in the Niagara Falls Water Board Wastewater Regulations shall be assessed a fine in the amount of up to \$10,000. This amount is available for each violation, and each day of a violation is a separate incident for which penalties may be sought.

The person violating any of the provisions of the Niagara Falls Water Board Wastewater Regulations will be liable for any expense, loss, or damage occasioned by reason of such violation. The expense, loss or damage will be taken to be to the extent determined by the Director.

In addition, any person who knowingly makes any false statements; representation or certification in any application, record, report, plan or other document filed or required to be maintained pursuant to the Niagara Falls Water Board Wastewater Regulations, or Wastewater Discharge Permit, or who falsifies, tampers with, or knowingly renders inaccurate any monitoring device or method required under the Niagara Falls Water Board — Wastewater Regulations will, upon conviction be punished by a fine up to \$5,000. Furthermore, the Niagara Falls Water Board may recover reasonable attorney's fees, court costs, court reporting fees, and other expenses of litigation by appropriate suit at law against the person found to have violated applicable laws, orders, rules and permits required by the Niagara Falls Water Board Wastewater Regulations.

- 7. In accordance with Federal Regulation CFR 40, Part 403.12(g), any exceedance of a numeric limitation noted by the SIU must be re-sampled, analyzed and resubmitted to the of Niagara Falls Water Board Wastewater Facilities within 30 days of becoming aware of the exceedance.
 - Specifically, if any limit that is <u>listed</u> in Section E of this permit is exceeded, then the permittee will undertake a short-term monitoring program for that pollutant. Samples will be collected identical to those required for routine monitoring purposes and will be collected on each of at least <u>two (2)</u> operating days and analyzed. Results will be reported in both concentration and mass, and will be submitted within <u>30</u> days of becoming aware of the exceedance.
- 8. Sampling frequency for any permitted compounds may be increased beyond the requirements set forth in Section E and F of this permit. If the permittee monitors (sample and analysis) more frequent than required under this permit, <u>all</u> results of this monitoring must be reported.
- 9. As noted in Section 1960.5g of the Niagara Falls Water Board Wastewater Regulations, "Personnel as designated by the Director will be permitted at any time for reasonable cause to enter upon all properties served by the Niagara Falls Water Board Wastewater Facilities for the purpose of, and to carry out, inspection of the premises, observation, measurement, sampling and testing, in accordance with provisions of the Regulations." (C. continued)

- C. <u>General Wastewater Discharge Permit Conditions</u> (continued)
- 10. As noted in Section 1960.5c of the Niagara Falls Water Board Wastewater Regulations, significant changes in discharge characteristics or volume must be reported immediately to the Niagara Falls Water Board Wastewater Facilities.
- 11. As noted in Section 1960.6b of the Niagara Falls Water Board Wastewater Regulations,
 - samples required to be collected via a 24-hour composite sampler must be retained refrigerated for an additional 24 hour plus unrefrigerated an additional 48 hours (total 72 hours).
- 12. As noted in Section 1960.5d of the Niagara Falls Water Board Wastewater Regulations, all "SIU's will keep on file for a minimum of three years, all records, flow charts, laboratory calculations or any other pertinent data on their discharge to the Niagara Falls Water Board Wastewater Facilities.
- 13. As noted in Section 1960.6g of the Niagara Falls Water Board Wastewater Regulations, "Permits are issued to a specific user for a specific monitoring station. A permit will not be reassigned or transferred without the approval of the Director which approval will not be unreasonably withheld. Any succeeding owner or user to which a permit has been transferred and approved will also comply with all the terms and conditions of the existing permit."
- 14. The Annual Average Limitation is equivalent to the specific SIU allocation, and will be defined as the permissible long-term average discharge of a particular pollutant. These limitations are listed in Section E of this permit. The computation of the Annual Average will be as follows; for each compound listed in Section F of this permit, the Annual Average will be the average of the present monitoring quarter and **three previous quarters' dat**a.
- 15. The Daily Maximum Limitation will be defined as the maximum allowable discharge on anyone day. The Daily Maximum Limitation will allow for periodic short-term discharge fluctuations. These specific limitations are listed in Section E of this permit.
- 16. Enforcement of the Annual Average Limitation will be based on the reported average of the last four quarters data vs. the Annual Average Limited listed in Section E of this permit. Enforcement of the Daily Maximum Limitation will be based on individual analysis results vs. the Daily Maximum Limit listed in Section E of this permit. These results may be obtained from self-monitoring (Section F), Niagara Falls Water Board Verification, incident investigation or billing samples. (C. continued)

C. <u>General Wastewater Discharge Permit Conditions</u> (continued)

17. The Niagara Falls Water Board Administrative Procedure Number 6 "Procedure for Determination and Use of Local Limits" lists all pollutants noted in the Niagara Falls Water Board – Wastewater Facilities SPDES Permit. The limits defined in the procedure are values which are based on the quantity of substances discharged which can be easily related to the Treatment Plant's removal capacity.

The pollutants listed in this procedure which are <u>not</u> specifically listed in Section E and F of this permit may be present in the permittee's wastewater discharge, but at levels which do not require specific permit limitations. Consequently, if any of the limits listed in this procedure, for pollutants <u>not</u> identified in Section E and F of this permit, are exceeded then the permittee will undertake a short-term, high intensity monitoring program for that pollutant.

Samples identical to those required for routine monitoring purposes will be collected on each of at least three operating days and analyzed. Results will be expressed in terms of both concentration and mass, and will be submitted no later than the end of the third month following the month when the limit was first exceeded.

If levels higher than the limit are confirmed, the permit may be reopened by the Niagara Falls Water Board for consideration of revised permit limits.

- 18. 40 CFR 403.17(c) contains the notification requirements for anticipated and unanticipated bypass. In the event of an <u>anticipated bypass</u>, the federal regulations at 403.17 (c)(1) require an industrial user to notify the Niagara Falls Water Board at least ten days prior to the date of the bypass. In the event of an <u>unanticipated bypass</u>, the federal regulations at 40 CFR 403.17(c)(2) require an industrial user to notify the Niagara Falls Waster Board within 24 hours from the time the industrial user becomes aware of the bypass and submit a written description of the bypass within five days of the time the industrial user becomes aware of the bypass. As defined at 403.17 (a)(1), a bypass is "the intentional diversion of wastestreams from any portion of an Industrials User's treatment facility."
- 19. 40 CFR 403.8(f)(2)(vi) requires the SIU to notify the Niagara Falls Water Board immediately of any changes at its facility affecting its potential for a slug discharge.

D. <u>Specific Wastewater Discharge Permit Conditions</u>

1. <u>Billing Agreement</u>:

- a) Flow quantities will be derived from the Wastewater Treatment Facility flow meter. The results of the daily flow readings will be compiled and submitted to the Niagara Falls Water Board in a Monthly Flow Report by the 15th day of the following month.
- b) Charges for TSS, SOC and Substances of Concern shall be developed based on Quarterly Self-Monitoring data.

2. <u>Love Canal Leachate Treatment Facility (LCLTF)</u>

The Niagara Falls Water Board agrees to accept wastewater processed from the Glenn Springs Holdings (GSH) LCLTF. These waters in addition to Love Canal wastewater shall include wastewater from the 102nd Street remedial site. This approval is subject to the following conditions:

- a) The LCLTF shall be properly operated and maintained at all times.
- b) To ensure proper operation GSH shall ensure sufficient feed, inter-stage (breakthrough), and effluent analysis to ensure timely carbon changes. Treatment levels of 10 **ug/l** shall be achieved and verified with quarterly composite sample analysis for the following compounds: trichloroethylene, tetrachloroethylene, monochlorotoluene, monochlorobenzenes, trichlorobenzenes, tetrachlorobenzenes, hexachlorocyclohexanes- alpha, beta, gamma and delta and hexachlorobenzene.
- c) The issuance of this approval if based on GSH's previous assertions that there is no reason to anticipate the presence of tetrachlorodibenzo-p-dioxins in the discharge from the treatment facility. The Niagara Falls Water Board hereby reserves the right to collect samples from the treatment facility effluent and analyze such wastewaters for their chemical constituents, including tetrachlorodibenzo-p-dioxins. If such analysis indicates the presence of tetrachlorodibenzo-p-dioxins, this approval may be withdrawn. If at any time, the Niagara Falls Water Board determines on any basis that the discharge of these wastewater to the POTW is interfering with the operation of that facility, the Niagara Falls Water Board will direct GSH to discontinue the discharge.
- d) These pretreated wastewaters shall be discharged to the POTW via Outfall MS # 1.

(D2. continued)

D. <u>Specific Wastewater Discharge Permit Conditions</u> (continued)

2. <u>Love Canal Leachate Treatment Facility (LCLTF)</u> (continued)

e) Periodically wet weather flow in the area around LCLTF results in surcharged sewers. The resultant surcharge requires overflow at combined sewer and storm sewer overflow points. Other points in the sewer shed require manual bypass pumping. Consequently, to minimize this overflow, the Niagara Falls Water Board will require the permittee to cease discharge from the LCLTF during these surcharge events.

A notification procedure has been established by the Niagara Falls Water Board to formalize the communication between the Niagara Falls Water Board and the permittee to halt and resume the LCLTF discharge. This procedure by reference is hereby incorporated as a condition of this permit.

3. <u>Slug Discharge Control Plan</u>:

Pursuant to the regulations contained in the Federal Industrial Pretreatment Program, 40CFR 403.8(f)(2), the NFWB is obligated to periodically review users for the need for a Slug Discharge Control Plan.

This permittee has been reviewed and is <u>NOT</u> required to develop and implement such a plan.

E. <u>Discharge Limitations & Monitoring Requirements</u>

During the Period beginning the effective date of this Permit and lasting until the expiration date, discharge from the permitted facility outfall(s) will be limited and monitored by the permittee as specified below.

	DISCHARGE LIMITATIONS			MINIMUM MONITORING REQUIREMENTS	
OUTFALL NUMBER/ EFFLUENT PARAMETER	ANNUAL AVERAGE	DAILY MAXIMUM	UNITS	MEASUREMENT FREQUENCY	SAMPLE TYPE
MS #1 Flow	0.3	0.3	MGD	Continuous	4
MS #1 Total Suspended Solids	25	50	lbs./d	1/Quarter	1
MS #1 Soluble Organic Carbon	50	75	lbs./d	1/Quarter	1
MS#1 Volatile Priority Pollutants (See Attached list (Section F-1)	MONITO	OR ONLY	lbs./d	1/Quarter	1
MS #1 Acid Extractable Priority Pollutants (See attached list Section F-2)	MONITO	OR ONLY	lbs./d	1/Quarter	1
MS #1 Base/Neutral Priority Pollutants (See attached list Section F-3)	MONITO	OR ONLY	lbs./d	1/Quarter	1
MS #1 Pesticides Hexachlorocyclohexanes- alpha, beta, gamma and delta (See attached list Section F-4)	MONITO	OR ONLY	lbs./d	1/Quarter	1
MS #1 Total Phenols	MONITO	OR ONLY	lbs./d	1/Quarter	1

SAMPLE TYPE FOOTNOTES

- (1) Each sample will consist of four (4) grabs collected spaced throughout the batch discharge, such that they are representative of the effluent being discharged pursuant to 40CFR 403.12.b5iii. The four (4) grabs will be composited in the laboratory and analyzed as one sample.
- (2) Each sample will consist of four (4) grabs collected spaced over the 24-hour period, such that they are representative of the effluent being discharged pursuant to 40CFR 403.12.b5iii. The four (4) grabs will be composited in the laboratory and analyzed as one sample.
- (3) Each sample will consist of a 24-hour, flow proportioned composite sample collected from the monitoring point.
- (4) Flow will be monitored continuously with the use of a water meter or another acceptable flow metering device.
- (5) Each sample will consist of a 24-hour, time proportioned composite sample collected from the monitoring point.
- (6) Reserved
- (7) Same as (3), however, five (5) samples will be collected per quarter from the monitoring point and analyzed by and at the Niagara Falls Water Board's expense.
- (8) Four (4) grab samples will be collected spaced over the 24-hour period, such that they are representative of the effluent being discharged pursuant to 40CFR 403.12.b5iii. Each grab will be analyzed and reported separately.
- (9) A grab sample is defined as an aliquot collected over a period of not more than 15 minutes.

F. <u>Discharge Monitoring Reporting Requirements</u>

During the period beginning the effective date of this permit and lasting until its expiration date, discharge monitoring results will be summarized and reported by the permittee; Monthly - 14 days after monitoring period, Quarterly - by the last day of the monitoring period = February 28, May 31, August 31, November 30. Semiannual reports will be submitted on the last day of the monitoring period = February 28, August 31. The annual average for each parameter listed in Section F, will be computed and reported quarterly. The individual sample analysis for present quarter will also be reported quarterly unless directed otherwise in this permit.

OUTFALL NO	PARAMETER	REPORTING FREQUENCY
MS #1	Flow	Monthly
MS #1	Total Suspended Solids	Quarterly
MS #1	Soluble Organic Carbon	Quarterly
MS #1	Volatile - Priority Pollutants (F-1) *	Quarterly
MS #1	Acid Extractables - Priority Pollutants (F-2) *	Quarterly
MS #1	Base/Neutral - Priority Pollutants (F-3) *	Quarterly
MS #1	Pesticides (F-4) *	Quarterly
MS #1	Total Phenols	Quarterly

^{*} See specific compounds listed on the following page.

(F. continued)

F. <u>Discharge Monitoring Reporting Requirements</u> (continued)

Discharge Monitoring Compounds

F1- Volatile Priority Pollutants					
Benzene	Bromoform	Trichloroethylene			
Carbon Tetrachloride	Dichloropropylenes	Methylene Chloride			
Chlorodibromethane	Ethylbenzene	Vinyl Chloride			
Monochlorobenzene	Tetrachloroethanes	Monochlorotoluenes			
Dichlorobromethane	Tetrachloroethylene	Monochlorobenzotrifluoride			
Chloroform	Toluene				
Dichloroethylenes	Trichloroethanes				

F2- Acid Extractables Priority Pollutants				
Monochlorophenol	Monochlorocresol	Pentachlorophenol		
Dichlorophenol	Trichlorophenol			

F3- Base/Neutrals Extractables Priority Pollutants					
Dimethyl Phthalate	Dichlorotoluene	Trichlorobenzene			
Butyl Benz Phthalate	Acenaphthlene	Trichlorotoluene			
Di-N-Butyl Phthalate	Fluoranthene	Hexachlorobutadiene			
Di-N-Octyl Phthalate	Chrysene	Tetrachlorobenzene			
Diethyl Phthalate	Napthalene	Hexachlorocyclopentadiene			
Nitrosodiphenylamine	Benzo (a) Anthracene	Hexachlorobenzene			
Dichlorobenzenes	Pyrene	Dichlorobenzotrifluoride			

F4- Pesticides
Hexachlorocyclohexanes- alpha, beta, delta, and gamma

G. <u>Comments/Revisions</u>

Appendix D Annual Groundwater Sampling Schedule



2055 Niagara Falls Blvd., Suite #3 Niagara Falls, New York 14304

Telephone: (716) 297-6150 Fax: (716) 297-2265

www.CRAworld.com

MEMORANDUM

To:

Clint Babcock, Ralph Schupp

REF. NO.:

009954

FROM:

Jane Pietraszek-Polovich/adh/8

DATE:

August 5, 2010

C.C.:

Darrell Crockett, Dennis Hoyt, John Pentilchuk,

Dave Tyran, Filing

RE:

Love Canal Annual Groundwater Sampling Schedule

At the request of Glenn Springs Holdings, Inc. (GSH), Conestoga-Rovers & Associates (CRA) has prepared the following memo to document the Annual Groundwater Sampling schedule for the Love Canal Facility in Niagara Falls, New York (Site).

Conservation (NYSDEC) sent to CRA and GSH on March 25, 2009 (email attached) states that it is no longer necessary for the NYSDEC to specifically list the wells to be sampled each year at the Site, since the annual and alternating (Group I and Group II) wells have remained the same throughout the years. From 1994 through 2008, the NYSDEC provided GSH with a list of wells to be sampled each year. The March 25, 2009 email from Mr. Sadowski stated that the NYSDEC will no longer provide such a list. Therefore, Mr. Sadowski suggested that the wells sampled during the 2007 annual groundwater monitoring event be used for the 2009 annual groundwater monitoring event, to remain consistent with the Long-Term Monitoring Program. The 2007 (and therefore 2009) monitoring wells represent the Group I wells (Table 1). The 2008 (and therefore 2010) monitoring wells represent the Group II wells. In addition, there are select overburden and bedrock wells that are to be sampled annually (Table 1).

Mr. Sadowski went on to further state that GSH must "ensure that the monitoring network and well selection provide adequate overburden and bedrock coverage that returns the data necessary for the evaluation of the remediation, and that the NYSDEC feels that the selection of the 2007 wells will meet those objectives." Mr. Sadowski indicated that GSH can enhance upon the objective by choosing other wells if they wish. Once the well selection is made for the annual event, GSH is to provide the NYSDEC with the monitoring well numbers. Any changes in the well selection must be accompanied with reasons for the addition/ deletion. Based on a review of the data for the wells suggested by NYSDEC, GSH agreed to sample the wells in Table 1 for future sampling events. This was communicated to the NYSDEC through a phone call to Mr. Sadowski on June 7, 2010, and documented in the attached email dated June 8, 2010. The NYSDEC is to be notified when the annual monitoring will take place for oversight purposes and to split samples if desired. A 2-week notice of the annual groundwater monitoring event is preferred by the NYSDEC.



TABLE 1

SAMPLE SCHEDULE LOVE CANAL FACILITY LONG-TERM MONITORING PROGRAM NIAGARA FALLS, NEW YORK

Annual Wells	Biannual Wells			
	Overburden Wells	Overburden Wells		
Bedrock Wells	Group I (2009)	Group II (2010)		
3257	3151	7115		
5221	7120	7125		
6209	7 155	8115		
7205	7161	8125		
8210	8110	9105		
9205	8120	9113		
9210	8130	9118		
10205	8140	10178 A		
10210A	9110			
10210B	9115			
10210C	9120			
10215	9125			
10225A	9130			
10225B	9140			
10225C	10105			
10270	10147			
10272	10174A			
10278				

Overburden Wells

7130 7132

8106

10135

Re: Love Canal Annual 2009 Sampling

Page 1 of 1

From: Crockett, Darrell

Sent: Wednesday, March 25, 2009 12:06 PM

To: Pentilchuk, John

Subject: 9954 FW: Love Canal Annual 2009 Sampling

John,

Please let me know how you'd like for me to proceed. I have the 2007 sampling event data.

Thanks Darrell

From: Brian Sadowski [mailto:bpsadows@gw.dec.state.ny.us]

Sent: Wed 3/25/2009 11:56 AM

To: Crockett, Darrell

Cc: Hoyt, Dennis; Clint_Babcock ext

Subject: Re: Love Canal Annual 2009 Sampling

Darrell,

Your contact and this response will be considered as our pre-sampling conference as stated on p.6. in Section 2.0 Monitoring Requirements of the February 19, 2001 Sampling Manual. Over the last fourteen years the Department has specifically listed the wells to sample and believe that is no longer necessary as the annual and alternating wells have stayed the same. GSHI and/or MSRM has clearly demonstrated their ability to operate, maintain and monitor the site. With the addition of CRA; there is an added layer of technological security and professional environmental judgement. With that said, the Department will not provide a specific list of wells to monitor. However, the suggested wells are the wells that were sampled in 2007 to remain consistent with the LTM program schedule listed on Table 2.2 of the Sampling Manual. The monitoring schedule is flexible. OXY and CRA is to ensure that the monitoring network and well selection provides adequate overburden and bedrock coverage that returns the data necessary for the evaluation of the remediation. The Department feels that the selection of the 2007 wells will meet those objectives. OXY and CRA can enhance upon the objectives by choosing other wells if they wish. When a decision is made by CRA on the well selection, please provide the Department with the well numbers. A simple return "as same as 2007 or 2007 with the addition, deletion or substitution of well #" will suffice. If changes are made, please provide reasoning. Finally, the Department will not be splitting this year. But, will need to be given notice when monitoring will take place for oversight purposes. Thank you.

>>> "Crockett, Darrell" <dcrockett@craworld.com> 3/18/2009 8:05 AM >>> Hello Brian,

At your convenience would you please provide me with a 2009 Annual sampling well list including the split samples.

Thank You Darrell Crockett 716/998-5804 From: Polovich, Jane

Sent: Tuesday, June 08, 2010 9:49 AM

To: Filing

Subject: 9954: Love Canal Annual Sampling - Conversation with Brian Sadowski NYSDEC I spoke with Brian Sadowski on Monday June 7, 2010 to confirm the annual groundwater sampling locations at Love Canal. As per Brians email of March 25 2009 to Darrell Crockett, the 2007 wells selected by the DEC were to be the ones sampled in 2009 (these represent the Group I wells). I confirmed with Mr. Sadowski that the 2008 wells selected by the DEC would be teh wells sampled in 2010 (representing the Group II wells). These Group I and Group II wells will from this point forward be the wells sampled for the alternating annual groundwater sampling events at Love Canal. Mr Sadowski went further to say that GSH may add or delete wells from the sampling but must provide a reason to the DEC prior to sampling for the additions / deletions. CRA documented this change in the sampling program in a internal memo.

Jane Polovich Conestoga-Rovers & Associates (CRA)

2055 Niagara Falls Blvd., Suite 3 Niagara Falls, New York 14304

Phone: 716.297,6150 Fax: 716.297.2265

Email: jpolovich@CRAworld.com

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Appendix E Laboratory Reports



Service Request No:R2005470

Ms. Kathy Willy GHD Services Inc. 2055 Niagara Falls Blvd., Niagara Falls, NY 14304

Laboratory Results for: Love Canal:292-402-D02-3100

Dear Ms. Willy,

Enclosed are the results of the sample(s) submitted to our laboratory June 25, 2020 For your reference, these analyses have been assigned our service request number R2005470.

All testing was performed according to our laboratory's quality assurance program and met the requirements of the TNI standards except as noted in the case narrative report. Any testing not included in the lab's accreditation is identified on a Non-Certified Analytes report. All results are intended to be considered in their entirety. ALS Environmental is not responsible for use of less than the complete report. Results apply only to the individual samples submitted to the lab for analysis, as listed in the report. The measurement uncertainty of the results included in this report is within that expected when using the prescribed method(s), and represented by Laboratory Control Sample control limits. Any events, such as QC failures or Holding Time exceedances, which may add to the uncertainty are explained in the report narrative or are flagged with qualifiers. The flags are explained in the Report Qualifiers and Definitions page of this report.

Please contact me if you have any questions. My extension is 7471. You may also contact me via email at Brady.Kalkman@alsglobal.com.

Respectfully submitted,

Goody Kullen

ALS Group USA, Corp. dba ALS Environmental

Brady Kalkman

Project Manager



Narrative Documents

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com



Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100 Date Received: 06/25/2020

Sample Matrix: Water

CASE NARRATIVE

All analyses were performed consistent with the quality assurance program of ALS Environmental. This report contains analytical results for samples for the Tier level IV requested by the client.

Sample Receipt:

Eight water samples were received for analysis at ALS Environmental on 06/25/2020. Any discrepancies upon initial sample inspection are annotated on the sample receipt and preservation form included within this report. The samples were stored at minimum in accordance with the analytical method requirements.

Semivolatiles by GC/MS:

Method 8270D, 06/30/2020: The lower control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). Since there were no detections of the analyte(s) above the MRL in the associated field samples, the quantitation is not affected. The data quality was not significantly affected and no further corrective action was taken. Method 8270D, 07/01/2020: The upper control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). The field samples analyzed in this sequence did not contain the analyte(s) in question above the Method Reporting Limit (MRL). Since the exceedance equates to a potential high bias, the data quality was not significantly affected and no further corrective action was taken.

Method 8270D, 07/01/2020: The lower control limit for the spike recovery of the Laboratory Control Sample Duplicate (LCSD) was exceeded for one or more analyte. Precision is also outside limits. There were no detections of the analyte(s) in the associated field samples. The LCS and batch MS/MSD were within limits for these anlytes. The analytes affected are flagged in the LCS Summary.

Method 8270D, 07/06/2020: The upper control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). The field samples analyzed in this sequence did not contain the analyte(s) in question above the Method Reporting Limit (MRL). Since the exceedance equates to a potential high bias, the data quality was not significantly affected and no further corrective action was taken.

Method 8270D, 07/06/2020: The control limit was exceeded for one or more surrogates in the Continuing Calibration Verification (CCV). The surrogates were within acceptance limits for the associated field samples. The data quality was not significantly affected and no further corrective action was taken.

Method 8270D, R2005470-003: The control limits were exceeded for one or more surrogates in the sample(s). Since the exceedance may indicate a potential bias in the analytical batch, all associated field samples were re-extracted and reanalyzed. The surrogates met control limits for the reanalysis. Since the results for the field samples were comparable for both determinations, the exceedance in the initial analysis was likely restricted to the surrogate recovery. Therefore, the results from the original analysis are reported and flagged.

Semivoa GC:

Method 8081B, 07/01/2020: The upper control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). The field samples analyzed in this sequence did not contain the analyte(s) in question above the Method Reporting Limit (MRL). Since the exceedance equates to a potential high bias, the data quality was not significantly affected and no further corrective action was taken.

Method 8081B, 07/01/2020: The lower control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). Since there were no detections of the analyte(s) above the MRL in the associated field samples, the quantitation is not affected. The data quality was not significantly affected and no further corrective action was taken.



Volatiles by GC/MS:

Method 8260C, 07/01/2020: The lower control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). Since there were no detections of the analyte(s) above the MRL in the associated field samples, the quantitation is not affected. The data quality was not significantly affected and no further corrective action was taken.

Method 8260C, 07/01/2020: The upper control criterion was exceeded for one or more analytes in the Laboratory Control Sample (LCS). There were no detections of the analyte(s) above the MRL in the associated field samples. The error associated with elevated recovery equates to a high bias. The sample data is not significantly affected. No further corrective action was appropriate.

Method 8260C, 07/02/2020: The upper control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). The field samples analyzed in this sequence did not contain the analyte(s) in question above the Method Reporting Limit (MRL). Since the exceedance equates to a potential high bias, the data quality was not significantly affected and no further corrective action was taken.

Method 8260C, 07/02/2020: The upper control criterion was exceeded for one or more analytes in the Laboratory Control Sample (LCS). There were no detections of the analyte(s) above the MRL in the associated field samples. The error associated with elevated recovery equates to a high bias. The sample data is not significantly affected. No further corrective action was appropriate.

Approved by _____

Date 07/14/2020



SAMPLE DETECTION SUMMARY

CLIENT ID: WG-9954-062420-SG-001		Lab	ID: R2005	5470-002		
Analyte	Results	Flag	MDL	MRL	Units	Method
Bis(2-ethylhexyl) Phthalate	1.7	J	0.91	9.1	ug/L	8270D
CLIENT ID: WG-9954-062420-SG-004		Lab	ID: R2005	5470-005		
Analyte	Results	Flag	MDL	MRL	Units	Method
Carbon Disulfide	0.83	J	0.42	10	ug/L	8260C



Sample Receipt Information

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request:R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

SAMPLE CROSS-REFERENCE

SAMPLE #	CLIENT SAMPLE ID	<u>DATE</u>	<u>TIME</u>
R2005470-001	TB-9954-062420-SG-001	6/24/2020	0000
R2005470-002	WG-9954-062420-SG-001	6/24/2020	0910
R2005470-003	WG-9954-062420-SG-002	6/24/2020	0955
R2005470-004	WG-9954-062420-SG-003	6/24/2020	1030
R2005470-005	WG-9954-062420-SG-004	6/24/2020	1105
R2005470-006	WG-9954-062420-SG-005	6/24/2020	1150
R2005470-007	WG-9954-062420-SG-006	6/24/2020	1245
R2005470-008	WG-9954-062420-SG-007	6/24/2020	1335

CHA!N-OF-CUSTODY/Analytica! Request Document The Chain-of-Custody is a LEGAL DOCUMENT. All relevant fields must be completed accurately.

Client In	formation
GLEN SPRINGS HOLDINGS INC	Report To: Kathy Willy
806 97TH STREET	Сору То:
LOVE CANAL	
NIAGARA FALLS, NEW YORK 14304	Invoice To:
Phone: 716-283-0111	PO:
Fax: 716-283-2866	Project Name: LOVE CANAL ANNUAL GW
Email: kathy.willy@ghd.com	Project Number: 9954

	Information
Laboratory: ALS	
Laboratory Location: 15 1565 JEFFERSON RD. BUILDING 300, SUITE 36 ROCHESTER, NY 14623	
Laboratory Contact: BR	ADY KALKMAN
Requested Due Date:	TAT: 10
QA/QC Requirements:	

	Event Information
	ID#: LC ANNUALGW SAMPLING 2020-01-1
	SSOW Ref#: 292-402-999-3100
Sam	oler Name: S GARDNER, DTYRAN

	Valid Matrix Code WG Groundwater WB Borehole Water WS Surface Water SO Soil SE Sediment	Matrix Code	Date Collected	Time Collected	PestPCBs(None)	SVOC(none)	VOA(HCI)		Sample Cond Temp in C Received on ice Sealed Cooler Samples Intact	Y/N Y/N Y/N
Sample Identification				-	ءَ ا	ίς	>	Remarks		
TB-9954-062420-SG-001		WG Q	06/24/2020	00:00	-	-	3			
WG-9954-062420-SG-001		WG	06/24/2020	09:10	2	2	3			
WG-9954-062420-SG-002		WG	06/24/2020	09:55	2	2	3		,	•
WG-9954-062420-SG-003		WG	06/24/2020	10:30	2	2	3			
WG-9954-062420-SG-004		WG	06/24/2020	11:05	2	2	3		161 3 1111 3	
WG-9954-062420-SG-005		WG	06/24/2020	11:50	2	2	3			
WG-9954-062420-SG-006		WG	06/24/2020	12:45	2	2	3			
WG-9954-062420-SG-007		WG	06/24/2020	13:35	2	2	3			
Total Bottles				4	14	14	24	Grand Tot	al:52	

SHIPMENT METHOD	NO. OF COOLERS	RELINQUISHED BY:	DATE	TIME	RECIEVED BY:	DATE	TIME
FedEx	2	Shaw Mardner	6/24/2	1500		695ab	20 11:00
AIRBILL#:							

R2005470 5
GHD Services Inc.
Love Canal:292-402-D02-3100

SR# CHAIN OF CUSTODY / LABORATORY ANALYSIS REQUEST FORM 004, 005, 006, 007, 008, 009, 010, 1565 Jefferson Road, Bldg 300, Suite 360, Rochester, NY 14623 T030477 011, 012, 013 Phone (585) 288-5380 / FAX (585) 288-8475 S) Environmental www.alsglobal.com **14**D 7 Love Canal:292-402-D02-3100 Project Number: 9954 Annual Long Term Monitoring Report To Kathy Willy NUMBER OF CONTAINERS Company / Address GHD Services Inc. 2055 Niagara Falls Bivd., Suite 3 Niagara Falls NY, 14304 8081B / Pest OC 3260C / VOC 3082A / PCB 3270D / SVO Phone # 716-297-2160 718-297-2265 Sampler Signature Sampler Printed Name Remarks SAMPLING Matrix **CLIENT SAMPLE ID** LABID Date Time Liquid Liquid Liquid Liquid Liquid 6. Liquid Liquid Liquid Liquid Liquid **Turnaround Requirements** Report Requirements Invoice Information Special Instructions/Comments: RUSH (SURCHARGES APPLY) I. Results Only P.O.#_ II. Results + QC Summaries (LCS, DUP, MS/MSD as required) Standard (3 weeks) III. Results + QC and Cilibration Bill To:___ Summaries X IV. Data Validation Report REQUESTED FAX DATE with Raw Data EData _____Yes ____No Requested Report Date Received By: Received By: Relinquished By: Relinquished By: Received By: Relinquished By: Signature Signature Signature Signature Signature Signature

Printed Name

Firm

Date/Time

Printed Name

Date/Time

Firm

Printed Name

Firm

Date/Time

Printed Name

Date/Time

Firm

Printed Name

Firm

Date/Time

Printed Name

Date/Time

Firm



Cooler Receipt and Preservation Check Form

Project/Clie	ent <i>GH</i>	<u>/</u>		<u>.</u>	_Fold	er Number_		_				
Cooler receiv	ed on <u>6/2</u>	5/2020	by:_	TO /1	5	COURIER:	ALS	UPS FEI	DEX) VEI	LOCITY CLII	ENT	
1 Were Cu	istody seals on	outside of coole	r?	1	Y) N	5a Percl	ılorate s	samples have	required h	eadspace?	Y N	(NA)
2 Custody	papers prope	rly completed (in	k, sign	ed)?	Y) N	5b Did V	OA via	ls, Alk,or Sul	fide have s	ig* bubbles?	Ø N	NA NA
3 Did all b	ottles arrive in	good condition ((unbrol	(en)?	Ý N	6 When	e did the	bottles origi	inate?	ALS/ROC	CLIE	NT
4 Circle:	Wet Ice Dry	Ice Gel packs	pres	sent?	Y N	7 Soil	/OA rec	eived as:	Bulk F	Encore 5035	iset N	NA
. Temperatu	re Readings	Date: 6/8/	2020	Time:	1122	<u>'</u> ID:	IR#7	(R#10)	From	: Temp Blank) Samj	ple Bottle
Observed Te	emp (°C)	3.8		1.1	0	2.4						
Within 0-6°	C?	Y) N		0	N	N	Y	N Y	'N	Y N	Y	N
If <0°C, wer	re samples froz				N	YN	Y	N Y	·· · •	YN	Y	N
If out of 1	Temperature.	note packing/ic	e cond	ition:		Ice mel	ted P	oorly Packet			Same D	Day Rule
	•	un Samples:						•	=	=		
All samples	held in storag	re location:	6/2	202 p	y <i>4</i> 2	on Ibe	at at	//30				
_	_	orage location:	Was /	טעט	, <u>7/</u> 2	on <i>\(\alpha \)</i>	at at		n 48 hours	of sampling?	V	N
									11 TO HOULD	or sumpling:		
					1-1		1.1-					
		ervation Check**					1451		by: <u>MO</u> NO			
		labels complete (bels and tags agr						VES VES	NO NO			
		ontainers used for				8!		E S	NO			
		s acceptable (no				ng)?		YES	NO		λ√/A	
		Cassettes / Tubes		-		anisters Pressu	rized		r® Bags In	flated	N/A	
pH	Lot of test	Reagent	Preser		Lot Re		Exp	Sample ID		Lot Adde	ed	Final
'	paper		Yes	No			1	Adjusted	Adde	d		pН
≥12		NaOH										
≤2		HNO ₃										
≤2		H ₂ SO ₄	T									ļ
<4		NaHSO ₄										
5-9		For 608pest			No=No	tify for 3day						
Residual		For CN,				ntact PM to add						
Chlorine		Phenol, 625,				3 (625, 608,	1					
(-)	1	608pest, 522			CN), as	corbic (phenol).		ļ				
		Na ₂ S ₂ O ₃						-				
		ZnAcetate	-	-						be tested before ar		_
		HC1	**	**				1		l samples with che	mical pre	servatives
-		1.0.										
·			1					are checked (not just repre	semanves).		
Bottle lot	numbers: 🖍		0 1	427%	VI- 181	MC .		are checked (not just repre	semanves).		
	numbers: <i>C</i>	luit coveres	l O	4276	V-181	MC		are checked (not just repre	semauves).		
Explain a	Il Discrepanci	lent covered es/Other Comm	ents:		V-1B1	MC		are checked (not just repre	scinauves).		
Explain a	Il Discrepanci	lent covered es/Other Comm	ents:		W-1B1	MC		are checked (not just repre	semauves).		
Explain a	Il Discrepanci	lent covered es/Other Comm	ents:		W-1181	мс		are checked (not just repre	semauves).		
Explain a	Il Discrepanci	lent covered es/Other Comm	ents:		W- 1181	MC		are checked (not just repre	semauves).		
Explain a	Il Discrepanci	luit coveres	ents:		V- 1B1	MC		are checked (not just repre	semanves).		

HPROD	BULK
HTR	FLDT
SUB	HGFB
ALS	LL3541

Labels secondary reviewed by:

PC Secondary Review:

*significant air bubbles: VOA \geq 5-6 mm : WC \geq 1 in. diameter Page 10 of 113



Miscellaneous Forms

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com



REPORT QUALIFIERS AND DEFINITIONS

- U Analyte was analyzed for but not detected. The sample quantitation limit has been corrected for dilution and for percent moisture, unless otherwise noted in the case narrative.
- J Estimated value due to either being a Tentatively Identified Compound (TIC) or that the concentration is between the MRL and the MDL. Concentrations are not verified within the linear range of the calibration. For DoD: concentration >40% difference between two GC columns (pesticides/Arclors).
- B Analyte was also detected in the associated method blank at a concentration that may have contributed to the sample result.
- E Inorganics- Concentration is estimated due to the serial dilution was outside control limits.
- E Organics- Concentration has exceeded the calibration range for that specific analysis.
- D Concentration is a result of a dilution, typically a secondary analysis of the sample due to exceeding the calibration range or that a surrogate has been diluted out of the sample and cannot be assessed.
- * Indicates that a quality control parameter has exceeded laboratory limits. Under the õNotesö column of the Form I, this qualifier denotes analysis was performed out of Holding Time.
- H Analysis was performed out of hold time for tests that have an õimmediateö hold time criteria.
- # Spike was diluted out.

- + Correlation coefficient for MSA is <0.995.
- N Inorganics- Matrix spike recovery was outside laboratory limits.
- N Organics- Presumptive evidence of a compound (reported as a TIC) based on the MS library search.
- S Concentration has been determined using Method of Standard Additions (MSA).
- W Post-Digestion Spike recovery is outside control limits and the sample absorbance is <50% of the spike absorbance.
- P Concentration >40% difference between the two GC columns.
- C Confirmed by GC/MS
- Q DoD reports: indicates a pesticide/Aroclor is not confirmed (×100% Difference between two GC columns).
- X See Case Narrative for discussion.
- MRL Method Reporting Limit. Also known as:
- LOQ Limit of Quantitation (LOQ)

 The lowest concentration at which the method analyte may be reliably quantified under the method conditions.
- MDL Method Detection Limit. A statistical value derived from a study designed to provide the lowest concentration that will be detected 99% of the time. Values between the MDL and MRL are estimated (see J qualifier).
- LOD Limit of Detection. A value at or above the MDL which has been verified to be detectable.
- ND Non-Detect. Analyte was not detected at the concentration listed. Same as U qualifier.



Rochester Lab ID # for State Certifications¹

Connecticut ID # PH0556	Maine ID #NY0032	Pennsylvania ID# 68-786
Delaware Approved	New Hampshire ID # 2941	Rhode Island ID # 158
DoD ELAP #65817	New York ID # 10145	Virginia #460167
Florida ID # E87674	North Carolina #676	

¹ Analyses were performed according to our laboratory

NELAP-approved quality assurance program and any applicable state or agency requirements. The test results meet requirements of the current NELAP/TNI standards or state or agency requirements, where applicable, except as noted in the case narrative. Since not all analyte/method/matrix combinations are offered for state/NELAC accreditation, this report may contain results which are not accredited. For a specific list of accredited analytes, contact the laboratory or go to https://www.alsglobal.com/locations/americas/north-america/usa/new-york/rochester-environmental

ALS Laboratory Group

Acronyms

ASTM American Society for Testing and Materials

A2LA American Association for Laboratory Accreditation

CARB California Air Resources Board

CAS Number Chemical Abstract Service registry Number

CFC Chlorofluorocarbon CFU Colony-Forming Unit

DEC Department of Environmental Conservation

DEQ Department of Environmental Quality

DHS Department of Health Services

DOE Department of Ecology DOH Department of Health

EPA U. S. Environmental Protection Agency

ELAP Environmental Laboratory Accreditation Program

GC Gas Chromatography

GC/MS Gas Chromatography/Mass Spectrometry

LUFT Leaking Underground Fuel Tank

M Modified

MCL Maximum Contaminant Level is the highest permissible concentration of a

substance allowed in drinking water as established by the USEPA.

MDL Method Detection Limit
MPN Most Probable Number
MRL Method Reporting Limit

NA Not Applicable NC Not Calculated

NCASI National Council of the Paper Industry for Air and Stream Improvement

ND Not Detected

NIOSH National Institute for Occupational Safety and Health

PQL Practical Quantitation Limit

RCRA Resource Conservation and Recovery Act

SIM Selected Ion Monitoring

TPH Total Petroleum Hydrocarbons

tr Trace level is the concentration of an analyte that is less than the PQL but

greater than or equal to the MDL.

Analyst Summary report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Monitoring

Sample Name: TB-9954-062420-SG-001

Lab Code: R2005470-001

Sample Matrix: Water

Analysis Method Extracted/Digested By Analyzed By

8260C FNAEGLER

Sample Name: WG-9954-062420-SG-001 **Date Collected:** 06/24/20

Lab Code: R2005470-002 **Date Received:** 06/25/20

Sample Matrix: Water

Analysis Method Extracted/Digested By Analyzed By

8081B KSERCU JMISIUREWICZ 8082A KSERCU BALLGEIER

8260C FNAEGLER

8270D KSERCU JMISIUREWICZ

Sample Name: WG-9954-062420-SG-002 **Date Collected:** 06/24/20

Lab Code: R2005470-003 **Date Received:** 06/25/20

Sample Matrix: Water

8270D

Analysis Method Extracted/Digested By Analyzed By

8081B KSERCU JMISIUREWICZ 8082A KSERCU BALLGEIER

8260C FNAEGLER

KSERCU

 Sample Name:
 WG-9954-062420-SG-002
 Date Collected: 06/24/20

 Lab Code:
 R2005470-003.R01
 Date Received: 06/25/20

Sample Matrix: Water

Analysis Method Extracted/Digested By Analyzed By

8270D KSERCU JMISIUREWICZ

JMISIUREWICZ

Service Request: R2005470

Date Collected: 06/24/20

Date Received: 06/25/20

Analyst Summary report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term **Project:**

Monitoring

Sample Name: WG-9954-062420-SG-003

Lab Code: R2005470-004

Sample Matrix: Water

8270D

Analyzed By Extracted/Digested By

Service Request: R2005470

Date Collected: 06/24/20

Date Received: 06/25/20

JMISIUREWICZ

Analysis Method 8081B **KSERCU JMISIUREWICZ** 8082A **KSERCU BALLGEIER**

KSERCU

8260C **FNAEGLER**

Sample Name: WG-9954-062420-SG-004 **Date Collected:** 06/24/20

Lab Code: R2005470-005 **Date Received:** 06/25/20 Sample Matrix: Water

Analyzed By Analysis Method Extracted/Digested By 8081B **KSERCU JMISIUREWICZ** 8082A **KSERCU BALLGEIER** 8260C **FNAEGLER**

8270D **KSERCU JMISIUREWICZ**

Sample Name: WG-9954-062420-SG-005 **Date Collected:** 06/24/20

Lab Code: R2005470-006 **Date Received:** 06/25/20

Sample Matrix: Water

Analyzed By Extracted/Digested By Analysis Method

8081B **KSERCU JMISIUREWICZ** 8082A **KSERCU BALLGEIER** 8260C **FNAEGLER KSERCU** 8270D **JMISIUREWICZ**

Sample Name: WG-9954-062420-SG-006

Date Collected: 06/24/20 Lab Code: R2005470-007 **Date Received:** 06/25/20

Sample Matrix: Water

Analyzed By Analysis Method Extracted/Digested By

8081B **KSERCU JMISIUREWICZ**

Printed 7/14/2020 2:17:48 PM Superset Reference:20-0000554597 rev 00 Page 15 of 113

Analyst Summary report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Monitoring

Sample Name: WG-9954-062420-SG-006

Lab Code: R2005470-007

Sample Matrix: Water

Date Collected: 06/24/20

Date Received: 06/25/20

Service Request: R2005470

Analysis Method Extracted/Digested By Analyzed By

8082A KSERCU BALLGEIER

8260C FNAEGLER

8270D KSERCU JMISIUREWICZ

Sample Name: WG-9954-062420-SG-007 **Date Collected:** 06/24/20

Lab Code: R2005470-008 **Date Received:** 06/25/20

Sample Matrix: Water

Analysis Method Extracted/Digested By Analyzed By

8081B KSERCU JMISIUREWICZ 8082A KSERCU BALLGEIER

8260C FNAEGLER

8270D KSERCU JMISIUREWICZ



INORGANIC PREPARATION METHODS

The preparation methods associated with this report are found in these tables unless discussed in the case narrative.

Water/Liquid Matrix

Analytical Method	Preparation Method
200.7	200.2
200.8	200.2
6010C	3005A/3010A
6020A	ILM05.3
9034 Sulfide Acid Soluble	9030B
SM 4500-CN-E Residual Cyanide	SM 4500-CN-G
SM 4500-CN-E WAD Cyanide	SM 4500-CN-I

Solid/Soil/Non-Aqueous Matrix

Analytical Method	Preparation Method			
6010C	3050B			
6020A	3050B			
6010C TCLP (1311)	3005A/3010A			
extract				
6010 SPLP (1312) extract	3005A/3010A			
7199	3060A			
300.0 Anions/ 350.1/	DI extraction			
353.2/ SM 2320B/ SM				
5210B/ 9056A Anions				
For analytical methods not listed, the preparation method is the same as the analytical method reference.				



Sample Results

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com



Volatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 00:00

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 TB-9954-062420-SG-001
 Units: ug/L

 Lab Code:
 R2005470-001
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/01/20 00:37	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/01/20 00:37	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/01/20 00:37	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/01/20 00:37	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/01/20 00:37	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/01/20 00:37	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/01/20 00:37	
2-Butanone (MEK)	10 U	10	0.78	1	07/01/20 00:37	
2-Hexanone	10 U	10	0.20	1	07/01/20 00:37	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/01/20 00:37	
Acetone	10 U	10	5.0	1	07/01/20 00:37	
Benzene	5.0 U	5.0	0.20	1	07/01/20 00:37	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/01/20 00:37	
Bromoform	5.0 U	5.0	0.25	1	07/01/20 00:37	
Bromomethane	5.0 U	5.0	0.70	1	07/01/20 00:37	
Carbon Disulfide	10 U	10	0.42	1	07/01/20 00:37	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/01/20 00:37	
Chlorobenzene	5.0 U	5.0	0.20	1	07/01/20 00:37	
Chloroethane	5.0 U	5.0	0.23	1	07/01/20 00:37	
Chloroform	5.0 U	5.0	0.24	1	07/01/20 00:37	
Chloromethane	5.0 U	5.0	0.28	1	07/01/20 00:37	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/01/20 00:37	
Dichloromethane	5.0 U	5.0	0.65	1	07/01/20 00:37	
Ethylbenzene	5.0 U	5.0	0.20	1	07/01/20 00:37	
Styrene	5.0 U	5.0	0.20	1	07/01/20 00:37	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/01/20 00:37	
Toluene	5.0 U	5.0	0.20	1	07/01/20 00:37	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/01/20 00:37	
Vinyl Acetate	10 U	10	1.1	1	07/01/20 00:37	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/01/20 00:37	
Xylenes, Total	5.0 U	5.0	0.23	1	07/01/20 00:37	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/01/20 00:37	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/01/20 00:37	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/01/20 00:37	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/01/20 00:37	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 00:00

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 TB-9954-062420-SG-001
 Units: ug/L

 Lab Code:
 R2005470-001
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	90	85 - 122	07/01/20 00:37	
Dibromofluoromethane	90	89 - 119	07/01/20 00:37	
Toluene-d8	92	87 - 121	07/01/20 00:37	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 00:00 **Project:**

Monitoring

Date Received: 06/25/20 11:00 **Sample Matrix:** Water

Sample Name: TB-9954-062420-SG-001 Units: ug/L Lab Code: R2005470-001 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

Result

CAS# **Compound Identification** Q RTug/L

No Tentatively Identified Compounds Detected

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 09:10

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-001
 Units: ug/L

 Lab Code:
 R2005470-002
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/01/20 00:59	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/01/20 00:59	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/01/20 00:59	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/01/20 00:59	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/01/20 00:59	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/01/20 00:59	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/01/20 00:59	
2-Butanone (MEK)	10 U	10	0.78	1	07/01/20 00:59	
2-Hexanone	10 U	10	0.20	1	07/01/20 00:59	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/01/20 00:59	
Acetone	10 U	10	5.0	1	07/01/20 00:59	
Benzene	5.0 U	5.0	0.20	1	07/01/20 00:59	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/01/20 00:59	
Bromoform	5.0 U	5.0	0.25	1	07/01/20 00:59	
Bromomethane	5.0 U	5.0	0.70	1	07/01/20 00:59	
Carbon Disulfide	10 U	10	0.42	1	07/01/20 00:59	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/01/20 00:59	
Chlorobenzene	5.0 U	5.0	0.20	1	07/01/20 00:59	
Chloroethane	5.0 U	5.0	0.23	1	07/01/20 00:59	
Chloroform	5.0 U	5.0	0.24	1	07/01/20 00:59	
Chloromethane	5.0 U	5.0	0.28	1	07/01/20 00:59	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/01/20 00:59	
Dichloromethane	5.0 U	5.0	0.65	1	07/01/20 00:59	
Ethylbenzene	5.0 U	5.0	0.20	1	07/01/20 00:59	
Styrene	5.0 U	5.0	0.20	1	07/01/20 00:59	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/01/20 00:59	
Toluene	5.0 U	5.0	0.20	1	07/01/20 00:59	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/01/20 00:59	
Vinyl Acetate	10 U	10	1.1	1	07/01/20 00:59	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/01/20 00:59	
Xylenes, Total	5.0 U	5.0	0.23	1	07/01/20 00:59	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/01/20 00:59	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/01/20 00:59	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/01/20 00:59	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/01/20 00:59	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 09:10

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-001
 Units: ug/L

 Lab Code:
 R2005470-002
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	90	85 - 122	07/01/20 00:59	
Dibromofluoromethane	92	89 - 119	07/01/20 00:59	
Toluene-d8	93	87 - 121	07/01/20 00:59	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 09:10 **Project:**

Monitoring

Date Received: 06/25/20 11:00 **Sample Matrix:** Water

Sample Name: WG-9954-062420-SG-001 Units: ug/L Lab Code: R2005470-002 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:**

EPA 5030C

Tentatively Identified Compounds

Result CAS# **Compound Identification** Q RTug/L

No Tentatively Identified Compounds

Detected

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 09:55

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-002
 Units: ug/L

 Lab Code:
 R2005470-003
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/02/20 19:13	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/02/20 19:13	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/02/20 19:13	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/02/20 19:13	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/02/20 19:13	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/02/20 19:13	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/02/20 19:13	
2-Butanone (MEK)	10 U	10	0.78	1	07/02/20 19:13	
2-Hexanone	10 U	10	0.20	1	07/02/20 19:13	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/02/20 19:13	
Acetone	10 U	10	5.0	1	07/02/20 19:13	
Benzene	5.0 U	5.0	0.20	1	07/02/20 19:13	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/02/20 19:13	
Bromoform	5.0 U	5.0	0.25	1	07/02/20 19:13	
Bromomethane	5.0 U	5.0	0.70	1	07/02/20 19:13	
Carbon Disulfide	10 U	10	0.42	1	07/02/20 19:13	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/02/20 19:13	
Chlorobenzene	5.0 U	5.0	0.20	1	07/02/20 19:13	
Chloroethane	5.0 U	5.0	0.23	1	07/02/20 19:13	
Chloroform	5.0 U	5.0	0.24	1	07/02/20 19:13	
Chloromethane	5.0 U	5.0	0.28	1	07/02/20 19:13	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/02/20 19:13	
Dichloromethane	5.0 U	5.0	0.65	1	07/02/20 19:13	
Ethylbenzene	5.0 U	5.0	0.20	1	07/02/20 19:13	
Styrene	5.0 U	5.0	0.20	1	07/02/20 19:13	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/02/20 19:13	
Toluene	5.0 U	5.0	0.20	1	07/02/20 19:13	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/02/20 19:13	
Vinyl Acetate	10 U	10	1.1	1	07/02/20 19:13	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/02/20 19:13	
Xylenes, Total	5.0 U	5.0	0.23	1	07/02/20 19:13	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/02/20 19:13	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/02/20 19:13	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/02/20 19:13	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/02/20 19:13	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 09:55

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-002
 Units: ug/L

 Lab Code:
 R2005470-003
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	89	85 - 122	07/02/20 19:13	
Dibromofluoromethane	92	89 - 119	07/02/20 19:13	
Toluene-d8	94	87 - 121	07/02/20 19:13	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 09:55

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-002
 Units: ug/L

 Lab Code:
 R2005470-003
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 10:30

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-003
 Units: ug/L

 Lab Code:
 R2005470-004
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/02/20 19:35	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/02/20 19:35	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/02/20 19:35	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/02/20 19:35	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/02/20 19:35	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/02/20 19:35	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/02/20 19:35	
2-Butanone (MEK)	10 U	10	0.78	1	07/02/20 19:35	
2-Hexanone	10 U	10	0.20	1	07/02/20 19:35	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/02/20 19:35	
Acetone	10 U	10	5.0	1	07/02/20 19:35	
Benzene	5.0 U	5.0	0.20	1	07/02/20 19:35	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/02/20 19:35	
Bromoform	5.0 U	5.0	0.25	1	07/02/20 19:35	
Bromomethane	5.0 U	5.0	0.70	1	07/02/20 19:35	
Carbon Disulfide	10 U	10	0.42	1	07/02/20 19:35	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/02/20 19:35	
Chlorobenzene	5.0 U	5.0	0.20	1	07/02/20 19:35	
Chloroethane	5.0 U	5.0	0.23	1	07/02/20 19:35	
Chloroform	5.0 U	5.0	0.24	1	07/02/20 19:35	
Chloromethane	5.0 U	5.0	0.28	1	07/02/20 19:35	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/02/20 19:35	
Dichloromethane	5.0 U	5.0	0.65	1	07/02/20 19:35	
Ethylbenzene	5.0 U	5.0	0.20	1	07/02/20 19:35	
Styrene	5.0 U	5.0	0.20	1	07/02/20 19:35	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/02/20 19:35	
Toluene	5.0 U	5.0	0.20	1	07/02/20 19:35	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/02/20 19:35	
Vinyl Acetate	10 U	10	1.1	1	07/02/20 19:35	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/02/20 19:35	
Xylenes, Total	5.0 U	5.0	0.23	1	07/02/20 19:35	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/02/20 19:35	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/02/20 19:35	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/02/20 19:35	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/02/20 19:35	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 10:30

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-003
 Units: ug/L

 Lab Code:
 R2005470-004
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	91	85 - 122	07/02/20 19:35	
Dibromofluoromethane	94	89 - 119	07/02/20 19:35	
Toluene-d8	96	87 - 121	07/02/20 19:35	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 10:30

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-003
 Units: ug/L

 Lab Code:
 R2005470-004
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

CAS# Compound Identification RT ug/L

No Tentatively Identified Compounds

Detected

Q

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 11:05

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-004
 Units: ug/L

 Lab Code:
 R2005470-005
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/02/20 19:58	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/02/20 19:58	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/02/20 19:58	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/02/20 19:58	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/02/20 19:58	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/02/20 19:58	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/02/20 19:58	
2-Butanone (MEK)	10 U	10	0.78	1	07/02/20 19:58	
2-Hexanone	10 U	10	0.20	1	07/02/20 19:58	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/02/20 19:58	
Acetone	10 U	10	5.0	1	07/02/20 19:58	
Benzene	5.0 U	5.0	0.20	1	07/02/20 19:58	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/02/20 19:58	
Bromoform	5.0 U	5.0	0.25	1	07/02/20 19:58	
Bromomethane	5.0 U	5.0	0.70	1	07/02/20 19:58	
Carbon Disulfide	0.83 J	10	0.42	1	07/02/20 19:58	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/02/20 19:58	
Chlorobenzene	5.0 U	5.0	0.20	1	07/02/20 19:58	
Chloroethane	5.0 U	5.0	0.23	1	07/02/20 19:58	
Chloroform	5.0 U	5.0	0.24	1	07/02/20 19:58	
Chloromethane	5.0 U	5.0	0.28	1	07/02/20 19:58	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/02/20 19:58	
Dichloromethane	5.0 U	5.0	0.65	1	07/02/20 19:58	
Ethylbenzene	5.0 U	5.0	0.20	1	07/02/20 19:58	
Styrene	5.0 U	5.0	0.20	1	07/02/20 19:58	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/02/20 19:58	
Toluene	5.0 U	5.0	0.20	1	07/02/20 19:58	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/02/20 19:58	
Vinyl Acetate	10 U	10	1.1	1	07/02/20 19:58	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/02/20 19:58	
Xylenes, Total	5.0 U	5.0	0.23	1	07/02/20 19:58	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/02/20 19:58	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/02/20 19:58	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/02/20 19:58	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/02/20 19:58	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 11:05

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-004
 Units: ug/L

 Lab Code:
 R2005470-005
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	90	85 - 122	07/02/20 19:58	
Dibromofluoromethane	94	89 - 119	07/02/20 19:58	
Toluene-d8	95	87 - 121	07/02/20 19:58	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 11:05

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-004
 Units: ug/L

 Lab Code:
 R2005470-005
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

Result

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 11:50

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-005
 Units: ug/L

 Lab Code:
 R2005470-006
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/01/20 02:27	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/01/20 02:27	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/01/20 02:27	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/01/20 02:27	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/01/20 02:27	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/01/20 02:27	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/01/20 02:27	
2-Butanone (MEK)	10 U	10	0.78	1	07/01/20 02:27	
2-Hexanone	10 U	10	0.20	1	07/01/20 02:27	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/01/20 02:27	
Acetone	10 U	10	5.0	1	07/01/20 02:27	
Benzene	5.0 U	5.0	0.20	1	07/01/20 02:27	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/01/20 02:27	
Bromoform	5.0 U	5.0	0.25	1	07/01/20 02:27	
Bromomethane	5.0 U	5.0	0.70	1	07/01/20 02:27	
Carbon Disulfide	10 U	10	0.42	1	07/01/20 02:27	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/01/20 02:27	
Chlorobenzene	5.0 U	5.0	0.20	1	07/01/20 02:27	
Chloroethane	5.0 U	5.0	0.23	1	07/01/20 02:27	
Chloroform	5.0 U	5.0	0.24	1	07/01/20 02:27	
Chloromethane	5.0 U	5.0	0.28	1	07/01/20 02:27	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/01/20 02:27	
Dichloromethane	5.0 U	5.0	0.65	1	07/01/20 02:27	
Ethylbenzene	5.0 U	5.0	0.20	1	07/01/20 02:27	
Styrene	5.0 U	5.0	0.20	1	07/01/20 02:27	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/01/20 02:27	
Toluene	5.0 U	5.0	0.20	1	07/01/20 02:27	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/01/20 02:27	
Vinyl Acetate	10 U	10	1.1	1	07/01/20 02:27	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/01/20 02:27	
Xylenes, Total	5.0 U	5.0	0.23	1	07/01/20 02:27	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/01/20 02:27	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/01/20 02:27	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/01/20 02:27	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/01/20 02:27	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 11:50

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-005
 Units: ug/L

 Lab Code:
 R2005470-006
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	93	85 - 122	07/01/20 02:27	
Dibromofluoromethane	93	89 - 119	07/01/20 02:27	
Toluene-d8	94	87 - 121	07/01/20 02:27	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 11:50

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-005
 Units: ug/L

 Lab Code:
 R2005470-006
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

Result

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 12:45

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-006
 Units: ug/L

 Lab Code:
 R2005470-007
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/01/20 02:50	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/01/20 02:50	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/01/20 02:50	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/01/20 02:50	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/01/20 02:50	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/01/20 02:50	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/01/20 02:50	
2-Butanone (MEK)	10 U	10	0.78	1	07/01/20 02:50	
2-Hexanone	10 U	10	0.20	1	07/01/20 02:50	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/01/20 02:50	
Acetone	10 U	10	5.0	1	07/01/20 02:50	
Benzene	5.0 U	5.0	0.20	1	07/01/20 02:50	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/01/20 02:50	
Bromoform	5.0 U	5.0	0.25	1	07/01/20 02:50	
Bromomethane	5.0 U	5.0	0.70	1	07/01/20 02:50	
Carbon Disulfide	10 U	10	0.42	1	07/01/20 02:50	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/01/20 02:50	
Chlorobenzene	5.0 U	5.0	0.20	1	07/01/20 02:50	
Chloroethane	5.0 U	5.0	0.23	1	07/01/20 02:50	
Chloroform	5.0 U	5.0	0.24	1	07/01/20 02:50	
Chloromethane	5.0 U	5.0	0.28	1	07/01/20 02:50	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/01/20 02:50	
Dichloromethane	5.0 U	5.0	0.65	1	07/01/20 02:50	
Ethylbenzene	5.0 U	5.0	0.20	1	07/01/20 02:50	
Styrene	5.0 U	5.0	0.20	1	07/01/20 02:50	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/01/20 02:50	
Toluene	5.0 U	5.0	0.20	1	07/01/20 02:50	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/01/20 02:50	
Vinyl Acetate	10 U	10	1.1	1	07/01/20 02:50	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/01/20 02:50	
Xylenes, Total	5.0 U	5.0	0.23	1	07/01/20 02:50	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/01/20 02:50	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/01/20 02:50	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/01/20 02:50	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/01/20 02:50	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 12:45

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-006
 Units: ug/L

 Lab Code:
 R2005470-007
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	89	85 - 122	07/01/20 02:50	
Dibromofluoromethane	89	89 - 119	07/01/20 02:50	
Toluene-d8	92	87 - 121	07/01/20 02:50	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 12:45

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-006
 Units: ug/L

 Lab Code:
 R2005470-007
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

Result

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 13:35

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-007
 Units: ug/L

 Lab Code:
 R2005470-008
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/01/20 03:12	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/01/20 03:12	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/01/20 03:12	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/01/20 03:12	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/01/20 03:12	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/01/20 03:12	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/01/20 03:12	
2-Butanone (MEK)	10 U	10	0.78	1	07/01/20 03:12	
2-Hexanone	10 U	10	0.20	1	07/01/20 03:12	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/01/20 03:12	
Acetone	10 U	10	5.0	1	07/01/20 03:12	
Benzene	5.0 U	5.0	0.20	1	07/01/20 03:12	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/01/20 03:12	
Bromoform	5.0 U	5.0	0.25	1	07/01/20 03:12	
Bromomethane	5.0 U	5.0	0.70	1	07/01/20 03:12	
Carbon Disulfide	10 U	10	0.42	1	07/01/20 03:12	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/01/20 03:12	
Chlorobenzene	5.0 U	5.0	0.20	1	07/01/20 03:12	
Chloroethane	5.0 U	5.0	0.23	1	07/01/20 03:12	
Chloroform	5.0 U	5.0	0.24	1	07/01/20 03:12	
Chloromethane	5.0 U	5.0	0.28	1	07/01/20 03:12	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/01/20 03:12	
Dichloromethane	5.0 U	5.0	0.65	1	07/01/20 03:12	
Ethylbenzene	5.0 U	5.0	0.20	1	07/01/20 03:12	
Styrene	5.0 U	5.0	0.20	1	07/01/20 03:12	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/01/20 03:12	
Toluene	5.0 U	5.0	0.20	1	07/01/20 03:12	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/01/20 03:12	
Vinyl Acetate	10 U	10	1.1	1	07/01/20 03:12	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/01/20 03:12	
Xylenes, Total	5.0 U	5.0	0.23	1	07/01/20 03:12	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/01/20 03:12	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/01/20 03:12	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/01/20 03:12	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/01/20 03:12	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 13:35

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-007
 Units: ug/L

 Lab Code:
 R2005470-008
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	90	85 - 122	07/01/20 03:12	
Dibromofluoromethane	90	89 - 119	07/01/20 03:12	
Toluene-d8	92	87 - 121	07/01/20 03:12	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 13:35

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-007
 Units: ug/L

 Lab Code:
 R2005470-008
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

Result

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected



Semivolatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 09:10

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-001
 Units: ug/L

 Lab Code:
 R2005470-002
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	06/30/20 18:29	6/29/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	06/30/20 18:29	6/29/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	06/30/20 18:29	6/29/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	06/30/20 18:29	6/29/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	06/30/20 18:29	6/29/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	06/30/20 18:29	6/29/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	06/30/20 18:29	6/29/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	06/30/20 18:29	6/29/20	
2,4-Dinitrophenol	45 U	45	19	1	06/30/20 18:29	6/29/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	06/30/20 18:29	6/29/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	06/30/20 18:29	6/29/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	06/30/20 18:29	6/29/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	06/30/20 18:29	6/29/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	06/30/20 18:29	6/29/20	
2-Methylphenol	9.1 U	9.1	0.91	1	06/30/20 18:29	6/29/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	06/30/20 18:29	6/29/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	06/30/20 18:29	6/29/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	06/30/20 18:29	6/29/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	06/30/20 18:29	6/29/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	06/30/20 18:29	6/29/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	06/30/20 18:29	6/29/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	06/30/20 18:29	6/29/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	06/30/20 18:29	6/29/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	06/30/20 18:29	6/29/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	06/30/20 18:29	6/29/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	06/30/20 18:29	6/29/20	
4-Nitrophenol	45 U	45	5.8	1	06/30/20 18:29	6/29/20	
Acenaphthene	9.1 U	9.1	1.3	1	06/30/20 18:29	6/29/20	
Acenaphthylene	9.1 U	9.1	1.3	1	06/30/20 18:29	6/29/20	
Anthracene	9.1 U	9.1	1.2	1	06/30/20 18:29	6/29/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	06/30/20 18:29	6/29/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	06/30/20 18:29	6/29/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	06/30/20 18:29	6/29/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	06/30/20 18:29	6/29/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	06/30/20 18:29	6/29/20	
Benzoic Acid	91 U	91	33	1	06/30/20 18:29	6/29/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	06/30/20 18:29	6/29/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	06/30/20 18:29	6/29/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	06/30/20 18:29	6/29/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	06/30/20 18:29	6/29/20	
Bis(2-ethylhexyl) Phthalate	1.7 J	9.1	0.91	1	06/30/20 18:29	6/29/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	06/30/20 18:29	6/29/20	
Chrysene	9.1 U	9.1	1.1	1	06/30/20 18:29	6/29/20	

Printed 7/14/2020 2:18:02 PM

Superset Reference:20-0000554597 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 09:10

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-001
 Units: ug/L

 Lab Code:
 R2005470-002
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	06/30/20 18:29	6/29/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	06/30/20 18:29	6/29/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	06/30/20 18:29	6/29/20	
Dibenzofuran	9.1 U	9.1	1.3	1	06/30/20 18:29	6/29/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	06/30/20 18:29	6/29/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	06/30/20 18:29	6/29/20	
Fluoranthene	9.1 U	9.1	1.4	1	06/30/20 18:29	6/29/20	
Fluorene	9.1 U	9.1	1.2	1	06/30/20 18:29	6/29/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	06/30/20 18:29	6/29/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	06/30/20 18:29	6/29/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	06/30/20 18:29	6/29/20	
Hexachloroethane	9.1 U	9.1	0.96	1	06/30/20 18:29	6/29/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	06/30/20 18:29	6/29/20	
Isophorone	9.1 U	9.1	1.3	1	06/30/20 18:29	6/29/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	06/30/20 18:29	6/29/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	06/30/20 18:29	6/29/20	
Naphthalene	9.1 U	9.1	1.1	1	06/30/20 18:29	6/29/20	
Nitrobenzene	9.1 U	9.1	1.4	1	06/30/20 18:29	6/29/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	06/30/20 18:29	6/29/20	
Phenanthrene	9.1 U	9.1	1.3	1	06/30/20 18:29	6/29/20	
Phenol	9.1 U	9.1	0.91	1	06/30/20 18:29	6/29/20	
Pyrene	9.1 U	9.1	1.3	1	06/30/20 18:29	6/29/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	73	35 - 141	06/30/20 18:29	
2-Fluorobiphenyl	53	31 - 118	06/30/20 18:29	
2-Fluorophenol	35	10 - 105	06/30/20 18:29	
Nitrobenzene-d5	43	31 - 110	06/30/20 18:29	
Phenol-d6	24	10 - 107	06/30/20 18:29	
p-Terphenyl-d14	84	10 - 165	06/30/20 18:29	

Tentatively Identified Compounds

			Result	
CAS#	Compound Identification	RT	ug/L	Q
1000126-80-6	Azacyclohexan-2-carboxylic acid, 1	8.33	5.1	JN

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470 **Date Collected:** 06/24/20 09:55

Love Canal:292-402-D02-3100/9954 Annual Long Term **Project:**

Monitoring

Date Received: 06/25/20 11:00 **Sample Matrix:** Water

Sample Name: Units: ug/L WG-9954-062420-SG-002 Basis: NA Lab Code: R2005470-003

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	06/30/20 18:57	6/29/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	06/30/20 18:57	6/29/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	06/30/20 18:57	6/29/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	06/30/20 18:57	6/29/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	06/30/20 18:57	6/29/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	06/30/20 18:57	6/29/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	06/30/20 18:57	6/29/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	06/30/20 18:57	6/29/20	
2,4-Dinitrophenol	45 U	45	19	1	06/30/20 18:57	6/29/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	06/30/20 18:57	6/29/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	06/30/20 18:57	6/29/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	06/30/20 18:57	6/29/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	06/30/20 18:57	6/29/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	06/30/20 18:57	6/29/20	
2-Methylphenol	9.1 U	9.1	0.91	1	06/30/20 18:57	6/29/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	06/30/20 18:57	6/29/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	06/30/20 18:57	6/29/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	06/30/20 18:57	6/29/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	06/30/20 18:57	6/29/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	06/30/20 18:57	6/29/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	06/30/20 18:57	6/29/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	06/30/20 18:57	6/29/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	06/30/20 18:57	6/29/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	06/30/20 18:57	6/29/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	06/30/20 18:57	6/29/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	06/30/20 18:57	6/29/20	
4-Nitrophenol	45 U	45	5.8	1	06/30/20 18:57	6/29/20	
Acenaphthene	9.1 U	9.1	1.3	1	06/30/20 18:57	6/29/20	
Acenaphthylene	9.1 U	9.1	1.3	1	06/30/20 18:57	6/29/20	
Anthracene	9.1 U	9.1	1.2	1	06/30/20 18:57	6/29/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	06/30/20 18:57	6/29/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	06/30/20 18:57	6/29/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	06/30/20 18:57	6/29/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	06/30/20 18:57	6/29/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	06/30/20 18:57	6/29/20	
Benzoic Acid	91 U	91	33	1	06/30/20 18:57	6/29/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	06/30/20 18:57	6/29/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	06/30/20 18:57	6/29/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	06/30/20 18:57	6/29/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	06/30/20 18:57	6/29/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	06/30/20 18:57	6/29/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	06/30/20 18:57	6/29/20	
Chrysene	9.1 U	9.1	1.1	1	06/30/20 18:57	6/29/20	

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Superset Reference:20-0000554597 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470 **Date Collected:** 06/24/20 09:55

Love Canal:292-402-D02-3100/9954 Annual Long Term **Project:**

Monitoring

Sample Matrix: Water **Date Received:** 06/25/20 11:00

Sample Name: WG-9954-062420-SG-002 Units: ug/L Lab Code: R2005470-003 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	06/30/20 18:57	6/29/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	06/30/20 18:57	6/29/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	06/30/20 18:57	6/29/20	
Dibenzofuran	9.1 U	9.1	1.3	1	06/30/20 18:57	6/29/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	06/30/20 18:57	6/29/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	06/30/20 18:57	6/29/20	
Fluoranthene	9.1 U	9.1	1.4	1	06/30/20 18:57	6/29/20	
Fluorene	9.1 U	9.1	1.2	1	06/30/20 18:57	6/29/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	06/30/20 18:57	6/29/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	06/30/20 18:57	6/29/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	06/30/20 18:57	6/29/20	
Hexachloroethane	9.1 U	9.1	0.96	1	06/30/20 18:57	6/29/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	06/30/20 18:57	6/29/20	
Isophorone	9.1 U	9.1	1.3	1	06/30/20 18:57	6/29/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	06/30/20 18:57	6/29/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	06/30/20 18:57	6/29/20	
Naphthalene	9.1 U	9.1	1.1	1	06/30/20 18:57	6/29/20	
Nitrobenzene	9.1 U	9.1	1.4	1	06/30/20 18:57	6/29/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	06/30/20 18:57	6/29/20	
Phenanthrene	9.1 U	9.1	1.3	1	06/30/20 18:57	6/29/20	
Phenol	9.1 U	9.1	0.91	1	06/30/20 18:57	6/29/20	
Pyrene	9.1 U	9.1	1.3	1	06/30/20 18:57	6/29/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	27 *	35 - 141	06/30/20 18:57	*
2-Fluorobiphenyl	26 *	31 - 118	06/30/20 18:57	*
2-Fluorophenol	18	10 - 105	06/30/20 18:57	
Nitrobenzene-d5	22 *	31 - 110	06/30/20 18:57	*
Phenol-d6	12	10 - 107	06/30/20 18:57	
p-Terphenyl-d14	35	10 - 165	06/30/20 18:57	

Tentatively Identified Compounds

			Result	
CAS#	Compound Identification	RT	ug/L	Q

No Tentatively Identified Compounds

Detected

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Love Canal:292-402-D02-3100/9954 Annual Long Term **Project:**

Date Collected: 06/24/20 09:55

Monitoring **Sample Matrix:** Water

Date Received: 06/25/20 11:00

Sample Name: WG-9954-062420-SG-002 Lab Code:

Units: ug/L R2005470-003 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 16:22	7/2/20	*
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 16:22	7/2/20	*
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/06/20 16:22	7/2/20	*
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 16:22	7/2/20	*
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/06/20 16:22	7/2/20	*
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/06/20 16:22	7/2/20	*
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/06/20 16:22	7/2/20	*
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/06/20 16:22	7/2/20	*
2,4-Dinitrophenol	45 U	45	19	1	07/06/20 16:22	7/2/20	*
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/06/20 16:22	7/2/20	*
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/06/20 16:22	7/2/20	*
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/06/20 16:22	7/2/20	*
2-Chlorophenol	9.1 U	9.1	0.97	1	07/06/20 16:22	7/2/20	*
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/06/20 16:22	7/2/20	*
2-Methylphenol	9.1 U	9.1	0.91	1	07/06/20 16:22	7/2/20	*
2-Nitroaniline	9.1 U	9.1	1.3	1	07/06/20 16:22	7/2/20	*
2-Nitrophenol	9.1 U	9.1	1.4	1	07/06/20 16:22	7/2/20	*
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/06/20 16:22	7/2/20	*
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/06/20 16:22	7/2/20	*
3-Nitroaniline	9.1 U	9.1	2.3	1	07/06/20 16:22	7/2/20	*
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/06/20 16:22	7/2/20	*
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/06/20 16:22	7/2/20	*
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/06/20 16:22	7/2/20	*
4-Chloroaniline	9.1 U	9.1	0.91	1	07/06/20 16:22	7/2/20	*
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/06/20 16:22	7/2/20	*
4-Nitroaniline	9.1 U	9.1	2.5	1	07/06/20 16:22	7/2/20	*
4-Nitrophenol	45 U	45	5.8	1	07/06/20 16:22	7/2/20	*
Acenaphthene	9.1 U	9.1	1.3	1	07/06/20 16:22	7/2/20	*
Acenaphthylene	9.1 U	9.1	1.3	1	07/06/20 16:22	7/2/20	*
Anthracene	9.1 U	9.1	1.2	1	07/06/20 16:22	7/2/20	*
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/06/20 16:22	7/2/20	*
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/06/20 16:22	7/2/20	*
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/06/20 16:22	7/2/20	*
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/06/20 16:22	7/2/20	*
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/06/20 16:22	7/2/20	*
Benzoic Acid	91 U	91	33	1	07/06/20 16:22	7/2/20	*
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/06/20 16:22	7/2/20	*
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/06/20 16:22	7/2/20	*
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/06/20 16:22	7/2/20	*
Bis(2-chloroethyl) Ether	9.1 U 9.1 U	9.1 9.1	1.8	1	07/06/20 16:22	7/2/20	*
Bis(2-ethylhexyl) Phthalate	9.1 U 9.1 U	9.1	0.91	1	07/06/20 16:22	7/2/20	*
	9.1 U 9.1 U	9.1 9.1	1.3	1	07/06/20 16:22	7/2/20 7/2/20	*
Butyl Benzyl Phthalate	9.1 U 9.1 U	9.1 9.1	1.3			7/2/20	*
Chrysene	9.1 U	9.1	1.1	1	07/06/20 16:22	1/2/20	~

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Superset Reference:20-0000554597 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 09:55

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-002
 Units: ug/L

 Lab Code:
 R2005470-003
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/06/20 16:22 7/2/20	*
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/06/20 16:22 7/2/20	*
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/06/20 16:22 7/2/20	*
Dibenzofuran	9.1 U	9.1	1.3	1	07/06/20 16:22 7/2/20	*
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/06/20 16:22 7/2/20	*
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/06/20 16:22 7/2/20	*
Fluoranthene	9.1 U	9.1	1.4	1	07/06/20 16:22 7/2/20	*
Fluorene	9.1 U	9.1	1.2	1	07/06/20 16:22 7/2/20	*
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/06/20 16:22 7/2/20	*
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/06/20 16:22 7/2/20	*
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/06/20 16:22 7/2/20	*
Hexachloroethane	9.1 U	9.1	0.96	1	07/06/20 16:22 7/2/20	*
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/06/20 16:22 7/2/20	*
Isophorone	9.1 U	9.1	1.3	1	07/06/20 16:22 7/2/20	*
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/06/20 16:22 7/2/20	*
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/06/20 16:22 7/2/20	*
Naphthalene	9.1 U	9.1	1.1	1	07/06/20 16:22 7/2/20	*
Nitrobenzene	9.1 U	9.1	1.4	1	07/06/20 16:22 7/2/20	*
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/06/20 16:22 7/2/20	*
Phenanthrene	9.1 U	9.1	1.3	1	07/06/20 16:22 7/2/20	*
Phenol	9.1 U	9.1	0.91	1	07/06/20 16:22 7/2/20	*
Pyrene	9.1 U	9.1	1.3	1	07/06/20 16:22 7/2/20	*

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	105	35 - 141	07/06/20 16:22	
2-Fluorobiphenyl	79	31 - 118	07/06/20 16:22	
2-Fluorophenol	51	10 - 105	07/06/20 16:22	
Nitrobenzene-d5	85	31 - 110	07/06/20 16:22	
Phenol-d6	32	10 - 107	07/06/20 16:22	
p-Terphenyl-d14	56	10 - 165	07/06/20 16:22	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown	12.07	4.7	J	
	unknown hydrocarbon	12.69	4.9	J	
	unknown hydrocarbon	13.36	6.5	J	
	unknown hydrocarbon	14.07	4.3	J	
	unknown	14.84	4.1	J	
	unknown	15.65	3.7	J	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 10:30

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-003
 Units: ug/L

 Lab Code:
 R2005470-004
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	06/30/20 19:26	6/29/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	06/30/20 19:26	6/29/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	06/30/20 19:26	6/29/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	06/30/20 19:26	6/29/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	06/30/20 19:26	6/29/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	06/30/20 19:26	6/29/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	06/30/20 19:26	6/29/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	06/30/20 19:26	6/29/20	
2,4-Dinitrophenol	45 U	45	19	1	06/30/20 19:26	6/29/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	06/30/20 19:26	6/29/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	06/30/20 19:26	6/29/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	06/30/20 19:26	6/29/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	06/30/20 19:26	6/29/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	06/30/20 19:26	6/29/20	
2-Methylphenol	9.1 U	9.1	0.91	1	06/30/20 19:26	6/29/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	06/30/20 19:26	6/29/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	06/30/20 19:26	6/29/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	06/30/20 19:26	6/29/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	06/30/20 19:26	6/29/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	06/30/20 19:26	6/29/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	06/30/20 19:26	6/29/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	06/30/20 19:26	6/29/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	06/30/20 19:26	6/29/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	06/30/20 19:26	6/29/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	06/30/20 19:26	6/29/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	06/30/20 19:26	6/29/20	
4-Nitrophenol	45 U	45	5.8	1	06/30/20 19:26	6/29/20	
Acenaphthene	9.1 U	9.1	1.3	1	06/30/20 19:26	6/29/20	
Acenaphthylene	9.1 U	9.1	1.3	1	06/30/20 19:26	6/29/20	
Anthracene	9.1 U	9.1	1.2	1	06/30/20 19:26	6/29/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	06/30/20 19:26	6/29/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	06/30/20 19:26	6/29/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	06/30/20 19:26	6/29/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	06/30/20 19:26	6/29/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	06/30/20 19:26	6/29/20	
Benzoic Acid	91 U	91	33	1	06/30/20 19:26	6/29/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	06/30/20 19:26	6/29/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	06/30/20 19:26	6/29/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	06/30/20 19:26	6/29/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	06/30/20 19:26	6/29/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	06/30/20 19:26	6/29/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	06/30/20 19:26	6/29/20	
Chrysene	9.1 U	9.1	1.1	1	06/30/20 19:26	6/29/20	

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Superset Reference: 20-0000554597 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 10:30

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-003
 Units: ug/L

 Lab Code:
 R2005470-004
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed I	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	06/30/20 19:26	6/29/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	06/30/20 19:26	6/29/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	06/30/20 19:26	6/29/20	
Dibenzofuran	9.1 U	9.1	1.3	1	06/30/20 19:26	6/29/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	06/30/20 19:26	6/29/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	06/30/20 19:26	6/29/20	
Fluoranthene	9.1 U	9.1	1.4	1	06/30/20 19:26	6/29/20	
Fluorene	9.1 U	9.1	1.2	1	06/30/20 19:26	6/29/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	06/30/20 19:26	6/29/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	06/30/20 19:26	6/29/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	06/30/20 19:26	6/29/20	
Hexachloroethane	9.1 U	9.1	0.96	1	06/30/20 19:26	6/29/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	06/30/20 19:26	6/29/20	
Isophorone	9.1 U	9.1	1.3	1	06/30/20 19:26	6/29/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	06/30/20 19:26	6/29/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	06/30/20 19:26	6/29/20	
Naphthalene	9.1 U	9.1	1.1	1	06/30/20 19:26	6/29/20	
Nitrobenzene	9.1 U	9.1	1.4	1	06/30/20 19:26	6/29/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	06/30/20 19:26	6/29/20	
Phenanthrene	9.1 U	9.1	1.3	1	06/30/20 19:26	6/29/20	
Phenol	9.1 U	9.1	0.91	1	06/30/20 19:26	6/29/20	
Pyrene	9.1 U	9.1	1.3	1	06/30/20 19:26	6/29/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	82	35 - 141	06/30/20 19:26	
2-Fluorobiphenyl	71	31 - 118	06/30/20 19:26	
2-Fluorophenol	46	10 - 105	06/30/20 19:26	
Nitrobenzene-d5	69	31 - 110	06/30/20 19:26	
Phenol-d6	31	10 - 107	06/30/20 19:26	
p-Terphenyl-d14	97	10 - 165	06/30/20 19:26	

Tentatively Identified Compounds

			Result	
CAS#	Compound Identification	RT	ug/L	Q
	unknown hydrocarbon	13.16	3.7	J

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 11:05

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-004
 Units: ug/L

 Lab Code:
 R2005470-005
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	06/30/20 19:54	6/29/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	06/30/20 19:54	6/29/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	06/30/20 19:54	6/29/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	06/30/20 19:54	6/29/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	06/30/20 19:54	6/29/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	06/30/20 19:54	6/29/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	06/30/20 19:54	6/29/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	06/30/20 19:54	6/29/20	
2,4-Dinitrophenol	45 U	45	19	1	06/30/20 19:54	6/29/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	06/30/20 19:54	6/29/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	06/30/20 19:54	6/29/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	06/30/20 19:54	6/29/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	06/30/20 19:54	6/29/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	06/30/20 19:54	6/29/20	
2-Methylphenol	9.1 U	9.1	0.91	1	06/30/20 19:54	6/29/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	06/30/20 19:54	6/29/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	06/30/20 19:54	6/29/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	06/30/20 19:54	6/29/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	06/30/20 19:54	6/29/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	06/30/20 19:54	6/29/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	06/30/20 19:54	6/29/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	06/30/20 19:54	6/29/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	06/30/20 19:54	6/29/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	06/30/20 19:54	6/29/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	06/30/20 19:54	6/29/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	06/30/20 19:54	6/29/20	
4-Nitrophenol	45 U	45	5.8	1	06/30/20 19:54	6/29/20	
Acenaphthene	9.1 U	9.1	1.3	1	06/30/20 19:54	6/29/20	
Acenaphthylene	9.1 U	9.1	1.3	1	06/30/20 19:54	6/29/20	
Anthracene	9.1 U	9.1	1.2	1	06/30/20 19:54	6/29/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	06/30/20 19:54	6/29/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	06/30/20 19:54	6/29/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	06/30/20 19:54	6/29/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	06/30/20 19:54	6/29/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	06/30/20 19:54	6/29/20	
Benzoic Acid	91 U	91	33	1	06/30/20 19:54	6/29/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	06/30/20 19:54	6/29/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	06/30/20 19:54	6/29/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	06/30/20 19:54	6/29/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	06/30/20 19:54	6/29/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	06/30/20 19:54	6/29/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	06/30/20 19:54	6/29/20	
Chrysene	9.1 U	9.1	1.1	1	06/30/20 19:54	6/29/20	

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Superset Reference: 20-0000554597 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 11:05

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-004
 Units: ug/L

 Lab Code:
 R2005470-005
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	06/30/20 19:54	6/29/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	06/30/20 19:54	6/29/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	06/30/20 19:54	6/29/20	
Dibenzofuran	9.1 U	9.1	1.3	1	06/30/20 19:54	6/29/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	06/30/20 19:54	6/29/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	06/30/20 19:54	6/29/20	
Fluoranthene	9.1 U	9.1	1.4	1	06/30/20 19:54	6/29/20	
Fluorene	9.1 U	9.1	1.2	1	06/30/20 19:54	6/29/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	06/30/20 19:54	6/29/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	06/30/20 19:54	6/29/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	06/30/20 19:54	6/29/20	
Hexachloroethane	9.1 U	9.1	0.96	1	06/30/20 19:54	6/29/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	06/30/20 19:54	6/29/20	
Isophorone	9.1 U	9.1	1.3	1	06/30/20 19:54	6/29/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	06/30/20 19:54	6/29/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	06/30/20 19:54	6/29/20	
Naphthalene	9.1 U	9.1	1.1	1	06/30/20 19:54	6/29/20	
Nitrobenzene	9.1 U	9.1	1.4	1	06/30/20 19:54	6/29/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	06/30/20 19:54	6/29/20	
Phenanthrene	9.1 U	9.1	1.3	1	06/30/20 19:54	6/29/20	
Phenol	9.1 U	9.1	0.91	1	06/30/20 19:54	6/29/20	
Pyrene	9.1 U	9.1	1.3	1	06/30/20 19:54	6/29/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	68	35 - 141	06/30/20 19:54	
2-Fluorobiphenyl	58	31 - 118	06/30/20 19:54	
2-Fluorophenol	34	10 - 105	06/30/20 19:54	
Nitrobenzene-d5	51	31 - 110	06/30/20 19:54	
Phenol-d6	23	10 - 107	06/30/20 19:54	
p-Terphenyl-d14	85	10 - 165	06/30/20 19:54	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	

No Tentatively Identified Compounds

Detected

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 11:50

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-005
 Units: ug/L

 Lab Code:
 R2005470-006
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	06/30/20 20:23	6/29/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	06/30/20 20:23	6/29/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	06/30/20 20:23	6/29/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	06/30/20 20:23	6/29/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	06/30/20 20:23	6/29/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	06/30/20 20:23	6/29/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	06/30/20 20:23	6/29/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	06/30/20 20:23	6/29/20	
2,4-Dinitrophenol	45 U	45	19	1	06/30/20 20:23	6/29/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	06/30/20 20:23	6/29/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	06/30/20 20:23	6/29/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	06/30/20 20:23	6/29/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	06/30/20 20:23	6/29/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	06/30/20 20:23	6/29/20	
2-Methylphenol	9.1 U	9.1	0.91	1	06/30/20 20:23	6/29/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	06/30/20 20:23	6/29/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	06/30/20 20:23	6/29/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	06/30/20 20:23	6/29/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	06/30/20 20:23	6/29/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	06/30/20 20:23	6/29/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	06/30/20 20:23	6/29/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	06/30/20 20:23	6/29/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	06/30/20 20:23	6/29/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	06/30/20 20:23	6/29/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	06/30/20 20:23	6/29/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	06/30/20 20:23	6/29/20	
4-Nitrophenol	45 U	45	5.8	1	06/30/20 20:23	6/29/20	
Acenaphthene	9.1 U	9.1	1.3	1	06/30/20 20:23	6/29/20	
Acenaphthylene	9.1 U	9.1	1.3	1	06/30/20 20:23	6/29/20	
Anthracene	9.1 U	9.1	1.2	1	06/30/20 20:23	6/29/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	06/30/20 20:23	6/29/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	06/30/20 20:23	6/29/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	06/30/20 20:23	6/29/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	06/30/20 20:23	6/29/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	06/30/20 20:23	6/29/20	
Benzoic Acid	91 U	91	33	1	06/30/20 20:23	6/29/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	06/30/20 20:23	6/29/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	06/30/20 20:23	6/29/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	06/30/20 20:23	6/29/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	06/30/20 20:23	6/29/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	06/30/20 20:23	6/29/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	06/30/20 20:23	6/29/20	
Chrysene	9.1 U	9.1	1.1	1	06/30/20 20:23	6/29/20	

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Superset Reference:20-0000554597 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 11:50

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-005
 Units: ug/L

 Lab Code:
 R2005470-006
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	06/30/20 20:23	6/29/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	06/30/20 20:23	6/29/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	06/30/20 20:23	6/29/20	
Dibenzofuran	9.1 U	9.1	1.3	1	06/30/20 20:23	6/29/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	06/30/20 20:23	6/29/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	06/30/20 20:23	6/29/20	
Fluoranthene	9.1 U	9.1	1.4	1	06/30/20 20:23	6/29/20	
Fluorene	9.1 U	9.1	1.2	1	06/30/20 20:23	6/29/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	06/30/20 20:23	6/29/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	06/30/20 20:23	6/29/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	06/30/20 20:23	6/29/20	
Hexachloroethane	9.1 U	9.1	0.96	1	06/30/20 20:23	6/29/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	06/30/20 20:23	6/29/20	
Isophorone	9.1 U	9.1	1.3	1	06/30/20 20:23	6/29/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	06/30/20 20:23	6/29/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	06/30/20 20:23	6/29/20	
Naphthalene	9.1 U	9.1	1.1	1	06/30/20 20:23	6/29/20	
Nitrobenzene	9.1 U	9.1	1.4	1	06/30/20 20:23	6/29/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	06/30/20 20:23	6/29/20	
Phenanthrene	9.1 U	9.1	1.3	1	06/30/20 20:23	6/29/20	
Phenol	9.1 U	9.1	0.91	1	06/30/20 20:23	6/29/20	
Pyrene	9.1 U	9.1	1.3	1	06/30/20 20:23	6/29/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	81	35 - 141	06/30/20 20:23	
2-Fluorobiphenyl	77	31 - 118	06/30/20 20:23	
2-Fluorophenol	41	10 - 105	06/30/20 20:23	
Nitrobenzene-d5	65	31 - 110	06/30/20 20:23	
Phenol-d6	29	10 - 107	06/30/20 20:23	
p-Terphenyl-d14	84	10 - 165	06/30/20 20:23	

Tentatively Identified Compounds

			Result	
CAS#	Compound Identification	RT	ug/L	Q

No Tentatively Identified Compounds

Detected

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 12:45

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-006
 Units: ug/L

 Lab Code:
 R2005470-007
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/01/20 18:26	6/30/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/01/20 18:26	6/30/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/01/20 18:26	6/30/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/01/20 18:26	6/30/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/01/20 18:26	6/30/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/01/20 18:26	6/30/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/01/20 18:26	6/30/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/01/20 18:26	6/30/20	
2,4-Dinitrophenol	45 U	45	19	1	07/01/20 18:26	6/30/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/01/20 18:26	6/30/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/01/20 18:26	6/30/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/01/20 18:26	6/30/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/01/20 18:26	6/30/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/01/20 18:26	6/30/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/01/20 18:26	6/30/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/01/20 18:26	6/30/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/01/20 18:26	6/30/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/01/20 18:26	6/30/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/01/20 18:26	6/30/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/01/20 18:26	6/30/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/01/20 18:26	6/30/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/01/20 18:26	6/30/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/01/20 18:26	6/30/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/01/20 18:26	6/30/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/01/20 18:26	6/30/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/01/20 18:26	6/30/20	
4-Nitrophenol	45 U	45	5.8	1	07/01/20 18:26	6/30/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/01/20 18:26	6/30/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/01/20 18:26	6/30/20	
Anthracene	9.1 U	9.1	1.2	1	07/01/20 18:26	6/30/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/01/20 18:26	6/30/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/01/20 18:26	6/30/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/01/20 18:26	6/30/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/01/20 18:26	6/30/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/01/20 18:26	6/30/20	
Benzoic Acid	91 U	91	33	1	07/01/20 18:26	6/30/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/01/20 18:26	6/30/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/01/20 18:26	6/30/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/01/20 18:26	6/30/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	07/01/20 18:26	6/30/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	07/01/20 18:26	6/30/20	
Butyl Benzyl Phthalate	9.1 U 9.1 U	9.1 9.1	1.3	1	07/01/20 18:26	6/30/20	
	9.1 U 9.1 U	9.1 9.1	1.5	1	07/01/20 18:26	6/30/20	
Chrysene	9.1 U	9.1	1.1	1	07/01/20 18:26	0/30/20	

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Superset Reference: 20-0000554597 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 12:45

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-006
 Units: ug/L

 Lab Code:
 R2005470-007
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/01/20 18:26	6/30/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/01/20 18:26	6/30/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/01/20 18:26	6/30/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/01/20 18:26	6/30/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/01/20 18:26	6/30/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/01/20 18:26	6/30/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/01/20 18:26	6/30/20	
Fluorene	9.1 U	9.1	1.2	1	07/01/20 18:26	6/30/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/01/20 18:26	6/30/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/01/20 18:26	6/30/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/01/20 18:26	6/30/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/01/20 18:26	6/30/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/01/20 18:26	6/30/20	
Isophorone	9.1 U	9.1	1.3	1	07/01/20 18:26	6/30/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/01/20 18:26	6/30/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/01/20 18:26	6/30/20	
Naphthalene	9.1 U	9.1	1.1	1	07/01/20 18:26	6/30/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/01/20 18:26	6/30/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/01/20 18:26	6/30/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/01/20 18:26	6/30/20	
Phenol	9.1 U	9.1	0.91	1	07/01/20 18:26	6/30/20	
Pyrene	9.1 U	9.1	1.3	1	07/01/20 18:26	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	71	35 - 141	07/01/20 18:26	
2-Fluorobiphenyl	59	31 - 118	07/01/20 18:26	
2-Fluorophenol	40	10 - 105	07/01/20 18:26	
Nitrobenzene-d5	58	31 - 110	07/01/20 18:26	
Phenol-d6	27	10 - 107	07/01/20 18:26	
p-Terphenyl-d14	82	10 - 165	07/01/20 18:26	

Tentatively Identified Compounds

			Result	
CAS#	Compound Identification	RT	ug/L	Q

No Tentatively Identified Compounds

Detected

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 13:35

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-007
 Units: ug/L

 Lab Code:
 R2005470-008
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/01/20 18:54	6/30/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/01/20 18:54	6/30/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/01/20 18:54	6/30/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/01/20 18:54	6/30/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/01/20 18:54	6/30/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/01/20 18:54	6/30/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/01/20 18:54	6/30/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/01/20 18:54	6/30/20	
2,4-Dinitrophenol	45 U	45	19	1	07/01/20 18:54	6/30/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/01/20 18:54	6/30/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/01/20 18:54	6/30/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/01/20 18:54	6/30/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/01/20 18:54	6/30/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/01/20 18:54	6/30/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/01/20 18:54	6/30/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/01/20 18:54	6/30/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/01/20 18:54	6/30/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/01/20 18:54	6/30/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/01/20 18:54	6/30/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/01/20 18:54	6/30/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/01/20 18:54	6/30/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/01/20 18:54	6/30/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/01/20 18:54	6/30/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/01/20 18:54	6/30/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/01/20 18:54	6/30/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/01/20 18:54	6/30/20	
4-Nitrophenol	45 U	45	5.8	1	07/01/20 18:54	6/30/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/01/20 18:54	6/30/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/01/20 18:54	6/30/20	
Anthracene	9.1 U	9.1	1.2	1	07/01/20 18:54	6/30/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/01/20 18:54	6/30/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/01/20 18:54	6/30/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/01/20 18:54	6/30/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/01/20 18:54	6/30/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/01/20 18:54	6/30/20	
Benzoic Acid	91 U	91	33	1	07/01/20 18:54	6/30/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/01/20 18:54	6/30/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/01/20 18:54	6/30/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/01/20 18:54	6/30/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	07/01/20 18:54	6/30/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	07/01/20 18:54	6/30/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	07/01/20 18:54	6/30/20	
Chrysene	9.1 U	9.1	1.1	1	07/01/20 18:54	6/30/20	

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Superset Reference: 20-0000554597 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 13:35

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-007
 Units: ug/L

 Lab Code:
 R2005470-008
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/01/20 18:54	6/30/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/01/20 18:54	6/30/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/01/20 18:54	6/30/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/01/20 18:54	6/30/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/01/20 18:54	6/30/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/01/20 18:54	6/30/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/01/20 18:54	6/30/20	
Fluorene	9.1 U	9.1	1.2	1	07/01/20 18:54	6/30/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/01/20 18:54	6/30/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/01/20 18:54	6/30/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/01/20 18:54	6/30/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/01/20 18:54	6/30/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/01/20 18:54	6/30/20	
Isophorone	9.1 U	9.1	1.3	1	07/01/20 18:54	6/30/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/01/20 18:54	6/30/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/01/20 18:54	6/30/20	
Naphthalene	9.1 U	9.1	1.1	1	07/01/20 18:54	6/30/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/01/20 18:54	6/30/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/01/20 18:54	6/30/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/01/20 18:54	6/30/20	
Phenol	9.1 U	9.1	0.91	1	07/01/20 18:54	6/30/20	
Pyrene	9.1 U	9.1	1.3	1	07/01/20 18:54	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	85	35 - 141	07/01/20 18:54	
2-Fluorobiphenyl	62	31 - 118	07/01/20 18:54	
2-Fluorophenol	41	10 - 105	07/01/20 18:54	
Nitrobenzene-d5	60	31 - 110	07/01/20 18:54	
Phenol-d6	30	10 - 107	07/01/20 18:54	
p-Terphenyl-d14	91	10 - 165	07/01/20 18:54	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown hydrocarbon	11.35	6.1	J	
	unknown hydrocarbon	11.89	10	J	
	unknown hydrocarbon	12.48	15	J	
	unknown hydrocarbon	13.13	15	J	
	unknown hydrocarbon	13.83	12	J	
	unknown hydrocarbon	14.58	11	J	
	unknown hydrocarbon	15.37	7.8	J	

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Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 13:35

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-007
 Units: ug/L

 Lab Code:
 R2005470-008
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Tentatively Identified Compounds

CAS# Compound Identification RT ug/L Q
unknown hydrocarbon 16.11 6.3 J



Semivolatile Organic Compounds by GC

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 09:10

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-001
 Units: ug/L

 Lab Code:
 R2005470-002
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/01/20 18:12	6/30/20	
4,4'-DDE	0.045 U	0.045	0.019	1	07/01/20 18:12	6/30/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/01/20 18:12	6/30/20	
Aldrin	0.045 U	0.045	0.019	1	07/01/20 18:12	6/30/20	
Dieldrin	0.045 U	0.045	0.019	1	07/01/20 18:12	6/30/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/01/20 18:12	6/30/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/01/20 18:12	6/30/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/01/20 18:12	6/30/20	
Endrin	0.045 U	0.045	0.019	1	07/01/20 18:12	6/30/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/01/20 18:12	6/30/20	
Heptachlor	0.045 U	0.045	0.019	1	07/01/20 18:12	6/30/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/01/20 18:12	6/30/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/01/20 18:12	6/30/20	
Toxaphene	0.46 U	0.46	0.46	1	07/01/20 18:12	6/30/20	
alpha-BHC	0.045 U	0.045	0.019	1	07/01/20 18:12	6/30/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/01/20 18:12	6/30/20	
beta-BHC	0.045 U	0.045	0.019	1	07/01/20 18:12	6/30/20	
delta-BHC	0.045 U	0.045	0.019	1	07/01/20 18:12	6/30/20	
gamma-BHC (Lindane)	0.045 U	0.045	0.019	1	07/01/20 18:12	6/30/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/01/20 18:12	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	19	10 - 164	07/01/20 18:12	
Tetrachloro-m-xylene	60	10 - 147	07/01/20 18:12	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 09:55

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-002
 Units: ug/L

 Lab Code:
 R2005470-003
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/01/20 18:31	6/30/20	_
4,4'-DDE	0.045 U	0.045	0.019	1	07/01/20 18:31	6/30/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/01/20 18:31	6/30/20	
Aldrin	0.045 U	0.045	0.019	1	07/01/20 18:31	6/30/20	
Dieldrin	0.045 U	0.045	0.019	1	07/01/20 18:31	6/30/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/01/20 18:31	6/30/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/01/20 18:31	6/30/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/01/20 18:31	6/30/20	
Endrin	0.045 U	0.045	0.019	1	07/01/20 18:31	6/30/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/01/20 18:31	6/30/20	
Heptachlor	0.045 U	0.045	0.019	1	07/01/20 18:31	6/30/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/01/20 18:31	6/30/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/01/20 18:31	6/30/20	
Toxaphene	0.46 U	0.46	0.46	1	07/01/20 18:31	6/30/20	
alpha-BHC	0.045 U	0.045	0.019	1	07/01/20 18:31	6/30/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/01/20 18:31	6/30/20	
beta-BHC	0.045 U	0.045	0.019	1	07/01/20 18:31	6/30/20	
delta-BHC	0.045 U	0.045	0.019	1	07/01/20 18:31	6/30/20	
gamma-BHC (Lindane)	0.045 U	0.045	0.019	1	07/01/20 18:31	6/30/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/01/20 18:31	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	44	10 - 164	07/01/20 18:31	
Tetrachloro-m-xylene	51	10 - 147	07/01/20 18:31	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 10:30

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-003
 Units: ug/L

 Lab Code:
 R2005470-004
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/01/20 18:50	6/30/20	
4,4'-DDE	0.045 U	0.045	0.019	1	07/01/20 18:50	6/30/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/01/20 18:50	6/30/20	
Aldrin	0.045 U	0.045	0.019	1	07/01/20 18:50	6/30/20	
Dieldrin	0.045 U	0.045	0.019	1	07/01/20 18:50	6/30/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/01/20 18:50	6/30/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/01/20 18:50	6/30/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/01/20 18:50	6/30/20	
Endrin	0.045 U	0.045	0.019	1	07/01/20 18:50	6/30/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/01/20 18:50	6/30/20	
Heptachlor	0.045 U	0.045	0.019	1	07/01/20 18:50	6/30/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/01/20 18:50	6/30/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/01/20 18:50	6/30/20	
Toxaphene	0.46 U	0.46	0.46	1	07/01/20 18:50	6/30/20	
alpha-BHC	0.045 U	0.045	0.019	1	07/01/20 18:50	6/30/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/01/20 18:50	6/30/20	
beta-BHC	0.045 U	0.045	0.019	1	07/01/20 18:50	6/30/20	
delta-BHC	0.045 U	0.045	0.019	1	07/01/20 18:50	6/30/20	
gamma-BHC (Lindane)	0.045 U	0.045	0.019	1	07/01/20 18:50	6/30/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/01/20 18:50	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	61	10 - 164	07/01/20 18:50	
Tetrachloro-m-xylene	50	10 - 147	07/01/20 18:50	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 11:05

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-004
 Units: ug/L

 Lab Code:
 R2005470-005
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/01/20 19:09	6/30/20	_
4,4'-DDE	0.045 U	0.045	0.019	1	07/01/20 19:09	6/30/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/01/20 19:09	6/30/20	
Aldrin	0.045 U	0.045	0.019	1	07/01/20 19:09	6/30/20	
Dieldrin	0.045 U	0.045	0.019	1	07/01/20 19:09	6/30/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/01/20 19:09	6/30/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/01/20 19:09	6/30/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/01/20 19:09	6/30/20	
Endrin	0.045 U	0.045	0.019	1	07/01/20 19:09	6/30/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/01/20 19:09	6/30/20	
Heptachlor	0.045 U	0.045	0.019	1	07/01/20 19:09	6/30/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/01/20 19:09	6/30/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/01/20 19:09	6/30/20	
Toxaphene	0.46 U	0.46	0.46	1	07/01/20 19:09	6/30/20	
alpha-BHC	0.045 U	0.045	0.019	1	07/01/20 19:09	6/30/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/01/20 19:09	6/30/20	
beta-BHC	0.045 U	0.045	0.019	1	07/01/20 19:09	6/30/20	
delta-BHC	0.045 U	0.045	0.019	1	07/01/20 19:09	6/30/20	
gamma-BHC (Lindane)	0.045 U	0.045	0.019	1	07/01/20 19:09	6/30/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/01/20 19:09	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	61	10 - 164	07/01/20 19:09	
Tetrachloro-m-xylene	55	10 - 147	07/01/20 19:09	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 11:50

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-005
 Units: ug/L

 Lab Code:
 R2005470-006
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/01/20 19:48	6/30/20	
4,4'-DDE	0.045 U	0.045	0.019	1	07/01/20 19:48	6/30/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/01/20 19:48	6/30/20	
Aldrin	0.045 U	0.045	0.019	1	07/01/20 19:48	6/30/20	
Dieldrin	0.045 U	0.045	0.019	1	07/01/20 19:48	6/30/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/01/20 19:48	6/30/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/01/20 19:48	6/30/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/01/20 19:48	6/30/20	
Endrin	0.045 U	0.045	0.019	1	07/01/20 19:48	6/30/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/01/20 19:48	6/30/20	
Heptachlor	0.045 U	0.045	0.019	1	07/01/20 19:48	6/30/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/01/20 19:48	6/30/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/01/20 19:48	6/30/20	
Toxaphene	0.46 U	0.46	0.46	1	07/01/20 19:48	6/30/20	
alpha-BHC	0.045 U	0.045	0.019	1	07/01/20 19:48	6/30/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/01/20 19:48	6/30/20	
beta-BHC	0.045 U	0.045	0.019	1	07/01/20 19:48	6/30/20	
delta-BHC	0.045 U	0.045	0.019	1	07/01/20 19:48	6/30/20	
gamma-BHC (Lindane)	0.045 U	0.045	0.019	1	07/01/20 19:48	6/30/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/01/20 19:48	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	20	10 - 164	07/01/20 19:48	
Tetrachloro-m-xylene	57	10 - 147	07/01/20 19:48	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 12:45

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-006
 Units: ug/L

 Lab Code:
 R2005470-007
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/01/20 20:07	6/30/20	_
4,4'-DDE	0.045 U	0.045	0.019	1	07/01/20 20:07	6/30/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/01/20 20:07	6/30/20	
Aldrin	0.045 U	0.045	0.019	1	07/01/20 20:07	6/30/20	
Dieldrin	0.045 U	0.045	0.019	1	07/01/20 20:07	6/30/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/01/20 20:07	6/30/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/01/20 20:07	6/30/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/01/20 20:07	6/30/20	
Endrin	0.045 U	0.045	0.019	1	07/01/20 20:07	6/30/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/01/20 20:07	6/30/20	
Heptachlor	0.045 U	0.045	0.019	1	07/01/20 20:07	6/30/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/01/20 20:07	6/30/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/01/20 20:07	6/30/20	
Toxaphene	0.46 U	0.46	0.46	1	07/01/20 20:07	6/30/20	
alpha-BHC	0.045 U	0.045	0.019	1	07/01/20 20:07	6/30/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/01/20 20:07	6/30/20	
beta-BHC	0.045 U	0.045	0.019	1	07/01/20 20:07	6/30/20	
delta-BHC	0.045 U	0.045	0.019	1	07/01/20 20:07	6/30/20	
gamma-BHC (Lindane)	0.045 U	0.045	0.019	1	07/01/20 20:07	6/30/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/01/20 20:07	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	35	10 - 164	07/01/20 20:07	
Tetrachloro-m-xylene	49	10 - 147	07/01/20 20:07	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 13:35

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-007
 Units: ug/L

 Lab Code:
 R2005470-008
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/01/20 20:26	6/30/20	_
4,4'-DDE	0.045 U	0.045	0.019	1	07/01/20 20:26	6/30/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/01/20 20:26	6/30/20	
Aldrin	0.045 U	0.045	0.019	1	07/01/20 20:26	6/30/20	
Dieldrin	0.045 U	0.045	0.019	1	07/01/20 20:26	6/30/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/01/20 20:26	6/30/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/01/20 20:26	6/30/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/01/20 20:26	6/30/20	
Endrin	0.045 U	0.045	0.019	1	07/01/20 20:26	6/30/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/01/20 20:26	6/30/20	
Heptachlor	0.045 U	0.045	0.019	1	07/01/20 20:26	6/30/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/01/20 20:26	6/30/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/01/20 20:26	6/30/20	
Toxaphene	0.46 U	0.46	0.46	1	07/01/20 20:26	6/30/20	
alpha-BHC	0.045 U	0.045	0.019	1	07/01/20 20:26	6/30/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/01/20 20:26	6/30/20	
beta-BHC	0.045 U	0.045	0.019	1	07/01/20 20:26	6/30/20	
delta-BHC	0.045 U	0.045	0.019	1	07/01/20 20:26	6/30/20	
gamma-BHC (Lindane)	0.045 U	0.045	0.019	1	07/01/20 20:26	6/30/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/01/20 20:26	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	28	10 - 164	07/01/20 20:26	
Tetrachloro-m-xylene	56	10 - 147	07/01/20 20:26	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 09:10

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-001
 Units: ug/L

 Lab Code:
 R2005470-002
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/01/20 19:08	6/30/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/01/20 19:08	6/30/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/01/20 19:08	6/30/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/01/20 19:08	6/30/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/01/20 19:08	6/30/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/01/20 19:08	6/30/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/01/20 19:08	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	21	10 - 152	07/01/20 19:08	
Tetrachloro-m-xylene	53	14 - 129	07/01/20 19:08	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 09:55

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-002
 Units: ug/L

 Lab Code:
 R2005470-003
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/01/20 19:28	6/30/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/01/20 19:28	6/30/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/01/20 19:28	6/30/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/01/20 19:28	6/30/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/01/20 19:28	6/30/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/01/20 19:28	6/30/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/01/20 19:28	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	44	10 - 152	07/01/20 19:28	
Tetrachloro-m-xylene	43	14 - 129	07/01/20 19:28	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 10:30

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-003
 Units: ug/L

 Lab Code:
 R2005470-004
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/01/20 19:48	6/30/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/01/20 19:48	6/30/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/01/20 19:48	6/30/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/01/20 19:48	6/30/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/01/20 19:48	6/30/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/01/20 19:48	6/30/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/01/20 19:48	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	62	10 - 152	07/01/20 19:48	
Tetrachloro-m-xylene	44	14 - 129	07/01/20 19:48	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 11:05

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-004
 Units: ug/L

 Lab Code:
 R2005470-005
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/01/20 20:08	6/30/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/01/20 20:08	6/30/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/01/20 20:08	6/30/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/01/20 20:08	6/30/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/01/20 20:08	6/30/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/01/20 20:08	6/30/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/01/20 20:08	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q	
Decachlorobiphenyl	59	10 - 152	07/01/20 20:08		
Tetrachloro-m-xvlene	47	14 - 129	07/01/20 20:08		

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 11:50

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-005
 Units: ug/L

 Lab Code:
 R2005470-006
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/01/20 20:29	6/30/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/01/20 20:29	6/30/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/01/20 20:29	6/30/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/01/20 20:29	6/30/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/01/20 20:29	6/30/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/01/20 20:29	6/30/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/01/20 20:29	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	20	10 - 152	07/01/20 20:29	
Tetrachloro-m-xylene	49	14 - 129	07/01/20 20:29	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 12:45

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-006
 Units: ug/L

 Lab Code:
 R2005470-007
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/01/20 21:10	6/30/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/01/20 21:10	6/30/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/01/20 21:10	6/30/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/01/20 21:10	6/30/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/01/20 21:10	6/30/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/01/20 21:10	6/30/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/01/20 21:10	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	33	10 - 152	07/01/20 21:10	
Tetrachloro-m-xylene	42	14 - 129	07/01/20 21:10	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/24/20 13:35

Monitoring

Sample Matrix: Water Date Received: 06/25/20 11:00

 Sample Name:
 WG-9954-062420-SG-007
 Units: ug/L

 Lab Code:
 R2005470-008
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/01/20 21:30	6/30/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/01/20 21:30	6/30/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/01/20 21:30	6/30/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/01/20 21:30	6/30/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/01/20 21:30	6/30/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/01/20 21:30	6/30/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/01/20 21:30	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	26	10 - 152	07/01/20 21:30	
Tetrachloro-m-xvlene	49	14 - 129	07/01/20 21:30	



QC Summary Forms

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com



Volatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

SURROGATE RECOVERY SUMMARYVolatile Organic Compounds by GC/MS

Analysis Method: 8260C

Extraction Method: EPA 5030C

		4-Bromofluorobenzene	Dibromofluoromethane	Toluene-d8
Sample Name	Lab Code	85-122	89-119	87-121
TB-9954-062420-SG-001	R2005470-001	90	90	92
WG-9954-062420-SG-001	R2005470-002	90	92	93
WG-9954-062420-SG-002	R2005470-003	89	92	94
WG-9954-062420-SG-003	R2005470-004	91	94	96
WG-9954-062420-SG-004	R2005470-005	90	94	95
WG-9954-062420-SG-005	R2005470-006	93	93	94
WG-9954-062420-SG-006	R2005470-007	89	89	92
WG-9954-062420-SG-007	R2005470-008	90	90	92
Method Blank	RQ2007077-06	92	95	97
Method Blank	RQ2007141-04	90	89	91
Lab Control Sample	RQ2007077-04	94	98	97
Lab Control Sample	RQ2007141-03	95	95	93

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007077-06
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/02/20 13:35	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/02/20 13:35	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/02/20 13:35	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/02/20 13:35	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/02/20 13:35	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/02/20 13:35	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/02/20 13:35	
2-Butanone (MEK)	10 U	10	0.78	1	07/02/20 13:35	
2-Hexanone	10 U	10	0.20	1	07/02/20 13:35	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/02/20 13:35	
Acetone	10 U	10	5.0	1	07/02/20 13:35	
Benzene	5.0 U	5.0	0.20	1	07/02/20 13:35	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/02/20 13:35	
Bromoform	5.0 U	5.0	0.25	1	07/02/20 13:35	
Bromomethane	5.0 U	5.0	0.70	1	07/02/20 13:35	
Carbon Disulfide	10 U	10	0.42	1	07/02/20 13:35	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/02/20 13:35	
Chlorobenzene	5.0 U	5.0	0.20	1	07/02/20 13:35	
Chloroethane	5.0 U	5.0	0.23	1	07/02/20 13:35	
Chloroform	5.0 U	5.0	0.24	1	07/02/20 13:35	
Chloromethane	5.0 U	5.0	0.28	1	07/02/20 13:35	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/02/20 13:35	
Dichloromethane	5.0 U	5.0	0.65	1	07/02/20 13:35	
Ethylbenzene	5.0 U	5.0	0.20	1	07/02/20 13:35	
Styrene	5.0 U	5.0	0.20	1	07/02/20 13:35	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/02/20 13:35	
Toluene	5.0 U	5.0	0.20	1	07/02/20 13:35	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/02/20 13:35	
Vinyl Acetate	10 U	10	1.1	1	07/02/20 13:35	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/02/20 13:35	
Xylenes, Total	5.0 U	5.0	0.23	1	07/02/20 13:35	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/02/20 13:35	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/02/20 13:35	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/02/20 13:35	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	11	07/02/20 13:35	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007077-06Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	92	85 - 122	07/02/20 13:35	
Dibromofluoromethane	95	89 - 119	07/02/20 13:35	
Toluene-d8	97	87 - 121	07/02/20 13:35	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007077-06Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Tentatively Identified Compounds

CAS# Compound Identification RT ug/L

No Tentatively Identified Compounds

Detected

Q

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007141-04
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	06/30/20 23:08	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	06/30/20 23:08	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	06/30/20 23:08	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	06/30/20 23:08	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	06/30/20 23:08	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	06/30/20 23:08	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	06/30/20 23:08	
2-Butanone (MEK)	10 U	10	0.78	1	06/30/20 23:08	
2-Hexanone	10 U	10	0.20	1	06/30/20 23:08	
4-Methyl-2-pentanone	10 U	10	0.20	1	06/30/20 23:08	
Acetone	10 U	10	5.0	1	06/30/20 23:08	
Benzene	5.0 U	5.0	0.20	1	06/30/20 23:08	
Bromodichloromethane	5.0 U	5.0	0.20	1	06/30/20 23:08	
Bromoform	5.0 U	5.0	0.25	1	06/30/20 23:08	
Bromomethane	5.0 U	5.0	0.70	1	06/30/20 23:08	
Carbon Disulfide	10 U	10	0.42	1	06/30/20 23:08	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	06/30/20 23:08	
Chlorobenzene	5.0 U	5.0	0.20	1	06/30/20 23:08	
Chloroethane	5.0 U	5.0	0.23	1	06/30/20 23:08	
Chloroform	5.0 U	5.0	0.24	1	06/30/20 23:08	
Chloromethane	5.0 U	5.0	0.28	1	06/30/20 23:08	
Dibromochloromethane	5.0 U	5.0	0.20	1	06/30/20 23:08	
Dichloromethane	5.0 U	5.0	0.65	1	06/30/20 23:08	
Ethylbenzene	5.0 U	5.0	0.20	1	06/30/20 23:08	
Styrene	5.0 U	5.0	0.20	1	06/30/20 23:08	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	06/30/20 23:08	
Toluene	5.0 U	5.0	0.20	1	06/30/20 23:08	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	06/30/20 23:08	
Vinyl Acetate	10 U	10	1.1	1	06/30/20 23:08	
Vinyl Chloride	5.0 U	5.0	0.20	1	06/30/20 23:08	
Xylenes, Total	5.0 U	5.0	0.23	1	06/30/20 23:08	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	06/30/20 23:08	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	06/30/20 23:08	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	06/30/20 23:08	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	06/30/20 23:08	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007141-04Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	90	85 - 122	06/30/20 23:08	
Dibromofluoromethane	89	89 - 119	06/30/20 23:08	
Toluene-d8	91	87 - 121	06/30/20 23:08	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007141-04Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

Result

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/02/20

Sample Matrix: Water

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005470

Lab Control Sample

RQ2007077-04

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
1,1,1-Trichloroethane (TCA)	8260C	22.5	20.0	113	75-125
1,1,2,2-Tetrachloroethane	8260C	28.8	20.0	144 *	78-126
1,1,2-Trichloroethane	8260C	21.8	20.0	109	82-121
1,1-Dichloroethane (1,1-DCA)	8260C	23.3	20.0	117	80-124
1,1-Dichloroethene (1,1-DCE)	8260C	21.1	20.0	105	71-118
1,2-Dichloroethane	8260C	20.4	20.0	102	71-127
1,2-Dichloropropane	8260C	22.9	20.0	114	80-119
2-Butanone (MEK)	8260C	23.7	20.0	118	61-137
2-Hexanone	8260C	22.8	20.0	114	63-124
4-Methyl-2-pentanone	8260C	22.8	20.0	114	66-124
Acetone	8260C	25.8	20.0	129	40-161
Benzene	8260C	21.7	20.0	109	79-119
Bromodichloromethane	8260C	21.7	20.0	109	81-123
Bromoform	8260C	21.9	20.0	110	65-146
Bromomethane	8260C	17.7	20.0	89	42-166
Carbon Disulfide	8260C	21.0	20.0	105	66-128
Carbon Tetrachloride	8260C	20.4	20.0	102	70-127
Chlorobenzene	8260C	21.4	20.0	107	80-121
Chloroethane	8260C	22.3	20.0	112	62-131
Chloroform	8260C	22.2	20.0	111	79-120
Chloromethane	8260C	22.1	20.0	110	65-135
Dibromochloromethane	8260C	22.5	20.0	112	72-128
Dichloromethane	8260C	21.8	20.0	109	73-122
Ethylbenzene	8260C	22.1	20.0	110	76-120
Styrene	8260C	22.0	20.0	110	80-124
Tetrachloroethene (PCE)	8260C	20.4	20.0	102	72-125
Toluene	8260C	21.8	20.0	109	79-119
Trichloroethene (TCE)	8260C	17.7	20.0	89	74-122
Vinyl Acetate	8260C	30.7	20.0	154	52-174
Vinyl Chloride	8260C	23.1	20.0	116	74-159
cis-1,2-Dichloroethene	8260C	22.8	20.0	114	80-121
cis-1,3-Dichloropropene	8260C	22.4	20.0	112	77-122
trans-1,2-Dichloroethene	8260C	21.9	20.0	110	73-118
Printed 7/14/2020 2:17:50 PM			Supers	et Reference:20-000	00554597 rev 00

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix:

Project:

Water

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

> Units:ug/L Basis:NA

Service Request: R2005470

Date Analyzed: 07/02/20

Lab Control Sample

RQ2007077-04

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
trans-1.3-Dichloropropene	8260C	23.4	20.0	117	71-133

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring **Date Analyzed:** 06/30/20

Sample Matrix: Water

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005470

Lab Control Sample

RQ2007141-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
1,1,1-Trichloroethane (TCA)	8260C	21.6	20.0	108	75-125
1,1,2,2-Tetrachloroethane	8260C	26.2	20.0	131 *	78-126
1,1,2-Trichloroethane	8260C	22.0	20.0	110	82-121
1,1-Dichloroethane (1,1-DCA)	8260C	20.8	20.0	104	80-124
1,1-Dichloroethene (1,1-DCE)	8260C	20.3	20.0	101	71-118
1,2-Dichloroethane	8260C	19.5	20.0	98	71-127
1,2-Dichloropropane	8260C	21.2	20.0	106	80-119
2-Butanone (MEK)	8260C	16.9	20.0	85	61-137
2-Hexanone	8260C	17.3	20.0	87	63-124
4-Methyl-2-pentanone	8260C	17.2	20.0	86	66-124
Acetone	8260C	20.1	20.0	101	40-161
Benzene	8260C	21.2	20.0	106	79-119
Bromodichloromethane	8260C	21.4	20.0	107	81-123
Bromoform	8260C	22.5	20.0	113	65-146
Bromomethane	8260C	14.4	20.0	72	42-166
Carbon Disulfide	8260C	17.2	20.0	86	66-128
Carbon Tetrachloride	8260C	21.4	20.0	107	70-127
Chlorobenzene	8260C	21.9	20.0	109	80-121
Chloroethane	8260C	16.2	20.0	81	62-131
Chloroform	8260C	21.1	20.0	105	79-120
Chloromethane	8260C	17.5	20.0	87	65-135
Dibromochloromethane	8260C	23.1	20.0	116	72-128
Dichloromethane	8260C	20.7	20.0	103	73-122
Ethylbenzene	8260C	22.5	20.0	113	76-120
Styrene	8260C	22.5	20.0	112	80-124
Tetrachloroethene (PCE)	8260C	21.6	20.0	108	72-125
Toluene	8260C	21.5	20.0	107	79-119
Trichloroethene (TCE)	8260C	18.3	20.0	91	74-122
Vinyl Acetate	8260C	23.8	20.0	119	52-174
Vinyl Chloride	8260C	17.0	20.0	85	74-159
cis-1,2-Dichloroethene	8260C	22.1	20.0	111	80-121
cis-1,3-Dichloropropene	8260C	21.4	20.0	107	77-122
trans-1,2-Dichloroethene	8260C	21.1	20.0	105	73-118
Printed 7/14/2020 2:17:51 PM			Supers	et Reference:20-000	00554597 rev 00

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix:

Project:

Water

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

> Units:ug/L Basis:NA

Service Request: R2005470

Date Analyzed: 06/30/20

Lab Control Sample

RQ2007141-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
trans-1 3-Dichloropropene	8260C	21.6	20.0	108	71-133



Semivolatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

SURROGATE RECOVERY SUMMARY Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Extraction Method:** EPA 3510C

		2,4,6-Tribromophenol	2-Fluorobiphenyl	2-Fluorophenol
Sample Name	Lab Code	35-141	31-118	10-105
WG-9954-062420-SG-001	R2005470-002	73	53	35
WG-9954-062420-SG-002	R2005470-003	27*	26*	18
WG-9954-062420-SG-002 RE	R2005470-003	105	79	51
WG-9954-062420-SG-003	R2005470-004	82	71	46
WG-9954-062420-SG-004	R2005470-005	68	58	34
WG-9954-062420-SG-005	R2005470-006	81	77	41
WG-9954-062420-SG-006	R2005470-007	71	59	40
WG-9954-062420-SG-007	R2005470-008	85	62	41
Method Blank	RQ2006888-01	74	59	45
Method Blank	RQ2006968-03	78	71	44
Method Blank	RQ2007073-03	101	71	48
Lab Control Sample Duplicate	RQ2006888-02	91	76	48
Lab Control Sample Lab	RQ2006888-03	84	81	50
Control Sample	RQ2006968-04	100	80	56
Duplicate Lab Control Sample	RQ2006968-05	64	54	39
Lab Control Sample	RQ2007073-04	115	78	47
Duplicate Lab Control Sample	RQ2007073-05	103	67	47

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

SURROGATE RECOVERY SUMMARY Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Extraction Method:** EPA 3510C

		Nitrobenzene-d5	Phenol-d6	p-Terphenyl-d14
Sample Name	Lab Code	31-110	10-107	10-165
WG-9954-062420-SG-001	R2005470-002	43	24	84
WG-9954-062420-SG-002	R2005470-003	22*	12	35
WG-9954-062420-SG-002 RE	R2005470-003	85	32	56
WG-9954-062420-SG-003	R2005470-004	69	31	97
WG-9954-062420-SG-004	R2005470-005	51	23	85
WG-9954-062420-SG-005	R2005470-006	65	29	84
WG-9954-062420-SG-006	R2005470-007	58	27	82
WG-9954-062420-SG-007	R2005470-008	60	30	91
Method Blank	RQ2006888-01	63	33	104
Method Blank	RQ2006968-03	65	31	95
Method Blank	RQ2007073-03	69	35	66
Lab Control Sample Duplicate	RQ2006888-02	67	35	105
Lab Control Sample Lab	RQ2006888-03	68	35	108
Control Sample	RQ2006968-04	85	41	104
Duplicate Lab Control Sample	RQ2006968-05	54	29	78
Lab Control Sample	RQ2007073-04	80	34	58
Duplicate Lab Control Sample	RQ2007073-05	66	34	59

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2006888-01
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

1,2,4-Trichlorobenzene	Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1.3-Dichlorobenzene 10 U 10 1.1 1 06/30/20 14:46 6/29/20 2.4.5-Trichlorophenol 10 U 10 1.2 1 06/30/20 14:46 6/29/20 2.4.5-Trichlorophenol 10 U 10 1.1 1 06/30/20 14:46 6/29/20 2.4.5-Trichlorophenol 10 U 10 1.4 1 06/30/20 14:46 6/29/20 2.4-Dichlorophenol 10 U 10 1.3 1 06/30/20 14:46 6/29/20 2.4-Dichlorophenol 10 U 10 1.4 1 06/30/20 14:46 6/29/20 2.4-Dinitrophenol 50 U 50 20 1 06/30/20 14:46 6/29/20 2.4-Dinitrophenol 50 U 50 20 1 06/30/20 14:46 6/29/20 2.4-Dinitrophenol 10 U 10 1.4 1 06/30/20 14:46 6/29/20 2.4-Dinitrophenol 10 U 10 1.4 1 06/30/20 14:46 6/29/20 2.4-Dinitrophenol 10 U 10 1.4 1 06/30/20 14:46 6/29/20 2.4-Dinitrophenol 10 U 10 1.4 1 06/30/20 14:46 6/29/20 2.4-Dinitrophenol 10 U 10 1.4 1 06/30/20 14:46 6/29/20 2.4-Dinitrophenol 10 U 10 1.4 1 06/30/20 14:46 6/29/20 2.4-Dinitrophenol 10 U 10 1.1 1 06/30/20 14:46 6/29/20 2.4-Dinitrophenol 10 U 10 1.1 1 06/30/20 14:46 6/29/20 2.4-Dinitrophenol 10 U 10 1.3 1 06/30/20 14:46 6/29/20 2.4-Dinitrophenol 10 U 10 1.4 1 06/30/20 14:46 6/29/20 2.4-Dinitrophenol 10 U 10 1.5 1 06/30/20 14:46 6/29/20 2.4-Dinitrophenol 10 U 10 1.5 1 06/30/20 14:46 6/29/20 2.4-Dinitrophenol 10 U 10 1.5 1 06/30/20 14:46 6/29/20 2.4-Dinitrophenol 10 U 10 1.2 1 06/30/20 14:46 6/29/20 3.3-Dichlorobenzidine 10 U 10 1.2 1 06/30/20 14:46 6/29/20 3.3-Dichlorobenzidine 10 U 10 1.2 1 06/30/20 14:46 6/29/20 3.3-Dichloro-2-methylphenol 50 U 50 20 U 10/63/20 14:46 6/29/20 3.3-Dichloro-3-methylphenol 50 U 50 20 U 10/63/20 14:46 6/29/20 4.4-Dinitro-3-methylphenol 50 U 50 6.4 1 06/30/20 14:46 6/29/20 4.4-Dinitro-3-methylphenol 50 U 50 6.4 1 06/30/20 14:46 6/29/20 4.4-Dinitro-3-methylphenol 50 U 50 6.4 1 06/30/2	1,2,4-Trichlorobenzene	10 U	10	1.2	1	06/30/20 14:46	6/29/20	
1,4-Dichlorobenzene	1,2-Dichlorobenzene	10 U	10	1.2	1	06/30/20 14:46	6/29/20	
2,4,5-Trichlorophenol 10 U 10	1,3-Dichlorobenzene	10 U	10	1.1	1	06/30/20 14:46	6/29/20	
2,4-Frichlorophenol	1,4-Dichlorobenzene	10 U	10	1.2	1	06/30/20 14:46	6/29/20	
2,4-Dintchlorophenol 10 U 10	2,4,5-Trichlorophenol	10 U	10	1.1	1	06/30/20 14:46	6/29/20	
2.4-Dimethylphenol 10 U 10 1.4 1 06/30/20 14:46 6/29/20 2.4-Dinitrophenol 50 U 50 20 1 06/30/20 14:46 6/29/20 2.4-Dinitrotoluene 10 U 10 1.4 1 06/30/20 14:46 6/29/20 2.Chlorophenol 10 U 10 1.4 1 06/30/20 14:46 6/29/20 2.Chlorophenol 10 U 10 1.4 1 06/30/20 14:46 6/29/20 2.Methylphenol 10 U 10 1.3 1 06/30/20 14:46 6/29/20 2.Nitroaniline 10 U 10 1.4 1 06/30/20 14:46 6/29/20 2.Nitrophenol 10 U 10 1.4 1 06/30/20 14:46 6/29/20 2.Nitrophenol 10 U 10 1.5 1 06/30/20 14:46 6/29/20 2.Nitrophenol 10 U 10 1.5 1 06/30/20 14:46 6/29/20 3.3*Dichlorobenzidine 10 U 10 1.2 1	2,4,6-Trichlorophenol	10 U	10	1.4	1	06/30/20 14:46	6/29/20	
2,4-Dimitrophenol 10 U 10	2,4-Dichlorophenol	10 U	10	1.3	1	06/30/20 14:46	6/29/20	
2.4-Dinitrotoluene 10 U 10 2.4 1 06/30/20 14:46 6/29/20 2.6-Dinitrotoluene 10 U 10 1.4 1 06/30/20 14:46 6/29/20 2-Chlorophenol 10 U 10 1.1 1 06/30/20 14:46 6/29/20 2-Methylnaphthalene 10 U 10 1.3 1 06/30/20 14:46 6/29/20 2-Methylphenol 10 U 10 1.0 1 06/30/20 14:46 6/29/20 2-Mitrophenol 10 U 10 1.4 1 06/30/20 14:46 6/29/20 2-Nitrophenol 10 U 10 1.5 1 06/30/20 14:46 6/29/20 3-Nitrophenol 10 U 10 1.2 1 06/30/20 14:46 6/29/20 3-Nitronilline 10 U 10 1.2 1 06/30/20 14:46 6/29/20 4-G-Dinitrobenol 50 U 50 20 1 06/30/20 14:46 6/29/20 4-Shironiline 10 U 10 1.7 1 <td< td=""><td></td><td>10 U</td><td>10</td><td>1.4</td><td>1</td><td>06/30/20 14:46</td><td>6/29/20</td><td></td></td<>		10 U	10	1.4	1	06/30/20 14:46	6/29/20	
2,6-Dinitrotoluene 10 U 11 U 10 U 10 U 11 U 10 U 11 U 10 U 10 U 11 U 10 U 11 U 10 U <td>2,4-Dinitrophenol</td> <td>50 U</td> <td>50</td> <td>20</td> <td>1</td> <td>06/30/20 14:46</td> <td>6/29/20</td> <td></td>	2,4-Dinitrophenol	50 U	50	20	1	06/30/20 14:46	6/29/20	
2-Chloronaphthalene 10 U 10 II 1.4 1 06/30/20 14:46 6/29/20 2-Chlorophenol 10 U 10 II 1.1 1 06/30/20 14:46 6/29/20 2-Methylaphthalene 10 U 10 1.3 1 06/30/20 14:46 6/29/20 2-Methylphenol 10 U 10 1.4 1 06/30/20 14:46 6/29/20 2-Nitrophenol 10 U 10 1.4 1 06/30/20 14:46 6/29/20 2-Nitrophenol 10 U 10 1.5 1 06/30/20 14:46 6/29/20 3.3-Dichlorobenzidine 10 U 10 1.2 1 06/30/20 14:46 6/29/20 3-and 4-Methylphenol Coelution 10 U 10 1.2 1 06/30/20 14:46 6/29/20 3-introaniline 10 U 10 1.2 1 06/30/20 14:46 6/29/20 4-G-Dinitro-2-methylphenol 50 U 50 20 1 06/30/20 14:46 6/29/20 4-G-Dinitro-2-methylphenol 10 U 10 1.7 1 06/30/20 14:46 6/29/20 4-Chloro-3-methylphenol 10 U <t< td=""><td>2,4-Dinitrotoluene</td><td>10 U</td><td>10</td><td>2.4</td><td>1</td><td>06/30/20 14:46</td><td>6/29/20</td><td></td></t<>	2,4-Dinitrotoluene	10 U	10	2.4	1	06/30/20 14:46	6/29/20	
2-Chlorophenol 10 U 10 1.1 1 06/30/20 14:46 6/29/20 2-Methylnaphthalene 10 U 10 1.3 1 06/30/20 14:46 6/29/20 2-Methylphenol 10 U 10 1.3 1 06/30/20 14:46 6/29/20 2-Methylphenol 10 U 10 1.4 1 06/30/20 14:46 6/29/20 2-Nitroaniline 10 U 10 1.4 1 06/30/20 14:46 6/29/20 2-Nitrophenol 10 U 10 1.5 1 06/30/20 14:46 6/29/20 3.3*-Dichlorobenzidine 10 U 10 1.5 1 06/30/20 14:46 6/29/20 3.3*-Dichlorobenzidine 10 U 10 1.2 1 06/30/20 14:46 6/29/20 3-Anitroaniline 10 U 10 1.2 1 06/30/20 14:46 6/29/20 3-Nitroaniline 10 U 10 1.2 1 06/30/20 14:46 6/29/20 3-Nitroaniline 10 U 10 2.5 1 06/30/20 14:46 6/29/20 4-Chloro-methylphenol 50 U 50 20 1 06/30/20 14:46 6/29/20 4-Chloro-methylphenol 50 U 50 20 1 06/30/20 14:46 6/29/20 4-Chloro-s-methylphenol 10 U 10 1.7 1 06/30/20 14:46 6/29/20 4-Chloro-s-methylphenol 10 U 10 1.1 1 06/30/20 14:46 6/29/20 4-Chloro-s-methylphenol 10 U 10 1.1 1 06/30/20 14:46 6/29/20 4-Chloro-s-methylphenol 10 U 10 1.1 1 06/30/20 14:46 6/29/20 4-Chloro-s-methylphenol 10 U 10 1.1 1 06/30/20 14:46 6/29/20 4-Chlorophenyl Phenyl Ether 10 U 10 1.5 1 06/30/20 14:46 6/29/20 4-Chlorophenyl Phenyl Ether 10 U 10 1.5 1 06/30/20 14:46 6/29/20 4-Chlorophenyl Phenyl Ether 10 U 10 1.5 1 06/30/20 14:46 6/29/20 4-Nitrophenol 50 U 50 6.4 1 06/30/20 14:46 6/29/20 4-Nitrophenol 50 U 50 6.4 1 06/30/20 14:46 6/29/20 4-Cenaphthene 10 U 10 1.4 1 06/30/20 14:46 6/29/20 4-Cenaphthylene 10 U 10 1.4 1 06/30/20 14:46 6/29/20 4-Cenaphthylene 10 U 10 1.3 1 06/30/20 14:46 6/29/20 Benz(a)amthracene 10 U 10 1.2 1 06/30/20 14:46 6/29/20 Benz(a)amthracene 10 U 10 1.2 1 06/30/20 14:46 6/29/20 Benz(a)mthracene 10 U 10 1.3 1 06/30/20 14:46 6/29/20 Benz(b)fluoranthene 10 U 10 1.3 1 06/30/20 14:46 6/29/20 Benz(b)fluoranthene 10 U 10 1.3 1 06/30/20 14:46 6/29/20 Benz(b)fluoranthene 10 U 10 1.4 1 06/30/20 14:46 6/29/20 Benz(b)fluoranthene 10 U 10 1.4 1 06/30/20 14:46 6/29/20 Benz(b)fluoranthene 10 U 10 1.4 1 06/30/20 14:46 6/29/20 Benz(b)fluoranthene 10 U 10 1.4 1 06/30/20 14:46 6/29/20 Benz(c)c-loroethyl) Ether 10 U 10 1.0 1.0 1.0 1.0 06/30/20 14:46 6/29/20	2,6-Dinitrotoluene	10 U	10	1.4	1	06/30/20 14:46	6/29/20	
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2-Methylnaphthalene		10 U	10	1.1	1	06/30/20 14:46	6/29/20	
2-Methylphenol 10 U 10 1.0 1.0 06/30/20 14:46 6/29/20 2-Nitropailline 10 U 10 1.4 1 06/30/20 14:46 6/29/20 3,3'-Dichlorobenzidine 10 U 10 1.2 1 06/30/20 14:46 6/29/20 3-and 4-Methylphenol Coelution 10 U 10 1.2 1 06/30/20 14:46 6/29/20 3-Nitroaniline 10 U 10 2.5 1 06/30/20 14:46 6/29/20 4,6-Dinitro-2-methylphenol 50 U 50 20 1 06/30/20 14:46 6/29/20 4-Bromophenyl Phenyl Ether 10 U 10 1.7 1 06/30/20 14:46 6/29/20 4-Chloro-3-methylphenol 10 U 10 1.7 1 06/30/20 14:46 6/29/20 4-Chloro-3-methylphenol 10 U 10 1.0 1.0 1.0 1.0 1.0 6/30/20 14:46 6/29/20 4-Chloro-3-methylphenol		10 U	10	1.3	1	06/30/20 14:46	6/29/20	
2-Nitronailline 10 U 10 II 1.4 1 06/30/20 14:46 6/29/20 2-Nitrophenol 10 U 10 II 1.5 1 06/30/20 14:46 6/29/20 33-'Dichlorobenzidine 10 U 10 II 1.2 1 06/30/20 14:46 6/29/20 3- and 4-Methylphenol Coelution 10 U 10 II 1.2 1 06/30/20 14:46 6/29/20 3-Nitroaniline 10 U 10 U 2.5 1 06/30/20 14:46 6/29/20 4,6-Dinitro-2-methylphenol 50 U 50 D 20 II 1 06/30/20 14:46 6/29/20 4-Bromophenyl Phenyl Ether 10 U 10 II 1.7 II 1 06/30/20 14:46 6/29/20 4-Chloro-3-methylphenol 10 U 10 II 1.1 II 1 06/30/20 14:46 6/29/20 4-Chlorophenyl Phenyl Ether 10 U 10 II 1.0 06/30/20 14:46 6/29/20 4-Chlorophenyl Phenyl Ether 10 U 10 II 1.5 II 1 06/30/20 14:46 6/29/20 4-Nitroaniline 10 U 10 II 2.7 II 1 06/30/20 14:46 6/29/20 4-Nitrophen		10 U		1.0	1			
2-Nitrophenol 10 U 10 1.5 1 06/30/20 14:46 6/29/20 3,3'-Dichlorobenzidine 10 U 10 1.2 1 06/30/20 14:46 6/29/20 3-and 4-Methylphenol Coelution 10 U 10 1.2 1 06/30/20 14:46 6/29/20 3-Nitroaniline 10 U 10 2.5 1 06/30/20 14:46 6/29/20 4,6-Dinitro-2-methylphenol 50 U 50 20 1 06/30/20 14:46 6/29/20 4-Bromophenyl Phenyl Ether 10 U 10 1.7 1 06/30/20 14:46 6/29/20 4-Bromophenyl Phenyl Ether 10 U 10 1.7 1 06/30/20 14:46 6/29/20 4-Chloro-3-methylphenol 10 U 10 1.1 1 06/30/20 14:46 6/29/20 4-Chloro-3-methylphenol 10 U 10 1.0 1 06/30/20 14:46 6/29/20 4-Chloroaniline 10 U 10 1.0 1 06/30/20 14:46 6/29/20 4-Chloroaniline 10 U 10 1.5 1 06/30/20 14:46 6/29/20 4-Chloroaniline 10 U 10 1.5 1 06/30/20 14:46 6/29/20 4-Nitroaniline 10 U 10 1.5 1 06/30/20 14:46 6/29/20 4-Nitrophenol 50 U 50 6.4 1 06/30/20 14:46 6/29/20 4-Nitrophenol 50 U 50 6.4 1 06/30/20 14:46 6/29/20 4-Cenaphthylene 10 U 10 1.4 1 06/30/20 14:46 6/29/20 4-Cenaphthylene 10 U 10 1.4 1 06/30/20 14:46 6/29/20 4-Cenaphthylene 10 U 10 1.4 1 06/30/20 14:46 6/29/20 4-Cenaphthylene 10 U 10 1.3 1 06/30/20 14:46 6/29/20 4-Cenaphthylene 10 U 10 1.5 1 06/30/20 14:46 6/29/20 4-Cenaphthylene 10 U 10 1.5 1 06/30/20 14:46 6/29/20 4-Cenaphthylene 10 U 10 1.2 1 06/30/20 14:46 6/29/20 4-Cenaphthylene 10 U 10 1.2 1 06/30/20 14:46 6/29/20 4-Cenaphthylene 10 U 10 1.2 1 06/30/20 14:46 6/29/20 4-Cenaphthylene 10 U 10 1.3 1 06/30/20 14:46 6/29/20 4-Cenaphthylene 10 U 10 1.3 1 06/30/20 14:46 6/29/20 4-Cenaphthylene 10 U 10 1.3 1 06/30/20 14:46 6/29/20 4-Cenaphthylene 10 U 10 1.4 1 06/30/20 14:46 6/29/20 4-Cenaphthylene 10 U 10 1.4 1 06/30/20 14:46 6/29/20 4-Cenaphthylene 10 U 10 1.4 1	2-Nitroaniline							
3,3'-Dichlorobenzidine 10 U 10 1.2 1 06/30/20 14:46 6/29/20		10 U	10	1.5	1	06/30/20 14:46	6/29/20	
3- and 4-Methylphenol Coelution 10 U 2.5 1 06/30/20 14:46 6/29/20 6/29/20 4,6-Dinitro-2-methylphenol 50 U 50 U 50 U 1 06/30/20 14:46 6/29/20 6/29/20 4-Chloro-3-methylphenol 10 U 10 U 11 0 06/30/20 14:46 6/29/20 6/29/20 4-Chloro-3-methylphenol 10 U 10 U 10 U 10 06/30/20 14:46 6/29/20 6/29/20 4-Chlorophenyl Phenyl Ether 10 U 10 U 1.5 1 06/30/20 14:46 6/29/20 6/29/20 4-Nitrophenyl Phenyl Ether 10 U 10 U 1.5 1 06/30/20 14:46 6/29/20 6/29/20 4-Nitrophenol 50 U 50 G-4 1 06/30/20 14:46 6/29/20 6/29/20 6/29/20 Acenaphthene 10 U 10 U 1.4 1 06/30/20 14:46 6/29/20 6/29/20 Acenaphthylene 10 U 10 U 1.4 1 06/30/20 14:46 6/29/20 6/29/20 Benzo(a)pyrene 10 U 10 U <t< td=""><td></td><td>10 U</td><td></td><td>1.2</td><td>1</td><td></td><td></td><td></td></t<>		10 U		1.2	1			
3-Nitroaniline 10 U 10 2.5 1 06/30/20 14:46 6/29/20								
4,6-Dinitro-2-methylphenol 50 U 50 U 20 I 1 06/30/20 14:46 6/29/20 4-Bromophenyl Phenyl Ether 10 U 10 I 1.7 1 06/30/20 14:46 6/29/20 4-Chloro-3-methylphenol 10 U 10 I 1.1 1 06/30/20 14:46 6/29/20 4-Chloroaniline 10 U 10 I 1.0 1 06/30/20 14:46 6/29/20 4-Chlorophenyl Phenyl Ether 10 U 10 I 1.5 1 06/30/20 14:46 6/29/20 4-Nitroaniline 10 U 10 U 2.7 1 06/30/20 14:46 6/29/20 4-Nitrophenol 50 U 50 G.4 1 06/30/20 14:46 6/29/20 4-Nitrophenol 50 U 50 G.4 1 06/30/20 14:46 6/29/20 Acenaphthene 10 U 10 I.4 1 06/30/20 14:46 6/29/20 Acenaphthylene 10 U 10 I.3 1 06/30/20 14:46 6/29/20 Anthracene 10 U 10 I.3 1 06/30/20 14:46 6/29/20 Benz(a)anthracene 10 U 10 I.2 1 06/30/20 14:46 6/29/20		10 U	10	2.5	1	06/30/20 14:46	6/29/20	
4-Bromophenyl Phenyl Ether 10 U 10 1.7 1 06/30/20 14:46 6/29/20 4-Chloro-3-methylphenol 10 U 10 1.1 1 06/30/20 14:46 6/29/20 4-Chlorophenyl Phenyl Ether 10 U 10 1.0 1 06/30/20 14:46 6/29/20 4-Chlorophenyl Phenyl Ether 10 U 10 1.5 1 06/30/20 14:46 6/29/20 4-Nitroaniline 10 U 10 2.7 1 06/30/20 14:46 6/29/20 4-Nitrophenol 50 U 50 6.4 1 06/30/20 14:46 6/29/20 Acenaphthene 10 U 10 1.4 1 06/30/20 14:46 6/29/20 Acenaphthylene 10 U 10 1.4 1 06/30/20 14:46 6/29/20 Acenaphthylene 10 U 10 1.3 1 06/30/20 14:46 6/29/20 Acenaphthylene 10 U 10 1.6 1 06/30/20 14:46 6/29/20 Benzo(a)aphtracene 10 U 10 1	4,6-Dinitro-2-methylphenol	50 U						
4-Chloro-3-methylphenol 10 U 10		10 U	10	1.7	1	06/30/20 14:46	6/29/20	
4-Chloroaniline 10 U 10 U 10 U 10 O 1.0 O		10 U	10	1.1	1	06/30/20 14:46	6/29/20	
4-Nitroaniline 10 U 10 U 2.7 1 06/30/20 14:46 6/29/20 4-Nitrophenol 50 U 50 G.4 1 06/30/20 14:46 6/29/20 Acenaphthene 10 U 10 U 1.4 1 06/30/20 14:46 6/29/20 Acenaphthylene 10 U 10 U 1.4 1 06/30/20 14:46 6/29/20 Anthracene 10 U 10 U 1.3 1 06/30/20 14:46 6/29/20 Benz(a)anthracene 10 U 10 U 1.6 1 06/30/20 14:46 6/29/20 Benzo(a)pyrene 10 U 10 U 1.2 1 06/30/20 14:46 6/29/20 Benzo(b)fluoranthene 10 U 10 U 1.2 1 06/30/20 14:46 6/29/20 Benzo(k)fluoranthene 10 U 10 U 1.0 1 06/30/20 14:46 6/29/20 Benzo(k)fluoranthene 10 U 10 U 1.0 1 06/30/20 14:46 6/29/20 Benzo(k)fluoranthene 10 U 10 U 1.0 1 06/30/20 14:46 6/29/20 Benzo(k)fluoranthene 10 U 10 U 1.0		10 U	10	1.0	1			
4-Nitroaniline 10 U 10 U 2.7 1 06/30/20 14:46 6/29/20 4-Nitrophenol 50 U 50 G.4 1 06/30/20 14:46 6/29/20 Acenaphthene 10 U 10 U 1.4 1 06/30/20 14:46 6/29/20 Acenaphthylene 10 U 10 U 1.4 1 06/30/20 14:46 6/29/20 Anthracene 10 U 10 U 1.3 1 06/30/20 14:46 6/29/20 Benz(a)anthracene 10 U 10 U 1.6 1 06/30/20 14:46 6/29/20 Benzo(a)pyrene 10 U 10 U 1.2 1 06/30/20 14:46 6/29/20 Benzo(b)fluoranthene 10 U 10 U 1.2 1 06/30/20 14:46 6/29/20 Benzo(k)fluoranthene 10 U 10 U 1.0 1 06/30/20 14:46 6/29/20 Benzo(k)fluoranthene 10 U 10 U 1.0 1 06/30/20 14:46 6/29/20 Benzo(k)fluoranthene 10 U 10 U 1.0 1 06/30/20 14:46 6/29/20 Benzo(k)fluoranthene 10 U 10 U 1.0	4-Chlorophenyl Phenyl Ether	10 U	10	1.5	1	06/30/20 14:46	6/29/20	
4-Nitrophenol 50 U 50 d 6.4 1 06/30/20 14:46 6/29/20 Acenaphthene 10 U 10 U 10 U 1.4 1 06/30/20 14:46 6/29/20 Acenaphthylene 10 U 10 U 1.4 1 06/30/20 14:46 6/29/20 Anthracene 10 U 10 U 1.3 1 06/30/20 14:46 6/29/20 Benzo(a)anthracene 10 U 10 U 1.6 1 06/30/20 14:46 6/29/20 Benzo(a)pyrene 10 U 10 U 1.2 1 06/30/20 14:46 6/29/20 Benzo(b)fluoranthene 10 U 10 U 1.0 1.2 1 06/30/20 14:46 6/29/20 Benzo(k)fluoranthene 10 U 10 U 1.0 1.0 1 06/30/20 14:46 6/29/20 Benzoic Acid 100 U 100 U 36 1 06/30/20 14:46 6/29/20 Benzyl Alcohol 10 U 10 U 1.6 1 06/30/20 14:46 6/29/20 2,2'-Oxybis(1-chloropropane) 10 U 10 U 1.4 1 06/30/20 14:46 6/29/20 Bis(2-chloro		10 U		2.7	1			
Acenaphthene 10 U 10 U 10 U 1.4 1 06/30/20 14:46 6/29/20 Acenaphthylene 10 U 10 U 1.4 1 06/30/20 14:46 6/29/20 Anthracene 10 U 10 U 1.3 1 06/30/20 14:46 6/29/20 Benz(a)anthracene 10 U 10 U 1.6 1 06/30/20 14:46 6/29/20 Benzo(a)pyrene 10 U 10 U 1.2 1 06/30/20 14:46 6/29/20 Benzo(b)fluoranthene 10 U 10 U 1.2 1 06/30/20 14:46 6/29/20 Benzo(g,h,i)perylene 10 U 10 U 1.0 1.0 1 06/30/20 14:46 6/29/20 Benzo(k)fluoranthene 10 U 10 U 1.0 1.3 1 06/30/20 14:46 6/29/20 Benzoic Acid 10 U 10 U 1.0 1.6 1 06/30/20 14:46 6/29/20 Benzyl Alcohol 10 U 10 U 1.0 1.6 1 06/30/20 14:46 6/29/20 2,2'-Oxybis(1-chloropropane) 10 U 10 U 1.4 1 06/30/20 14:46 6/29/20 </td <td>4-Nitrophenol</td> <td>50 U</td> <td>50</td> <td>6.4</td> <td>1</td> <td></td> <td></td> <td></td>	4-Nitrophenol	50 U	50	6.4	1			
Acenaphthylene 10 U		10 U	10	1.4	1	06/30/20 14:46	6/29/20	
Anthracene 10 U		10 U	10	1.4	1	06/30/20 14:46	6/29/20	
Benz(a)anthracene 10 U 10 U <td></td> <td>10 U</td> <td>10</td> <td>1.3</td> <td>1</td> <td>06/30/20 14:46</td> <td>6/29/20</td> <td></td>		10 U	10	1.3	1	06/30/20 14:46	6/29/20	
Benzo(b)fluoranthene 10 U 10 U<		10 U	10	1.6	1		6/29/20	
Benzo(b)fluoranthene 10 U 10 U<	* *		10	1.2	1		6/29/20	
Benzo(g,h,i)perylene 10 U 10 U<		10 U						
Benzo(k)fluoranthene 10 U 10 U<			10		1			
Benzoic Acid 100 U 100 36 1 06/30/20 14:46 6/29/20 Benzyl Alcohol 10 U 10 U 1.6 1 06/30/20 14:46 6/29/20 2,2'-Oxybis(1-chloropropane) 10 U 10 U 1.4 1 06/30/20 14:46 6/29/20 Bis(2-chloroethoxy)methane 10 U 10 U 1.9 1 06/30/20 14:46 6/29/20 Bis(2-chloroethyl) Ether 10 U 10 U 1.3 1 06/30/20 14:46 6/29/20 Bis(2-ethylhexyl) Phthalate 10 U 10 U 1.0 1 06/30/20 14:46 6/29/20		10 U	10	1.3	1	06/30/20 14:46	6/29/20	
Benzyl Alcohol 10 U								
2,2'-Oxybis(1-chloropropane) 10 U 10 U 1.4 1 06/30/20 14:46 6/29/20 Bis(2-chloroethoxy)methane 10 U 10 U 1.9 1 06/30/20 14:46 6/29/20 Bis(2-chloroethyl) Ether 10 U 10 U 1.3 1 06/30/20 14:46 6/29/20 Bis(2-ethylhexyl) Phthalate 10 U 10 U 1.0 1 06/30/20 14:46 6/29/20			10	1.6	1			
Bis(2-chloroethoxy)methane 10 U 10 U <th< td=""><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td></th<>								
Bis(2-chloroethyl) Ether 10 U 10 U 1.3 1 06/30/20 14:46 6/29/20 Bis(2-ethylhexyl) Phthalate 10 U 10 U 1.0 1 06/30/20 14:46 6/29/20								
Bis(2-ethylhexyl) Phthalate 10 U 10 1.0 1 06/30/20 14:46 6/29/20								
Dulyi Dehzyi f huididle 10 U 10 1.4 1 00/30/20 14.40 0/29/20	Butyl Benzyl Phthalate	10 U	10	1.4	1	06/30/20 14:46	6/29/20	
Chrysene 10 U 10 1.2 1 06/30/20 14:46 6/29/20								

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Superset Reference: 20-0000554597 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2006888-01
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	10 U	10	2.0	1	06/30/20 14:46	6/29/20	
Di-n-octyl Phthalate	10 U	10	3.3	1	06/30/20 14:46	6/29/20	
Dibenz(a,h)anthracene	10 U	10	1.1	1	06/30/20 14:46	6/29/20	
Dibenzofuran	10 U	10	1.4	1	06/30/20 14:46	6/29/20	
Diethyl Phthalate	10 U	10	1.1	1	06/30/20 14:46	6/29/20	
Dimethyl Phthalate	10 U	10	1.3	1	06/30/20 14:46	6/29/20	
Fluoranthene	10 U	10	1.5	1	06/30/20 14:46	6/29/20	
Fluorene	10 U	10	1.3	1	06/30/20 14:46	6/29/20	
Hexachlorobenzene	10 U	10	1.6	1	06/30/20 14:46	6/29/20	
Hexachlorobutadiene	10 U	10	1.0	1	06/30/20 14:46	6/29/20	
Hexachlorocyclopentadiene	10 U	10	2.2	1	06/30/20 14:46	6/29/20	
Hexachloroethane	10 U	10	1.1	1	06/30/20 14:46	6/29/20	
Indeno(1,2,3-cd)pyrene	10 U	10	1.8	1	06/30/20 14:46	6/29/20	
Isophorone	10 U	10	1.4	1	06/30/20 14:46	6/29/20	
N-Nitrosodi-n-propylamine	10 U	10	1.2	1	06/30/20 14:46	6/29/20	
N-Nitrosodiphenylamine	10 U	10	2.7	1	06/30/20 14:46	6/29/20	
Naphthalene	10 U	10	1.2	1	06/30/20 14:46	6/29/20	
Nitrobenzene	10 U	10	1.5	1	06/30/20 14:46	6/29/20	
Pentachlorophenol (PCP)	50 U	50	9.8	1	06/30/20 14:46	6/29/20	
Phenanthrene	10 U	10	1.4	1	06/30/20 14:46	6/29/20	
Phenol	10 U	10	1.0	1	06/30/20 14:46	6/29/20	
Pyrene	10 U	10	1.5	1	06/30/20 14:46	6/29/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	74	35 - 141	06/30/20 14:46	
2-Fluorobiphenyl	59	31 - 118	06/30/20 14:46	
2-Fluorophenol	45	10 - 105	06/30/20 14:46	
Nitrobenzene-d5	63	31 - 110	06/30/20 14:46	
Phenol-d6	33	10 - 107	06/30/20 14:46	
p-Terphenyl-d14	104	10 - 165	06/30/20 14:46	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown hydrocarbon	11.37	4.9	J	
	unknown hydrocarbon	11.92	7.5	J	
	unknown hydrocarbon	12.51	11	J	
	unknown hydrocarbon	13.16	11	J	
	unknown hydrocarbon	13.87	10	J	
	unknown hydrocarbon	14.61	8.2	J	
	unknown hydrocarbon	15.40	6.5	J	

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Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2006888-01Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown	16.14	4.1	J	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2006968-03
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	10 U	10	1.2	1	07/01/20 17:03	6/30/20	
1,2-Dichlorobenzene	10 U	10	1.2	1	07/01/20 17:03	6/30/20	
1,3-Dichlorobenzene	10 U	10	1.1	1	07/01/20 17:03	6/30/20	
1,4-Dichlorobenzene	10 U	10	1.2	1	07/01/20 17:03	6/30/20	
2,4,5-Trichlorophenol	10 U	10	1.1	1	07/01/20 17:03	6/30/20	
2,4,6-Trichlorophenol	10 U	10	1.4	1	07/01/20 17:03	6/30/20	
2,4-Dichlorophenol	10 U	10	1.3	1	07/01/20 17:03	6/30/20	
2,4-Dimethylphenol	10 U	10	1.4	1	07/01/20 17:03	6/30/20	
2,4-Dinitrophenol	50 U	50	20	1	07/01/20 17:03	6/30/20	
2,4-Dinitrotoluene	10 U	10	2.4	1	07/01/20 17:03	6/30/20	
2,6-Dinitrotoluene	10 U	10	1.4	1	07/01/20 17:03	6/30/20	-
2-Chloronaphthalene	10 U	10	1.4	1	07/01/20 17:03	6/30/20	
2-Chlorophenol	10 U	10	1.1	1	07/01/20 17:03	6/30/20	
2-Methylnaphthalene	10 U	10	1.3	1	07/01/20 17:03	6/30/20	
2-Methylphenol	10 U	10	1.0	1	07/01/20 17:03	6/30/20	
2-Nitroaniline	10 U	10	1.4	1	07/01/20 17:03	6/30/20	-
2-Nitrophenol	10 U	10	1.5	1	07/01/20 17:03	6/30/20	
3,3'-Dichlorobenzidine	10 U	10	1.2	1	07/01/20 17:03	6/30/20	
3- and 4-Methylphenol Coelution	10 U	10	1.2	1	07/01/20 17:03	6/30/20	
3-Nitroaniline	10 U	10	2.5	1	07/01/20 17:03	6/30/20	
4,6-Dinitro-2-methylphenol	50 U	50	20	1	07/01/20 17:03	6/30/20	
4-Bromophenyl Phenyl Ether	10 U	10	1.7	1	07/01/20 17:03	6/30/20	
4-Chloro-3-methylphenol	10 U	10	1.1	1	07/01/20 17:03	6/30/20	
4-Chloroaniline	10 U	10	1.0	1	07/01/20 17:03	6/30/20	
4-Chlorophenyl Phenyl Ether	10 U	10	1.5	1	07/01/20 17:03	6/30/20	
4-Nitroaniline	10 U	10	2.7	1	07/01/20 17:03	6/30/20	
4-Nitrophenol	50 U	50	6.4	1	07/01/20 17:03	6/30/20	
Acenaphthene	10 U	10	1.4	1	07/01/20 17:03	6/30/20	
Acenaphthylene	10 U	10	1.4	1	07/01/20 17:03	6/30/20	
Anthracene	10 U	10	1.3	1	07/01/20 17:03	6/30/20	
Benz(a)anthracene	10 U	10	1.6	1	07/01/20 17:03	6/30/20	
Benzo(a)pyrene	10 U	10	1.2	1	07/01/20 17:03	6/30/20	
Benzo(b)fluoranthene	10 U	10	1.2	1	07/01/20 17:03	6/30/20	
Benzo(g,h,i)perylene	10 U	10	1.0	1	07/01/20 17:03	6/30/20	
Benzo(k)fluoranthene	10 U	10	1.3	1	07/01/20 17:03	6/30/20	
Benzoic Acid	100 U	100	36	1	07/01/20 17:03	6/30/20	
Benzyl Alcohol	10 U	10	1.6	1	07/01/20 17:03	6/30/20	
2,2'-Oxybis(1-chloropropane)	10 U	10	1.4	1	07/01/20 17:03	6/30/20	
Bis(2-chloroethoxy)methane	10 U	10	1.9	1	07/01/20 17:03	6/30/20	
Bis(2-chloroethyl) Ether	10 U	10	1.3	1	07/01/20 17:03	6/30/20	
Bis(2-ethylhexyl) Phthalate	10 U	10	1.0	1	07/01/20 17:03	6/30/20	
Butyl Benzyl Phthalate	10 U	10	1.4	1	07/01/20 17:03	6/30/20	
Chrysene	10 U	10	1.2	1	07/01/20 17:03	6/30/20	

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Superset Reference: 20-0000554597 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2006968-03Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	10 U	10	2.0	1	07/01/20 17:03	6/30/20	
Di-n-octyl Phthalate	10 U	10	3.3	1	07/01/20 17:03	6/30/20	
Dibenz(a,h)anthracene	10 U	10	1.1	1	07/01/20 17:03	6/30/20	
Dibenzofuran	10 U	10	1.4	1	07/01/20 17:03	6/30/20	
Diethyl Phthalate	10 U	10	1.1	1	07/01/20 17:03	6/30/20	
Dimethyl Phthalate	10 U	10	1.3	1	07/01/20 17:03	6/30/20	
Fluoranthene	10 U	10	1.5	1	07/01/20 17:03	6/30/20	
Fluorene	10 U	10	1.3	1	07/01/20 17:03	6/30/20	
Hexachlorobenzene	10 U	10	1.6	1	07/01/20 17:03	6/30/20	
Hexachlorobutadiene	10 U	10	1.0	1	07/01/20 17:03	6/30/20	
Hexachlorocyclopentadiene	10 U	10	2.2	1	07/01/20 17:03	6/30/20	
Hexachloroethane	10 U	10	1.1	1	07/01/20 17:03	6/30/20	
Indeno(1,2,3-cd)pyrene	10 U	10	1.8	1	07/01/20 17:03	6/30/20	
Isophorone	10 U	10	1.4	1	07/01/20 17:03	6/30/20	
N-Nitrosodi-n-propylamine	10 U	10	1.2	1	07/01/20 17:03	6/30/20	
N-Nitrosodiphenylamine	10 U	10	2.7	1	07/01/20 17:03	6/30/20	
Naphthalene	10 U	10	1.2	1	07/01/20 17:03	6/30/20	
Nitrobenzene	10 U	10	1.5	1	07/01/20 17:03	6/30/20	
Pentachlorophenol (PCP)	50 U	50	9.8	1	07/01/20 17:03	6/30/20	
Phenanthrene	10 U	10	1.4	1	07/01/20 17:03	6/30/20	
Phenol	10 U	10	1.0	1	07/01/20 17:03	6/30/20	
Pyrene	10 U	10	1.5	1	07/01/20 17:03	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	78	35 - 141	07/01/20 17:03	
2-Fluorobiphenyl	71	31 - 118	07/01/20 17:03	
2-Fluorophenol	44	10 - 105	07/01/20 17:03	
Nitrobenzene-d5	65	31 - 110	07/01/20 17:03	
Phenol-d6	31	10 - 107	07/01/20 17:03	
p-Terphenyl-d14	95	10 - 165	07/01/20 17:03	

Tentatively Identified Compounds

			Result	
CAS#	Compound Identification	RT	ug/L	Q

No Tentatively Identified Compounds

Detected

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007073-03Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

1,24-Trichlorobenzene	Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1.3-Dichlorobenzene	1,2,4-Trichlorobenzene	10 U	10	1.2	1	07/06/20 14:53	7/2/20	
1.4-Dichlorobenzene	1,2-Dichlorobenzene	10 U	10	1.2	1	07/06/20 14:53	7/2/20	
2.4.5-Trichlorophenol 10 U 10	1,3-Dichlorobenzene	10 U	10	1.1	1	07/06/20 14:53	7/2/20	
2.4Trichlorophenol 10 U 10	1,4-Dichlorobenzene	10 U	10	1.2	1	07/06/20 14:53	7/2/20	
2.4-Dichlorophenol 10 U 10	2,4,5-Trichlorophenol	10 U	10	1.1	1	07/06/20 14:53	7/2/20	
2,4-Dimitrylphenol 10 U 10 U 1.4 I 07/06/20 14:53 7/2/20 2,4-Dinitrophenol 50 U 50 20 I 1 07/06/20 14:53 7/2/20 2,4-Dinitrotoluene 10 U 10 U 10 I.4 I 1 07/06/20 14:53 7/2/20 2,6-Dinitrotoluene 10 U 10 II.4 I 1 07/06/20 14:53 7/2/20 2-Chloropaphthalene 10 U 10 II.4 I 1 07/06/20 14:53 7/2/20 2-Chlorophenol 10 U 10 II.1 II.1 II.1 II.1 II.1 07/06/20 14:53 7/2/20 2-Methylphenol 10 U II.1 II.1 II.1 II.1 II.1 07/06/20 14:53 7/2/20 2-Methylphenol 10 U II.1 II.1 II.1 II.1 07/06/20 14:53 7/2/20 2-Nitrophenol 10 U II.1 II.1 II.1 07/06/20 14:53 7/2/20 2-Nitrophenol 10 U II.1 II.1 II.1 07/06/20 14:53 7/2/20 3,3-Dichlorobenzidine 10 U II.1 II.1 II.1 07/06/20 14:53 7/2/20 3-and 4-Methylphenol Coclution 10 U II.1 II.1 II.1 07/06/20 14:53 7/2/20 3-Shitroaniline 10 U II.1 II.1 II.1 07/06/20 14:53 7/2/20 4-Bromophenyl Phenyl Ether 10 U II.1 II.1 II.1 07/06/20 14:53 <t< td=""><td>2,4,6-Trichlorophenol</td><td>10 U</td><td>10</td><td>1.4</td><td>1</td><td>07/06/20 14:53</td><td>7/2/20</td><td></td></t<>	2,4,6-Trichlorophenol	10 U	10	1.4	1	07/06/20 14:53	7/2/20	
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2-Methylnaphthalene		10 U	10	1.1	1	07/06/20 14:53	7/2/20	
2-Methylphenol		10 U	10	1.3	1		7/2/20	
2-Nitroaniline								
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3,3'-Dichlorobenzidine								
3- and 4-Methylphenol Coelution 10 U 10 1.2 1 07/06/20 14:53 7/2/20 3-Nitroaniline 10 U 10 2.5 1 07/06/20 14:53 7/2/20 4-6-Dinitro-2-methylphenol 50 U 50 20 1 07/06/20 14:53 7/2/20 4-Bromophenyl Phenyl Ether 10 U 10 1.7 1 07/06/20 14:53 7/2/20 4-Chloro-3-methylphenol 10 U 10 1.1 1 07/06/20 14:53 7/2/20 4-Chloroaniline 10 U 10 1.1 1 07/06/20 14:53 7/2/20 4-Chlorophenyl Phenyl Ether 10 U 10 1.5 1 07/06/20 14:53 7/2/20 4-Chlorophenyl Phenyl Ether 10 U 10 1.5 1 07/06/20 14:53 7/2/20 4-Chlorophenyl Phenyl Ether 10 U 10 1.5 1 07/06/20 14:53 7/2/20 4-Nitroaniline 10 U 10 2.7 1 07/06/20 14:53 7/2/20 4-Nitroaniline 10 U 10 2.7 1 07/06/20 14:53 7/2/20 4-Nitrophenol 50 U 50 6.4 1 07/06/20 14:53 7/2/20 4-Nitrophenol 50 U 50 6.4 1 07/06/20 14:53 7/2/20 4-Nitrophenol 10 U 10 1.4 1 07/06/20 14:53 7/2/20 4-Nitracene 10 U 10 1.4 1 07/06/20 14:53 7/2/20 4-Nitracene 10 U 10 1.3 1 07/06/20 14:53 7/2/20 4-Nitracene 10 U 10 1.3 1 07/06/20 14:53 7/2/20 4-Nitracene 10 U 10 1.6 1 07/06/20 14:53 7/2/20 4-Nitracene 10 U 10 1.6 1 07/06/20 14:53 7/2/20 4-Nitracene 10 U 10 1.2 1 07/06/20 14:53 7/2/20 4-Nitracene 10 U 10 1.2 1 07/06/20 14:53 7/2/20 4-Nitracene 10 U 10 1.2 1 07/06/20 14:53 7/2/20 4-Nitracene 10 U 10 1.3 1 07/06/20 14:53 7/2/20 4-Nitracene 10 U 10 1.3 1 07/06/20 14:53 7/2/20 4-Nitracene 10 U 10 1.3 1 07/06/20 14:53 7/2/20 4-Nitracene 10 U 10 1.3 1 07/06/20 14:53 7/2/20 4-Nitracene 10 U 10 1.3 1 07/06/20 14:53 7/2/20 4-Nitracene 10 U 10 1.6 1 07/06/20 14:53 7/2/20 4-Nitracene 10 U 10 1.6 1 07/06/20 14:53 7/2/20 4-Nitracene 10 U 10 1.4 1 07/06/20 14:53 7/2/20 4-Nitracene 10 U 10 1.6 1 07/06/20 14:53 7/2								
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Butyl Benzyl Phthalate 10 U 10 1.4 1 07/06/20 14:53 7/2/20								
	Chrysene	10 U	10	1.2	1	07/06/20 14:53	7/2/20	

Printed 7/14/2020 2:18:07 PM

Superset Reference: 20-0000554597 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007073-03
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	10 U	10	2.0	1	07/06/20 14:53	7/2/20	
Di-n-octyl Phthalate	10 U	10	3.3	1	07/06/20 14:53	7/2/20	
Dibenz(a,h)anthracene	10 U	10	1.1	1	07/06/20 14:53	7/2/20	
Dibenzofuran	10 U	10	1.4	1	07/06/20 14:53	7/2/20	
Diethyl Phthalate	10 U	10	1.1	1	07/06/20 14:53	7/2/20	
Dimethyl Phthalate	10 U	10	1.3	1	07/06/20 14:53	7/2/20	
Fluoranthene	10 U	10	1.5	1	07/06/20 14:53	7/2/20	
Fluorene	10 U	10	1.3	1	07/06/20 14:53	7/2/20	
Hexachlorobenzene	10 U	10	1.6	1	07/06/20 14:53	7/2/20	
Hexachlorobutadiene	10 U	10	1.0	1	07/06/20 14:53	7/2/20	
Hexachlorocyclopentadiene	10 U	10	2.2	1	07/06/20 14:53	7/2/20	
Hexachloroethane	10 U	10	1.1	1	07/06/20 14:53	7/2/20	
Indeno(1,2,3-cd)pyrene	10 U	10	1.8	1	07/06/20 14:53	7/2/20	
Isophorone	10 U	10	1.4	1	07/06/20 14:53	7/2/20	
N-Nitrosodi-n-propylamine	10 U	10	1.2	1	07/06/20 14:53	7/2/20	
N-Nitrosodiphenylamine	10 U	10	2.7	1	07/06/20 14:53	7/2/20	
Naphthalene	10 U	10	1.2	1	07/06/20 14:53	7/2/20	
Nitrobenzene	10 U	10	1.5	1	07/06/20 14:53	7/2/20	
Pentachlorophenol (PCP)	50 U	50	9.8	1	07/06/20 14:53	7/2/20	
Phenanthrene	10 U	10	1.4	1	07/06/20 14:53	7/2/20	
Phenol	10 U	10	1.0	1	07/06/20 14:53	7/2/20	
Pyrene	10 U	10	1.5	1	07/06/20 14:53	7/2/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	101	35 - 141	07/06/20 14:53	
2-Fluorobiphenyl	71	31 - 118	07/06/20 14:53	
2-Fluorophenol	48	10 - 105	07/06/20 14:53	
Nitrobenzene-d5	69	31 - 110	07/06/20 14:53	
Phenol-d6	35	10 - 107	07/06/20 14:53	
p-Terphenyl-d14	66	10 - 165	07/06/20 14:53	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown	11.52	7.9	J	
	unknown	11.93	8.1	J	
	unknown hydrocarbon	12.08	12	J	
	unknown hydrocarbon	12.69	17	J	
	unknown hydrocarbon	13.36	17	J	
	unknown hydrocarbon	14.08	14	J	
	unknown hydrocarbon	14.85	13	J	

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Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007073-03Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown hydrocarbon	15.65	11	J	
	unknown	16.34	8.5	J	

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring **Date Analyzed:** 06/30/20

Sample Matrix: Water

Duplicate Lab Control Sample Summary Semivolatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005470

Lab Control Sample

Duplicate Lab Control Sample

RQ2006888-02

RQ2006888-03

Analyte Name	Analytica l Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit		
1,2,4-Trichlorobenzene	8270D	48.2	80.0	60	51.2	80.0	64	10-127	6	30		
1,2-Dichlorobenzene	8270D	47.7	80.0	60	48.4	80.0	61	23-130	2	30		
1,3-Dichlorobenzene	8270D	47.1	80.0	59	47.8	80.0	60	21-90	2	30		
1,4-Dichlorobenzene	8270D	46.4	80.0	58	48.6	80.0	61	10-124	5	30		
2,4,5-Trichlorophenol	8270D	61.3	80.0	77	65.0	80.0	81	48-134	5	30		
2,4,6-Trichlorophenol	8270D	60.0	80.0	75	64.1	80.0	80	44-135	6	30		
2,4-Dichlorophenol	8270D	49.4	80.0	62	52.5	80.0	66	48-127	6	30		
2,4-Dimethylphenol	8270D	57.0	80.0	71	57.9	80.0	72	59-113	1	30		
2,4-Dinitrophenol	8270D	52.1	80.0	65	51.0	80.0	64	21-154	2	30		
2,4-Dinitrotoluene	8270D	70.5	80.0	88	66.1	80.0	83	54-130	6	30		
2,6-Dinitrotoluene	8270D	76.2	80.0	95	72.4	80.0	91	51-127	4	30		
2-Chloronaphthalene	8270D	61.8	80.0	77	65.6	80.0	82	40-108	6	30		
2-Chlorophenol	8270D	48.1	80.0	60	50.5	80.0	63	42-112	5	30		
2-Methylnaphthalene	8270D	50.8	80.0	63	55.2	80.0	69	34-102	9	30		
2-Methylphenol	8270D	51.5	80.0	64	55.0	80.0	69	47-100	8	30		
2-Nitroaniline	8270D	68.1	80.0	85	70.3	80.0	88	52-133	3	30		
2-Nitrophenol	8270D	53.6	80.0	67	53.8	80.0	67	43-131	<1	30		
3,3'-Dichlorobenzidine	8270D	64.6	80.0	81	64.1	80.0	80	43-126	1	30		
3- and 4-Methylphenol Coelution	8270D	45.9	80.0	57	49.4	80.0	62	40-92	8	30		
3-Nitroaniline	8270D	63.4	80.0	79	64.3	80.0	80	42-111	1	30		
4,6-Dinitro-2-methylphenol	8270D	55.5	80.0	69	54.9	80.0	69	36-152	<1	30		
4-Bromophenyl Phenyl Ether	8270D	66.9	80.0	84	64.2	80.0	80	48-114	5	30		
4-Chloro-3-methylphenol	8270D	53.5	80.0	67	59.9	80.0	75	52-113	11	30		
4-Chloroaniline	8270D	61.1	80.0	76	61.7	80.0	77	44-109	1	30		
4-Chlorophenyl Phenyl Ether	8270D	58.0	80.0	72	56.7	80.0	71	51-107	1	30		
4-Nitroaniline	8270D	59.6	80.0	74	58.9	80.0	74	54-133	<1	30		
4-Nitrophenol	8270D	20.2 J	80.0	25	19.3 J	80.0	24	10-126	4	30		
Acenaphthene	8270D	64.0	80.0	80	65.4	80.0	82	52-107	2	30		
Acenaphthylene	8270D	68.6	80.0	86	68.5	80.0	86	55-109	<1	30		
Anthracene	8270D	71.3	80.0	89	70.7	80.0	88	55-116	1	30		
Benz(a)anthracene	8270D	68.0	80.0	85	68.8	80.0	86	61-121	1	30		
Benzo(a)pyrene	8270D	72.8	80.0	91	73.1	80.0	91	44-114	<1	30		
Benzo(b)fluoranthene	8270D	68.1	80.0	85	69.8	80.0	87	62-115	2	30		
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QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request: R2005470 **Project:** Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring **Date Analyzed:** 06/30/20

Sample Matrix: Water

Duplicate Lab Control Sample Summary Semivolatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Lab Control Sample

Duplicate Lab Control Sample

RQ2006888-02

RQ2006888-03

Analyte Name	Analytica l Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
Benzo(g,h,i)perylene	8270D	81.3	80.0	102	80.9	80.0	101	63-136	<1	30
Benzo(k)fluoranthene	8270D	76.6	80.0	96	73.8	80.0	92	49-133	4	30
Benzoic Acid	8270D	66.5 J	120	55	70.2 J	120	58	10-94	5	30
Benzyl Alcohol	8270D	58.9	80.0	74	61.3	80.0	77	31-109	4	30
2,2'-Oxybis(1-chloropropane)	8270D	53.3	80.0	67	55.4	80.0	69	32-122	3	30
Bis(2-chloroethoxy)methane	8270D	57.7	80.0	72	59.7	80.0	75	55-110	4	30
Bis(2-chloroethyl) Ether	8270D	52.8	80.0	66	53.8	80.0	67	46-102	2	30
Bis(2-ethylhexyl) Phthalate	8270D	73.5	80.0	92	73.7	80.0	92	51-132	<1	30
Butyl Benzyl Phthalate	8270D	70.0	80.0	88	68.3	80.0	85	41-148	3	30
Chrysene	8270D	73.2	80.0	91	75.1	80.0	94	57-118	3	30
Di-n-butyl Phthalate	8270D	79.3	80.0	99	71.5	80.0	89	57-128	11	30
Di-n-octyl Phthalate	8270D	73.8	80.0	92	73.5	80.0	92	62-124	<1	30
Dibenz(a,h)anthracene	8270D	87.1	80.0	109	89.3	80.0	112	54-135	3	30
Dibenzofuran	8270D	68.5	80.0	86	67.1	80.0	84	55-110	2	30
Diethyl Phthalate	8270D	67.4	80.0	84	63.4	80.0	79	53-113	6	30
Dimethyl Phthalate	8270D	70.5	80.0	88	68.0	80.0	85	51-112	3	30
Fluoranthene	8270D	74.8	80.0	93	72.6	80.0	91	66-127	2	30
Fluorene	8270D	67.0	80.0	84	66.6	80.0	83	54-106	1	30
Hexachlorobenzene	8270D	79.1	80.0	99	77.3	80.0	97	53-123	2	30
Hexachlorobutadiene	8270D	52.4	80.0	66	55.8	80.0	70	16-95	6	30
Hexachlorocyclopentadiene	8270D	24.2	80.0	30	26.0	80.0	32	10-99	6	30
Hexachloroethane	8270D	47.0	80.0	59	47.2	80.0	59	15-92	<1	30
Indeno(1,2,3-cd)pyrene	8270D	71.8	80.0	90	73.4	80.0	92	62-137	2	30
Isophorone	8270D	50.0	80.0	63	49.9	80.0	62	50-116	2	30
N-Nitrosodi-n-propylamine	8270D	58.9	80.0	74	62.8	80.0	78	49-115	5	30
N-Nitrosodiphenylamine	8270D	80.4	80.0	101	78.3	80.0	98	45-123	3	30
Naphthalene	8270D	55.4	80.0	69	54.5	80.0	68	38-99	1	30
Nitrobenzene	8270D	50.8	80.0	64	53.6	80.0	67	46-108	5	30
Pentachlorophenol (PCP)	8270D	61.8	80.0	77	61.0	80.0	76	29-164	1	30
Phenanthrene	8270D	71.2	80.0	89	68.7	80.0	86	58-118	3	30
Phenol	8270D	32.3	80.0	40	32.6	80.0	41	10-113	2	30
Pyrene	8270D	77.6	80.0	97	74.2	80.0	93	61-122	4	30

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QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/01/20

Sample Matrix: Water

Duplicate Lab Control Sample Summary Semivolatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005470

Lab Control Sample

Duplicate Lab Control Sample

RQ2006968-04

RQ2006968-05

Analyte Name	Analytica l Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
1,2,4-Trichlorobenzene	8270D	60.4	80.0	76	39.0	80.0	49	10-127	43*	30
1,2-Dichlorobenzene	8270D	57.5	80.0	72	37.6	80.0	47	23-130	42*	30
1,3-Dichlorobenzene	8270D	57.0	80.0	71	36.9	80.0	46	21-90	43*	30
1,4-Dichlorobenzene	8270D	55.7	80.0	70	36.5	80.0	46	10-124	41*	30
2,4,5-Trichlorophenol	8270D	69.0	80.0	86	46.8	80.0	58	48-134	39*	30
2,4,6-Trichlorophenol	8270D	65.6	80.0	82	42.1	80.0	53	44-135	43*	30
2,4-Dichlorophenol	8270D	62.6	80.0	78	41.4	80.0	52	48-127	40*	30
2,4-Dimethylphenol	8270D	65.7	80.0	82	44.7	80.0	56 *	59-113	38*	30
2,4-Dinitrophenol	8270D	60.5	80.0	76	38.9 J	80.0	49	21-154	43*	30
2,4-Dinitrotoluene	8270D	70.4	80.0	88	46.4	80.0	58	54-130	41*	30
2,6-Dinitrotoluene	8270D	79.2	80.0	99	54.5	80.0	68	51-127	37*	30
2-Chloronaphthalene	8270D	66.2	80.0	83	44.5	80.0	56	40-108	39*	30
2-Chlorophenol	8270D	58.8	80.0	74	37.5	80.0	47	42-112	45*	30
2-Methylnaphthalene	8270D	62.8	80.0	79	39.8	80.0	50	34-102	45*	30
2-Methylphenol	8270D	61.4	80.0	77	44.0	80.0	55	47-100	33*	30
2-Nitroaniline	8270D	71.5	80.0	89	47.7	80.0	60	52-133	39*	30
2-Nitrophenol	8270D	64.1	80.0	80	41.7	80.0	52	43-131	42*	30
3,3'-Dichlorobenzidine	8270D	66.6	80.0	83	43.9	80.0	55	43-126	41*	30
3- and 4-Methylphenol Coelution	8270D	55.1	80.0	69	38.5	80.0	48	40-92	36*	30
3-Nitroaniline	8270D	67.2	80.0	84	55.6	80.0	69	42-111	20	30
4,6-Dinitro-2-methylphenol	8270D	60.3	80.0	75	41.1 J	80.0	51	36-152	38*	30
4-Bromophenyl Phenyl Ether	8270D	69.7	80.0	87	50.5	80.0	63	48-114	32*	30
4-Chloro-3-methylphenol	8270D	66.1	80.0	83	40.6	80.0	51 *	52-113	48*	30
4-Chloroaniline	8270D	64.5	80.0	81	54.9	80.0	69	44-109	16	30
4-Chlorophenyl Phenyl Ether	8270D	62.5	80.0	78	43.8	80.0	55	51-107	35*	30
4-Nitroaniline	8270D	63.5	80.0	79	42.6	80.0	53 *	54-133	39*	30
4-Nitrophenol	8270D	27.1 J	80.0	34	20.8 J	80.0	26	10-126	27	30
Acenaphthene	8270D	69.8	80.0	87	46.0	80.0	57	52-107	42*	30
Acenaphthylene	8270D	73.7	80.0	92	49.7	80.0	62	55-109	39*	30
Anthracene	8270D	74.4	80.0	93	49.5	80.0	62	55-116	40*	30
Benz(a)anthracene	8270D	69.6	80.0	87	49.6	80.0	62	61-121	34*	30
Benzo(a)pyrene	8270D	75.2	80.0	94	52.1	80.0	65	44-114	36*	30
Benzo(b)fluoranthene	8270D	72.6	80.0	91	50.6	80.0	63	62-115	36*	30

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QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/01/20

Sample Matrix: Water

Duplicate Lab Control Sample Summary Semivolatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005470

Lab Control Sample

Duplicate Lab Control Sample

RQ2006968-04

RQ2006968-05

Analyte Name	Analytica l Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
Benzo(g,h,i)perylene	8270D	76.4	80.0	95	56.5	80.0	71	63-136	29	30
Benzo(k)fluoranthene	8270D	77.9	80.0	97	56.2	80.0	70	49-133	32*	30
Benzoic Acid	8270D	65.8 J	120	55	56.4 J	120	47	10-94	16	30
Benzyl Alcohol	8270D	63.3	80.0	79	47.2	80.0	59	31-109	29	30
2,2'-Oxybis(1-chloropropane)	8270D	64.2	80.0	80	41.7	80.0	52	32-122	42*	30
Bis(2-chloroethoxy)methane	8270D	65.9	80.0	82	45.6	80.0	57	55-110	36*	30
Bis(2-chloroethyl) Ether	8270D	61.5	80.0	77	39.9	80.0	50	46-102	43*	30
Bis(2-ethylhexyl) Phthalate	8270D	73.4	80.0	92	51.7	80.0	65	51-132	34*	30
Butyl Benzyl Phthalate	8270D	75.3	80.0	94	49.3	80.0	62	41-148	41*	30
Chrysene	8270D	73.1	80.0	91	53.9	80.0	67	57-118	30	30
Di-n-butyl Phthalate	8270D	83.3	80.0	104	52.5	80.0	66	57-128	45*	30
Di-n-octyl Phthalate	8270D	80.2	80.0	100	53.2	80.0	66	62-124	41*	30
Dibenz(a,h)anthracene	8270D	82.6	80.0	103	62.0	80.0	78	54-135	28	30
Dibenzofuran	8270D	72.8	80.0	91	49.0	80.0	61	55-110	39*	30
Diethyl Phthalate	8270D	65.2	80.0	82	45.6	80.0	57	53-113	36*	30
Dimethyl Phthalate	8270D	75.9	80.0	95	50.9	80.0	64	51-112	39*	30
Fluoranthene	8270D	80.7	80.0	101	54.0	80.0	68	66-127	39*	30
Fluorene	8270D	72.2	80.0	90	49.5	80.0	62	54-106	37*	30
Hexachlorobenzene	8270D	79.9	80.0	100	58.9	80.0	74	53-123	30	30
Hexachlorobutadiene	8270D	62.6	80.0	78	40.2	80.0	50	16-95	44*	30
Hexachlorocyclopentadiene	8270D	26.8	80.0	33	18.1	80.0	23	10-99	36*	30
Hexachloroethane	8270D	55.2	80.0	69	35.9	80.0	45	15-92	42*	30
Indeno(1,2,3-cd)pyrene	8270D	70.3	80.0	88	49.6	80.0	62	62-137	35*	30
Isophorone	8270D	61.1	80.0	76	37.2	80.0	46 *	50-116	49*	30
N-Nitrosodi-n-propylamine	8270D	70.8	80.0	89	45.4	80.0	57	49-115	44*	30
N-Nitrosodiphenylamine	8270D	82.8	80.0	103	60.2	80.0	75	45-123	31*	30
Naphthalene	8270D	64.9	80.0	81	42.8	80.0	53	38-99	42*	30
Nitrobenzene	8270D	68.2	80.0	85	42.7	80.0	53	46-108	46*	30
Pentachlorophenol (PCP)	8270D	78.0	80.0	97	49.9 J	80.0	62	29-164	44*	30
Phenanthrene	8270D	72.6	80.0	91	49.0	80.0	61	58-118	39*	30
Phenol	8270D	36.4	80.0	46	27.1	80.0	34	10-113	30	30
Pyrene	8270D	78.0	80.0	98	52.0	80.0	65	61-122	40*	30

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QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix: Water

Duplicate Lab Control Sample Summary Semivolatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Date Analyzed: 07/06/20 - 07/08/20

Service Request: R2005470

Lab Control Sample

Duplicate Lab Control Sample

RQ2007073-04

RQ2007073-05

Analyte Name	Analytica l Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
1,2,4-Trichlorobenzene	8270D	56.0	80.0	70 10	50.2	80.0	63	10-127	11	30
1,2-Dichlorobenzene	8270D	56.6	80.0	71	47.1	80.0	59	23-130	18	30
1,3-Dichlorobenzene	8270D	55.5	80.0	69	48.6	80.0	61	21-90	12	30
1,4-Dichlorobenzene	8270D	53.4	80.0	67	47.6	80.0	60	10-124	11	30
2,4,5-Trichlorophenol	8270D	69.5	80.0	87	57.7	80.0	72	48-134	19	30
2,4,6-Trichlorophenol	8270D	65.0	80.0	81	55.5	80.0	69	44-135	16	30
2,4-Dichlorophenol	8270D	58.0	80.0	73	50.9	80.0	64	48-127	13	30
2,4-Dimethylphenol	8270D	58.8	80.0	74	56.3	80.0	70	59-113	6	30
2,4-Dinitrophenol	8270D	48.1 J	80.0	60	43.6 J	80.0	55	21-154	9	30
2,4-Dinitrotoluene	8270D	64.1	80.0	80	55.3	80.0	69	54-130	15	30
2,6-Dinitrotoluene	8270D	73.4	80.0	92	65.1	80.0	81	51-127	13	30
2-Chloronaphthalene	8270D	65.8	80.0	82	56.4	80.0	71	40-108	14	30
2-Chlorophenol	8270D	52.9	80.0	66	49.9	80.0	62	42-112	6	30
2-Methylnaphthalene	8270D	58.8	80.0	74	52.4	80.0	65	34-102	13	30
2-Methylphenol	8270D	55.4	80.0	69	50.8	80.0	64	47-100	8	30
2-Nitroaniline	8270D	67.4	80.0	84	65.2	80.0	82	52-133	2	30
2-Nitrophenol	8270D	58.6	80.0	73	51.8	80.0	65	43-131	12	30
3,3'-Dichlorobenzidine	8270D	59.9	80.0	75	56.1	80.0	70	43-126	7	30
3- and 4-Methylphenol Coelution	8270D	48.9	80.0	61	47.2	80.0	59	40-92	3	30
3-Nitroaniline	8270D	58.4	80.0	73	62.2	80.0	78	42-111	7	30
4,6-Dinitro-2-methylphenol	8270D	54.0	80.0	68	45.9 J	80.0	57	36-152	18	30
4-Bromophenyl Phenyl Ether	8270D	66.2	80.0	83	60.4	80.0	75	48-114	10	30
4-Chloro-3-methylphenol	8270D	62.1	80.0	78	55.8	80.0	70	52-113	11	30
4-Chloroaniline	8270D	57.6	80.0	72	61.9	80.0	77	44-109	7	30
4-Chlorophenyl Phenyl Ether	8270D	58.4	80.0	73	49.4	80.0	62	51-107	16	30
4-Nitroaniline	8270D	58.1	80.0	73	50.1	80.0	63	54-133	15	30
4-Nitrophenol	8270D	20.2 J	80.0	25	18.7 J	80.0	23	10-126	8	30
Acenaphthene	8270D	64.3	80.0	80	58.2	80.0	73	52-107	9	30
Acenaphthylene	8270D	70.3	80.0	88	62.1	80.0	78	55-109	12	30
Anthracene	8270D	67.2	80.0	84	59.9	80.0	75	55-116	11	30
Benz(a)anthracene	8270D	60.6	80.0	76	58.1	80.0	73	61-121	4	30
Benzo(a)pyrene	8270D	62.2	80.0	78	61.3	80.0	77	44-114	1	30
Benzo(b)fluoranthene	8270D	59.2	80.0	74	58.3	80.0	73	62-115	1	30
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QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix: Water

Service Request: R2005470 **Date Analyzed:** 07/06/20 - 07/08/20

Duplicate Lab Control Sample Summary Semivolatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Lab Control Sample

Duplicate Lab Control Sample

RQ2007073-04

RQ2007073-05

Analyte Name	Analytica l Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
Benzo(g,h,i)perylene	8270D	70.4	80.0	88	68.4	80.0	85	63-136	3	30
Benzo(k)fluoranthene	8270D	68.5	80.0	86	64.2	80.0	80	49-133	7	30
Benzoic Acid	8270D	70.9 J	120	59	67.3 J	120	56	10-94	5	30
Benzyl Alcohol	8270D	60.0	80.0	75	58.9	80.0	74	31-109	1	30
2,2'-Oxybis(1-chloropropane)	8270D	63.4	80.0	79	54.4	80.0	68	32-122	15	30
Bis(2-chloroethoxy)methane	8270D	66.2	80.0	83	57.4	80.0	72	55-110	14	30
Bis(2-chloroethyl) Ether	8270D	61.0	80.0	76	51.7	80.0	65	46-102	16	30
Bis(2-ethylhexyl) Phthalate	8270D	62.1	80.0	78	60.8	80.0	76	51-132	3	30
Butyl Benzyl Phthalate	8270D	62.3	80.0	78	60.3	80.0	75	41-148	4	30
Chrysene	8270D	65.3	80.0	82	63.8	80.0	80	57-118	2	30
Di-n-butyl Phthalate	8270D	72.4	80.0	91	66.3	80.0	83	57-128	9	30
Di-n-octyl Phthalate	8270D	63.7	80.0	80	61.0	80.0	76	62-124	5	30
Dibenz(a,h)anthracene	8270D	75.6	80.0	95	71.4	80.0	89	54-135	7	30
Dibenzofuran	8270D	69.7	80.0	87	57.6	80.0	72	55-110	19	30
Diethyl Phthalate	8270D	62.9	80.0	79	53.9	80.0	67	53-113	16	30
Dimethyl Phthalate	8270D	72.4	80.0	90	64.0	80.0	80	51-112	12	30
Fluoranthene	8270D	72.0	80.0	90	66.4	80.0	83	66-127	8	30
Fluorene	8270D	69.1	80.0	86	56.8	80.0	71	54-106	19	30
Hexachlorobenzene	8270D	73.1	80.0	91	69.0	80.0	86	53-123	6	30
Hexachlorobutadiene	8270D	58.0	80.0	73	53.1	80.0	66	16-95	10	30
Hexachlorocyclopentadiene	8270D	19.1	80.0	24	18.0	80.0	23	10-99	4	30
Hexachloroethane	8270D	55.7	80.0	70	49.0	80.0	61	15-92	14	30
Indeno(1,2,3-cd)pyrene	8270D	62.2	80.0	78	61.9	80.0	77	62-137	1	30
Isophorone	8270D	54.6	80.0	68	48.0	80.0	60	50-116	13	30
N-Nitrosodi-n-propylamine	8270D	66.7	80.0	83	60.7	80.0	76	49-115	9	30
N-Nitrosodiphenylamine	8270D	85.1	80.0	106	75.5	80.0	94	45-123	12	30
Naphthalene	8270D	62.2	80.0	78	54.1	80.0	68	38-99	14	30
Nitrobenzene	8270D	60.1	80.0	75	54.5	80.0	68	46-108	10	30
Pentachlorophenol (PCP)	8270D	67.6	80.0	85	59.1	80.0	74	29-164	14	30
Phenanthrene	8270D	67.1	80.0	84	61.2	80.0	77	58-118	9	30
Phenol	8270D	31.5	80.0	39	32.0	80.0	40	10-113	3	30
Pyrene	8270D	69.2	80.0	86	65.3	80.0	82	61-122	5	30

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Semivolatile Organic Compounds by GC

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

SURROGATE RECOVERY SUMMARY

Organochlorine Pesticides by Gas Chromatography

Analysis Method: 8081B **Extraction Method:** EPA 3510C

		Decachlorobiphenyl	Tetrachloro-m-xylene	
Sample Name	Lab Code	10-164	10-147	
WG-9954-062420-SG-001	R2005470-002	19	60	
WG-9954-062420-SG-002	R2005470-003	44	51	
WG-9954-062420-SG-003	R2005470-004	61	50	
WG-9954-062420-SG-004	R2005470-005	61	55	
WG-9954-062420-SG-005	R2005470-006	20	57	
WG-9954-062420-SG-006	R2005470-007	35	49	
WG-9954-062420-SG-007	R2005470-008	28	56	
Method Blank	RQ2006967-03	58	59	
Lab Control Sample	RQ2006967-04	71	67	
Duplicate Lab Control Sample	RQ2006967-05	60	57	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2006967-03Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analysis Method: 8081B **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	_
4,4'-DDE	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
4,4'-DDT	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
Aldrin	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
Dieldrin	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
Endosulfan I	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
Endosulfan II	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
Endosulfan Sulfate	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
Endrin	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
Endrin Ketone	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
Heptachlor	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
Heptachlor Epoxide	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
Methoxychlor	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
Toxaphene	0.50 U	0.50	0.50	1	07/01/20 16:36	6/30/20	
alpha-BHC	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
alpha-Chlordane	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
beta-BHC	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
delta-BHC	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
gamma-BHC (Lindane)	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
gamma-Chlordane	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	58	10 - 164	07/01/20 16:36	
Tetrachloro-m-xylene	59	10 - 147	07/01/20 16:36	

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/01/20

Sample Matrix: Water

Project:

Duplicate Lab Control Sample Summary Organochlorine Pesticides by Gas Chromatography

Units:ug/L Basis:NA

Service Request: R2005470

Lab Control Sample

Duplicate Lab Control Sample

RQ2006967-04

RQ2006967-05

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
4,4'-DDD	8081B	0.289	0.400	72	0.242	0.400	61	42-159	17	30
4,4'-DDE	8081B	0.287	0.400	72	0.245	0.400	61	47-147	16	30
4,4'-DDT	8081B	0.267	0.400	67	0.224	0.400	56	41-149	18	30
Aldrin	8081B	0.222	0.400	56	0.195	0.400	49	22-137	13	30
Dieldrin	8081B	0.320	0.400	80	0.269	0.400	67	52-144	17	30
Endosulfan I	8081B	0.316	0.400	79	0.268	0.400	67	52-136	16	30
Endosulfan II	8081B	0.323	0.400	81	0.276	0.400	69	57-138	16	30
Endosulfan Sulfate	8081B	0.271	0.400	68	0.229	0.400	57	34-156	17	30
Endrin	8081B	0.310	0.400	78	0.264	0.400	66	56-143	16	30
Endrin Ketone	8081B	0.317	0.400	79	0.269	0.400	67	59-143	16	30
Heptachlor	8081B	0.222	0.400	55	0.196	0.400	49	32-141	12	30
Heptachlor Epoxide	8081B	0.313	0.400	78	0.266	0.400	66	51-143	16	30
Methoxychlor	8081B	0.269	0.400	67	0.240	0.400	60	56-149	11	30
alpha-BHC	8081B	0.301	0.400	75	0.250	0.400	62	36-151	19	30
alpha-Chlordane	8081B	0.303	0.400	76	0.259	0.400	65	50-139	15	30
beta-BHC	8081B	0.326	0.400	82	0.270	0.400	68	55-149	19	30
delta-BHC	8081B	0.287	0.400	72	0.237	0.400	59	29-159	19	30
gamma-BHC (Lindane)	8081B	0.303	0.400	76	0.251	0.400	63	41-149	19	30
gamma-Chlordane	8081B	0.295	0.400	74	0.253	0.400	63	50-140	15	30

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

SURROGATE RECOVERY SUMMARY Polychlorinated Biphenyls (PCBs) by GC

Analysis Method: 8082A **Extraction Method:** EPA 3510C

		Decachlorobiphenyl	Tetrachloro-m-xylene	
Sample Name	Lab Code	10-152	14-129	
WG-9954-062420-SG-001	R2005470-002	21	53	
WG-9954-062420-SG-002	R2005470-003	44	43	
WG-9954-062420-SG-003	R2005470-004	62	44	
WG-9954-062420-SG-004	R2005470-005	59	47	
WG-9954-062420-SG-005	R2005470-006	20	49	
WG-9954-062420-SG-006	R2005470-007	33	42	
WG-9954-062420-SG-007	R2005470-008	26	49	
Method Blank	RQ2006967-03	56	51	
Lab Control Sample	RQ2006967-04	56	49	
Duplicate Lab Control Sample	RQ2006967-05	58	49	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005470

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2006967-03Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analysis Method: 8082A **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	1.0 U	1.0	0.50	1	07/01/20 17:28	6/30/20	
Aroclor 1221	2.0 U	2.0	1.0	1	07/01/20 17:28	6/30/20	
Aroclor 1232	1.0 U	1.0	0.50	1	07/01/20 17:28	6/30/20	
Aroclor 1242	1.0 U	1.0	0.50	1	07/01/20 17:28	6/30/20	
Aroclor 1248	1.0 U	1.0	0.50	1	07/01/20 17:28	6/30/20	
Aroclor 1254	1.0 U	1.0	0.50	1	07/01/20 17:28	6/30/20	
Aroclor 1260	1.0 U	1.0	0.50	1	07/01/20 17:28	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q	
Decachlorobiphenyl	56	10 - 152	07/01/20 17:28		
Tetrachloro-m-xvlene	51	14 - 129	07/01/20 17:28		

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix:

Project:

Water

Duplicate Lab Control Sample Summary Polychlorinated Biphenyls (PCBs) by GC

> Units:ug/L Basis:NA

Service Request: R2005470

Date Analyzed: 07/01/20

Lab Control Sample

Duplicate Lab Control Sample

RQ2006967-04

RQ2006967-05

	Analytical		Spike			Spike		% Rec		RPD
Analyte Name	Method	Result	Amount	% Rec	Result	Amount	% Rec	Limits	RPD	Limit
Aroclor 1016	8082A	2.57	4.00	64	2.65	4.00	66	49-123	3	30
Aroclor 1260	8082A	2.77	4.00	69	2.87	4.00	72	30-120	4	30



Service Request No:R2005520

Ms. Kathy Willy GHD Services Inc. 2055 Niagara Falls Blvd., Niagara Falls, NY 14304

Laboratory Results for: Love Canal:292-402-D02-3100

Dear Ms. Willy,

Enclosed are the results of the sample(s) submitted to our laboratory June 26, 2020 For your reference, these analyses have been assigned our service request number **R2005520**.

All testing was performed according to our laboratory's quality assurance program and met the requirements of the TNI standards except as noted in the case narrative report. Any testing not included in the lab's accreditation is identified on a Non-Certified Analytes report. All results are intended to be considered in their entirety. ALS Environmental is not responsible for use of less than the complete report. Results apply only to the individual samples submitted to the lab for analysis, as listed in the report. The measurement uncertainty of the results included in this report is within that expected when using the prescribed method(s), and represented by Laboratory Control Sample control limits. Any events, such as QC failures or Holding Time exceedances, which may add to the uncertainty are explained in the report narrative or are flagged with qualifiers. The flags are explained in the Report Qualifiers and Definitions page of this report.

Please contact me if you have any questions. My extension is 7471. You may also contact me via email at Brady.Kalkman@alsglobal.com.

Respectfully submitted,

Goody Kullen

ALS Group USA, Corp. dba ALS Environmental

Brady Kalkman Project Manager



Narrative Documents

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com



Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100 Date Received: 06/26/2020

Sample Matrix: Water

CASE NARRATIVE

All analyses were performed consistent with the quality assurance program of ALS Environmental. This report contains analytical results for samples for the Tier level IV requested by the client.

Sample Receipt:

Eight water samples were received for analysis at ALS Environmental on 06/26/2020. Any discrepancies upon initial sample inspection are annotated on the sample receipt and preservation form included within this report. The samples were stored at minimum in accordance with the analytical method requirements.

Semivolatiles by GC/MS:

Method 8270D, 07/01/2020: The upper control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). The field samples analyzed in this sequence did not contain the analyte(s) in question above the Method Reporting Limit (MRL). Since the exceedance equates to a potential high bias, the data quality was not significantly affected and no further corrective action was taken.

Method 8270D, 07/01/2020: The lower control limit for the spike recovery of the Laboratory Control Sample Duplicate (LCSD) was exceeded for one or more analyte. Precision is also outside limits. There were no detections of the analyte(s) in the associated field samples. The LCS/MS/MSD were within limits for these anlytes. The analytes affected are flagged in the LCS Summary.

Method 8270D, 07/01/2020: The lower control limit for the spike recovery of the Matrix Spike/Matrix Spike Duplicate (MS/MSD) was exceeded for one or more analyte. There were no detections of the analyte(s) in the associated field samples. The LCS/LCSD were within limits for theses analytes. The analytes affected are flagged in the MS Summary.

Semivoa GC:

Method 8081B, 07/08/2020: The lower control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). Since there were no detections of the analyte(s) above the MRL in the associated field samples, the quantitation is not affected. The data quality was not significantly affected and no further corrective action was taken. Method 8081B, 07/06/2020: The lower control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). Since there were no detections of the analyte(s) above the MRL in the associated field samples, the quantitation is not affected. The data quality was not significantly affected and no further corrective action was taken. Method 8082A, 07/06/2020: The control limits were exceeded for one or more surrogates in one or more QC samples associated with samples in this report. The associated recoveries of target compounds were in control, indicating the analysis was in control. The surrogate outlier is flagged accordingly. The sample and ms/msd confirm each other for surrogate low. No further corrective action was appropriate.

Volatiles by GC/MS:

Method 8260C, 06/30/2020: The lower control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). Since there were no detections of the analyte(s) above the MRL in the associated field samples, the quantitation is not affected. The data quality was not significantly affected and no further corrective action was taken.

Method 8260C, 06/30/2020: The upper control criterion was exceeded for one or more analytes in the Laboratory Control Sample (LCS). There were no detections of the analyte(s) above the MRL in the associated field samples. The error associated with elevated recovery equates to a high bias. The sample data is not significantly affected. No further corrective action was appropriate.

Method 8260C, 07/02/2020: The upper control criterion was exceeded for one or more analytes in the Laboratory Control Sample (LCS). There were no detections of the analyte(s) above the MRL in the associated field samples. The error associated with



elevated recovery equates to a high bias. The sample data is not significantly affected. No further corrective action was appropriate.

Method 8260C, 07/02/2020: The upper control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). The field samples analyzed in this sequence did not contain the analyte(s) in question above the Method Reporting Limit (MRL). Since the exceedance equates to a potential high bias, the data quality was not significantly affected and no further corrective action was taken.

Method 8260C, 07/07/2020: The upper control criterion was exceeded for one or more analytes in the Laboratory Control Sample (LCS). There were no detections of the analyte(s) above the MRL in the associated field samples. The error associated with elevated recovery equates to a high bias. The sample data is not significantly affected. No further corrective action was appropriate.

Approved by Approved by

Date 07/16/2020



SAMPLE DETECTION SUMMARY

CLIENT ID: WG-9954-062520-SG-013		Lab	ID: R2005	5520-001		
Analyte	Results	Flag	MDL	MRL	Units	Method
Carbon Disulfide	5.4	J	0.42	10	ug/L	8260C
Bis(2-ethylhexyl) Phthalate	3.7	J	0.91	9.1	ug/L	8270D
Aldrin	0.019	J	0.019	0.045	ug/L	8081B
alpha-BHC	0.18		0.019	0.045	ug/L	8081B
delta-BHC	0.28		0.019	0.045	ug/L	8081B
gamma-BHC (Lindane)	0.21		0.019	0.045	ug/L	8081B
CLIENT ID: WG-9954-062520-SG-011		Lab	ID: R2005	5520-002		
Analyte	Results	Flag	MDL	MRL	Units	Method
Carbon Disulfide	2.0	J	0.42	10	ug/L	8260C
alpha-BHC	0.31		0.019	0.045	ug/L	8081B
delta-BHC	0.13		0.019	0.045	ug/L	8081B
gamma-BHC (Lindane)	0.25		0.019	0.045	ug/L	8081B
CLIENT ID: WG-9954-062520-SG-008		Lab	D: R2005	5520-003		
Analyte	Results	Flag	MDL	MRL	Units	Method
4-Methyl-2-pentanone	0.55	J	0.20	10	ug/L	8260C
Chloromethane	0.32	J	0.28	5.0	ug/L	8260C
CLIENT ID: WG-9954-062520-SG-009		Lab	ID: R2005	5520-004		
Analyte	Results	Flag	MDL	MRL	Units	Method
Carbon Disulfide	0.48	J	0.42	10	ug/L	8260C
Chloromethane	0.38	J	0.28	5.0	ug/L	8260C
CLIENT ID: WG-9954-062520-SG-012		Lab	ID: R2005	5520-005		
Analyte	Results	Flag	MDL	MRL	Units	Method
Carbon Disulfide	2.6	J	0.42	10	ug/L	8260C
Chloromethane	0.30	J	0.28	5.0	ug/L	8260C
alpha-BHC	0.033	J	0.019	0.045	ug/L	8081B
delta-BHC	0.068		0.019	0.045	ug/L	8081B
gamma-BHC (Lindane)	0.037	J	0.019	0.045	ug/L	8081B
gamma-Chlordane	0.028	JP	0.019	0.045	ug/L	8081B
CLIENT ID: RB-9954-062520-SG-001		Lab	ID: R2005	5520-006		
Analyte	Results	Flag	MDL	MRL	Units	Method
alpha-BHC	0.079		0.019	0.045	ug/L	8081B
delta-BHC	0.044	J	0.019	0.045	ug/L	8081B
gamma-BHC (Lindane)	0.075		0.019	0.045	ug/L	8081B
CLIENT ID: WG-9954-062520-SG-010		Lab	ID: R2005	5520-008		
Analyte	Results	Flag	MDL	MRL	Units	Method
Carbon Disulfide	0.75	J	0.42	10	ug/L	8260C



Sample Receipt Information

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request:R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

SAMPLE CROSS-REFERENCE

SAMPLE #	CLIENT SAMPLE ID	<u>DATE</u>	<u>TIME</u>
R2005520-001	WG-9954-062520-SG-013	6/25/2020	1315
R2005520-002	WG-9954-062520-SG-011	6/25/2020	1110
R2005520-003	WG-9954-062520-SG-008	6/25/2020	0915
R2005520-004	WG-9954-062520-SG-009	6/25/2020	1005
R2005520-005	WG-9954-062520-SG-012	6/25/2020	1225
R2005520-006	RB-9954-062520-SG-001	6/25/2020	1350
R2005520-007	TB-9954-062520-SG-002	6/25/2020	0000
R2005520-008	WG-9954-062520-SG-010	6/25/2020	1005

CHAIN-OF-CUSTODY/Analytical Request Document The Chain-of-Custody is a LEGAL DOCUMENT. All relevant fields must be completed accurately.

Client to	formation
GLEN SPRINGS HOLDINGS INC	Report To: Kathy Willy
805 97TH STREET	Сору То:
LOVE CANAL	
NIAGARA FALLS, NEW YORK 14304	Invoice To:
Phone: 716-283-0111	PO:
Fax: 716-283-2866	Project Name: LOVE CANAL ANNUAL GW
Email: kathy.willy@ghd.com	Project Number: 9954

Laboratory: ALS	b Information
Laboratory Location: 1 BUILDING 300, SUITE : ROCHESTER, NY 1462	360
ROCHESTER, NY 1462	3
Laboratory Contact: Bi	RADY KALKMAN
Requested Due Date:	TAT: 10
QA/QC Regulrements:	

Event Information ID#: LC ANNUAL GW SAMPLING 2020-02-1 SSOW Ref#: 292-402-999-3100 Sampler Name: S GARDNER,

	Valid Matrix Code WG Groundwater WB Borehote Water					<u> </u>			Sample Cond	iltion
	WS Surface Water		10	ō	Je)				Temp in C	
	SO Soil SE Sediment	ode	ecte	acte .	Ž,	ĝ			Received on ice	Y/N
		ပ	Collected	Collected	PestPCBs(None)	SVOC(none)	ହି		Sealed Cooler	Y/N
		Matrix	Date (Time) st	g	VOA(HCI)	,	Samples Intact	Y/N
Sample Identification		ž	ă	4	" ا	(6)	>	Remarks		
WG-9954-062520-SG-013		wg	06/25/2020	13:15	2	2	3			
WG-9954-062520-SG-011		WG	06/25/2020	11:10	2	2	3			•
WG-9954-062520-SG-008		WG	06/25/2020	09:15	2	2	3			
WG-9954-062520-SG-009		WG	06/25/2020	10:05	2	2	3			
WG-9954-062520-SG-012		WG	06/25/2020	12:25	6	6	9	MS/MSD		
RB-9954-062520-SG-001		WG Q	06/25/2020	1 3 :50	2	2	3			
TB-9954-062520-SG-002		WG Q	06/25/2020	00:00	0	0	3			
WG-9954-062520-SG-010		WG	06/25/2020	10:05	2	2	3			
Total Bottles			18	18	30	Grand Tot	al:66			

SHIPMENT METHOD	NO. OF COOLERS	RELINQUISHED BY:	DATE	TIME	RECIEVED BY:	DATE	TIME
FedEx	2	Shawn Hardner_	6/25/2	150	MINIMUM AVS	696200	10 0125
AIRBILL#:							

CHAIN OF CUSTODY / LABORATORY ANALYSIS REQUEST FORM SR# 004, 005, 006, 007, 008, 009, 010, 1565 Jefferson Road, Bldg 300, Suite 360, Rochester, NY 14623 011, 012, 013 T030477 Phone (585) 288-5380 / FAX (585) 268-8475 S) Enutronmental www.alsglobal.com Project Name: 14D Love Canal:292-402-D02-3100 Project Number: 9954 Annual Long Term Monitoring Company / Address GHD Services Inc. Report To Kathy Willy NUMBER OF CONTAINERS 2055 Niagara Falts Blvd., Suite 3 Niagara Falis NY, 14304 3081B / Pest OC 260C / VOC FP 082A / PCB 270D / SVO Phone # 716-297-2160 Sampler Signature 716-297-2265 Sampler Printed Name Remarks SAMPLING Matrix **CLIENT SAMPLE ID** LABID Date Time Liquid Llquid Liquid Liquid Liquid Liquid Liquid 8. Liquid Liquid 10. Liquid Special Instructions/Comments: Report Requirements Turnaround Requirements Invoice Information RUSH (SURCHARGES APPLY) I. Results Only II. Results + QC Summaries (LCS, DUP, MS/MSD as required) P.O.#_ Standard (3 weeks) III. Results + QC and Cilibration Bill To: Summaries X IV. Data Validation Report REQUESTED FAX DATE with Raw Data EData ______No Requested Report Date Relinquished By: Received By: Relinquished By: Received By: Relinquished By: Received By: Signature Signature Signature Signature Signature Signature Printed Name Printed Name Printed Name Printed Name Printed Name Printed Name

Firm

Date/Time

Firm

Date/Time

Firm

Date/Time

Firm

Date/Time

Date/Time

Firm

Date/Time

AL	s) ~	Cooler 1	Receip	t and P	reservatio	n Che	eck F	orm	Love C	irvices inc. Insi:292-402-00; I	?-3100	
Project/Clie	1 (1)	D		/ Fol	der Number							
-	•	<u>ン</u>	113	// / 101	dei Nuinbei_							
Cooler receive	ed on <u>60262</u>	<u>vao</u>	by: 111/	46	COURÍER:	ALS	UPS (FEDEX	, AETO	CITY CLI	ENT	_
1 Were Cu	stody seals on	outside of coole	r?	N (Y)	5a Perch	lorate s	samples	have requ	ired head	space?	Y N	NA
2 Custody	papers proper	ly completed (in	k, signed)'	? (Y) N	5b Did V	OA via	ls, Alk,o	r Sulfide	have sig*	bubbles?	YN	I NA
3 Did all be	ottles arrive in	good condition (unbroken)? (Ý) N	6 Where	did the	bottles	originate'	· (1	LS/ROC	CLIE	
4 Circle:	Wet Ice Dry	Ice Gel packs	present	? Ŷ N	7 Soil V	OA rec	eived as	: Bull	c Enc	ore 5035	set N	ĪĀ)
3. Temperatur	e Readings	Date: 626d	DAO Tin	ne: <u> ()13(</u>	D:	IR#7(IR#10)	From	Temp Blank	Samp	ole Bottle
Observed Te	mp (°C)	2,8	0.6	o '								
Within 0-6°0	C?	(Y) N	Y	N	YN	Y	N	YN	1	YN	Y	N
If <0°C, wer	e samples froz	en? Y N	Y	N	YN	Y	N	Y	1	YN	Y	N
If out of I	Temperature,	note packing/ic	e conditio	n:	Ice melt	ed P	oorly Pa	cked (des	scribed be	elow)	Same D	ay Rule
	-				pproval Client		-			-		-
•	held in storages placed in storage	e location: orage location:	<i>R-0</i> 02	by definition by	on Graph on	2020 at _ at _	1038		hours of	sampling?	Y	N
· Cooler Bro	eakdown/Prese	rvation Check**	· Date ·	1./24	7020 Time:	10	1416	by:	Đ)			
		labels complete (7-6			NO			_
		bels and tags agr					₹	ES .	NO			
		ontainers used for						ES)	NO			
12. V	Vere 5035 vial	s acceptable (no	extra label	ls, not leak	ring)?		Y	ES	NO		<i>X</i> ₩ <i>A</i>	
		assettes / Tubes			Canisters Pressu	ized		edlar® B		ed	N/A)	
pН	Lot of test	Reagent	Preserved		Received	Exp	Sampl	i i	Vol.	Lot Add	ed	Final
	paper		Yes N	io		1	Adjust	ed	Added			pН
≥12		NaOH				<u> </u>						
_≤2		HNO ₃								<u> </u>		
≤2		H ₂ SO ₄										
<4		NaHSO ₄					1					
5-9		For 608pest			lotify for 3day		ļ			1		
Residual		For CN,			ontact PM to add O ₃ (625, 608,			ŀ				
Chlorine		Phenol, 625,	1 1	1,0202	.03 (023, 000,		l					1 1

CN), ascorbic (phenol).

Bottle lot numbers: Chil Course 040720-18mc Explain all Discrepancies/ Other Comments:

608pest, 522 Na₂S₂O₃

ZnAcetate

HCl

* Trip Blank: 20f3 V/al3

HPROD	BULK
HTR	FLDT ·
SUB	HGFB
ALS	LL3541

Labels secondary reviewed by:_	Ø
PC Secondary Review:	

*significant air bubbles: VOA > 5-6 mm : WC > 1 in. diameter Page 10 of 118

**VOAs and 1664 Not to be tested before analysis.

are checked (not just representatives).

Otherwise, all bottles of all samples with chemical preservatives



Miscellaneous Forms

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com



REPORT QUALIFIERS AND DEFINITIONS

- U Analyte was analyzed for but not detected. The sample quantitation limit has been corrected for dilution and for percent moisture, unless otherwise noted in the case narrative.
- J Estimated value due to either being a Tentatively Identified Compound (TIC) or that the concentration is between the MRL and the MDL. Concentrations are not verified within the linear range of the calibration. For DoD: concentration >40% difference between two GC columns (pesticides/Arclors).
- B Analyte was also detected in the associated method blank at a concentration that may have contributed to the sample result.
- E Inorganics- Concentration is estimated due to the serial dilution was outside control limits.
- E Organics- Concentration has exceeded the calibration range for that specific analysis.
- D Concentration is a result of a dilution, typically a secondary analysis of the sample due to exceeding the calibration range or that a surrogate has been diluted out of the sample and cannot be assessed.
- * Indicates that a quality control parameter has exceeded laboratory limits. Under the õNotesö column of the Form I, this qualifier denotes analysis was performed out of Holding Time.
- H Analysis was performed out of hold time for tests that have an õimmediateö hold time criteria.
- # Spike was diluted out.

- + Correlation coefficient for MSA is <0.995.
- N Inorganics- Matrix spike recovery was outside laboratory limits.
- N Organics- Presumptive evidence of a compound (reported as a TIC) based on the MS library search.
- S Concentration has been determined using Method of Standard Additions (MSA).
- W Post-Digestion Spike recovery is outside control limits and the sample absorbance is <50% of the spike absorbance.
- P Concentration >40% difference between the two GC columns.
- C Confirmed by GC/MS
- Q DoD reports: indicates a pesticide/Aroclor is not confirmed (×100% Difference between two GC columns).
- X See Case Narrative for discussion.
- MRL Method Reporting Limit. Also known as:
- LOQ Limit of Quantitation (LOQ)

 The lowest concentration at which the method analyte may be reliably quantified under the method conditions.
- MDL Method Detection Limit. A statistical value derived from a study designed to provide the lowest concentration that will be detected 99% of the time. Values between the MDL and MRL are estimated (see J qualifier).
- LOD Limit of Detection. A value at or above the MDL which has been verified to be detectable.
- ND Non-Detect. Analyte was not detected at the concentration listed. Same as U qualifier.



Rochester Lab ID # for State Certifications¹

Connecticut ID # PH0556	Maine ID #NY0032	Pennsylvania ID# 68-786
Delaware Approved	New Hampshire ID # 2941	Rhode Island ID # 158
DoD ELAP #65817	New York ID # 10145	Virginia #460167
Florida ID # E87674	North Carolina #676	

¹ Analyses were performed according to our laboratory

NELAP-approved quality assurance program and any applicable state or agency requirements. The test results meet requirements of the current NELAP/TNI standards or state or agency requirements, where applicable, except as noted in the case narrative. Since not all analyte/method/matrix combinations are offered for state/NELAC accreditation, this report may contain results which are not accredited. For a specific list of accredited analytes, contact the laboratory or go to https://www.alsglobal.com/locations/americas/north-america/usa/new-york/rochester-environmental

ALS Laboratory Group

Acronyms

ASTM American Society for Testing and Materials

A2LA American Association for Laboratory Accreditation

CARB California Air Resources Board

CAS Number Chemical Abstract Service registry Number

CFC Chlorofluorocarbon CFU Colony-Forming Unit

DEC Department of Environmental Conservation

DEQ Department of Environmental Quality

DHS Department of Health Services

DOE Department of Ecology DOH Department of Health

EPA U. S. Environmental Protection Agency

ELAP Environmental Laboratory Accreditation Program

GC Gas Chromatography

GC/MS Gas Chromatography/Mass Spectrometry

LUFT Leaking Underground Fuel Tank

M Modified

MCL Maximum Contaminant Level is the highest permissible concentration of a

substance allowed in drinking water as established by the USEPA.

MDL Method Detection Limit
MPN Most Probable Number
MRL Method Reporting Limit

NA Not Applicable NC Not Calculated

NCASI National Council of the Paper Industry for Air and Stream Improvement

ND Not Detected

NIOSH National Institute for Occupational Safety and Health

PQL Practical Quantitation Limit

RCRA Resource Conservation and Recovery Act

SIM Selected Ion Monitoring

TPH Total Petroleum Hydrocarbons

tr Trace level is the concentration of an analyte that is less than the PQL but

greater than or equal to the MDL.

Analyst Summary report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term **Project:**

Monitoring

Sample Name: WG-9954-062520-SG-013

Lab Code: R2005520-001

Sample Matrix: Water

Analysis Method

Analyzed By Extracted/Digested By

Service Request: R2005520

Date Collected: 06/25/20

Date Received: 06/26/20

8081B **KSERCU JMISIUREWICZ** 8082A **KSERCU BALLGEIER** 8260C **FNAEGLER**

8270D **KSERCU JMISIUREWICZ**

Sample Name: WG-9954-062520-SG-011 **Date Collected:** 06/25/20

Lab Code: R2005520-002 **Date Received:** 06/26/20

Sample Matrix: Water

Extracted/Digested By Analyzed By Analysis Method

8081B **KSERCU JMISIUREWICZ** 8082A **KSERCU BALLGEIER** 8260C **FNAEGLER** 8270D **KSERCU JMISIUREWICZ**

Sample Name: WG-9954-062520-SG-008 **Date Collected:** 06/25/20

Lab Code: R2005520-003 **Date Received:** 06/26/20 **Sample Matrix:** Water

Analyzed By Extracted/Digested By Analysis Method

8081B **KSERCU JMISIUREWICZ** 8082A **KSERCU BALLGEIER** 8260C **FNAEGLER KSERCU** 8270D **JMISIUREWICZ**

Sample Name: WG-9954-062520-SG-009 Date Collected: 06/25/20 Lab Code: R2005520-004 **Date Received:** 06/26/20

Sample Matrix: Water

Analyzed By Extracted/Digested By Analysis Method 8081B **KSERCU JMISIUREWICZ**

Printed 7/16/2020 11:08:07 AM

Page 14 of 118

Superset Reference:20-0000554598 rev 00

Analyst Summary report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Monitoring

Sample Name: WG-9954-062520-SG-009

Lab Code: R2005520-004

Sample Matrix: Water

Analysis Method

Extracted/Digested By Analyzed By

Service Request: R2005520

Date Collected: 06/25/20

Date Received: 06/26/20

8082A KSERCU BALLGEIER

8260C FNAEGLER

8270D KSERCU JMISIUREWICZ

Sample Name: WG-9954-062520-SG-012 **Date Collected:** 06/25/20

Lab Code: R2005520-005 **Date Received:** 06/26/20

Sample Matrix: Water

Analysis Method Extracted/Digested By Analyzed By

8081B KSERCU JMISIUREWICZ 8082A KSERCU BALLGEIER

8260C RSERCU BALLUEIER
FNAEGLER

8270D KSERCU JMISIUREWICZ

Sample Name: RB-9954-062520-SG-001 **Date Collected:** 06/25/20

Lab Code: R2005520-006 Date Received: 06/26/20 Sample Matrix: Water

Analysis Method Extracted/Digested By Analyzed By

8081B KSERCU BALLGEIER 8082A KSERCU BALLGEIER

8260C RSERCE BALEGEER
FNAEGLER

8270D KSERCU JMISIUREWICZ

Sample Name: TB-9954-062520-SG-002 **Date Collected:** 06/25/20

Lab Code: R2005520-007 **Date Received:** 06/26/20

Sample Matrix: Water

Analysis Method Extracted/Digested By Analyzed By

8260C FNAEGLER

Analyst Summary report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Monitoring

Sample Name: WG-9954-062520-SG-010

Lab Code: R2005520-008

Sample Matrix: Water

Service Request: R2005520

Date Collected: 06/25/20

Date Received: 06/26/20

Analysis Method	Extracted/Digested By	Analyzed By
8081B	KSERCU	BALLGEIER
8082A	KSERCU	BALLGEIER
8260C		FNAEGLER
8270D	KSERCU	JMISIUREWICZ



INORGANIC PREPARATION METHODS

The preparation methods associated with this report are found in these tables unless discussed in the case narrative.

Water/Liquid Matrix

Analytical Method	Preparation Method
200.7	200.2
200.8	200.2
6010C	3005A/3010A
6020A	ILM05.3
9034 Sulfide Acid Soluble	9030B
SM 4500-CN-E Residual Cyanide	SM 4500-CN-G
SM 4500-CN-E WAD Cyanide	SM 4500-CN-I

Solid/Soil/Non-Aqueous Matrix

Analytical Method	Preparation
	Method
6010C	3050B
6020A	3050B
6010C TCLP (1311)	3005A/3010A
extract	
6010 SPLP (1312) extract	3005A/3010A
7199	3060A
300.0 Anions/ 350.1/	DI extraction
353.2/ SM 2320B/ SM	
5210B/ 9056A Anions	
For analytical methods not listed,	
method is the same as the analytic	cal method
I ICICICILC.	



Sample Results

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com



Volatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 13:15

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-013
 Units: ug/L

 Lab Code:
 R2005520-001
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/02/20 20:20	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/02/20 20:20	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/02/20 20:20	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/02/20 20:20	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/02/20 20:20	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/02/20 20:20	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/02/20 20:20	
2-Butanone (MEK)	10 U	10	0.78	1	07/02/20 20:20	
2-Hexanone	10 U	10	0.20	1	07/02/20 20:20	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/02/20 20:20	
Acetone	10 U	10	5.0	1	07/02/20 20:20	
Benzene	5.0 U	5.0	0.20	1	07/02/20 20:20	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/02/20 20:20	
Bromoform	5.0 U	5.0	0.25	1	07/02/20 20:20	
Bromomethane	5.0 U	5.0	0.70	1	07/02/20 20:20	
Carbon Disulfide	5.4 J	10	0.42	1	07/02/20 20:20	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/02/20 20:20	
Chlorobenzene	5.0 U	5.0	0.20	1	07/02/20 20:20	
Chloroethane	5.0 U	5.0	0.23	1	07/02/20 20:20	
Chloroform	5.0 U	5.0	0.24	1	07/02/20 20:20	
Chloromethane	5.0 U	5.0	0.28	1	07/02/20 20:20	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/02/20 20:20	
Dichloromethane	5.0 U	5.0	0.65	1	07/02/20 20:20	
Ethylbenzene	5.0 U	5.0	0.20	1	07/02/20 20:20	
Styrene	5.0 U	5.0	0.20	1	07/02/20 20:20	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/02/20 20:20	
Toluene	5.0 U	5.0	0.20	1	07/02/20 20:20	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/02/20 20:20	
Vinyl Acetate	10 U	10	1.1	1	07/02/20 20:20	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/02/20 20:20	
Xylenes, Total	5.0 U	5.0	0.23	1	07/02/20 20:20	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/02/20 20:20	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/02/20 20:20	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/02/20 20:20	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/02/20 20:20	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 13:15

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-013
 Units: ug/L

 Lab Code:
 R2005520-001
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	92	85 - 122	07/02/20 20:20	
Dibromofluoromethane	97	89 - 119	07/02/20 20:20	
Toluene-d8	97	87 - 121	07/02/20 20:20	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 13:15

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-013
 Units: ug/L

 Lab Code:
 R2005520-001
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 11:10

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-011
 Units: ug/L

 Lab Code:
 R2005520-002
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/02/20 20:42	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/02/20 20:42	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/02/20 20:42	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/02/20 20:42	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/02/20 20:42	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/02/20 20:42	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/02/20 20:42	
2-Butanone (MEK)	10 U	10	0.78	1	07/02/20 20:42	
2-Hexanone	10 U	10	0.20	1	07/02/20 20:42	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/02/20 20:42	
Acetone	10 U	10	5.0	1	07/02/20 20:42	
Benzene	5.0 U	5.0	0.20	1	07/02/20 20:42	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/02/20 20:42	
Bromoform	5.0 U	5.0	0.25	1	07/02/20 20:42	
Bromomethane	5.0 U	5.0	0.70	1	07/02/20 20:42	
Carbon Disulfide	2.0 J	10	0.42	1	07/02/20 20:42	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/02/20 20:42	
Chlorobenzene	5.0 U	5.0	0.20	1	07/02/20 20:42	
Chloroethane	5.0 U	5.0	0.23	1	07/02/20 20:42	
Chloroform	5.0 U	5.0	0.24	1	07/02/20 20:42	
Chloromethane	5.0 U	5.0	0.28	1	07/02/20 20:42	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/02/20 20:42	
Dichloromethane	5.0 U	5.0	0.65	1	07/02/20 20:42	
Ethylbenzene	5.0 U	5.0	0.20	1	07/02/20 20:42	
Styrene	5.0 U	5.0	0.20	1	07/02/20 20:42	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/02/20 20:42	
Toluene	5.0 U	5.0	0.20	1	07/02/20 20:42	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/02/20 20:42	
Vinyl Acetate	10 U	10	1.1	1	07/02/20 20:42	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/02/20 20:42	
Xylenes, Total	5.0 U	5.0	0.23	1	07/02/20 20:42	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/02/20 20:42	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/02/20 20:42	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/02/20 20:42	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/02/20 20:42	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 11:10

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-011
 Units: ug/L

 Lab Code:
 R2005520-002
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	92	85 - 122	07/02/20 20:42	
Dibromofluoromethane	100	89 - 119	07/02/20 20:42	
Toluene-d8	98	87 - 121	07/02/20 20:42	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 11:10

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

Sample Name: WG-9954-062520-SG-011 **Units:** ug/L

Lab Code: R2005520-002 **Basis:** NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 09:15

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-008
 Units: ug/L

 Lab Code:
 R2005520-003
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/01/20 06:09	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/01/20 06:09	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/01/20 06:09	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/01/20 06:09	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/01/20 06:09	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/01/20 06:09	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/01/20 06:09	
2-Butanone (MEK)	10 U	10	0.78	1	07/01/20 06:09	
2-Hexanone	10 U	10	0.20	1	07/01/20 06:09	
4-Methyl-2-pentanone	0.55 J	10	0.20	1	07/01/20 06:09	
Acetone	10 U	10	5.0	1	07/01/20 06:09	
Benzene	5.0 U	5.0	0.20	1	07/01/20 06:09	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/01/20 06:09	
Bromoform	5.0 U	5.0	0.25	1	07/01/20 06:09	
Bromomethane	5.0 U	5.0	0.70	1	07/01/20 06:09	
Carbon Disulfide	10 U	10	0.42	1	07/01/20 06:09	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/01/20 06:09	
Chlorobenzene	5.0 U	5.0	0.20	1	07/01/20 06:09	
Chloroethane	5.0 U	5.0	0.23	1	07/01/20 06:09	
Chloroform	5.0 U	5.0	0.24	1	07/01/20 06:09	
Chloromethane	0.32 J	5.0	0.28	1	07/01/20 06:09	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/01/20 06:09	
Dichloromethane	5.0 U	5.0	0.65	1	07/01/20 06:09	
Ethylbenzene	5.0 U	5.0	0.20	1	07/01/20 06:09	
Styrene	5.0 U	5.0	0.20	1	07/01/20 06:09	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/01/20 06:09	
Toluene	5.0 U	5.0	0.20	1	07/01/20 06:09	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/01/20 06:09	
Vinyl Acetate	10 U	10	1.1	1	07/01/20 06:09	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/01/20 06:09	
Xylenes, Total	5.0 U	5.0	0.23	1	07/01/20 06:09	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/01/20 06:09	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/01/20 06:09	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/01/20 06:09	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/01/20 06:09	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 09:15

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-008
 Units: ug/L

 Lab Code:
 R2005520-003
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	90	85 - 122	07/01/20 06:09	
Dibromofluoromethane	92	89 - 119	07/01/20 06:09	
Toluene-d8	93	87 - 121	07/01/20 06:09	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 09:15 **Project:**

Monitoring

Date Received: 06/26/20 10:25 **Sample Matrix:** Water

Sample Name: WG-9954-062520-SG-008 Units: ug/L Lab Code:

R2005520-003 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

Result CAS# **Compound Identification** Q RTug/L

No Tentatively Identified Compounds

Detected

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 10:05

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-009
 Units: ug/L

 Lab Code:
 R2005520-004
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/02/20 21:04	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/02/20 21:04	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/02/20 21:04	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/02/20 21:04	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/02/20 21:04	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/02/20 21:04	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/02/20 21:04	
2-Butanone (MEK)	10 U	10	0.78	1	07/02/20 21:04	
2-Hexanone	10 U	10	0.20	1	07/02/20 21:04	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/02/20 21:04	
Acetone	10 U	10	5.0	1	07/02/20 21:04	
Benzene	5.0 U	5.0	0.20	1	07/02/20 21:04	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/02/20 21:04	
Bromoform	5.0 U	5.0	0.25	1	07/02/20 21:04	
Bromomethane	5.0 U	5.0	0.70	1	07/02/20 21:04	
Carbon Disulfide	0.48 J	10	0.42	1	07/02/20 21:04	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/02/20 21:04	
Chlorobenzene	5.0 U	5.0	0.20	1	07/02/20 21:04	
Chloroethane	5.0 U	5.0	0.23	1	07/02/20 21:04	
Chloroform	5.0 U	5.0	0.24	1	07/02/20 21:04	
Chloromethane	0.38 J	5.0	0.28	1	07/02/20 21:04	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/02/20 21:04	
Dichloromethane	5.0 U	5.0	0.65	1	07/02/20 21:04	
Ethylbenzene	5.0 U	5.0	0.20	1	07/02/20 21:04	
Styrene	5.0 U	5.0	0.20	1	07/02/20 21:04	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/02/20 21:04	
Toluene	5.0 U	5.0	0.20	1	07/02/20 21:04	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/02/20 21:04	
Vinyl Acetate	10 U	10	1.1	1	07/02/20 21:04	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/02/20 21:04	
Xylenes, Total	5.0 U	5.0	0.23	1	07/02/20 21:04	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/02/20 21:04	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/02/20 21:04	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/02/20 21:04	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/02/20 21:04	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 10:05

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-009
 Units: ug/L

 Lab Code:
 R2005520-004
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	91	85 - 122	07/02/20 21:04	
Dibromofluoromethane	96	89 - 119	07/02/20 21:04	
Toluene-d8	96	87 - 121	07/02/20 21:04	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 10:05

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-009
 Units: ug/L

 Lab Code:
 R2005520-004
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Tentatively Identified Compounds

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 12:25

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-012
 Units: ug/L

 Lab Code:
 R2005520-005
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/02/20 21:26	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/02/20 21:26	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/02/20 21:26	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/02/20 21:26	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/02/20 21:26	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/02/20 21:26	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/02/20 21:26	
2-Butanone (MEK)	10 U	10	0.78	1	07/02/20 21:26	
2-Hexanone	10 U	10	0.20	1	07/02/20 21:26	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/02/20 21:26	
Acetone	10 U	10	5.0	1	07/02/20 21:26	
Benzene	5.0 U	5.0	0.20	1	07/02/20 21:26	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/02/20 21:26	
Bromoform	5.0 U	5.0	0.25	1	07/02/20 21:26	
Bromomethane	5.0 U	5.0	0.70	1	07/02/20 21:26	
Carbon Disulfide	2.6 J	10	0.42	1	07/02/20 21:26	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/02/20 21:26	
Chlorobenzene	5.0 U	5.0	0.20	1	07/02/20 21:26	
Chloroethane	5.0 U	5.0	0.23	1	07/02/20 21:26	
Chloroform	5.0 U	5.0	0.24	1	07/02/20 21:26	
Chloromethane	0.30 J	5.0	0.28	1	07/02/20 21:26	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/02/20 21:26	
Dichloromethane	5.0 U	5.0	0.65	1	07/02/20 21:26	
Ethylbenzene	5.0 U	5.0	0.20	1	07/02/20 21:26	
Styrene	5.0 U	5.0	0.20	1	07/02/20 21:26	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/02/20 21:26	
Toluene	5.0 U	5.0	0.20	1	07/02/20 21:26	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/02/20 21:26	
Vinyl Acetate	10 U	10	1.1	1	07/02/20 21:26	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/02/20 21:26	
Xylenes, Total	5.0 U	5.0	0.23	1	07/02/20 21:26	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/02/20 21:26	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/02/20 21:26	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/02/20 21:26	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	11	07/02/20 21:26	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 12:25

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-012
 Units: ug/L

 Lab Code:
 R2005520-005
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	93	85 - 122	07/02/20 21:26	
Dibromofluoromethane	98	89 - 119	07/02/20 21:26	
Toluene-d8	98	87 - 121	07/02/20 21:26	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 12:25 **Project:**

Monitoring

Sample Matrix: Date Received: 06/26/20 10:25 Water

Sample Name: WG-9954-062520-SG-012 Units: ug/L

Lab Code: R2005520-005 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown	1.32	7.5	J	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 13:50

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 RB-9954-062520-SG-001
 Units: ug/L

 Lab Code:
 R2005520-006
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	06/30/20 23:52	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	06/30/20 23:52	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	06/30/20 23:52	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	06/30/20 23:52	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	06/30/20 23:52	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	06/30/20 23:52	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	06/30/20 23:52	
2-Butanone (MEK)	10 U	10	0.78	1	06/30/20 23:52	
2-Hexanone	10 U	10	0.20	1	06/30/20 23:52	
4-Methyl-2-pentanone	10 U	10	0.20	1	06/30/20 23:52	
Acetone	10 U	10	5.0	1	06/30/20 23:52	
Benzene	5.0 U	5.0	0.20	1	06/30/20 23:52	
Bromodichloromethane	5.0 U	5.0	0.20	1	06/30/20 23:52	
Bromoform	5.0 U	5.0	0.25	1	06/30/20 23:52	
Bromomethane	5.0 U	5.0	0.70	1	06/30/20 23:52	
Carbon Disulfide	10 U	10	0.42	1	06/30/20 23:52	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	06/30/20 23:52	
Chlorobenzene	5.0 U	5.0	0.20	1	06/30/20 23:52	
Chloroethane	5.0 U	5.0	0.23	1	06/30/20 23:52	
Chloroform	5.0 U	5.0	0.24	1	06/30/20 23:52	
Chloromethane	5.0 U	5.0	0.28	1	06/30/20 23:52	
Dibromochloromethane	5.0 U	5.0	0.20	1	06/30/20 23:52	
Dichloromethane	5.0 U	5.0	0.65	1	06/30/20 23:52	
Ethylbenzene	5.0 U	5.0	0.20	1	06/30/20 23:52	
Styrene	5.0 U	5.0	0.20	1	06/30/20 23:52	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	06/30/20 23:52	
Toluene	5.0 U	5.0	0.20	1	06/30/20 23:52	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	06/30/20 23:52	
Vinyl Acetate	10 U	10	1.1	1	06/30/20 23:52	
Vinyl Chloride	5.0 U	5.0	0.20	1	06/30/20 23:52	
Xylenes, Total	5.0 U	5.0	0.23	1	06/30/20 23:52	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	06/30/20 23:52	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	06/30/20 23:52	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	06/30/20 23:52	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	06/30/20 23:52	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 13:50

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 RB-9954-062520-SG-001
 Units: ug/L

 Lab Code:
 R2005520-006
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	90	85 - 122	06/30/20 23:52	
Dibromofluoromethane	89	89 - 119	06/30/20 23:52	
Toluene-d8	92	87 - 121	06/30/20 23:52	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 13:50 **Project:**

Monitoring

Date Received: 06/26/20 10:25 **Sample Matrix:** Water

Sample Name: RB-9954-062520-SG-001 Units: ug/L Lab Code: Basis: NA R2005520-006

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

Result

CAS# **Compound Identification** Q RTug/L

No Tentatively Identified Compounds

Detected

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 00:00

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 TB-9954-062520-SG-002
 Units: ug/L

 Lab Code:
 R2005520-007
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	06/30/20 23:30	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	06/30/20 23:30	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	06/30/20 23:30	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	06/30/20 23:30	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	06/30/20 23:30	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	06/30/20 23:30	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	06/30/20 23:30	
2-Butanone (MEK)	10 U	10	0.78	1	06/30/20 23:30	
2-Hexanone	10 U	10	0.20	1	06/30/20 23:30	
4-Methyl-2-pentanone	10 U	10	0.20	1	06/30/20 23:30	
Acetone	10 U	10	5.0	1	06/30/20 23:30	
Benzene	5.0 U	5.0	0.20	1	06/30/20 23:30	
Bromodichloromethane	5.0 U	5.0	0.20	1	06/30/20 23:30	
Bromoform	5.0 U	5.0	0.25	1	06/30/20 23:30	
Bromomethane	5.0 U	5.0	0.70	1	06/30/20 23:30	
Carbon Disulfide	10 U	10	0.42	1	06/30/20 23:30	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	06/30/20 23:30	
Chlorobenzene	5.0 U	5.0	0.20	1	06/30/20 23:30	
Chloroethane	5.0 U	5.0	0.23	1	06/30/20 23:30	
Chloroform	5.0 U	5.0	0.24	1	06/30/20 23:30	
Chloromethane	5.0 U	5.0	0.28	1	06/30/20 23:30	
Dibromochloromethane	5.0 U	5.0	0.20	1	06/30/20 23:30	
Dichloromethane	5.0 U	5.0	0.65	1	06/30/20 23:30	
Ethylbenzene	5.0 U	5.0	0.20	1	06/30/20 23:30	
Styrene	5.0 U	5.0	0.20	1	06/30/20 23:30	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	06/30/20 23:30	
Toluene	5.0 U	5.0	0.20	1	06/30/20 23:30	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	06/30/20 23:30	
Vinyl Acetate	10 U	10	1.1	1	06/30/20 23:30	
Vinyl Chloride	5.0 U	5.0	0.20	1	06/30/20 23:30	
Xylenes, Total	5.0 U	5.0	0.23	1	06/30/20 23:30	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	06/30/20 23:30	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	06/30/20 23:30	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	06/30/20 23:30	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	11	06/30/20 23:30	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 00:00

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 TB-9954-062520-SG-002
 Units: ug/L

 Lab Code:
 R2005520-007
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	91	85 - 122	06/30/20 23:30	
Dibromofluoromethane	92	89 - 119	06/30/20 23:30	
Toluene-d8	95	87 - 121	06/30/20 23:30	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 00:00

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

Sample Name: TB-9954-062520-SG-002 **Units:** ug/L

Lab Code: R2005520-007 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

Result

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 10:05

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-010
 Units: ug/L

 Lab Code:
 R2005520-008
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/07/20 22:08	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/07/20 22:08	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/07/20 22:08	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/07/20 22:08	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/07/20 22:08	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/07/20 22:08	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/07/20 22:08	
2-Butanone (MEK)	10 U	10	0.78	1	07/07/20 22:08	
2-Hexanone	10 U	10	0.20	1	07/07/20 22:08	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/07/20 22:08	
Acetone	10 U	10	5.0	1	07/07/20 22:08	
Benzene	5.0 U	5.0	0.20	1	07/07/20 22:08	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/07/20 22:08	
Bromoform	5.0 U	5.0	0.25	1	07/07/20 22:08	
Bromomethane	5.0 U	5.0	0.70	1	07/07/20 22:08	
Carbon Disulfide	0.75 J	10	0.42	1	07/07/20 22:08	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/07/20 22:08	
Chlorobenzene	5.0 U	5.0	0.20	1	07/07/20 22:08	
Chloroethane	5.0 U	5.0	0.23	1	07/07/20 22:08	
Chloroform	5.0 U	5.0	0.24	1	07/07/20 22:08	
Chloromethane	5.0 U	5.0	0.28	1	07/07/20 22:08	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/07/20 22:08	
Dichloromethane	5.0 U	5.0	0.65	1	07/07/20 22:08	
Ethylbenzene	5.0 U	5.0	0.20	1	07/07/20 22:08	
Styrene	5.0 U	5.0	0.20	1	07/07/20 22:08	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/07/20 22:08	
Toluene	5.0 U	5.0	0.20	1	07/07/20 22:08	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/07/20 22:08	
Vinyl Acetate	10 U	10	1.1	1	07/07/20 22:08	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/07/20 22:08	
Xylenes, Total	5.0 U	5.0	0.23	1	07/07/20 22:08	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/07/20 22:08	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/07/20 22:08	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/07/20 22:08	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/07/20 22:08	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 10:05

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-010
 Units: ug/L

 Lab Code:
 R2005520-008
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	107	85 - 122	07/07/20 22:08	
Dibromofluoromethane	109	89 - 119	07/07/20 22:08	
Toluene-d8	107	87 - 121	07/07/20 22:08	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 10:05

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-010
 Units: ug/L

 Lab Code:
 R2005520-008
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

			Result	
CAS#	Compound Identification	RT	ug/L	Q
007446-09-5	Sulfur dioxide	1.23	28.9	JN



Semivolatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 13:15

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-013
 Units: ug/L

 Lab Code:
 R2005520-001
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/01/20 19:49	6/30/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/01/20 19:49	6/30/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/01/20 19:49	6/30/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/01/20 19:49	6/30/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/01/20 19:49	6/30/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/01/20 19:49	6/30/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/01/20 19:49	6/30/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/01/20 19:49	6/30/20	
2,4-Dinitrophenol	45 U	45	19	1	07/01/20 19:49	6/30/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/01/20 19:49	6/30/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/01/20 19:49	6/30/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/01/20 19:49	6/30/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/01/20 19:49	6/30/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/01/20 19:49	6/30/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/01/20 19:49	6/30/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/01/20 19:49	6/30/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/01/20 19:49	6/30/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/01/20 19:49	6/30/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/01/20 19:49	6/30/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/01/20 19:49	6/30/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/01/20 19:49	6/30/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/01/20 19:49	6/30/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/01/20 19:49	6/30/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/01/20 19:49	6/30/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/01/20 19:49	6/30/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/01/20 19:49	6/30/20	
4-Nitrophenol	45 U	45	5.8	1	07/01/20 19:49	6/30/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/01/20 19:49	6/30/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/01/20 19:49	6/30/20	
Anthracene	9.1 U	9.1	1.2	1	07/01/20 19:49	6/30/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/01/20 19:49	6/30/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/01/20 19:49	6/30/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/01/20 19:49	6/30/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/01/20 19:49	6/30/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/01/20 19:49	6/30/20	
Benzoic Acid	91 U	91	33	1	07/01/20 19:49	6/30/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/01/20 19:49	6/30/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/01/20 19:49	6/30/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/01/20 19:49	6/30/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	07/01/20 19:49	6/30/20	
Bis(2-ethylhexyl) Phthalate	3.7 J	9.1	0.91	1	07/01/20 19:49	6/30/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	07/01/20 19:49	6/30/20	
Chrysene	9.1 U	9.1	1.1	1	07/01/20 19:49	6/30/20	

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Superset Reference: 20-0000554598 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 13:15

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-013
 Units: ug/L

 Lab Code:
 R2005520-001
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/01/20 19:49	6/30/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/01/20 19:49	6/30/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/01/20 19:49	6/30/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/01/20 19:49	6/30/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/01/20 19:49	6/30/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/01/20 19:49	6/30/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/01/20 19:49	6/30/20	
Fluorene	9.1 U	9.1	1.2	1	07/01/20 19:49	6/30/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/01/20 19:49	6/30/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/01/20 19:49	6/30/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/01/20 19:49	6/30/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/01/20 19:49	6/30/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/01/20 19:49	6/30/20	
Isophorone	9.1 U	9.1	1.3	1	07/01/20 19:49	6/30/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/01/20 19:49	6/30/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/01/20 19:49	6/30/20	
Naphthalene	9.1 U	9.1	1.1	1	07/01/20 19:49	6/30/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/01/20 19:49	6/30/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/01/20 19:49	6/30/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/01/20 19:49	6/30/20	
Phenol	9.1 U	9.1	0.91	1	07/01/20 19:49	6/30/20	
Pyrene	9.1 U	9.1	1.3	1	07/01/20 19:49	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	104	35 - 141	07/01/20 19:49	
2-Fluorobiphenyl	69	31 - 118	07/01/20 19:49	
2-Fluorophenol	38	10 - 105	07/01/20 19:49	
Nitrobenzene-d5	61	31 - 110	07/01/20 19:49	
Phenol-d6	32	10 - 107	07/01/20 19:49	
p-Terphenyl-d14	92	10 - 165	07/01/20 19:49	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown hydrocarbon	11.35	6.0	J	
	unknown hydrocarbon	11.89	9.6	J	
	unknown hydrocarbon	12.48	12	J	
	unknown hydrocarbon	13.13	13	J	
	unknown hydrocarbon	13.83	11	J	
	unknown hydrocarbon	14.58	13	J	
	unknown hydrocarbon	15.37	8.5	J	

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Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 13:15

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-013
 Units: ug/L

 Lab Code:
 R2005520-001
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown hydrocarbon	16.11	8.0	J	
000112-05-0	Nonanoic acid	6.09	4.5	JN	
013798-23-7	Sulfur	7.83	42	JN	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 11:10

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-011
 Units: ug/L

 Lab Code:
 R2005520-002
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/01/20 20:16	6/30/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/01/20 20:16	6/30/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/01/20 20:16	6/30/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/01/20 20:16	6/30/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/01/20 20:16	6/30/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/01/20 20:16	6/30/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/01/20 20:16	6/30/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/01/20 20:16	6/30/20	
2,4-Dinitrophenol	45 U	45	19	1	07/01/20 20:16	6/30/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/01/20 20:16	6/30/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/01/20 20:16	6/30/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/01/20 20:16	6/30/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/01/20 20:16	6/30/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/01/20 20:16	6/30/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/01/20 20:16	6/30/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/01/20 20:16	6/30/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/01/20 20:16	6/30/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/01/20 20:16	6/30/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/01/20 20:16	6/30/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/01/20 20:16	6/30/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/01/20 20:16	6/30/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/01/20 20:16	6/30/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/01/20 20:16	6/30/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/01/20 20:16	6/30/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/01/20 20:16	6/30/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/01/20 20:16	6/30/20	
4-Nitrophenol	45 U	45	5.8	1	07/01/20 20:16	6/30/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/01/20 20:16	6/30/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/01/20 20:16	6/30/20	
Anthracene	9.1 U	9.1	1.2	1	07/01/20 20:16	6/30/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/01/20 20:16	6/30/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/01/20 20:16	6/30/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/01/20 20:16	6/30/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/01/20 20:16	6/30/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/01/20 20:16	6/30/20	
Benzoic Acid	91 U	91	33	1	07/01/20 20:16	6/30/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/01/20 20:16	6/30/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/01/20 20:16	6/30/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/01/20 20:16	6/30/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	07/01/20 20:16	6/30/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	07/01/20 20:16	6/30/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	07/01/20 20:16	6/30/20	
Chrysene	9.1 U	9.1	1.1	1	07/01/20 20:16	6/30/20	

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Superset Reference:20-0000554598 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 11:10

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-011
 Units: ug/L

 Lab Code:
 R2005520-002
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/01/20 20:16	6/30/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/01/20 20:16	6/30/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/01/20 20:16	6/30/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/01/20 20:16	6/30/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/01/20 20:16	6/30/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/01/20 20:16	6/30/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/01/20 20:16	6/30/20	
Fluorene	9.1 U	9.1	1.2	1	07/01/20 20:16	6/30/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/01/20 20:16	6/30/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/01/20 20:16	6/30/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/01/20 20:16	6/30/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/01/20 20:16	6/30/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/01/20 20:16	6/30/20	
Isophorone	9.1 U	9.1	1.3	1	07/01/20 20:16	6/30/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/01/20 20:16	6/30/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/01/20 20:16	6/30/20	
Naphthalene	9.1 U	9.1	1.1	1	07/01/20 20:16	6/30/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/01/20 20:16	6/30/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/01/20 20:16	6/30/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/01/20 20:16	6/30/20	
Phenol	9.1 U	9.1	0.91	1	07/01/20 20:16	6/30/20	
Pyrene	9.1 U	9.1	1.3	1	07/01/20 20:16	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	93	35 - 141	07/01/20 20:16	
2-Fluorobiphenyl	70	31 - 118	07/01/20 20:16	
2-Fluorophenol	43	10 - 105	07/01/20 20:16	
Nitrobenzene-d5	68	31 - 110	07/01/20 20:16	
Phenol-d6	30	10 - 107	07/01/20 20:16	
p-Terphenyl-d14	93	10 - 165	07/01/20 20:16	

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
013798-23-7	Sulfur	7.77	8.0	JN	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 09:15

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-008
 Units: ug/L

 Lab Code:
 R2005520-003
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/01/20 20:44	6/30/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/01/20 20:44	6/30/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/01/20 20:44	6/30/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/01/20 20:44	6/30/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/01/20 20:44	6/30/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/01/20 20:44	6/30/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/01/20 20:44	6/30/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/01/20 20:44	6/30/20	
2,4-Dinitrophenol	45 U	45	19	1	07/01/20 20:44	6/30/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/01/20 20:44	6/30/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/01/20 20:44	6/30/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/01/20 20:44	6/30/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/01/20 20:44	6/30/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/01/20 20:44	6/30/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/01/20 20:44	6/30/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/01/20 20:44	6/30/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/01/20 20:44	6/30/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/01/20 20:44	6/30/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/01/20 20:44	6/30/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/01/20 20:44	6/30/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/01/20 20:44	6/30/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/01/20 20:44	6/30/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/01/20 20:44	6/30/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/01/20 20:44	6/30/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/01/20 20:44	6/30/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/01/20 20:44	6/30/20	
4-Nitrophenol	45 U	45	5.8	1	07/01/20 20:44	6/30/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/01/20 20:44	6/30/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/01/20 20:44	6/30/20	
Anthracene	9.1 U	9.1	1.2	1	07/01/20 20:44	6/30/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/01/20 20:44	6/30/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/01/20 20:44	6/30/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/01/20 20:44	6/30/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/01/20 20:44	6/30/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/01/20 20:44	6/30/20	
Benzoic Acid	91 U	91	33	1	07/01/20 20:44	6/30/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/01/20 20:44	6/30/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/01/20 20:44	6/30/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/01/20 20:44	6/30/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	07/01/20 20:44	6/30/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	07/01/20 20:44	6/30/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	07/01/20 20:44	6/30/20	
Chrysene	9.1 U	9.1	1.1	1	07/01/20 20:44	6/30/20	

Printed 7/16/2020 11:08:28 AM

Superset Reference: 20-0000554598 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 09:15

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-008
 Units: ug/L

 Lab Code:
 R2005520-003
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed D	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/01/20 20:44	6/30/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/01/20 20:44	6/30/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/01/20 20:44	6/30/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/01/20 20:44	6/30/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/01/20 20:44	6/30/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/01/20 20:44	6/30/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/01/20 20:44	6/30/20	
Fluorene	9.1 U	9.1	1.2	1	07/01/20 20:44	6/30/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/01/20 20:44	6/30/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/01/20 20:44	6/30/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/01/20 20:44	6/30/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/01/20 20:44	6/30/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/01/20 20:44	6/30/20	
Isophorone	9.1 U	9.1	1.3	1	07/01/20 20:44	6/30/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/01/20 20:44	6/30/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/01/20 20:44	6/30/20	
Naphthalene	9.1 U	9.1	1.1	1	07/01/20 20:44	6/30/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/01/20 20:44	6/30/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/01/20 20:44	6/30/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/01/20 20:44	6/30/20	
Phenol	9.1 U	9.1	0.91	1	07/01/20 20:44	6/30/20	
Pyrene	9.1 U	9.1	1.3	1	07/01/20 20:44	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	95	35 - 141	07/01/20 20:44	
2-Fluorobiphenyl	74	31 - 118	07/01/20 20:44	
2-Fluorophenol	45	10 - 105	07/01/20 20:44	
Nitrobenzene-d5	69	31 - 110	07/01/20 20:44	
Phenol-d6	31	10 - 107	07/01/20 20:44	
p-Terphenyl-d14	94	10 - 165	07/01/20 20:44	

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
1000336-66-8	6-Octadecenoic acid	10.16	11	JN	
	unknown hydrocarbon	13.13	4.6	J	
	unknown	6.36	3.7	J	
	unknown	6.51	12	J	
000057-10-3	n-Hexadecanoic acid	9.44	5.8	JN	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 10:05

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-009
 Units: ug/L

 Lab Code:
 R2005520-004
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/01/20 21:12	6/30/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/01/20 21:12	6/30/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/01/20 21:12	6/30/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/01/20 21:12	6/30/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/01/20 21:12	6/30/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/01/20 21:12	6/30/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/01/20 21:12	6/30/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/01/20 21:12	6/30/20	
2,4-Dinitrophenol	45 U	45	19	1	07/01/20 21:12	6/30/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/01/20 21:12	6/30/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/01/20 21:12	6/30/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/01/20 21:12	6/30/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/01/20 21:12	6/30/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/01/20 21:12	6/30/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/01/20 21:12	6/30/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/01/20 21:12	6/30/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/01/20 21:12	6/30/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/01/20 21:12	6/30/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/01/20 21:12	6/30/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/01/20 21:12	6/30/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/01/20 21:12	6/30/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/01/20 21:12	6/30/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/01/20 21:12	6/30/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/01/20 21:12	6/30/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/01/20 21:12	6/30/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/01/20 21:12	6/30/20	
4-Nitrophenol	45 U	45	5.8	1	07/01/20 21:12	6/30/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/01/20 21:12	6/30/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/01/20 21:12	6/30/20	
Anthracene	9.1 U	9.1	1.2	1	07/01/20 21:12	6/30/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/01/20 21:12	6/30/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/01/20 21:12	6/30/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/01/20 21:12	6/30/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/01/20 21:12	6/30/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/01/20 21:12	6/30/20	
Benzoic Acid	91 U	9.1	33	1	07/01/20 21:12	6/30/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/01/20 21:12	6/30/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/01/20 21:12	6/30/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/01/20 21:12	6/30/20	
Bis(2-chloroethyl) Ether	9.1 U 9.1 U	9.1	1.8	1	07/01/20 21:12	6/30/20	
Bis(2-ethylhexyl) Phthalate	9.1 U 9.1 U	9.1	0.91	1	07/01/20 21:12	6/30/20	
	9.1 U 9.1 U	9.1 9.1	1.3		07/01/20 21:12	6/30/20	
Butyl Benzyl Phthalate				1			
Chrysene	9.1 U	9.1	1.1	1	07/01/20 21:12	6/30/20	

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Superset Reference:20-0000554598 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 10:05

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-009
 Units: ug/L

 Lab Code:
 R2005520-004
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/01/20 21:12	6/30/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/01/20 21:12	6/30/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/01/20 21:12	6/30/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/01/20 21:12	6/30/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/01/20 21:12	6/30/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/01/20 21:12	6/30/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/01/20 21:12	6/30/20	
Fluorene	9.1 U	9.1	1.2	1	07/01/20 21:12	6/30/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/01/20 21:12	6/30/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/01/20 21:12	6/30/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/01/20 21:12	6/30/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/01/20 21:12	6/30/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/01/20 21:12	6/30/20	
Isophorone	9.1 U	9.1	1.3	1	07/01/20 21:12	6/30/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/01/20 21:12	6/30/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/01/20 21:12	6/30/20	
Naphthalene	9.1 U	9.1	1.1	1	07/01/20 21:12	6/30/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/01/20 21:12	6/30/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/01/20 21:12	6/30/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/01/20 21:12	6/30/20	
Phenol	9.1 U	9.1	0.91	1	07/01/20 21:12	6/30/20	
Pyrene	9.1 U	9.1	1.3	1	07/01/20 21:12	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	81	35 - 141	07/01/20 21:12	
2-Fluorobiphenyl	64	31 - 118	07/01/20 21:12	
2-Fluorophenol	38	10 - 105	07/01/20 21:12	
Nitrobenzene-d5	66	31 - 110	07/01/20 21:12	
Phenol-d6	28	10 - 107	07/01/20 21:12	
p-Terphenyl-d14	87	10 - 165	07/01/20 21:12	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown hydrocarbon	11.35	4.5	J	
	unknown hydrocarbon	11.89	7.7	J	
	unknown hydrocarbon	12.48	10	J	
	unknown hydrocarbon	13.13	11	J	
	unknown hydrocarbon	13.83	8.7	J	
	unknown hydrocarbon	14.58	7.4	J	
	unknown hydrocarbon	15.36	5.9	J	

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Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 10:05

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-009
 Units: ug/L

 Lab Code:
 R2005520-004
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown hydrocarbon	16.11	4.0	J	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 12:25

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-012
 Units: ug/L

 Lab Code:
 R2005520-005
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/01/20 21:40	6/30/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/01/20 21:40	6/30/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/01/20 21:40	6/30/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/01/20 21:40	6/30/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/01/20 21:40	6/30/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/01/20 21:40	6/30/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/01/20 21:40	6/30/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/01/20 21:40	6/30/20	
2,4-Dinitrophenol	45 U	45	19	1	07/01/20 21:40	6/30/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/01/20 21:40	6/30/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/01/20 21:40	6/30/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/01/20 21:40	6/30/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/01/20 21:40	6/30/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/01/20 21:40	6/30/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/01/20 21:40	6/30/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/01/20 21:40	6/30/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/01/20 21:40	6/30/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/01/20 21:40	6/30/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/01/20 21:40	6/30/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/01/20 21:40	6/30/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/01/20 21:40	6/30/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/01/20 21:40	6/30/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/01/20 21:40	6/30/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/01/20 21:40	6/30/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/01/20 21:40	6/30/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/01/20 21:40	6/30/20	
4-Nitrophenol	45 U	45	5.8	1	07/01/20 21:40	6/30/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/01/20 21:40	6/30/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/01/20 21:40	6/30/20	
Anthracene	9.1 U	9.1	1.2	1	07/01/20 21:40	6/30/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/01/20 21:40	6/30/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/01/20 21:40	6/30/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/01/20 21:40	6/30/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/01/20 21:40	6/30/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/01/20 21:40	6/30/20	
Benzoic Acid	91 U	91	33	1	07/01/20 21:40	6/30/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/01/20 21:40	6/30/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/01/20 21:40	6/30/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/01/20 21:40	6/30/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	07/01/20 21:40	6/30/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	07/01/20 21:40	6/30/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	07/01/20 21:40	6/30/20	
Chrysene	9.1 U	9.1	1.1	1	07/01/20 21:40	6/30/20	

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Superset Reference:20-0000554598 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 12:25

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-012
 Units: ug/L

 Lab Code:
 R2005520-005
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/01/20 21:40	6/30/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/01/20 21:40	6/30/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/01/20 21:40	6/30/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/01/20 21:40	6/30/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/01/20 21:40	6/30/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/01/20 21:40	6/30/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/01/20 21:40	6/30/20	
Fluorene	9.1 U	9.1	1.2	1	07/01/20 21:40	6/30/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/01/20 21:40	6/30/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/01/20 21:40	6/30/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/01/20 21:40	6/30/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/01/20 21:40	6/30/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/01/20 21:40	6/30/20	
Isophorone	9.1 U	9.1	1.3	1	07/01/20 21:40	6/30/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/01/20 21:40	6/30/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/01/20 21:40	6/30/20	
Naphthalene	9.1 U	9.1	1.1	1	07/01/20 21:40	6/30/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/01/20 21:40	6/30/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/01/20 21:40	6/30/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/01/20 21:40	6/30/20	
Phenol	9.1 U	9.1	0.91	1	07/01/20 21:40	6/30/20	
Pyrene	9.1 U	9.1	1.3	1	07/01/20 21:40	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	83	35 - 141	07/01/20 21:40	
2-Fluorobiphenyl	59	31 - 118	07/01/20 21:40	
2-Fluorophenol	40	10 - 105	07/01/20 21:40	
Nitrobenzene-d5	59	31 - 110	07/01/20 21:40	
Phenol-d6	28	10 - 107	07/01/20 21:40	
p-Terphenyl-d14	93	10 - 165	07/01/20 21:40	

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
013798-23-7	Sulfur	7.79	9.6	JN	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 13:50

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 RB-9954-062520-SG-001
 Units: ug/L

 Lab Code:
 R2005520-006
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/01/20 23:03	6/30/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/01/20 23:03	6/30/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/01/20 23:03	6/30/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/01/20 23:03	6/30/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/01/20 23:03	6/30/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/01/20 23:03	6/30/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/01/20 23:03	6/30/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/01/20 23:03	6/30/20	
2,4-Dinitrophenol	45 U	45	19	1	07/01/20 23:03	6/30/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/01/20 23:03	6/30/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/01/20 23:03	6/30/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/01/20 23:03	6/30/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/01/20 23:03	6/30/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/01/20 23:03	6/30/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/01/20 23:03	6/30/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/01/20 23:03	6/30/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/01/20 23:03	6/30/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/01/20 23:03	6/30/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/01/20 23:03	6/30/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/01/20 23:03	6/30/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/01/20 23:03	6/30/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/01/20 23:03	6/30/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/01/20 23:03	6/30/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/01/20 23:03	6/30/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/01/20 23:03	6/30/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/01/20 23:03	6/30/20	
4-Nitrophenol	45 U	45	5.8	1	07/01/20 23:03	6/30/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/01/20 23:03	6/30/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/01/20 23:03	6/30/20	
Anthracene	9.1 U	9.1	1.2	1	07/01/20 23:03	6/30/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/01/20 23:03	6/30/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/01/20 23:03	6/30/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/01/20 23:03	6/30/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/01/20 23:03	6/30/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/01/20 23:03	6/30/20	
Benzoic Acid	91 U	91	33	1	07/01/20 23:03	6/30/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/01/20 23:03	6/30/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/01/20 23:03	6/30/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/01/20 23:03	6/30/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	07/01/20 23:03	6/30/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	07/01/20 23:03	6/30/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	07/01/20 23:03	6/30/20	
Chrysene	9.1 U	9.1	1.1	1	07/01/20 23:03	6/30/20	
2111 3 50110).i U	/.1	1.1	1	57,01,20 25.05	0/20/20	

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Superset Reference: 20-0000554598 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 13:50

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 RB-9954-062520-SG-001
 Units: ug/L

 Lab Code:
 R2005520-006
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/01/20 23:03	6/30/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/01/20 23:03	6/30/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/01/20 23:03	6/30/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/01/20 23:03	6/30/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/01/20 23:03	6/30/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/01/20 23:03	6/30/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/01/20 23:03	6/30/20	
Fluorene	9.1 U	9.1	1.2	1	07/01/20 23:03	6/30/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/01/20 23:03	6/30/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/01/20 23:03	6/30/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/01/20 23:03	6/30/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/01/20 23:03	6/30/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/01/20 23:03	6/30/20	
Isophorone	9.1 U	9.1	1.3	1	07/01/20 23:03	6/30/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/01/20 23:03	6/30/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/01/20 23:03	6/30/20	
Naphthalene	9.1 U	9.1	1.1	1	07/01/20 23:03	6/30/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/01/20 23:03	6/30/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/01/20 23:03	6/30/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/01/20 23:03	6/30/20	
Phenol	9.1 U	9.1	0.91	1	07/01/20 23:03	6/30/20	
Pyrene	9.1 U	9.1	1.3	1	07/01/20 23:03	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	89	35 - 141	07/01/20 23:03	
2-Fluorobiphenyl	78	31 - 118	07/01/20 23:03	
2-Fluorophenol	39	10 - 105	07/01/20 23:03	
Nitrobenzene-d5	68	31 - 110	07/01/20 23:03	
Phenol-d6	28	10 - 107	07/01/20 23:03	
p-Terphenyl-d14	94	10 - 165	07/01/20 23:03	

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown hydrocarbon	11.89	4.0	J	
	unknown hydrocarbon	12.48	4.9	J	
	unknown hydrocarbon	13.13	4.9	J	
	unknown hydrocarbon	13.83	4.1	J	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 10:05

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-010
 Units: ug/L

 Lab Code:
 R2005520-008
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/01/20 23:32	6/30/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/01/20 23:32	6/30/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/01/20 23:32	6/30/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/01/20 23:32	6/30/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/01/20 23:32	6/30/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/01/20 23:32	6/30/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/01/20 23:32	6/30/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/01/20 23:32	6/30/20	
2,4-Dinitrophenol	45 U	45	19	1	07/01/20 23:32	6/30/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/01/20 23:32	6/30/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/01/20 23:32	6/30/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/01/20 23:32	6/30/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/01/20 23:32	6/30/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/01/20 23:32	6/30/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/01/20 23:32	6/30/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/01/20 23:32	6/30/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/01/20 23:32	6/30/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/01/20 23:32	6/30/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/01/20 23:32	6/30/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/01/20 23:32	6/30/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/01/20 23:32	6/30/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/01/20 23:32	6/30/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/01/20 23:32	6/30/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/01/20 23:32	6/30/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/01/20 23:32	6/30/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/01/20 23:32	6/30/20	
4-Nitrophenol	45 U	45	5.8	1	07/01/20 23:32	6/30/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/01/20 23:32	6/30/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/01/20 23:32	6/30/20	
Anthracene	9.1 U	9.1	1.2	1	07/01/20 23:32	6/30/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/01/20 23:32	6/30/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/01/20 23:32	6/30/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/01/20 23:32	6/30/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/01/20 23:32	6/30/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/01/20 23:32	6/30/20	
Benzoic Acid	91 U	91	33	1	07/01/20 23:32	6/30/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/01/20 23:32	6/30/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/01/20 23:32	6/30/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/01/20 23:32	6/30/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	07/01/20 23:32	6/30/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	07/01/20 23:32	6/30/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	07/01/20 23:32	6/30/20	
Chrysene	9.1 U	9.1	1.1	1	07/01/20 23:32	6/30/20	

Printed 7/16/2020 11:08:29 AM

Superset Reference: 20-0000554598 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 10:05

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-010
 Units: ug/L

 Lab Code:
 R2005520-008
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed Date Extra	acted Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/01/20 23:32 6/30/20	0
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/01/20 23:32 6/30/20	C
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/01/20 23:32 6/30/20)
Dibenzofuran	9.1 U	9.1	1.3	1	07/01/20 23:32 6/30/20)
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/01/20 23:32 6/30/20)
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/01/20 23:32 6/30/20)
Fluoranthene	9.1 U	9.1	1.4	1	07/01/20 23:32 6/30/20	0
Fluorene	9.1 U	9.1	1.2	1	07/01/20 23:32 6/30/20)
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/01/20 23:32 6/30/20)
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/01/20 23:32 6/30/20)
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/01/20 23:32 6/30/20)
Hexachloroethane	9.1 U	9.1	0.96	1	07/01/20 23:32 6/30/20)
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/01/20 23:32 6/30/20)
Isophorone	9.1 U	9.1	1.3	1	07/01/20 23:32 6/30/20)
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/01/20 23:32 6/30/20)
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/01/20 23:32 6/30/20)
Naphthalene	9.1 U	9.1	1.1	1	07/01/20 23:32 6/30/20)
Nitrobenzene	9.1 U	9.1	1.4	1	07/01/20 23:32 6/30/20)
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/01/20 23:32 6/30/20)
Phenanthrene	9.1 U	9.1	1.3	1	07/01/20 23:32 6/30/20)
Phenol	9.1 U	9.1	0.91	1	07/01/20 23:32 6/30/20)
Pyrene	9.1 U	9.1	1.3	1	07/01/20 23:32 6/30/20)

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	79	35 - 141	07/01/20 23:32	
2-Fluorobiphenyl	67	31 - 118	07/01/20 23:32	
2-Fluorophenol	44	10 - 105	07/01/20 23:32	
Nitrobenzene-d5	68	31 - 110	07/01/20 23:32	
Phenol-d6	29	10 - 107	07/01/20 23:32	
p-Terphenyl-d14	92	10 - 165	07/01/20 23:32	

Tentatively Identified Compounds

			Result	
CAS#	Compound Identification	RT	ug/L	Q

No Tentatively Identified Compounds

Detected



Semivolatile Organic Compounds by GC

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 13:15

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-013
 Units: ug/L

 Lab Code:
 R2005520-001
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/08/20 09:54	6/30/20	
4,4'-DDE	0.045 U	0.045	0.019	1	07/08/20 09:54	6/30/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/08/20 09:54	6/30/20	
Aldrin	0.019 J	0.045	0.019	1	07/08/20 09:54	6/30/20	
Dieldrin	0.045 U	0.045	0.019	1	07/08/20 09:54	6/30/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/08/20 09:54	6/30/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/08/20 09:54	6/30/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/08/20 09:54	6/30/20	
Endrin	0.045 U	0.045	0.019	1	07/08/20 09:54	6/30/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/08/20 09:54	6/30/20	
Heptachlor	0.045 U	0.045	0.019	1	07/08/20 09:54	6/30/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/08/20 09:54	6/30/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/08/20 09:54	6/30/20	
Toxaphene	0.46 U	0.46	0.46	1	07/08/20 09:54	6/30/20	
alpha-BHC	0.18	0.045	0.019	1	07/08/20 09:54	6/30/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/08/20 09:54	6/30/20	
beta-BHC	0.045 U	0.045	0.019	1	07/08/20 09:54	6/30/20	
delta-BHC	0.28	0.045	0.019	1	07/08/20 09:54	6/30/20	
gamma-BHC (Lindane)	0.21	0.045	0.019	1	07/08/20 09:54	6/30/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/08/20 09:54	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	67	10 - 164	07/08/20 09:54	
Tetrachloro-m-xylene	80	10 - 147	07/08/20 09:54	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 11:10

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-011
 Units: ug/L

 Lab Code:
 R2005520-002
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/08/20 10:13	6/30/20	
4,4'-DDE	0.045 U	0.045	0.019	1	07/08/20 10:13	6/30/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/08/20 10:13	6/30/20	
Aldrin	0.045 U	0.045	0.019	1	07/08/20 10:13	6/30/20	
Dieldrin	0.045 U	0.045	0.019	1	07/08/20 10:13	6/30/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/08/20 10:13	6/30/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/08/20 10:13	6/30/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/08/20 10:13	6/30/20	
Endrin	0.045 U	0.045	0.019	1	07/08/20 10:13	6/30/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/08/20 10:13	6/30/20	
Heptachlor	0.045 U	0.045	0.019	1	07/08/20 10:13	6/30/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/08/20 10:13	6/30/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/08/20 10:13	6/30/20	
Toxaphene	0.46 U	0.46	0.46	1	07/08/20 10:13	6/30/20	
alpha-BHC	0.31	0.045	0.019	1	07/08/20 10:13	6/30/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/08/20 10:13	6/30/20	
beta-BHC	0.045 U	0.045	0.019	1	07/08/20 10:13	6/30/20	
delta-BHC	0.13	0.045	0.019	1	07/08/20 10:13	6/30/20	
gamma-BHC (Lindane)	0.25	0.045	0.019	1	07/08/20 10:13	6/30/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/08/20 10:13	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	56	10 - 164	07/08/20 10:13	
Tetrachloro-m-xylene	71	10 - 147	07/08/20 10:13	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 09:15

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-008
 Units: ug/L

 Lab Code:
 R2005520-003
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/08/20 10:32	6/30/20	_
4,4'-DDE	0.045 U	0.045	0.019	1	07/08/20 10:32	6/30/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/08/20 10:32	6/30/20	
Aldrin	0.045 U	0.045	0.019	1	07/08/20 10:32	6/30/20	
Dieldrin	0.045 U	0.045	0.019	1	07/08/20 10:32	6/30/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/08/20 10:32	6/30/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/08/20 10:32	6/30/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/08/20 10:32	6/30/20	
Endrin	0.045 U	0.045	0.019	1	07/08/20 10:32	6/30/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/08/20 10:32	6/30/20	
Heptachlor	0.045 U	0.045	0.019	1	07/08/20 10:32	6/30/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/08/20 10:32	6/30/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/08/20 10:32	6/30/20	
Toxaphene	0.46 U	0.46	0.46	1	07/08/20 10:32	6/30/20	
alpha-BHC	0.045 U	0.045	0.019	1	07/08/20 10:32	6/30/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/08/20 10:32	6/30/20	
beta-BHC	0.045 U	0.045	0.019	1	07/08/20 10:32	6/30/20	
delta-BHC	0.045 U	0.045	0.019	1	07/08/20 10:32	6/30/20	
gamma-BHC (Lindane)	0.045 U	0.045	0.019	1	07/08/20 10:32	6/30/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/08/20 10:32	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	65	10 - 164	07/08/20 10:32	
Tetrachloro-m-xylene	74	10 - 147	07/08/20 10:32	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 10:05

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-009
 Units: ug/L

 Lab Code:
 R2005520-004
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/08/20 10:51	6/30/20	
4,4'-DDE	0.045 U	0.045	0.019	1	07/08/20 10:51	6/30/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/08/20 10:51	6/30/20	
Aldrin	0.045 U	0.045	0.019	1	07/08/20 10:51	6/30/20	
Dieldrin	0.045 U	0.045	0.019	1	07/08/20 10:51	6/30/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/08/20 10:51	6/30/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/08/20 10:51	6/30/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/08/20 10:51	6/30/20	
Endrin	0.045 U	0.045	0.019	1	07/08/20 10:51	6/30/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/08/20 10:51	6/30/20	
Heptachlor	0.045 U	0.045	0.019	1	07/08/20 10:51	6/30/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/08/20 10:51	6/30/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/08/20 10:51	6/30/20	
Toxaphene	0.46 U	0.46	0.46	1	07/08/20 10:51	6/30/20	
alpha-BHC	0.045 U	0.045	0.019	1	07/08/20 10:51	6/30/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/08/20 10:51	6/30/20	
beta-BHC	0.045 U	0.045	0.019	1	07/08/20 10:51	6/30/20	
delta-BHC	0.045 U	0.045	0.019	1	07/08/20 10:51	6/30/20	
gamma-BHC (Lindane)	0.045 U	0.045	0.019	1	07/08/20 10:51	6/30/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/08/20 10:51	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	37	10 - 164	07/08/20 10:51	
Tetrachloro-m-xylene	57	10 - 147	07/08/20 10:51	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 12:25

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-012
 Units: ug/L

 Lab Code:
 R2005520-005
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/08/20 18:43	6/30/20	
4,4'-DDE	0.045 U	0.045	0.019	1	07/08/20 18:43	6/30/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/08/20 18:43	6/30/20	
Aldrin	0.045 U	0.045	0.019	1	07/08/20 18:43	6/30/20	
Dieldrin	0.045 U	0.045	0.019	1	07/08/20 18:43	6/30/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/08/20 18:43	6/30/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/08/20 18:43	6/30/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/08/20 18:43	6/30/20	
Endrin	0.045 U	0.045	0.019	1	07/08/20 18:43	6/30/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/08/20 18:43	6/30/20	
Heptachlor	0.045 U	0.045	0.019	1	07/08/20 18:43	6/30/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/08/20 18:43	6/30/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/08/20 18:43	6/30/20	
Toxaphene	0.46 U	0.46	0.46	1	07/08/20 18:43	6/30/20	
alpha-BHC	0.033 J	0.045	0.019	1	07/08/20 18:43	6/30/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/08/20 18:43	6/30/20	
beta-BHC	0.045 U	0.045	0.019	1	07/08/20 18:43	6/30/20	
delta-BHC	0.068	0.045	0.019	1	07/08/20 18:43	6/30/20	
gamma-BHC (Lindane)	0.037 J	0.045	0.019	1	07/08/20 18:43	6/30/20	
gamma-Chlordane	0.028 JP	0.045	0.019	1	07/08/20 18:43	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	49	10 - 164	07/08/20 18:43	
Tetrachloro-m-xylene	85	10 - 147	07/08/20 18:43	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 13:50

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 RB-9954-062520-SG-001
 Units: ug/L

 Lab Code:
 R2005520-006
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/06/20 11:38	7/1/20	
4,4'-DDE	0.045 U	0.045	0.019	1	07/06/20 11:38	7/1/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/06/20 11:38	7/1/20	
Aldrin	0.045 U	0.045	0.019	1	07/06/20 11:38	7/1/20	
Dieldrin	0.045 U	0.045	0.019	1	07/06/20 11:38	7/1/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/06/20 11:38	7/1/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/06/20 11:38	7/1/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/06/20 11:38	7/1/20	
Endrin	0.045 U	0.045	0.019	1	07/06/20 11:38	7/1/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/06/20 11:38	7/1/20	
Heptachlor	0.045 U	0.045	0.019	1	07/06/20 11:38	7/1/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/06/20 11:38	7/1/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/06/20 11:38	7/1/20	
Toxaphene	0.46 U	0.46	0.46	1	07/06/20 11:38	7/1/20	
alpha-BHC	0.079	0.045	0.019	1	07/06/20 11:38	7/1/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/06/20 11:38	7/1/20	
beta-BHC	0.045 U	0.045	0.019	1	07/06/20 11:38	7/1/20	
delta-BHC	0.044 J	0.045	0.019	1	07/06/20 11:38	7/1/20	
gamma-BHC (Lindane)	0.075	0.045	0.019	1	07/06/20 11:38	7/1/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/06/20 11:38	7/1/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	10	10 - 164	07/06/20 11:38	
Tetrachloro-m-xylene	53	10 - 147	07/06/20 11:38	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 10:05

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-010
 Units: ug/L

 Lab Code:
 R2005520-008
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/06/20 11:57	7/1/20	_
4,4'-DDE	0.045 U	0.045	0.019	1	07/06/20 11:57	7/1/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/06/20 11:57	7/1/20	
Aldrin	0.045 U	0.045	0.019	1	07/06/20 11:57	7/1/20	
Dieldrin	0.045 U	0.045	0.019	1	07/06/20 11:57	7/1/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/06/20 11:57	7/1/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/06/20 11:57	7/1/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/06/20 11:57	7/1/20	
Endrin	0.045 U	0.045	0.019	1	07/06/20 11:57	7/1/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/06/20 11:57	7/1/20	
Heptachlor	0.045 U	0.045	0.019	1	07/06/20 11:57	7/1/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/06/20 11:57	7/1/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/06/20 11:57	7/1/20	
Toxaphene	0.46 U	0.46	0.46	1	07/06/20 11:57	7/1/20	
alpha-BHC	0.045 U	0.045	0.019	1	07/06/20 11:57	7/1/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/06/20 11:57	7/1/20	
beta-BHC	0.045 U	0.045	0.019	1	07/06/20 11:57	7/1/20	
delta-BHC	0.045 U	0.045	0.019	1	07/06/20 11:57	7/1/20	
gamma-BHC (Lindane)	0.045 U	0.045	0.019	1	07/06/20 11:57	7/1/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/06/20 11:57	7/1/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	27	10 - 164	07/06/20 11:57	
Tetrachloro-m-xylene	45	10 - 147	07/06/20 11:57	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 13:15

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-013
 Units: ug/L

 Lab Code:
 R2005520-001
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/07/20 18:23	6/30/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/07/20 18:23	6/30/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/07/20 18:23	6/30/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/07/20 18:23	6/30/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/07/20 18:23	6/30/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/07/20 18:23	6/30/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/07/20 18:23	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	53	10 - 152	07/07/20 18:23	
Tetrachloro-m-xvlene	47	14 - 129	07/07/20 18:23	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 11:10

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-011
 Units: ug/L

 Lab Code:
 R2005520-002
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/07/20 18:42	6/30/20	_
Aroclor 1221	1.8 U	1.8	0.91	1	07/07/20 18:42	6/30/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/07/20 18:42	6/30/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/07/20 18:42	6/30/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/07/20 18:42	6/30/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/07/20 18:42	6/30/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/07/20 18:42	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	41	10 - 152	07/07/20 18:42	
Tetrachloro-m-xylene	45	14 - 129	07/07/20 18:42	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 09:15

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-008
 Units: ug/L

 Lab Code:
 R2005520-003
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/01/20 23:12	6/30/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/01/20 23:12	6/30/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/01/20 23:12	6/30/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/01/20 23:12	6/30/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/01/20 23:12	6/30/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/01/20 23:12	6/30/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/01/20 23:12	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	54	10 - 152	07/01/20 23:12	
Tetrachloro-m-xylene	57	14 - 129	07/01/20 23:12	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 10:05

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-009
 Units: ug/L

 Lab Code:
 R2005520-004
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/01/20 23:32	6/30/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/01/20 23:32	6/30/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/01/20 23:32	6/30/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/01/20 23:32	6/30/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/01/20 23:32	6/30/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/01/20 23:32	6/30/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/01/20 23:32	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	30	10 - 152	07/01/20 23:32	
Tetrachloro-m-xvlene	59	14 - 129	07/01/20 23:32	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 12:25

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-012
 Units: ug/L

 Lab Code:
 R2005520-005
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/07/20 19:02	6/30/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/07/20 19:02	6/30/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/07/20 19:02	6/30/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/07/20 19:02	6/30/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/07/20 19:02	6/30/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/07/20 19:02	6/30/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/07/20 19:02	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q	
Decachlorobiphenyl	60	10 - 152	07/07/20 19:02		
Tetrachloro-m-xvlene	46	14 - 129	07/07/20 19:02		

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 13:50

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 RB-9954-062520-SG-001
 Units: ug/L

 Lab Code:
 R2005520-006
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/06/20 12:41	7/1/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/06/20 12:41	7/1/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/06/20 12:41	7/1/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/06/20 12:41	7/1/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/06/20 12:41	7/1/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/06/20 12:41	7/1/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/06/20 12:41	7/1/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	10	10 - 152	07/06/20 12:41	
Tetrachloro-m-xylene	47	14 - 129	07/06/20 12:41	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/25/20 10:05

Monitoring

Sample Matrix: Water Date Received: 06/26/20 10:25

 Sample Name:
 WG-9954-062520-SG-010
 Units: ug/L

 Lab Code:
 R2005520-008
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/06/20 13:42	7/1/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/06/20 13:42	7/1/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/06/20 13:42	7/1/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/06/20 13:42	7/1/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/06/20 13:42	7/1/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/06/20 13:42	7/1/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/06/20 13:42	7/1/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	28	10 - 152	07/06/20 13:42	
Tetrachloro-m-xylene	39	14 - 129	07/06/20 13:42	



QC Summary Forms

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com



Volatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

QA/QC Report

Service Request: R2005520

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

Project:

SURROGATE RECOVERY SUMMARY Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Extraction Method: EPA 5030C

		4-Bromofluorobenzene	Dibromofluoromethane	Toluene-d8
Sample Name	Lab Code	85-122	89-119	87-121
WG-9954-062520-SG-013	R2005520-001	92	97	97
WG-9954-062520-SG-011	R2005520-002	92	100	98
WG-9954-062520-SG-008	R2005520-003	90	92	93
WG-9954-062520-SG-009	R2005520-004	91	96	96
WG-9954-062520-SG-012	R2005520-005	93	98	98
RB-9954-062520-SG-001	R2005520-006	90	89	92
TB-9954-062520-SG-002	R2005520-007	91	92	95
WG-9954-062520-SG-010	R2005520-008	107	109	107
Method Blank	RQ2007077-06	92	95	97
Method Blank	RQ2007141-04	90	89	91
Method Blank	RQ2007251-04	101	102	100
Lab Control Sample	RQ2007077-04	94	98	97
Lab Control Sample	RQ2007141-03	95	95	93
Lab Control Sample	RQ2007251-03	102	103	101
WG-9954-062520-SG-012 MS	RQ2007077-07	92	98	97
WG-9954-062520-SG-012 DMS	RQ2007077-08	97	102	100

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Date Collected: 06/25/20 **Date Received:** 06/26/20

Sample Matrix: Water

Date Analyzed: 07/2/20

NA

Date Extracted:

Duplicate Matrix Spike Summary Volatile Organic Compounds by GC/MS

Sample Name: WG-9954-062520-SG-012

Units: ug/L

Lab Code: R2005520-005

Basis: NA

Analysis Method: 8260C **Prep Method:** EPA 5030C

Matrix SpikeDuplicate Matrix SpikeRQ2007077-07RQ2007077-08

4 1 4 27	Sample	D 1/	Spike	0/ D	D 14	Spike	0 / TD	% Rec	DDD	RPD
Analyte Name	Result	Result	Amount	% Rec	Result	Amount	% Rec	Limits	RPD	Limit
1,1,1-Trichloroethane (TCA)	5.0 U	57.7	50.0	115	63.7	50.0	127	74-127	10	30
1,1,2,2-Tetrachloroethane	5.0 U	64.9	50.0	130 *	70.7	50.0	141 *	72-122	9	30
1,1,2-Trichloroethane	5.0 U	51.8	50.0	104	57.8	50.0	116	82-121	11	30
1,1-Dichloroethane (1,1-DCA)	5.0 U	58.8	50.0	118	63.7	50.0	127	74-132	8	30
1,1-Dichloroethene (1,1-DCE)	5.0 U	55.1	50.0	110	60.6	50.0	121 *	71-118	9	30
1,2-Dichloroethane	5.0 U	48.5	50.0	97	53.8	50.0	108	68-130	10	30
1,2-Dichloropropane	5.0 U	56.1	50.0	112	61.8	50.0	124	79-124	10	30
2-Butanone (MEK)	10 U	50.5	50.0	101	58.0	50.0	116	61-137	14	30
2-Hexanone	10 U	53.8	50.0	108	61.7	50.0	123	56-132	14	30
4-Methyl-2-pentanone	10 U	54.2	50.0	108	62.8	50.0	126	60-141	15	30
Acetone	10 U	47.5	50.0	95	53.5	50.0	107	35-183	12	30
Benzene	5.0 U	54.1	50.0	108	59.3	50.0	119	76-129	9	30
Bromodichloromethane	5.0 U	55.3	50.0	111	61.2	50.0	122	78-133	10	30
Bromoform	5.0 U	51.1	50.0	102	56.1	50.0	112	58-133	9	30
Bromomethane	5.0 U	23.8	50.0	48	24.8	50.0	50	10-184	4	30
Carbon Disulfide	2.6 J	59.6	50.0	114	68.9	50.0	133	59-140	14	30
Carbon Tetrachloride	5.0 U	55.5	50.0	111	61.1	50.0	122	65-135	10	30
Chlorobenzene	5.0 U	51.1	50.0	102	55.9	50.0	112	76-125	9	30
Chloroethane	5.0 U	59.8	50.0	120	67.1	50.0	134	48-146	12	30
Chloroform	5.0 U	55.4	50.0	111	59.9	50.0	120	75-130	8	30
Chloromethane	0.30 J	51.9	50.0	103	56.4	50.0	112	55-160	8	30
Dibromochloromethane	5.0 U	53.0	50.0	106	58.0	50.0	116	72-128	9	30
Dichloromethane	5.0 U	53.3	50.0	107	56.9	50.0	114	73-122	7	30
Ethylbenzene	5.0 U	55.1	50.0	110	59.8	50.0	120	72-134	8	30
Styrene	5.0 U	54.3	50.0	109	58.9	50.0	118	74-136	8	30
Tetrachloroethene (PCE)	5.0 U	50.9	50.0	102	54.8	50.0	110	72-125	7	30
Toluene	5.0 U	54.2	50.0	108	59.5	50.0	119	79-119	9	30
Trichloroethene (TCE)	5.0 U	42.0	50.0	84	47.1	50.0	94	74-122	11	30
Vinyl Acetate	10 U	62.8	50.0	126	70.6	50.0	141	48-172	12	30
Vinyl Chloride	5.0 U	61.8	50.0	124	66.7	50.0	133	74-159	8	30
cis-1,2-Dichloroethene	5.0 U	56.3	50.0	113	62.1	50.0	124	77-127	10	30
cis-1,3-Dichloropropene	5.0 U	52.8	50.0	106	58.7	50.0	117	52-134	11	30
trans-1,2-Dichloroethene	5.0 U	55.6	50.0	111	61.3	50.0	123 *	73-118	10	30

Results flagged with an asterisk (*) indicate values outside control criteria.

Results flagged with a pound (#) indicate the control criteria is not applicable.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

Matrix Spike and Matrix Spike Duplicate Data is presented for information purposes only. The matrix may or may not be relevant to samples reported in this report. The laboratory evaluates system performance based on the LCS and LCSD control limits.

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates) **Service Request:**

R2005520

Project:

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Date Collected:

06/25/20

Sample Matrix:

Water

Date Received:

06/26/20

Date Analyzed:

07/2/20

Date Extracted:

NA

Duplicate Matrix Spike Summary Volatile Organic Compounds by GC/MS

Sample Name:

WG-9954-062520-SG-012

Units:

ug/L

Lab Code:

Prep Method:

R2005520-005

Basis:

NA

Analysis Method:

8260C

EPA 5030C

Matrix Spike

Duplicate Matrix Spike

RQ2007077-08

RQ2007077-07

RPD Sample **Spike** Spike % Rec Analyte Name Result Amount **RPD** Result **Amount** % Rec Result % Rec Limits Limit 52.2 trans-1,3-Dichloropropene 5.0 U 50.0 104 58.6 50.0 30

Results flagged with an asterisk (*) indicate values outside control criteria.

Results flagged with a pound (#) indicate the control criteria is not applicable.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

Matrix Spike and Matrix Spike Duplicate Data is presented for information purposes only. The matrix may or may not be relevant to samples reported in this report. The laboratory evaluates system performance based on the LCS and LCSD control limits.

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007077-06
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/02/20 13:35	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/02/20 13:35	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/02/20 13:35	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/02/20 13:35	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/02/20 13:35	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/02/20 13:35	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/02/20 13:35	
2-Butanone (MEK)	10 U	10	0.78	1	07/02/20 13:35	
2-Hexanone	10 U	10	0.20	1	07/02/20 13:35	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/02/20 13:35	
Acetone	10 U	10	5.0	1	07/02/20 13:35	
Benzene	5.0 U	5.0	0.20	1	07/02/20 13:35	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/02/20 13:35	
Bromoform	5.0 U	5.0	0.25	1	07/02/20 13:35	
Bromomethane	5.0 U	5.0	0.70	1	07/02/20 13:35	
Carbon Disulfide	10 U	10	0.42	1	07/02/20 13:35	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/02/20 13:35	
Chlorobenzene	5.0 U	5.0	0.20	1	07/02/20 13:35	
Chloroethane	5.0 U	5.0	0.23	1	07/02/20 13:35	
Chloroform	5.0 U	5.0	0.24	1	07/02/20 13:35	
Chloromethane	5.0 U	5.0	0.28	1	07/02/20 13:35	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/02/20 13:35	
Dichloromethane	5.0 U	5.0	0.65	1	07/02/20 13:35	
Ethylbenzene	5.0 U	5.0	0.20	1	07/02/20 13:35	
Styrene	5.0 U	5.0	0.20	1	07/02/20 13:35	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/02/20 13:35	
Toluene	5.0 U	5.0	0.20	1	07/02/20 13:35	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/02/20 13:35	
Vinyl Acetate	10 U	10	1.1	1	07/02/20 13:35	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/02/20 13:35	
Xylenes, Total	5.0 U	5.0	0.23	1	07/02/20 13:35	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/02/20 13:35	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/02/20 13:35	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/02/20 13:35	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/02/20 13:35	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007077-06Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	92	85 - 122	07/02/20 13:35	
Dibromofluoromethane	95	89 - 119	07/02/20 13:35	
Toluene-d8	97	87 - 121	07/02/20 13:35	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Love Canal:292-402-D02-3100/9954 Annual Long Term **Project: Date Collected:** NA

Monitoring

Date Received: NA **Sample Matrix:** Water

Sample Name: Method Blank Units: ug/L Lab Code:

RQ2007077-06 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

Result CAS# **Compound Identification** RTug/L

No Tentatively Identified Compounds

Detected

Q

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007141-04
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	06/30/20 23:08	_
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	06/30/20 23:08	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	06/30/20 23:08	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	06/30/20 23:08	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	06/30/20 23:08	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	06/30/20 23:08	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	06/30/20 23:08	
2-Butanone (MEK)	10 U	10	0.78	1	06/30/20 23:08	
2-Hexanone	10 U	10	0.20	1	06/30/20 23:08	
4-Methyl-2-pentanone	10 U	10	0.20	1	06/30/20 23:08	
Acetone	10 U	10	5.0	1	06/30/20 23:08	
Benzene	5.0 U	5.0	0.20	1	06/30/20 23:08	
Bromodichloromethane	5.0 U	5.0	0.20	1	06/30/20 23:08	
Bromoform	5.0 U	5.0	0.25	1	06/30/20 23:08	
Bromomethane	5.0 U	5.0	0.70	1	06/30/20 23:08	
Carbon Disulfide	10 U	10	0.42	1	06/30/20 23:08	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	06/30/20 23:08	
Chlorobenzene	5.0 U	5.0	0.20	1	06/30/20 23:08	
Chloroethane	5.0 U	5.0	0.23	1	06/30/20 23:08	
Chloroform	5.0 U	5.0	0.24	1	06/30/20 23:08	
Chloromethane	5.0 U	5.0	0.28	1	06/30/20 23:08	
Dibromochloromethane	5.0 U	5.0	0.20	1	06/30/20 23:08	
Dichloromethane	5.0 U	5.0	0.65	1	06/30/20 23:08	
Ethylbenzene	5.0 U	5.0	0.20	1	06/30/20 23:08	
Styrene	5.0 U	5.0	0.20	1	06/30/20 23:08	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	06/30/20 23:08	
Toluene	5.0 U	5.0	0.20	1	06/30/20 23:08	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	06/30/20 23:08	
Vinyl Acetate	10 U	10	1.1	1	06/30/20 23:08	
Vinyl Chloride	5.0 U	5.0	0.20	1	06/30/20 23:08	
Xylenes, Total	5.0 U	5.0	0.23	1	06/30/20 23:08	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	06/30/20 23:08	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	06/30/20 23:08	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	06/30/20 23:08	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	06/30/20 23:08	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007141-04Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	90	85 - 122	06/30/20 23:08	
Dibromofluoromethane	89	89 - 119	06/30/20 23:08	
Toluene-d8	91	87 - 121	06/30/20 23:08	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007141-04Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

Result

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007251-04Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/07/20 17:23	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/07/20 17:23	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/07/20 17:23	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/07/20 17:23	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/07/20 17:23	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/07/20 17:23	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/07/20 17:23	
2-Butanone (MEK)	10 U	10	0.78	1	07/07/20 17:23	
2-Hexanone	10 U	10	0.20	1	07/07/20 17:23	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/07/20 17:23	
Acetone	10 U	10	5.0	1	07/07/20 17:23	
Benzene	5.0 U	5.0	0.20	1	07/07/20 17:23	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/07/20 17:23	
Bromoform	5.0 U	5.0	0.25	1	07/07/20 17:23	
Bromomethane	5.0 U	5.0	0.70	1	07/07/20 17:23	
Carbon Disulfide	10 U	10	0.42	1	07/07/20 17:23	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/07/20 17:23	
Chlorobenzene	5.0 U	5.0	0.20	1	07/07/20 17:23	
Chloroethane	5.0 U	5.0	0.23	1	07/07/20 17:23	
Chloroform	5.0 U	5.0	0.24	1	07/07/20 17:23	
Chloromethane	5.0 U	5.0	0.28	1	07/07/20 17:23	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/07/20 17:23	
Dichloromethane	5.0 U	5.0	0.65	1	07/07/20 17:23	
Ethylbenzene	5.0 U	5.0	0.20	1	07/07/20 17:23	
Styrene	5.0 U	5.0	0.20	1	07/07/20 17:23	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/07/20 17:23	
Toluene	5.0 U	5.0	0.20	1	07/07/20 17:23	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/07/20 17:23	
Vinyl Acetate	10 U	10	1.1	1	07/07/20 17:23	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/07/20 17:23	
Xylenes, Total	5.0 U	5.0	0.23	1	07/07/20 17:23	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/07/20 17:23	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/07/20 17:23	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/07/20 17:23	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/07/20 17:23	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007251-04Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	101	85 - 122	07/07/20 17:23	
Dibromofluoromethane	102	89 - 119	07/07/20 17:23	
Toluene-d8	100	87 - 121	07/07/20 17:23	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007251-04Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

Result

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/02/20

Sample Matrix: Water

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005520

Lab Control Sample

RQ2007077-04

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
1,1,1-Trichloroethane (TCA)	8260C	22.5	20.0	113	75-125
1,1,2,2-Tetrachloroethane	8260C	28.8	20.0	144 *	78-126
1,1,2-Trichloroethane	8260C	21.8	20.0	109	82-121
1,1-Dichloroethane (1,1-DCA)	8260C	23.3	20.0	117	80-124
1,1-Dichloroethene (1,1-DCE)	8260C	21.1	20.0	105	71-118
1,2-Dichloroethane	8260C	20.4	20.0	102	71-127
1,2-Dichloropropane	8260C	22.9	20.0	114	80-119
2-Butanone (MEK)	8260C	23.7	20.0	118	61-137
2-Hexanone	8260C	22.8	20.0	114	63-124
4-Methyl-2-pentanone	8260C	22.8	20.0	114	66-124
Acetone	8260C	25.8	20.0	129	40-161
Benzene	8260C	21.7	20.0	109	79-119
Bromodichloromethane	8260C	21.7	20.0	109	81-123
Bromoform	8260C	21.9	20.0	110	65-146
Bromomethane	8260C	17.7	20.0	89	42-166
Carbon Disulfide	8260C	21.0	20.0	105	66-128
Carbon Tetrachloride	8260C	20.4	20.0	102	70-127
Chlorobenzene	8260C	21.4	20.0	107	80-121
Chloroethane	8260C	22.3	20.0	112	62-131
Chloroform	8260C	22.2	20.0	111	79-120
Chloromethane	8260C	22.1	20.0	110	65-135
Dibromochloromethane	8260C	22.5	20.0	112	72-128
Dichloromethane	8260C	21.8	20.0	109	73-122
Ethylbenzene	8260C	22.1	20.0	110	76-120
Styrene	8260C	22.0	20.0	110	80-124
Tetrachloroethene (PCE)	8260C	20.4	20.0	102	72-125
Toluene	8260C	21.8	20.0	109	79-119
Trichloroethene (TCE)	8260C	17.7	20.0	89	74-122
Vinyl Acetate	8260C	30.7	20.0	154	52-174
Vinyl Chloride	8260C	23.1	20.0	116	74-159
cis-1,2-Dichloroethene	8260C	22.8	20.0	114	80-121
cis-1,3-Dichloropropene	8260C	22.4	20.0	112	77-122
trans-1,2-Dichloroethene	8260C	21.9	20.0	110	73-118
Printed 7/16/2020 11:08:09 AM			Supers	et Reference:20-000	00554598 rev 00

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix:

Project:

Water

Service Request: R2005520 Date Analyzed: 07/02/20

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

> Units:ug/L Basis:NA

Lab Control Sample

RQ2007077-04

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
trans-1 3-Dichloropropene	8260C	23.4	20.0	117	71-133

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring **Date Analyzed:** 06/30/20

Sample Matrix: Water

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005520

Lab Control Sample

RQ2007141-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
1,1,1-Trichloroethane (TCA)	8260C	21.6	20.0	108	75-125
1,1,2,2-Tetrachloroethane	8260C	26.2	20.0	131 *	78-126
1,1,2-Trichloroethane	8260C	22.0	20.0	110	82-121
1,1-Dichloroethane (1,1-DCA)	8260C	20.8	20.0	104	80-124
1,1-Dichloroethene (1,1-DCE)	8260C	20.3	20.0	101	71-118
1,2-Dichloroethane	8260C	19.5	20.0	98	71-127
1,2-Dichloropropane	8260C	21.2	20.0	106	80-119
2-Butanone (MEK)	8260C	16.9	20.0	85	61-137
2-Hexanone	8260C	17.3	20.0	87	63-124
4-Methyl-2-pentanone	8260C	17.2	20.0	86	66-124
Acetone	8260C	20.1	20.0	101	40-161
Benzene	8260C	21.2	20.0	106	79-119
Bromodichloromethane	8260C	21.4	20.0	107	81-123
Bromoform	8260C	22.5	20.0	113	65-146
Bromomethane	8260C	14.4	20.0	72	42-166
Carbon Disulfide	8260C	17.2	20.0	86	66-128
Carbon Tetrachloride	8260C	21.4	20.0	107	70-127
Chlorobenzene	8260C	21.9	20.0	109	80-121
Chloroethane	8260C	16.2	20.0	81	62-131
Chloroform	8260C	21.1	20.0	105	79-120
Chloromethane	8260C	17.5	20.0	87	65-135
Dibromochloromethane	8260C	23.1	20.0	116	72-128
Dichloromethane	8260C	20.7	20.0	103	73-122
Ethylbenzene	8260C	22.5	20.0	113	76-120
Styrene	8260C	22.5	20.0	112	80-124
Tetrachloroethene (PCE)	8260C	21.6	20.0	108	72-125
Toluene	8260C	21.5	20.0	107	79-119
Trichloroethene (TCE)	8260C	18.3	20.0	91	74-122
Vinyl Acetate	8260C	23.8	20.0	119	52-174
Vinyl Chloride	8260C	17.0	20.0	85	74-159
cis-1,2-Dichloroethene	8260C	22.1	20.0	111	80-121
cis-1,3-Dichloropropene	8260C	21.4	20.0	107	77-122
trans-1,2-Dichloroethene	8260C	21.1	20.0	105	73-118
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QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix:

Project:

Water

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

> Units:ug/L Basis:NA

Service Request: R2005520

Date Analyzed: 06/30/20

Lab Control Sample

RQ2007141-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
trans-1.3-Dichloropropene	8260C	21.6	20.0	108	71-133

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring **Date Analyzed:** 07/07/20

Sample Matrix: Water

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005520

Lab Control Sample

RQ2007251-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
1,1,1-Trichloroethane (TCA)	8260C	19.1	20.0	95	75-125
1,1,2,2-Tetrachloroethane	8260C	21.7	20.0	108	78-126
1,1,2-Trichloroethane	8260C	20.6	20.0	103	82-121
1,1-Dichloroethane (1,1-DCA)	8260C	18.6	20.0	93	80-124
1,1-Dichloroethene (1,1-DCE)	8260C	22.3	20.0	112	71-118
1,2-Dichloroethane	8260C	19.0	20.0	95	71-127
1,2-Dichloropropane	8260C	18.3	20.0	92	80-119
2-Butanone (MEK)	8260C	21.7	20.0	109	61-137
2-Hexanone	8260C	19.5	20.0	98	63-124
4-Methyl-2-pentanone	8260C	20.4	20.0	102	66-124
Acetone	8260C	26.2	20.0	131	40-161
Benzene	8260C	19.2	20.0	96	79-119
Bromodichloromethane	8260C	19.2	20.0	96	81-123
Bromoform	8260C	21.7	20.0	109	65-146
Bromomethane	8260C	19.2	20.0	96	42-166
Carbon Disulfide	8260C	23.7	20.0	118	66-128
Carbon Tetrachloride	8260C	19.7	20.0	99	70-127
Chlorobenzene	8260C	19.2	20.0	96	80-121
Chloroethane	8260C	16.7	20.0	84	62-131
Chloroform	8260C	20.1	20.0	100	79-120
Chloromethane	8260C	17.5	20.0	88	65-135
Dibromochloromethane	8260C	20.9	20.0	104	72-128
Dichloromethane	8260C	19.6	20.0	98	73-122
Ethylbenzene	8260C	19.2	20.0	96	76-120
Styrene	8260C	19.4	20.0	97	80-124
Tetrachloroethene (PCE)	8260C	18.2	20.0	91	72-125
Toluene	8260C	19.4	20.0	97	79-119
Trichloroethene (TCE)	8260C	18.5	20.0	93	74-122
Vinyl Acetate	8260C	37.2	20.0	186 *	52-174
Vinyl Chloride	8260C	17.2	20.0	86	74-159
cis-1,2-Dichloroethene	8260C	20.8	20.0	104	80-121
cis-1,3-Dichloropropene	8260C	19.1	20.0	96	77-122
trans-1,2-Dichloroethene	8260C	20.5	20.0	102	73-118
Printed 7/16/2020 11:08:11 AM			Supers	et Reference:20-000	00554598 rev 00

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix: Water

Project:

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

> Units:ug/L Basis:NA

Service Request: R2005520

Date Analyzed: 07/07/20

Lab Control Sample

RQ2007251-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
trans-1 3-Dichloropropene	8260C	20.9	20.0	104	71-133



Semivolatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

SURROGATE RECOVERY SUMMARY Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Extraction Method:** EPA 3510C

		2,4,6-Tribromophenol	2-Fluorobiphenyl	2-Fluorophenol
Sample Name	Lab Code	35-141	31-118	10-105
WG-9954-062520-SG-013	R2005520-001	104	69	38
WG-9954-062520-SG-011	R2005520-002	93	70	43
WG-9954-062520-SG-008	R2005520-003	95	74	45
WG-9954-062520-SG-009	R2005520-004	81	64	38
WG-9954-062520-SG-012	R2005520-005	83	59	40
RB-9954-062520-SG-001	R2005520-006	89	78	39
WG-9954-062520-SG-010	R2005520-008	79	67	44
Method Blank	RQ2006968-03	78	71	44
Lab Control Sample	RQ2006968-04	100	80	56
Duplicate Lab Control Sample	RQ2006968-05	64	54	39
WG-9954-062520-SG-012 MS	RQ2006968-01	107	78	43
WG-9954-062520-SG-012 DMS	RQ2006968-02	97	71	43

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

SURROGATE RECOVERY SUMMARY Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Extraction Method:** EPA 3510C

		Nitrobenzene-d5	Phenol-d6	p-Terphenyl-d14
Sample Name	Lab Code	31-110	10-107	10-165
WG-9954-062520-SG-013	R2005520-001	61	32	92
WG-9954-062520-SG-011	R2005520-002	68	30	93
WG-9954-062520-SG-008	R2005520-003	69	31	94
WG-9954-062520-SG-009	R2005520-004	66	28	87
WG-9954-062520-SG-012	R2005520-005	59	28	93
RB-9954-062520-SG-001	R2005520-006	68	28	94
WG-9954-062520-SG-010	R2005520-008	68	29	92
Method Blank	RQ2006968-03	65	31	95
Lab Control Sample	RQ2006968-04	85	41	104
Duplicate Lab Control Sample	RQ2006968-05	54	29	78
WG-9954-062520-SG-012 MS	RQ2006968-01	77	32	94
WG-9954-062520-SG-012 DMS	RQ2006968-02	68	34	92

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Date Collected: 06/25/20

Sample Matrix: Water

Date Received: 06/26/20 **Date Analyzed:** 07/1/20

Date Extracted: 06/30/20

Duplicate Matrix Spike Summary Semivolatile Organic Compounds by GC/MS

Sample Name: WG-9954-062520-SG-012

Units: ug/L

Lab Code: R2005520-005

Basis: NA

Analysis Method: 8270D **Prep Method:** EPA 3510C

		Matrix Spike Duplicate Matrix Spike								
		RQ2006968-01			R	Q2006968-0	02			
	Sample		Spike			Spike		% Rec		RPD
Analyte Name	Result	Result	Amount	% Rec	Result	Amount	% Rec	Limits	RPD	Limit
1,2,4-Trichlorobenzene	9.1 U	44.7	72.7	61	42.3	72.7	58	10-127	5	30
1,2-Dichlorobenzene	9.1 U	39.9	72.7	55	40.1	72.7	55	17-105	<1	30
1,3-Dichlorobenzene	9.1 U	38.5	72.7	53	39.5	72.7	54	21-99	2	30
1,4-Dichlorobenzene	9.1 U	41.1	72.7	56	37.8	72.7	52	10-124	7	30
2,4,5-Trichlorophenol	9.1 U	62.7	72.7	86	59.6	72.7	82	48-134	5	30
2,4,6-Trichlorophenol	9.1 U	59.8	72.7	82	55.5	72.7	76	44-135	8	30
2,4-Dichlorophenol	9.1 U	48.4	72.7	67	48.0	72.7	66	40-130	2	30
2,4-Dimethylphenol	9.1 U	54.2	72.7	75	55.4	72.7	76	42-121	1	30
2,4-Dinitrophenol	45 U	53.5	72.7	74	49.6	72.7	68	21-168	8	30
2,4-Dinitrotoluene	9.1 U	72.1	72.7	99	68.2	72.7	94	37-143	5	30
2,6-Dinitrotoluene	9.1 U	78.8	72.7	108	72.3	72.7	99	39-136	9	30
2-Chloronaphthalene	9.1 U	58.5	72.7	80	54.1	72.7	74	40-108	8	30
2-Chlorophenol	9.1 U	41.7	72.7	57	42.2	72.7	58	37-112	2	30
2-Methylnaphthalene	9.1 U	52.7	72.7	72	50.2	72.7	69	34-102	4	30
2-Methylphenol	9.1 U	47.7	72.7	66	47.0	72.7	65	37-102	2	30
2-Nitroaniline	9.1 U	73.3	72.7	101	73.9	72.7	102	40-136	<1	30
2-Nitrophenol	9.1 U	48.8	72.7	67	49.7	72.7	68	27-143	1	30
3,3'-Dichlorobenzidine	9.1 U	65.2	72.7	90	62.7	72.7	86	11-131	5	30
3- and 4-Methylphenol Coelution	9.1 U	45.1	72.7	62	44.4	72.7	61	30-95	2	30
3-Nitroaniline	9.1 U	69.3	72.7	95	69.3	72.7	95	19-117	<1	30
4,6-Dinitro-2-methylphenol	45 U	58.3	72.7	80	57.5	72.7	79	25-154	1	30
4-Bromophenyl Phenyl Ether	9.1 U	64.3	72.7	88	57.1	72.7	78	39-115	12	30
4-Chloro-3-methylphenol	9.1 U	60.9	72.7	84	59.9	72.7	82	41-126	2	30
4-Chloroaniline	9.1 U	58.3	72.7	80	63.6	72.7	87	19-111	8	30
4-Chlorophenyl Phenyl Ether	9.1 U	57.4	72.7	79	52.0	72.7	72	41-111	9	30
4-Nitroaniline	9.1 U	76.4	72.7	105	85.0	72.7	117	18-143	11	30
4-Nitrophenol	45 U	26.2 J	72.7	36	28.7 J	72.7	39	10-126	8	30
Acenaphthene	9.1 U	63.6	72.7	87	58.4	72.7	80	43-117	8	30
Acenaphthylene	9.1 U	66.8	72.7	92	61.4	72.7	84	45-119	9	30
Anthracene	9.1 U	70.5	72.7	97	63.5	72.7	87	45-127	11	30
Benz(a)anthracene	9.1 U	56.7	72.7	78	53.0	72.7	73	46-126	7	30
Benzo(a)pyrene	9.1 U	58.0	72.7	80	56.8	72.7	78	44-114	3	30
Benzo(b)fluoranthene	9.1 U	51.5	72.7	71	48.5	72.7	67	41-127	6	30

Results flagged with an asterisk (*) indicate values outside control criteria.

Results flagged with a pound (#) indicate the control criteria is not applicable.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

Matrix Spike and Matrix Spike Duplicate Data is presented for information purposes only. The matrix may or may not be relevant to samples reported in this report. The laboratory evaluates system performance based on the LCS and LCSD control limits.

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates) **Service Request:**

R2005520

Project:

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Date Collected:

06/25/20

Sample Matrix:

Water

Date Received:

06/26/20

Date Analyzed: **Date Extracted:**

07/1/20 06/30/20

Duplicate Matrix Spike Summary Semivolatile Organic Compounds by GC/MS

Matrix Spike

Sample Name: WG-9954-062520-SG-012 **Units:**

Basis:

ug/L NA

Lab Code:

R2005520-005

Analysis Method: 8270D **Prep Method:** EPA 3510C

Duplicate Matrix Spike

		RQ2006968-01 RQ				RQ2006968-02				
	Sample		Spike			Spike		% Rec		RPD
Analyte Name	Result	Result	Amount	% Rec	Result	Amount	% Rec	Limits	RPD	Limit
Benzo(g,h,i)perylene	9.1 U	56.3	72.7	77	55.5	72.7	76	50-143	1	30
Benzo(k)fluoranthene	9.1 U	55.7	72.7	77	55.0	72.7	76	46-139	1	30
Benzoic Acid	91 U	74.4 J	109	68	69.3 J	109	64	10-94	6	30
Benzyl Alcohol	9.1 U	56.4	72.7	78	57.0	72.7	78	31-109	<1	30
2,2'-Oxybis(1-chloropropane)	9.1 U	49.9	72.7	69	49.0	72.7	67	21-126	3	30
Bis(2-chloroethoxy)methane	9.1 U	53.7	72.7	74	54.6	72.7	75	41-118	1	30
Bis(2-chloroethyl) Ether	9.1 U	44.9	72.7	62	45.3	72.7	62	33-108	<1	30
Bis(2-ethylhexyl) Phthalate	9.1 U	52.9	72.7	73	50.7	72.7	70	41-132	4	30
Butyl Benzyl Phthalate	9.1 U	62.9	72.7	86	61.7	72.7	85	41-148	1	30
Chrysene	9.1 U	59.0	72.7	81	55.6	72.7	76	47-126	6	30
Di-n-butyl Phthalate	9.1 U	73.3	72.7	101	66.4	72.7	91	43-130	10	30
Di-n-octyl Phthalate	9.1 U	50.7	72.7	70	54.0	72.7	74	40-139	6	30
Dibenz(a,h)anthracene	9.1 U	60.8	72.7	84	59.2	72.7	81	43-136	4	30
Dibenzofuran	9.1 U	68.3	72.7	94	61.8	72.7	85	46-119	10	30
Diethyl Phthalate	9.1 U	65.9	72.7	91	63.1	72.7	87	36-122	4	30
Dimethyl Phthalate	9.1 U	72.3	72.7	99	67.4	72.7	93	33-123	6	30
Fluoranthene	9.1 U	73.9	72.7	102	65.8	72.7	90	43-135	13	30
Fluorene	9.1 U	69.5	72.7	96	63.1	72.7	87	43-113	10	30
Hexachlorobenzene	9.1 U	67.8	72.7	93	60.1	72.7	83	42-125	11	30
Hexachlorobutadiene	9.1 U	44.8	72.7	62	44.9	72.7	62	10-111	<1	30
Hexachlorocyclopentadiene	9.1 U	4.05 J	72.7	6 *	5.07 J	72.7	7 *	10-103	15	30
Hexachloroethane	9.1 U	39.8	72.7	55	39.6	72.7	54	12-101	2	30
Indeno(1,2,3-cd)pyrene	9.1 U	50.9	72.7	70	48.4	72.7	67	49-140	4	30
Isophorone	9.1 U	49.1	72.7	68	46.4	72.7	64	40-111	6	30
N-Nitrosodi-n-propylamine	9.1 U	61.9	72.7	85	58.9	72.7	81	35-108	5	30
N-Nitrosodiphenylamine	9.1 U	83.7	72.7	115	77.1	72.7	106	43-127	8	30
Naphthalene	9.1 U	49.1	72.7	68	48.5	72.7	67	37-108	1	30
Nitrobenzene	9.1 U	52.1	72.7	72	48.4	72.7	67	35-112	7	30
Pentachlorophenol (PCP)	45 U	71.7	72.7	99	61.1	72.7	84	29-164	16	30
Phenanthrene	9.1 U	69.0	72.7	95	63.6	72.7	87	46-123	9	30
Phenol	9.1 U	26.4	72.7	36	28.4	72.7	39	10-113	8	30
Pyrene	9.1 U	72.3	72.7	99	70.5	72.7	97	44-129	2	30

Results flagged with an asterisk (*) indicate values outside control criteria.

Results flagged with a pound (#) indicate the control criteria is not applicable.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

Matrix Spike and Matrix Spike Duplicate Data is presented for information purposes only. The matrix may or may not be relevant to samples reported in this report. The laboratory evaluates system performance based on the LCS and LCSD control limits.

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2006968-03
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	10 U	10	1.2	1	07/01/20 17:03	6/30/20	
1,2-Dichlorobenzene	10 U	10	1.2	1	07/01/20 17:03	6/30/20	
1,3-Dichlorobenzene	10 U	10	1.1	1	07/01/20 17:03	6/30/20	
1,4-Dichlorobenzene	10 U	10	1.2	1	07/01/20 17:03	6/30/20	
2,4,5-Trichlorophenol	10 U	10	1.1	1	07/01/20 17:03	6/30/20	
2,4,6-Trichlorophenol	10 U	10	1.4	1	07/01/20 17:03	6/30/20	
2,4-Dichlorophenol	10 U	10	1.3	1	07/01/20 17:03	6/30/20	
2,4-Dimethylphenol	10 U	10	1.4	1	07/01/20 17:03	6/30/20	
2,4-Dinitrophenol	50 U	50	20	1	07/01/20 17:03	6/30/20	
2,4-Dinitrotoluene	10 U	10	2.4	1	07/01/20 17:03	6/30/20	
2,6-Dinitrotoluene	10 U	10	1.4	1	07/01/20 17:03	6/30/20	
2-Chloronaphthalene	10 U	10	1.4	1	07/01/20 17:03	6/30/20	
2-Chlorophenol	10 U	10	1.1	1	07/01/20 17:03	6/30/20	
2-Methylnaphthalene	10 U	10	1.3	1	07/01/20 17:03	6/30/20	
2-Methylphenol	10 U	10	1.0	1	07/01/20 17:03	6/30/20	
2-Nitroaniline	10 U	10	1.4	1	07/01/20 17:03	6/30/20	
2-Nitrophenol	10 U	10	1.5	1	07/01/20 17:03	6/30/20	
3,3'-Dichlorobenzidine	10 U	10	1.2	1	07/01/20 17:03	6/30/20	
3- and 4-Methylphenol Coelution	10 U	10	1.2	1	07/01/20 17:03	6/30/20	
3-Nitroaniline	10 U	10	2.5	1	07/01/20 17:03	6/30/20	
4,6-Dinitro-2-methylphenol	50 U	50	20	1	07/01/20 17:03	6/30/20	
4-Bromophenyl Phenyl Ether	10 U	10	1.7	1	07/01/20 17:03	6/30/20	
4-Chloro-3-methylphenol	10 U	10	1.1	1	07/01/20 17:03	6/30/20	
4-Chloroaniline	10 U	10	1.0	1	07/01/20 17:03	6/30/20	
4-Chlorophenyl Phenyl Ether	10 U	10	1.5	1	07/01/20 17:03	6/30/20	
4-Nitroaniline	10 U	10	2.7	1	07/01/20 17:03	6/30/20	
4-Nitrophenol	50 U	50	6.4	1	07/01/20 17:03	6/30/20	
Acenaphthene	10 U	10	1.4	1	07/01/20 17:03	6/30/20	
Acenaphthylene	10 U	10	1.4	1	07/01/20 17:03	6/30/20	
Anthracene	10 U	10	1.3	1	07/01/20 17:03	6/30/20	
Benz(a)anthracene	10 U	10	1.6	1	07/01/20 17:03	6/30/20	
Benzo(a)pyrene	10 U	10	1.2	1	07/01/20 17:03	6/30/20	
Benzo(b)fluoranthene	10 U	10	1.2	1	07/01/20 17:03	6/30/20	
Benzo(g,h,i)perylene	10 U	10	1.0	1	07/01/20 17:03	6/30/20	
Benzo(k)fluoranthene	10 U	10	1.3	1	07/01/20 17:03	6/30/20	
Benzoic Acid	100 U	100	36	1	07/01/20 17:03	6/30/20	
Benzyl Alcohol	10 U	10	1.6	1	07/01/20 17:03	6/30/20	
2,2'-Oxybis(1-chloropropane)	10 U	10	1.4	1	07/01/20 17:03	6/30/20	
Bis(2-chloroethoxy)methane	10 U	10	1.9	1	07/01/20 17:03	6/30/20	
Bis(2-chloroethyl) Ether	10 U	10	1.3	1	07/01/20 17:03	6/30/20	
Bis(2-ethylhexyl) Phthalate	10 U	10	1.0	1	07/01/20 17:03	6/30/20	
Butyl Benzyl Phthalate	10 U	10	1.4	1	07/01/20 17:03	6/30/20	
Chrysene	10 U	10	1.2	1	07/01/20 17:03	6/30/20	
,	10 0	10		-	001.20 17.03	0, 2 0, 20	

Printed 7/16/2020 11:08:30 AM

Superset Reference: 20-0000554598 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2006968-03Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	10 U	10	2.0	1	07/01/20 17:03	6/30/20	
Di-n-octyl Phthalate	10 U	10	3.3	1	07/01/20 17:03	6/30/20	
Dibenz(a,h)anthracene	10 U	10	1.1	1	07/01/20 17:03	6/30/20	
Dibenzofuran	10 U	10	1.4	1	07/01/20 17:03	6/30/20	
Diethyl Phthalate	10 U	10	1.1	1	07/01/20 17:03	6/30/20	
Dimethyl Phthalate	10 U	10	1.3	1	07/01/20 17:03	6/30/20	
Fluoranthene	10 U	10	1.5	1	07/01/20 17:03	6/30/20	
Fluorene	10 U	10	1.3	1	07/01/20 17:03	6/30/20	
Hexachlorobenzene	10 U	10	1.6	1	07/01/20 17:03	6/30/20	
Hexachlorobutadiene	10 U	10	1.0	1	07/01/20 17:03	6/30/20	
Hexachlorocyclopentadiene	10 U	10	2.2	1	07/01/20 17:03	6/30/20	
Hexachloroethane	10 U	10	1.1	1	07/01/20 17:03	6/30/20	
Indeno(1,2,3-cd)pyrene	10 U	10	1.8	1	07/01/20 17:03	6/30/20	
Isophorone	10 U	10	1.4	1	07/01/20 17:03	6/30/20	
N-Nitrosodi-n-propylamine	10 U	10	1.2	1	07/01/20 17:03	6/30/20	
N-Nitrosodiphenylamine	10 U	10	2.7	1	07/01/20 17:03	6/30/20	
Naphthalene	10 U	10	1.2	1	07/01/20 17:03	6/30/20	
Nitrobenzene	10 U	10	1.5	1	07/01/20 17:03	6/30/20	
Pentachlorophenol (PCP)	50 U	50	9.8	1	07/01/20 17:03	6/30/20	
Phenanthrene	10 U	10	1.4	1	07/01/20 17:03	6/30/20	
Phenol	10 U	10	1.0	1	07/01/20 17:03	6/30/20	
Pyrene	10 U	10	1.5	1	07/01/20 17:03	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	78	35 - 141	07/01/20 17:03	
2-Fluorobiphenyl	71	31 - 118	07/01/20 17:03	
2-Fluorophenol	44	10 - 105	07/01/20 17:03	
Nitrobenzene-d5	65	31 - 110	07/01/20 17:03	
Phenol-d6	31	10 - 107	07/01/20 17:03	
p-Terphenyl-d14	95	10 - 165	07/01/20 17:03	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	

No Tentatively Identified Compounds

Detected

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/01/20

Sample Matrix: Water

Duplicate Lab Control Sample Summary Semivolatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005520

Lab Control Sample

Duplicate Lab Control Sample

RQ2006968-04

RQ2006968-05

Analyte Name	Analytica l Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
1,2,4-Trichlorobenzene	8270D	60.4	80.0	76	39.0	80.0	49	10-127	43*	30
1,2-Dichlorobenzene	8270D	57.5	80.0	72	37.6	80.0	47	23-130	42*	30
1,3-Dichlorobenzene	8270D	57.0	80.0	71	36.9	80.0	46	21-90	43*	30
1,4-Dichlorobenzene	8270D	55.7	80.0	70	36.5	80.0	46	10-124	41*	30
2,4,5-Trichlorophenol	8270D	69.0	80.0	86	46.8	80.0	58	48-134	39*	30
2,4,6-Trichlorophenol	8270D	65.6	80.0	82	42.1	80.0	53	44-135	43*	30
2,4-Dichlorophenol	8270D	62.6	80.0	78	41.4	80.0	52	48-127	40*	30
2,4-Dimethylphenol	8270D	65.7	80.0	82	44.7	80.0	56 *	59-113	38*	30
2,4-Dinitrophenol	8270D	60.5	80.0	76	38.9 J	80.0	49	21-154	43*	30
2,4-Dinitrotoluene	8270D	70.4	80.0	88	46.4	80.0	58	54-130	41*	30
2,6-Dinitrotoluene	8270D	79.2	80.0	99	54.5	80.0	68	51-127	37*	30
2-Chloronaphthalene	8270D	66.2	80.0	83	44.5	80.0	56	40-108	39*	30
2-Chlorophenol	8270D	58.8	80.0	74	37.5	80.0	47	42-112	45*	30
2-Methylnaphthalene	8270D	62.8	80.0	79	39.8	80.0	50	34-102	45*	30
2-Methylphenol	8270D	61.4	80.0	77	44.0	80.0	55	47-100	33*	30
2-Nitroaniline	8270D	71.5	80.0	89	47.7	80.0	60	52-133	39*	30
2-Nitrophenol	8270D	64.1	80.0	80	41.7	80.0	52	43-131	42*	30
3,3'-Dichlorobenzidine	8270D	66.6	80.0	83	43.9	80.0	55	43-126	41*	30
3- and 4-Methylphenol Coelution	8270D	55.1	80.0	69	38.5	80.0	48	40-92	36*	30
3-Nitroaniline	8270D	67.2	80.0	84	55.6	80.0	69	42-111	20	30
4,6-Dinitro-2-methylphenol	8270D	60.3	80.0	75	41.1 J	80.0	51	36-152	38*	30
4-Bromophenyl Phenyl Ether	8270D	69.7	80.0	87	50.5	80.0	63	48-114	32*	30
4-Chloro-3-methylphenol	8270D	66.1	80.0	83	40.6	80.0	51 *	52-113	48*	30
4-Chloroaniline	8270D	64.5	80.0	81	54.9	80.0	69	44-109	16	30
4-Chlorophenyl Phenyl Ether	8270D	62.5	80.0	78	43.8	80.0	55	51-107	35*	30
4-Nitroaniline	8270D	63.5	80.0	79	42.6	80.0	53 *	54-133	39*	30
4-Nitrophenol	8270D	27.1 J	80.0	34	20.8 J	80.0	26	10-126	27	30
Acenaphthene	8270D	69.8	80.0	87	46.0	80.0	57	52-107	42*	30
Acenaphthylene	8270D	73.7	80.0	92	49.7	80.0	62	55-109	39*	30
Anthracene	8270D	74.4	80.0	93	49.5	80.0	62	55-116	40*	30
Benz(a)anthracene	8270D	69.6	80.0	87	49.6	80.0	62	61-121	34*	30
Benzo(a)pyrene	8270D	75.2	80.0	94	52.1	80.0	65	44-114	36*	30
Benzo(b)fluoranthene	8270D	72.6	80.0	91	50.6	80.0	63	62-115	36*	30
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Superset Reference: 20-0000554598 rev 00

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/01/20

Sample Matrix: Water

Duplicate Lab Control Sample Summary Semivolatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005520

Lab Control Sample

Duplicate Lab Control Sample

RQ2006968-04

RQ2006968-05

Analyte Name	Analytica l Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
Benzo(g,h,i)perylene	8270D	76.4	80.0	95	56.5	80.0	71	63-136	29	30
Benzo(k)fluoranthene	8270D	77.9	80.0	97	56.2	80.0	70	49-133	32*	30
Benzoic Acid	8270D	65.8 J	120	55	56.4 J	120	47	10-94	16	30
Benzyl Alcohol	8270D	63.3	80.0	79	47.2	80.0	59	31-109	29	30
2,2'-Oxybis(1-chloropropane)	8270D	64.2	80.0	80	41.7	80.0	52	32-122	42*	30
Bis(2-chloroethoxy)methane	8270D	65.9	80.0	82	45.6	80.0	57	55-110	36*	30
Bis(2-chloroethyl) Ether	8270D	61.5	80.0	77	39.9	80.0	50	46-102	43*	30
Bis(2-ethylhexyl) Phthalate	8270D	73.4	80.0	92	51.7	80.0	65	51-132	34*	30
Butyl Benzyl Phthalate	8270D	75.3	80.0	94	49.3	80.0	62	41-148	41*	30
Chrysene	8270D	73.1	80.0	91	53.9	80.0	67	57-118	30	30
Di-n-butyl Phthalate	8270D	83.3	80.0	104	52.5	80.0	66	57-128	45*	30
Di-n-octyl Phthalate	8270D	80.2	80.0	100	53.2	80.0	66	62-124	41*	30
Dibenz(a,h)anthracene	8270D	82.6	80.0	103	62.0	80.0	78	54-135	28	30
Dibenzofuran	8270D	72.8	80.0	91	49.0	80.0	61	55-110	39*	30
Diethyl Phthalate	8270D	65.2	80.0	82	45.6	80.0	57	53-113	36*	30
Dimethyl Phthalate	8270D	75.9	80.0	95	50.9	80.0	64	51-112	39*	30
Fluoranthene	8270D	80.7	80.0	101	54.0	80.0	68	66-127	39*	30
Fluorene	8270D	72.2	80.0	90	49.5	80.0	62	54-106	37*	30
Hexachlorobenzene	8270D	79.9	80.0	100	58.9	80.0	74	53-123	30	30
Hexachlorobutadiene	8270D	62.6	80.0	78	40.2	80.0	50	16-95	44*	30
Hexachlorocyclopentadiene	8270D	26.8	80.0	33	18.1	80.0	23	10-99	36*	30
Hexachloroethane	8270D	55.2	80.0	69	35.9	80.0	45	15-92	42*	30
Indeno(1,2,3-cd)pyrene	8270D	70.3	80.0	88	49.6	80.0	62	62-137	35*	30
Isophorone	8270D	61.1	80.0	76	37.2	80.0	46 *	50-116	49*	30
N-Nitrosodi-n-propylamine	8270D	70.8	80.0	89	45.4	80.0	57	49-115	44*	30
N-Nitrosodiphenylamine	8270D	82.8	80.0	103	60.2	80.0	75	45-123	31*	30
Naphthalene	8270D	64.9	80.0	81	42.8	80.0	53	38-99	42*	30
Nitrobenzene	8270D	68.2	80.0	85	42.7	80.0	53	46-108	46*	30
Pentachlorophenol (PCP)	8270D	78.0	80.0	97	49.9 J	80.0	62	29-164	44*	30
Phenanthrene	8270D	72.6	80.0	91	49.0	80.0	61	58-118	39*	30
Phenol	8270D	36.4	80.0	46	27.1	80.0	34	10-113	30	30
Pyrene	8270D	78.0	80.0	98	52.0	80.0	65	61-122	40*	30

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Semivolatile Organic Compounds by GC

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

SURROGATE RECOVERY SUMMARY

Organochlorine Pesticides by Gas Chromatography

Analysis Method: 8081B **Extraction Method:** EPA 3510C

		Decachlorobiphenyl	Tetrachloro-m-xylene	
Sample Name	Lab Code	10-164	10-147	
WG-9954-062520-SG-013	R2005520-001	67	80	
WG-9954-062520-SG-011	R2005520-002	56	71	
WG-9954-062520-SG-008	R2005520-003	65	74	
WG-9954-062520-SG-009	R2005520-004	37	57	
WG-9954-062520-SG-012	R2005520-005	49	85	
RB-9954-062520-SG-001	R2005520-006	10	53	
WG-9954-062520-SG-010	R2005520-008	27	45	
Method Blank	RQ2006967-03	58	59	
Method Blank	RQ2007029-01	84	75	
Lab Control Sample	RQ2006967-04	71	67	
Duplicate Lab Control Sample	RQ2006967-05	60	57	
Lab Control Sample	RQ2007029-02	63	60	
Duplicate Lab Control Sample	RQ2007029-03	68	58	
WG-9954-062520-SG-012 MS	RQ2006967-01	42	67	
WG-9954-062520-SG-012 DMS	RQ2006967-02	43	66	

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Date Collected: 06/25/20 **Date Received:** 06/26/20

Water Date Received:

07/8/20

Date Analyzed:
Date Extracted:

06/30/20

Duplicate Matrix Spike Summary Organochlorine Pesticides by Gas Chromatography

Sample Name: WG-9954-062520-SG-012

Units: ug/L

Lab Code: R2005520-005

Basis: NA

Analysis Method: 8081B **Prep Method:** EPA 3510C

Sample Matrix:

			Matrix	Spike		Duplicate Ma	ıtrix Spike	!		
			RQ20069	967-01		RQ20069	67-02			
	Sample		Spike			Spike		% Rec		RPD
Analyte Name	Result	Result	Amount	% Rec	Result	Amount	% Rec	Limits	RPD	Limit
4,4'-DDD	0.045 U	0.244	0.364	67	0.270	0.364	74	38-157	10	30
4,4'-DDE	0.045 U	0.236	0.364	65	0.261	0.364	72	10-200	10	30
4,4'-DDT	0.045 U	0.237	0.364	65	0.260	0.364	72	19-154	10	30
Aldrin	0.045 U	0.214	0.364	59	0.239	0.364	66	26-149	11	30
Dieldrin	0.045 U	0.249	0.364	68	0.279	0.364	77	41-164	11	30
Endosulfan I	0.045 U	0.251	0.364	69	0.280	0.364	77	47-149	11	30
Endosulfan II	0.045 U	0.254	0.364	70	0.285	0.364	78	51-148	12	30
Endosulfan Sulfate	0.045 U	0.234	0.364	64	0.266	0.364	73	10-170	13	30
Endrin	0.045 U	0.255	0.364	70	0.285	0.364	78	48-165	11	30
Endrin Ketone	0.045 U	0.259	0.364	71	0.290	0.364	80	48-162	11	30
Heptachlor	0.045 U	0.181	0.364	50	0.200	0.364	55	29-168	10	30
Heptachlor Epoxide	0.045 U	0.249	0.364	69	0.280	0.364	77	29-180	12	30
Methoxychlor	0.045 U	0.263	0.364	72	0.289	0.364	80	38-162	10	30
alpha-BHC	0.033 J	0.268	0.364	65	0.303	0.364	74	27-154	12	30
alpha-Chlordane	0.045 U	0.243	0.364	67	0.272	0.364	75	35-160	11	30
beta-BHC	0.045 U	0.271	0.364	75	0.303	0.364	83	32-184	11	30
delta-BHC	0.068	0.272	0.364	56	0.317	0.364	68	10-182	15	30
gamma-BHC (Lindane)	0.037 J	0.268	0.364	64	0.305	0.364	74	43-164	13	30
gamma-Chlordane	0.028 JP	0.246	0.364	60	0.272	0.364	67	35-165	10	30

Results flagged with an asterisk (*) indicate values outside control criteria.

Results flagged with a pound (#) indicate the control criteria is not applicable.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

Matrix Spike and Matrix Spike Duplicate Data is presented for information purposes only. The matrix may or may not be relevant to samples reported in this report. The laboratory evaluates system performance based on the LCS and LCSD control limits.

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2006967-03Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analysis Method: 8081B **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	_
4,4'-DDE	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
4,4'-DDT	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
Aldrin	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
Dieldrin	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
Endosulfan I	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
Endosulfan II	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
Endosulfan Sulfate	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
Endrin	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
Endrin Ketone	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
Heptachlor	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
Heptachlor Epoxide	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
Methoxychlor	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
Toxaphene	0.50 U	0.50	0.50	1	07/01/20 16:36	6/30/20	
alpha-BHC	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
alpha-Chlordane	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
beta-BHC	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
delta-BHC	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
gamma-BHC (Lindane)	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	
gamma-Chlordane	0.050 U	0.050	0.020	1	07/01/20 16:36	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	58	10 - 164	07/01/20 16:36	
Tetrachloro-m-xylene	59	10 - 147	07/01/20 16:36	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007029-01
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analysis Method: 8081B **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
4,4'-DDE	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
4,4'-DDT	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
Aldrin	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
Dieldrin	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
Endosulfan I	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
Endosulfan II	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
Endosulfan Sulfate	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
Endrin	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
Endrin Ketone	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
Heptachlor	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
Heptachlor Epoxide	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
Methoxychlor	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
Toxaphene	0.50 U	0.50	0.50	1	07/06/20 10:40	7/1/20	
alpha-BHC	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
alpha-Chlordane	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
beta-BHC	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
delta-BHC	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
gamma-BHC (Lindane)	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
gamma-Chlordane	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	84	10 - 164	07/06/20 10:40	
Tetrachloro-m-xylene	75	10 - 147	07/06/20 10:40	

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/01/20

Sample Matrix: Water

Project:

Duplicate Lab Control Sample Summary Organochlorine Pesticides by Gas Chromatography

Units:ug/L Basis:NA

Service Request: R2005520

Lab Control Sample

Duplicate Lab Control Sample

RQ2006967-04

RQ2006967-05

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
4,4'-DDD	8081B	0.289	0.400	72	0.242	0.400	61	42-159	17	30
4,4'-DDE	8081B	0.287	0.400	72	0.245	0.400	61	47-147	16	30
4,4'-DDT	8081B	0.267	0.400	67	0.224	0.400	56	41-149	18	30
Aldrin	8081B	0.222	0.400	56	0.195	0.400	49	22-137	13	30
Dieldrin	8081B	0.320	0.400	80	0.269	0.400	67	52-144	17	30
Endosulfan I	8081B	0.316	0.400	79	0.268	0.400	67	52-136	16	30
Endosulfan II	8081B	0.323	0.400	81	0.276	0.400	69	57-138	16	30
Endosulfan Sulfate	8081B	0.271	0.400	68	0.229	0.400	57	34-156	17	30
Endrin	8081B	0.310	0.400	78	0.264	0.400	66	56-143	16	30
Endrin Ketone	8081B	0.317	0.400	79	0.269	0.400	67	59-143	16	30
Heptachlor	8081B	0.222	0.400	55	0.196	0.400	49	32-141	12	30
Heptachlor Epoxide	8081B	0.313	0.400	78	0.266	0.400	66	51-143	16	30
Methoxychlor	8081B	0.269	0.400	67	0.240	0.400	60	56-149	11	30
alpha-BHC	8081B	0.301	0.400	75	0.250	0.400	62	36-151	19	30
alpha-Chlordane	8081B	0.303	0.400	76	0.259	0.400	65	50-139	15	30
beta-BHC	8081B	0.326	0.400	82	0.270	0.400	68	55-149	19	30
delta-BHC	8081B	0.287	0.400	72	0.237	0.400	59	29-159	19	30
gamma-BHC (Lindane)	8081B	0.303	0.400	76	0.251	0.400	63	41-149	19	30
gamma-Chlordane	8081B	0.295	0.400	74	0.253	0.400	63	50-140	15	30

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/06/20

Sample Matrix: Water

Duplicate Lab Control Sample Summary Organochlorine Pesticides by Gas Chromatography

Units:ug/L Basis:NA

Service Request: R2005520

Lab Control Sample

Duplicate Lab Control Sample

RQ2007029-02

RQ2007029-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
4,4'-DDD	8081B	0.260	0.400	65	0.268	0.400	67	42-159	3	30
4,4'-DDE	8081B	0.262	0.400	65	0.278	0.400	69	47-147	6	30
4,4'-DDT	8081B	0.251	0.400	63	0.275	0.400	69	41-149	9	30
Aldrin	8081B	0.216	0.400	54	0.219	0.400	55	22-137	1	30
Dieldrin	8081B	0.295	0.400	74	0.307	0.400	77	52-144	4	30
Endosulfan I	8081B	0.289	0.400	72	0.298	0.400	75	52-136	3	30
Endosulfan II	8081B	0.297	0.400	74	0.318	0.400	80	57-138	7	30
Endosulfan Sulfate	8081B	0.247	0.400	62	0.261	0.400	65	34-156	6	30
Endrin	8081B	0.282	0.400	71	0.300	0.400	75	56-143	6	30
Endrin Ketone	8081B	0.279	0.400	70	0.303	0.400	76	59-143	8	30
Heptachlor	8081B	0.190	0.400	47	0.189	0.400	47	32-141	<1	30
Heptachlor Epoxide	8081B	0.286	0.400	71	0.293	0.400	73	51-143	3	30
Methoxychlor	8081B	0.234	0.400	58	0.260	0.400	65	56-149	11	30
alpha-BHC	8081B	0.263	0.400	66	0.258	0.400	65	36-151	2	30
alpha-Chlordane	8081B	0.281	0.400	70	0.290	0.400	72	50-139	3	30
beta-BHC	8081B	0.290	0.400	72	0.293	0.400	73	55-149	1	30
delta-BHC	8081B	0.244	0.400	61	0.249	0.400	62	29-159	2	30
gamma-BHC (Lindane)	8081B	0.263	0.400	66	0.262	0.400	66	41-149	<1	30
gamma-Chlordane	8081B	0.275	0.400	69	0.283	0.400	71	50-140	3	30

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

SURROGATE RECOVERY SUMMARY Polychlorinated Biphenyls (PCBs) by GC

Analysis Method: 8082A **Extraction Method:** EPA 3510C

		Decachlorobiphenyl	Tetrachloro-m-xylene	
Sample Name	Lab Code	10-152	14-129	
WG-9954-062520-SG-013	R2005520-001	53	47	
WG-9954-062520-SG-011	R2005520-002	41	45	
WG-9954-062520-SG-008	R2005520-003	54	57	
WG-9954-062520-SG-009	R2005520-004	30	59	
WG-9954-062520-SG-012	R2005520-005	60	46	
RB-9954-062520-SG-001	R2005520-006	10	47	_
WG-9954-062520-SG-010	R2005520-008	28	39	
Method Blank	RQ2006967-03	56	51	
Method Blank	RQ2007029-01	89	70	
Lab Control Sample	RQ2006967-04	56	49	
Duplicate Lab Control Sample	RQ2006967-05	58	49	_
Lab Control Sample	RQ2007029-02	71	53	
Duplicate Lab Control Sample	RQ2007029-03	68	49	
WG-9954-062520-SG-012 MS	RQ2006967-01	54	36	
WG-9954-062520-SG-012 DMS	RQ2006967-02	84	50	
RB-9954-062520-SG-001 MS	RQ2007029-04	7*	34	
RB-9954-062520-SG-001 DMS	RQ2007029-05	9*	39	

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates) **Service Request:** R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring 06/25/20

Sample Matrix:

06/26/20

Water **Date Received:** Date Analyzed:

07/8/20

Date Extracted:

Date Collected:

06/30/20

Duplicate Matrix Spike Summary Polychlorinated Biphenyls (PCBs) by GC

Sample Name: WG-9954-062520-SG-012 **Units:**

R2005520-005

Basis:

NA

ug/L

Analysis Method: 8082A **Prep Method:** EPA 3510C

Lab Code:

Duplicate Matrix Spike

Matrix Spike RQ2006967-01

RQ2006967-02

	Sample		Spike			Spike		% Rec		RPD
Analyte Name	Result	Result	Amount	% Rec	Result	Amount	% Rec	Limits	RPD	Limit
Aroclor 1016	0.91 U	2.23	3.64	61	2.49	3.64	68	32-142	11	30
Aroclor 1260	0.91 U	2.24	3.64	62	2.96	3.64	81	28-142	28	30

Results flagged with an asterisk (*) indicate values outside control criteria.

Results flagged with a pound (#) indicate the control criteria is not applicable.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

Matrix Spike and Matrix Spike Duplicate Data is presented for information purposes only. The matrix may or may not be relevant to samples reported in this report. The laboratory evaluates system performance based on the LCS and LCSD control limits.

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates) **Service Request:** R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Date Collected: 06/25/20

Sample Matrix: Water **Date Received:**

06/26/20

Date Analyzed:

07/6/20

Date Extracted:

07/1/20

Duplicate Matrix Spike Summary Polychlorinated Biphenyls (PCBs) by GC

Sample Name:

RB-9954-062520-SG-001

Units:

ug/L

Lab Code:

R2005520-006

Basis:

NA

Analysis Method: Prep Method:

8082A

EPA 3510C

Matrix Spike RQ2007029-04 **Duplicate Matrix Spike**

RQ2007029-05

	Sample		Spike			Spike		% Rec		RPD
Analyte Name	Result	Result	Amount	% Rec	Result	Amount	% Rec	Limits	RPD	Limit
Aroclor 1016	0.91 U	1.73	3.64	48	2.26	3.64	62	32-142	26	30
Aroclor 1260	0.91 U	1.14	3.64	31	1.64	3.64	45	28-142	36*	30

Results flagged with an asterisk (*) indicate values outside control criteria.

Results flagged with a pound (#) indicate the control criteria is not applicable.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

Matrix Spike and Matrix Spike Duplicate Data is presented for information purposes only. The matrix may or may not be relevant to samples reported in this report. The laboratory evaluates system performance based on the LCS and LCSD control limits.

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2006967-03Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analysis Method: 8082A **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	1.0 U	1.0	0.50	1	07/01/20 17:28	6/30/20	
Aroclor 1221	2.0 U	2.0	1.0	1	07/01/20 17:28	6/30/20	
Aroclor 1232	1.0 U	1.0	0.50	1	07/01/20 17:28	6/30/20	
Aroclor 1242	1.0 U	1.0	0.50	1	07/01/20 17:28	6/30/20	
Aroclor 1248	1.0 U	1.0	0.50	1	07/01/20 17:28	6/30/20	
Aroclor 1254	1.0 U	1.0	0.50	1	07/01/20 17:28	6/30/20	
Aroclor 1260	1.0 U	1.0	0.50	1	07/01/20 17:28	6/30/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	56	10 - 152	07/01/20 17:28	
Tetrachloro-m-xylene	51	14 - 129	07/01/20 17:28	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005520

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007029-01Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analysis Method: 8082A **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	1.0 U	1.0	0.50	1	07/06/20 11:40	7/1/20	
Aroclor 1221	2.0 U	2.0	1.0	1	07/06/20 11:40	7/1/20	
Aroclor 1232	1.0 U	1.0	0.50	1	07/06/20 11:40	7/1/20	
Aroclor 1242	1.0 U	1.0	0.50	1	07/06/20 11:40	7/1/20	
Aroclor 1248	1.0 U	1.0	0.50	1	07/06/20 11:40	7/1/20	
Aroclor 1254	1.0 U	1.0	0.50	1	07/06/20 11:40	7/1/20	
Aroclor 1260	1.0 U	1.0	0.50	1	07/06/20 11:40	7/1/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	89	10 - 152	07/06/20 11:40	
Tetrachloro-m-xylene	70	14 - 129	07/06/20 11:40	

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix:

Project:

Water

Duplicate Lab Control Sample Summary Polychlorinated Biphenyls (PCBs) by GC

Units:ug/L Basis:NA

Service Request: R2005520

Date Analyzed: 07/01/20

Lab Control Sample

Duplicate Lab Control Sample

RQ2006967-04

RQ2006967-05

	Analytical		Spike			Spike		% Rec		RPD
Analyte Name	Method	Result	Amount	% Rec	Result	Amount	% Rec	Limits	RPD	Limit
Aroclor 1016	8082A	2.57	4.00	64	2.65	4.00	66	49-123	3	30
Aroclor 1260	8082A	2.77	4.00	69	2.87	4.00	72	30-120	4	30

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix: Water

Project:

Duplicate Lab Control Sample Summary Polychlorinated Biphenyls (PCBs) by GC

> Units:ug/L Basis:NA

Service Request: R2005520

Date Analyzed: 07/06/20

Lab Control Sample

Duplicate Lab Control Sample

RQ2007029-02

RQ2007029-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
Aroclor 1016	8082A	2.88	4.00	72	2.96	4.00	74	49-123	2	30
Aroclor 1260	8082A	3.52	4.00	88	3.35	4.00	84	30-120	5	30



Service Request No:R2005539

Ms. Kathy Willy GHD Services Inc. 2055 Niagara Falls Blvd., Niagara Falls, NY 14304

Laboratory Results for: Love Canal:292-402-D02-3100

Dear Ms. Willy,

Enclosed are the results of the sample(s) submitted to our laboratory June 27, 2020 For your reference, these analyses have been assigned our service request number **R2005539**.

All testing was performed according to our laboratory's quality assurance program and met the requirements of the TNI standards except as noted in the case narrative report. Any testing not included in the lab's accreditation is identified on a Non-Certified Analytes report. All results are intended to be considered in their entirety. ALS Environmental is not responsible for use of less than the complete report. Results apply only to the individual samples submitted to the lab for analysis, as listed in the report. The measurement uncertainty of the results included in this report is within that expected when using the prescribed method(s), and represented by Laboratory Control Sample control limits. Any events, such as QC failures or Holding Time exceedances, which may add to the uncertainty are explained in the report narrative or are flagged with qualifiers. The flags are explained in the Report Qualifiers and Definitions page of this report.

Please contact me if you have any questions. My extension is 7471. You may also contact me via email at Brady.Kalkman@alsglobal.com.

Respectfully submitted,

Goody Kullen

ALS Group USA, Corp. dba ALS Environmental

Brady Kalkman Project Manager



Narrative Documents

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com



Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100 Date Received: 06/27/2020

Sample Matrix: Water

CASE NARRATIVE

All analyses were performed consistent with the quality assurance program of ALS Environmental. This report contains analytical results for samples for the Tier level IV requested by the client.

Sample Receipt:

Six water samples were received for analysis at ALS Environmental on 06/27/2020. Any discrepancies upon initial sample inspection are annotated on the sample receipt and preservation form included within this report. The samples were stored at minimum in accordance with the analytical method requirements.

Semivolatiles by GC/MS:

Method 8270D, 07/06/2020: The upper control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). The field samples analyzed in this sequence did not contain the analyte(s) in question above the Method Reporting Limit (MRL). Since the exceedance equates to a potential high bias, the data quality was not significantly affected and no further corrective action was taken.

Method 8270D, R2005539-005: The control limits were exceeded for one or more surrogates. A reanalysis was not performed because insufficient sample was available. No further corrective action was possible.

Semivoa GC:

Method 8081B, 07/07/2020: The upper control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). The field samples analyzed in this sequence did not contain the analyte(s) in question above the Method Reporting Limit (MRL). Since the exceedance equates to a potential high bias, the data quality was not significantly affected and no further corrective action was taken.

Method 8081B, 07/07/2020: The lower control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). Since there were no detections of the analyte(s) above the MRL in the associated field samples, the quantitation is not affected. The data quality was not significantly affected and no further corrective action was taken.

Method 8081B, 07/06/2020: The lower control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). Since there were no detections of the analyte(s) above the MRL in the associated field samples, the quantitation is not affected. The data quality was not significantly affected and no further corrective action was taken. **Volatiles by GC/MS:**

Method 8260C, 07/01/2020: The upper control criterion was exceeded for one or more analytes in the Laboratory Control Sample (LCS). There were no detections of the analyte(s) above the MRL in the associated field samples. The error associated with elevated recovery equates to a high bias. The sample data is not significantly affected. No further corrective action was appropriate.

Method 8260C, 07/01/2020: The lower control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). Since there were no detections of the analyte(s) above the MRL in the associated field samples, the quantitation is not affected. The data quality was not significantly affected and no further corrective action was taken.

Method 8260C, 07/07/2020: The upper control criterion was exceeded for one or more analytes in the Laboratory Control Sample (LCS). There were no detections of the analyte(s) above the MRL in the associated field samples. The error associated with elevated recovery equates to a high bias. The sample data is not significantly affected. No further corrective action was appropriate.



SAMPLE DETECTION SUMMARY

CLIENT ID: WG-9954-062620-RM-014		Lab	ID: R2005	5539-001		
Analyte	Results	Flag	MDL	MRL	Units	Method
4-Methyl-2-pentanone	2.4	J	0.20	10	ug/L	8260C
Carbon Disulfide	4.0	J	0.42	10	ug/L	8260C
Chloromethane	0.30	J	0.28	5.0	ug/L	8260C
Bis(2-ethylhexyl) Phthalate	2.2	J	0.91	9.1	ug/L	8270D
alpha-BHC	0.088		0.019	0.045	ug/L	8081B
beta-BHC	0.036	J	0.019	0.045	ug/L	8081B
delta-BHC	0.32		0.019	0.045	ug/L	8081B
gamma-BHC (Lindane)	0.11		0.019	0.045	ug/L	8081B
CLIENT ID: WG-9954-062620-RM-015		Lab	ID: R2005	5539-002		
Analyte	Results	Flag	MDL	MRL	Units	Method
Carbon Disulfide	2.7	J	0.42	10	ug/L	8260C
Chloromethane	0.35	J	0.28	5.0	ug/L	8260C
alpha-BHC	0.064		0.019	0.045	ug/L	8081B
delta-BHC	0.042	J	0.019	0.045	ug/L	8081B
gamma-BHC (Lindane)	0.060		0.019	0.045	ug/L	8081B
CLIENT ID: WG-9954-062620-RM-016		Lab	ID: R2005	5539-003		
Analyte	Results	Flag	MDL	MRL	Units	Method
Carbon Disulfide	1.9	J	0.42	10	ug/L	8260C
alpha-BHC	0.097		0.019	0.045	ug/L	8081B
delta-BHC	0.10		0.019	0.045	ug/L	8081B
gamma-BHC (Lindane)	0.096		0.019	0.045	ug/L	8081B
CLIENT ID: WG-9954-062620-RM-017		Lab	ID: R2005	5539-004		
Analyte	Results	Flag	MDL	MRL	Units	Method
Chloromethane	0.38	J	0.28	5.0	ug/L	8260C
CLIENT ID: WG-9954-062620-RM-018		Lab	ID: R2005	5539-005		
Analyte	Results	Flag	MDL	MRL	Units	Method
Carbon Disulfide	1.2	J	0.42	10	ug/L	8260C
4,4'-DDE	0.029	J	0.019	0.045	ug/L	8081B
Endosulfan I	0.037	J	0.019	0.045	ug/L	8081B
Endosulian i	0.00.					
beta-BHC	0.045	JP	0.019	0.045	ug/L	8081B
		JP J	0.019 0.019	0.045 0.045	ug/L ug/L	8081B 8081B



Sample Receipt Information

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request:R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

SAMPLE CROSS-REFERENCE

SAMPLE #	CLIENT SAMPLE ID	<u>DATE</u>	<u>TIME</u>
R2005539-001	WG-9954-062620-RM-014	6/26/2020	0850
R2005539-002	WG-9954-062620-RM-015	6/26/2020	1000
R2005539-003	WG-9954-062620-RM-016	6/26/2020	1105
R2005539-004	WG-9954-062620-RM-017	6/26/2020	1155
R2005539-005	WG-9954-062620-RM-018	6/26/2020	1240
R2005539-006	TB-9954-062620-RM-003	6/26/2020	0000



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	(Containers for each sample may be combined on one PRESERVATION - (S					L.,	HC										 			<u> </u>			•
1	WG 9954 - 062620-RH	014 6/26/2	0 0850) WG	9	N	X	×	K	X							 	7	T				•
2	WG-9954-062620 RM-0								X.	X					7			7	T		· · · · · · · · · · · · · · · · · · ·		•
3	WG-9954-062620-RM-C	16 6/26/2	0 1105	WG	G	N	X	Х	×	X								7	T				•
4	WG-9954-0626-20-RM-0	17 6/26/2	0 1155	W4	G	N	X	X	X	X								7	T	-			•
5	WG-9954-062620-RM-C	18 6 26/2	0 124	o W6	6	N	Х	Х	X	X								7			2(51)		•
6	TB-9954-062620-RM-0			TB											\neg			3			-6		•
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	T Required in business days (use separate 1 Day □ 2 Days □ 3 Days □ 1 Week ↓		•			No	tes/ S	Speci	iai Re	equin	emei	nts:				_	_						
	RELINQUISHED BY	COMPANY		DATE		TIM	E				_	F	RECEN	VED B	<u>Y</u>				C	OMPANY	DATE	TIME	
1.	Duc ayıan	GHD	61	26/2	0	14	10	5	1			1				K	2	A	L	<i>-</i>	6-27-2020	08128	-
د. ع.	V								3.				<u></u>		<u>.</u>					R2005	5539	 5	-

THE CHAIN OF CUSTODY IS A LEGAL DOCUMENT - ALL FIELDS MUST BE COMPLETED ACCURATELY

Distribution: WHITE - Fully Executed Copy (CRA) YELLOW - Receiving Laboratory Copy

PINK - Shipper

GOLDENROD - Samplin

GHD Services inc. Love Canal:292-402-002-3100

CHAIN OF CUSTODY / LABORATORY ANALYSIS REQUEST FORM SR# 004, 005, 006, 007, 008, 009, 010, 011, 012, 013 1565 Jefferson Road, Bidg 300, Suite 380, Rochester, NY 14623 T030477 Phone (585) 288-5380 / FAX (585) 288-8475 S) Environmental www.alsglobal.com 14D Love Canal:292-402-D02-3100 Project Number: 9954 Annual Long Term Monitoring Report To Kathy Willy NUMBER OF CONTAINERS Company / Address GHD Services Inc. 2055 Niagara Falls Blvd., Suite 3 Niagara Falls NY, 14304 3081B / Pest OC svo 260C / VOC 082A / PCB Phone # 716-297-2160 Sampler Signature 716-297-2265 Sampler Printed Name 3270D/ Remarks SAMPLING Matrix LABID Date Time CLIENT SAMPLE ID Liquid Llquid Liquid ∐quid Liquid ∐auld Liquid Llquid ∐guid 10. Liquid Turnaround Requirements Report Requirements Invoice Information Special Instructions/Comments: RUSH (SURCHARGES APPLY) I. Results Only _II. Results + QC Summaries (LCS, DUP, MS/MSD as required) P.O.# Standard (3 weeks) ItI. Results + OC and Cilibration Bill To: Summaries X IV. Data Validation Report REQUESTED FAX DATE with Raw Data EData ____Yes ____No Requested Report Date Received By: Relinquished By: Received By: Relinquished By: Received By: Relinguished By: Signature Signature Signature Signature Signature Signature Printed Name Printed Name Printed Name Printed Name Printed Name Printed Name Firm Firm Firm Firm Firm Firm Date/Time Date/Time Date/Time Date/Time Date/Time Date/Time

(ALS)

Cooler Receipt and Preservation Check Form

\mathcal{L}	R2005539	5
]	GHD Services Inc. Love Canal:292-402-D02-3100	

Project/Clie		ر ــــــــــــــــــــــــــــــــــــ		1	_Fold	ler Number			· · · · · ·				
Cooler receive	ed on 627	<u>2020</u>	by:	<u>U</u>	· 201	COURIE	R: ALS	UPS	FEDEX	VEL	OCITY C	LIENT	
1 Were Cu	stody seals on	outside of coole	r?		Y/N	5a Per	chlorate :	samples	have requ	uired hea	adspace?	Y	N (NA
2 Custody	papers proper	ly completed (in	k, sign	ed)?	Y) N	5b Did	VOA via	ls, Alk,	or Sulfide	have sig	g* bubbles	? Y	N NA
3 Did all b	ottles arrive in	good condition (unbrok	(en)?	Ý) N	6 Wh	ere did th	e bottle	s originate	? (ALS/RO	CLI	ENT
4 Circle:	Wet Ice Dry	Ice Gel packs	pres	ent?	Y)N	7 Soil	VOA rec	eived a	s: Bu	lk Er	icore 5	035set (NA NA
3. Temperatur	e Readings	Date: 6 2)-2	1000	Time:	08:5	ם ל	D: IR#7	1R#10	\supset	From:	Temp Bla	ank San	iple Bottle
Observed To	emp (°C)	1.4									-		
Within 0-6%	C?	(Y) N		Y	N	Y N	Y	N	Y	N	Y N	Y	N
If <0°C, wer	e samples froz	en? Y N		Y	N	Y N	Y	N	Y	N	Y N	Y	N .
If out of 7	Temperature,	note packing/ic	e cond	ition:		Ice m	elted I	Poorly I	acked (de	scribed	below)	Same l	Day Rule
&Client A	Approval to R	un Samples:		_ Stan	ding Ap	proval Clie	ent aware	at drop	-off Cli	ent noti	fied by:		
All samples	held in storag	e location:	R-OV	<u> </u>	y K	on Go	720 at (19100					
-	_	orage location:			у	on	at		within 48	hours o	of samplin	g? Y	N
Cooler Breakdown/Preservation Check**: Date: 6/29/2020 Time: /2/8 by: 6 9. Were all bottle labels complete (i.e. analysis, preservation, etc.)? YES NO 10. Did all bottle labels and tags agree with custody papers? YES NO 11. Were correct containers used for the tests indicated? YES NO 12. Were 5035 vials acceptable (no extra labels, not leaking)? YES NO 13. Air Samples: Cassettes / Tubes Intact with MS? Canisters Pressurized Tedlar® Bags Inflated N/8													
pН	Lot of test paper	Reagent	Preser Yes	No	Lock	eceived	Exp	Adju	ole ID sted	Vol. Added	Lot A	daca	Final pH
≥12		NaOH											
≤2		HNO ₃											
≤2		H₂SO₄											
<4		NaHSO ₄								·			
5-9		For 608pest				otify for 3day							
Residual		For CN,		1		ontact PM to add	i				ľ		
Chlorine		Phenol, 625,				D ₃ (625, 608, scorbic (phenol	.						
(-)		608pest, 522			Civy, a.	scorbic (pilenoi	,.						-
		Na ₂ S ₂ O ₃	ļ					*****	1166	137.44.1			
	•	ZnAcetate HCl	**	**		_		Other		tles of all	e tested befor samples with	•	eservatives
Bottle lot Explain a	numbers: II Discrepanci	05//20-/18 es/Other Comm UNK: All	smc lents:	clu	is a	weed		T are ell	vonou (mor Ji	un roprosi			
15	ITI VI	1116 XIII	ノ'	•									

HPROD	BULK
HTR	FLDT
SUB	HGFB
ALS	LL3541

Labels secondary reviewed by:_	(Q)
PC Secondary Review:	



Miscellaneous Forms

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com



REPORT QUALIFIERS AND DEFINITIONS

- U Analyte was analyzed for but not detected. The sample quantitation limit has been corrected for dilution and for percent moisture, unless otherwise noted in the case narrative.
- J Estimated value due to either being a Tentatively Identified Compound (TIC) or that the concentration is between the MRL and the MDL. Concentrations are not verified within the linear range of the calibration. For DoD: concentration >40% difference between two GC columns (pesticides/Arclors).
- B Analyte was also detected in the associated method blank at a concentration that may have contributed to the sample result.
- E Inorganics- Concentration is estimated due to the serial dilution was outside control limits.
- E Organics- Concentration has exceeded the calibration range for that specific analysis.
- D Concentration is a result of a dilution, typically a secondary analysis of the sample due to exceeding the calibration range or that a surrogate has been diluted out of the sample and cannot be assessed.
- * Indicates that a quality control parameter has exceeded laboratory limits. Under the õNotesö column of the Form I, this qualifier denotes analysis was performed out of Holding Time.
- H Analysis was performed out of hold time for tests that have an õimmediateö hold time criteria.
- # Spike was diluted out.

- + Correlation coefficient for MSA is <0.995.
- N Inorganics- Matrix spike recovery was outside laboratory limits.
- N Organics- Presumptive evidence of a compound (reported as a TIC) based on the MS library search.
- S Concentration has been determined using Method of Standard Additions (MSA).
- W Post-Digestion Spike recovery is outside control limits and the sample absorbance is <50% of the spike absorbance.
- P Concentration >40% difference between the two GC columns.
- C Confirmed by GC/MS
- Q DoD reports: indicates a pesticide/Aroclor is not confirmed (×100% Difference between two GC columns).
- X See Case Narrative for discussion.
- MRL Method Reporting Limit. Also known as:
- LOQ Limit of Quantitation (LOQ)

 The lowest concentration at which the method analyte may be reliably quantified under the method conditions.
- MDL Method Detection Limit. A statistical value derived from a study designed to provide the lowest concentration that will be detected 99% of the time. Values between the MDL and MRL are estimated (see J qualifier).
- LOD Limit of Detection. A value at or above the MDL which has been verified to be detectable.
- ND Non-Detect. Analyte was not detected at the concentration listed. Same as U qualifier.



Rochester Lab ID # for State Certifications¹

Connecticut ID # PH0556	Maine ID #NY0032	Pennsylvania ID# 68-786
Delaware Approved	New Hampshire ID # 2941	Rhode Island ID # 158
DoD ELAP #65817	New York ID # 10145	Virginia #460167
Florida ID # E87674	North Carolina #676	

¹ Analyses were performed according to our laboratory

NELAP-approved quality assurance program and any applicable state or agency requirements. The test results meet requirements of the current NELAP/TNI standards or state or agency requirements, where applicable, except as noted in the case narrative. Since not all analyte/method/matrix combinations are offered for state/NELAC accreditation, this report may contain results which are not accredited. For a specific list of accredited analytes, contact the laboratory or go to https://www.alsglobal.com/locations/americas/north-america/usa/new-york/rochester-environmental

ALS Laboratory Group

Acronyms

ASTM American Society for Testing and Materials

A2LA American Association for Laboratory Accreditation

CARB California Air Resources Board

CAS Number Chemical Abstract Service registry Number

CFC Chlorofluorocarbon CFU Colony-Forming Unit

DEC Department of Environmental Conservation

DEQ Department of Environmental Quality

DHS Department of Health Services

DOE Department of Ecology DOH Department of Health

EPA U. S. Environmental Protection Agency

ELAP Environmental Laboratory Accreditation Program

GC Gas Chromatography

GC/MS Gas Chromatography/Mass Spectrometry

LUFT Leaking Underground Fuel Tank

M Modified

MCL Maximum Contaminant Level is the highest permissible concentration of a

substance allowed in drinking water as established by the USEPA.

MDL Method Detection Limit
MPN Most Probable Number
MRL Method Reporting Limit

NA Not Applicable NC Not Calculated

NCASI National Council of the Paper Industry for Air and Stream Improvement

ND Not Detected

NIOSH National Institute for Occupational Safety and Health

PQL Practical Quantitation Limit

RCRA Resource Conservation and Recovery Act

SIM Selected Ion Monitoring

TPH Total Petroleum Hydrocarbons

tr Trace level is the concentration of an analyte that is less than the PQL but

greater than or equal to the MDL.

Analyst Summary report

Service Request: R2005539

Date Collected: 06/26/20

Date Received: 06/27/20

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Monitoring

Sample Name: WG-9954-062620-RM-014

Lab Code: R2005539-001

Sample Matrix: Water

Analysis Method Extracted/Digested By Analyzed By

8081B KSERCU BALLGEIER 8082A KSERCU BALLGEIER 8260C FNAEGLER

8270D KSERCU JMISIUREWICZ

Sample Name: WG-9954-062620-RM-015 Date Collected: 06/26/20

Lab Code: R2005539-002 **Date Received:** 06/27/20

Sample Matrix: Water

Analysis Method Extracted/Digested By Analyzed By

8081B KSERCU JMISIUREWICZ 8082A KSERCU BALLGEIER 8260C FNAEGLER 8270D KSERCU JMISIUREWICZ

 Sample Name:
 WG-9954-062620-RM-016
 Date Collected: 06/26/20

 Lab Code:
 R2005539-003
 Date Received: 06/27/20

Lab Code:R2005539-003Date Received: 06/27/20Sample Matrix:Water

Analysis Method Extracted/Digested By Analyzed By

8081BKSERCUJMISIUREWICZ8082AKSERCUBALLGEIER8260CFNAEGLER8270DKSERCUJMISIUREWICZ

Sample Name: WG-9954-062620-RM-017 **Date Collected:** 06/26/20

Lab Code: R2005539-004 **Date Received:** 06/27/20

Sample Matrix: Water

Analysis Method Extracted/Digested By Analyzed By

8081B KSERCU JMISIUREWICZ

Printed 7/15/2020 12:37:46 PM Superset Reference:20-0000554991 rev 00 Page 13 of 90

Analyst Summary report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term **Project:**

Monitoring

Sample Name: WG-9954-062620-RM-017

Lab Code: R2005539-004

Sample Matrix: Water

Analyzed By Analysis Method Extracted/Digested By

KSERCU 8082A **BALLGEIER**

8260C **FNAEGLER**

8270D **KSERCU JMISIUREWICZ**

Sample Name: WG-9954-062620-RM-018 **Date Collected:** 06/26/20

Lab Code: R2005539-005 **Date Received:** 06/27/20

Sample Matrix: Water

Analyzed By Analysis Method Extracted/Digested By

8081B **KSERCU JMISIUREWICZ** 8082A **KSERCU BALLGEIER**

8260C **FNAEGLER**

8270D **KSERCU JMISIUREWICZ**

TB-9954-062620-RM-003

Sample Name: Date Collected: 06/26/20 Lab Code: R2005539-006 **Date Received:** 06/27/20

Sample Matrix: Water

Analyzed By **Analysis Method Extracted/Digested By**

8260C **FNAEGLER**

Service Request: R2005539

Date Collected: 06/26/20

Date Received: 06/27/20



INORGANIC PREPARATION METHODS

The preparation methods associated with this report are found in these tables unless discussed in the case narrative.

Water/Liquid Matrix

Analytical Method	Preparation Method
200.7	200.2
200.8	200.2
6010C	3005A/3010A
6020A	ILM05.3
9034 Sulfide Acid Soluble	9030B
SM 4500-CN-E Residual Cyanide	SM 4500-CN-G
SM 4500-CN-E WAD Cyanide	SM 4500-CN-I

Solid/Soil/Non-Aqueous Matrix

Analytical Method	Preparation				
	Method				
6010C	3050B				
6020A	3050B				
6010C TCLP (1311)	3005A/3010A				
extract					
6010 SPLP (1312) extract	3005A/3010A				
7199	3060A				
300.0 Anions/ 350.1/	DI extraction				
353.2/ SM 2320B/ SM					
5210B/ 9056A Anions					
method is the same as the analytical method					
5210B/ 9056A Anions For analytical methods not listed, the preparation					



Sample Results

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475

www.alsglobal.com



Volatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 08:50

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-014
 Units: ug/L

 Lab Code:
 R2005539-001
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/01/20 03:34	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/01/20 03:34	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/01/20 03:34	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/01/20 03:34	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/01/20 03:34	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/01/20 03:34	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/01/20 03:34	
2-Butanone (MEK)	10 U	10	0.78	1	07/01/20 03:34	
2-Hexanone	10 U	10	0.20	1	07/01/20 03:34	
4-Methyl-2-pentanone	2.4 J	10	0.20	1	07/01/20 03:34	
Acetone	10 U	10	5.0	1	07/01/20 03:34	
Benzene	5.0 U	5.0	0.20	1	07/01/20 03:34	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/01/20 03:34	
Bromoform	5.0 U	5.0	0.25	1	07/01/20 03:34	
Bromomethane	5.0 U	5.0	0.70	1	07/01/20 03:34	
Carbon Disulfide	4.0 J	10	0.42	1	07/01/20 03:34	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/01/20 03:34	
Chlorobenzene	5.0 U	5.0	0.20	1	07/01/20 03:34	
Chloroethane	5.0 U	5.0	0.23	1	07/01/20 03:34	
Chloroform	5.0 U	5.0	0.24	1	07/01/20 03:34	
Chloromethane	0.30 J	5.0	0.28	1	07/01/20 03:34	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/01/20 03:34	
Dichloromethane	5.0 U	5.0	0.65	1	07/01/20 03:34	
Ethylbenzene	5.0 U	5.0	0.20	1	07/01/20 03:34	
Styrene	5.0 U	5.0	0.20	1	07/01/20 03:34	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/01/20 03:34	
Toluene	5.0 U	5.0	0.20	1	07/01/20 03:34	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/01/20 03:34	
Vinyl Acetate	10 U	10	1.1	1	07/01/20 03:34	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/01/20 03:34	
Xylenes, Total	5.0 U	5.0	0.23	1	07/01/20 03:34	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/01/20 03:34	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/01/20 03:34	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/01/20 03:34	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/01/20 03:34	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 08:50

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-014
 Units: ug/L

 Lab Code:
 R2005539-001
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	93	85 - 122	07/01/20 03:34	
Dibromofluoromethane	94	89 - 119	07/01/20 03:34	
Toluene-d8	94	87 - 121	07/01/20 03:34	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 08:50

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-014
 Units: ug/L

 Lab Code:
 R2005539-001
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

			Result	
CAS#	Compound Identification	RT	ug/L	Q
	unknown	1.32	8.5	J
000110-93-0	5-Hepten-2-one, 6-methyl-	11.53	10.2	JN

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 10:00

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-015
 Units: ug/L

 Lab Code:
 R2005539-002
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/01/20 03:56	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/01/20 03:56	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/01/20 03:56	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/01/20 03:56	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/01/20 03:56	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/01/20 03:56	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/01/20 03:56	
2-Butanone (MEK)	10 U	10	0.78	1	07/01/20 03:56	
2-Hexanone	10 U	10	0.20	1	07/01/20 03:56	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/01/20 03:56	
Acetone	10 U	10	5.0	1	07/01/20 03:56	
Benzene	5.0 U	5.0	0.20	1	07/01/20 03:56	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/01/20 03:56	
Bromoform	5.0 U	5.0	0.25	1	07/01/20 03:56	
Bromomethane	5.0 U	5.0	0.70	1	07/01/20 03:56	
Carbon Disulfide	2.7 J	10	0.42	1	07/01/20 03:56	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/01/20 03:56	
Chlorobenzene	5.0 U	5.0	0.20	1	07/01/20 03:56	
Chloroethane	5.0 U	5.0	0.23	1	07/01/20 03:56	
Chloroform	5.0 U	5.0	0.24	1	07/01/20 03:56	
Chloromethane	0.35 J	5.0	0.28	1	07/01/20 03:56	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/01/20 03:56	
Dichloromethane	5.0 U	5.0	0.65	1	07/01/20 03:56	
Ethylbenzene	5.0 U	5.0	0.20	1	07/01/20 03:56	
Styrene	5.0 U	5.0	0.20	1	07/01/20 03:56	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/01/20 03:56	
Toluene	5.0 U	5.0	0.20	1	07/01/20 03:56	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/01/20 03:56	
Vinyl Acetate	10 U	10	1.1	1	07/01/20 03:56	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/01/20 03:56	
Xylenes, Total	5.0 U	5.0	0.23	1	07/01/20 03:56	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/01/20 03:56	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/01/20 03:56	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/01/20 03:56	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/01/20 03:56	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 10:00

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-015
 Units: ug/L

 Lab Code:
 R2005539-002
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	98	85 - 122	07/01/20 03:56	
Dibromofluoromethane	100	89 - 119	07/01/20 03:56	
Toluene-d8	98	87 - 121	07/01/20 03:56	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 10:00

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-015
 Units: ug/L

 Lab Code:
 R2005539-002
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Tentatively Identified Compounds

			Result	
CAS#	Compound Identification	RT	ug/L	Q
000420-56-4	Silane, fluorotrimethyl-	1.32	10.2	JN

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 11:05

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-016
 Units: ug/L

 Lab Code:
 R2005539-003
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/07/20 21:25	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/07/20 21:25	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/07/20 21:25	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/07/20 21:25	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/07/20 21:25	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/07/20 21:25	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/07/20 21:25	
2-Butanone (MEK)	10 U	10	0.78	1	07/07/20 21:25	
2-Hexanone	10 U	10	0.20	1	07/07/20 21:25	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/07/20 21:25	
Acetone	10 U	10	5.0	1	07/07/20 21:25	
Benzene	5.0 U	5.0	0.20	1	07/07/20 21:25	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/07/20 21:25	
Bromoform	5.0 U	5.0	0.25	1	07/07/20 21:25	
Bromomethane	5.0 U	5.0	0.70	1	07/07/20 21:25	
Carbon Disulfide	1.9 J	10	0.42	1	07/07/20 21:25	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/07/20 21:25	
Chlorobenzene	5.0 U	5.0	0.20	1	07/07/20 21:25	
Chloroethane	5.0 U	5.0	0.23	1	07/07/20 21:25	
Chloroform	5.0 U	5.0	0.24	1	07/07/20 21:25	
Chloromethane	5.0 U	5.0	0.28	1	07/07/20 21:25	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/07/20 21:25	
Dichloromethane	5.0 U	5.0	0.65	1	07/07/20 21:25	
Ethylbenzene	5.0 U	5.0	0.20	1	07/07/20 21:25	
Styrene	5.0 U	5.0	0.20	1	07/07/20 21:25	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/07/20 21:25	
Toluene	5.0 U	5.0	0.20	1	07/07/20 21:25	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/07/20 21:25	
Vinyl Acetate	10 U	10	1.1	1	07/07/20 21:25	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/07/20 21:25	
Xylenes, Total	5.0 U	5.0	0.23	1	07/07/20 21:25	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/07/20 21:25	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/07/20 21:25	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/07/20 21:25	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/07/20 21:25	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 11:05

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-016
 Units: ug/L

 Lab Code:
 R2005539-003
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	105	85 - 122	07/07/20 21:25	
Dibromofluoromethane	104	89 - 119	07/07/20 21:25	
Toluene-d8	104	87 - 121	07/07/20 21:25	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 11:05

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-016
 Units: ug/L

 Lab Code:
 R2005539-003
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Tentatively Identified Compounds

			Result	
CAS#	Compound Identification	RT	ug/L	Q
007446-09-5	Sulfur dioxide	1.26	71.1	JN

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539 **Date Collected:** 06/26/20 11:55

Love Canal:292-402-D02-3100/9954 Annual Long Term **Project:**

Monitoring

Sample Matrix: Date Received: 06/27/20 08:50 Water

Sample Name: WG-9954-062620-RM-017 Units: ug/L Basis: NA Lab Code: R2005539-004

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/01/20 04:40	_
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/01/20 04:40	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/01/20 04:40	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/01/20 04:40	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/01/20 04:40	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/01/20 04:40	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/01/20 04:40	
2-Butanone (MEK)	10 U	10	0.78	1	07/01/20 04:40	
2-Hexanone	10 U	10	0.20	1	07/01/20 04:40	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/01/20 04:40	
Acetone	10 U	10	5.0	1	07/01/20 04:40	
Benzene	5.0 U	5.0	0.20	1	07/01/20 04:40	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/01/20 04:40	
Bromoform	5.0 U	5.0	0.25	1	07/01/20 04:40	
Bromomethane	5.0 U	5.0	0.70	1	07/01/20 04:40	
Carbon Disulfide	10 U	10	0.42	1	07/01/20 04:40	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/01/20 04:40	
Chlorobenzene	5.0 U	5.0	0.20	1	07/01/20 04:40	
Chloroethane	5.0 U	5.0	0.23	1	07/01/20 04:40	
Chloroform	5.0 U	5.0	0.24	1	07/01/20 04:40	
Chloromethane	0.38 J	5.0	0.28	1	07/01/20 04:40	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/01/20 04:40	
Dichloromethane	5.0 U	5.0	0.65	1	07/01/20 04:40	
Ethylbenzene	5.0 U	5.0	0.20	1	07/01/20 04:40	
Styrene	5.0 U	5.0	0.20	1	07/01/20 04:40	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/01/20 04:40	
Toluene	5.0 U	5.0	0.20	1	07/01/20 04:40	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/01/20 04:40	
Vinyl Acetate	10 U	10	1.1	1	07/01/20 04:40	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/01/20 04:40	
Xylenes, Total	5.0 U	5.0	0.23	1	07/01/20 04:40	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/01/20 04:40	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/01/20 04:40	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/01/20 04:40	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/01/20 04:40	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 11:55

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-017
 Units: ug/L

 Lab Code:
 R2005539-004
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	92	85 - 122	07/01/20 04:40	
Dibromofluoromethane	93	89 - 119	07/01/20 04:40	
Toluene-d8	94	87 - 121	07/01/20 04:40	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 11:55

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-017
 Units: ug/L

 Lab Code:
 R2005539-004
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

Result

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 12:40

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-018
 Units: ug/L

 Lab Code:
 R2005539-005
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/07/20 21:46	_
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/07/20 21:46	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/07/20 21:46	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/07/20 21:46	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/07/20 21:46	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/07/20 21:46	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/07/20 21:46	
2-Butanone (MEK)	10 U	10	0.78	1	07/07/20 21:46	
2-Hexanone	10 U	10	0.20	1	07/07/20 21:46	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/07/20 21:46	
Acetone	10 U	10	5.0	1	07/07/20 21:46	
Benzene	5.0 U	5.0	0.20	1	07/07/20 21:46	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/07/20 21:46	
Bromoform	5.0 U	5.0	0.25	1	07/07/20 21:46	
Bromomethane	5.0 U	5.0	0.70	1	07/07/20 21:46	
Carbon Disulfide	1.2 J	10	0.42	1	07/07/20 21:46	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/07/20 21:46	
Chlorobenzene	5.0 U	5.0	0.20	1	07/07/20 21:46	
Chloroethane	5.0 U	5.0	0.23	1	07/07/20 21:46	
Chloroform	5.0 U	5.0	0.24	1	07/07/20 21:46	
Chloromethane	5.0 U	5.0	0.28	1	07/07/20 21:46	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/07/20 21:46	
Dichloromethane	5.0 U	5.0	0.65	1	07/07/20 21:46	
Ethylbenzene	5.0 U	5.0	0.20	1	07/07/20 21:46	
Styrene	5.0 U	5.0	0.20	1	07/07/20 21:46	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/07/20 21:46	
Toluene	5.0 U	5.0	0.20	1	07/07/20 21:46	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/07/20 21:46	
Vinyl Acetate	10 U	10	1.1	1	07/07/20 21:46	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/07/20 21:46	
Xylenes, Total	5.0 U	5.0	0.23	1	07/07/20 21:46	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/07/20 21:46	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/07/20 21:46	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/07/20 21:46	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/07/20 21:46	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 12:40

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-018
 Units: ug/L

 Lab Code:
 R2005539-005
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	102	85 - 122	07/07/20 21:46	
Dibromofluoromethane	103	89 - 119	07/07/20 21:46	
Toluene-d8	103	87 - 121	07/07/20 21:46	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 12:40

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-018
 Units: ug/L

 Lab Code:
 R2005539-005
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Tentatively Identified Compounds

 CAS#
 Compound Identification
 RT
 ug/L
 Q

 007446-09-5
 Sulfur dioxide
 1.23
 35.9
 JN

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 00:00

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 TB-9954-062620-RM-003
 Units: ug/L

 Lab Code:
 R2005539-006
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/01/20 00:15	_
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/01/20 00:15	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/01/20 00:15	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/01/20 00:15	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/01/20 00:15	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/01/20 00:15	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/01/20 00:15	
2-Butanone (MEK)	10 U	10	0.78	1	07/01/20 00:15	
2-Hexanone	10 U	10	0.20	1	07/01/20 00:15	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/01/20 00:15	
Acetone	10 U	10	5.0	1	07/01/20 00:15	
Benzene	5.0 U	5.0	0.20	1	07/01/20 00:15	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/01/20 00:15	
Bromoform	5.0 U	5.0	0.25	1	07/01/20 00:15	
Bromomethane	5.0 U	5.0	0.70	1	07/01/20 00:15	
Carbon Disulfide	10 U	10	0.42	1	07/01/20 00:15	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/01/20 00:15	
Chlorobenzene	5.0 U	5.0	0.20	1	07/01/20 00:15	
Chloroethane	5.0 U	5.0	0.23	1	07/01/20 00:15	
Chloroform	5.0 U	5.0	0.24	1	07/01/20 00:15	
Chloromethane	5.0 U	5.0	0.28	1	07/01/20 00:15	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/01/20 00:15	
Dichloromethane	5.0 U	5.0	0.65	1	07/01/20 00:15	
Ethylbenzene	5.0 U	5.0	0.20	1	07/01/20 00:15	
Styrene	5.0 U	5.0	0.20	1	07/01/20 00:15	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/01/20 00:15	
Toluene	5.0 U	5.0	0.20	1	07/01/20 00:15	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/01/20 00:15	
Vinyl Acetate	10 U	10	1.1	1	07/01/20 00:15	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/01/20 00:15	
Xylenes, Total	5.0 U	5.0	0.23	1	07/01/20 00:15	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/01/20 00:15	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/01/20 00:15	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/01/20 00:15	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/01/20 00:15	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 00:00

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 TB-9954-062620-RM-003
 Units: ug/L

 Lab Code:
 R2005539-006
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	91	85 - 122	07/01/20 00:15	
Dibromofluoromethane	89	89 - 119	07/01/20 00:15	
Toluene-d8	93	87 - 121	07/01/20 00:15	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 00:00

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 TB-9954-062620-RM-003
 Units: ug/L

 Lab Code:
 R2005539-006
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Tentatively Identified Compounds

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected



Semivolatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 08:50

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-014
 Units: ug/L

 Lab Code:
 R2005539-001
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 10:32	7/1/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 10:32	7/1/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/06/20 10:32	7/1/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 10:32	7/1/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/06/20 10:32	7/1/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/06/20 10:32	7/1/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/06/20 10:32	7/1/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/06/20 10:32	7/1/20	
2,4-Dinitrophenol	45 U	45	19	1	07/06/20 10:32	7/1/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/06/20 10:32	7/1/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/06/20 10:32	7/1/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/06/20 10:32	7/1/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/06/20 10:32	7/1/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/06/20 10:32	7/1/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/06/20 10:32	7/1/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/06/20 10:32	7/1/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/06/20 10:32	7/1/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/06/20 10:32	7/1/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/06/20 10:32	7/1/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/06/20 10:32	7/1/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/06/20 10:32	7/1/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/06/20 10:32	7/1/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/06/20 10:32	7/1/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/06/20 10:32	7/1/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/06/20 10:32	7/1/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/06/20 10:32	7/1/20	
4-Nitrophenol	45 U	45	5.8	1	07/06/20 10:32	7/1/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/06/20 10:32	7/1/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/06/20 10:32	7/1/20	
Anthracene	9.1 U	9.1	1.2	1	07/06/20 10:32	7/1/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/06/20 10:32	7/1/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/06/20 10:32	7/1/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/06/20 10:32	7/1/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/06/20 10:32	7/1/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/06/20 10:32	7/1/20	
Benzoic Acid	91 U	91	33	1	07/06/20 10:32	7/1/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/06/20 10:32	7/1/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/06/20 10:32	7/1/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/06/20 10:32	7/1/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	07/06/20 10:32	7/1/20	
Bis(2-ethylhexyl) Phthalate	2.2 J	9.1	0.91	1	07/06/20 10:32	7/1/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	07/06/20 10:32	7/1/20	
Chrysene	9.1 U	9.1	1.1	1	07/06/20 10:32	7/1/20	

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Superset Reference:20-0000554991 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 08:50

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-014
 Units: ug/L

 Lab Code:
 R2005539-001
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/06/20 10:32	7/1/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/06/20 10:32	7/1/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/06/20 10:32	7/1/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/06/20 10:32	7/1/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/06/20 10:32	7/1/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/06/20 10:32	7/1/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/06/20 10:32	7/1/20	
Fluorene	9.1 U	9.1	1.2	1	07/06/20 10:32	7/1/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/06/20 10:32	7/1/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/06/20 10:32	7/1/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/06/20 10:32	7/1/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/06/20 10:32	7/1/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/06/20 10:32	7/1/20	
Isophorone	9.1 U	9.1	1.3	1	07/06/20 10:32	7/1/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/06/20 10:32	7/1/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/06/20 10:32	7/1/20	
Naphthalene	9.1 U	9.1	1.1	1	07/06/20 10:32	7/1/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/06/20 10:32	7/1/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/06/20 10:32	7/1/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/06/20 10:32	7/1/20	
Phenol	9.1 U	9.1	0.91	1	07/06/20 10:32	7/1/20	
Pyrene	9.1 U	9.1	1.3	1	07/06/20 10:32	7/1/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	84	35 - 141	07/06/20 10:32	
2-Fluorobiphenyl	63	31 - 118	07/06/20 10:32	
2-Fluorophenol	42	10 - 105	07/06/20 10:32	
Nitrobenzene-d5	60	31 - 110	07/06/20 10:32	
Phenol-d6	30	10 - 107	07/06/20 10:32	
p-Terphenyl-d14	48	10 - 165	07/06/20 10:32	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown	6.30	5.2	J	
	unknown	6.51	24	J	
013798-23-7	Sulfur	7.81	14	JN	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 10:00

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-015
 Units: ug/L

 Lab Code:
 R2005539-002
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 11:00	7/1/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 11:00	7/1/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/06/20 11:00	7/1/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 11:00	7/1/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/06/20 11:00	7/1/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/06/20 11:00	7/1/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/06/20 11:00	7/1/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/06/20 11:00	7/1/20	
2,4-Dinitrophenol	45 U	45	19	1	07/06/20 11:00	7/1/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/06/20 11:00	7/1/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/06/20 11:00	7/1/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/06/20 11:00	7/1/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/06/20 11:00	7/1/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/06/20 11:00	7/1/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/06/20 11:00	7/1/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/06/20 11:00	7/1/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/06/20 11:00	7/1/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/06/20 11:00	7/1/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/06/20 11:00	7/1/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/06/20 11:00	7/1/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/06/20 11:00	7/1/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/06/20 11:00	7/1/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/06/20 11:00	7/1/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/06/20 11:00	7/1/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/06/20 11:00	7/1/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/06/20 11:00	7/1/20	
4-Nitrophenol	45 U	45	5.8	1	07/06/20 11:00	7/1/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/06/20 11:00	7/1/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/06/20 11:00	7/1/20	
Anthracene	9.1 U	9.1	1.2	1	07/06/20 11:00	7/1/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/06/20 11:00	7/1/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/06/20 11:00	7/1/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/06/20 11:00	7/1/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/06/20 11:00	7/1/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/06/20 11:00	7/1/20	
Benzoic Acid	91 U	91	33	1	07/06/20 11:00	7/1/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/06/20 11:00	7/1/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/06/20 11:00	7/1/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/06/20 11:00	7/1/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	07/06/20 11:00	7/1/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	07/06/20 11:00	7/1/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	07/06/20 11:00	7/1/20	
Chrysene	9.1 U	9.1	1.1	1	07/06/20 11:00	7/1/20	
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Superset Reference:20-0000554991 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 10:00

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-015
 Units: ug/L

 Lab Code:
 R2005539-002
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed 1	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/06/20 11:00	7/1/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/06/20 11:00	7/1/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/06/20 11:00	7/1/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/06/20 11:00	7/1/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/06/20 11:00	7/1/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/06/20 11:00	7/1/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/06/20 11:00	7/1/20	
Fluorene	9.1 U	9.1	1.2	1	07/06/20 11:00	7/1/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/06/20 11:00	7/1/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/06/20 11:00	7/1/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/06/20 11:00	7/1/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/06/20 11:00	7/1/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/06/20 11:00	7/1/20	
Isophorone	9.1 U	9.1	1.3	1	07/06/20 11:00	7/1/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/06/20 11:00	7/1/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/06/20 11:00	7/1/20	
Naphthalene	9.1 U	9.1	1.1	1	07/06/20 11:00	7/1/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/06/20 11:00	7/1/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/06/20 11:00	7/1/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/06/20 11:00	7/1/20	
Phenol	9.1 U	9.1	0.91	1	07/06/20 11:00	7/1/20	
Pyrene	9.1 U	9.1	1.3	1	07/06/20 11:00	7/1/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	86	35 - 141	07/06/20 11:00	
2-Fluorobiphenyl	57	31 - 118	07/06/20 11:00	
2-Fluorophenol	40	10 - 105	07/06/20 11:00	
Nitrobenzene-d5	56	31 - 110	07/06/20 11:00	
Phenol-d6	30	10 - 107	07/06/20 11:00	
p-Terphenyl-d14	59	10 - 165	07/06/20 11:00	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown hydrocarbon	11.35	5.9	J	
	unknown hydrocarbon	11.88	8.6	J	
	unknown hydrocarbon	12.48	10	J	
	unknown hydrocarbon	13.12	10	J	
	unknown hydrocarbon	13.82	9.0	J	
	unknown hydrocarbon	14.57	7.5	J	
	unknown hydrocarbon	15.36	5.7	J	

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Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 10:00

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-015
 Units: ug/L

 Lab Code:
 R2005539-002
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown hydrocarbon	16.11	4.3	J	
013798-23-7	Sulfur	7.76	5.1	JN	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 11:05

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-016
 Units: ug/L

 Lab Code:
 R2005539-003
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 11:27	7/1/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 11:27	7/1/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/06/20 11:27	7/1/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 11:27	7/1/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/06/20 11:27	7/1/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/06/20 11:27	7/1/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/06/20 11:27	7/1/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/06/20 11:27	7/1/20	
2,4-Dinitrophenol	45 U	45	19	1	07/06/20 11:27	7/1/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/06/20 11:27	7/1/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/06/20 11:27	7/1/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/06/20 11:27	7/1/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/06/20 11:27	7/1/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/06/20 11:27	7/1/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/06/20 11:27	7/1/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/06/20 11:27	7/1/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/06/20 11:27	7/1/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/06/20 11:27	7/1/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/06/20 11:27	7/1/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/06/20 11:27	7/1/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/06/20 11:27	7/1/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/06/20 11:27	7/1/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/06/20 11:27	7/1/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/06/20 11:27	7/1/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/06/20 11:27	7/1/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/06/20 11:27	7/1/20	
4-Nitrophenol	45 U	45	5.8	1	07/06/20 11:27	7/1/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/06/20 11:27	7/1/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/06/20 11:27	7/1/20	
Anthracene	9.1 U	9.1	1.2	1	07/06/20 11:27	7/1/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/06/20 11:27	7/1/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/06/20 11:27	7/1/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/06/20 11:27	7/1/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/06/20 11:27	7/1/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/06/20 11:27	7/1/20	
Benzoic Acid	91 U	91	33	1	07/06/20 11:27	7/1/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/06/20 11:27	7/1/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/06/20 11:27	7/1/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/06/20 11:27	7/1/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	07/06/20 11:27	7/1/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	07/06/20 11:27	7/1/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	07/06/20 11:27	7/1/20	
Chrysene	9.1 U	9.1	1.1	1	07/06/20 11:27	7/1/20	

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Superset Reference:20-0000554991 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 11:05

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-016
 Units: ug/L

 Lab Code:
 R2005539-003
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/06/20 11:27	7/1/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/06/20 11:27	7/1/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/06/20 11:27	7/1/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/06/20 11:27	7/1/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/06/20 11:27	7/1/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/06/20 11:27	7/1/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/06/20 11:27	7/1/20	
Fluorene	9.1 U	9.1	1.2	1	07/06/20 11:27	7/1/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/06/20 11:27	7/1/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/06/20 11:27	7/1/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/06/20 11:27	7/1/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/06/20 11:27	7/1/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/06/20 11:27	7/1/20	
Isophorone	9.1 U	9.1	1.3	1	07/06/20 11:27	7/1/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/06/20 11:27	7/1/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/06/20 11:27	7/1/20	
Naphthalene	9.1 U	9.1	1.1	1	07/06/20 11:27	7/1/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/06/20 11:27	7/1/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/06/20 11:27	7/1/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/06/20 11:27	7/1/20	
Phenol	9.1 U	9.1	0.91	1	07/06/20 11:27	7/1/20	
Pyrene	9.1 U	9.1	1.3	1	07/06/20 11:27	7/1/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	84	35 - 141	07/06/20 11:27	
2-Fluorobiphenyl	64	31 - 118	07/06/20 11:27	
2-Fluorophenol	35	10 - 105	07/06/20 11:27	
Nitrobenzene-d5	57	31 - 110	07/06/20 11:27	
Phenol-d6	27	10 - 107	07/06/20 11:27	
p-Terphenyl-d14	58	10 - 165	07/06/20 11:27	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown hydrocarbon	11.35	6.9	J	
	unknown hydrocarbon	11.88	11	J	
	unknown hydrocarbon	12.48	15	J	
	unknown hydrocarbon	13.12	15	J	
	unknown hydrocarbon	13.83	12	J	
	unknown hydrocarbon	14.57	12	J	
	unknown hydrocarbon	15.36	8.9	J	

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Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 11:05

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-016
 Units: ug/L

 Lab Code:
 R2005539-003
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown hydrocarbon	16.11	5.6	J	
013798-23-7	Sulfur	7.75	4.0	JN	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Date Collected: 06/26/20 11:55

Date Received: 06/27/20 08:50

Sample Matrix: Water

Sample Name: WG-9954-062620-RM-017 **Lab Code:** R2005539-004

Units: ug/L Basis: NA

Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

1,24-Trichlorobenzene	Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1.3-Dichlorobenzene	1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 11:55	7/1/20	
1,4-Dichlorobenzene	1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 11:55	7/1/20	
2,4.5-Trichlorophenol	1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/06/20 11:55	7/1/20	
2.4-Frichlorophenol	1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 11:55	7/1/20	
2.4-Dintchlorophenol	2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/06/20 11:55	7/1/20	
2.4-Dimethylphenol 9.1 U 9.1 U 9.1 D 1.3 D 1.07/06/20 11:55 7/1/20 2.4-Dinitrophenol 45 U 45 U 9.1 U 9.1 D 0.706/20 11:55 7/1/20 2.4-Dinitrophenol 9.1 U 9.1 L 1.2 L 1.07/06/20 11:55 7/1/20 2.6-Dinitrotoluene 9.1 U 9.1 L 1.2 L 1.07/06/20 11:55 7/1/20 2Chlorophenol 9.1 U 9.1 U 9.1 0.97 1.07/06/20 11:55 7/1/20 2-Methylaphthalene 9.1 U 9.1 U.9.1 0.97 1.07/06/20 11:55 7/1/20 2-Methylaphenol 9.1 U.9.1 0.91 1.07/06/20 11:55 7/1/20 2-Mitrophenol 9.1 U.9.1 1.3 1.07/06/20 11:55 7/1/20 2-Nitrophenol 9.1 U.9.1 1.4 1.07/06/20 11:55 7/1/20 3Aitrophenol Coelution 9.1 U.9.1 1.4 1.07/06/20 11:55 7/1/20 3Aitrophenol Coelution 9.1 U.9.1 1.1 1.07/06/20 11:55 7/1/20 4.6-Dinitro-2-methylphenol Coelution 45 U.9.1 1.5 1.07/06/20 11:55 7/1/20 4.Bromophenyl Phenyl Ether 9.1 U.9.1 1.5 1.07/06/20 11:55 7/1/20 <	2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/06/20 11:55	7/1/20	
2.4-Dinitrophenol 45 U 45 ID 1 07/06/20 11:55 7/1/20 2.4-Dinitrotoluene 9.1 U 9.1 L2 1 07/06/20 11:55 7/1/20 2.6-Dinitrotoluene 9.1 U 9.1 L3 1 07/06/20 11:55 7/1/20 2-Chloroaphthalene 9.1 U 9.1 0.97 1 07/06/20 11:55 7/1/20 2-Methylaphthalene 9.1 U 9.1 0.91 1.7 1 07/06/20 11:55 7/1/20 2-Methylaphthalene 9.1 U 9.1 0.91 1 07/06/20 11:55 7/1/20 2-Methylaphthalene 9.1 U 9.1 0.91 1 07/06/20 11:55 7/1/20 2-Methylaphthalene 9.1 U 9.1 0.91 1 07/06/20 11:55 7/1/20 2-Nitroaniline 9.1 U 9.1 1.3 1 07/06/20 11:55 7/1/20 2-Nitroaniline 9.1 U 9.1 1.1 1 07/06/20 11:55 7/1/20 3.3-Dichlorobenzidine 9.1 U 9.1 1.1 1 07/06/20 11:55 7/1/20 3.3-Dichlorobenzidine 9.1 U 9.1 U 9.1 1.1 1 07/06/20 11:55 7/1/20 3-Nitroaniline 9.1 U	2,4-Dichlorophenol		9.1		1	07/06/20 11:55	7/1/20	
2.4-Dinitrotoluene 9.1 U 9.1 U 9.1 L2 1 O7/06/20 11:55 7/1/20 2.6-Dinitrotoluene 9.1 U 9.1 L3 1 O7/06/20 11:55 7/1/20 2-Chlorophenol 9.1 U 9.1 U 9.1 U 9.1 O7/06/20 11:55 7/1/20 2-Methylnaphthalene 9.1 U 9.1 U 9.1 L2 1 O7/06/20 11:55 7/1/20 2-Methylphenol 9.1 U 9.1 U 9.1 U 9.1 O/91 1 O7/06/20 11:55 7/1/20 2-Mitrophenol 9.1 U 9.1 U 9.1 U 9.1 U 9.1 U 7/1/20 2-Nitrophenol 9.1 U 9.1 U 9.1 U 9.1 U 9.1 U 7/1/20 3.3'-Dichlorobenzidine 9.1 U 9.1	2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/06/20 11:55	7/1/20	
2,6-Dinitrotoluene 9.1 U 9.1 I.2 1 07/06/20 11:55 7/1/20 2-Chlorophenol 9.1 U 9.1 U.9.1 1.3 1 07/06/20 11:55 7/1/20 2-Chlorophenol 9.1 U.9.1 0.97 1 07/06/20 11:55 7/1/20 2-Methylaphthalene 9.1 U.9.1 1.2 1 07/06/20 11:55 7/1/20 2-Methylphenol 9.1 U.9.1 1.2 1 07/06/20 11:55 7/1/20 2-Nitroaniline 9.1 U.9.1 1.4 1 07/06/20 11:55 7/1/20 2-Nitrophenol 9.1 U.9.1 1.4 1 07/06/20 11:55 7/1/20 3-Sichlorobenzidine 9.1 U.9.1 1.1 1 07/06/20 11:55 7/1/20 3-Ad 4-Methylphenol Coelution 9.1 U.9.1 2.3 1 07/06/20 11:55 7/1/20 4-G-Dinitro-2-methylphenol 45 U.9.1 45 I.8 1 07/06/20 11:55 7/1/20 4-Bromophenyl Phenyl Ether 9.1 U.9.1 9.1 U.9.1 0.98 I.07/06/20 11:55 7/1/20 4-Chloro-3-methylphenol 9.1 U.9.1 0.98 I.07/06/20 11:55 7/1/20 4-Chlorophenyl Phenyl Ether	2,4-Dinitrophenol	45 U	45	19	1	07/06/20 11:55	7/1/20	
2-Chloronaphthalene	2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/06/20 11:55	7/1/20	
2-Chlorophenol 9.1 U	2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/06/20 11:55	7/1/20	
2-Methylnaphthalene	2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/06/20 11:55	7/1/20	
2-Methylphenol		9.1 U	9.1	0.97	1	07/06/20 11:55	7/1/20	
2-Nitroaniline	2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/06/20 11:55	7/1/20	
2-Nitroaniline	2-Methylphenol	9.1 U	9.1	0.91	1	07/06/20 11:55	7/1/20	
3,3'-Dichlorobenzidine			9.1	1.3	1		7/1/20	
3,3'-Dichlorobenzidine	2-Nitrophenol	9.1 U	9.1	1.4	1	07/06/20 11:55	7/1/20	
3- and 4-Methylphenol Coelution 9.1 U 9.1 U 9.1 D 1.1 D 07/06/20 11:55 D 7/1/20 D 3-Nitroaniline 9.1 U 9.1 U 9.1 E 2.3 1 D 07/06/20 11:55 D 7/1/20 D 4,6-Dinitro-2-methylphenol 45 U 45 I8 I 1 07/06/20 11:55 D 7/1/20 D 4-Bromophenyl Phenyl Ether 9.1 U 9.1 U 9.1 D 9.1 U 9.1 D 9.1 U 9.		9.1 U	9.1	1.1	1		7/1/20	
3-Nitroaniline								
4,6-Dinitro-2-methylphenol 45 U 45 18 1 07/06/20 11:55 7/1/20 4-Bromophenyl Phenyl Ether 9.1 U 9.1 1.5 1 07/06/20 11:55 7/1/20 4-Chloro-3-methylphenol 9.1 U 9.1 0.98 1 07/06/20 11:55 7/1/20 4-Chloroaniline 9.1 U 9.1 0.91 1 07/06/20 11:55 7/1/20 4-Chlorophenyl Phenyl Ether 9.1 U 9.1 1.4 1 07/06/20 11:55 7/1/20 4-Nitroaniline 9.1 U 9.1 2.5 1 07/06/20 11:55 7/1/20 4-Nitrophenol 45 U 45 5.8 1 07/06/20 11:55 7/1/20 4-Nitrophenol 45 U 45 5.8 1 07/06/20 11:55 7/1/20 Acenaphthene 9.1 U 9.1 1.3 1 07/06/20 11:55 7/1/20 Acenaphthylene 9.1 U 9.1 1.3 1 07/06/20 11:55 7/1/20 Anthracene 9.1 U 9.1 1.2 1 07/06/20 11:55 7/1/20 Benz(a)apyrene 9.1 U 9.1 1.5 1 07/06/20 11:55 7/1/20 Benzo(b)fluoranthene 9.1 U 9.1 1.1 1 07/06/20 11:55 <td></td> <td>9.1 U</td> <td>9.1</td> <td>2.3</td> <td>1</td> <td>07/06/20 11:55</td> <td>7/1/20</td> <td></td>		9.1 U	9.1	2.3	1	07/06/20 11:55	7/1/20	
4-Bromophenyl Phenyl Ether 9.1 U 9.1 U 9.1 0.98 1 07/06/20 11:55 7/1/20 4-Chloro-3-methylphenol 9.1 U 9.1 0.91 0.98 1 07/06/20 11:55 7/1/20 4-Chloroaniline 9.1 U 9.1 0.91 1 07/06/20 11:55 7/1/20 4-Chlorophenyl Phenyl Ether 9.1 U 9.1 1.4 1 07/06/20 11:55 7/1/20 4-Nitrophenol 45 U 45 5.8 1 07/06/20 11:55 7/1/20 4-Nitrophenol 45 U 45 5.8 1 07/06/20 11:55 7/1/20 Acenaphthene 9.1 U 9.1 1.3 1 07/06/20 11:55 7/1/20 Acenaphthene 9.1 U 9.1 1.3 1 07/06/20 11:55 7/1/20 Achtracene 9.1 U 9.1 1.3 1 07/06/20 11:55 7/1/20 Benz(a)anthracene 9.1 U 9.1 1.5 1 07/06/20 11:55 7/1/20 Benzo(a)pyrene 9.1 U 9.1 1.1 1 07/06/20 11:55 7/1/20 Benzo(b)fluoranthene 9.1 U 9.1 1.1 1 07/06/20 11:55 7/1/20 Benzo(b)fluoranthene 9.1 U 9.1 0.91 1 07/06/20 11:55 7/1/20 Benzo(k)fluoranthene <td>4,6-Dinitro-2-methylphenol</td> <td>45 U</td> <td>45</td> <td>18</td> <td>1</td> <td></td> <td>7/1/20</td> <td></td>	4,6-Dinitro-2-methylphenol	45 U	45	18	1		7/1/20	
4-Chloro-3-methylphenol 9.1 U 9.1 D 9.1	4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/06/20 11:55	7/1/20	
4-Chlorophenyl Phenyl Ether 9.1 U 9.1 1.4 1 07/06/20 11:55 7/1/20 4-Nitroaniline 9.1 U 9.1 U 9.1 U 2.5 1 07/06/20 11:55 7/1/20 4-Nitrophenol 45 U 45 S.8 1 07/06/20 11:55 7/1/20 Acenaphthene 9.1 U 9.1 I.3 1 07/06/20 11:55 7/1/20 Acenaphthylene 9.1 U 9.1 I.3 1 07/06/20 11:55 7/1/20 Anthracene 9.1 U 9.1 I.2 1 07/06/20 11:55 7/1/20 Benz(a)anthracene 9.1 U 9.1 I.5 1 07/06/20 11:55 7/1/20 Benzo(a)pyrene 9.1 U 9.1 I.1 1 07/06/20 11:55 7/1/20 Benzo(b)fluoranthene 9.1 U 9.1 I.1 1 07/06/20 11:55 7/1/20 Benzo(c)g,h,i)perylene 9.1 U 9.1 I.1 1 07/06/20 11:55 7/1/20 Benzo(k)fluoranthene 9.1 U 9.1 I.1 1 07/06/20 11:55 7/1/20 Benzolc Acid 91 U 9.1 I.1 1 07/06/20 11:55 7/1/20 Benzyl Alcohol 9.1 U 9.1 I.5		9.1 U	9.1	0.98	1	07/06/20 11:55	7/1/20	
4-Chlorophenyl Phenyl Ether 9.1 U 9.1 1.4 1 07/06/20 11:55 7/1/20 4-Nitroaniline 9.1 U 9.1 U 9.1 U 2.5 1 07/06/20 11:55 7/1/20 4-Nitrophenol 45 U 45 S.8 1 07/06/20 11:55 7/1/20 Acenaphthene 9.1 U 9.1 I.3 1 07/06/20 11:55 7/1/20 Acenaphthylene 9.1 U 9.1 I.3 1 07/06/20 11:55 7/1/20 Anthracene 9.1 U 9.1 I.2 1 07/06/20 11:55 7/1/20 Benz(a)anthracene 9.1 U 9.1 I.5 1 07/06/20 11:55 7/1/20 Benzo(a)pyrene 9.1 U 9.1 I.1 1 07/06/20 11:55 7/1/20 Benzo(b)fluoranthene 9.1 U 9.1 I.1 1 07/06/20 11:55 7/1/20 Benzo(c)g,h,i)perylene 9.1 U 9.1 I.1 1 07/06/20 11:55 7/1/20 Benzo(k)fluoranthene 9.1 U 9.1 I.1 1 07/06/20 11:55 7/1/20 Benzolc Acid 91 U 9.1 I.1 1 07/06/20 11:55 7/1/20 Benzyl Alcohol 9.1 U 9.1 I.5	4-Chloroaniline	9.1 U	9.1	0.91	1	07/06/20 11:55	7/1/20	
4-Nitrophenol 45 U 45 5.8 1 07/06/20 11:55 7/1/20 Acenaphthene 9.1 U 9.1 1.3 1 07/06/20 11:55 7/1/20 Acenaphthylene 9.1 U 9.1 1.3 1 07/06/20 11:55 7/1/20 Anthracene 9.1 U 9.1 1.2 1 07/06/20 11:55 7/1/20 Benz(a)anthracene 9.1 U 9.1 1.5 1 07/06/20 11:55 7/1/20 Benzo(a)pyrene 9.1 U 9.1 1.1 1 07/06/20 11:55 7/1/20 Benzo(b)fluoranthene 9.1 U 9.1 1.1 1 07/06/20 11:55 7/1/20 Benzo(g,h,i)perylene 9.1 U 9.1 0.91 1 07/06/20 11:55 7/1/20 Benzo(k)fluoranthene 9.1 U 9.1 1.1 1 07/06/20 11:55 7/1/20 Benzoic Acid 9.1 U 9.1 1.1 1 07/06/20 11:55 7/1/20 Benzyl Alcohol 9.1 U 9.1 1.5 1 07/06/20 11:55 7/1/20 2,2'-Oxybis(1-chloropropane) 9.1 U 9.1 1.3 1 07/06/20 11:55 7/1/20 Bis(2-chloroethoxy)methane 9.1 U 9.1 1.8 1 07/06/20 11:55 7/1/20 Bis(2-chloroethyl) Ether 9.1 U 9.1 U.			9.1	1.4	1	07/06/20 11:55	7/1/20	
Acenaphthene 9.1 U 9.1 I.3 1 07/06/20 11:55 7/1/20 Acenaphthylene 9.1 U 9.1 I.3 1 07/06/20 11:55 7/1/20 Anthracene 9.1 U 9.1 I.2 1 07/06/20 11:55 7/1/20 Benz(a)anthracene 9.1 U 9.1 I.5 1 07/06/20 11:55 7/1/20 Benzo(a)pyrene 9.1 U 9.1 I.1 1 07/06/20 11:55 7/1/20 Benzo(b)fluoranthene 9.1 U 9.1 I.1 1 07/06/20 11:55 7/1/20 Benzo(g,h,i)perylene 9.1 U 9.1 I.1 1 07/06/20 11:55 7/1/20 Benzo(k)fluoranthene 9.1 U 9.1 I.1 1 07/06/20 11:55 7/1/20 Benzoic Acid 91 U 9.1 I.1 1 07/06/20 11:55 7/1/20 Benzyl Alcohol 9.1 U 9.1 I.5 1 07/06/20 11:55 7/1/20 2,2'-Oxybis(1-chloropropane) 9.1 U 9.1 I.3 1 07/06/20 11:55 7/1/20 Bis(2-chloroethoxy)methane 9.1 U 9.1 I.8 1 07/06/20 11:55 7/1/20 Bis(2-chloroethyl) Ether 9.1 U 9.1 I.2 1 07/06/20 11:55 7/1/20 Butyl Benzyl Phthalate 9.1 U <	4-Nitroaniline	9.1 U	9.1	2.5	1	07/06/20 11:55	7/1/20	
Acenaphthene 9.1 U 9.1 1.3 1 07/06/20 11:55 7/1/20 Acenaphthylene 9.1 U 9.1 1.3 1 07/06/20 11:55 7/1/20 Anthracene 9.1 U 9.1 1.2 1 07/06/20 11:55 7/1/20 Benz(a)anthracene 9.1 U 9.1 1.5 1 07/06/20 11:55 7/1/20 Benzo(a)pyrene 9.1 U 9.1 1.1 1 07/06/20 11:55 7/1/20 Benzo(b)fluoranthene 9.1 U 9.1 1.1 1 07/06/20 11:55 7/1/20 Benzo(g,h,i)perylene 9.1 U 9.1 0.91 1 07/06/20 11:55 7/1/20 Benzo(k)fluoranthene 9.1 U 9.1 1.1 1 07/06/20 11:55 7/1/20 Benzoic Acid 9.1 U 9.1 33 1 07/06/20 11:55 7/1/20 Benzyl Alcohol 9.1 U 9.1 1.5 1 07/06/20 11:55 7/1/20 2,2'-Oxybis(1-chloropropane) 9.1 U 9.1 1.3 1 07/06/20 11:55 7/1/20 Bis(2-chloroethoxy)methane 9.1 U 9.1 1.8 1 07/06/20 11:55 7/1/20 Bis(2-chloroethyl) Ether 9.1 U 9.1 1.2 1 07/06/20 11:55 7/1/20 Butyl Benzyl Phthalate 9.1 U	4-Nitrophenol	45 U	45	5.8	1	07/06/20 11:55	7/1/20	
Acenaphthylene 9.1 U 9.1 U 9.1 I.3 1 07/06/20 11:55 7/1/20 Anthracene 9.1 U 9.1 U 9.1 I.2 1 07/06/20 11:55 7/1/20 Benz(a)anthracene 9.1 U 9.1 I.5 1 07/06/20 11:55 7/1/20 Benzo(a)pyrene 9.1 U 9.1 I.1 1 07/06/20 11:55 7/1/20 Benzo(b)fluoranthene 9.1 U 9.1 I.1 1 07/06/20 11:55 7/1/20 Benzo(g,h,i)perylene 9.1 U		9.1 U	9.1	1.3	1	07/06/20 11:55	7/1/20	
Anthracene 9.1 U 9.1 1.2 1 07/06/20 11:55 7/1/20 Benz(a)anthracene 9.1 U 9.1 1.5 1 07/06/20 11:55 7/1/20 Benzo(a)pyrene 9.1 U 9.1 1.1 1 07/06/20 11:55 7/1/20 Benzo(b)fluoranthene 9.1 U 9.1 1.1 1 07/06/20 11:55 7/1/20 Benzo(g,h,i)perylene 9.1 U 9.1 0.91 1 07/06/20 11:55 7/1/20 Benzo(k)fluoranthene 9.1 U 9.1 1.1 1 07/06/20 11:55 7/1/20 Benzoic Acid 9.1 U 9.1 33 1 07/06/20 11:55 7/1/20 Benzyl Alcohol 9.1 U 9.1 1.5 1 07/06/20 11:55 7/1/20 2,2'-Oxybis(1-chloropropane) 9.1 U 9.1 1.3 1 07/06/20 11:55 7/1/20 Bis(2-chloroethoxy)methane 9.1 U 9.1 1.8 1 07/06/20 11:55 7/1/20 Bis(2-chloroethyl) Ether 9.1 U 9.1 1.2 1 07/06/20 11:55 7/1/20 Butyl Benzyl Phthalate 9.1 U 9.1 0.91 1 0.91 1 07/06/20 11:55 7/1/20		9.1 U	9.1	1.3	1	07/06/20 11:55	7/1/20	
Benz(a)anthracene 9.1 U 9.1 1.5 1 07/06/20 11:55 7/1/20 Benzo(a)pyrene 9.1 U 9.1 1.1 1 07/06/20 11:55 7/1/20 Benzo(b)fluoranthene 9.1 U 9.1 1.1 1 07/06/20 11:55 7/1/20 Benzo(g,h,i)perylene 9.1 U 9.1 0.91 1 07/06/20 11:55 7/1/20 Benzo(k)fluoranthene 9.1 U 9.1 1.1 1 07/06/20 11:55 7/1/20 Benzoic Acid 91 U 91 33 1 07/06/20 11:55 7/1/20 Benzyl Alcohol 9.1 U 9.1 1.5 1 07/06/20 11:55 7/1/20 2,2'-Oxybis(1-chloropropane) 9.1 U 9.1 1.3 1 07/06/20 11:55 7/1/20 Bis(2-chloroethoxy)methane 9.1 U 9.1 1.8 1 07/06/20 11:55 7/1/20 Bis(2-chloroethyl) Ether 9.1 U 9.1 1.2 1 07/06/20 11:55 7/1/20 Bis(2-ethylhexyl) Phthalate 9.1 U 9.1 0.91 1 07/06/20 11:55 7/1/20 Butyl Benzyl Phthalate 9.1 U 9.1 1.3 1 07/06/20 11:55 7/1/20		9.1 U	9.1	1.2	1	07/06/20 11:55	7/1/20	
Benzo(b)fluoranthene 9.1 U 9.1 D 9.1 D </td <td></td> <td>9.1 U</td> <td>9.1</td> <td></td> <td>1</td> <td></td> <td>7/1/20</td> <td></td>		9.1 U	9.1		1		7/1/20	
Benzo(b)fluoranthene 9.1 U 9.1 D 9.1 D </td <td>Benzo(a)pyrene</td> <td>9.1 U</td> <td>9.1</td> <td>1.1</td> <td>1</td> <td>07/06/20 11:55</td> <td>7/1/20</td> <td></td>	Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/06/20 11:55	7/1/20	
Benzo(g,h,i)perylene 9.1 U 9.1 D 0.91 D 1 07/06/20 11:55 7/1/20 Benzo(k)fluoranthene 9.1 U 9.1 D 1.1 D 1 07/06/20 11:55 7/1/20 Benzoic Acid 91 U 91 S 33 D 1 07/06/20 11:55 7/1/20 Benzyl Alcohol 9.1 U 9.1 D 1.5 D 1 07/06/20 11:55 7/1/20 2,2'-Oxybis(1-chloropropane) 9.1 U 9.1 D 1.3 D 1 07/06/20 11:55 7/1/20 Bis(2-chloroethoxy)methane 9.1 U 9.1 D 1.8 D 1 07/06/20 11:55 7/1/20 Bis(2-chloroethyl) Ether 9.1 U 9.1 D 9.1 D 1.2 D 1 07/06/20 11:55 7/1/20 Bis(2-ethylhexyl) Phthalate 9.1 U 9.1 D 0.91 D 1 07/06/20 11:55 7/1/20 Butyl Benzyl Phthalate 9.1 U 9.1 D 1.3 D 1 07/06/20 11:55 7/1/20		9.1 U	9.1	1.1	1		7/1/20	
Benzo(k)fluoranthene 9.1 U 9.1 1.1 1 07/06/20 11:55 7/1/20 Benzoic Acid 91 U 91 33 1 07/06/20 11:55 7/1/20 Benzyl Alcohol 9.1 U 9.1 1.5 1 07/06/20 11:55 7/1/20 2,2'-Oxybis(1-chloropropane) 9.1 U 9.1 1.3 1 07/06/20 11:55 7/1/20 Bis(2-chloroethoxy)methane 9.1 U 9.1 1.8 1 07/06/20 11:55 7/1/20 Bis(2-chloroethyl) Ether 9.1 U 9.1 1.2 1 07/06/20 11:55 7/1/20 Bis(2-ethylhexyl) Phthalate 9.1 U 9.1 0.91 1 07/06/20 11:55 7/1/20 Butyl Benzyl Phthalate 9.1 U 9.1 1.3 1 07/06/20 11:55 7/1/20		9.1 U	9.1	0.91	1	07/06/20 11:55	7/1/20	
Benzoic Acid 91 U 91 33 1 07/06/20 11:55 7/1/20 Benzyl Alcohol 9.1 U 9.1 1.5 1 07/06/20 11:55 7/1/20 2,2'-Oxybis(1-chloropropane) 9.1 U 9.1 1.3 1 07/06/20 11:55 7/1/20 Bis(2-chloroethoxy)methane 9.1 U 9.1 1.8 1 07/06/20 11:55 7/1/20 Bis(2-chloroethyl) Ether 9.1 U 9.1 1.2 1 07/06/20 11:55 7/1/20 Bis(2-ethylhexyl) Phthalate 9.1 U 9.1 0.91 1 07/06/20 11:55 7/1/20 Butyl Benzyl Phthalate 9.1 U 9.1 1.3 1 07/06/20 11:55 7/1/20		9.1 U	9.1	1.1	1	07/06/20 11:55	7/1/20	
Benzyl Alcohol 9.1 U 9.1 1.5 1 07/06/20 11:55 7/1/20 2,2'-Oxybis(1-chloropropane) 9.1 U 9.1 1.3 1 07/06/20 11:55 7/1/20 Bis(2-chloroethoxy)methane 9.1 U 9.1 1.8 1 07/06/20 11:55 7/1/20 Bis(2-chloroethyl) Ether 9.1 U 9.1 1.2 1 07/06/20 11:55 7/1/20 Bis(2-ethylhexyl) Phthalate 9.1 U 9.1 0.91 1 07/06/20 11:55 7/1/20 Butyl Benzyl Phthalate 9.1 U 9.1 1.3 1 07/06/20 11:55 7/1/20					1			
2,2'-Oxybis(1-chloropropane) 9.1 U 9.1 1.3 1 07/06/20 11:55 7/1/20 Bis(2-chloroethoxy)methane 9.1 U 9.1 1.8 1 07/06/20 11:55 7/1/20 Bis(2-chloroethyl) Ether 9.1 U 9.1 1.2 1 07/06/20 11:55 7/1/20 Bis(2-ethylhexyl) Phthalate 9.1 U 9.1 0.91 1 07/06/20 11:55 7/1/20 Butyl Benzyl Phthalate 9.1 U 9.1 1.3 1 07/06/20 11:55 7/1/20		9.1 U	9.1	1.5	1	07/06/20 11:55	7/1/20	
Bis(2-chloroethoxy)methane 9.1 U 9.1 1.8 1 07/06/20 11:55 7/1/20 Bis(2-chloroethyl) Ether 9.1 U 9.1 1.2 1 07/06/20 11:55 7/1/20 Bis(2-ethylhexyl) Phthalate 9.1 U 9.1 0.91 1 07/06/20 11:55 7/1/20 Butyl Benzyl Phthalate 9.1 U 9.1 1.3 1 07/06/20 11:55 7/1/20		9.1 U	9.1	1.3	1	07/06/20 11:55	7/1/20	
Bis(2-chloroethyl) Ether 9.1 U 9.1 1.2 1 07/06/20 11:55 7/1/20 Bis(2-ethylhexyl) Phthalate 9.1 U 9.1 0.91 1 07/06/20 11:55 7/1/20 Butyl Benzyl Phthalate 9.1 U 9.1 1.3 1 07/06/20 11:55 7/1/20								
Bis(2-ethylhexyl) Phthalate 9.1 U 9.1 U 9.1 U 0.91 I 0.91 U		9.1 U	9.1	1.2	1	07/06/20 11:55	7/1/20	
Butyl Benzyl Phthalate 9.1 U 9.1 1.3 1 07/06/20 11:55 7/1/20								
	Chrysene							

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Superset Reference:20-0000554991 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 11:55

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-017
 Units: ug/L

 Lab Code:
 R2005539-004
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/06/20 11:55	7/1/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/06/20 11:55	7/1/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/06/20 11:55	7/1/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/06/20 11:55	7/1/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/06/20 11:55	7/1/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/06/20 11:55	7/1/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/06/20 11:55	7/1/20	
Fluorene	9.1 U	9.1	1.2	1	07/06/20 11:55	7/1/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/06/20 11:55	7/1/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/06/20 11:55	7/1/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/06/20 11:55	7/1/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/06/20 11:55	7/1/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/06/20 11:55	7/1/20	
Isophorone	9.1 U	9.1	1.3	1	07/06/20 11:55	7/1/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/06/20 11:55	7/1/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/06/20 11:55	7/1/20	
Naphthalene	9.1 U	9.1	1.1	1	07/06/20 11:55	7/1/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/06/20 11:55	7/1/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/06/20 11:55	7/1/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/06/20 11:55	7/1/20	
Phenol	9.1 U	9.1	0.91	1	07/06/20 11:55	7/1/20	
Pyrene	9.1 U	9.1	1.3	1	07/06/20 11:55	7/1/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	83	35 - 141	07/06/20 11:55	
2-Fluorobiphenyl	70	31 - 118	07/06/20 11:55	
2-Fluorophenol	41	10 - 105	07/06/20 11:55	
Nitrobenzene-d5	68	31 - 110	07/06/20 11:55	
Phenol-d6	32	10 - 107	07/06/20 11:55	
p-Terphenyl-d14	57	10 - 165	07/06/20 11:55	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown hydrocarbon	11.35	4.8	J	
	unknown hydrocarbon	11.88	7.6	J	
	unknown	12.48	11	J	
	unknown hydrocarbon	13.12	11	J	
	unknown hydrocarbon	13.83	8.4	J	
	unknown hydrocarbon	14.58	8.2	J	
	unknown hydrocarbon	15.36	5.9	J	

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Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 11:55

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-017
 Units: ug/L

 Lab Code:
 R2005539-004
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown	16.11	5.1	J	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 12:40

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-018
 Units: ug/L

 Lab Code:
 R2005539-005
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 12:23	7/1/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 12:23	7/1/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/06/20 12:23	7/1/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 12:23	7/1/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/06/20 12:23	7/1/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/06/20 12:23	7/1/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/06/20 12:23	7/1/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/06/20 12:23	7/1/20	
2,4-Dinitrophenol	45 U	45	19	1	07/06/20 12:23	7/1/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/06/20 12:23	7/1/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/06/20 12:23	7/1/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/06/20 12:23	7/1/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/06/20 12:23	7/1/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/06/20 12:23	7/1/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/06/20 12:23	7/1/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/06/20 12:23	7/1/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/06/20 12:23	7/1/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/06/20 12:23	7/1/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/06/20 12:23	7/1/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/06/20 12:23	7/1/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/06/20 12:23	7/1/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/06/20 12:23	7/1/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/06/20 12:23	7/1/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/06/20 12:23	7/1/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/06/20 12:23	7/1/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/06/20 12:23	7/1/20	
4-Nitrophenol	45 U	45	5.8	1	07/06/20 12:23	7/1/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/06/20 12:23	7/1/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/06/20 12:23	7/1/20	
Anthracene	9.1 U	9.1	1.2	1	07/06/20 12:23	7/1/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/06/20 12:23	7/1/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/06/20 12:23	7/1/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/06/20 12:23	7/1/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/06/20 12:23	7/1/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/06/20 12:23	7/1/20	
Benzoic Acid	91 U	91	33	1	07/06/20 12:23	7/1/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/06/20 12:23	7/1/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/06/20 12:23	7/1/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/06/20 12:23	7/1/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	07/06/20 12:23	7/1/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	07/06/20 12:23	7/1/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	07/06/20 12:23	7/1/20	
Chrysene	9.1 U	9.1	1.1	1	07/06/20 12:23	7/1/20	

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Superset Reference:20-0000554991 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 12:40

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-018
 Units: ug/L

 Lab Code:
 R2005539-005
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/06/20 12:23	7/1/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/06/20 12:23	7/1/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/06/20 12:23	7/1/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/06/20 12:23	7/1/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/06/20 12:23	7/1/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/06/20 12:23	7/1/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/06/20 12:23	7/1/20	
Fluorene	9.1 U	9.1	1.2	1	07/06/20 12:23	7/1/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/06/20 12:23	7/1/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/06/20 12:23	7/1/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/06/20 12:23	7/1/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/06/20 12:23	7/1/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/06/20 12:23	7/1/20	
Isophorone	9.1 U	9.1	1.3	1	07/06/20 12:23	7/1/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/06/20 12:23	7/1/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/06/20 12:23	7/1/20	
Naphthalene	9.1 U	9.1	1.1	1	07/06/20 12:23	7/1/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/06/20 12:23	7/1/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/06/20 12:23	7/1/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/06/20 12:23	7/1/20	
Phenol	9.1 U	9.1	0.91	1	07/06/20 12:23	7/1/20	
Pyrene	9.1 U	9.1	1.3	1	07/06/20 12:23	7/1/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	86	35 - 141	07/06/20 12:23	
2-Fluorobiphenyl	62	31 - 118	07/06/20 12:23	
2-Fluorophenol	40	10 - 105	07/06/20 12:23	
Nitrobenzene-d5	60	31 - 110	07/06/20 12:23	
Phenol-d6	30	10 - 107	07/06/20 12:23	
p-Terphenyl-d14	69	10 - 165	07/06/20 12:23	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown hydrocarbon	11.35	5.0	J	
	unknown hydrocarbon	11.89	7.2	J	
	unknown hydrocarbon	12.48	9.9	J	
	unknown hydrocarbon	13.13	9.6	J	
	unknown hydrocarbon	13.83	8.3	J	
	unknown hydrocarbon	14.58	7.3	J	
	unknown	15.36	5.3	J	

Printed 7/15/2020 12:38:04 PM

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 12:40

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-018
 Units: ug/L

 Lab Code:
 R2005539-005
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Tentatively Identified Compounds

 CAS#
 Compound Identification
 RT
 ug/L
 Q

 013798-23-7
 Sulfur
 7.79
 9.1
 JN



Semivolatile Organic Compounds by GC

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 08:50

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-014
 Units: ug/L

 Lab Code:
 R2005539-001
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/06/20 12:16	7/1/20	_
4,4'-DDE	0.045 U	0.045	0.019	1	07/06/20 12:16	7/1/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/06/20 12:16	7/1/20	
Aldrin	0.045 U	0.045	0.019	1	07/06/20 12:16	7/1/20	
Dieldrin	0.045 U	0.045	0.019	1	07/06/20 12:16	7/1/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/06/20 12:16	7/1/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/06/20 12:16	7/1/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/06/20 12:16	7/1/20	
Endrin	0.045 U	0.045	0.019	1	07/06/20 12:16	7/1/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/06/20 12:16	7/1/20	
Heptachlor	0.045 U	0.045	0.019	1	07/06/20 12:16	7/1/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/06/20 12:16	7/1/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/06/20 12:16	7/1/20	
Toxaphene	0.46 U	0.46	0.46	1	07/06/20 12:16	7/1/20	
alpha-BHC	0.088	0.045	0.019	1	07/06/20 12:16	7/1/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/06/20 12:16	7/1/20	
beta-BHC	0.036 J	0.045	0.019	1	07/06/20 12:16	7/1/20	
delta-BHC	0.32	0.045	0.019	1	07/06/20 12:16	7/1/20	
gamma-BHC (Lindane)	0.11	0.045	0.019	1	07/06/20 12:16	7/1/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/06/20 12:16	7/1/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	42	10 - 164	07/06/20 12:16	
Tetrachloro-m-xylene	52	10 - 147	07/06/20 12:16	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 10:00

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-015
 Units: ug/L

 Lab Code:
 R2005539-002
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/07/20 16:01	7/1/20	_
4,4'-DDE	0.045 U	0.045	0.019	1	07/07/20 16:01	7/1/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/07/20 16:01	7/1/20	
Aldrin	0.045 U	0.045	0.019	1	07/07/20 16:01	7/1/20	
Dieldrin	0.045 U	0.045	0.019	1	07/07/20 16:01	7/1/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/07/20 16:01	7/1/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/07/20 16:01	7/1/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/07/20 16:01	7/1/20	
Endrin	0.045 U	0.045	0.019	1	07/07/20 16:01	7/1/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/07/20 16:01	7/1/20	
Heptachlor	0.045 U	0.045	0.019	1	07/07/20 16:01	7/1/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/07/20 16:01	7/1/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/07/20 16:01	7/1/20	
Toxaphene	0.46 U	0.46	0.46	1	07/07/20 16:01	7/1/20	
alpha-BHC	0.064	0.045	0.019	1	07/07/20 16:01	7/1/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/07/20 16:01	7/1/20	
beta-BHC	0.045 U	0.045	0.019	1	07/07/20 16:01	7/1/20	
delta-BHC	0.042 J	0.045	0.019	1	07/07/20 16:01	7/1/20	
gamma-BHC (Lindane)	0.060	0.045	0.019	1	07/07/20 16:01	7/1/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/07/20 16:01	7/1/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	67	10 - 164	07/07/20 16:01	
Tetrachloro-m-xylene	59	10 - 147	07/07/20 16:01	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 11:05

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-016
 Units: ug/L

 Lab Code:
 R2005539-003
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/07/20 16:20	7/1/20	_
4,4'-DDE	0.045 U	0.045	0.019	1	07/07/20 16:20	7/1/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/07/20 16:20	7/1/20	
Aldrin	0.045 U	0.045	0.019	1	07/07/20 16:20	7/1/20	
Dieldrin	0.045 U	0.045	0.019	1	07/07/20 16:20	7/1/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/07/20 16:20	7/1/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/07/20 16:20	7/1/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/07/20 16:20	7/1/20	
Endrin	0.045 U	0.045	0.019	1	07/07/20 16:20	7/1/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/07/20 16:20	7/1/20	
Heptachlor	0.045 U	0.045	0.019	1	07/07/20 16:20	7/1/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/07/20 16:20	7/1/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/07/20 16:20	7/1/20	
Toxaphene	0.46 U	0.46	0.46	1	07/07/20 16:20	7/1/20	
alpha-BHC	0.097	0.045	0.019	1	07/07/20 16:20	7/1/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/07/20 16:20	7/1/20	
beta-BHC	0.045 U	0.045	0.019	1	07/07/20 16:20	7/1/20	
delta-BHC	0.10	0.045	0.019	1	07/07/20 16:20	7/1/20	
gamma-BHC (Lindane)	0.096	0.045	0.019	1	07/07/20 16:20	7/1/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/07/20 16:20	7/1/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	55	10 - 164	07/07/20 16:20	
Tetrachloro-m-xylene	44	10 - 147	07/07/20 16:20	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 11:55

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-017
 Units: ug/L

 Lab Code:
 R2005539-004
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/07/20 16:39	7/1/20	_
4,4'-DDE	0.045 U	0.045	0.019	1	07/07/20 16:39	7/1/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/07/20 16:39	7/1/20	
Aldrin	0.045 U	0.045	0.019	1	07/07/20 16:39	7/1/20	
Dieldrin	0.045 U	0.045	0.019	1	07/07/20 16:39	7/1/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/07/20 16:39	7/1/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/07/20 16:39	7/1/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/07/20 16:39	7/1/20	
Endrin	0.045 U	0.045	0.019	1	07/07/20 16:39	7/1/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/07/20 16:39	7/1/20	
Heptachlor	0.045 U	0.045	0.019	1	07/07/20 16:39	7/1/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/07/20 16:39	7/1/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/07/20 16:39	7/1/20	
Toxaphene	0.46 U	0.46	0.46	1	07/07/20 16:39	7/1/20	
alpha-BHC	0.045 U	0.045	0.019	1	07/07/20 16:39	7/1/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/07/20 16:39	7/1/20	
beta-BHC	0.045 U	0.045	0.019	1	07/07/20 16:39	7/1/20	
delta-BHC	0.045 U	0.045	0.019	1	07/07/20 16:39	7/1/20	
gamma-BHC (Lindane)	0.045 U	0.045	0.019	1	07/07/20 16:39	7/1/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/07/20 16:39	7/1/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	33	10 - 164	07/07/20 16:39	
Tetrachloro-m-xylene	54	10 - 147	07/07/20 16:39	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 12:40

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-018
 Units: ug/L

 Lab Code:
 R2005539-005
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/08/20 18:05	7/1/20	_
4,4'-DDE	0.029 J	0.045	0.019	1	07/08/20 18:05	7/1/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/08/20 18:05	7/1/20	
Aldrin	0.045 U	0.045	0.019	1	07/08/20 18:05	7/1/20	
Dieldrin	0.045 U	0.045	0.019	1	07/08/20 18:05	7/1/20	
Endosulfan I	0.037 J	0.045	0.019	1	07/08/20 18:05	7/1/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/08/20 18:05	7/1/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/08/20 18:05	7/1/20	
Endrin	0.045 U	0.045	0.019	1	07/08/20 18:05	7/1/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/08/20 18:05	7/1/20	
Heptachlor	0.045 U	0.045	0.019	1	07/08/20 18:05	7/1/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/08/20 18:05	7/1/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/08/20 18:05	7/1/20	
Toxaphene	0.46 U	0.46	0.46	1	07/08/20 18:05	7/1/20	
alpha-BHC	0.045 U	0.045	0.019	1	07/08/20 18:05	7/1/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/08/20 18:05	7/1/20	
beta-BHC	0.045 JP	0.045	0.019	1	07/08/20 18:05	7/1/20	
delta-BHC	0.045 U	0.045	0.019	1	07/08/20 18:05	7/1/20	
gamma-BHC (Lindane)	0.026 J	0.045	0.019	1	07/08/20 18:05	7/1/20	
gamma-Chlordane	0.020 JP	0.045	0.019	1	07/08/20 18:05	7/1/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q	
Decachlorobiphenyl	3 *	10 - 164	07/08/20 18:05	*	
Tetrachloro-m-xylene	34	10 - 147	07/08/20 18:05		

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 08:50

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-014
 Units: ug/L

 Lab Code:
 R2005539-001
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/07/20 20:03	7/1/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/07/20 20:03	7/1/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/07/20 20:03	7/1/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/07/20 20:03	7/1/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/07/20 20:03	7/1/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/07/20 20:03	7/1/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/07/20 20:03	7/1/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	42	10 - 152	07/07/20 20:03	
Tetrachloro-m-xylene	44	14 - 129	07/07/20 20:03	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 10:00

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-015
 Units: ug/L

 Lab Code:
 R2005539-002
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/07/20 20:23	7/1/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/07/20 20:23	7/1/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/07/20 20:23	7/1/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/07/20 20:23	7/1/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/07/20 20:23	7/1/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/07/20 20:23	7/1/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/07/20 20:23	7/1/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	54	10 - 152	07/07/20 20:23	
Tetrachloro-m-xylene	43	14 - 129	07/07/20 20:23	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 11:05

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-016
 Units: ug/L

 Lab Code:
 R2005539-003
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/07/20 20:43	7/1/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/07/20 20:43	7/1/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/07/20 20:43	7/1/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/07/20 20:43	7/1/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/07/20 20:43	7/1/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/07/20 20:43	7/1/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/07/20 20:43	7/1/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	64	10 - 152	07/07/20 20:43	
Tetrachloro-m-xvlene	29	14 - 129	07/07/20 20:43	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 11:55

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-017
 Units: ug/L

 Lab Code:
 R2005539-004
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/06/20 15:26	7/1/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/06/20 15:26	7/1/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/06/20 15:26	7/1/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/06/20 15:26	7/1/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/06/20 15:26	7/1/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/06/20 15:26	7/1/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/06/20 15:26	7/1/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	27	10 - 152	07/06/20 15:26	
Tetrachloro-m-xylene	37	14 - 129	07/06/20 15:26	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/26/20 12:40

Monitoring

Sample Matrix: Water Date Received: 06/27/20 08:50

 Sample Name:
 WG-9954-062620-RM-018
 Units: ug/L

 Lab Code:
 R2005539-005
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/06/20 15:45	7/1/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/06/20 15:45	7/1/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/06/20 15:45	7/1/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/06/20 15:45	7/1/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/06/20 15:45	7/1/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/06/20 15:45	7/1/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/06/20 15:45	7/1/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	11	10 - 152	07/06/20 15:45	
Tetrachloro-m-xylene	43	14 - 129	07/06/20 15:45	



QC Summary Forms

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com



Volatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

QA/QC Report

Service Request: R2005539

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

Project:

SURROGATE RECOVERY SUMMARYVolatile Organic Compounds by GC/MS

Analysis Method: 8260C

Extraction Method: EPA 5030C

		4-Bromofluorobenzene	Dibromofluoromethane	Toluene-d8
Sample Name	Lab Code	85-122	89-119	87-121
WG-9954-062620-RM-014	R2005539-001	93	94	94
WG-9954-062620-RM-015	R2005539-002	98	100	98
WG-9954-062620-RM-016	R2005539-003	105	104	104
WG-9954-062620-RM-017	R2005539-004	92	93	94
WG-9954-062620-RM-018	R2005539-005	102	103	103
TB-9954-062620-RM-003	R2005539-006	91	89	93
Method Blank	RQ2007141-04	90	89	91
Method Blank	RQ2007251-04	101	102	100
Lab Control Sample	RQ2007141-03	95	95	93
Lab Control Sample	RQ2007251-03	102	103	101

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007141-04Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	06/30/20 23:08	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	06/30/20 23:08	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	06/30/20 23:08	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	06/30/20 23:08	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	06/30/20 23:08	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	06/30/20 23:08	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	06/30/20 23:08	
2-Butanone (MEK)	10 U	10	0.78	1	06/30/20 23:08	
2-Hexanone	10 U	10	0.20	1	06/30/20 23:08	
4-Methyl-2-pentanone	10 U	10	0.20	1	06/30/20 23:08	
Acetone	10 U	10	5.0	1	06/30/20 23:08	
Benzene	5.0 U	5.0	0.20	1	06/30/20 23:08	
Bromodichloromethane	5.0 U	5.0	0.20	1	06/30/20 23:08	
Bromoform	5.0 U	5.0	0.25	1	06/30/20 23:08	
Bromomethane	5.0 U	5.0	0.70	1	06/30/20 23:08	
Carbon Disulfide	10 U	10	0.42	1	06/30/20 23:08	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	06/30/20 23:08	
Chlorobenzene	5.0 U	5.0	0.20	1	06/30/20 23:08	
Chloroethane	5.0 U	5.0	0.23	1	06/30/20 23:08	
Chloroform	5.0 U	5.0	0.24	1	06/30/20 23:08	
Chloromethane	5.0 U	5.0	0.28	1	06/30/20 23:08	
Dibromochloromethane	5.0 U	5.0	0.20	1	06/30/20 23:08	
Dichloromethane	5.0 U	5.0	0.65	1	06/30/20 23:08	
Ethylbenzene	5.0 U	5.0	0.20	1	06/30/20 23:08	
Styrene	5.0 U	5.0	0.20	1	06/30/20 23:08	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	06/30/20 23:08	
Toluene	5.0 U	5.0	0.20	1	06/30/20 23:08	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	06/30/20 23:08	
Vinyl Acetate	10 U	10	1.1	1	06/30/20 23:08	
Vinyl Chloride	5.0 U	5.0	0.20	1	06/30/20 23:08	
Xylenes, Total	5.0 U	5.0	0.23	1	06/30/20 23:08	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	06/30/20 23:08	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	06/30/20 23:08	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	06/30/20 23:08	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	06/30/20 23:08	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007141-04Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	90	85 - 122	06/30/20 23:08	
Dibromofluoromethane	89	89 - 119	06/30/20 23:08	
Toluene-d8	91	87 - 121	06/30/20 23:08	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007141-04Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Tentatively Identified Compounds

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007251-04Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/07/20 17:23	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/07/20 17:23	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/07/20 17:23	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/07/20 17:23	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/07/20 17:23	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/07/20 17:23	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/07/20 17:23	
2-Butanone (MEK)	10 U	10	0.78	1	07/07/20 17:23	
2-Hexanone	10 U	10	0.20	1	07/07/20 17:23	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/07/20 17:23	
Acetone	10 U	10	5.0	1	07/07/20 17:23	
Benzene	5.0 U	5.0	0.20	1	07/07/20 17:23	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/07/20 17:23	
Bromoform	5.0 U	5.0	0.25	1	07/07/20 17:23	
Bromomethane	5.0 U	5.0	0.70	1	07/07/20 17:23	
Carbon Disulfide	10 U	10	0.42	1	07/07/20 17:23	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/07/20 17:23	
Chlorobenzene	5.0 U	5.0	0.20	1	07/07/20 17:23	
Chloroethane	5.0 U	5.0	0.23	1	07/07/20 17:23	
Chloroform	5.0 U	5.0	0.24	1	07/07/20 17:23	
Chloromethane	5.0 U	5.0	0.28	1	07/07/20 17:23	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/07/20 17:23	
Dichloromethane	5.0 U	5.0	0.65	1	07/07/20 17:23	
Ethylbenzene	5.0 U	5.0	0.20	1	07/07/20 17:23	
Styrene	5.0 U	5.0	0.20	1	07/07/20 17:23	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/07/20 17:23	
Toluene	5.0 U	5.0	0.20	1	07/07/20 17:23	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/07/20 17:23	
Vinyl Acetate	10 U	10	1.1	1	07/07/20 17:23	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/07/20 17:23	
Xylenes, Total	5.0 U	5.0	0.23	1	07/07/20 17:23	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/07/20 17:23	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/07/20 17:23	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/07/20 17:23	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/07/20 17:23	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007251-04Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Surrogate Name % Rec Q **Control Limits Date Analyzed** 4-Bromofluorobenzene 101 85 - 122 07/07/20 17:23 Dibromofluoromethane 102 89 - 119 07/07/20 17:23 100 87 - 121 07/07/20 17:23 Toluene-d8

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007251-04Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Tentatively Identified Compounds

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring **Date Analyzed:** 06/30/20

Sample Matrix: Water

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005539

Lab Control Sample

RQ2007141-03

1,1,1-Trichloroethane (TCA) 8260C 21.6 20.0 1,1,2,2-Tetrachloroethane 8260C 26.2 20.0 1,1,2-Trichloroethane 8260C 22.0 20.0 1,1-Dichloroethane (1,1-DCA) 8260C 20.8 20.0 1,1-Dichloroethane (1,1-DCE) 8260C 20.3 20.0 1,2-Dichloroethane 8260C 19.5 20.0 1,2-Dichloropropane 8260C 21.2 20.0 2-Butanone (MEK) 8260C 16.9 20.0 2-Hexanone 8260C 17.3 20.0 4-Methyl-2-pentanone 8260C 17.2 20.0 Acetone 8260C 21.1 20.0 Benzene 8260C 21.2 20.0 Bromodichloromethane 8260C 21.2 20.0 Bromoform 8260C 21.4 20.0 Bromomethane 8260C 21.4 20.0 Carbon Disulfide 8260C 17.2 20.0 Chloroethane 8260C 21.4 2	ount % Rec	% Rec Limits
1,1,2-Trichloroethane 8260C 22.0 20.0 1,1-Dichloroethane (1,1-DCA) 8260C 20.8 20.0 1,1-Dichloroethene (1,1-DCE) 8260C 20.3 20.0 1,2-Dichloroethane 8260C 19.5 20.0 1,2-Dichloropropane 8260C 21.2 20.0 2-Butanone (MEK) 8260C 16.9 20.0 2-Hexanone 8260C 17.3 20.0 4-Methyl-2-pentanone 8260C 17.2 20.0 Acetone 8260C 20.1 20.0 Benzene 8260C 21.2 20.0 Bromodichloromethane 8260C 21.4 20.0 Bromoform 8260C 21.4 20.0 Bromomethane 8260C 17.2 20.0 Carbon Disulfide 8260C 17.2 20.0 Chlorobenzene 8260C 21.4 20.0 Chloroethane 8260C 21.4 20.0 Chloroform 8260C 21.1 20.0 Chloromethane 8260C 21.1 20.0 Dibromoc	108	75-125
1,1-Dichloroethane (1,1-DCA) 8260C 20.8 20.0 1,1-Dichloroethene (1,1-DCE) 8260C 20.3 20.0 1,2-Dichloroethane 8260C 19.5 20.0 1,2-Dichloropropane 8260C 21.2 20.0 2-Butanone (MEK) 8260C 16.9 20.0 2-Hexanone 8260C 17.3 20.0 4-Methyl-2-pentanone 8260C 17.2 20.0 Acetone 8260C 20.1 20.0 Benzene 8260C 21.2 20.0 Bromodichloromethane 8260C 21.4 20.0 Bromoform 8260C 21.4 20.0 Bromomethane 8260C 14.4 20.0 Carbon Disulfide 8260C 17.2 20.0 Carbon Tetrachloride 8260C 21.4 20.0 Chlorobenzene 8260C 21.4 20.0 Chloroform 8260C 21.1 20.0 Chloromethane 8260C 21.1 20.0 Dibromochloromethane 8260C 23.1 20.0 E	131 *	78-126
1,1-Dichloroethene (1,1-DCE) 8260C 20.3 20.0 1,2-Dichloroethane 8260C 19.5 20.0 1,2-Dichloropropane 8260C 21.2 20.0 2-Butanone (MEK) 8260C 16.9 20.0 2-Hexanone 8260C 17.3 20.0 4-Methyl-2-pentanone 8260C 17.2 20.0 Acetone 8260C 20.1 20.0 Benzene 8260C 21.2 20.0 Bromodichloromethane 8260C 21.4 20.0 Bromoform 8260C 22.5 20.0 Bromomethane 8260C 14.4 20.0 Carbon Disulfide 8260C 17.2 20.0 Carbon Tetrachloride 8260C 21.4 20.0 Chlorobenzene 8260C 21.4 20.0 Chloroform 8260C 21.1 20.0 Chloroform 8260C 21.1 20.0 Chloromethane 8260C 21.1 20.0 Dichloromethane 8260C 23.1 20.0 Styrene 8	110	82-121
1,2-Dichloroethane 8260C 19.5 20.0 1,2-Dichloropropane 8260C 21.2 20.0 2-Butanone (MEK) 8260C 16.9 20.0 2-Hexanone 8260C 17.3 20.0 4-Methyl-2-pentanone 8260C 17.2 20.0 Acetone 8260C 20.1 20.0 Benzene 8260C 21.2 20.0 Bromodichloromethane 8260C 21.4 20.0 Bromoform 8260C 22.5 20.0 Bromomethane 8260C 14.4 20.0 Carbon Disulfide 8260C 17.2 20.0 Chlorobenzene 8260C 21.4 20.0 Chlorobenzene 8260C 21.4 20.0 Chloroform 8260C 21.1 20.0 Chloroform 8260C 21.1 20.0 Chloromethane 8260C 21.1 20.0 Dibromochloromethane 8260C 20.7 20.0 Ethylbenzene	104	80-124
1,2-Dichloropropane 8260C 21.2 20.0 2-Butanone (MEK) 8260C 16.9 20.0 2-Hexanone 8260C 17.3 20.0 4-Methyl-2-pentanone 8260C 17.2 20.0 Acetone 8260C 20.1 20.0 Benzene 8260C 21.2 20.0 Bromodichloromethane 8260C 21.4 20.0 Bromoform 8260C 22.5 20.0 Bromomethane 8260C 14.4 20.0 Carbon Disulfide 8260C 17.2 20.0 Carbon Tetrachloride 8260C 21.4 20.0 Chlorobenzene 8260C 21.4 20.0 Chloroform 8260C 21.9 20.0 Chloroform 8260C 21.1 20.0 Chloromethane 8260C 21.1 20.0 Dibromochloromethane 8260C 23.1 20.0 Ethylbenzene 8260C 22.5 20.0 Styrene	101	71-118
2-Butanone (MEK) 8260C 16.9 20.0 2-Hexanone 8260C 17.3 20.0 4-Methyl-2-pentanone 8260C 17.2 20.0 Acetone 8260C 20.1 20.0 Benzene 8260C 21.2 20.0 Bromodichloromethane 8260C 21.4 20.0 Bromoform 8260C 22.5 20.0 Bromomethane 8260C 14.4 20.0 Carbon Disulfide 8260C 17.2 20.0 Carbon Tetrachloride 8260C 21.4 20.0 Chlorobenzene 8260C 21.4 20.0 Chlorothane 8260C 21.9 20.0 Chloroform 8260C 21.1 20.0 Chloromethane 8260C 21.1 20.0 Dibromochloromethane 8260C 23.1 20.0 Ethylbenzene 8260C 22.5 20.0 Styrene 8260C 21.6 20.0 Toluene 8260C </td <td>98</td> <td>71-127</td>	98	71-127
2-Hexanone 8260C 17.3 20.0 4-Methyl-2-pentanone 8260C 17.2 20.0 Acetone 8260C 20.1 20.0 Benzene 8260C 21.2 20.0 Bromodichloromethane 8260C 21.4 20.0 Bromoform 8260C 22.5 20.0 Bromomethane 8260C 14.4 20.0 Carbon Disulfide 8260C 17.2 20.0 Carbon Tetrachloride 8260C 21.4 20.0 Chlorobenzene 8260C 21.4 20.0 Chlorothane 8260C 21.9 20.0 Chloroform 8260C 21.1 20.0 Chloromethane 8260C 21.1 20.0 Dibrloromethane 8260C 23.1 20.0 Ethylbenzene 8260C 22.5 20.0 Styrene 8260C 22.5 20.0 Tetrachloroethene (PCE) 8260C 21.6 20.0 Trichloroethene (TCE) 8260C 18.3 20.0 Vinyl Acetate 8260C	106	80-119
4-Methyl-2-pentanone 8260C 17.2 20.0 Acetone 8260C 20.1 20.0 Benzene 8260C 21.2 20.0 Bromodichloromethane 8260C 21.4 20.0 Bromoform 8260C 22.5 20.0 Bromomethane 8260C 14.4 20.0 Carbon Disulfide 8260C 17.2 20.0 Carbon Tetrachloride 8260C 21.4 20.0 Chlorobenzene 8260C 21.9 20.0 Chlorothane 8260C 16.2 20.0 Chloroform 8260C 21.1 20.0 Chloromethane 8260C 21.1 20.0 Dibromochloromethane 8260C 23.1 20.0 Ethylbenzene 8260C 22.5 20.0 Styrene 8260C 22.5 20.0 Tetrachloroethene (PCE) 8260C 21.6 20.0 Trichloroethene (TCE) 8260C 18.3 20.0 Vinyl Acetate<	85	61-137
Acetone 8260C 20.1 20.0 Benzene 8260C 21.2 20.0 Bromodichloromethane 8260C 21.4 20.0 Bromoform 8260C 22.5 20.0 Bromomethane 8260C 14.4 20.0 Carbon Disulfide 8260C 17.2 20.0 Carbon Tetrachloride 8260C 21.4 20.0 Chlorobenzene 8260C 21.9 20.0 Chlorotethane 8260C 21.9 20.0 Chloroform 8260C 21.1 20.0 Chloromethane 8260C 21.1 20.0 Dibromochloromethane 8260C 23.1 20.0 Ethylbenzene 8260C 20.7 20.0 Styrene 8260C 22.5 20.0 Tetrachloroethene (PCE) 8260C 21.6 20.0 Trichloroethene (TCE) 8260C 21.5 20.0 Trichloroethene (TCE) 8260C 23.8 20.0 Vinyl Chlor	87	63-124
Benzene 8260C 21.2 20.0 Bromodichloromethane 8260C 21.4 20.0 Bromoform 8260C 22.5 20.0 Bromomethane 8260C 14.4 20.0 Carbon Disulfide 8260C 17.2 20.0 Carbon Tetrachloride 8260C 21.4 20.0 Chlorobenzene 8260C 21.9 20.0 Chlorotethane 8260C 21.9 20.0 Chloroform 8260C 21.1 20.0 Chloromethane 8260C 21.1 20.0 Dibromochloromethane 8260C 23.1 20.0 Ethylbenzene 8260C 20.7 20.0 Styrene 8260C 22.5 20.0 Tetrachloroethene (PCE) 8260C 21.6 20.0 Trichloroethene (TCE) 8260C 21.5 20.0 Vinyl Acetate 8260C 23.8 20.0 Vinyl Chloride 8260C 17.0 20.0 cis-1,2-Dich	86	66-124
Bromodichloromethane 8260C 21.4 20.0 Bromoform 8260C 22.5 20.0 Bromomethane 8260C 14.4 20.0 Carbon Disulfide 8260C 17.2 20.0 Carbon Tetrachloride 8260C 21.4 20.0 Chlorobenzene 8260C 21.9 20.0 Chlorotethane 8260C 16.2 20.0 Chloroform 8260C 21.1 20.0 Chloromethane 8260C 21.1 20.0 Dibromochloromethane 8260C 23.1 20.0 Dichloromethane 8260C 20.7 20.0 Ethylbenzene 8260C 22.5 20.0 Styrene 8260C 22.5 20.0 Tetrachloroethene (PCE) 8260C 21.6 20.0 Trichloroethene (TCE) 8260C 18.3 20.0 Vinyl Acetate 8260C 17.0 20.0 Cis-1,2-Dichloroethene 8260C 22.1 20.0	101	40-161
Bromoform 8260C 22.5 20.0 Bromomethane 8260C 14.4 20.0 Carbon Disulfide 8260C 17.2 20.0 Carbon Tetrachloride 8260C 21.4 20.0 Chlorobenzene 8260C 21.9 20.0 Chloroethane 8260C 16.2 20.0 Chloroform 8260C 21.1 20.0 Chloromethane 8260C 21.1 20.0 Dibromochloromethane 8260C 23.1 20.0 Ethylbenzene 8260C 20.7 20.0 Styrene 8260C 22.5 20.0 Tetrachloroethene (PCE) 8260C 21.6 20.0 Toluene 8260C 21.5 20.0 Vinyl Acetate 8260C 18.3 20.0 Vinyl Chloride 8260C 17.0 20.0 cis-1,2-Dichloroethene 8260C 22.1 20.0	106	79-119
Bromomethane 8260C 14.4 20.0 Carbon Disulfide 8260C 17.2 20.0 Carbon Tetrachloride 8260C 21.4 20.0 Chlorobenzene 8260C 21.9 20.0 Chloroethane 8260C 16.2 20.0 Chloroform 8260C 21.1 20.0 Chloromethane 8260C 17.5 20.0 Dibromochloromethane 8260C 23.1 20.0 Dichloromethane 8260C 20.7 20.0 Ethylbenzene 8260C 22.5 20.0 Styrene 8260C 22.5 20.0 Toluene 8260C 21.6 20.0 Trichloroethene (PCE) 8260C 21.5 20.0 Vinyl Acetate 8260C 18.3 20.0 Vinyl Chloride 8260C 17.0 20.0 cis-1,2-Dichloroethene 8260C 22.1 20.0	107	81-123
Carbon Disulfide 8260C 17.2 20.0 Carbon Tetrachloride 8260C 21.4 20.0 Chlorobenzene 8260C 21.9 20.0 Chloroethane 8260C 16.2 20.0 Chloroform 8260C 21.1 20.0 Chloromethane 8260C 17.5 20.0 Dibromochloromethane 8260C 23.1 20.0 Dichloromethane 8260C 20.7 20.0 Ethylbenzene 8260C 22.5 20.0 Styrene 8260C 22.5 20.0 Toluene 8260C 21.5 20.0 Trichloroethene (PCE) 8260C 21.5 20.0 Vinyl Acetate 8260C 18.3 20.0 Vinyl Chloride 8260C 17.0 20.0 cis-1,2-Dichloroethene 8260C 22.1 20.0	113	65-146
Carbon Tetrachloride 8260C 21.4 20.0 Chlorobenzene 8260C 21.9 20.0 Chloroethane 8260C 16.2 20.0 Chloroform 8260C 21.1 20.0 Chloromethane 8260C 17.5 20.0 Dibromochloromethane 8260C 23.1 20.0 Dichloromethane 8260C 20.7 20.0 Ethylbenzene 8260C 22.5 20.0 Styrene 8260C 22.5 20.0 Tetrachloroethene (PCE) 8260C 21.6 20.0 Toluene 8260C 21.5 20.0 Vinyl Acetate 8260C 18.3 20.0 Vinyl Chloride 8260C 17.0 20.0 Vinyl Chloride 8260C 17.0 20.0 cis-1,2-Dichloroethene 8260C 22.1 20.0	72	42-166
Chlorobenzene 8260C 21.9 20.0 Chloroethane 8260C 16.2 20.0 Chloroform 8260C 21.1 20.0 Chloromethane 8260C 17.5 20.0 Dibromochloromethane 8260C 23.1 20.0 Dichloromethane 8260C 20.7 20.0 Ethylbenzene 8260C 22.5 20.0 Styrene 8260C 22.5 20.0 Tetrachloroethene (PCE) 8260C 21.6 20.0 Toluene 8260C 21.5 20.0 Vinyl Acetate 8260C 18.3 20.0 Vinyl Chloride 8260C 23.8 20.0 Vinyl Chloride 8260C 17.0 20.0 cis-1,2-Dichloroethene 8260C 22.1 20.0	86	66-128
Chloroethane 8260C 16.2 20.0 Chloroform 8260C 21.1 20.0 Chloromethane 8260C 17.5 20.0 Dibromochloromethane 8260C 23.1 20.0 Dichloromethane 8260C 20.7 20.0 Ethylbenzene 8260C 22.5 20.0 Styrene 8260C 22.5 20.0 Tetrachloroethene (PCE) 8260C 21.6 20.0 Toluene 8260C 21.5 20.0 Trichloroethene (TCE) 8260C 18.3 20.0 Vinyl Acetate 8260C 23.8 20.0 Vinyl Chloride 8260C 17.0 20.0 cis-1,2-Dichloroethene 8260C 22.1 20.0	107	70-127
Chloroform 8260C 21.1 20.0 Chloromethane 8260C 17.5 20.0 Dibromochloromethane 8260C 23.1 20.0 Dichloromethane 8260C 20.7 20.0 Ethylbenzene 8260C 22.5 20.0 Styrene 8260C 22.5 20.0 Tetrachloroethene (PCE) 8260C 21.6 20.0 Toluene 8260C 21.5 20.0 Trichloroethene (TCE) 8260C 18.3 20.0 Vinyl Acetate 8260C 23.8 20.0 Vinyl Chloride 8260C 17.0 20.0 cis-1,2-Dichloroethene 8260C 22.1 20.0	109	80-121
Chloromethane 8260C 17.5 20.0 Dibromochloromethane 8260C 23.1 20.0 Dichloromethane 8260C 20.7 20.0 Ethylbenzene 8260C 22.5 20.0 Styrene 8260C 22.5 20.0 Tetrachloroethene (PCE) 8260C 21.6 20.0 Toluene 8260C 21.5 20.0 Trichloroethene (TCE) 8260C 18.3 20.0 Vinyl Acetate 8260C 23.8 20.0 Vinyl Chloride 8260C 17.0 20.0 cis-1,2-Dichloroethene 8260C 22.1 20.0	81	62-131
Dibromochloromethane 8260C 23.1 20.0 Dichloromethane 8260C 20.7 20.0 Ethylbenzene 8260C 22.5 20.0 Styrene 8260C 22.5 20.0 Tetrachloroethene (PCE) 8260C 21.6 20.0 Toluene 8260C 21.5 20.0 Trichloroethene (TCE) 8260C 18.3 20.0 Vinyl Acetate 8260C 23.8 20.0 Vinyl Chloride 8260C 17.0 20.0 cis-1,2-Dichloroethene 8260C 22.1 20.0	105	79-120
Dichloromethane 8260C 20.7 20.0 Ethylbenzene 8260C 22.5 20.0 Styrene 8260C 22.5 20.0 Tetrachloroethene (PCE) 8260C 21.6 20.0 Toluene 8260C 21.5 20.0 Trichloroethene (TCE) 8260C 18.3 20.0 Vinyl Acetate 8260C 23.8 20.0 Vinyl Chloride 8260C 17.0 20.0 cis-1,2-Dichloroethene 8260C 22.1 20.0	87	65-135
Ethylbenzene 8260C 22.5 20.0 Styrene 8260C 22.5 20.0 Tetrachloroethene (PCE) 8260C 21.6 20.0 Toluene 8260C 21.5 20.0 Trichloroethene (TCE) 8260C 18.3 20.0 Vinyl Acetate 8260C 23.8 20.0 Vinyl Chloride 8260C 17.0 20.0 cis-1,2-Dichloroethene 8260C 22.1 20.0	116	72-128
Styrene 8260C 22.5 20.0 Tetrachloroethene (PCE) 8260C 21.6 20.0 Toluene 8260C 21.5 20.0 Trichloroethene (TCE) 8260C 18.3 20.0 Vinyl Acetate 8260C 23.8 20.0 Vinyl Chloride 8260C 17.0 20.0 cis-1,2-Dichloroethene 8260C 22.1 20.0	103	73-122
Tetrachloroethene (PCE) 8260C 21.6 20.0 Toluene 8260C 21.5 20.0 Trichloroethene (TCE) 8260C 18.3 20.0 Vinyl Acetate 8260C 23.8 20.0 Vinyl Chloride 8260C 17.0 20.0 cis-1,2-Dichloroethene 8260C 22.1 20.0	113	76-120
Toluene 8260C 21.5 20.0 Trichloroethene (TCE) 8260C 18.3 20.0 Vinyl Acetate 8260C 23.8 20.0 Vinyl Chloride 8260C 17.0 20.0 cis-1,2-Dichloroethene 8260C 22.1 20.0	112	80-124
Trichloroethene (TCE) 8260C 18.3 20.0 Vinyl Acetate 8260C 23.8 20.0 Vinyl Chloride 8260C 17.0 20.0 cis-1,2-Dichloroethene 8260C 22.1 20.0	108	72-125
Vinyl Acetate 8260C 23.8 20.0 Vinyl Chloride 8260C 17.0 20.0 cis-1,2-Dichloroethene 8260C 22.1 20.0	107	79-119
Vinyl Chloride 8260C 17.0 20.0 cis-1,2-Dichloroethene 8260C 22.1 20.0	91	74-122
cis-1,2-Dichloroethene 8260C 22.1 20.0	119	52-174
	85	74-159
cis 1.2 Dichloromenono 9260C 21.4 20.0	111	80-121
cis-1,3-Dichloropropene 8260C 21.4 20.0	107	77-122
trans-1,2-Dichloroethene 8260C 21.1 20.0	105	73-118
Printed 7/15/2020 12:37:48 PM	Superset Reference:20-00	00554991 rev 00

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix:

Project:

Water

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

> Units:ug/L Basis:NA

Service Request: R2005539

Date Analyzed: 06/30/20

Lab Control Sample

RQ2007141-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
trans-1 3-Dichloropropene	8260C	21.6	20.0	108	71-133

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring **Date Analyzed:** 07/07/20

Sample Matrix: Water

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005539

Lab Control Sample

RQ2007251-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
1,1,1-Trichloroethane (TCA)	8260C	19.1	20.0	95	75-125
1,1,2,2-Tetrachloroethane	8260C	21.7	20.0	108	78-126
1,1,2-Trichloroethane	8260C	20.6	20.0	103	82-121
1,1-Dichloroethane (1,1-DCA)	8260C	18.6	20.0	93	80-124
1,1-Dichloroethene (1,1-DCE)	8260C	22.3	20.0	112	71-118
1,2-Dichloroethane	8260C	19.0	20.0	95	71-127
1,2-Dichloropropane	8260C	18.3	20.0	92	80-119
2-Butanone (MEK)	8260C	21.7	20.0	109	61-137
2-Hexanone	8260C	19.5	20.0	98	63-124
4-Methyl-2-pentanone	8260C	20.4	20.0	102	66-124
Acetone	8260C	26.2	20.0	131	40-161
Benzene	8260C	19.2	20.0	96	79-119
Bromodichloromethane	8260C	19.2	20.0	96	81-123
Bromoform	8260C	21.7	20.0	109	65-146
Bromomethane	8260C	19.2	20.0	96	42-166
Carbon Disulfide	8260C	23.7	20.0	118	66-128
Carbon Tetrachloride	8260C	19.7	20.0	99	70-127
Chlorobenzene	8260C	19.2	20.0	96	80-121
Chloroethane	8260C	16.7	20.0	84	62-131
Chloroform	8260C	20.1	20.0	100	79-120
Chloromethane	8260C	17.5	20.0	88	65-135
Dibromochloromethane	8260C	20.9	20.0	104	72-128
Dichloromethane	8260C	19.6	20.0	98	73-122
Ethylbenzene	8260C	19.2	20.0	96	76-120
Styrene	8260C	19.4	20.0	97	80-124
Tetrachloroethene (PCE)	8260C	18.2	20.0	91	72-125
Toluene	8260C	19.4	20.0	97	79-119
Trichloroethene (TCE)	8260C	18.5	20.0	93	74-122
Vinyl Acetate	8260C	37.2	20.0	186 *	52-174
Vinyl Chloride	8260C	17.2	20.0	86	74-159
cis-1,2-Dichloroethene	8260C	20.8	20.0	104	80-121
cis-1,3-Dichloropropene	8260C	19.1	20.0	96	77-122
trans-1,2-Dichloroethene	8260C	20.5	20.0	102	73-118
Printed 7/15/2020 12:37:49 PM			Supers	et Reference:20-000	00554991 rev 00

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix:

Project:

Water

Service Request: R2005539 **Date Analyzed:** 07/07/20

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

> Units:ug/L Basis:NA

Lab Control Sample

RQ2007251-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
trans-1.3-Dichloropropene	8260C	20.9	20.0	104	71-133



Semivolatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

SURROGATE RECOVERY SUMMARY Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Extraction Method:** EPA 3510C

		2,4,6-Tribromophenol	2-Fluorobiphenyl	2-Fluorophenol
Sample Name Lab Code		35-141	31-118	10-105
WG-9954-062620-RM-014	R2005539-001	84	63	42
WG-9954-062620-RM-015	R2005539-002	86	57	40
WG-9954-062620-RM-016	R2005539-003	84	64	35
WG-9954-062620-RM-017	R2005539-004	83	70	41
WG-9954-062620-RM-018	R2005539-005	86	62	40
Method Blank	RQ2007030-01	83	72	45
Lab Control Sample	RQ2007030-02	85	67	42
Duplicate Lab Control Sample	RQ2007030-03	81	67	40
WG-9954-062620-RM-018 MS	RQ2007030-04	88	78	47
WG-9954-062620-RM-018 DMS	RQ2007030-05	93	78	45

QA/QC Report

Service Request: R2005539

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

SURROGATE RECOVERY SUMMARY Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Extraction Method:** EPA 3510C

		Nitrobenzene-d5	Phenol-d6	p-Terphenyl-d14
Sample Name	Lab Code	31-110	10-107	10-165
WG-9954-062620-RM-014	R2005539-001	60	30	48
WG-9954-062620-RM-015	R2005539-002	56	30	59
WG-9954-062620-RM-016	R2005539-003	57	27	58
WG-9954-062620-RM-017	R2005539-004	68	32	57
WG-9954-062620-RM-018	R2005539-005	60	30	69
Method Blank	RQ2007030-01	72	35	70
Lab Control Sample	RQ2007030-02	70	31	50
Duplicate Lab Control Sample	RQ2007030-03	67	30	58
WG-9954-062620-RM-018 MS	RQ2007030-04	77	37	60
WG-9954-062620-RM-018 DMS	RQ2007030-05	70	34	56

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates) **Service Request:**

R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring **Sample Matrix:**

Date Collected:

06/26/20

Water

Date Received:

06/27/20

Date Analyzed: **Date Extracted:**

07/6/20 07/1/20

ug/L

Duplicate Matrix Spike Summary Semivolatile Organic Compounds by GC/MS

Sample Name: WG-9954-062620-RM-018 **Units:**

Lab Code:

R2005539-005

8270D

Basis: NA

Analysis Method: Prep Method: EPA 3510C

> **Matrix Spike Duplicate Matrix Spike** RQ2007030-04 RQ2007030-05

		IQ2	2007030-04		r	Q2007030-1	33			
	Sample		Spike			Spike		% Rec		RPD
Analyte Name	Result	Result	Amount	% Rec	Result	Amount	% Rec	Limits	RPD	Limit
1,2,4-Trichlorobenzene	9.1 U	45.0	72.7	62	40.7	72.7	56	10-127	10	30
1,2-Dichlorobenzene	9.1 U	40.6	72.7	56	40.0	72.7	55	17-105	2	30
1,3-Dichlorobenzene	9.1 U	39.7	72.7	55	38.5	72.7	53	21-99	4	30
1,4-Dichlorobenzene	9.1 U	40.6	72.7	56	38.5	72.7	53	10-124	6	30
2,4,5-Trichlorophenol	9.1 U	61.9	72.7	85	62.7	72.7	86	48-134	1	30
2,4,6-Trichlorophenol	9.1 U	58.8	72.7	81	61.6	72.7	85	44-135	5	30
2,4-Dichlorophenol	9.1 U	51.1	72.7	70	48.7	72.7	67	40-130	4	30
2,4-Dimethylphenol	9.1 U	56.0	72.7	77	55.3	72.7	76	42-121	1	30
2,4-Dinitrophenol	45 U	45.6	72.7	63	47.7	72.7	66	21-168	5	30
2,4-Dinitrotoluene	9.1 U	62.8	72.7	86	67.4	72.7	93	37-143	8	30
2,6-Dinitrotoluene	9.1 U	73.8	72.7	101	75.9	72.7	104	39-136	3	30
2-Chloronaphthalene	9.1 U	59.4	72.7	82	60.0	72.7	83	40-108	1	30
2-Chlorophenol	9.1 U	43.8	72.7	60	42.5	72.7	58	37-112	3	30
2-Methylnaphthalene	9.1 U	56.4	72.7	78	52.4	72.7	72	34-102	8	30
2-Methylphenol	9.1 U	50.7	72.7	70	50.3	72.7	69	37-102	1	30
2-Nitroaniline	9.1 U	71.3	72.7	98	75.8	72.7	104	40-136	6	30
2-Nitrophenol	9.1 U	54.3	72.7	75	50.6	72.7	70	27-143	7	30
3,3'-Dichlorobenzidine	9.1 U	62.3	72.7	86	62.5	72.7	86	11-131	<1	30
3- and 4-Methylphenol Coelution	9.1 U	48.5	72.7	67	45.1	72.7	62	30-95	8	30
3-Nitroaniline	9.1 U	66.2	72.7	91	59.8	72.7	82	19-117	10	30
4,6-Dinitro-2-methylphenol	45 U	56.5	72.7	78	56.9	72.7	78	25-154	<1	30
4-Bromophenyl Phenyl Ether	9.1 U	62.4	72.7	86	61.0	72.7	84	39-115	2	30
4-Chloro-3-methylphenol	9.1 U	61.5	72.7	85	60.9	72.7	84	41-126	1	30
4-Chloroaniline	9.1 U	59.7	72.7	82	53.1	72.7	73	19-111	12	30
4-Chlorophenyl Phenyl Ether	9.1 U	53.9	72.7	74	55.7	72.7	77	41-111	4	30
4-Nitroaniline	9.1 U	62.9	72.7	86	62.9	72.7	87	18-143	1	30
4-Nitrophenol	45 U	24.9 J	72.7	34	25.6 J	72.7	35	10-126	3	30
Acenaphthene	9.1 U	61.7	72.7	85	60.3	72.7	83	43-117	2	30
Acenaphthylene	9.1 U	66.7	72.7	92	65.2	72.7	90	45-119	2	30
Anthracene	9.1 U	66.4	72.7	91	65.3	72.7	90	45-127	1	30
Benz(a)anthracene	9.1 U	56.5	72.7	78	53.9	72.7	74	46-126	5	30
Benzo(a)pyrene	9.1 U	60.4	72.7	83	56.3	72.7	77	44-114	8	30
Benzo(b)fluoranthene	9.1 U	52.2	72.7	72	49.1	72.7	68	41-127	6	30

Results flagged with an asterisk (*) indicate values outside control criteria.

Results flagged with a pound (#) indicate the control criteria is not applicable.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

Matrix Spike and Matrix Spike Duplicate Data is presented for information purposes only. The matrix may or may not be relevant to samples reported in this report. The laboratory evaluates system performance based on the LCS and LCSD control limits.

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Date Collected: 06/26/20 **Date Received:** 06/27/20

Sample Matrix: Water

Date Analyzed: 07/6/20

Date Extracted: 07/1/20

Duplicate Matrix Spike

Duplicate Matrix Spike Summary Semivolatile Organic Compounds by GC/MS

Matrix Spike

Sample Name: WG-9954-062620-RM-018

Units: ug/L

Lab Code: R2005539-005

Analysis Method: 8270D

Basis: NA

Prep Method: EPA 3510C

		RQ2	007030-04		R	Q2007030-0)5			
	Sample		Spike			Spike		% Rec		RPD
Analyte Name	Result	Result	Amount	% Rec	Result	Amount	% Rec	Limits	RPD	Limit
Benzo(g,h,i)perylene	9.1 U	63.5	72.7	87	57.5	72.7	79	50-143	10	30
Benzo(k)fluoranthene	9.1 U	60.4	72.7	83	56.3	72.7	77	46-139	8	30
Benzoic Acid	91 U	71.8 J	109	66	73.9 J	109	68	10-94	3	30
Benzyl Alcohol	9.1 U	58.1	72.7	80	55.0	72.7	76	31-109	5	30
2,2'-Oxybis(1-chloropropane)	9.1 U	51.9	72.7	71	49.1	72.7	68	21-126	4	30
Bis(2-chloroethoxy)methane	9.1 U	58.3	72.7	80	56.3	72.7	77	41-118	4	30
Bis(2-chloroethyl) Ether	9.1 U	49.7	72.7	68	45.4	72.7	62	33-108	9	30
Bis(2-ethylhexyl) Phthalate	9.1 U	55.0	72.7	76	51.7	72.7	71	41-132	7	30
Butyl Benzyl Phthalate	9.1 U	61.6	72.7	85	59.8	72.7	82	41-148	4	30
Chrysene	9.1 U	59.4	72.7	82	56.7	72.7	78	47-126	5	30
Di-n-butyl Phthalate	9.1 U	71.7	72.7	99	70.8	72.7	97	43-130	2	30
Di-n-octyl Phthalate	9.1 U	54.1	72.7	74	50.6	72.7	70	40-139	6	30
Dibenz(a,h)anthracene	9.1 U	67.7	72.7	93	59.3	72.7	82	43-136	13	30
Dibenzofuran	9.1 U	65.4	72.7	90	64.7	72.7	89	46-119	1	30
Diethyl Phthalate	9.1 U	58.0	72.7	80	59.1	72.7	81	36-122	1	30
Dimethyl Phthalate	9.1 U	67.1	72.7	92	69.7	72.7	96	33-123	4	30
Fluoranthene	9.1 U	73.6	72.7	101	71.5	72.7	98	43-135	3	30
Fluorene	9.1 U	64.5	72.7	89	65.7	72.7	90	43-113	1	30
Hexachlorobenzene	9.1 U	68.2	72.7	94	66.7	72.7	92	42-125	2	30
Hexachlorobutadiene	9.1 U	47.4	72.7	65	42.2	72.7	58	10-111	11	30
Hexachlorocyclopentadiene	9.1 U	9.50	72.7	13	10.2	72.7	14	10-103	7	30
Hexachloroethane	9.1 U	39.9	72.7	55	37.9	72.7	52	12-101	6	30
Indeno(1,2,3-cd)pyrene	9.1 U	54.2	72.7	75	48.5	72.7	67	49-140	11	30
Isophorone	9.1 U	50.7	72.7	70	49.0	72.7	67	40-111	4	30
N-Nitrosodi-n-propylamine	9.1 U	63.9	72.7	88	59.6	72.7	82	35-108	7	30
N-Nitrosodiphenylamine	9.1 U	81.0	72.7	111	78.4	72.7	108	43-127	3	30
Naphthalene	9.1 U	52.9	72.7	73	48.4	72.7	67	37-108	9	30
Nitrobenzene	9.1 U	53.8	72.7	74	48.7	72.7	67	35-112	10	30
Pentachlorophenol (PCP)	45 U	63.5	72.7	87	63.9	72.7	88	29-164	1	30
Phenanthrene	9.1 U	64.6	72.7	89	64.7	72.7	89	46-123	<1	30
Phenol	9.1 U	29.2	72.7	40	28.4	72.7	39	10-113	3	30
Pyrene	9.1 U	69.5	72.7	96	68.9	72.7	95	44-129	1	30

Results flagged with an asterisk (*) indicate values outside control criteria.

Results flagged with a pound (#) indicate the control criteria is not applicable.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

Matrix Spike and Matrix Spike Duplicate Data is presented for information purposes only. The matrix may or may not be relevant to samples reported in this report. The laboratory evaluates system performance based on the LCS and LCSD control limits.

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007030-01
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

1,24-Trichlorobenzene	Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1.3-Dichlorobenzene	1,2,4-Trichlorobenzene	10 U	10	1.2	1	07/06/20 09:10	7/1/20	
1.4-Dichlorobenzene	1,2-Dichlorobenzene	10 U	10	1.2	1	07/06/20 09:10	7/1/20	
2.4.5-Trichlorophenol 10 U 10	1,3-Dichlorobenzene	10 U	10	1.1	1	07/06/20 09:10	7/1/20	
2.4Trichlorophenol	1,4-Dichlorobenzene	10 U	10	1.2	1	07/06/20 09:10	7/1/20	
2.4-Dichlorophenol 10 U 10	2,4,5-Trichlorophenol	10 U	10	1.1	1	07/06/20 09:10	7/1/20	
2.4-Dimethylphenol 10 U 10 1.4 1 07/06/20 09:10 7/1/20 2.4-Dimitrophenol 50 U 50 20 1 07/06/20 09:10 7/1/20 2.4-Dimitrotoluene 10 U 10 2.4 1 07/06/20 09:10 7/1/20 2.6-Dimitrotoluene 10 U 10 1.4 1 07/06/20 09:10 7/1/20 2-Chloropaphthalene 10 U 10 1.4 1 07/06/20 09:10 7/1/20 2-Methylaphthalene 10 U 10 1.3 1 07/06/20 09:10 7/1/20 2-Methylaphenol 10 U 10 1.0 1 07/06/20 09:10 7/1/20 2-Mitrophenol 10 U 10 1.6 1.4 1 07/06/20 09:10 7/1/20 2-Nitrophenol 10 U 10 1.5 1 07/06/20 09:10 7/1/20 3-and 4-Methylphenol Coelution 10 U 10 1.2 1 07/06/20 09:10 7/1/20 3-Nitroaniline 10 U 10 <td< td=""><td>2,4,6-Trichlorophenol</td><td>10 U</td><td>10</td><td>1.4</td><td>1</td><td>07/06/20 09:10</td><td>7/1/20</td><td></td></td<>	2,4,6-Trichlorophenol	10 U	10	1.4	1	07/06/20 09:10	7/1/20	
2.4-Dimitrophenol 50 U 50 U 20 U 1 07/06/20 09:10 7/1/20 7/1/20 2.4-Dimitrotoluene 10 U 10 U 10 U 10 U 10 U 7/1/20 2.6-Dimitrotoluene 10 U 10 U 10 U 10 U 10 U 10 U 7/1/20 2-Chlorophenol 10 U 10 U 11 U 10 U <t< td=""><td>2,4-Dichlorophenol</td><td>10 U</td><td>10</td><td>1.3</td><td>1</td><td>07/06/20 09:10</td><td>7/1/20</td><td></td></t<>	2,4-Dichlorophenol	10 U	10	1.3	1	07/06/20 09:10	7/1/20	
2.4-Dinitrotoluene 10 U 10 2.4 1 07/06/20 09:10 7/1/20 2.6-Dinitrotoluene 10 U 10 1.4 1 07/06/20 09:10 7/1/20 2-Chlorophenol 10 U 10 1.4 1 07/06/20 09:10 7/1/20 2-Methylmaphthalene 10 U 10 1.3 1 07/06/20 09:10 7/1/20 2-Methylphenol 10 U 10 1.3 1 07/06/20 09:10 7/1/20 2-Nitrophinol 10 U 10 1.4 1 07/06/20 09:10 7/1/20 2-Nitrophinol 10 U 10 1.5 1 07/06/20 09:10 7/1/20 3-Shitroaniline 10 U 10 1.5 1 07/06/20 09:10 7/1/20 3-Shitroaniline 10 U 10 1.2 1 07/06/20 09:10 7/1/20 3-Silchorobenzidine 10 U 10 1.2 1 07/06/20 09:10 7/1/20 3-Silchorobenzidine 10 U 10 1.2 1	2,4-Dimethylphenol	10 U	10	1.4	1	07/06/20 09:10	7/1/20	
2.6-Dinitrotoluene 10 U 10 U 10 U 10 U 10 U 71/20 2-Chloronaphthalene 10 U 10 U 10 U 11 U 10 U/20 71/20 2-Chlorophenol 10 U 10 U 10 U 11 U/20 71/20 2-Methylphenol 10 U 10 U 10 U 10 U/20 71/20 2-Methylphenol 10 U 10 U 10 U 10 U/20 71/20 2-Nitroaniline 10 U 10 U 10 U/20 10 U/20 71/20 2-Nitrophenol 10 U 10 U/20 10 U/20 71/20 71/20 3-3'-Dichlorobenzidine 10 U 10 U/20 10 T/06/20 09:10 T/1/20 71/20 4-Methylphenol 10 U 10 U/20 10 T/06/20 09:10 T/1/20 71/20 4-G-Dinitro-2-methylphenol 50 U 50 U/20 10 T/06/20 09:10 T/1/20 4-Bromophenyl Phenyl Ether 10 U 10 U/20 10 T/06/20 09:10 T/1/20 4-Chloro-3-methylphenol 10 U 10 U/20 10 T/06/20 09:10 T/1/20 4-Chlorophenyl Phenyl Ether <td>2,4-Dinitrophenol</td> <td>50 U</td> <td>50</td> <td>20</td> <td>1</td> <td>07/06/20 09:10</td> <td>7/1/20</td> <td></td>	2,4-Dinitrophenol	50 U	50	20	1	07/06/20 09:10	7/1/20	
2-Chloronaphthalene 10 U 10 U 10 U 11 OU 11 OU 11 OU 711/20 2-Chlorophenol 10 U 10 U 10 U 10 U 10 OU 71/120 2-Methylaphthalene 10 U 10 U 10 U 10 OU 10	2,4-Dinitrotoluene	10 U	10	2.4	1	07/06/20 09:10	7/1/20	
2-Chlorophenol		10 U	10	1.4	1	07/06/20 09:10	7/1/20	
2-Chlorophenol 10 U 10 1.1 1 07/06/20 09:10 7/1/20 2-Methylnaphthalene 10 U 10 1.3 1 07/06/20 09:10 7/1/20 2-Methylphenol 10 U 10 1.4 1 07/06/20 09:10 7/1/20 2-Nitroaniline 10 U 10 1.4 1 07/06/20 09:10 7/1/20 2-Nitrophenol 10 U 10 1.5 1 07/06/20 09:10 7/1/20 2-Nitrophenol 10 U 10 1.5 1 07/06/20 09:10 7/1/20 3-37-Dichlorobenzidine 10 U 10 1.5 1 07/06/20 09:10 7/1/20 3-and 4-Methylphenol Coelution 10 U 10 1.2 1 07/06/20 09:10 7/1/20 3-Aitroaniline 10 U 10 1.2 1 07/06/20 09:10 7/1/20 3-Nitroaniline 10 U 10 1.2 1 07/06/20 09:10 7/1/20 3-Nitroaniline 10 U 10 1.2 1 07/06/20 09:10 7/1/20 4-Bromophenyl Phenyl Ether 10 U 10 1.7 1 07/06/20 09:10 7/1/20 4-Bromophenyl Phenyl Ether 10 U 10 1.7 1 07/06/20 09:10 7/1/20 4-Chloro-3-methylphenol 10 U 10 1.1 1 07/06/20 09:10 7/1/20 4-Chloro-3-methylphenol 10 U 10 1.1 1 07/06/20 09:10 7/1/20 4-Chlorophenyl Phenyl Ether 10 U 10 1.1 1 07/06/20 09:10 7/1/20 4-Chlorophenyl Phenyl Ether 10 U 10 1.5 1 07/06/20 09:10 7/1/20 4-Chlorophenyl Phenyl Ether 10 U 10 1.5 1 07/06/20 09:10 7/1/20 4-Chlorophenyl Phenyl Ether 10 U 10 1.5 1 07/06/20 09:10 7/1/20 4-Nitroaniline 10 U 10 1.5 1 07/06/20 09:10 7/1/20 4-Nitroaniline 10 U 10 1.5 1 07/06/20 09:10 7/1/20 4-Nitroaniline 10 U 10 1.4 1 07/06/20 09:10 7/1/20 Acenaphthene 10 U 10 1.4 1 07/06/20 09:10 7/1/20 Acenaphthylene 10 U 10 1.4 1 07/06/20 09:10 7/1/20 Benzo(a)myrene 10 U 10 1.4 1 07/06/20 09:10 7/1/20 Benzo(a)myrene 10 U 10 1.6 1 07/06/20 09:10 7/1/20 Benzo(a)myrene 10 U 10 1.6 1 07/06/20 09:10 7/1/20 Benzo(a)myrene 10 U 10 1.3 1 07/06/20 09:10 7/1/20 Benzo(b)fluoranthene 10 U 10 1.3 1 07/06/20 09:10 7/1/20 Benzo(b)fluoranthene 10 U 10 1.4 1 07/06/20 09:10 7/1/20 Benzo(c)myrene 10 U 10 1.4 1 07/06/20 09:10 7/1/20 Benzo(c)myrene 10 U 10 1.4 1 07/06/20 09:10 7/1/20 Benzo(c)myrene 10 U 10 1.4 1 07/06/20 09:10 7/1/20 Benzo(c)myrene 10 U 10 1.4 1 07/06/20 09:10 7/1/20 Benzo(c)myrene 10 U 10 1.4 1 07/06/20 09:10 7/1/20 Benzo(c)myrene 10 U 10 1.4 1 07/06/20 09:10 7/1/20 Benzo(c)myrene 10 U 10 1.4 1 07/06/20 09:10 7/1/20 Benzo(c)myrene 10 U 10 1.4 1	2-Chloronaphthalene	10 U	10	1.4	1	07/06/20 09:10	7/1/20	
2-Methylnaphthalene		10 U	10	1.1	1	07/06/20 09:10	7/1/20	
2-Methylphenol		10 U	10	1.3	1	07/06/20 09:10	7/1/20	
2-Nitroaniline		10 U		1.0	1		7/1/20	
2-Nitrophenol 10 U 10 1.5 1 07/06/20 09:10 7/1/20 3,3"-Dichlorobenzidine 10 U 10 1.2 1 07/06/20 09:10 7/1/20 3-and 4-Methylphenol Coelution 10 U 10 1.2 1 07/06/20 09:10 7/1/20 3-Nitroaniline 10 U 10 2.5 1 07/06/20 09:10 7/1/20 4-Dinitro-2-methylphenol 50 U 50 20 1 07/06/20 09:10 7/1/20 4-Bromophenyl Phenyl Ether 10 U 10 1.7 1 07/06/20 09:10 7/1/20 4-Chloro-3-methylphenol 10 U 10 1.7 1 07/06/20 09:10 7/1/20 4-Chloro-3-methylphenol 10 U 10 1.1 1 07/06/20 09:10 7/1/20 4-Chloro-3-methylphenol 10 U 10 1.5 1 07/06/20 09:10 7/1/20 4-Chloro-3-methylphenol 10 U 10 1.5 1 07/06/20 09:10 7/1/20 4-Chloro-3-methylphenol 10 U 10 1.5 1 07/06/20 09:10 7/1/20 4-Chloro-3-methylphenol 10 U 10 1.5 1 07/06/20 09:10 7/1/20 4-Chloro-benyl Phenyl Ether 10 U 10 1.5 1 07/06/20 09:10 7/1/20 4-Nitroaniline 10 U 10 1.5 1 07/06/20 09:10 7/1/20 4-Nitrophenol 50 U 50 6.4 1 07/06/20 09:10 7/1/20 4-Nitrophenol 50 U 50 6.4 1 07/06/20 09:10 7/1/20 4-Nitrophenol 10 U 10 1.4 1 07/06/20 09:10 7/1/20 4-Nitrophenol 10 U 10 1.4 1 07/06/20 09:10 7/1/20 4-Nitrophenol 10 U 10 1.4 1 07/06/20 09:10 7/1/20 4-Nitrophenol 10 U 10 1.4 1 07/06/20 09:10 7/1/20 4-Nitrophenol 10 U 10 1.4 1 07/06/20 09:10 7/1/20 4-Nitrophenol 10 U 10 1.5 1 07/06/20 09:10 7/1/20 4-Nitrophenol 10 U 10 1.5 1 07/06/20 09:10 7/1/20 4-Nitrophenol 10 U 10 1.5 1 07/06/20 09:10 7/1/20 4-Nitrophenol 10 U 10 1.5 1 07/06/20 09:10 7/1/20 4-Nitrophenol 10 U 10 1.5 1 07/06/20 09:10 7/1/20 4-Nitrophenol 10 U 10 1.5 1 07/06/20 09:10 7/1/20 4-Nitrophenol 10 U 10 1.5 1 07/06/20 09:10 7/1/20 4-Nitrophenol 10 U 10 1.5 1 07/06/20 09:10 7/1/20 4-Nitrophenol 10 U 10 1.5 1 07/06/20 09:10 7/1/20	2-Nitroaniline							
3,3'-Dichlorobenzidine		10 U	10	1.5	1	07/06/20 09:10	7/1/20	
3- and 4-Methylphenol Coelution 10 U 10 1.2 1 07/06/20 09:10 7/1/20 3-Nitroaniline 10 U 10 2.5 1 07/06/20 09:10 7/1/20 4.6-Dinitro-2-methylphenol 50 U 50 20 1 07/06/20 09:10 7/1/20 4-Bromophenyl Phenyl Ether 10 U 10 1.7 1 07/06/20 09:10 7/1/20 4-Chloro-3-methylphenol 10 U 10 1.1 1 07/06/20 09:10 7/1/20 4-Chloroaniline 10 U 10 1.1 1 07/06/20 09:10 7/1/20 4-Chlorophenyl Phenyl Ether 10 U 10 1.5 1 07/06/20 09:10 7/1/20 4-Chlorophenyl Phenyl Ether 10 U 10 1.5 1 07/06/20 09:10 7/1/20 4-Chlorophenyl Phenyl Ether 10 U 10 1.5 1 07/06/20 09:10 7/1/20 4-Nitroaniline 10 U 10 2.7 1 07/06/20 09:10 7/1/20 4-Nitroaniline 10 U 10 2.7 1 07/06/20 09:10 7/1/20 4-Nitroaniline 10 U 10 1.4 1 07/06/20 09:10 7/1/20 4-Nitroaniline 10 U 10 1.4 1 07/06/20 09:10 7/1/20 4-Nitroaniline 10 U 10 1.4 1 07/06/20 09:10 7/1/20 4-Nitroaniline 10 U 10 1.4 1 07/06/20 09:10 7/1/20 4-Nitroaniline 10 U 10 1.4 1 07/06/20 09:10 7/1/20 4-Nitroaniline 10 U 10 1.4 1 07/06/20 09:10 7/1/20 4-Nitroaniline 10 U 10 1.4 1 07/06/20 09:10 7/1/20 4-Nitroaniline 10 U 10 1.3 1 07/06/20 09:10 7/1/20 4-Nitroaniline 10 U 10 1.6 1 07/06/20 09:10 7/1/20 4-Nitroaniline 10 U 10 1.2 1 07/06/20 09:10 7/1/20 4-Nitroaniline 10 U 10 1.2 1 07/06/20 09:10 7/1/20 4-Nitroaniline 10 U 10 1.3 1 07/06/20 09:10 7/1/20 4-Nitroaniline 10 U 10 1.3 1 07/06/20 09:10 7/1/20 4-Nitroaniline 10 U 10 1.4 1 07/06/20 09:10 7/1/20 4-Nitroaniline 10 U 10 1.4 1 07/06/20 09:10 7/1/20 4-Nitroaniline 10 U 10 1.4 1 07/06/20 09:10 7/1/20 4-Nitroaniline 10 U 10 1.4 1 07/06/20 09:10 7/1/20 4-Nitroaniline 10 U 10 1.4 1 07/06/20 09:10 7/1/20 4-Nitroaniline 10 U 10 1.4 1 07/06/20 09:10 7/1/20 4-Nit		10 U	10	1.2	1		7/1/20	
3-Nitroaniline 10 U 10 2.5 1 07/06/20 09:10 7/1/20								
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4-Bromophenyl Phenyl Ether 10 U 10 1.7 1 07/06/20 09:10 7/1/20 4-Chloro-3-methylphenol 10 U 10 1.1 1 07/06/20 09:10 7/1/20 4-Chloroaniline 10 U 10 1.0 1 07/06/20 09:10 7/1/20 4-Chlorophenyl Phenyl Ether 10 U 10 1.5 1 07/06/20 09:10 7/1/20 4-Nitrophenol 50 U 50 6.4 1 07/06/20 09:10 7/1/20 4-Nitrophenol 50 U 50 6.4 1 07/06/20 09:10 7/1/20 4-Chlorophenyl Phenyl Ether 10 U 10 1.4 1 07/06/20 09:10 7/1/20 4-Nitrophenol 50 U 50 6.4 1 07/06/20 09:10 7/1/20 4-Nitrophenol 50 U 50 6.4 1 07/06/20 09:10 7/1/20 Accapaththene 10 U 10 1.4 1 07/06/20 09:10 7/1/20 Benzo(a)ptrace 10 U 10 1.6		50 U	50					
4-Chloro-3-methylphenol 10 U 10 U 10 U 1.1 1 07/06/20 09:10 7/1/20 4-Chloroaniline 10 U 10 U 10 U 1.0 1 07/06/20 09:10 7/1/20 4-Chlorophenyl Phenyl Ether 10 U 10 U 1.5 1 07/06/20 09:10 7/1/20 4-Nitrophenol 50 U 50 G.4 1 07/06/20 09:10 7/1/20 4-Chlorophenol 50 U 50 G.4 1 07/06/20 09:10 7/1/20 4-Chlorophenol 10 U 10 I.4 1 07/06/20 09:10 7/1/20 4-Chlorophenol 50 U 50 G.4 1 07/06/20 09:10 7/1/20 4-Nitrophenol 10 U 10 I.4 1 07/06/20 09:10 7/1/20 Acenaphthene 10 U 10 I.4 1 07/06/20 09:10 7/1/20 Acenaphthylene 10 U 10 I.3 1 07/06/20 09:10 7/1/20 Benz(a)anthracene 10 U 10 I.5 1 07/06/20 09:10 7/1/20 Benz(a)aphyrene 10 U 10 I.5 1 07/06/20 09:10 7/1/20 Benzo(b)fluoranthene 1		10 U	10	1.7	1	07/06/20 09:10	7/1/20	
4-Chloroaniline 10 U 10 10 1.0 1 07/06/20 09:10 7/1/20 4-Chlorophenyl Phenyl Ether 10 U 10 1.5 1 07/06/20 09:10 7/1/20 4-Nitroaniline 10 U 10 U 2.7 1 07/06/20 09:10 7/1/20 4-Nitrophenol 50 U 50 G.4 1 07/06/20 09:10 7/1/20 Acenaphthene 10 U 10 II 1.4 1 07/06/20 09:10 7/1/20 Acenaphthylene 10 U 10 II 1.4 1 07/06/20 09:10 7/1/20 Acenaphthylene 10 U 10 II 1.4 1 07/06/20 09:10 7/1/20 Acenaphthylene 10 U 10 II 1.4 1 07/06/20 09:10 7/1/20 Acenaphthylene 10 U 10 II 1.3 1 07/06/20 09:10 7/1/20 Acenaphthylene 10 U 10 II 1.3 1 07/06/20 09:10 7/1/20 Benzo(a)anthracene 10 U 10 II 1.2 1 07/06/20 09:10 7/1/20 Benzo(b)fluoranthene 10 U 10 II 1.2 1 0		10 U	10	1.1	1	07/06/20 09:10	7/1/20	
4-Nitroaniline 10 U 10 U 2.7 1 07/06/20 09:10 7/1/20 4-Nitrophenol 50 U 50 G.4 1 07/06/20 09:10 7/1/20 Acenaphthene 10 U 10 I.4 1 07/06/20 09:10 7/1/20 Acenaphthylene 10 U 10 I.4 1 07/06/20 09:10 7/1/20 Anthracene 10 U 10 I.3 1 07/06/20 09:10 7/1/20 Benz(a)anthracene 10 U 10 I.6 1 07/06/20 09:10 7/1/20 Benzo(a)pyrene 10 U 10 I.2 1 07/06/20 09:10 7/1/20 Benzo(b)fluoranthene 10 U 10 I.2 1 07/06/20 09:10 7/1/20 Benzo(g,h,i)perylene 10 U 10 I.3 1 07/06/20 09:10 7/1/20 Benzo(k)fluoranthene 10 U 10 I.3 1 07/06/20 09:10 7/1/20 Benzoic Acid 10 U 10 I.3 1 07/06/20 09:10 7/1/20 Benzyl Alcohol 10 U 10 I.6 1 07/06/20 09:10 7/1/20 2,2'-Oxybis(1-chloropropane) 10 U 10 I.4 1 07/06/20 09:10 7/1/20		10 U	10	1.0	1		7/1/20	
4-Nitroaniline 10 U 10 U 2.7 1 07/06/20 09:10 7/1/20 4-Nitrophenol 50 U 50 G.4 1 07/06/20 09:10 7/1/20 Acenaphthene 10 U 10 I.4 1 07/06/20 09:10 7/1/20 Acenaphthylene 10 U 10 I.4 1 07/06/20 09:10 7/1/20 Anthracene 10 U 10 I.3 1 07/06/20 09:10 7/1/20 Benz(a)anthracene 10 U 10 I.6 1 07/06/20 09:10 7/1/20 Benzo(a)pyrene 10 U 10 I.2 1 07/06/20 09:10 7/1/20 Benzo(b)fluoranthene 10 U 10 I.2 1 07/06/20 09:10 7/1/20 Benzo(g,h,i)perylene 10 U 10 I.3 1 07/06/20 09:10 7/1/20 Benzo(k)fluoranthene 10 U 10 I.3 1 07/06/20 09:10 7/1/20 Benzoic Acid 10 U 10 I.3 1 07/06/20 09:10 7/1/20 Benzyl Alcohol 10 U 10 I.6 1 07/06/20 09:10 7/1/20 2,2'-Oxybis(1-chloropropane) 10 U 10 I.4 1 07/06/20 09:10 7/1/20	4-Chlorophenyl Phenyl Ether	10 U	10	1.5	1	07/06/20 09:10	7/1/20	
4-Nitrophenol 50 U 50 d 6.4 1 07/06/20 09:10 7/1/20 Acenaphthene 10 U		10 U		2.7	1			
Acenaphthene 10 U 10 U 10 U 1.4 1 07/06/20 09:10 7/1/20 Acenaphthylene 10 U 10 U 10 U 1.4 1 07/06/20 09:10 7/1/20 Anthracene 10 U 10 U 10 U 1.3 1 07/06/20 09:10 7/1/20 Benzo(a)anthracene 10 U 10 U 10 U 1.6 1 07/06/20 09:10 7/1/20 Benzo(a)pyrene 10 U 10 U 10 U 1.2 1 07/06/20 09:10 7/1/20 Benzo(b)fluoranthene 10 U 10 U 10 U 1.0 1 07/06/20 09:10 7/1/20 Benzo(k)fluoranthene 10 U 10 U 10 U 1.0 1 07/06/20 09:10 7/1/20 Benzoic Acid 100 U 100 U 36 I 1 07/06/20 09:10 7/1/20 Benzyl Alcohol 10 U 10 U 1.6 1 07/06/20 09:10 7/1/20 Bis(2-chloroethoxy)methane 10 U 10 U 1.4 1 07/06/20 09:10 7/1/20 Bis(2-chloroethyl) Ether 10 U 10 U 1.3 1 07/06/20 09:10 7/1/20 Butyl Benzyl Phthalate 10 U 10	4-Nitrophenol	50 U	50	6.4	1		7/1/20	
Acenaphthylene 10 U		10 U	10	1.4	1	07/06/20 09:10	7/1/20	
Anthracene 10 U		10 U	10	1.4	1	07/06/20 09:10	7/1/20	
Benz(a)anthracene 10 U 10 U <td></td> <td>10 U</td> <td>10</td> <td>1.3</td> <td>1</td> <td>07/06/20 09:10</td> <td>7/1/20</td> <td></td>		10 U	10	1.3	1	07/06/20 09:10	7/1/20	
Benzo(b)fluoranthene 10 U 10 U<		10 U	10	1.6	1		7/1/20	
Benzo(b)fluoranthene 10 U 10 U<	* *	10 U	10	1.2	1		7/1/20	
Benzo(g,h,i)perylene 10 U 10 U<		10 U				07/06/20 09:10	7/1/20	
Benzo(k)fluoranthene 10 U 10 U<			10		1			
Benzoic Acid 100 U 100 36 1 07/06/20 09:10 7/1/20 Benzyl Alcohol 10 U		10 U	10	1.3	1	07/06/20 09:10	7/1/20	
Benzyl Alcohol 10 U								
2,2'-Oxybis(1-chloropropane) 10 U 10 U 1.4 1 07/06/20 09:10 7/1/20 Bis(2-chloroethoxy)methane 10 U 10 U 1.9 1 07/06/20 09:10 7/1/20 Bis(2-chloroethyl) Ether 10 U 10 U 1.3 1 07/06/20 09:10 7/1/20 Bis(2-ethylhexyl) Phthalate 10 U 10 U 1.0 1 07/06/20 09:10 7/1/20 Butyl Benzyl Phthalate 10 U 10 U 1.4 1 07/06/20 09:10 7/1/20			10	1.6	1		7/1/20	
Bis(2-chloroethoxy)methane 10 U 10 U <th< td=""><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td></th<>								
Bis(2-chloroethyl) Ether 10 U 10 U 1.3 1 07/06/20 09:10 7/1/20 Bis(2-ethylhexyl) Phthalate 10 U 10 I.0 1 07/06/20 09:10 7/1/20 Butyl Benzyl Phthalate 10 U 10 I.4 1 07/06/20 09:10 7/1/20								
Bis(2-ethylhexyl) Phthalate 10 U 10 1.0 1 07/06/20 09:10 7/1/20 Butyl Benzyl Phthalate 10 U 10 1.4 1 07/06/20 09:10 7/1/20								
Butyl Benzyl Phthalate 10 U 10 1.4 1 07/06/20 09:10 7/1/20								
	Chrysene	10 U	10	1.2	1	07/06/20 09:10	7/1/20	

Printed 7/15/2020 12:38:04 PM

Superset Reference:20-0000554991 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007030-01
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	10 U	10	2.0	1	07/06/20 09:10	7/1/20	
Di-n-octyl Phthalate	10 U	10	3.3	1	07/06/20 09:10	7/1/20	
Dibenz(a,h)anthracene	10 U	10	1.1	1	07/06/20 09:10	7/1/20	
Dibenzofuran	10 U	10	1.4	1	07/06/20 09:10	7/1/20	
Diethyl Phthalate	10 U	10	1.1	1	07/06/20 09:10	7/1/20	
Dimethyl Phthalate	10 U	10	1.3	1	07/06/20 09:10	7/1/20	
Fluoranthene	10 U	10	1.5	1	07/06/20 09:10	7/1/20	
Fluorene	10 U	10	1.3	1	07/06/20 09:10	7/1/20	
Hexachlorobenzene	10 U	10	1.6	1	07/06/20 09:10	7/1/20	
Hexachlorobutadiene	10 U	10	1.0	1	07/06/20 09:10	7/1/20	
Hexachlorocyclopentadiene	10 U	10	2.2	1	07/06/20 09:10	7/1/20	
Hexachloroethane	10 U	10	1.1	1	07/06/20 09:10	7/1/20	
Indeno(1,2,3-cd)pyrene	10 U	10	1.8	1	07/06/20 09:10	7/1/20	
Isophorone	10 U	10	1.4	1	07/06/20 09:10	7/1/20	
N-Nitrosodi-n-propylamine	10 U	10	1.2	1	07/06/20 09:10	7/1/20	
N-Nitrosodiphenylamine	10 U	10	2.7	1	07/06/20 09:10	7/1/20	
Naphthalene	10 U	10	1.2	1	07/06/20 09:10	7/1/20	
Nitrobenzene	10 U	10	1.5	1	07/06/20 09:10	7/1/20	
Pentachlorophenol (PCP)	50 U	50	9.8	1	07/06/20 09:10	7/1/20	
Phenanthrene	10 U	10	1.4	1	07/06/20 09:10	7/1/20	
Phenol	10 U	10	1.0	1	07/06/20 09:10	7/1/20	
Pyrene	10 U	10	1.5	1	07/06/20 09:10	7/1/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	83	35 - 141	07/06/20 09:10	
2-Fluorobiphenyl	72	31 - 118	07/06/20 09:10	
2-Fluorophenol	45	10 - 105	07/06/20 09:10	
Nitrobenzene-d5	72	31 - 110	07/06/20 09:10	
Phenol-d6	35	10 - 107	07/06/20 09:10	
p-Terphenyl-d14	70	10 - 165	07/06/20 09:10	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown	11.89	4.8	J	
	unknown hydrocarbon	12.48	6.6	J	
	unknown hydrocarbon	13.13	7.2	J	
	unknown hydrocarbon	13.83	6.6	J	
	unknown hydrocarbon	14.58	5.0	J	

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/06/20

Sample Matrix: Water

Duplicate Lab Control Sample Summary Semivolatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005539

Lab Control Sample

Duplicate Lab Control Sample

RQ2007030-02

RQ2007030-03

Analyte Name	Analytica l Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
1,2,4-Trichlorobenzene	8270D	44.6	80.0	56	46.6	80.0	58	10-127	4	30
1,2-Dichlorobenzene	8270D	43.6	80.0	54	44.1	80.0	55	23-130	2	30
1,3-Dichlorobenzene	8270D	41.9	80.0	52	42.1	80.0	53	21-90	2	30
1,4-Dichlorobenzene	8270D	42.9	80.0	54	41.2	80.0	52	10-124	4	30
2,4,5-Trichlorophenol	8270D	58.5	80.0	73	59.7	80.0	75	48-134	3	30
2,4,6-Trichlorophenol	8270D	54.1	80.0	68	54.5	80.0	68	44-135	<1	30
2,4-Dichlorophenol	8270D	50.2	80.0	63	49.4	80.0	62	48-127	2	30
2,4-Dimethylphenol	8270D	58.5	80.0	73	54.3	80.0	68	59-113	7	30
2,4-Dinitrophenol	8270D	46.5 J	80.0	58	45.7 J	80.0	57	21-154	2	30
2,4-Dinitrotoluene	8270D	62.1	80.0	78	60.5	80.0	76	54-130	3	30
2,6-Dinitrotoluene	8270D	66.6	80.0	83	69.2	80.0	86	51-127	4	30
2-Chloronaphthalene	8270D	56.4	80.0	70	55.8	80.0	70	40-108	<1	30
2-Chlorophenol	8270D	43.8	80.0	55	44.0	80.0	55	42-112	<1	30
2-Methylnaphthalene	8270D	52.7	80.0	66	49.9	80.0	62	34-102	6	30
2-Methylphenol	8270D	49.9	80.0	62	47.9	80.0	60	47-100	3	30
2-Nitroaniline	8270D	63.4	80.0	79	61.0	80.0	76	52-133	4	30
2-Nitrophenol	8270D	54.4	80.0	68	51.1	80.0	64	43-131	6	30
3,3'-Dichlorobenzidine	8270D	59.2	80.0	74	62.3	80.0	78	43-126	5	30
3- and 4-Methylphenol Coelution	8270D	44.8	80.0	56	43.0	80.0	54	40-92	4	30
3-Nitroaniline	8270D	60.6	80.0	76	57.6	80.0	72	42-111	5	30
4,6-Dinitro-2-methylphenol	8270D	53.0	80.0	66	52.9	80.0	66	36-152	<1	30
4-Bromophenyl Phenyl Ether	8270D	63.2	80.0	79	61.6	80.0	77	48-114	3	30
4-Chloro-3-methylphenol	8270D	55.8	80.0	70	51.0	80.0	64	52-113	9	30
4-Chloroaniline	8270D	55.2	80.0	69	49.1	80.0	61	44-109	12	30
4-Chlorophenyl Phenyl Ether	8270D	54.7	80.0	68	52.4	80.0	65	51-107	5	30
4-Nitroaniline	8270D	56.1	80.0	70	55.1	80.0	69	54-133	1	30
4-Nitrophenol	8270D	21.6 J	80.0	27	21.1 J	80.0	26	10-126	4	30
Acenaphthene	8270D	59.5	80.0	74	57.9	80.0	72	52-107	3	30
Acenaphthylene	8270D	62.6	80.0	78	62.2	80.0	78	55-109	<1	30
Anthracene	8270D	62.7	80.0	78	63.0	80.0	79	55-116	1	30
Benz(a)anthracene	8270D	54.4	80.0	68	60.3	80.0	75	61-121	10	30
Benzo(a)pyrene	8270D	55.5	80.0	69	62.5	80.0	78	44-114	12	30
Benzo(b)fluoranthene	8270D	52.4	80.0	65	58.6	80.0	73	62-115	12	30
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Superset Reference: 20-0000554991 rev 00

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request: R2005539 **Project:** Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/06/20

Sample Matrix: Water

Duplicate Lab Control Sample Summary Semivolatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Lab Control Sample

Duplicate Lab Control Sample

RQ2007030-02

RQ2007030-03

Analyte Name	Analytica l Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
Benzo(g,h,i)perylene	8270D	59.3	80.0	74	70.7	80.0	88	63-136	17	30
Benzo(k)fluoranthene	8270D	57.0	80.0	71	62.8	80.0	79	49-133	11	30
Benzoic Acid	8270D	68.8 J	120	57	71.9 J	120	60	10-94	5	30
Benzyl Alcohol	8270D	55.3	80.0	69	53.0	80.0	66	31-109	4	30
2,2'-Oxybis(1-chloropropane)	8270D	51.3	80.0	64	51.1	80.0	64	32-122	<1	30
Bis(2-chloroethoxy)methane	8270D	60.4	80.0	76	57.2	80.0	71	55-110	7	30
Bis(2-chloroethyl) Ether	8270D	46.4	80.0	58	47.2	80.0	59	46-102	2	30
Bis(2-ethylhexyl) Phthalate	8270D	53.6	80.0	67	64.5	80.0	81	51-132	19	30
Butyl Benzyl Phthalate	8270D	57.7	80.0	72	64.7	80.0	81	41-148	12	30
Chrysene	8270D	58.9	80.0	74	63.5	80.0	79	57-118	7	30
Di-n-butyl Phthalate	8270D	66.9	80.0	84	70.6	80.0	88	57-128	5	30
Di-n-octyl Phthalate	8270D	55.2	80.0	69	64.4	80.0	81	62-124	16	30
Dibenz(a,h)anthracene	8270D	62.4	80.0	78	72.2	80.0	90	54-135	14	30
Dibenzofuran	8270D	63.0	80.0	79	61.7	80.0	77	55-110	3	30
Diethyl Phthalate	8270D	58.7	80.0	73	57.0	80.0	71	53-113	3	30
Dimethyl Phthalate	8270D	66.0	80.0	83	65.5	80.0	82	51-112	1	30
Fluoranthene	8270D	68.4	80.0	86	69.8	80.0	87	66-127	1	30
Fluorene	8270D	63.4	80.0	79	63.6	80.0	80	54-106	1	30
Hexachlorobenzene	8270D	68.5	80.0	86	68.2	80.0	85	53-123	1	30
Hexachlorobutadiene	8270D	48.2	80.0	60	48.5	80.0	61	16-95	2	30
Hexachlorocyclopentadiene	8270D	20.5	80.0	26	21.4	80.0	27	10-99	4	30
Hexachloroethane	8270D	43.5	80.0	54	41.8	80.0	52	15-92	4	30
Indeno(1,2,3-cd)pyrene	8270D	53.8	80.0	67	65.1	80.0	81	62-137	19	30
Isophorone	8270D	51.8	80.0	65	48.6	80.0	61	50-116	6	30
N-Nitrosodi-n-propylamine	8270D	60.2	80.0	75	59.5	80.0	74	49-115	1	30
N-Nitrosodiphenylamine	8270D	79.3	80.0	99	78.7	80.0	98	45-123	1	30
Naphthalene	8270D	50.6	80.0	63	52.4	80.0	65	38-99	3	30
Nitrobenzene	8270D	54.2	80.0	68	52.7	80.0	66	46-108	3	30
Pentachlorophenol (PCP)	8270D	62.5	80.0	78	64.7	80.0	81	29-164	4	30
Phenanthrene	8270D	63.0	80.0	79	63.4	80.0	79	58-118	<1	30
Phenol	8270D	27.9	80.0	35	27.2	80.0	34	10-113	3	30
Pyrene	8270D	63.1	80.0	79	67.8	80.0	85	61-122	7	30

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Semivolatile Organic Compounds by GC

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

QA/QC Report

Service Request: R2005539

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

SURROGATE RECOVERY SUMMARY

Organochlorine Pesticides by Gas Chromatography

Analysis Method: 8081B **Extraction Method:** EPA 3510C

		Decachlorobiphenyl	Tetrachloro-m-xylene	
Sample Name	Lab Code	10-164	10-147	
WG-9954-062620-RM-014	R2005539-001	42	52	
WG-9954-062620-RM-015	R2005539-002	67	59	
WG-9954-062620-RM-016	R2005539-003	55	44	
WG-9954-062620-RM-017	R2005539-004	33	54	
WG-9954-062620-RM-018	R2005539-005	3*	34	
Method Blank	RQ2007029-01	84	75	
Lab Control Sample	RQ2007029-02	63	60	
Duplicate Lab Control Sample	RQ2007029-03	68	58	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007029-01
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	_
4,4'-DDE	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
4,4'-DDT	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
Aldrin	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
Dieldrin	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
Endosulfan I	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
Endosulfan II	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
Endosulfan Sulfate	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
Endrin	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
Endrin Ketone	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
Heptachlor	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
Heptachlor Epoxide	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
Methoxychlor	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
Toxaphene	0.50 U	0.50	0.50	1	07/06/20 10:40	7/1/20	
alpha-BHC	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
alpha-Chlordane	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
beta-BHC	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
delta-BHC	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
gamma-BHC (Lindane)	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	
gamma-Chlordane	0.050 U	0.050	0.020	1	07/06/20 10:40	7/1/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	84	10 - 164	07/06/20 10:40	
Tetrachloro-m-xylene	75	10 - 147	07/06/20 10:40	

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request: R2005539 Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/06/20

Sample Matrix: Water

Project:

Duplicate Lab Control Sample Summary Organochlorine Pesticides by Gas Chromatography

Units:ug/L Basis:NA

Lab Control Sample

Duplicate Lab Control Sample

RQ2007029-02

RQ2007029-03

	Analytical		Spike			Spike		% Rec		RPD
Analyte Name	Method	Result	Amount	% Rec	Result	Amount	% Rec	Limits	RPD	Limit
4,4'-DDD	8081B	0.260	0.400	65	0.268	0.400	67	42-159	3	30
4,4'-DDE	8081B	0.262	0.400	65	0.278	0.400	69	47-147	6	30
4,4'-DDT	8081B	0.251	0.400	63	0.275	0.400	69	41-149	9	30
Aldrin	8081B	0.216	0.400	54	0.219	0.400	55	22-137	1	30
Dieldrin	8081B	0.295	0.400	74	0.307	0.400	77	52-144	4	30
Endosulfan I	8081B	0.289	0.400	72	0.298	0.400	75	52-136	3	30
Endosulfan II	8081B	0.297	0.400	74	0.318	0.400	80	57-138	7	30
Endosulfan Sulfate	8081B	0.247	0.400	62	0.261	0.400	65	34-156	6	30
Endrin	8081B	0.282	0.400	71	0.300	0.400	75	56-143	6	30
Endrin Ketone	8081B	0.279	0.400	70	0.303	0.400	76	59-143	8	30
Heptachlor	8081B	0.190	0.400	47	0.189	0.400	47	32-141	<1	30
Heptachlor Epoxide	8081B	0.286	0.400	71	0.293	0.400	73	51-143	3	30
Methoxychlor	8081B	0.234	0.400	58	0.260	0.400	65	56-149	11	30
alpha-BHC	8081B	0.263	0.400	66	0.258	0.400	65	36-151	2	30
alpha-Chlordane	8081B	0.281	0.400	70	0.290	0.400	72	50-139	3	30
beta-BHC	8081B	0.290	0.400	72	0.293	0.400	73	55-149	1	30
delta-BHC	8081B	0.244	0.400	61	0.249	0.400	62	29-159	2	30
gamma-BHC (Lindane)	8081B	0.263	0.400	66	0.262	0.400	66	41-149	<1	30
gamma-Chlordane	8081B	0.275	0.400	69	0.283	0.400	71	50-140	3	30

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

SURROGATE RECOVERY SUMMARY Polychlorinated Biphenyls (PCBs) by GC

Analysis Method: 8082A **Extraction Method:** EPA 3510C

		Decachlorobiphenyl	Tetrachloro-m-xylene	
Sample Name	Lab Code	10-152	14-129	
WG-9954-062620-RM-014	R2005539-001	42	44	
WG-9954-062620-RM-015	R2005539-002	54	43	
WG-9954-062620-RM-016	R2005539-003	64	29	
WG-9954-062620-RM-017	R2005539-004	27	37	
WG-9954-062620-RM-018	R2005539-005	11	43	
Method Blank	RQ2007029-01	89	70	
Lab Control Sample	RQ2007029-02	71	53	
Duplicate Lab Control Sample	RQ2007029-03	68	49	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005539

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007029-01Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	1.0 U	1.0	0.50	1	07/06/20 11:40	7/1/20	
Aroclor 1221	2.0 U	2.0	1.0	1	07/06/20 11:40	7/1/20	
Aroclor 1232	1.0 U	1.0	0.50	1	07/06/20 11:40	7/1/20	
Aroclor 1242	1.0 U	1.0	0.50	1	07/06/20 11:40	7/1/20	
Aroclor 1248	1.0 U	1.0	0.50	1	07/06/20 11:40	7/1/20	
Aroclor 1254	1.0 U	1.0	0.50	1	07/06/20 11:40	7/1/20	
Aroclor 1260	1.0 U	1.0	0.50	1	07/06/20 11:40	7/1/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	89	10 - 152	07/06/20 11:40	
Tetrachloro-m-xylene	70	14 - 129	07/06/20 11:40	

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix:

Project:

Water

Duplicate Lab Control Sample Summary Polychlorinated Biphenyls (PCBs) by GC

Units:ug/L Basis:NA

Service Request: R2005539

Date Analyzed: 07/06/20

Lab Control Sample

Duplicate Lab Control Sample

RQ2007029-02

RQ2007029-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
Aroclor 1016	8082A	2.88	4.00	72	2.96	4.00	74	49-123	2	30
Aroclor 1260	8082A	3.52	4.00	88	3.35	4.00	84	30-120	5	30



Service Request No:R2005635

Ms. Kathy Willy GHD Services Inc. 2055 Niagara Falls Blvd., Niagara Falls, NY 14304

Laboratory Results for: Love Canal:292-402-D02-3100

Dear Ms. Willy,

Enclosed are the results of the sample(s) submitted to our laboratory July 01, 2020 For your reference, these analyses have been assigned our service request number **R2005635**.

All testing was performed according to our laboratory's quality assurance program and met the requirements of the TNI standards except as noted in the case narrative report. Any testing not included in the lab's accreditation is identified on a Non-Certified Analytes report. All results are intended to be considered in their entirety. ALS Environmental is not responsible for use of less than the complete report. Results apply only to the individual samples submitted to the lab for analysis, as listed in the report. The measurement uncertainty of the results included in this report is within that expected when using the prescribed method(s), and represented by Laboratory Control Sample control limits. Any events, such as QC failures or Holding Time exceedances, which may add to the uncertainty are explained in the report narrative or are flagged with qualifiers. The flags are explained in the Report Qualifiers and Definitions page of this report.

Please contact me if you have any questions. My extension is 7471. You may also contact me via email at Brady.Kalkman@alsglobal.com.

Respectfully submitted,

Goody Kullin

ALS Group USA, Corp. dba ALS Environmental

Brady Kalkman Project Manager

ADDRESS



Narrative Documents

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com



Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100

Date Received: 07/01/2020

Sample Matrix: Water

CASE NARRATIVE

All analyses were performed consistent with the quality assurance program of ALS Environmental. This report contains analytical results for samples for the Tier level IV requested by the client.

Sample Receipt:

Seven water samples were received for analysis at ALS Environmental on 07/01/2020. Any discrepancies upon initial sample inspection are annotated on the sample receipt and preservation form included within this report. The samples were stored at minimum in accordance with the analytical method requirements.

Semivolatiles by GC/MS:

Method 8270D, 07/06/2020: The upper control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). The field samples analyzed in this sequence did not contain the analyte(s) in question above the Method Reporting Limit (MRL). Since the exceedance equates to a potential high bias, the data quality was not significantly affected and no further corrective action was taken.

Method 8270D, 07/06/2020: The control limit was exceeded for one or more surrogates in the Continuing Calibration Verification (CCV). The surrogates were within acceptance limits for the associated field samples. The data quality was not significantly affected and no further corrective action was taken.

Method 8270D, 07/08/2020: The upper control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). The field samples analyzed in this sequence did not contain the analyte(s) in question above the Method Reporting Limit (MRL). Since the exceedance equates to a potential high bias, the data quality was not significantly affected and no further corrective action was taken.

Method 8270D, 07/08/2020: The control limit was exceeded for one or more surrogates in the Continuing Calibration Verification (CCV). The surrogates were within acceptance limits for the associated field samples. The data quality was not significantly affected and no further corrective action was taken.

Semivoa GC:

Method 8081B, 07/07/2020: The upper control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). The field samples analyzed in this sequence did not contain the analyte(s) in question above the Method Reporting Limit (MRL). Since the exceedance equates to a potential high bias, the data quality was not significantly affected and no further corrective action was taken.

Method 8081B, 07/07/2020: The lower control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). Since there were no detections of the analyte(s) above the MRL in the associated field samples, the quantitation is not affected. The data quality was not significantly affected and no further corrective action was taken.

Method 8081B, R2005635-005: The control limits were exceeded for one or more surrogates due to matrix interferences. A reextraction and reanalysis was performed, but produced similar results. The re-extraction was performed out of holding time. No further corrective action was required.

Method 8081B, 07/06/2020: The lower control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). Since there were no detections of the analyte(s) above the MRL in the associated field samples, the quantitation is not affected. The data quality was not significantly affected and no further corrective action was taken. Method 8082A, 07/08/2020: The upper control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). The field samples analyzed in this sequence did not contain the analyte(s) in question above the Method Reporting Limit (MRL). Since the exceedance equates to a potential high bias, the data quality was not significantly affected and



no further corrective action was taken.

Method 8082A, 07/13/2020: The upper control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). The field samples analyzed in this sequence did not contain the analyte(s) in question above the Method Reporting Limit (MRL). Since the exceedance equates to a potential high bias, the data quality was not significantly affected and no further corrective action was taken.

Method 8082A, r2005635-005: The control limits were exceeded for one or more surrogates due to matrix interferences. A reextraction and reanalysis was performed, but produced similar results. No further corrective action was required.

Method 8081B, R2005635-005: Due to matrix interference, the MRL and MDL were elevated for alpha-BHC in both the primary and secondary analysis. Analyte is flagged with an "i" for both analyses.

Volatiles by GC/MS:

Method 8260C, : The upper control criterion was exceeded for one or more analytes in the Laboratory Control Sample (LCS). There were no detections of the analyte(s) above the MRL in the associated field samples. The error associated with elevated recovery equates to a high bias. The sample data is not significantly affected. No further corrective action was appropriate.

Approved by _____

Date 07/24/2020



SAMPLE DETECTION SUMMARY

CLIENT ID: WG-9954-063020-SG-019		Lab	ID: R2005	635-001									
Analyte	Results	Flag	MDL	MRL	Units	Method							
Carbon Disulfide	4.3	J	0.42	10	ug/L	8260C							
Toluene	0.34	BJ	0.20	5.0	ug/L	8260C							
alpha-BHC	0.12		0.019	0.045	ug/L	8081B							
delta-BHC	0.076		0.019	0.045	ug/L	8081B							
gamma-BHC (Lindane)	0.10		0.019	0.045	ug/L	8081B							
CLIENT ID: WG-9954-063020-SG-020	Lab ID: R2005635-002												
Analyte	Results	Flag	MDL	MRL	Units	Method							
Carbon Disulfide	4.2	J	0.42	10	ug/L	8260C							
Toluene	0.35	BJ	0.20	5.0	ug/L	8260C							
alpha-BHC	0.081		0.019	0.045	ug/L	8081B							
delta-BHC	0.068		0.019	0.045	ug/L	8081B							
gamma-BHC (Lindane)	0.089		0.019	0.045	ug/L	8081B							
LIENT ID: WG-9954-063020-SG-022		Lab	ID: R2005	635-003									
Analyte	Results	Flag	MDL	MRL	Units	Method							
Carbon Disulfide	3.5	J	0.42	10	ug/L	8260C							
Toluene	0.38	BJ	0.20	5.0	ug/L	8260C							
alpha-BHC	0.064		0.019	0.045	ug/L	8081B							
beta-BHC	0.027	J	0.019	0.045	ug/L	8081B							
delta-BHC	0.29		0.019	0.045	ug/L	8081B							
gamma-BHC (Lindane)	0.074		0.019	0.045	ug/L	8081B							
LIENT ID: WG-9954-063020-RM-023		Lab	ID: R2005	635-004									
Analyte	Results	Flag	MDL	MRL	Units	Method							
Carbon Disulfide	2.9	J	0.42	10	ug/L	8260C							
Toluene	0.31	BJ	0.20	5.0	ug/L	8260C							
LIENT ID: RB-9954-063020-SG-002		Lab	ID: R2005	635-005									
Analyte	Results	Flag	MDL	MRL	Units	Method							
Toluene	0.25	BJ	0.20	5.0	ug/L	8260C							
Phenol	0.94	J	0.91	9.1	ug/L	8270D							
delta-BHC	0.037	J	0.019	0.045	ug/L	8081B							
gamma-BHC (Lindane)	0.058		0.019	0.045	ug/L	8081B							
delta-BHC	0.026	J	0.019	0.045	ug/L	8081B							
gamma-BHC (Lindane)	0.053		0.019	0.045	ug/L	8081B							
LIENT ID: TB-9954-063020-SG-004		Lab	ID: R2005	5635-006									
Analyte	Results	Flag	MDL	MRL	Units	Method							
Toluene	0.30	BJ	0.20	5.0	ug/L	8260C							
CLIENT ID: WG-9954-063020-SG-021		Lab	ID: R2005	5635-007									
Analyte	Results	Flag	MDL	MRL	Units	Method							
Carbon Disulfide	3.1	J	0.42	10	ug/L	8260C							



SAMPLE DETECTION SUMMARY

CLIENT ID: WG-9954-063020-SG-021	Lab ID: R2005635-007										
Analyte	Results	Flag	MDL	MRL	Units	Method					
Toluene	0.31	BJ	0.20	5.0	ug/L	8260C					
Phenol	1.1	J	0.91	9.1	ug/L	8270D					
alpha-BHC	0.10		0.019	0.045	ug/L	8081B					
delta-BHC	0.12		0.019	0.045	ug/L	8081B					
gamma-BHC (Lindane)	0.11		0.019	0.045	ug/L	8081B					



Sample Receipt Information

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request:R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

SAMPLE CROSS-REFERENCE

SAMPLE #	CLIENT SAMPLE ID	<u>DATE</u>	<u>TIME</u>
R2005635-001	WG-9954-063020-SG-019	6/30/2020	0905
R2005635-002	WG-9954-063020-SG-020	6/30/2020	1010
R2005635-003	WG-9954-063020-SG-022	6/30/2020	1100
R2005635-004	WG-9954-063020-RM-023	6/30/2020	1215
R2005635-005	RB-9954-063020-SG-002	6/30/2020	1300
R2005635-006	TB-9954-063020-SG-004	6/30/2020	0000
R2005635-007	WG-9954-063020-SG-021	6/30/2020	1010

CHAIN-OF-CUSTODY/Analytical Request Document The Chain-of-Custody is a LEGAL DOCUMENT. All relevant fields must be completed accurately.

Client Information						
GLEN SPRINGS HOLDINGS INC	Report To: Kathy Willy					
805 97TH STREET	Сору То:					
LOVE CANAL						
NIAGARA FALLS, NEW YORK 14304	Invoice To:					
Phone: 716-283-0111	PO:					
Fax: 716-283-2866	Project Name: LOVE CANAL ANNUAL GW					
Email: kathy.willy@ghd.com	Project Number: 9954					

•	Information	
Laboratory: ALS		
	-	
Laboratory Location: 15		
BUILDING 300, SUITE 36 ROCHESTER, NY 14623	3 U	
NOCHES IER, NT 14623		
ROCHESTER, NY 14623		
Laboratory Contact: BR	ADV KAI KMAN	
	ADT NACHWAR	
Requested Due Date:	TAT: 10	
QA/QC Requirements:		

Ever	nt Information
	GW SAMPLING 2020-04-1
SSOW Ref#	: 292-402-999-3100
Sampler Name: S GAR	RONER, DIVRAN

	Valid Matrix Code WG Groundwater WB Borehole Water								Sample Cond	ltion
	WS Surface Water		-	,	e e				Тетр іл С	
	SO Soil SE Sediment	90	Collected	ollected	s(None)	<u> </u>			Received on ice	Y/N
		U	3 1 1 1 1 1 1 1 1	8		Ē	₹		Sealed Cooler	Y/N
		Matrix	Date (Time (PestPCB	SVOC(none)	VOA(HCI)		Samples Intact	Y/N
Sample Identification		Ž	۵	₽	ے ا	6	۶	Remarks		
WG-9954-063020-SG-019		WG	06/30/2020	09:05	2	2	3			
WG-9954-063020-SG-020		WG	06/30/2020	10:10	2	2	3			
WG-9954-063020-SG-022		WG	06/30/2020	11:00	2	2	3			
WG-9954-063020-RM-023		WG	06/30/2020	12:15	6	6	9	MS/MSD		
RB-9954-063020-SG-002		WG	06/30/2020	13:00	2	2	3			
TB-9954-063020-SG-004		WG Q	06/30/2020	00:00	0	0	3			
WG-9954-063020-SG-021		WG	06/30/2020	10:10	2	2	3			
Total Bottles					16	16	27	Grand Tot	al:69	

SHIPMENT METHOD	NO. OF COOLERS	RELINQUISHED BY:	DATE	TIME	RECIEVED BY:	al. D.	DATE	TIME
FedEx	2	Sharen Pardner	6 3dz	1405	-11/10	MANY AUS	7-1-2020	11:05
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Project Name: Love Canal:292-402-D02-3100						<u>5</u>		14D			. <u> </u>					П	P. 6. 1. 1. 1. 1. 1. 1. 1. 1. 1. 1. 1. 1. 1.	,	
Project Number: 9954 Annual Long Term Monitoring	Report To Kathy With	· · · · · · · · · · · · · · · · · · ·	•	1	<u> </u>	_		14		-						ŀ			
Company / Address GHD Services Inc. 2055 Niagara Falls Blvd., Suite 3 Niagara Falls NY, 14304				CONTAINERS	သ			; FP											•
Phone # 716-297-2160	FAX # 716-297-2	265		l ö	est	8	200	Įξ											
Sampler Signature		rinted Name		NUMBER OF	3081B / Pest	8082A / PCB	8270D / SVO	8260C / VOC			اي	_		Rema	arks		·		•
CLIENT SAMPLE ID	LABID	SAMPLING Date Time	Matrix			8		3		I.A.	,	4					•		
1.			Liquid																
2.			Uquid																
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9.			Liquid																
10.			Liquid									L.,							
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Signature	Signature		Sign	nature			•			s	igna	ture					Signature	\neg	Signature
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Date/Time	Date/Tim	0	Dat	Date/Time						Ē	Date/Time						Date/Time		Date/Time



Cooler Receipt and Preservation Check Form



Project/Clie	ent (5 H				Fol	der Numbe	r			•		A 10 10 10 10 10 10 10 10 10 10 10 10 10	
Cooler receive	4 17	020	by:	Z	<u> </u>	COURIE	R: ALS	UPS (FEDE	VELO	OCITY CLI	ENT	
1 Were Cu	stody seals on	outside of coole	τ?		YM) 5a Pe	rchlorate s	samples	have req	uired he	adspace?	Y N	INA
2 Custody	papers proper	ly completed (in	k, sign	ed)? (Y) N	5b Di	VOA via	جا Alk,و	or Sulfide	have sig	g* bubbles?	Y	NA V
1		good condition (_		Ŷ) N		here did the				ALS/ROC	CLIE	
		Ice Gel packs	-		Y) N	7 So	il VOA rec	eived a	s: Bu	lk Er	ncore 503	Sset N	VA >
3. Temperatur		Date: 7-1-2			<u> </u>					From	Temp Blank	Sam	ple Bottle
Observed Te	emp (°C)	-124	Tá	23						ľ			
Within 0-6°	C?	(V) N		**	N	Y N	Y	N	Y	N	Y N	Y	N
If <0°C, wer	e samples froz	en? Y N		Y	N	Y N	Y	N	Y	N	Y N	Y	N
If out of T	Temperature,	note packing/ic	e cond	ition:		Ice r	nelted F	oorly P	acked (d	escribed	below)	Same I	Day Rule
&Client A	Approval to R	un Samples:		_ Stan	ding A	pproval Cl	ient aware	at drop	-off Cl	ient notii	fied by:		
All comples	held in storag	re location:	R-00) 2 t	y X	on 7	1-20 at 1	1:17					
		orage location:	17-00	<u> </u>) //c	on '	at	· · ·	within 4	R hours a	of sampling?	Y	N
											,, samping,		
10. I 11. V 12. V	Did all bottle la Were correct co Were 5035 vial	labels complete of the labels and tags ago ontainers used for sacceptable (no Cassettes / Tubes	ree with r the tes extra la	n custo sts ind abels, n	dy pape icated? not leak	ers?	ssurized	₹ Y	ES ES ES Tedlar®	NO NO NO Bags Inf	lated	₹7 <u>A</u> N/A	
pH	Lot of test	Reagent	Preser			eceived	Exp	Samp		Vol.	Lot Add	ed	Final
	paper		Yes	No				Adjus	sted	Added	•		рН
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≤2		HNO ₃			-								
≤2		H ₂ SO ₄		-	 	•		 			-		,
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(-)		608pest, 522			CN), a	scorbic (pheno	d).						
		Na ₂ S ₂ O ₃				<u> </u>							
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		HCl	**	**					vise, all bor ecked (not		samples with che entatives).	emical pro	eservatives
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Explain a	ll Discrepanci	es/ Other Comm	ents:										
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* N	id Hank?	A113 V/	a15										

	HPROD	BULK
ļ	HTR	FLDT
	SUB	HGFB
	ALS	LL3541

Labels secondary reviewed	l by: <i>@</i>	
PC Secondary Review:		

*significant air bubbles: VOA > 5-6 mm : WC > 1 in. diameter

Internal Chain of Custody Report

Client: GHD Services Inc. Service Request: R2005635

Bottle ID	Methods	Date	Time	Sample Location / User	Disposed On
R2005635-001.01					
	8081B				
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
R2005635-001.03					
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1843	R-001 / GLAFORCE	
D400#<2# 004 04		7/1/2020	1043	R-001 / GLAFORCE	
R2005635-001.04					
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1843	R-001 / GLAFORCE	
		7/13/2020	1751	In Lab / KRUEST	
		7/13/2020	1805	R-001-S12 / KRUEST	
R2005635-001.05					
	8260C				
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1843	R-001 / GLAFORCE	
		7/14/2020	1314	In Lab / KRUEST	
		7/14/2020	1451	R-001-S12 / KRUEST	
R2005635-001.06					
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
R2005635-001.08					
	8082A				
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
		7/2/2020	0816	In Lab / VSTAUFFER	
R2005635-001.09					
	8270D	7/1/2020	10.40	GMO / GLA FORGE	
		7/1/2020	1843	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
		7/2/2020	0815	In Lab / VSTAUFFER	
R2005635-002.01	00015				
	8081B	7/1/2020	1942	SMO / CLAEODCE	
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020 7/2/2020	1844 0816	R-002 / GLAFORCE In Lab / VSTAUFFER	
R2005635-002.03		1/2/2020	0010	III Lau / VSIAUFFEK	

Internal Chain of Custody Report

Client: GHD Services Inc. Service Request: R2005635

Bottle ID	Methods	Date	Time	Sample Location / User	Disposed On
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1843	R-001 / GLAFORCE	
R2005635-002.04					
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1842	R-001 / GLAFORCE	
		7/13/2020	1751	In Lab / KRUEST	
		7/13/2020	1805	R-001-S12 / KRUEST	
R2005635-002.05		7/13/2020	1003	R-001-512 / RROLS1	
	8260C				
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1843	R-001 / GLAFORCE	
		7/14/2020	1314	In Lab / KRUEST	
		7/14/2020	1451	R-001-S12 / KRUEST	
R2005635-002.06					
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
R2005635-002.08					
	8082A				
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
R2005635-002.09					
	8270D				
		7/1/2020	1843	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
		7/2/2020	0815	In Lab / VSTAUFFER	
R2005635-003.01	8081B				
	0001D	7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
R2005635-003.03					
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1843	R-001 / GLAFORCE	
R2005635-003.04		., 1, 2020		TOUT, CELH ONCE	_
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1843	R-001 / GLAFORCE	
		7/13/2020	1751	In Lab / KRUEST	
Printed 7/24/2020 2:55	:37 PM		Page 13 of 12	1	

Internal Chain of Custody Report

Client: GHD Services Inc. Service Request: R2005635

Bottle ID	Methods	Date	Time	Sample Location / User	Disposed On
		7/13/2020	1805	R-001-S12 / KRUEST	
R2005635-003.05					
	8260C				
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1843	R-001 / GLAFORCE	
		7/14/2020	1314	In Lab / KRUEST	
		7/14/2020	1451	R-001-S12 / KRUEST	
R2005635-003.06					
	8270D				
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
		7/2/2020	0815	In Lab / VSTAUFFER	
R2005635-003.08					
	8082A	7/1/2020	1042	CMO / CLAEODOE	
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
		7/2/2020	0816	In Lab / VSTAUFFER	
R2005635-003.09					
		7/1/2020	1843	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
R2005635-004.01					
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
R2005635-004.07		7,1,2020	1011	R 0027 GEIN GROE	
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1843	R-001 / GLAFORCE	
R2005635-004.08	92606				
	8260C	7/1/2020	1042	SMO / CLAEODCE	
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1843	R-001 / GLAFORCE	
		7/14/2020	1315	In Lab / KRUEST	
D2005/25 004 00		7/14/2020	1451	R-001-S12 / KRUEST	_
R2005635-004.09					
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1843	R-001 / GLAFORCE	

Internal Chain of Custody Report

Client: GHD Services Inc. Service Request: R2005635

Bottle ID	Methods	Date	Time	Sample Location / User	Disposed On
R2005635-004.10					
		7/1/2020	1843	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
R2005635-004.11					
		7/1/2020	1843	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
R2005635-004.12	8270D				
	8270D	7/1/2020	1843	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
		7/2/2020	0815	In Lab / VSTAUFFER	
R2005635-004.13					
	8081B,8082A				
		7/1/2020	1843	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
		7/2/2020	0815	In Lab / VSTAUFFER	
R2005635-004.14					
		7/1/2020	1843	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
R2005635-004.15		., ., _ , _ ,			
		7/1/2020	1843	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
		7/2/2020	0815	In Lab / VSTAUFFER	
R2005635-004.16					
		7/1/2020	1843	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
		7/2/2020	0815	In Lab / VSTAUFFER	
R2005635-004.17					
		7/1/2020	1843	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
D2005/25 004 40		7/2/2020	0815	In Lab / VSTAUFFER	
R2005635-004.18					
		7/1/2020	1843	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
D 1 5/21/2022 5 5 5 5 5	DI 4		Page 15 of 121		
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Internal Chain of Custody Report

Client: GHD Services Inc. Service Request: R2005635

Bottle ID	Methods	Date	Time	Sample Location / User	Disposed On
		7/2/2020	0815	In Lab / VSTAUFFER	
R2005635-004.19		,, _, _,			
		7/1/2020	1843	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
		7/2/2020	0815	In Lab / VSTAUFFER	
R2005635-004.20					
		7/1/2020	1843	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
		7/2/2020	0815	In Lab / VSTAUFFER	
R2005635-004.21		,			
		7/1/2020	1843	R-001 / GLAFORCE	
		7/1/2020	1843	SMO / GLAFORCE	
		7/1/2020	1751	In Lab / KRUEST	
		7/13/2020	1805	R-001-S12 / KRUEST	
R2005635-004.22		.,,,			
		7/1/2020	1843	R-001 / GLAFORCE	
		7/1/2020	1843	SMO / GLAFORCE	
R2005635-004.23					
		7/1/2020	1843	R-001 / GLAFORCE	
		7/1/2020	1843	SMO / GLAFORCE	
R2005635-004.24					
		7/1/2020	1843	R-001 / GLAFORCE	
		7/1/2020	1843	SMO / GLAFORCE	
R2005635-004.25					
		7/1/2020	1843	R-001 / GLAFORCE	
		7/1/2020	1843	SMO / GLAFORCE	
R2005635-004.26					
		7/1/2020	1843	R-001 / GLAFORCE	
		7/1/2020	1843	SMO / GLAFORCE	

Internal Chain of Custody Report

Client: GHD Services Inc. Service Request: R2005635

Bottle ID	Methods	Date	Time	Sample Location / User	Disposed On
	8081B,8081B				
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
		7/2/2020	0815	In Lab / VSTAUFFER	
R2005635-005.03					
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1843	R-001 / GLAFORCE	
R2005635-005.04					
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1843	R-001 / GLAFORCE	
		7/13/2020	1751	In Lab / KRUEST	
		7/13/2020	1805	R-001-S12 / KRUEST	
R2005635-005.05	0.5.10.5				
	8260C	7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1842 1843	R-001 / GLAFORCE	
		7/1/2020	1314	In Lab / KRUEST	
		7/14/2020	1452	R-001-S12 / KRUEST	
R2005635-005.06		7/14/2020	1432	K-001-512 / KKOL51	
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
D2005/25 005 00		7/1/2020	1044	K-0027 GLAFORCE	
R2005635-005.08	8082A,8082A				
	0002/1,0002/1	7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
		7/10/2020	0804	In Lab / VSTAUFFER	
R2005635-005.09					
	8270D				
		7/1/2020	1843	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
		7/2/2020	0815	In Lab / VSTAUFFER	
R2005635-006.01					
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1843	R-001 / GLAFORCE	
		7/13/2020	1751	In Lab / KRUEST	
		7/13/2020	1805	R-001-S12 / KRUEST	

Internal Chain of Custody Report

Client: GHD Services Inc. Service Request: R2005635

Bottle ID	Methods	Date	Time	Sample Location / User	Disposed On
	8260C				
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1843	R-001 / GLAFORCE	
		7/14/2020	1314	In Lab / KRUEST	
		7/14/2020	1451	R-001-S12 / KRUEST	
R2005635-006.03					
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1843	R-001 / GLAFORCE	
R2005635-007.01					
	8081B				
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
R2005635-007.03					
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1843	R-001 / GLAFORCE	
R2005635-007.04					
	8260C				
	02000	7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1843	R-001 / GLAFORCE	
		7/13/2020	1751	In Lab / KRUEST	
		7/14/2020	1451	R-001-S12 / KRUEST	
R2005635-007.05					
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1843	R-001 / GLAFORCE	
		7/13/2020	1805	R-001-S12 / KRUEST	
		7/14/2020	1314	In Lab / KRUEST	
R2005635-007.06					
		7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
R2005635-007.08					
K4003033=007.00	8082A				
	000211	7/1/2020	1842	SMO / GLAFORCE	
		7/1/2020	1844	R-002 / GLAFORCE	
		7/2/2020	0815	In Lab / VSTAUFFER	
D2005625 007 00		1,2,2020	0013	III Duo / TOTTIOTI DIX	
R2005635-007.09	8270D				
	02701	7/1/2020	1843	SMO / GLAFORCE	
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Internal Chain of Custody Report

Client: GHD Services Inc. Service Request: R2005635

Bottle ID	Methods	Date	Time	Sample Location / User	Disposed On
	8270D				_
		7/1/2020	1844	R-002 / GLAFORCE	
		7/2/2020	0815	In Lab / VSTAUFFER	



Miscellaneous Forms

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com



REPORT QUALIFIERS AND DEFINITIONS

- U Analyte was analyzed for but not detected. The sample quantitation limit has been corrected for dilution and for percent moisture, unless otherwise noted in the case narrative.
- J Estimated value due to either being a Tentatively Identified Compound (TIC) or that the concentration is between the MRL and the MDL. Concentrations are not verified within the linear range of the calibration. For DoD: concentration >40% difference between two GC columns (pesticides/Arclors).
- B Analyte was also detected in the associated method blank at a concentration that may have contributed to the sample result.
- E Inorganics- Concentration is estimated due to the serial dilution was outside control limits.
- E Organics- Concentration has exceeded the calibration range for that specific analysis.
- D Concentration is a result of a dilution, typically a secondary analysis of the sample due to exceeding the calibration range or that a surrogate has been diluted out of the sample and cannot be assessed.
- * Indicates that a quality control parameter has exceeded laboratory limits. Under the õNotesö column of the Form I, this qualifier denotes analysis was performed out of Holding Time.
- H Analysis was performed out of hold time for tests that have an õimmediateö hold time criteria.
- # Spike was diluted out.

- + Correlation coefficient for MSA is <0.995.
- N Inorganics- Matrix spike recovery was outside laboratory limits.
- N Organics- Presumptive evidence of a compound (reported as a TIC) based on the MS library search.
- S Concentration has been determined using Method of Standard Additions (MSA).
- W Post-Digestion Spike recovery is outside control limits and the sample absorbance is <50% of the spike absorbance.
- P Concentration >40% difference between the two GC columns.
- C Confirmed by GC/MS
- Q DoD reports: indicates a pesticide/Aroclor is not confirmed (×100% Difference between two GC columns).
- X See Case Narrative for discussion.
- MRL Method Reporting Limit. Also known as:
- LOQ Limit of Quantitation (LOQ)

 The lowest concentration at which the method analyte may be reliably quantified under the method conditions.
- MDL Method Detection Limit. A statistical value derived from a study designed to provide the lowest concentration that will be detected 99% of the time. Values between the MDL and MRL are estimated (see J qualifier).
- LOD Limit of Detection. A value at or above the MDL which has been verified to be detectable.
- ND Non-Detect. Analyte was not detected at the concentration listed. Same as U qualifier.



Rochester Lab ID # for State Certifications¹

Connecticut ID # PH0556	Maine ID #NY0032	Pennsylvania ID# 68-786
Delaware Approved	New Hampshire ID # 2941	Rhode Island ID # 158
DoD ELAP #65817	New York ID # 10145	Virginia #460167
Florida ID # E87674	North Carolina #676	

¹ Analyses were performed according to our laboratory

NELAP-approved quality assurance program and any applicable state or agency requirements. The test results meet requirements of the current NELAP/TNI standards or state or agency requirements, where applicable, except as noted in the case narrative. Since not all analyte/method/matrix combinations are offered for state/NELAC accreditation, this report may contain results which are not accredited. For a specific list of accredited analytes, contact the laboratory or go to https://www.alsglobal.com/locations/americas/north-america/usa/new-york/rochester-environmental

9/28/18

ALS Laboratory Group

Acronyms

ASTM American Society for Testing and Materials

A2LA American Association for Laboratory Accreditation

CARB California Air Resources Board

CAS Number Chemical Abstract Service registry Number

CFC Chlorofluorocarbon CFU Colony-Forming Unit

DEC Department of Environmental Conservation

DEQ Department of Environmental Quality

DHS Department of Health Services

DOE Department of Ecology DOH Department of Health

EPA U. S. Environmental Protection Agency

ELAP Environmental Laboratory Accreditation Program

GC Gas Chromatography

GC/MS Gas Chromatography/Mass Spectrometry

LUFT Leaking Underground Fuel Tank

M Modified

MCL Maximum Contaminant Level is the highest permissible concentration of a

substance allowed in drinking water as established by the USEPA.

MDL Method Detection Limit
MPN Most Probable Number
MRL Method Reporting Limit

NA Not Applicable NC Not Calculated

NCASI National Council of the Paper Industry for Air and Stream Improvement

ND Not Detected

NIOSH National Institute for Occupational Safety and Health

PQL Practical Quantitation Limit

RCRA Resource Conservation and Recovery Act

SIM Selected Ion Monitoring

TPH Total Petroleum Hydrocarbons

tr Trace level is the concentration of an analyte that is less than the PQL but

greater than or equal to the MDL.

Analyst Summary report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term **Project:**

Monitoring

Sample Name: WG-9954-063020-SG-019

Lab Code: R2005635-001

Sample Matrix: Water

Analyzed By

Service Request: R2005635

Date Collected: 06/30/20

Date Received: 07/1/20

Analysis Method Extracted/Digested By 8081B **KSERCU JMISIUREWICZ** 8082A **KSERCU BALLGEIER**

8260C KRUEST

8270D **KSERCU JMISIUREWICZ**

WG-9954-063020-SG-020 **Sample Name: Date Collected:** 06/30/20

Lab Code: R2005635-002 **Date Received:** 07/1/20

Sample Matrix: Water

Extracted/Digested By Analyzed By Analysis Method

8081B **KSERCU JMISIUREWICZ** 8082A **KSERCU BALLGEIER** 8260C **KRUEST**

8270D **KSERCU JMISIUREWICZ**

Sample Name: WG-9954-063020-SG-022 **Date Collected:** 06/30/20

Lab Code: R2005635-003 Date Received: 07/1/20

Sample Matrix: Water

Analyzed By Extracted/Digested By Analysis Method

8081B **KSERCU JMISIUREWICZ** 8082A **KSERCU BALLGEIER** 8260C **KRUEST**

KSERCU 8270D **JMISIUREWICZ**

Sample Name: WG-9954-063020-RM-023 Date Collected: 06/30/20

Lab Code: R2005635-004 **Date Received:** 07/1/20 Sample Matrix: Water

Analyzed By Extracted/Digested By Analysis Method

8081B **KSERCU JMISIUREWICZ**

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Analyst Summary report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Monitoring

Sample Name: WG-9954-063020-RM-023

Lab Code: R2005635-004

Sample Matrix: Water

Analysis Method

Extracted/Digested By Analyzed By

Service Request: R2005635

Date Collected: 06/30/20

Date Received: 07/1/20

8082A KSERCU BALLGEIER

8260C KRUEST

8270D KSERCU JMISIUREWICZ

ozrob wildere wie

Sample Name: RB-9954-063020-SG-002 **Date Collected:** 06/30/20

Lab Code: R2005635-005 **Date Received:** 07/1/20

Sample Matrix: Water

Analysis Method Extracted/Digested By Analyzed By

8081BKSERCUJMISIUREWICZ8082AKSERCUBALLGEIER8260CKRUEST

8270D KSERCU JMISIUREWICZ

Sample Name: RB-9954-063020-SG-002 **Date Collected:** 06/30/20

Lab Code: R2005635-005.R01 Date Received: 07/1/20 Sample Matrix: Water

Analysis Method Extracted/Digested By Analyzed By

8081B KSERCU BALLGEIER 8082A KSERCU BALLGEIER

Sample Name: TB-9954-063020-SG-004 **Date Collected:** 06/30/20

Lab Code:R2005635-006Date Received: 07/1/20Sample Matrix:Water

Analysis Method Extracted/Digested By Analyzed By 8260C KRUEST

Analyst Summary report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Monitoring

Sample Name: WG-9954-063020-SG-021

Lab Code: R2005635-007

Sample Matrix: Water

Service Request: R2005635

Date Collected: 06/30/20

Date Received: 07/1/20

Analysis MethodExtracted/Digested ByAnalyzed By8081BKSERCUJMISIUREWICZ8082AKSERCUBALLGEIER8260CKRUEST8270DKSERCUJMISIUREWICZ

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INORGANIC PREPARATION METHODS

The preparation methods associated with this report are found in these tables unless discussed in the case narrative.

Water/Liquid Matrix

Analytical Method	Preparation Method
200.7	200.2
200.8	200.2
6010C	3005A/3010A
6020A	ILM05.3
9034 Sulfide Acid Soluble	9030B
SM 4500-CN-E Residual Cyanide	SM 4500-CN-G
SM 4500-CN-E WAD Cyanide	SM 4500-CN-I

Solid/Soil/Non-Aqueous Matrix

Analytical Method	Preparation Method			
6010C	3050B			
6020A	3050B			
6010C TCLP (1311)	3005A/3010A			
extract				
6010 SPLP (1312) extract	3005A/3010A			
7199	3060A			
300.0 Anions/ 350.1/	DI extraction			
353.2/ SM 2320B/ SM				
5210B/ 9056A Anions				
For analytical methods not listed, the preparation method is the same as the analytical method reference.				



Sample Results

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com



Volatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 09:05

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-019
 Units: ug/L

 Lab Code:
 R2005635-001
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/14/20 15:20	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/14/20 15:20	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/14/20 15:20	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/14/20 15:20	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/14/20 15:20	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/14/20 15:20	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/14/20 15:20	
2-Butanone (MEK)	10 U	10	0.78	1	07/14/20 15:20	
2-Hexanone	10 U	10	0.20	1	07/14/20 15:20	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/14/20 15:20	
Acetone	10 U	10	5.0	1	07/14/20 15:20	
Benzene	5.0 U	5.0	0.20	1	07/14/20 15:20	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/14/20 15:20	
Bromoform	5.0 U	5.0	0.25	1	07/14/20 15:20	
Bromomethane	5.0 U	5.0	0.70	1	07/14/20 15:20	
Carbon Disulfide	4.3 J	10	0.42	1	07/14/20 15:20	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/14/20 15:20	
Chlorobenzene	5.0 U	5.0	0.20	1	07/14/20 15:20	
Chloroethane	5.0 U	5.0	0.23	1	07/14/20 15:20	
Chloroform	5.0 U	5.0	0.24	1	07/14/20 15:20	
Chloromethane	5.0 U	5.0	0.28	1	07/14/20 15:20	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/14/20 15:20	
Dichloromethane	5.0 U	5.0	0.65	1	07/14/20 15:20	
Ethylbenzene	5.0 U	5.0	0.20	1	07/14/20 15:20	
Styrene	5.0 U	5.0	0.20	1	07/14/20 15:20	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/14/20 15:20	
Toluene	0.34 BJ	5.0	0.20	1	07/14/20 15:20	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/14/20 15:20	
Vinyl Acetate	10 U	10	1.1	1	07/14/20 15:20	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/14/20 15:20	
Xylenes, Total	5.0 U	5.0	0.23	1	07/14/20 15:20	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/14/20 15:20	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/14/20 15:20	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/14/20 15:20	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/14/20 15:20	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 09:05

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-019
 Units: ug/L

 Lab Code:
 R2005635-001
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	94	85 - 122	07/14/20 15:20	
Dibromofluoromethane	100	89 - 119	07/14/20 15:20	
Toluene-d8	101	87 - 121	07/14/20 15:20	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 09:05

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-019
 Units: ug/L

 Lab Code:
 R2005635-001
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

CAS#	Compound Identification	RT	ug/L	Q	
	unknown	1.21	9.0	J	
	unknown	1.32	190.6	J	
	unknown	1.54	6.4	J	
007446-09-5	Sulfur dioxide	1.62	78.4	JN	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 10:10

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-020
 Units: ug/L

 Lab Code:
 R2005635-002
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/14/20 15:42	_
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/14/20 15:42	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/14/20 15:42	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/14/20 15:42	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/14/20 15:42	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/14/20 15:42	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/14/20 15:42	
2-Butanone (MEK)	10 U	10	0.78	1	07/14/20 15:42	
2-Hexanone	10 U	10	0.20	1	07/14/20 15:42	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/14/20 15:42	
Acetone	10 U	10	5.0	1	07/14/20 15:42	
Benzene	5.0 U	5.0	0.20	1	07/14/20 15:42	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/14/20 15:42	
Bromoform	5.0 U	5.0	0.25	1	07/14/20 15:42	
Bromomethane	5.0 U	5.0	0.70	1	07/14/20 15:42	
Carbon Disulfide	4.2 J	10	0.42	1	07/14/20 15:42	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/14/20 15:42	
Chlorobenzene	5.0 U	5.0	0.20	1	07/14/20 15:42	
Chloroethane	5.0 U	5.0	0.23	1	07/14/20 15:42	
Chloroform	5.0 U	5.0	0.24	1	07/14/20 15:42	
Chloromethane	5.0 U	5.0	0.28	1	07/14/20 15:42	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/14/20 15:42	
Dichloromethane	5.0 U	5.0	0.65	1	07/14/20 15:42	
Ethylbenzene	5.0 U	5.0	0.20	1	07/14/20 15:42	
Styrene	5.0 U	5.0	0.20	1	07/14/20 15:42	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/14/20 15:42	
Toluene	0.35 BJ	5.0	0.20	1	07/14/20 15:42	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/14/20 15:42	
Vinyl Acetate	10 U	10	1.1	1	07/14/20 15:42	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/14/20 15:42	
Xylenes, Total	5.0 U	5.0	0.23	1	07/14/20 15:42	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/14/20 15:42	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/14/20 15:42	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/14/20 15:42	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/14/20 15:42	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 10:10

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-020
 Units: ug/L

 Lab Code:
 R2005635-002
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	97	85 - 122	07/14/20 15:42	
Dibromofluoromethane	105	89 - 119	07/14/20 15:42	
Toluene-d8	103	87 - 121	07/14/20 15:42	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 10:10

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-020
 Units: ug/L

 Lab Code:
 R2005635-002
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Tentatively Identified Compounds

 CAS#
 Compound Identification
 RT
 ug/L
 Q

 007446-09-5
 Sulfur dioxide unknown
 1.29
 41.3
 JN

 47.4
 J

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 11:00

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-022
 Units: ug/L

 Lab Code:
 R2005635-003
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/14/20 16:04	_
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/14/20 16:04	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/14/20 16:04	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/14/20 16:04	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/14/20 16:04	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/14/20 16:04	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/14/20 16:04	
2-Butanone (MEK)	10 U	10	0.78	1	07/14/20 16:04	
2-Hexanone	10 U	10	0.20	1	07/14/20 16:04	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/14/20 16:04	
Acetone	10 U	10	5.0	1	07/14/20 16:04	
Benzene	5.0 U	5.0	0.20	1	07/14/20 16:04	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/14/20 16:04	
Bromoform	5.0 U	5.0	0.25	1	07/14/20 16:04	
Bromomethane	5.0 U	5.0	0.70	1	07/14/20 16:04	
Carbon Disulfide	3.5 J	10	0.42	1	07/14/20 16:04	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/14/20 16:04	
Chlorobenzene	5.0 U	5.0	0.20	1	07/14/20 16:04	
Chloroethane	5.0 U	5.0	0.23	1	07/14/20 16:04	
Chloroform	5.0 U	5.0	0.24	1	07/14/20 16:04	
Chloromethane	5.0 U	5.0	0.28	1	07/14/20 16:04	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/14/20 16:04	
Dichloromethane	5.0 U	5.0	0.65	1	07/14/20 16:04	
Ethylbenzene	5.0 U	5.0	0.20	1	07/14/20 16:04	
Styrene	5.0 U	5.0	0.20	1	07/14/20 16:04	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/14/20 16:04	
Toluene	0.38 BJ	5.0	0.20	1	07/14/20 16:04	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/14/20 16:04	
Vinyl Acetate	10 U	10	1.1	1	07/14/20 16:04	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/14/20 16:04	
Xylenes, Total	5.0 U	5.0	0.23	1	07/14/20 16:04	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/14/20 16:04	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/14/20 16:04	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/14/20 16:04	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/14/20 16:04	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 11:00

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-022
 Units: ug/L

 Lab Code:
 R2005635-003
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	95	85 - 122	07/14/20 16:04	
Dibromofluoromethane	102	89 - 119	07/14/20 16:04	
Toluene-d8	102	87 - 121	07/14/20 16:04	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 11:00

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-022
 Units: ug/L

 Lab Code:
 R2005635-003
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

CAS#			Result		
	Compound Identification	RT	ug/L	Q	
007446-09-5	Sulfur dioxide	1.29	51.5	JN	
	unknown	1.62	19.6	J	
	unknown	9.24	5.1	J	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 12:15

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-RM-023
 Units: ug/L

 Lab Code:
 R2005635-004
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/14/20 16:25	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/14/20 16:25	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/14/20 16:25	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/14/20 16:25	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/14/20 16:25	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/14/20 16:25	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/14/20 16:25	
2-Butanone (MEK)	10 U	10	0.78	1	07/14/20 16:25	
2-Hexanone	10 U	10	0.20	1	07/14/20 16:25	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/14/20 16:25	
Acetone	10 U	10	5.0	1	07/14/20 16:25	
Benzene	5.0 U	5.0	0.20	1	07/14/20 16:25	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/14/20 16:25	
Bromoform	5.0 U	5.0	0.25	1	07/14/20 16:25	
Bromomethane	5.0 U	5.0	0.70	1	07/14/20 16:25	
Carbon Disulfide	2.9 J	10	0.42	1	07/14/20 16:25	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/14/20 16:25	
Chlorobenzene	5.0 U	5.0	0.20	1	07/14/20 16:25	
Chloroethane	5.0 U	5.0	0.23	1	07/14/20 16:25	
Chloroform	5.0 U	5.0	0.24	1	07/14/20 16:25	
Chloromethane	5.0 U	5.0	0.28	1	07/14/20 16:25	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/14/20 16:25	
Dichloromethane	5.0 U	5.0	0.65	1	07/14/20 16:25	
Ethylbenzene	5.0 U	5.0	0.20	1	07/14/20 16:25	
Styrene	5.0 U	5.0	0.20	1	07/14/20 16:25	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/14/20 16:25	
Toluene	0.31 BJ	5.0	0.20	1	07/14/20 16:25	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/14/20 16:25	
Vinyl Acetate	10 U	10	1.1	1	07/14/20 16:25	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/14/20 16:25	
Xylenes, Total	5.0 U	5.0	0.23	1	07/14/20 16:25	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/14/20 16:25	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/14/20 16:25	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/14/20 16:25	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/14/20 16:25	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 12:15

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-RM-023
 Units: ug/L

 Lab Code:
 R2005635-004
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	90	85 - 122	07/14/20 16:25	
Dibromofluoromethane	99	89 - 119	07/14/20 16:25	
Toluene-d8	96	87 - 121	07/14/20 16:25	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 12:15

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-RM-023
 Units: ug/L

 Lab Code:
 R2005635-004
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
007446-09-5	Sulfur dioxide	1.29	47.6	JN	
	unknown	1.62	22.2	J	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 13:00

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 RB-9954-063020-SG-002
 Units: ug/L

 Lab Code:
 R2005635-005
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/14/20 16:47	_
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/14/20 16:47	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/14/20 16:47	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/14/20 16:47	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/14/20 16:47	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/14/20 16:47	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/14/20 16:47	
2-Butanone (MEK)	10 U	10	0.78	1	07/14/20 16:47	
2-Hexanone	10 U	10	0.20	1	07/14/20 16:47	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/14/20 16:47	
Acetone	10 U	10	5.0	1	07/14/20 16:47	
Benzene	5.0 U	5.0	0.20	1	07/14/20 16:47	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/14/20 16:47	
Bromoform	5.0 U	5.0	0.25	1	07/14/20 16:47	
Bromomethane	5.0 U	5.0	0.70	1	07/14/20 16:47	
Carbon Disulfide	10 U	10	0.42	1	07/14/20 16:47	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/14/20 16:47	
Chlorobenzene	5.0 U	5.0	0.20	1	07/14/20 16:47	
Chloroethane	5.0 U	5.0	0.23	1	07/14/20 16:47	
Chloroform	5.0 U	5.0	0.24	1	07/14/20 16:47	
Chloromethane	5.0 U	5.0	0.28	1	07/14/20 16:47	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/14/20 16:47	
Dichloromethane	5.0 U	5.0	0.65	1	07/14/20 16:47	
Ethylbenzene	5.0 U	5.0	0.20	1	07/14/20 16:47	
Styrene	5.0 U	5.0	0.20	1	07/14/20 16:47	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/14/20 16:47	
Toluene	0.25 BJ	5.0	0.20	1	07/14/20 16:47	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/14/20 16:47	
Vinyl Acetate	10 U	10	1.1	1	07/14/20 16:47	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/14/20 16:47	
Xylenes, Total	5.0 U	5.0	0.23	1	07/14/20 16:47	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/14/20 16:47	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/14/20 16:47	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/14/20 16:47	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/14/20 16:47	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 13:00

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 RB-9954-063020-SG-002
 Units: ug/L

 Lab Code:
 R2005635-005
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	94	85 - 122	07/14/20 16:47	
Dibromofluoromethane	99	89 - 119	07/14/20 16:47	
Toluene-d8	101	87 - 121	07/14/20 16:47	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 13:00

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

Sample Name: RB-9954-063020-SG-002 **Units:** ug/L

Lab Code: R2005635-005 **Basis:** NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Tentatively Identified Compounds

 CAS#
 Compound Identification
 RT
 ug/L
 Q

 007446-09-5
 Sulfur dioxide
 1.28
 84.1
 JN

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 00:00

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 TB-9954-063020-SG-004
 Units: ug/L

 Lab Code:
 R2005635-006
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/14/20 14:58	_
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/14/20 14:58	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/14/20 14:58	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/14/20 14:58	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/14/20 14:58	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/14/20 14:58	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/14/20 14:58	
2-Butanone (MEK)	10 U	10	0.78	1	07/14/20 14:58	
2-Hexanone	10 U	10	0.20	1	07/14/20 14:58	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/14/20 14:58	
Acetone	10 U	10	5.0	1	07/14/20 14:58	
Benzene	5.0 U	5.0	0.20	1	07/14/20 14:58	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/14/20 14:58	
Bromoform	5.0 U	5.0	0.25	1	07/14/20 14:58	
Bromomethane	5.0 U	5.0	0.70	1	07/14/20 14:58	
Carbon Disulfide	10 U	10	0.42	1	07/14/20 14:58	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/14/20 14:58	
Chlorobenzene	5.0 U	5.0	0.20	1	07/14/20 14:58	
Chloroethane	5.0 U	5.0	0.23	1	07/14/20 14:58	
Chloroform	5.0 U	5.0	0.24	1	07/14/20 14:58	
Chloromethane	5.0 U	5.0	0.28	1	07/14/20 14:58	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/14/20 14:58	
Dichloromethane	5.0 U	5.0	0.65	1	07/14/20 14:58	
Ethylbenzene	5.0 U	5.0	0.20	1	07/14/20 14:58	
Styrene	5.0 U	5.0	0.20	1	07/14/20 14:58	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/14/20 14:58	
Toluene	0.30 BJ	5.0	0.20	1	07/14/20 14:58	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/14/20 14:58	
Vinyl Acetate	10 U	10	1.1	1	07/14/20 14:58	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/14/20 14:58	
Xylenes, Total	5.0 U	5.0	0.23	1	07/14/20 14:58	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/14/20 14:58	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/14/20 14:58	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/14/20 14:58	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/14/20 14:58	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 00:00

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 TB-9954-063020-SG-004
 Units: ug/L

 Lab Code:
 R2005635-006
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	94	85 - 122	07/14/20 14:58	
Dibromofluoromethane	98	89 - 119	07/14/20 14:58	
Toluene-d8	102	87 - 121	07/14/20 14:58	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 00:00

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 TB-9954-063020-SG-004
 Units: ug/L

 Lab Code:
 R2005635-006
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown	1.60	6.5	J	
001825-61-2	Silane, methoxytrimethyl-	2.92	11.9	JN	
	unknown	4.49	10.8	J	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 10:10

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-021
 Units: ug/L

 Lab Code:
 R2005635-007
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/14/20 17:09	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/14/20 17:09	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/14/20 17:09	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/14/20 17:09	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/14/20 17:09	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/14/20 17:09	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/14/20 17:09	
2-Butanone (MEK)	10 U	10	0.78	1	07/14/20 17:09	
2-Hexanone	10 U	10	0.20	1	07/14/20 17:09	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/14/20 17:09	
Acetone	10 U	10	5.0	1	07/14/20 17:09	
Benzene	5.0 U	5.0	0.20	1	07/14/20 17:09	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/14/20 17:09	
Bromoform	5.0 U	5.0	0.25	1	07/14/20 17:09	
Bromomethane	5.0 U	5.0	0.70	1	07/14/20 17:09	
Carbon Disulfide	3.1 J	10	0.42	1	07/14/20 17:09	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/14/20 17:09	
Chlorobenzene	5.0 U	5.0	0.20	1	07/14/20 17:09	
Chloroethane	5.0 U	5.0	0.23	1	07/14/20 17:09	
Chloroform	5.0 U	5.0	0.24	1	07/14/20 17:09	
Chloromethane	5.0 U	5.0	0.28	1	07/14/20 17:09	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/14/20 17:09	
Dichloromethane	5.0 U	5.0	0.65	1	07/14/20 17:09	
Ethylbenzene	5.0 U	5.0	0.20	1	07/14/20 17:09	
Styrene	5.0 U	5.0	0.20	1	07/14/20 17:09	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/14/20 17:09	
Toluene	0.31 BJ	5.0	0.20	1	07/14/20 17:09	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/14/20 17:09	
Vinyl Acetate	10 U	10	1.1	1	07/14/20 17:09	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/14/20 17:09	
Xylenes, Total	5.0 U	5.0	0.23	1	07/14/20 17:09	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/14/20 17:09	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/14/20 17:09	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/14/20 17:09	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/14/20 17:09	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 10:10

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-021
 Units: ug/L

 Lab Code:
 R2005635-007
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	95	85 - 122	07/14/20 17:09	
Dibromofluoromethane	100	89 - 119	07/14/20 17:09	
Toluene-d8	101	87 - 121	07/14/20 17:09	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 10:10

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-021
 Units: ug/L

 Lab Code:
 R2005635-007
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
007446-09-5	Sulfur dioxide	1.29	46.7	JN	
	unknown	1.62	24.0	J	



Semivolatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 09:05

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-019
 Units: ug/L

 Lab Code:
 R2005635-001
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed D	ate Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 16:52	7/2/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 16:52	7/2/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/06/20 16:52	7/2/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 16:52	7/2/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/06/20 16:52	7/2/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/06/20 16:52	7/2/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/06/20 16:52	7/2/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/06/20 16:52	7/2/20	
2,4-Dinitrophenol	45 U	45	19	1	07/06/20 16:52	7/2/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/06/20 16:52	7/2/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/06/20 16:52	7/2/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/06/20 16:52	7/2/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/06/20 16:52	7/2/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/06/20 16:52	7/2/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/06/20 16:52	7/2/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/06/20 16:52	7/2/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/06/20 16:52	7/2/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/06/20 16:52	7/2/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/06/20 16:52	7/2/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/06/20 16:52	7/2/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/06/20 16:52	7/2/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/06/20 16:52	7/2/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/06/20 16:52	7/2/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/06/20 16:52	7/2/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/06/20 16:52	7/2/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/06/20 16:52	7/2/20	
4-Nitrophenol	45 U	45	5.8	1	07/06/20 16:52	7/2/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/06/20 16:52	7/2/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/06/20 16:52	7/2/20	
Anthracene	9.1 U	9.1	1.2	1	07/06/20 16:52	7/2/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/06/20 16:52	7/2/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/06/20 16:52	7/2/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/06/20 16:52	7/2/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/06/20 16:52	7/2/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/06/20 16:52	7/2/20	
Benzoic Acid	91 U	91	33	1	07/06/20 16:52	7/2/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/06/20 16:52	7/2/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/06/20 16:52	7/2/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/06/20 16:52	7/2/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.3	1	07/06/20 16:52	7/2/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	07/06/20 16:52	7/2/20	
Butyl Benzyl Phthalate	9.1 U 9.1 U	9.1 9.1	1.3	1	07/06/20 16:52	7/2/20	
	9.1 U 9.1 U	9.1 9.1	1.3	1	07/06/20 16:52	7/2/20	
Chrysene	9.1 U	9.1	1.1	1	07/00/20 10:52	1/2/2U	

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Superset Reference: 20-0000555197 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 09:05

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-019
 Units: ug/L

 Lab Code:
 R2005635-001
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/06/20 16:52	7/2/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/06/20 16:52	7/2/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/06/20 16:52	7/2/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/06/20 16:52	7/2/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/06/20 16:52	7/2/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/06/20 16:52	7/2/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/06/20 16:52	7/2/20	
Fluorene	9.1 U	9.1	1.2	1	07/06/20 16:52	7/2/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/06/20 16:52	7/2/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/06/20 16:52	7/2/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/06/20 16:52	7/2/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/06/20 16:52	7/2/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/06/20 16:52	7/2/20	
Isophorone	9.1 U	9.1	1.3	1	07/06/20 16:52	7/2/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/06/20 16:52	7/2/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/06/20 16:52	7/2/20	
Naphthalene	9.1 U	9.1	1.1	1	07/06/20 16:52	7/2/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/06/20 16:52	7/2/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/06/20 16:52	7/2/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/06/20 16:52	7/2/20	
Phenol	9.1 U	9.1	0.91	1	07/06/20 16:52	7/2/20	
Pyrene	9.1 U	9.1	1.3	1	07/06/20 16:52	7/2/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	121	35 - 141	07/06/20 16:52	
2-Fluorobiphenyl	73	31 - 118	07/06/20 16:52	
2-Fluorophenol	45	10 - 105	07/06/20 16:52	
Nitrobenzene-d5	71	31 - 110	07/06/20 16:52	
Phenol-d6	29	10 - 107	07/06/20 16:52	
p-Terphenyl-d14	51	10 - 165	07/06/20 16:52	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown	12.08	5.0	J	
	unknown	12.69	4.9	J	
	unknown	13.36	5.2	J	
	unknown	13.69	4.4	J	
	unknown	14.85	4.6	J	
	unknown hydrocarbon	16.34	4.0	J	
	unknown	7.21	3.8	J	

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Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 10:10

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-020
 Units: ug/L

 Lab Code:
 R2005635-002
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 17:22	7/2/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 17:22	7/2/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/06/20 17:22	7/2/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 17:22	7/2/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/06/20 17:22	7/2/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/06/20 17:22	7/2/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/06/20 17:22	7/2/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/06/20 17:22	7/2/20	
2,4-Dinitrophenol	45 U	45	19	1	07/06/20 17:22	7/2/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/06/20 17:22	7/2/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/06/20 17:22	7/2/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/06/20 17:22	7/2/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/06/20 17:22	7/2/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/06/20 17:22	7/2/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/06/20 17:22	7/2/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/06/20 17:22	7/2/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/06/20 17:22	7/2/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/06/20 17:22	7/2/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/06/20 17:22	7/2/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/06/20 17:22	7/2/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/06/20 17:22	7/2/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/06/20 17:22	7/2/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/06/20 17:22	7/2/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/06/20 17:22	7/2/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/06/20 17:22	7/2/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/06/20 17:22	7/2/20	
4-Nitrophenol	45 U	45	5.8	1	07/06/20 17:22	7/2/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/06/20 17:22	7/2/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/06/20 17:22	7/2/20	
Anthracene	9.1 U	9.1	1.2	1	07/06/20 17:22	7/2/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/06/20 17:22	7/2/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/06/20 17:22	7/2/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/06/20 17:22	7/2/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/06/20 17:22	7/2/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/06/20 17:22	7/2/20	
Benzoic Acid	91 U	91	33	1	07/06/20 17:22	7/2/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/06/20 17:22	7/2/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/06/20 17:22	7/2/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/06/20 17:22	7/2/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	07/06/20 17:22	7/2/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	07/06/20 17:22	7/2/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	07/06/20 17:22	7/2/20	
Chrysene	9.1 U	9.1	1.1	1	07/06/20 17:22	7/2/20	
,	<i>7.1</i> C	/.1		•	5., 55, 25 1,.22	20	

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Superset Reference: 20-0000555197 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 10:10

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-020
 Units: ug/L

 Lab Code:
 R2005635-002
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/06/20 17:22	7/2/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/06/20 17:22	7/2/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/06/20 17:22	7/2/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/06/20 17:22	7/2/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/06/20 17:22	7/2/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/06/20 17:22	7/2/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/06/20 17:22	7/2/20	
Fluorene	9.1 U	9.1	1.2	1	07/06/20 17:22	7/2/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/06/20 17:22	7/2/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/06/20 17:22	7/2/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/06/20 17:22	7/2/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/06/20 17:22	7/2/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/06/20 17:22	7/2/20	
Isophorone	9.1 U	9.1	1.3	1	07/06/20 17:22	7/2/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/06/20 17:22	7/2/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/06/20 17:22	7/2/20	
Naphthalene	9.1 U	9.1	1.1	1	07/06/20 17:22	7/2/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/06/20 17:22	7/2/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/06/20 17:22	7/2/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/06/20 17:22	7/2/20	
Phenol	9.1 U	9.1	0.91	1	07/06/20 17:22	7/2/20	
Pyrene	9.1 U	9.1	1.3	1	07/06/20 17:22	7/2/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	102	35 - 141	07/06/20 17:22	
2-Fluorobiphenyl	69	31 - 118	07/06/20 17:22	
2-Fluorophenol	44	10 - 105	07/06/20 17:22	
Nitrobenzene-d5	71	31 - 110	07/06/20 17:22	
Phenol-d6	31	10 - 107	07/06/20 17:22	
p-Terphenyl-d14	61	10 - 165	07/06/20 17:22	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown	12.08	6.1	J	
	unknown	12.69	7.2	J	
	unknown	13.36	7.2	J	
	unknown	14.08	6.3	J	
	unknown hydrocarbon	14.85	5.8	J	
	unknown	15.65	4.3	J	
013798-23-7	Sulfur	7.85	6.4	JN	

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Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 11:00

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-022
 Units: ug/L

 Lab Code:
 R2005635-003
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 17:52	7/2/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 17:52	7/2/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/06/20 17:52	7/2/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 17:52	7/2/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/06/20 17:52	7/2/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/06/20 17:52	7/2/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/06/20 17:52	7/2/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/06/20 17:52	7/2/20	
2,4-Dinitrophenol	45 U	45	19	1	07/06/20 17:52	7/2/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/06/20 17:52	7/2/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/06/20 17:52	7/2/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/06/20 17:52	7/2/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/06/20 17:52	7/2/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/06/20 17:52	7/2/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/06/20 17:52	7/2/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/06/20 17:52	7/2/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/06/20 17:52	7/2/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/06/20 17:52	7/2/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/06/20 17:52	7/2/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/06/20 17:52	7/2/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/06/20 17:52	7/2/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/06/20 17:52	7/2/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/06/20 17:52	7/2/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/06/20 17:52	7/2/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/06/20 17:52	7/2/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/06/20 17:52	7/2/20	
4-Nitrophenol	45 U	45	5.8	1	07/06/20 17:52	7/2/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/06/20 17:52	7/2/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/06/20 17:52	7/2/20	
Anthracene	9.1 U	9.1	1.2	1	07/06/20 17:52	7/2/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/06/20 17:52	7/2/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/06/20 17:52	7/2/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/06/20 17:52	7/2/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/06/20 17:52	7/2/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/06/20 17:52	7/2/20	
Benzoic Acid	91 U	91	33	1	07/06/20 17:52	7/2/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/06/20 17:52	7/2/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/06/20 17:52	7/2/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/06/20 17:52	7/2/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	07/06/20 17:52	7/2/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	07/06/20 17:52	7/2/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	07/06/20 17:52	7/2/20	
Chrysene	9.1 U	9.1	1.1	1	07/06/20 17:52	7/2/20	

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Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 11:00

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-022
 Units: ug/L

 Lab Code:
 R2005635-003
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/06/20 17:52	7/2/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/06/20 17:52	7/2/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/06/20 17:52	7/2/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/06/20 17:52	7/2/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/06/20 17:52	7/2/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/06/20 17:52	7/2/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/06/20 17:52	7/2/20	
Fluorene	9.1 U	9.1	1.2	1	07/06/20 17:52	7/2/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/06/20 17:52	7/2/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/06/20 17:52	7/2/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/06/20 17:52	7/2/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/06/20 17:52	7/2/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/06/20 17:52	7/2/20	
Isophorone	9.1 U	9.1	1.3	1	07/06/20 17:52	7/2/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/06/20 17:52	7/2/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/06/20 17:52	7/2/20	
Naphthalene	9.1 U	9.1	1.1	1	07/06/20 17:52	7/2/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/06/20 17:52	7/2/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/06/20 17:52	7/2/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/06/20 17:52	7/2/20	
Phenol	9.1 U	9.1	0.91	1	07/06/20 17:52	7/2/20	
Pyrene	9.1 U	9.1	1.3	1	07/06/20 17:52	7/2/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	118	35 - 141	07/06/20 17:52	
2-Fluorobiphenyl	80	31 - 118	07/06/20 17:52	
2-Fluorophenol	44	10 - 105	07/06/20 17:52	
Nitrobenzene-d5	71	31 - 110	07/06/20 17:52	
Phenol-d6	32	10 - 107	07/06/20 17:52	
p-Terphenyl-d14	64	10 - 165	07/06/20 17:52	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown	12.07	7.1	J	
	unknown hydrocarbon	12.69	8.2	J	
	unknown	13.11	4.6	J	
	unknown	13.36	8.7	J	
	unknown	14.08	7.5	J	
	unknown hydrocarbon	14.85	6.3	J	
	unknown	15.64	4.9	J	

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Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 11:00

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-022
 Units: ug/L

 Lab Code:
 R2005635-003
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown	16.34	4.4	J	
	unknown	6.18	4.2	J	
	unknown	7.84	4.4	J	
	unknown	7.92	3.9	J	
052253-93-7	Homomenthyl salicylate	9.37	4.5	JN	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 12:15

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-RM-023
 Units: ug/L

 Lab Code:
 R2005635-004
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 18:21	7/2/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 18:21	7/2/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/06/20 18:21	7/2/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/06/20 18:21	7/2/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/06/20 18:21	7/2/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/06/20 18:21	7/2/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/06/20 18:21	7/2/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/06/20 18:21	7/2/20	
2,4-Dinitrophenol	45 U	45	19	1	07/06/20 18:21	7/2/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/06/20 18:21	7/2/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/06/20 18:21	7/2/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/06/20 18:21	7/2/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/06/20 18:21	7/2/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/06/20 18:21	7/2/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/06/20 18:21	7/2/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/06/20 18:21	7/2/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/06/20 18:21	7/2/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/06/20 18:21	7/2/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/06/20 18:21	7/2/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/06/20 18:21	7/2/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/06/20 18:21	7/2/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/06/20 18:21	7/2/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/06/20 18:21	7/2/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/06/20 18:21	7/2/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/06/20 18:21	7/2/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/06/20 18:21	7/2/20	
4-Nitrophenol	45 U	45	5.8	1	07/06/20 18:21	7/2/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/06/20 18:21	7/2/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/06/20 18:21	7/2/20	
Anthracene	9.1 U	9.1	1.2	1	07/06/20 18:21	7/2/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/06/20 18:21	7/2/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/06/20 18:21	7/2/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/06/20 18:21	7/2/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/06/20 18:21	7/2/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/06/20 18:21	7/2/20	
Benzoic Acid	91 U	91	33	1	07/06/20 18:21	7/2/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/06/20 18:21	7/2/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/06/20 18:21	7/2/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/06/20 18:21	7/2/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	07/06/20 18:21	7/2/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	07/06/20 18:21	7/2/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	07/06/20 18:21	7/2/20	
Chrysene	9.1 U	9.1	1.1	1	07/06/20 18:21	7/2/20	
Cinybone	<i>)</i> .1 0	7.1	1.1	1	01/00/20 10.21	112120	

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Superset Reference: 20-0000555197 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 12:15

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-RM-023
 Units: ug/L

 Lab Code:
 R2005635-004
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/06/20 18:21	7/2/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/06/20 18:21	7/2/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/06/20 18:21	7/2/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/06/20 18:21	7/2/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/06/20 18:21	7/2/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/06/20 18:21	7/2/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/06/20 18:21	7/2/20	
Fluorene	9.1 U	9.1	1.2	1	07/06/20 18:21	7/2/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/06/20 18:21	7/2/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/06/20 18:21	7/2/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/06/20 18:21	7/2/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/06/20 18:21	7/2/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/06/20 18:21	7/2/20	
Isophorone	9.1 U	9.1	1.3	1	07/06/20 18:21	7/2/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/06/20 18:21	7/2/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/06/20 18:21	7/2/20	
Naphthalene	9.1 U	9.1	1.1	1	07/06/20 18:21	7/2/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/06/20 18:21	7/2/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/06/20 18:21	7/2/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/06/20 18:21	7/2/20	
Phenol	9.1 U	9.1	0.91	1	07/06/20 18:21	7/2/20	
Pyrene	9.1 U	9.1	1.3	1	07/06/20 18:21	7/2/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	102	35 - 141	07/06/20 18:21	
2-Fluorobiphenyl	74	31 - 118	07/06/20 18:21	
2-Fluorophenol	40	10 - 105	07/06/20 18:21	
Nitrobenzene-d5	72	31 - 110	07/06/20 18:21	
Phenol-d6	25	10 - 107	07/06/20 18:21	
p-Terphenyl-d14	67	10 - 165	07/06/20 18:21	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown	11.52	4.9	J	
	unknown hydrocarbon	12.07	7.1	J	
	unknown hydrocarbon	12.69	9.1	J	
	unknown hydrocarbon	13.35	10	J	
	unknown hydrocarbon	14.08	9.0	J	
	unknown	14.85	7.6	J	
	unknown hydrocarbon	15.65	6.3	J	

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Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 12:15

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-RM-023
 Units: ug/L

 Lab Code:
 R2005635-004
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Tentatively Identified Compounds

CAS# Compound Identification RT ug/L Q
unknown 16.34 5.1 J

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 13:00

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 RB-9954-063020-SG-002
 Units: ug/L

 Lab Code:
 R2005635-005
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/08/20 12:43	7/2/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/08/20 12:43	7/2/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/08/20 12:43	7/2/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/08/20 12:43	7/2/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/08/20 12:43	7/2/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/08/20 12:43	7/2/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/08/20 12:43	7/2/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/08/20 12:43	7/2/20	
2,4-Dinitrophenol	45 U	45	19	1	07/08/20 12:43	7/2/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/08/20 12:43	7/2/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/08/20 12:43	7/2/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/08/20 12:43	7/2/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/08/20 12:43	7/2/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/08/20 12:43	7/2/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/08/20 12:43	7/2/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/08/20 12:43	7/2/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/08/20 12:43	7/2/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/08/20 12:43	7/2/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/08/20 12:43	7/2/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/08/20 12:43	7/2/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/08/20 12:43	7/2/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/08/20 12:43	7/2/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/08/20 12:43	7/2/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/08/20 12:43	7/2/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/08/20 12:43	7/2/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/08/20 12:43	7/2/20	
4-Nitrophenol	45 U	45	5.8	1	07/08/20 12:43	7/2/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/08/20 12:43	7/2/20	
Acenaphthele Acenaphthylene	9.1 U	9.1	1.3	1	07/08/20 12:43	7/2/20	
Anthracene	9.1 U	9.1	1.3	1	07/08/20 12:43	7/2/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/08/20 12:43	7/2/20	
Benzo(a)pyrene	9.1 U	9.1	1.3	1	07/08/20 12:43	7/2/20	
Benzo(a)pyrene Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/08/20 12:43	7/2/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/08/20 12:43	7/2/20	
	9.1 U 9.1 U	9.1	1.1	1		7/2/20	
Benzo(k)fluoranthene Benzoic Acid	9.1 U 91 U	9.1	33	1	07/08/20 12:43 07/08/20 12:43	7/2/20	
	9.1 U	91 9.1	33 1.5	1	07/08/20 12:43	7/2/20	
Benzyl Alcohol			1.3				
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1		1	07/08/20 12:43	7/2/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/08/20 12:43	7/2/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	07/08/20 12:43	7/2/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	07/08/20 12:43	7/2/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	07/08/20 12:43	7/2/20	
Chrysene	9.1 U	9.1	1.1	1	07/08/20 12:43	7/2/20	

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Superset Reference: 20-0000555197 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 13:00

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 RB-9954-063020-SG-002
 Units: ug/L

 Lab Code:
 R2005635-005
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/08/20 12:43	7/2/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/08/20 12:43	7/2/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/08/20 12:43	7/2/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/08/20 12:43	7/2/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/08/20 12:43	7/2/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/08/20 12:43	7/2/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/08/20 12:43	7/2/20	
Fluorene	9.1 U	9.1	1.2	1	07/08/20 12:43	7/2/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/08/20 12:43	7/2/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/08/20 12:43	7/2/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/08/20 12:43	7/2/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/08/20 12:43	7/2/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/08/20 12:43	7/2/20	
Isophorone	9.1 U	9.1	1.3	1	07/08/20 12:43	7/2/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/08/20 12:43	7/2/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/08/20 12:43	7/2/20	
Naphthalene	9.1 U	9.1	1.1	1	07/08/20 12:43	7/2/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/08/20 12:43	7/2/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/08/20 12:43	7/2/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/08/20 12:43	7/2/20	
Phenol	0.94 J	9.1	0.91	1	07/08/20 12:43	7/2/20	
Pyrene	9.1 U	9.1	1.3	1	07/08/20 12:43	7/2/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	98	35 - 141	07/08/20 12:43	
2-Fluorobiphenyl	68	31 - 118	07/08/20 12:43	
2-Fluorophenol	39	10 - 105	07/08/20 12:43	
Nitrobenzene-d5	65	31 - 110	07/08/20 12:43	
Phenol-d6	27	10 - 107	07/08/20 12:43	
p-Terphenyl-d14	62	10 - 165	07/08/20 12:43	

Tentatively Identified Compounds

	CAS# Compound Identification		Result		
CAS#		RT	ug/L	Q	
	unknown	12.69	5.7	J	
	unknown	13.35	6.4	J	
	unknown	14.08	4.3	J	
	unknown	14.85	4.8	J	
	unknown	15.65	4.0	J	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 10:10

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-021
 Units: ug/L

 Lab Code:
 R2005635-007
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed Da	te Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/08/20 13:12	7/2/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/08/20 13:12	7/2/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/08/20 13:12	7/2/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/08/20 13:12	7/2/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/08/20 13:12	7/2/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/08/20 13:12	7/2/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/08/20 13:12	7/2/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/08/20 13:12	7/2/20	
2,4-Dinitrophenol	45 U	45	19	1	07/08/20 13:12	7/2/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/08/20 13:12	7/2/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/08/20 13:12	7/2/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/08/20 13:12	7/2/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/08/20 13:12	7/2/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/08/20 13:12	7/2/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/08/20 13:12	7/2/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/08/20 13:12	7/2/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/08/20 13:12	7/2/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/08/20 13:12	7/2/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/08/20 13:12	7/2/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/08/20 13:12	7/2/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/08/20 13:12	7/2/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/08/20 13:12	7/2/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/08/20 13:12	7/2/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/08/20 13:12	7/2/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/08/20 13:12	7/2/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/08/20 13:12	7/2/20	
4-Nitrophenol	45 U	45	5.8	1	07/08/20 13:12	7/2/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/08/20 13:12	7/2/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/08/20 13:12	7/2/20	
Anthracene	9.1 U	9.1	1.2	1	07/08/20 13:12	7/2/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/08/20 13:12	7/2/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/08/20 13:12	7/2/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/08/20 13:12	7/2/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/08/20 13:12	7/2/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/08/20 13:12	7/2/20	
Benzoic Acid	91 U	91	33	1	07/08/20 13:12	7/2/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/08/20 13:12	7/2/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/08/20 13:12	7/2/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/08/20 13:12	7/2/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.3	1	07/08/20 13:12	7/2/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	07/08/20 13:12	7/2/20	-
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	07/08/20 13:12	7/2/20	
Chrysene	9.1 U 9.1 U	9.1	1.3	1	07/08/20 13:12	7/2/20	
CIII y SCIIC	9.1 U	7.1	1.1	1	07/00/20 13.12	112120	

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Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 10:10

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-021
 Units: ug/L

 Lab Code:
 R2005635-007
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed I	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/08/20 13:12	7/2/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/08/20 13:12	7/2/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/08/20 13:12	7/2/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/08/20 13:12	7/2/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/08/20 13:12	7/2/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/08/20 13:12	7/2/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/08/20 13:12	7/2/20	
Fluorene	9.1 U	9.1	1.2	1	07/08/20 13:12	7/2/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/08/20 13:12	7/2/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/08/20 13:12	7/2/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/08/20 13:12	7/2/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/08/20 13:12	7/2/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/08/20 13:12	7/2/20	
Isophorone	9.1 U	9.1	1.3	1	07/08/20 13:12	7/2/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/08/20 13:12	7/2/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/08/20 13:12	7/2/20	
Naphthalene	9.1 U	9.1	1.1	1	07/08/20 13:12	7/2/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/08/20 13:12	7/2/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/08/20 13:12	7/2/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/08/20 13:12	7/2/20	
Phenol	1.1 J	9.1	0.91	1	07/08/20 13:12	7/2/20	
Pyrene	9.1 U	9.1	1.3	1	07/08/20 13:12	7/2/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	114	35 - 141	07/08/20 13:12	
2-Fluorobiphenyl	83	31 - 118	07/08/20 13:12	
2-Fluorophenol	50	10 - 105	07/08/20 13:12	
Nitrobenzene-d5	83	31 - 110	07/08/20 13:12	
Phenol-d6	33	10 - 107	07/08/20 13:12	
p-Terphenyl-d14	54	10 - 165	07/08/20 13:12	

Tentatively Identified Compounds

	Result		
RT	ug/L	Q	
12.69	4.1	J	
13.11	6.9	J	
7.86	4.1	J	
7.93	5.5	J	
	12.69 13.11 7.86	12.69 4.1 13.11 6.9 7.86 4.1	RT ug/L Q 12.69 4.1 J 13.11 6.9 J 7.86 4.1 J

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Semivolatile Organic Compounds by GC

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Date Collected: 06/30/20 09:05

Monitoring Sample Matrix: Water

Lab Code:

Date Received: 07/01/20 11:05

Sample Name: WG-9954-063020-SG-019

R2005635-001

Units: ug/L Basis: NA

Dusis

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/08/20 17:27	7/2/20	
4,4'-DDE	0.045 U	0.045	0.019	1	07/08/20 17:27	7/2/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/08/20 17:27	7/2/20	
Aldrin	0.045 U	0.045	0.019	1	07/08/20 17:27	7/2/20	
Dieldrin	0.045 U	0.045	0.019	1	07/08/20 17:27	7/2/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/08/20 17:27	7/2/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/08/20 17:27	7/2/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/08/20 17:27	7/2/20	
Endrin	0.045 U	0.045	0.019	1	07/08/20 17:27	7/2/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/08/20 17:27	7/2/20	
Heptachlor	0.045 U	0.045	0.019	1	07/08/20 17:27	7/2/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/08/20 17:27	7/2/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/08/20 17:27	7/2/20	
Toxaphene	0.46 U	0.46	0.46	1	07/08/20 17:27	7/2/20	
alpha-BHC	0.12	0.045	0.019	1	07/08/20 17:27	7/2/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/08/20 17:27	7/2/20	
beta-BHC	0.045 U	0.045	0.019	1	07/08/20 17:27	7/2/20	
delta-BHC	0.076	0.045	0.019	1	07/08/20 17:27	7/2/20	
gamma-BHC (Lindane)	0.10	0.045	0.019	1	07/08/20 17:27	7/2/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/08/20 17:27	7/2/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	72	10 - 164	07/08/20 17:27	
Tetrachloro-m-xylene	75	10 - 147	07/08/20 17:27	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 10:10

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-020
 Units: ug/L

 Lab Code:
 R2005635-002
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/07/20 13:08	7/2/20	
4,4'-DDE	0.045 U	0.045	0.019	1	07/07/20 13:08	7/2/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/07/20 13:08	7/2/20	
Aldrin	0.045 U	0.045	0.019	1	07/07/20 13:08	7/2/20	
Dieldrin	0.045 U	0.045	0.019	1	07/07/20 13:08	7/2/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/07/20 13:08	7/2/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/07/20 13:08	7/2/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/07/20 13:08	7/2/20	
Endrin	0.045 U	0.045	0.019	1	07/07/20 13:08	7/2/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/07/20 13:08	7/2/20	
Heptachlor	0.045 U	0.045	0.019	1	07/07/20 13:08	7/2/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/07/20 13:08	7/2/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/07/20 13:08	7/2/20	
Toxaphene	0.46 U	0.46	0.46	1	07/07/20 13:08	7/2/20	
alpha-BHC	0.081	0.045	0.019	1	07/07/20 13:08	7/2/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/07/20 13:08	7/2/20	
beta-BHC	0.045 U	0.045	0.019	1	07/07/20 13:08	7/2/20	
delta-BHC	0.068	0.045	0.019	1	07/07/20 13:08	7/2/20	
gamma-BHC (Lindane)	0.089	0.045	0.019	1	07/07/20 13:08	7/2/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/07/20 13:08	7/2/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	74	10 - 164	07/07/20 13:08	
Tetrachloro-m-xylene	59	10 - 147	07/07/20 13:08	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 11:00

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-022
 Units: ug/L

 Lab Code:
 R2005635-003
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/07/20 13:27	7/2/20	_
4,4'-DDE	0.045 U	0.045	0.019	1	07/07/20 13:27	7/2/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/07/20 13:27	7/2/20	
Aldrin	0.045 U	0.045	0.019	1	07/07/20 13:27	7/2/20	
Dieldrin	0.045 U	0.045	0.019	1	07/07/20 13:27	7/2/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/07/20 13:27	7/2/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/07/20 13:27	7/2/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/07/20 13:27	7/2/20	
Endrin	0.045 U	0.045	0.019	1	07/07/20 13:27	7/2/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/07/20 13:27	7/2/20	
Heptachlor	0.045 U	0.045	0.019	1	07/07/20 13:27	7/2/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/07/20 13:27	7/2/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/07/20 13:27	7/2/20	
Toxaphene	0.46 U	0.46	0.46	1	07/07/20 13:27	7/2/20	
alpha-BHC	0.064	0.045	0.019	1	07/07/20 13:27	7/2/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/07/20 13:27	7/2/20	
beta-BHC	0.027 J	0.045	0.019	1	07/07/20 13:27	7/2/20	
delta-BHC	0.29	0.045	0.019	1	07/07/20 13:27	7/2/20	
gamma-BHC (Lindane)	0.074	0.045	0.019	1	07/07/20 13:27	7/2/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/07/20 13:27	7/2/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	52	10 - 164	07/07/20 13:27	
Tetrachloro-m-xylene	66	10 - 147	07/07/20 13:27	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 12:15

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-RM-023
 Units: ug/L

 Lab Code:
 R2005635-004
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/07/20 13:46	7/2/20	_
4,4'-DDE	0.045 U	0.045	0.019	1	07/07/20 13:46	7/2/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/07/20 13:46	7/2/20	
Aldrin	0.045 U	0.045	0.019	1	07/07/20 13:46	7/2/20	
Dieldrin	0.045 U	0.045	0.019	1	07/07/20 13:46	7/2/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/07/20 13:46	7/2/20	<u>.</u>
Endosulfan II	0.045 U	0.045	0.019	1	07/07/20 13:46	7/2/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/07/20 13:46	7/2/20	
Endrin	0.045 U	0.045	0.019	1	07/07/20 13:46	7/2/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/07/20 13:46	7/2/20	
Heptachlor	0.045 U	0.045	0.019	1	07/07/20 13:46	7/2/20	<u>.</u>
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/07/20 13:46	7/2/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/07/20 13:46	7/2/20	
Toxaphene	0.46 U	0.46	0.46	1	07/07/20 13:46	7/2/20	
alpha-BHC	0.045 U	0.045	0.019	1	07/07/20 13:46	7/2/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/07/20 13:46	7/2/20	
beta-BHC	0.045 U	0.045	0.019	1	07/07/20 13:46	7/2/20	
delta-BHC	0.045 U	0.045	0.019	1	07/07/20 13:46	7/2/20	
gamma-BHC (Lindane)	0.045 U	0.045	0.019	1	07/07/20 13:46	7/2/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/07/20 13:46	7/2/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	18	10 - 164	07/07/20 13:46	
Tetrachloro-m-xylene	58	10 - 147	07/07/20 13:46	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 13:00

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 RB-9954-063020-SG-002
 Units: ug/L

 Lab Code:
 R2005635-005
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/06/20 17:42	7/2/20	
4,4'-DDE	0.045 U	0.045	0.019	1	07/06/20 17:42	7/2/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/06/20 17:42	7/2/20	
Aldrin	0.045 U	0.045	0.019	1	07/06/20 17:42	7/2/20	
Dieldrin	0.045 U	0.045	0.019	1	07/06/20 17:42	7/2/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/06/20 17:42	7/2/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/06/20 17:42	7/2/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/06/20 17:42	7/2/20	
Endrin	0.045 U	0.045	0.019	1	07/06/20 17:42	7/2/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/06/20 17:42	7/2/20	
Heptachlor	0.045 U	0.045	0.019	1	07/06/20 17:42	7/2/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/06/20 17:42	7/2/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/06/20 17:42	7/2/20	
Toxaphene	0.46 U	0.46	0.46	1	07/06/20 17:42	7/2/20	
alpha-BHC	0.063 Ui	0.063	0.063	1	07/06/20 17:42	7/2/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/06/20 17:42	7/2/20	
beta-BHC	0.045 U	0.045	0.019	1	07/06/20 17:42	7/2/20	
delta-BHC	0.037 J	0.045	0.019	1	07/06/20 17:42	7/2/20	
gamma-BHC (Lindane)	0.058	0.045	0.019	1	07/06/20 17:42	7/2/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/06/20 17:42	7/2/20	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 13:00

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 RB-9954-063020-SG-002
 Units: ug/L

 Lab Code:
 R2005635-005
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	9 *	10 - 164	07/06/20 17:42	*
Tetrachloro-m-xylene	55	10 - 147	07/06/20 17:42	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 13:00

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 RB-9954-063020-SG-002
 Units: ug/L

 Lab Code:
 R2005635-005
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/13/20 15:58	7/10/20	*
4,4'-DDE	0.045 U	0.045	0.019	1	07/13/20 15:58	7/10/20	*
4,4'-DDT	0.045 U	0.045	0.019	1	07/13/20 15:58	7/10/20	*
Aldrin	0.045 U	0.045	0.019	1	07/13/20 15:58	7/10/20	*
Dieldrin	0.045 U	0.045	0.019	1	07/13/20 15:58	7/10/20	*
Endosulfan I	0.045 U	0.045	0.019	1	07/13/20 15:58	7/10/20	*
Endosulfan II	0.045 U	0.045	0.019	1	07/13/20 15:58	7/10/20	*
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/13/20 15:58	7/10/20	*
Endrin	0.045 U	0.045	0.019	1	07/13/20 15:58	7/10/20	*
Endrin Ketone	0.045 U	0.045	0.019	1	07/13/20 15:58	7/10/20	*
Heptachlor	0.045 U	0.045	0.019	1	07/13/20 15:58	7/10/20	*
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/13/20 15:58	7/10/20	*
Methoxychlor	0.045 U	0.045	0.019	1	07/13/20 15:58	7/10/20	*
Toxaphene	0.46 U	0.46	0.46	1	07/13/20 15:58	7/10/20	*
alpha-BHC	0.059 Ui	0.059	0.059	1	07/13/20 15:58	7/10/20	*
alpha-Chlordane	0.045 U	0.045	0.019	1	07/13/20 15:58	7/10/20	*
beta-BHC	0.045 U	0.045	0.019	1	07/13/20 15:58	7/10/20	*
delta-BHC	0.026 J	0.045	0.019	1	07/13/20 15:58	7/10/20	*
gamma-BHC (Lindane)	0.053	0.045	0.019	1	07/13/20 15:58	7/10/20	*
gamma-Chlordane	0.045 U	0.045	0.019	1	07/13/20 15:58	7/10/20	*

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 13:00

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 RB-9954-063020-SG-002
 Units: ug/L

 Lab Code:
 R2005635-005
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analysis Method: 8081B

Prep Method: EPA 3510C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	9 *	10 - 164	07/13/20 15:58	*
Tetrachloro-m-xylene	36	10 - 147	07/13/20 15:58	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 10:10

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-021
 Units: ug/L

 Lab Code:
 R2005635-007
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/08/20 18:24	7/2/20	
4,4'-DDE	0.045 U	0.045	0.019	1	07/08/20 18:24	7/2/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/08/20 18:24	7/2/20	
Aldrin	0.045 U	0.045	0.019	1	07/08/20 18:24	7/2/20	
Dieldrin	0.045 U	0.045	0.019	1	07/08/20 18:24	7/2/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/08/20 18:24	7/2/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/08/20 18:24	7/2/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/08/20 18:24	7/2/20	
Endrin	0.045 U	0.045	0.019	1	07/08/20 18:24	7/2/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/08/20 18:24	7/2/20	
Heptachlor	0.045 U	0.045	0.019	1	07/08/20 18:24	7/2/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/08/20 18:24	7/2/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/08/20 18:24	7/2/20	
Toxaphene	0.46 U	0.46	0.46	1	07/08/20 18:24	7/2/20	
alpha-BHC	0.10	0.045	0.019	1	07/08/20 18:24	7/2/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/08/20 18:24	7/2/20	
beta-BHC	0.045 U	0.045	0.019	1	07/08/20 18:24	7/2/20	
delta-BHC	0.12	0.045	0.019	1	07/08/20 18:24	7/2/20	
gamma-BHC (Lindane)	0.11	0.045	0.019	1	07/08/20 18:24	7/2/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/08/20 18:24	7/2/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	85	10 - 164	07/08/20 18:24	
Tetrachloro-m-xylene	79	10 - 147	07/08/20 18:24	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 09:05

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-019
 Units: ug/L

 Lab Code:
 R2005635-001
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/08/20 21:12	7/2/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/08/20 21:12	7/2/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/08/20 21:12	7/2/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/08/20 21:12	7/2/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/08/20 21:12	7/2/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/08/20 21:12	7/2/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/08/20 21:12	7/2/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	103	10 - 152	07/08/20 21:12	
Tetrachloro-m-xvlene	52	14 - 129	07/08/20 21:12	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 10:10

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-020
 Units: ug/L

 Lab Code:
 R2005635-002
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/08/20 21:33	7/2/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/08/20 21:33	7/2/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/08/20 21:33	7/2/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/08/20 21:33	7/2/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/08/20 21:33	7/2/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/08/20 21:33	7/2/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/08/20 21:33	7/2/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	78	10 - 152	07/08/20 21:33	
Tetrachloro-m-xylene	41	14 - 129	07/08/20 21:33	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 11:00

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-022
 Units: ug/L

 Lab Code:
 R2005635-003
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/08/20 11:16	7/2/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/08/20 11:16	7/2/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/08/20 11:16	7/2/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/08/20 11:16	7/2/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/08/20 11:16	7/2/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/08/20 11:16	7/2/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/08/20 11:16	7/2/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q	
Decachlorobiphenyl	42	10 - 152	07/08/20 11:16		
Tetrachloro-m-xvlene	45	14 - 129	07/08/20 11:16		

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 12:15

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-RM-023
 Units: ug/L

 Lab Code:
 R2005635-004
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/08/20 11:36	7/2/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/08/20 11:36	7/2/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/08/20 11:36	7/2/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/08/20 11:36	7/2/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/08/20 11:36	7/2/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/08/20 11:36	7/2/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/08/20 11:36	7/2/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	15	10 - 152	07/08/20 11:36	
Tetrachloro-m-xylene	40	14 - 129	07/08/20 11:36	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 13:00

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 RB-9954-063020-SG-002
 Units: ug/L

 Lab Code:
 R2005635-005
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/08/20 12:37	7/2/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/08/20 12:37	7/2/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/08/20 12:37	7/2/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/08/20 12:37	7/2/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/08/20 12:37	7/2/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/08/20 12:37	7/2/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/08/20 12:37	7/2/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q	
Decachlorobiphenyl	9 *	10 - 152	07/08/20 12:37	*	
Tetrachloro-m-xylene	40	14 - 129	07/08/20 12:37		

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 13:00

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 RB-9954-063020-SG-002
 Units: ug/L

 Lab Code:
 R2005635-005
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/13/20 15:14	7/10/20	*
Aroclor 1221	1.8 U	1.8	0.91	1	07/13/20 15:14	7/10/20	*
Aroclor 1232	0.91 U	0.91	0.46	1	07/13/20 15:14	7/10/20	*
Aroclor 1242	0.91 U	0.91	0.46	1	07/13/20 15:14	7/10/20	*
Aroclor 1248	0.91 U	0.91	0.46	1	07/13/20 15:14	7/10/20	*
Aroclor 1254	0.91 U	0.91	0.46	1	07/13/20 15:14	7/10/20	*
Aroclor 1260	0.91 U	0.91	0.46	1	07/13/20 15:14	7/10/20	*

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q	
Decachlorobiphenyl	9 *	10 - 152	07/13/20 15:14	*	
Tetrachloro-m-xylene	31	14 - 129	07/13/20 15:14		

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 06/30/20 10:10

Monitoring

Sample Matrix: Water Date Received: 07/01/20 11:05

 Sample Name:
 WG-9954-063020-SG-021
 Units: ug/L

 Lab Code:
 R2005635-007
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/08/20 22:33	7/2/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/08/20 22:33	7/2/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/08/20 22:33	7/2/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/08/20 22:33	7/2/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/08/20 22:33	7/2/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/08/20 22:33	7/2/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/08/20 22:33	7/2/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	95	10 - 152	07/08/20 22:33	
Tetrachloro-m-xvlene	54	14 - 129	07/08/20 22:33	



QC Summary Forms

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com



Volatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

QA/QC Report

Service Request: R2005635

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

Project:

SURROGATE RECOVERY SUMMARY Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Extraction Method: EPA 5030C

		4-Bromofluorobenzene	Dibromofluoromethane	Toluene-d8
Sample Name	Sample Name Lab Code		89-119	87-121
WG-9954-063020-SG-019	R2005635-001	94	100	101
WG-9954-063020-SG-020	R2005635-002	97	105	103
WG-9954-063020-SG-022	R2005635-003	95	102	102
WG-9954-063020-RM-023	R2005635-004	90	99	96
RB-9954-063020-SG-002	R2005635-005	94	99	101
TB-9954-063020-SG-004	R2005635-006	94	98	102
WG-9954-063020-SG-021	R2005635-007	95	100	101
Method Blank	RQ2007501-04	95	99	101
Lab Control Sample	RQ2007501-03	103	103	105
Method Blank	RQ2007451-04	95	99	102
Lab Control Sample	RQ2007451-03	99	100	101
WG-9954-063020-RM-023 MS	RQ2007451-05	96	96	98
WG-9954-063020-RM-023 DMS	RQ2007451-06	104	107	104

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Date Collected:
Date Received:

06/30/20 07/01/20

Sample Matrix: Water

Data Analyza

Date Analyzed: 07/14/20

Duplicate Matrix Spike Summary Volatile Organic Compounds by GC/MS

Sample Name: WG-9954-063020-RM-023

Units: ug/L

Lab Code: R2005635-004

Basis: NA

Analysis Method: 8260C

		M	atrix Spike		Dupl	icate Matrix	k Spike			
		RQ	2007451-05]	RQ2007451-	06			
	Sample		Spike			Spike		% Rec		RPD
Analyte Name	Result	Result	Amount	% Rec	Result	Amount	% Rec	Limits	RPD	Limit
1,1,1-Trichloroethane (TCA)	5.0 U	51.1	50.0	102	56.3	50.0	113	74-127	10	30
1,1,2,2-Tetrachloroethane	5.0 U	51.3	50.0	103	56.5	50.0	113	72-122	10	30
1,1,2-Trichloroethane	5.0 U	48.8	50.0	98	52.9	50.0	106	82-121	8	30
1,1-Dichloroethane (1,1-DCA)	5.0 U	49.1	50.0	98	53.8	50.0	108	74-132	9	30
1,1-Dichloroethene (1,1-DCE)	5.0 U	57.1	50.0	114	64.9	50.0	130 *	71-118	13	30
1,2-Dichloroethane	5.0 U	46.2	50.0	92	51.1	50.0	102	68-130	10	30
1,2-Dichloropropane	5.0 U	47.9	50.0	96	54.3	50.0	109	79-124	12	30
2-Butanone (MEK)	10 U	51.9	50.0	104	54.5	50.0	109	61-137	5	30
2-Hexanone	10 U	46.0	50.0	92	48.7	50.0	97	56-132	6	30
4-Methyl-2-pentanone	10 U	46.8	50.0	94	53.4	50.0	107	60-141	13	30
Acetone	10 U	50.4	50.0	101	53.3	50.0	107	35-183	6	30
Benzene	5.0 U	48.9	50.0	98	54.3	50.0	109	76-129	10	30
Bromodichloromethane	5.0 U	48.4	50.0	97	53.5	50.0	107	78-133	10	30
Bromoform	5.0 U	46.1	50.0	92	51.3	50.0	103	58-133	11	30
Bromomethane	5.0 U	41.9	50.0	84	39.3	50.0	79	10-184	6	30
Carbon Disulfide	2.9 J	57.7	50.0	110	63.6	50.0	121	59-140	10	30
Carbon Tetrachloride	5.0 U	49.8	50.0	100	57.5	50.0	115	65-135	14	30
Chlorobenzene	5.0 U	47.5	50.0	95	52.1	50.0	104	76-125	9	30
Chloroethane	5.0 U	43.6	50.0	87	46.5	50.0	93	48-146	6	30
Chloroform	5.0 U	49.0	50.0	98	55.4	50.0	111	75-130	12	30
Chloromethane	5.0 U	54.0	50.0	108	57.1	50.0	114	55-160	6	30
Dibromochloromethane	5.0 U	51.2	50.0	102	57.6	50.0	115	72-128	12	30
Dichloromethane	5.0 U	47.1	50.0	94	51.6	50.0	103	73-122	9	30
Ethylbenzene	5.0 U	49.5	50.0	99	53.2	50.0	106	72-134	7	30
Styrene	5.0 U	45.0	50.0	90	48.9	50.0	98	74-136	8	30
Tetrachloroethene (PCE)	5.0 U	45.5	50.0	91	50.8	50.0	102	72-125	11	30
Toluene	0.31 BJ	52.1	50.0	104	56.6	50.0	113	79-119	8	30
Trichloroethene (TCE)	5.0 U	43.9	50.0	88	48.8	50.0	98	74-122	10	30
Vinyl Acetate	10 U	67.3	50.0	135	75.4	50.0	151	48-172	11	30
Vinyl Chloride	5.0 U	49.6	50.0	99	52.8	50.0	106	74-159	6	30
cis-1,2-Dichloroethene	5.0 U	49.3	50.0	99	54.0	50.0	108	77-127	9	30
cis-1,3-Dichloropropene	5.0 U	46.7	50.0	93	51.2	50.0	102	52-134	9	30
trans-1,2-Dichloroethene	5.0 U	54.2	50.0	108	61.0	50.0	122 *	73-118	12	30

Results flagged with an asterisk (*) indicate values outside control criteria.

Results flagged with a pound (#) indicate the control criteria is not applicable.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

Matrix Spike and Matrix Spike Duplicate Data is presented for information purposes only. The matrix may or may not be relevant to samples reported in this report. The laboratory evaluates system performance based on the LCS and LCSD control limits.

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request:

R2005635

Project: Sample Matrix: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Date Collected:

06/30/20

Water

Date Received:

07/01/20 07/14/20

Date Analyzed:

·

Duplicate Matrix Spike Summary Volatile Organic Compounds by GC/MS

WG-9954-063020-RM-023

Units:

ug/L

Lab Code:

Sample Name:

R2005635-004

Basis:

NA

Analysis Method:

8260C

Matrix Spike RQ2007451-05 **Duplicate Matrix Spike**

RQ2007451-06

	Sample		Spike			Spike		% Rec		RPD
Analyte Name	Result	Result	Amount	% Rec	Result	Amount	% Rec	Limits	RPD	Limit
trans-1.3-Dichloropropene	5.0 U	45.2	50.0	90	50.8	50.0	102	71-133	12	30

Results flagged with an asterisk (*) indicate values outside control criteria.

Results flagged with a pound (#) indicate the control criteria is not applicable.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

Matrix Spike and Matrix Spike Duplicate Data is presented for information purposes only. The matrix may or may not be relevant to samples reported in this report. The laboratory evaluates system performance based on the LCS and LCSD control limits.

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007451-04
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/13/20 19:02	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/13/20 19:02	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/13/20 19:02	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/13/20 19:02	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/13/20 19:02	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/13/20 19:02	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/13/20 19:02	
2-Butanone (MEK)	10 U	10	0.78	1	07/13/20 19:02	
2-Hexanone	10 U	10	0.20	1	07/13/20 19:02	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/13/20 19:02	
Acetone	10 U	10	5.0	1	07/13/20 19:02	
Benzene	5.0 U	5.0	0.20	1	07/13/20 19:02	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/13/20 19:02	
Bromoform	5.0 U	5.0	0.25	1	07/13/20 19:02	
Bromomethane	5.0 U	5.0	0.70	1	07/13/20 19:02	
Carbon Disulfide	10 U	10	0.42	1	07/13/20 19:02	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/13/20 19:02	
Chlorobenzene	5.0 U	5.0	0.20	1	07/13/20 19:02	
Chloroethane	5.0 U	5.0	0.23	1	07/13/20 19:02	
Chloroform	5.0 U	5.0	0.24	1	07/13/20 19:02	
Chloromethane	0.32 J	5.0	0.28	1	07/13/20 19:02	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/13/20 19:02	
Dichloromethane	5.0 U	5.0	0.65	1	07/13/20 19:02	
Ethylbenzene	5.0 U	5.0	0.20	1	07/13/20 19:02	
Styrene	5.0 U	5.0	0.20	1	07/13/20 19:02	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/13/20 19:02	
Toluene	5.0 U	5.0	0.20	1	07/13/20 19:02	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/13/20 19:02	
Vinyl Acetate	10 U	10	1.1	1	07/13/20 19:02	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/13/20 19:02	
Xylenes, Total	5.0 U	5.0	0.23	1	07/13/20 19:02	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/13/20 19:02	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/13/20 19:02	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/13/20 19:02	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/13/20 19:02	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007451-04Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	95	85 - 122	07/13/20 19:02	
Dibromofluoromethane	99	89 - 119	07/13/20 19:02	
Toluene-d8	102	87 - 121	07/13/20 19:02	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007451-04Basis: NA

Volatile Organic Compounds by GC/MS

, onto organic compounds of contra

Analysis Method: 8260C

Tentatively Identified Compounds

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007501-04
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/14/20 14:25	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/14/20 14:25	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/14/20 14:25	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/14/20 14:25	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/14/20 14:25	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/14/20 14:25	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/14/20 14:25	
2-Butanone (MEK)	10 U	10	0.78	1	07/14/20 14:25	
2-Hexanone	10 U	10	0.20	1	07/14/20 14:25	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/14/20 14:25	
Acetone	10 U	10	5.0	1	07/14/20 14:25	
Benzene	5.0 U	5.0	0.20	1	07/14/20 14:25	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/14/20 14:25	
Bromoform	5.0 U	5.0	0.25	1	07/14/20 14:25	
Bromomethane	5.0 U	5.0	0.70	1	07/14/20 14:25	
Carbon Disulfide	10 U	10	0.42	1	07/14/20 14:25	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/14/20 14:25	
Chlorobenzene	5.0 U	5.0	0.20	1	07/14/20 14:25	
Chloroethane	5.0 U	5.0	0.23	1	07/14/20 14:25	
Chloroform	5.0 U	5.0	0.24	1	07/14/20 14:25	
Chloromethane	5.0 U	5.0	0.28	1	07/14/20 14:25	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/14/20 14:25	
Dichloromethane	5.0 U	5.0	0.65	1	07/14/20 14:25	
Ethylbenzene	5.0 U	5.0	0.20	1	07/14/20 14:25	
Styrene	5.0 U	5.0	0.20	1	07/14/20 14:25	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/14/20 14:25	
Toluene	0.41 J	5.0	0.20	1	07/14/20 14:25	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/14/20 14:25	
Vinyl Acetate	10 U	10	1.1	1	07/14/20 14:25	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/14/20 14:25	
Xylenes, Total	5.0 U	5.0	0.23	1	07/14/20 14:25	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/14/20 14:25	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/14/20 14:25	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/14/20 14:25	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/14/20 14:25	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007501-04Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	95	85 - 122	07/14/20 14:25	
Dibromofluoromethane	99	89 - 119	07/14/20 14:25	
Toluene-d8	101	87 - 121	07/14/20 14:25	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Love Canal:292-402-D02-3100/9954 Annual Long Term **Project: Date Collected:** NA

Monitoring

Date Received: NA **Sample Matrix:** Water

Sample Name: Method Blank Units: ug/L Lab Code:

RQ2007501-04 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

Result CAS# **Compound Identification** RTug/L unknown 1.61

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/13/20

Sample Matrix: Water

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005635

Lab Control Sample

RQ2007451-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
1,1,1-Trichloroethane (TCA)	8260C	21.0	20.0	105	75-125
1,1,2,2-Tetrachloroethane	8260C	23.6	20.0	118	78-126
1,1,2-Trichloroethane	8260C	21.4	20.0	107	82-121
1,1-Dichloroethane (1,1-DCA)	8260C	20.6	20.0	103	80-124
1,1-Dichloroethene (1,1-DCE)	8260C	24.0	20.0	120 *	71-118
1,2-Dichloroethane	8260C	20.8	20.0	104	71-127
1,2-Dichloropropane	8260C	21.3	20.0	106	80-119
2-Butanone (MEK)	8260C	23.3	20.0	116	61-137
2-Hexanone	8260C	20.5	20.0	102	63-124
4-Methyl-2-pentanone	8260C	21.0	20.0	105	66-124
Acetone	8260C	23.4	20.0	117	40-161
Benzene	8260C	21.2	20.0	106	79-119
Bromodichloromethane	8260C	21.1	20.0	105	81-123
Bromoform	8260C	21.4	20.0	107	65-146
Bromomethane	8260C	23.4	20.0	117	42-166
Carbon Disulfide	8260C	21.0	20.0	105	66-128
Carbon Tetrachloride	8260C	21.5	20.0	108	70-127
Chlorobenzene	8260C	21.1	20.0	106	80-121
Chloroethane	8260C	18.6	20.0	93	62-131
Chloroform	8260C	20.5	20.0	103	79-120
Chloromethane	8260C	22.9	20.0	115	65-135
Dibromochloromethane	8260C	22.6	20.0	113	72-128
Dichloromethane	8260C	20.3	20.0	102	73-122
Ethylbenzene	8260C	20.7	20.0	103	76-120
Styrene	8260C	21.1	20.0	106	80-124
Tetrachloroethene (PCE)	8260C	19.1	20.0	95	72-125
Toluene	8260C	21.4	20.0	107	79-119
Trichloroethene (TCE)	8260C	19.6	20.0	98	74-122
Vinyl Acetate	8260C	30.4	20.0	152	52-174
Vinyl Chloride	8260C	22.9	20.0	115	74-159
cis-1,2-Dichloroethene	8260C	21.2	20.0	106	80-121
cis-1,3-Dichloropropene	8260C	21.0	20.0	105	77-122
trans-1,2-Dichloroethene	8260C	23.2	20.0	116	73-118
Printed 7/24/2020 2:55:41 PM			Superso	et Reference:20-000	00555197 rev 00

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix:

Project:

Water

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

> Units:ug/L Basis:NA

Service Request: R2005635

Date Analyzed: 07/13/20

Lab Control Sample

RQ2007451-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
trans-1,3-Dichloropropene	8260C	20.5	20.0	103	71-133

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring **Date Analyzed:** 07/14/20

Sample Matrix: Water

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005635

Lab Control Sample

RQ2007501-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
1,1,1-Trichloroethane (TCA)	8260C	20.1	20.0	100	75-125
1,1,2,2-Tetrachloroethane	8260C	22.4	20.0	112	78-126
1,1,2-Trichloroethane	8260C	21.9	20.0	110	82-121
1,1-Dichloroethane (1,1-DCA)	8260C	20.0	20.0	100	80-124
1,1-Dichloroethene (1,1-DCE)	8260C	22.5	20.0	112	71-118
1,2-Dichloroethane	8260C	19.9	20.0	99	71-127
1,2-Dichloropropane	8260C	19.8	20.0	99	80-119
2-Butanone (MEK)	8260C	22.7	20.0	114	61-137
2-Hexanone	8260C	19.9	20.0	99	63-124
4-Methyl-2-pentanone	8260C	20.8	20.0	104	66-124
Acetone	8260C	21.8	20.0	109	40-161
Benzene	8260C	20.3	20.0	101	79-119
Bromodichloromethane	8260C	21.0	20.0	105	81-123
Bromoform	8260C	19.8	20.0	99	65-146
Bromomethane	8260C	19.2	20.0	96	42-166
Carbon Disulfide	8260C	21.2	20.0	106	66-128
Carbon Tetrachloride	8260C	21.0	20.0	105	70-127
Chlorobenzene	8260C	20.7	20.0	104	80-121
Chloroethane	8260C	18.2	20.0	91	62-131
Chloroform	8260C	20.1	20.0	101	79-120
Chloromethane	8260C	19.9	20.0	100	65-135
Dibromochloromethane	8260C	22.1	20.0	111	72-128
Dichloromethane	8260C	19.2	20.0	96	73-122
Ethylbenzene	8260C	21.4	20.0	107	76-120
Styrene	8260C	20.7	20.0	103	80-124
Tetrachloroethene (PCE)	8260C	19.9	20.0	100	72-125
Γoluene	8260C	21.2	20.0	106	79-119
Trichloroethene (TCE)	8260C	18.0	20.0	90	74-122
Vinyl Acetate	8260C	35.0	20.0	175 *	52-174
Vinyl Chloride	8260C	18.8	20.0	94	74-159
cis-1,2-Dichloroethene	8260C	19.8	20.0	99	80-121
cis-1,3-Dichloropropene	8260C	20.6	20.0	103	77-122
trans-1,2-Dichloroethene	8260C	22.7	20.0	114	73-118
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QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix: Water

Project:

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

> Units:ug/L Basis:NA

Service Request: R2005635

Date Analyzed: 07/14/20

Lab Control Sample

RQ2007501-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
trans-1.3-Dichloropropene	8260C	19 9	20.0	100	71-133



Semivolatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

SURROGATE RECOVERY SUMMARY Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Extraction Method:** EPA 3510C

		2,4,6-Tribromophenol	2-Fluorobiphenyl	2-Fluorophenol
Sample Name	Lab Code	35-141	31-118	10-105
WG-9954-063020-SG-019	R2005635-001	121	73	45
WG-9954-063020-SG-020	R2005635-002	102	69	44
WG-9954-063020-SG-022	R2005635-003	118	80	44
WG-9954-063020-RM-023	R2005635-004	102	74	40
RB-9954-063020-SG-002	R2005635-005	98	68	39
WG-9954-063020-SG-021	R2005635-007	114	83	50
Method Blank	RQ2007073-03	101	71	48
Lab Control Sample	RQ2007073-04	115	78	47
Duplicate Lab Control Sample	RQ2007073-05	103	67	47
WG-9954-063020-RM-023 MS	RQ2007073-01	121	86	50
WG-9954-063020-RM-023 DMS	RQ2007073-02	105	83	40

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

SURROGATE RECOVERY SUMMARY Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Extraction Method:** EPA 3510C

		Nitrobenzene-d5	Phenol-d6	p-Terphenyl-d14
Sample Name	Lab Code	31-110	10-107	10-165
WG-9954-063020-SG-019	R2005635-001	71	29	51
WG-9954-063020-SG-020	R2005635-002	71	31	61
WG-9954-063020-SG-022	R2005635-003	71	32	64
WG-9954-063020-RM-023	R2005635-004	72	25	67
RB-9954-063020-SG-002	R2005635-005	65	27	62
WG-9954-063020-SG-021	R2005635-007	83	33	54
Method Blank	RQ2007073-03	69	35	66
Lab Control Sample	RQ2007073-04	80	34	58
Duplicate Lab Control Sample	RQ2007073-05	66	34	59
WG-9954-063020-RM-023 MS	RQ2007073-01	78	34	55
WG-9954-063020-RM-023 DMS	RO2007073-02	73	29	45

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Matrix Spike

Sample Matrix: Water

Project:

Service Request: R2005635

Date Collected: 06/30/20

Date Received: 07/01/20

Date Analyzed: 07/6/20

Date Extracted: 07/2/20

Basis:

Duplicate Matrix Spike Summary Semivolatile Organic Compounds by GC/MS

Sample Name: WG-9954-063020-RM-023

Units: ug/L

NA

Lab Code: R2005635-004

Analysis Method: 8270D **Prep Method:** EPA 3510C

Duplicate Matrix Spike

		RQ2	007073-01		R	Q2007073-0)2			
	Sample		Spike			Spike		% Rec		RPD
Analyte Name	Result	Result	Amount	% Rec	Result	Amount	% Rec	Limits	RPD	Limit
1,2,4-Trichlorobenzene	9.1 U	54.7	72.7	75	49.2	72.7	68	10-127	10	30
1,2-Dichlorobenzene	9.1 U	50.1	72.7	69	40.9	72.7	56	17-105	21	30
1,3-Dichlorobenzene	9.1 U	51.3	72.7	71	41.0	72.7	56	21-99	24	30
1,4-Dichlorobenzene	9.1 U	51.3	72.7	71	41.3	72.7	57	10-124	22	30
2,4,5-Trichlorophenol	9.1 U	65.0	72.7	89	59.6	72.7	82	48-134	8	30
2,4,6-Trichlorophenol	9.1 U	63.7	72.7	88	53.5	72.7	74	44-135	17	30
2,4-Dichlorophenol	9.1 U	59.9	72.7	82	50.3	72.7	69	40-130	17	30
2,4-Dimethylphenol	9.1 U	61.9	72.7	85	61.6	72.7	85	42-121	<1	30
2,4-Dinitrophenol	45 U	32.7 J	72.7	45	28.3 J	72.7	39	21-168	14	30
2,4-Dinitrotoluene	9.1 U	70.5	72.7	97	64.5	72.7	89	37-143	9	30
2,6-Dinitrotoluene	9.1 U	75.1	72.7	103	65.4	72.7	90	39-136	13	30
2-Chloronaphthalene	9.1 U	63.8	72.7	88	59.8	72.7	82	40-108	7	30
2-Chlorophenol	9.1 U	50.3	72.7	69	42.2	72.7	58	37-112	17	30
2-Methylnaphthalene	9.1 U	59.0	72.7	81	55.9	72.7	77	34-102	5	30
2-Methylphenol	9.1 U	52.5	72.7	72	47.1	72.7	65	37-102	10	30
2-Nitroaniline	9.1 U	74.3	72.7	102	73.2	72.7	101	40-136	<1	30
2-Nitrophenol	9.1 U	56.6	72.7	78	49.9	72.7	69	27-143	12	30
3,3'-Dichlorobenzidine	9.1 U	71.2	72.7	98	70.0	72.7	96	11-131	2	30
3- and 4-Methylphenol Coelution	9.1 U	49.0	72.7	67	46.0	72.7	63	30-95	6	30
3-Nitroaniline	9.1 U	60.1	72.7	83	62.0	72.7	85	19-117	2	30
4,6-Dinitro-2-methylphenol	45 U	52.1	72.7	72	49.5	72.7	68	25-154	6	30
4-Bromophenyl Phenyl Ether	9.1 U	65.3	72.7	90	64.2	72.7	88	39-115	2	30
4-Chloro-3-methylphenol	9.1 U	64.7	72.7	89	63.9	72.7	88	41-126	1	30
4-Chloroaniline	9.1 U	60.3	72.7	83	64.8	72.7	89	19-111	7	30
4-Chlorophenyl Phenyl Ether	9.1 U	61.7	72.7	85	62.9	72.7	86	41-111	1	30
4-Nitroaniline	9.1 U	65.4	72.7	90	68.2	72.7	94	18-143	4	30
4-Nitrophenol	45 U	30.6 J	72.7	42	26.4 J	72.7	36	10-126	15	30
Acenaphthene	9.1 U	64.5	72.7	89	62.5	72.7	86	43-117	3	30
Acenaphthylene	9.1 U	68.7	72.7	94	67.6	72.7	93	45-119	1	30
Anthracene	9.1 U	67.7	72.7	93	65.8	72.7	90	45-127	3	30
Benz(a)anthracene	9.1 U	50.7	72.7	70	45.3	72.7	62	46-126	12	30
Benzo(a)pyrene	9.1 U	52.1	72.7	72	45.5	72.7	63	44-114	13	30
Benzo(b)fluoranthene	9.1 U	46.3	72.7	64	42.0	72.7	58	41-127	10	30

Results flagged with an asterisk (*) indicate values outside control criteria.

Results flagged with a pound (#) indicate the control criteria is not applicable.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

Matrix Spike and Matrix Spike Duplicate Data is presented for information purposes only. The matrix may or may not be relevant to samples reported in this report. The laboratory evaluates system performance based on the LCS and LCSD control limits.

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates) **Service Request:**

R2005635

Project:

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Date Collected:

06/30/20

Sample Matrix: Water

Date Received:

07/01/20

Date Analyzed: **Date Extracted:**

07/6/20 07/2/20

Duplicate Matrix Spike Summary Semivolatile Organic Compounds by GC/MS

Sample Name: WG-9954-063020-RM-023 **Units:**

ug/L

Lab Code:

R2005635-004

Basis:

NA

Analysis Method: Prep Method:

8270D

EPA 3510C

Matrix Spike

Duplicate Matrix Spike

		RQ2	007073-01		R	.Q2007073-0)2			
	Sample		Spike			Spike		% Rec		RPD
Analyte Name	Result	Result	Amount	% Rec	Result	Amount	% Rec	Limits	RPD	Limit
Benzo(g,h,i)perylene	9.1 U	53.0	72.7	73	47.7	72.7	66	50-143	10	30
Benzo(k)fluoranthene	9.1 U	49.5	72.7	68	42.9	72.7	59	46-139	14	30
Benzoic Acid	91 U	63.2 J	109	58	58.1 J	109	53	10-94	9	30
Benzyl Alcohol	9.1 U	64.8	72.7	89	62.8	72.7	86	31-109	3	30
2,2'-Oxybis(1-chloropropane)	9.1 U	50.4	72.7	69	41.9	72.7	58	21-126	17	30
Bis(2-chloroethoxy)methane	9.1 U	55.2	72.7	76	55.0	72.7	76	41-118	<1	30
Bis(2-chloroethyl) Ether	9.1 U	48.6	72.7	67	42.6	72.7	59	33-108	13	30
Bis(2-ethylhexyl) Phthalate	9.1 U	44.3	72.7	61	38.8	72.7	53	41-132	14	30
Butyl Benzyl Phthalate	9.1 U	55.3	72.7	76	53.2	72.7	73	41-148	4	30
Chrysene	9.1 U	53.2	72.7	73	46.8	72.7	64	47-126	13	30
Di-n-butyl Phthalate	9.1 U	72.2	72.7	99	67.8	72.7	93	43-130	6	30
Di-n-octyl Phthalate	9.1 U	43.5	72.7	60	37.9	72.7	52	40-139	14	30
Dibenz(a,h)anthracene	9.1 U	57.9	72.7	80	50.2	72.7	69	43-136	15	30
Dibenzofuran	9.1 U	69.3	72.7	95	67.1	72.7	92	46-119	3	30
Diethyl Phthalate	9.1 U	64.9	72.7	89	61.4	72.7	84	36-122	6	30
Dimethyl Phthalate	9.1 U	73.9	72.7	102	72.8	72.7	100	33-123	2	30
Fluoranthene	9.1 U	74.6	72.7	103	71.5	72.7	98	43-135	5	30
Fluorene	9.1 U	70.5	72.7	97	67.4	72.7	93	43-113	4	30
Hexachlorobenzene	9.1 U	70.8	72.7	97	65.4	72.7	90	42-125	7	30
Hexachlorobutadiene	9.1 U	63.6	72.7	87	59.6	72.7	82	10-111	6	30
Hexachlorocyclopentadiene	9.1 U	23.5	72.7	32	19.6	72.7	27	10-103	17	30
Hexachloroethane	9.1 U	54.0	72.7	74	41.8	72.7	57	12-101	26	30
Indeno(1,2,3-cd)pyrene	9.1 U	53.4	72.7	73	47.3	72.7	65	49-140	12	30
Isophorone	9.1 U	50.4	72.7	69	50.1	72.7	69	40-111	<1	30
N-Nitrosodi-n-propylamine	9.1 U	63.5	72.7	87	59.7	72.7	82	35-108	6	30
N-Nitrosodiphenylamine	9.1 U	80.4	72.7	111	78.7	72.7	108	43-127	3	30
Naphthalene	9.1 U	59.5	72.7	82	51.3	72.7	71	37-108	14	30
Nitrobenzene	9.1 U	59.1	72.7	81	51.5	72.7	71	35-112	13	30
Pentachlorophenol (PCP)	45 U	60.6	72.7	83	53.0	72.7	73	29-164	13	30
Phenanthrene	9.1 U	66.9	72.7	92	65.0	72.7	89	46-123	3	30
Phenol	9.1 U	28.2	72.7	39	27.2	72.7	37	10-113	5	30
Pyrene	9.1 U	64.3	72.7	88	60.3	72.7	83	44-129	6	30

Results flagged with an asterisk (*) indicate values outside control criteria.

Results flagged with a pound (#) indicate the control criteria is not applicable.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

Matrix Spike and Matrix Spike Duplicate Data is presented for information purposes only. The matrix may or may not be relevant to samples reported in this report. The laboratory evaluates system performance based on the LCS and LCSD control limits.

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007073-03
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	10 U	10	1.2	1	07/06/20 14:53	7/2/20	
1,2-Dichlorobenzene	10 U	10	1.2	1	07/06/20 14:53	7/2/20	
1,3-Dichlorobenzene	10 U	10	1.1	1	07/06/20 14:53	7/2/20	
1,4-Dichlorobenzene	10 U	10	1.2	1	07/06/20 14:53	7/2/20	
2,4,5-Trichlorophenol	10 U	10	1.1	1	07/06/20 14:53	7/2/20	
2,4,6-Trichlorophenol	10 U	10	1.4	1	07/06/20 14:53	7/2/20	
2,4-Dichlorophenol	10 U	10	1.3	1	07/06/20 14:53	7/2/20	
2,4-Dimethylphenol	10 U	10	1.4	1	07/06/20 14:53	7/2/20	
2,4-Dinitrophenol	50 U	50	20	1	07/06/20 14:53	7/2/20	
2,4-Dinitrotoluene	10 U	10	2.4	1	07/06/20 14:53	7/2/20	
2,6-Dinitrotoluene	10 U	10	1.4	1	07/06/20 14:53	7/2/20	
2-Chloronaphthalene	10 U	10	1.4	1	07/06/20 14:53	7/2/20	
2-Chlorophenol	10 U	10	1.1	1	07/06/20 14:53	7/2/20	
2-Methylnaphthalene	10 U	10	1.3	1	07/06/20 14:53	7/2/20	
2-Methylphenol	10 U	10	1.0	1	07/06/20 14:53	7/2/20	
2-Nitroaniline	10 U	10	1.4	1	07/06/20 14:53	7/2/20	
2-Nitrophenol	10 U	10	1.5	1	07/06/20 14:53	7/2/20	
3,3'-Dichlorobenzidine	10 U	10	1.2	1	07/06/20 14:53	7/2/20	
3- and 4-Methylphenol Coelution	10 U	10	1.2	1	07/06/20 14:53	7/2/20	
3-Nitroaniline	10 U	10	2.5	1	07/06/20 14:53	7/2/20	
4,6-Dinitro-2-methylphenol	50 U	50	2.3	1	07/06/20 14:53	7/2/20	
4-Bromophenyl Phenyl Ether	10 U	10	1.7	1	07/06/20 14:53	7/2/20	
4-Chloro-3-methylphenol	10 U	10	1.7	1	07/06/20 14:53	7/2/20	
4-Chloroaniline	10 U	10	1.1	1	07/06/20 14:53	7/2/20	
4-Chlorophenyl Phenyl Ether	10 U	10	1.5	1	07/06/20 14:53	7/2/20	
4-Chlorophenyl Fhenyl Ethel 4-Nitroaniline	10 U	10	2.7	1	07/06/20 14:53	7/2/20	
	50 U	50	2.7 6.4			7/2/20	
4-Nitrophenol				1	07/06/20 14:53		
Acenaphthene	10 U	10	1.4	1	07/06/20 14:53	7/2/20	
Acenaphthylene	10 U	10	1.4	1	07/06/20 14:53	7/2/20	
Anthracene	10 U	10	1.3	1	07/06/20 14:53	7/2/20	
Benz(a)anthracene	10 U	10	1.6	1	07/06/20 14:53	7/2/20	
Benzo(a)pyrene	10 U	10	1.2	1	07/06/20 14:53	7/2/20	
Benzo(b)fluoranthene	10 U	10	1.2	1	07/06/20 14:53	7/2/20	
Benzo(g,h,i)perylene	10 U	10	1.0	1	07/06/20 14:53	7/2/20	
Benzo(k)fluoranthene	10 U	10	1.3	1	07/06/20 14:53	7/2/20	
Benzoic Acid	100 U	100	36	1	07/06/20 14:53	7/2/20	
Benzyl Alcohol	10 U	10	1.6	1	07/06/20 14:53	7/2/20	
2,2'-Oxybis(1-chloropropane)	10 U	10	1.4	1	07/06/20 14:53	7/2/20	
Bis(2-chloroethoxy)methane	10 U	10	1.9	1	07/06/20 14:53	7/2/20	
Bis(2-chloroethyl) Ether	10 U	10	1.3	1	07/06/20 14:53	7/2/20	
Bis(2-ethylhexyl) Phthalate	10 U	10	1.0	1	07/06/20 14:53	7/2/20	
Butyl Benzyl Phthalate	10 U	10	1.4	1	07/06/20 14:53	7/2/20	
Chrysene	10 U	10	1.2	1	07/06/20 14:53	7/2/20	

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Superset Reference: 20-0000555197 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007073-03
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	10 U	10	2.0	1	07/06/20 14:53	7/2/20	
Di-n-octyl Phthalate	10 U	10	3.3	1	07/06/20 14:53	7/2/20	
Dibenz(a,h)anthracene	10 U	10	1.1	1	07/06/20 14:53	7/2/20	
Dibenzofuran	10 U	10	1.4	1	07/06/20 14:53	7/2/20	
Diethyl Phthalate	10 U	10	1.1	1	07/06/20 14:53	7/2/20	
Dimethyl Phthalate	10 U	10	1.3	1	07/06/20 14:53	7/2/20	
Fluoranthene	10 U	10	1.5	1	07/06/20 14:53	7/2/20	
Fluorene	10 U	10	1.3	1	07/06/20 14:53	7/2/20	
Hexachlorobenzene	10 U	10	1.6	1	07/06/20 14:53	7/2/20	
Hexachlorobutadiene	10 U	10	1.0	1	07/06/20 14:53	7/2/20	
Hexachlorocyclopentadiene	10 U	10	2.2	1	07/06/20 14:53	7/2/20	
Hexachloroethane	10 U	10	1.1	1	07/06/20 14:53	7/2/20	
Indeno(1,2,3-cd)pyrene	10 U	10	1.8	1	07/06/20 14:53	7/2/20	
Isophorone	10 U	10	1.4	1	07/06/20 14:53	7/2/20	
N-Nitrosodi-n-propylamine	10 U	10	1.2	1	07/06/20 14:53	7/2/20	
N-Nitrosodiphenylamine	10 U	10	2.7	1	07/06/20 14:53	7/2/20	
Naphthalene	10 U	10	1.2	1	07/06/20 14:53	7/2/20	
Nitrobenzene	10 U	10	1.5	1	07/06/20 14:53	7/2/20	
Pentachlorophenol (PCP)	50 U	50	9.8	1	07/06/20 14:53	7/2/20	
Phenanthrene	10 U	10	1.4	1	07/06/20 14:53	7/2/20	
Phenol	10 U	10	1.0	1	07/06/20 14:53	7/2/20	
Pyrene	10 U	10	1.5	1	07/06/20 14:53	7/2/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	101	35 - 141	07/06/20 14:53	
2-Fluorobiphenyl	71	31 - 118	07/06/20 14:53	
2-Fluorophenol	48	10 - 105	07/06/20 14:53	
Nitrobenzene-d5	69	31 - 110	07/06/20 14:53	
Phenol-d6	35	10 - 107	07/06/20 14:53	
p-Terphenyl-d14	66	10 - 165	07/06/20 14:53	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown	11.52	7.9	J	
	unknown	11.93	8.1	J	
	unknown hydrocarbon	12.08	12	J	
	unknown hydrocarbon	12.69	17	J	
	unknown hydrocarbon	13.36	17	J	
	unknown hydrocarbon	14.08	14	J	
	unknown hydrocarbon	14.85	13	J	

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Superset Reference:20-0000555197 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA **Project:**

Monitoring

Sample Matrix: Date Received: NA Water

Sample Name: Method Blank Units: ug/L

Lab Code: RQ2007073-03 Basis: NA

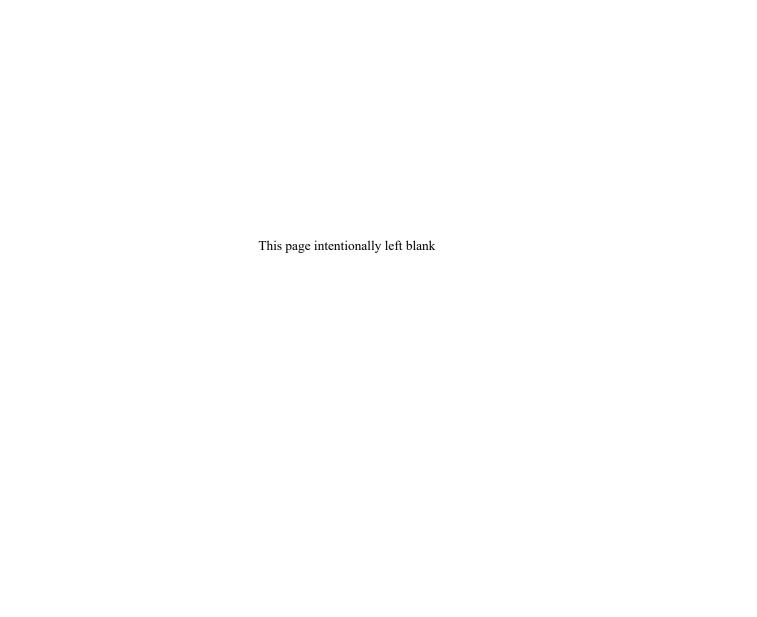
Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Tentatively Identified Compounds

CAS#	Compound Identification	RT	ug/L	Q	
	unknown hydrocarbon	15.65	11	J	
	unknown	16.34	8.5	J	





QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix: Water

Date Analyzed: 07/06/20 - 07/08/20

Service Request: R2005635

Duplicate Lab Control Sample Summary Semivolatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Lab Control Sample

Duplicate Lab Control Sample

RQ2007073-04

RQ2007073-05

Analyte Name	Analytica l Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
1,2,4-Trichlorobenzene	8270D	56.0	80.0	70	50.2	80.0	63	10-127	11	30
1,2-Dichlorobenzene	8270D	56.6	80.0	71	47.1	80.0	59	23-130	18	30
1,3-Dichlorobenzene	8270D	55.5	80.0	69	48.6	80.0	61	21-90	12	30
1,4-Dichlorobenzene	8270D	53.4	80.0	67	47.6	80.0	60	10-124	11	30
2,4,5-Trichlorophenol	8270D	69.5	80.0	87	57.7	80.0	72	48-134	19	30
2,4,6-Trichlorophenol	8270D	65.0	80.0	81	55.5	80.0	69	44-135	16	30
2,4-Dichlorophenol	8270D	58.0	80.0	73	50.9	80.0	64	48-127	13	30
2,4-Dimethylphenol	8270D	58.8	80.0	74	56.3	80.0	70	59-113	6	30
2,4-Dinitrophenol	8270D	48.1 J	80.0	60	43.6 J	80.0	55	21-154	9	30
2,4-Dinitrotoluene	8270D	64.1	80.0	80	55.3	80.0	69	54-130	15	30
2,6-Dinitrotoluene	8270D	73.4	80.0	92	65.1	80.0	81	51-127	13	30
2-Chloronaphthalene	8270D	65.8	80.0	82	56.4	80.0	71	40-108	14	30
2-Chlorophenol	8270D	52.9	80.0	66	49.9	80.0	62	42-112	6	30
2-Methylnaphthalene	8270D	58.8	80.0	74	52.4	80.0	65	34-102	13	30
2-Methylphenol	8270D	55.4	80.0	69	50.8	80.0	64	47-100	8	30
2-Nitroaniline	8270D	67.4	80.0	84	65.2	80.0	82	52-133	2	30
2-Nitrophenol	8270D	58.6	80.0	73	51.8	80.0	65	43-131	12	30
3,3'-Dichlorobenzidine	8270D	59.9	80.0	75	56.1	80.0	70	43-126	7	30
3- and 4-Methylphenol Coelution	8270D	48.9	80.0	61	47.2	80.0	59	40-92	3	30
3-Nitroaniline	8270D	58.4	80.0	73	62.2	80.0	78	42-111	7	30
4,6-Dinitro-2-methylphenol	8270D	54.0	80.0	68	45.9 J	80.0	57	36-152	18	30
4-Bromophenyl Phenyl Ether	8270D	66.2	80.0	83	60.4	80.0	75	48-114	10	30
4-Chloro-3-methylphenol	8270D	62.1	80.0	78	55.8	80.0	70	52-113	11	30
4-Chloroaniline	8270D	57.6	80.0	72	61.9	80.0	77	44-109	7	30
4-Chlorophenyl Phenyl Ether	8270D	58.4	80.0	73	49.4	80.0	62	51-107	16	30
4-Nitroaniline	8270D	58.1	80.0	73	50.1	80.0	63	54-133	15	30
4-Nitrophenol	8270D	20.2 J	80.0	25	18.7 J	80.0	23	10-126	8	30
Acenaphthene	8270D	64.3	80.0	80	58.2	80.0	73	52-107	9	30
Acenaphthylene	8270D	70.3	80.0	88	62.1	80.0	78	55-109	12	30
Anthracene	8270D	67.2	80.0	84	59.9	80.0	75	55-116	11	30
Benz(a)anthracene	8270D	60.6	80.0	76	58.1	80.0	73	61-121	4	30
Benzo(a)pyrene	8270D	62.2	80.0	78	61.3	80.0	77	44-114	1	30
Benzo(b)fluoranthene	8270D	59.2	80.0	74	58.3	80.0	73	62-115	1	30
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QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix: Water

Service Request: R2005635 **Date Analyzed:** 07/06/20 - 07/08/20

Duplicate Lab Control Sample Summary Semivolatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Lab Control Sample

Duplicate Lab Control Sample

RQ2007073-04

RQ2007073-05

Analyte Name	Analytica l Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
Benzo(g,h,i)perylene	8270D	70.4	80.0	88	68.4	80.0	85	63-136	3	30
Benzo(k)fluoranthene	8270D	68.5	80.0	86	64.2	80.0	80	49-133	7	30
Benzoic Acid	8270D	70.9 J	120	59	67.3 J	120	56	10-94	5	30
Benzyl Alcohol	8270D	60.0	80.0	75	58.9	80.0	74	31-109	1	30
2,2'-Oxybis(1-chloropropane)	8270D	63.4	80.0	79	54.4	80.0	68	32-122	15	30
Bis(2-chloroethoxy)methane	8270D	66.2	80.0	83	57.4	80.0	72	55-110	14	30
Bis(2-chloroethyl) Ether	8270D	61.0	80.0	76	51.7	80.0	65	46-102	16	30
Bis(2-ethylhexyl) Phthalate	8270D	62.1	80.0	78	60.8	80.0	76	51-132	3	30
Butyl Benzyl Phthalate	8270D	62.3	80.0	78	60.3	80.0	75	41-148	4	30
Chrysene	8270D	65.3	80.0	82	63.8	80.0	80	57-118	2	30
Di-n-butyl Phthalate	8270D	72.4	80.0	91	66.3	80.0	83	57-128	9	30
Di-n-octyl Phthalate	8270D	63.7	80.0	80	61.0	80.0	76	62-124	5	30
Dibenz(a,h)anthracene	8270D	75.6	80.0	95	71.4	80.0	89	54-135	7	30
Dibenzofuran	8270D	69.7	80.0	87	57.6	80.0	72	55-110	19	30
Diethyl Phthalate	8270D	62.9	80.0	79	53.9	80.0	67	53-113	16	30
Dimethyl Phthalate	8270D	72.4	80.0	90	64.0	80.0	80	51-112	12	30
Fluoranthene	8270D	72.0	80.0	90	66.4	80.0	83	66-127	8	30
Fluorene	8270D	69.1	80.0	86	56.8	80.0	71	54-106	19	30
Hexachlorobenzene	8270D	73.1	80.0	91	69.0	80.0	86	53-123	6	30
Hexachlorobutadiene	8270D	58.0	80.0	73	53.1	80.0	66	16-95	10	30
Hexachlorocyclopentadiene	8270D	19.1	80.0	24	18.0	80.0	23	10-99	4	30
Hexachloroethane	8270D	55.7	80.0	70	49.0	80.0	61	15-92	14	30
Indeno(1,2,3-cd)pyrene	8270D	62.2	80.0	78	61.9	80.0	77	62-137	1	30
Isophorone	8270D	54.6	80.0	68	48.0	80.0	60	50-116	13	30
N-Nitrosodi-n-propylamine	8270D	66.7	80.0	83	60.7	80.0	76	49-115	9	30
N-Nitrosodiphenylamine	8270D	85.1	80.0	106	75.5	80.0	94	45-123	12	30
Naphthalene	8270D	62.2	80.0	78	54.1	80.0	68	38-99	14	30
Nitrobenzene	8270D	60.1	80.0	75	54.5	80.0	68	46-108	10	30
Pentachlorophenol (PCP)	8270D	67.6	80.0	85	59.1	80.0	74	29-164	14	30
Phenanthrene	8270D	67.1	80.0	84	61.2	80.0	77	58-118	9	30
Phenol	8270D	31.5	80.0	39	32.0	80.0	40	10-113	3	30
Pyrene	8270D	69.2	80.0	86	65.3	80.0	82	61-122	5	30

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Semivolatile Organic Compounds by GC

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

SURROGATE RECOVERY SUMMARY

Organochlorine Pesticides by Gas Chromatography

Analysis Method: 8081B **Extraction Method:** EPA 3510C

		Decachlorobiphenyl	Tetrachloro-m-xylene	
Sample Name	Lab Code	10-164	10-147	
WG-9954-063020-SG-019	R2005635-001	72	75	
WG-9954-063020-SG-020	R2005635-002	74	59	
WG-9954-063020-SG-022	R2005635-003	52	66	
WG-9954-063020-RM-023	R2005635-004	18	58	
RB-9954-063020-SG-002	R2005635-005	9*	55	
RB-9954-063020-SG-002 RE	R2005635-005	9*	36	
WG-9954-063020-SG-021	R2005635-007	85	79	
Method Blank	RQ2007072-05	63	55	
Method Blank	RQ2007347-01	63	65	
Lab Control Sample	RQ2007072-06	55	49	
Duplicate Lab Control Sample	RQ2007072-07	76	60	
Lab Control Sample	RQ2007347-02	56	58	
Duplicate Lab Control Sample	RQ2007347-03	62	55	
WG-9954-063020-RM-023 MS	RQ2007072-01	19	63	
WG-9954-063020-RM-023 DMS	RQ2007072-02	32	66	

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Date Collected: 06/30/20 **Date Received:** 07/01/20

Water

07/01/20 07/7/20

Date Analyzed: Date Extracted:

07/2/20

Duplicate Matrix Spike Summary Organochlorine Pesticides by Gas Chromatography

Sample Name: WG-9954-063020-RM-023

Units: ug/L

Lab Code: R2005635-004

Basis: NA

Analysis Method: 8081B **Prep Method:** EPA 3510C

Sample Matrix:

			Matrix S RQ20070	-		Duplicate Ma RQ20070	_			
	Sample		Spike			Spike		% Rec		RPD
Analyte Name	Result	Result	Amount	% Rec	Result	Amount	% Rec	Limits	RPD	Limit
4,4'-DDD	0.045 U	0.261	0.364	72	0.308	0.364	85	38-157	17	30
4,4'-DDE	0.045 U	0.261	0.364	72	0.303	0.364	83	10-200	15	30
4,4'-DDT	0.045 U	0.246	0.364	68	0.293	0.364	81	19-154	18	30
Aldrin	0.045 U	0.212	0.364	58	0.235	0.364	65	26-149	10	30
Dieldrin	0.045 U	0.297	0.364	82	0.354	0.364	97	41-164	18	30
Endosulfan I	0.045 U	0.294	0.364	81	0.352	0.364	97	47-149	18	30
Endosulfan II	0.045 U	0.306	0.364	84	0.360	0.364	99	51-148	16	30
Endosulfan Sulfate	0.045 U	0.259	0.364	71	0.301	0.364	83	10-170	15	30
Endrin	0.045 U	0.291	0.364	80	0.347	0.364	96	48-165	18	30
Endrin Ketone	0.045 U	0.301	0.364	83	0.361	0.364	99	48-162	18	30
Heptachlor	0.045 U	0.157	0.364	43	0.181	0.364	50	29-168	15	30
Heptachlor Epoxide	0.045 U	0.296	0.364	81	0.351	0.364	97	29-180	17	30
Methoxychlor	0.045 U	0.286	0.364	79	0.338	0.364	93	38-162	17	30
alpha-BHC	0.045 U	0.283	0.364	78	0.340	0.364	94	27-154	18	30
alpha-Chlordane	0.045 U	0.277	0.364	76	0.326	0.364	90	35-160	16	30
beta-BHC	0.045 U	0.309	0.364	85	0.362	0.364	100	32-184	16	30
delta-BHC	0.045 U	0.278	0.364	76	0.331	0.364	91	10-182	17	30
gamma-BHC (Lindane)	0.045 U	0.289	0.364	79	0.345	0.364	95	43-164	18	30
gamma-Chlordane	0.045 U	0.292	0.364	80	0.341	0.364	94	35-165	15	30

Results flagged with an asterisk (*) indicate values outside control criteria.

Results flagged with a pound (#) indicate the control criteria is not applicable.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

Matrix Spike and Matrix Spike Duplicate Data is presented for information purposes only. The matrix may or may not be relevant to samples reported in this report. The laboratory evaluates system performance based on the LCS and LCSD control limits.

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007072-05Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analysis Method: 8081B **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.050 U	0.050	0.020	1	07/06/20 14:30	7/2/20	
4,4'-DDE	0.050 U	0.050	0.020	1	07/06/20 14:30	7/2/20	
4,4'-DDT	0.050 U	0.050	0.020	1	07/06/20 14:30	7/2/20	
Aldrin	0.050 U	0.050	0.020	1	07/06/20 14:30	7/2/20	
Dieldrin	0.050 U	0.050	0.020	1	07/06/20 14:30	7/2/20	
Endosulfan I	0.050 U	0.050	0.020	1	07/06/20 14:30	7/2/20	
Endosulfan II	0.050 U	0.050	0.020	1	07/06/20 14:30	7/2/20	
Endosulfan Sulfate	0.050 U	0.050	0.020	1	07/06/20 14:30	7/2/20	
Endrin	0.050 U	0.050	0.020	1	07/06/20 14:30	7/2/20	
Endrin Ketone	0.050 U	0.050	0.020	1	07/06/20 14:30	7/2/20	
Heptachlor	0.050 U	0.050	0.020	1	07/06/20 14:30	7/2/20	
Heptachlor Epoxide	0.050 U	0.050	0.020	1	07/06/20 14:30	7/2/20	
Methoxychlor	0.050 U	0.050	0.020	1	07/06/20 14:30	7/2/20	
Toxaphene	0.50 U	0.50	0.50	1	07/06/20 14:30	7/2/20	
alpha-BHC	0.050 U	0.050	0.020	1	07/06/20 14:30	7/2/20	
alpha-Chlordane	0.050 U	0.050	0.020	1	07/06/20 14:30	7/2/20	
beta-BHC	0.050 U	0.050	0.020	1	07/06/20 14:30	7/2/20	
delta-BHC	0.050 U	0.050	0.020	1	07/06/20 14:30	7/2/20	
gamma-BHC (Lindane)	0.050 U	0.050	0.020	1	07/06/20 14:30	7/2/20	
gamma-Chlordane	0.050 U	0.050	0.020	1	07/06/20 14:30	7/2/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	63	10 - 164	07/06/20 14:30	
Tetrachloro-m-xylene	55	10 - 147	07/06/20 14:30	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007347-01Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analysis Method: 8081B **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	_
4,4'-DDE	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
4,4'-DDT	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
Aldrin	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
Dieldrin	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
Endosulfan I	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
Endosulfan II	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
Endosulfan Sulfate	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
Endrin	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
Endrin Ketone	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
Heptachlor	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
Heptachlor Epoxide	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
Methoxychlor	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
Toxaphene	0.50 U	0.50	0.50	1	07/13/20 14:42	7/10/20	
alpha-BHC	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
alpha-Chlordane	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
beta-BHC	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
delta-BHC	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
gamma-BHC (Lindane)	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
gamma-Chlordane	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	63	10 - 164	07/13/20 14:42	
Tetrachloro-m-xylene	65	10 - 147	07/13/20 14:42	

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/06/20

Sample Matrix: Water

Duplicate Lab Control Sample Summary Organochlorine Pesticides by Gas Chromatography

Units:ug/L Basis:NA

Service Request: R2005635

Lab Control Sample

Duplicate Lab Control Sample

RQ2007072-06

RQ2007072-07

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
4,4'-DDD	8081B	0.249	0.400	62	0.314	0.400	7 9 79	42-159	23	30
4,4'-DDE	8081B	0.235	0.400	59	0.316	0.400	79	47-147	29	30
4,4'-DDT	8081B	0.234	0.400	58	0.309	0.400	77	41-149	28	30
Aldrin	8081B	0.190	0.400	48	0.222	0.400	55	22-137	15	30
Dieldrin	8081B	0.259	0.400	65	0.348	0.400	87	52-144	30	30
Endosulfan I	8081B	0.254	0.400	63	0.336	0.400	84	52-136	28	30
Endosulfan II	8081B	0.261	0.400	65	0.360	0.400	90	57-138	32*	30
Endosulfan Sulfate	8081B	0.222	0.400	56	0.302	0.400	75	34-156	30	30
Endrin	8081B	0.255	0.400	64	0.343	0.400	86	56-143	29	30
Endrin Ketone	8081B	0.262	0.400	66	0.345	0.400	86	59-143	27	30
Heptachlor	8081B	0.158	0.400	39	0.195	0.400	49	32-141	21	30
Heptachlor Epoxide	8081B	0.251	0.400	63	0.331	0.400	83	51-143	28	30
Methoxychlor	8081B	0.232	0.400	58	0.302	0.400	76	56-149	26	30
alpha-BHC	8081B	0.232	0.400	58	0.289	0.400	72	36-151	22	30
alpha-Chlordane	8081B	0.246	0.400	62	0.323	0.400	81	50-139	27	30
beta-BHC	8081B	0.255	0.400	64	0.326	0.400	82	55-149	24	30
delta-BHC	8081B	0.223	0.400	56	0.298	0.400	74	29-159	29	30
gamma-BHC (Lindane)	8081B	0.236	0.400	59	0.298	0.400	74	41-149	23	30
gamma-Chlordane	8081B	0.260	0.400	65	0.320	0.400	80	50-140	21	30

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/13/20

Sample Matrix: Water

Duplicate Lab Control Sample Summary Organochlorine Pesticides by Gas Chromatography

Units:ug/L Basis:NA

Service Request: R2005635

Lab Control Sample

Duplicate Lab Control Sample

RQ2007347-02

RQ2007347-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
4,4'-DDD	8081B	0.303	0.400	76	0.342	0.400	86	42-159	12	30
4,4'-DDE	8081B	0.280	0.400	70	0.322	0.400	81	47-147	14	30
4,4'-DDT	8081B	0.309	0.400	77	0.353	0.400	88	41-149	13	30
Aldrin	8081B	0.236	0.400	59	0.234	0.400	59	22-137	<1	30
Dieldrin	8081B	0.307	0.400	77	0.346	0.400	86	52-144	12	30
Endosulfan I	8081B	0.296	0.400	74	0.329	0.400	82	52-136	11	30
Endosulfan II	8081B	0.236	0.400	59	0.283	0.400	71	57-138	18	30
Endosulfan Sulfate	8081B	0.267	0.400	67	0.310	0.400	78	34-156	15	30
Endrin	8081B	0.319	0.400	80	0.357	0.400	89	56-143	11	30
Endrin Ketone	8081B	0.301	0.400	75	0.346	0.400	87	59-143	14	30
Heptachlor	8081B	0.274	0.400	68	0.283	0.400	71	32-141	3	30
Heptachlor Epoxide	8081B	0.307	0.400	77	0.338	0.400	84	51-143	10	30
Methoxychlor	8081B	0.318	0.400	80	0.359	0.400	90	56-149	12	30
alpha-BHC	8081B	0.297	0.400	74	0.313	0.400	78	36-151	5	30
alpha-Chlordane	8081B	0.293	0.400	73	0.322	0.400	80	50-139	9	30
beta-BHC	8081B	0.315	0.400	79	0.341	0.400	85	55-149	8	30
delta-BHC	8081B	0.292	0.400	73	0.328	0.400	82	29-159	12	30
gamma-BHC (Lindane)	8081B	0.305	0.400	76	0.326	0.400	82	41-149	7	30
gamma-Chlordane	8081B	0.296	0.400	74	0.323	0.400	81	50-140	9	30

QA/QC Report

Service Request: R2005635

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

SURROGATE RECOVERY SUMMARY Polychlorinated Biphenyls (PCBs) by GC

Analysis Method: 8082A **Extraction Method:** EPA 3510C

		Decachlorobiphenyl	Tetrachloro-m-xylene	
Sample Name	Lab Code	10-152	14-129	
WG-9954-063020-SG-019	R2005635-001	103	52	
WG-9954-063020-SG-020	R2005635-002	78	41	
WG-9954-063020-SG-022	R2005635-003	42	45	
WG-9954-063020-RM-023	R2005635-004	15	40	
RB-9954-063020-SG-002	R2005635-005	9*	40	
RB-9954-063020-SG-002 RE	R2005635-005	9*	31	
WG-9954-063020-SG-021	R2005635-007	95	54	
Method Blank	RQ2007072-05	57	44	
Method Blank	RQ2007347-01	69	59	
Lab Control Sample	RQ2007072-06	63	48	
Duplicate Lab Control Sample	RQ2007072-07	66	48	
Lab Control Sample	RQ2007347-02	70	48	
Duplicate Lab Control Sample	RQ2007347-03	68	53	
WG-9954-063020-RM-023 MS	RQ2007072-03	29	38	
WG-9954-063020-RM-023 DMS	RQ2007072-04	19	35	

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates) **Service Request:**

R2005635

Project:

Date Collected:

06/30/20

Sample Matrix: Water

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Date Received:

07/01/20

Date Analyzed: **Date Extracted:**

07/8/20 07/2/20

Duplicate Matrix Spike Summary Polychlorinated Biphenyls (PCBs) by GC

Sample Name:

WG-9954-063020-RM-023

Units:

ug/L

Lab Code:

R2005635-004

Basis:

NA

Analysis Method: Prep Method:

8082A

EPA 3510C

Duplicate Matrix Spike

Matrix Spike RQ2007072-03

RQ2007072-04

	Sample		Spike			Spike		% Rec		RPD
Analyte Name	Result	Result	Amount	% Rec	Result	Amount	% Rec	Limits	RPD	Limit
Aroclor 1016	0.91 U	2.30	3.64	63	2.36	3.64	65	32-142	2	30
Aroclor 1260	0.91 U	2.38	3.64	65	2.43	3.64	67	28-142	2	30

Results flagged with an asterisk (*) indicate values outside control criteria.

Results flagged with a pound (#) indicate the control criteria is not applicable.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

Matrix Spike and Matrix Spike Duplicate Data is presented for information purposes only. The matrix may or may not be relevant to samples reported in this report. The laboratory evaluates system performance based on the LCS and LCSD control limits.

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007072-05Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analysis Method: 8082A **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	1.0 U	1.0	0.50	1	07/06/20 19:25	7/2/20	
Aroclor 1221	2.0 U	2.0	1.0	1	07/06/20 19:25	7/2/20	
Aroclor 1232	1.0 U	1.0	0.50	1	07/06/20 19:25	7/2/20	
Aroclor 1242	1.0 U	1.0	0.50	1	07/06/20 19:25	7/2/20	
Aroclor 1248	1.0 U	1.0	0.50	1	07/06/20 19:25	7/2/20	
Aroclor 1254	1.0 U	1.0	0.50	1	07/06/20 19:25	7/2/20	
Aroclor 1260	1.0 U	1.0	0.50	1	07/06/20 19:25	7/2/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	57	10 - 152	07/06/20 19:25	
Tetrachloro-m-xylene	44	14 - 129	07/06/20 19:25	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005635

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007347-01Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analysis Method: 8082A **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	1.0 U	1.0	0.50	1	07/13/20 14:13	7/10/20	
Aroclor 1221	2.0 U	2.0	1.0	1	07/13/20 14:13	7/10/20	
Aroclor 1232	1.0 U	1.0	0.50	1	07/13/20 14:13	7/10/20	
Aroclor 1242	1.0 U	1.0	0.50	1	07/13/20 14:13	7/10/20	
Aroclor 1248	1.0 U	1.0	0.50	1	07/13/20 14:13	7/10/20	
Aroclor 1254	1.0 U	1.0	0.50	1	07/13/20 14:13	7/10/20	
Aroclor 1260	1.0 U	1.0	0.50	1	07/13/20 14:13	7/10/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	69	10 - 152	07/13/20 14:13	
Tetrachloro-m-xylene	59	14 - 129	07/13/20 14:13	

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix:

Project:

Water

Duplicate Lab Control Sample Summary Polychlorinated Biphenyls (PCBs) by GC

Units:ug/L Basis:NA

Service Request: R2005635

Date Analyzed: 07/06/20

Lab Control Sample

Duplicate Lab Control Sample

RQ2007072-06

RQ2007072-07

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
Aroclor 1016	8082A	2.83	4.00	71	2.66	4.00	67	49-123	6	30
Aroclor 1260	8082A	3.17	4.00	79	3.22	4.00	80	30-120	2	30

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix:

Project:

Water

Duplicate Lab Control Sample Summary Polychlorinated Biphenyls (PCBs) by GC

Units:ug/L Basis:NA

Service Request: R2005635

Date Analyzed: 07/13/20

Lab Control Sample

Duplicate Lab Control Sample

RQ2007347-02

RQ2007347-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
Aroclor 1016	8082A	3.35	4.00	84	3.40	4.00	85	49-123	2	30
Aroclor 1260	8082A	3.98	4.00	100	3.79	4.00	95	30-120	5	30



Service Request No:R2005701

Ms. Kathy Willy GHD Services Inc. 2055 Niagara Falls Blvd., Niagara Falls, NY 14304

Laboratory Results for: Love Canal:292-402-D02-3100

Dear Ms. Willy,

Enclosed are the results of the sample(s) submitted to our laboratory July 02, 2020 For your reference, these analyses have been assigned our service request number **R2005701**.

All testing was performed according to our laboratory's quality assurance program and met the requirements of the TNI standards except as noted in the case narrative report. Any testing not included in the lab's accreditation is identified on a Non-Certified Analytes report. All results are intended to be considered in their entirety. ALS Environmental is not responsible for use of less than the complete report. Results apply only to the individual samples submitted to the lab for analysis, as listed in the report. The measurement uncertainty of the results included in this report is within that expected when using the prescribed method(s), and represented by Laboratory Control Sample control limits. Any events, such as QC failures or Holding Time exceedances, which may add to the uncertainty are explained in the report narrative or are flagged with qualifiers. The flags are explained in the Report Qualifiers and Definitions page of this report.

Please contact me if you have any questions. My extension is 7471. You may also contact me via email at Brady.Kalkman@alsglobal.com.

Respectfully submitted,

Goody Kullen

ALS Group USA, Corp. dba ALS Environmental

Brady Kalkman Project Manager

ADDRESS



Narrative Documents

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com



Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100 Date Received: 07/02/2020

Sample Matrix: Water

CASE NARRATIVE

All analyses were performed consistent with the quality assurance program of ALS Environmental. This report contains analytical results for samples for the Tier level IV requested by the client.

Sample Receipt:

Six water samples were received for analysis at ALS Environmental on 07/02/2020. Any discrepancies upon initial sample inspection are annotated on the sample receipt and preservation form included within this report. The samples were stored at minimum in accordance with the analytical method requirements.

Semivolatiles by GC/MS:

Method 8270D, 07/09/2020: The upper control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). The field samples analyzed in this sequence did not contain the analyte(s) in question above the Method Reporting Limit (MRL). Since the exceedance equates to a potential high bias, the data quality was not significantly affected and no further corrective action was taken.

Method 8270D, 07/09/2020: The lower control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). Since there were no detections of the analyte(s) above the MRL in the associated field samples, the quantitation is not affected. The data quality was not significantly affected and no further corrective action was taken.

Method 8270D, 07/09/2020: The lower control limit for the spike recovery of the Matrix Spike/Matrix Spike Duplicate (MS/MSD) was exceeded for one or more analyte due to sample matrix. The LCS/LCSD was within limits for all analytes. The analytes affected are flagged in the MS Summary.

Method 8270D, 07/13/2020: The upper control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). The field samples analyzed in this sequence did not contain the analyte(s) in question above the Method Reporting Limit (MRL). Since the exceedance equates to a potential high bias, the data quality was not significantly affected and no further corrective action was taken.

Method 8270D, R2005701-001, -006: The control limits for one or more surrogates in the sample are not applicable. The analysis of the sample required a dilution, which resulted in a surrogate concentration below the Method Reporting Limit (MRL). No further corrective action was appropriate.

Semivoa GC:

Method 8081B, R2005701-001: The control limits were exceeded for one or more surrogates. A reanalysis was not performed because insufficient sample was available. No further corrective action was possible.

Method 8081B, r2005701-006: The control limits were exceeded for one or more surrogates due to matrix interferences. A reextraction and reanalysis was performed, but produced similar results. The re-extraction was performed out of holding time. No further corrective action was required.

Method 8082A, 07/08/2020: The upper control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). The field samples analyzed in this sequence did not contain the analyte(s) in question above the Method Reporting Limit (MRL). Since the exceedance equates to a potential high bias, the data quality was not significantly affected and no further corrective action was taken.

Method 8082A, 07/10/2020: The upper control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). The field samples analyzed in this sequence did not contain the analyte(s) in question above the Method Reporting Limit (MRL). Since the exceedance equates to a potential high bias, the data quality was not significantly affected and no further corrective action was taken.

Volatiles by GC/MS:

Approved by	 Date	07/30/2020

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Method 8260C, 07/13/2020: The upper control criterion was exceeded for one or more analytes in the Laboratory Control Sample (LCS). There were no detections of the analyte(s) above the MRL in the associated field samples. The error associated with elevated recovery equates to a high bias. The sample data is not significantly affected. No further corrective action was appropriate.

Method 8260C, 07/15/2020: The lower control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). Since there were no detections of the analyte(s) above the MRL in the associated field samples, the quantitation is not affected. The data quality was not significantly affected and no further corrective action was taken. Method 8260C, 07/14/2020: The upper control criterion was exceeded for one or more analytes in the Laboratory Control Sample (LCS). There were no detections of the analyte(s) above the MRL in the associated field samples. The error associated with elevated recovery equates to a high bias. The sample data is not significantly affected. No further corrective action was appropriate.

Approved by ______

Date 07/30/2020



CLIENT ID: WG-9954-070120-SG-027		Lab	ID: R2005	701-001		
Analyte	Results	Flag	MDL	MRL	Units	Method
1,1,2,2-Tetrachloroethane	6.1		0.20	5.0	ug/L	8260C
1,1,2-Trichloroethane	12		0.20	5.0	ug/L	8260C
1,1-Dichloroethene (1,1-DCE)	0.63	J	0.20	5.0	ug/L	8260C
2-Butanone (MEK)	12		0.78	10	ug/L	8260C
Acetone	79		5.0	10	ug/L	8260C
Benzene	1300	E	0.20	5.0	ug/L	8260C
Chlorobenzene	770	E	0.20	5.0	ug/L	8260C
Chloroform	140		0.24	5.0	ug/L	8260C
Dichloromethane	6.4		0.65	5.0	ug/L	8260C
Ethylbenzene	12		0.20	5.0	ug/L	8260C
Tetrachloroethene (PCE)	20		0.21	5.0	ug/L	8260C
Toluene	1700	Е	0.20	5.0	ug/L	8260C
Trichloroethene (TCE)	120		0.20	5.0	ug/L	8260C
Vinyl Chloride	10		0.20	5.0	ug/L	8260C
Xylenes, Total	58		0.23	5.0	ug/L	8260C
cis-1,2-Dichloroethene	26		0.23	5.0	ug/L	8260C
trans-1,2-Dichloroethene	29		0.20	5.0	ug/L	8260C
Benzene	6500	D	40	1000	ug/L	8260C
Bromodichloromethane	60	DJ	40	1000	ug/L	8260C
Chlorobenzene	2500	D	40	1000	ug/L	8260C
Chloroform	400	DJ	48	1000	ug/L	8260C
Toluene	22000	D	40	1000	ug/L	8260C
Trichloroethene (TCE)	100	DJ	40	1000	ug/L	8260C
1,2,4-Trichlorobenzene	93		1.1	9.1	ug/L	8270D
1,2-Dichlorobenzene	25		1.1	9.1	ug/L	8270D
1,3-Dichlorobenzene	3.6	J	0.92	9.1	ug/L	8270D
1,4-Dichlorobenzene	73		1.1	9.1	ug/L	8270D
2,4,5-Trichlorophenol	37		0.99	9.1	ug/L	8270D
2,4-Dichlorophenol	550	E	1.2	9.1	ug/L	8270D
2,4-Dimethylphenol	8.9	J	1.3	9.1	ug/L	8270D
2-Chlorophenol	27		0.97	9.1	ug/L	8270D
2-Methylnaphthalene	1.4	J	1.2	9.1	ug/L	8270D
2-Methylphenol	32		0.91	9.1	ug/L	8270D
3- and 4-Methylphenol Coelution	71		1.1	9.1	ug/L	8270D
4-Chloro-3-methylphenol	46		0.98	9.1	ug/L	8270D
Benzoic Acid	4500	Е	33	91	ug/L	8270D
Benzyl Alcohol	290		1.5	9.1	ug/L	8270D
Bis(2-chloroethyl) Ether	20		1.2	9.1	ug/L	8270D
Phenol			0.04	0.4	ua/l	8270D
	40		0.91	9.1	ug/L	02/00
1,2,4-Trichlorobenzene	40 59	J	26	230	ug/L ug/L	8270D 8270D

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CLIENT ID: WG-9954-070120-SG-027		Lab	ID: R2005	701-001		
Analyte	Results	Flag	MDL	MRL	Units	Method
2,4,5-Trichlorophenol	50	J	25	230	ug/L	8270D
2,4-Dichlorophenol	330	D	29	230	ug/L	8270D
2-Methylphenol	32	J	23	230	ug/L	8270D
3- and 4-Methylphenol Coelution	60	J	27	230	ug/L	8270D
4-Chloro-3-methylphenol	31	J	25	230	ug/L	8270D
Benzoic Acid	9200	Е	820	2300	ug/L	8270D
Benzyl Alcohol	240		36	230	ug/L	8270D
Phenol	32	J	23	230	ug/L	8270D
1,2,4-Trichlorobenzene	67	J	51	450	ug/L	8270D
1,4-Dichlorobenzene	73	J	52	450	ug/L	8270D
2,4-Dichlorophenol	340	J	57	450	ug/L	8270D
3- and 4-Methylphenol Coelution	61	J	53	450	ug/L	8270D
Benzoic Acid	11000	D	1700	4500	ug/L	8270D
Benzyl Alcohol	220	J	71	450	ug/L	8270D
Aldrin	1.0	Р	0.19	0.45	ug/L	8081B
alpha-BHC	27		0.19	0.45	ug/L	8081B
beta-BHC	6.8	Р	0.19	0.45	ug/L	8081B
delta-BHC	9.2		0.19	0.45	ug/L	8081B
gamma-BHC (Lindane)	6.5		0.19	0.45	ug/L	8081B
CLIENT ID: WG-9954-070120-SG-026		Lab	ID: R2005	701-002		
Analyte	Results	Flag	MDL	MRL	Units	Method
Carbon Disulfide	3.3	J	0.42	10	ug/L	8260C
Chlorobenzene	0.89	J	0.20	5.0	ug/L	8260C
Toluene	0.26	BJ	0.20	5.0	ug/L	8260C
Trichloroethene (TCE)	6.5		0.20	5.0	ug/L	8260C
cis-1,2-Dichloroethene	3.3	J	0.23	5.0	ug/L	8260C
1,2,4-Trichlorobenzene	7.7	J	1.1	9.1	ug/L	8270D
alpha-BHC	0.27		0.019	0.045	ug/L	8081B
beta-BHC	0.040	J	0.019	0.045	ug/L	8081B
delta-BHC	0.17		0.019	0.045	ug/L	8081B
gamma-BHC (Lindane)	0.21		0.019	0.045	ug/L	8081B
CLIENT ID: WG-9954-070120-SG-025		Lab	ID: R2005	701-003		
Analyte	Results	Flag	MDL	MRL	Units	Method
Toluene	0.27	BJ	0.20	5.0	ug/L	8260C
alpha-BHC	0.13		0.019	0.045	ug/L	8081B
delta-BHC	0.029	J	0.019	0.045	ug/L	8081B
gamma-BHC (Lindane)	0.076		0.019	0.045	ug/L	8081B
CLIENT ID: WG-9954-070120-SG-024		Lab	ID: R2005	701-004		
Analyte	Results	Flag	MDL	MRL	Units	Method
Toluene	0.26 Page	s of 115	0.20	5.0	ug/L	8260C



Analyte Results Flag MDL MRL Units Method	CLIENT ID: WG-9954-070120-SG-024		Lab	ID: R2005	5701-004			
Analyte	Analyte	Results	Flag	MDL	MRL	Units	Method	
1,1,2,2-Tictrloroethane	CLIENT ID: WG-9954-070120-SG-028	Lab ID: R2005701-006						
1,1,2-Trichloroethane 14 0,20 5.0 ug/L 8260C 1,1-Dichloroethane (1,1-DCA) 0,27 J 0,20 5.0 ug/L 8260C 1,1-Dichloroethene (1,1-DCE) 0,78 J 0,20 5.0 ug/L 8260C 2-Butanone (MEK) 6.8 J 0,78 10 ug/L 8260C 2-Hexanone 2.1 J 0,20 10 ug/L 8260C Acetone 71 5.0 10 ug/L 8260C Benzene 1500 E 0,20 5.0 ug/L 8260C Chloroform 140 0,24 5.0 ug/L 8260C Chloroform 140 0,24 5.0 ug/L 8260C Dichloromethane 7.2 0,65 5.0 ug/L 8260C Ethylbenzene 12 0,20 5.0 ug/L 8260C Ethylbenzene 12 0,20 5.0 ug/L 8260C Tolue	Analyte	Results	Flag	MDL	MRL	Units	Method	
1,1-Dichloroethane (1,1-DCA)	1,1,2,2-Tetrachloroethane	7.9		0.20	5.0	ug/L	8260C	
1,1-Dichloroethene (1,1-DCE)	1,1,2-Trichloroethane	14		0.20	5.0	ug/L	8260C	
2-Butanone (MEK) 6.8 J 0.78 10 ug/L 8260C 2-Hexanone 2.1 J 0.20 10 ug/L 8260C Acetone 71 5.0 10 ug/L 8260C Benzene 1500 E 0.20 5.0 ug/L 8260C Chloroform 140 0.24 5.0 ug/L 8260C Dichloromethane 7.2 0.65 5.0 ug/L 8260C Ethylbenzene 12 0.20 5.0 ug/L 8260C Toluene 2000 E 0.20 5.0 ug/L 8260C Trichloroethene (PCE) 130 0.20 5.0 ug/L 8260C Vinyl Chloride 11 0.20 5.0 ug/L 8260C Xylenes, Total 58 0.23 5.0 ug/L 8260C Kylenes, Total 58 0.23 5.0 ug/L 8260C Enzene 6300 D	1,1-Dichloroethane (1,1-DCA)	0.27	J	0.20	5.0	ug/L	8260C	
2-Hexanone	1,1-Dichloroethene (1,1-DCE)	0.78	J	0.20	5.0	ug/L	8260C	
Acetone	2-Butanone (MEK)	6.8	J	0.78	10	ug/L	8260C	
Benzene	2-Hexanone	2.1	J	0.20	10	ug/L	8260C	
Chlorobenzene	Acetone	71		5.0	10	ug/L	8260C	
Chloroform 140 0.24 5.0 ug/L 8260C Dichloromethane 7.2 0.65 5.0 ug/L 8260C Ethylbenzene 12 0.20 5.0 ug/L 8260C Toluene 2000 E 0.21 5.0 ug/L 8260C Trichloroethene (TCE) 130 0.20 5.0 ug/L 8260C Vinyl Chloride 11 0.20 5.0 ug/L 8260C Xylenes, Total 58 0.23 5.0 ug/L 8260C Xylenes, Total 58 0.23 5.0 ug/L 8260C trans-1,2-Dichloroethene 28 0.23 5.0 ug/L 8260C trans-1,2-Dichloroethene 35 0.20 5.0 ug/L 8260C Chloroform 270 D J 48 1000 ug/L 8260C Chloroform 270 D J 48 1000 ug/L 8260C Trichlor	Benzene	1500	Ε	0.20	5.0	ug/L	8260C	
Dichloromethane	Chlorobenzene	850	Ε	0.20	5.0	ug/L	8260C	
Ethylbenzene	Chloroform	140		0.24	5.0	ug/L	8260C	
Tetrachloroethene (PCE) 23 0.21 5.0 ug/L 8260C Toluene 2000 E 0.20 5.0 ug/L 8260C Trichloroethene (TCE) 130 0.20 5.0 ug/L 8260C Vinyl Chloride 11 0.20 5.0 ug/L 8260C Xylenes, Total 58 0.23 5.0 ug/L 8260C cis-1,2-Dichloroethene 28 0.23 5.0 ug/L 8260C trans-1,2-Dichloroethene 35 0.20 5.0 ug/L 8260C Chlorobenzene 6300 D 40 1000 ug/L 8260C Chloroform 270 DJ 48 1000 ug/L 8260C Chlorotene 2400 D 40 1000 ug/L 8260C Chlorotene 270 DJ 48 1000 ug/L 8260C Chlorotene 270 DJ 48 1000 ug/L 8260C Trichloroethene (TCE) 120 DJ 40 1000 ug/L 8260C Trichloroethene (TCE) 120 DJ 40 1000 ug/L 8260C Trichloroethene (TCE) 120 DJ 40 1000 ug/L 8260C Trichloroethene 34 1.1 9.1 ug/L 8270D 1,2-Dichlorobenzene 34 1.1 9.1 ug/L 8270D 1,3-Dichlorobenzene 4.1 J 0.92 9.1 ug/L 8270D 1,4-Dichlorobenzene 97 1.1 9.1 ug/L 8270D 2,4-Dichlorophenol 26 0.99 9.1 ug/L 8270D 2,4-Dichlorophenol 520 E 1.2 9.1 ug/L 8270D 2,4-Dichlorophenol 32 0.91 9.1 ug/L 8270D 2,4-Dimethylphenol 32 0.91 9.1 ug/L 8270D 2-Chlorophenol 31 0.97 9.1 ug/L 8270D 2-Chlorophenol 32 0.91 9.1 ug/L 8270D 2-Methylphenol 32 0.91 9.1 ug/L 8270D 3- and 4-Methylphenol Coelution 71 1.1 9.1 ug/L 8270D Benzoic Acid 4100 E 33 91 ug/L 8270D Benzoic Acid 4100 E 33 91 ug/L 8270D Benzoic Acid 4100 E 33 91 ug/L 8270D	Dichloromethane	7.2		0.65	5.0	ug/L	8260C	
Toluene 2000 E 0.20 5.0 ug/L 8260C Trichloroethene (TCE) 130 0.20 5.0 ug/L 8260C Vinyl Chloride 11 0.20 5.0 ug/L 8260C Xylenes, Total 58 0.23 5.0 ug/L 8260C cis-1,2-Dichloroethene 28 0.23 5.0 ug/L 8260C trans-1,2-Dichloroethene 35 0.20 5.0 ug/L 8260C Benzene 6300 D 40 1000 ug/L 8260C Chloroform 270 DJ 48 1000 ug/L 8260C Chloroform 270 DJ 48 1000 ug/L 8260C Trichloroethene (TCE) 120 DJ 40 1000 ug/L 8260C Trichloroethene (TCE) 120 DJ 40 1000 ug/L 8260C Trichloroethene (TCE) 120 DJ 40 1000 ug/L	Ethylbenzene	12		0.20	5.0	ug/L	8260C	
Trichloroethene (TCE) 130 0.20 5.0 ug/L 8260C Vinyl Chloride 11 0.20 5.0 ug/L 8260C Xylenes, Total 58 0.23 5.0 ug/L 8260C cis-1,2-Dichloroethene 28 0.23 5.0 ug/L 8260C trans-1,2-Dichloroethene 35 0.20 5.0 ug/L 8260C Benzene 6300 D 40 1000 ug/L 8260C Chloroberzene 2400 D 40 1000 ug/L 8260C Chloroberzene 2400 D 40 1000 ug/L 8260C Chloroberzene 21000 D 40 1000 ug/L 8260C Trichloroethene (TCE) 120 DJ 40 1000 ug/L 8260C Trichlorobenzene (TCE) 120 DJ 40 1000 ug/L 8260C Trichlorobenzene (TCE) 120 DJ 40 1000 <td< td=""><td>Tetrachloroethene (PCE)</td><td>23</td><td></td><td>0.21</td><td>5.0</td><td>ug/L</td><td>8260C</td></td<>	Tetrachloroethene (PCE)	23		0.21	5.0	ug/L	8260C	
Vinyl Chloride 11 0.20 5.0 ug/L 8260C Xylenes, Total 58 0.23 5.0 ug/L 8260C cis-1,2-Dichloroethene 28 0.23 5.0 ug/L 8260C trans-1,2-Dichloroethene 35 0.20 5.0 ug/L 8260C Benzene 6300 D 40 1000 ug/L 8260C Chlorobenzene 2400 D 40 1000 ug/L 8260C Chloroform 270 DJ 48 1000 ug/L 8260C Toluene 21000 D 40 1000 ug/L 8260C Trichloroethene (TCE) 120 DJ 40 1000 ug/L 8260C trans-1,2-Dichloroethene 51 DJ 40 1000 ug/L 8260C trans-1,2-Dichlorobenzene 100 1.1 9.1 ug/L 8270D 1,2-Trichlorobenzene 34 1.1 9.1 ug/L 8270D	Toluene	2000	Е	0.20	5.0	ug/L	8260C	
Xylenes, Total 58 0.23 5.0 ug/L 8260C cis-1,2-Dichloroethene 28 0.23 5.0 ug/L 8260C trans-1,2-Dichloroethene 35 0.20 5.0 ug/L 8260C Benzene 6300 D 40 1000 ug/L 8260C Chlorobenzene 2400 D 40 1000 ug/L 8260C Chloroform 270 DJ 48 1000 ug/L 8260C Toluene 21000 D 40 1000 ug/L 8260C Trichloroethene (TCE) 120 DJ 40 1000 ug/L 8260C Trichloroethene (TCE) 120 DJ 40 1000 ug/L 8260C Trichloroethene (TCE) 120 DJ 40 1000 ug/L 8260C trans-1,2-Dichloroethene 51 DJ 40 1000 ug/L 8260C trans-1,2-Dichlorobenzene 34 1.1	Trichloroethene (TCE)	130		0.20	5.0	ug/L	8260C	
Xylenes, Total 58 0.23 5.0 ug/L 8260C cis-1,2-Dichloroethene 28 0.23 5.0 ug/L 8260C trans-1,2-Dichloroethene 35 0.20 5.0 ug/L 8260C Benzene 6300 D 40 1000 ug/L 8260C Chlorobenzene 2400 D 40 1000 ug/L 8260C Chloroform 270 DJ 48 1000 ug/L 8260C Toluene 21000 D 40 1000 ug/L 8260C Trichloroethene (TCE) 120 DJ 40 1000 ug/L 8260C Trichlorobenzene (TCE) 120 DJ 40 <td>Vinyl Chloride</td> <td>11</td> <td></td> <td>0.20</td> <td>5.0</td> <td>_</td> <td>8260C</td>	Vinyl Chloride	11		0.20	5.0	_	8260C	
cis-1,2-Dichloroethene 28 0.23 5.0 ug/L 8260C trans-1,2-Dichloroethene 35 0.20 5.0 ug/L 8260C Benzene 6300 D 40 1000 ug/L 8260C Chlorobenzene 2400 D 40 1000 ug/L 8260C Chloroform 270 DJ 48 1000 ug/L 8260C Toluene 21000 D 40 1000 ug/L 8260C Trichloroethene (TCE) 120 DJ 40 1000 ug/L 8260C 1,2,4-Trichloroethene (TCE) 34	Xylenes, Total	58		0.23	5.0	_	8260C	
trans-1,2-Dichloroethene 35 0.20 5.0 ug/L 8260C Benzene 6300 D 40 1000 ug/L 8260C Chlorobenzene 2400 D 40 1000 ug/L 8260C Chloroform 270 DJ 48 1000 ug/L 8260C Toluene 21000 D 40 1000 ug/L 8260C Trichloroethene (TCE) 120 DJ 40 1000 ug/L 8270D 1,2-4-Trichloroethene (TCE)	cis-1,2-Dichloroethene	28		0.23	5.0	_	8260C	
Benzene 6300 D 40 1000 ug/L 8260C Chlorobenzene 2400 D 40 1000 ug/L 8260C Chloroform 270 DJ 48 1000 ug/L 8260C Toluene 21000 D 40 1000 ug/L 8260C Trichloroethene (TCE) 120 DJ 40 1000 ug/L 8260C trans-1,2-Dichloroethene 51 DJ 40 1000 ug/L 8260C 1,2,4-Trichlorobenzene 100 1.1 9.1 ug/L 8270D 1,2-Dichlorobenzene 34 1.1 9.1 ug/L 8270D 1,3-Dichlorobenzene 4.1 J 0.92 9.1 ug/L 8270D 1,4-Dichlorobenzene 97 1.1 9.1 ug/L 8270D 2,4,5-Trichlorophenol 26 0.99 9.1 ug/L 8270D 2,4-Dinethylphenol 520 E 1.2 9.1	trans-1,2-Dichloroethene	35		0.20	5.0	_	8260C	
Chlorobenzene 2400 D 40 1000 ug/L 8260C Chloroform 270 DJ 48 1000 ug/L 8260C Toluene 21000 D 40 1000 ug/L 8260C Trichloroethene (TCE) 120 DJ 40 1000 ug/L 8260C trans-1,2-Dichloroethene 51 DJ 40 1000 ug/L 8260C 1,2,4-Trichlorobenzene 100 1.1 9.1 ug/L 8270D 1,2-Dichlorobenzene 34 1.1 9.1 ug/L 8270D 1,3-Dichlorobenzene 4.1 J 0.92 9.1 ug/L 8270D 1,4-Dichlorobenzene 97 1.1 9.1 ug/L 8270D 2,4,5-Trichlorophenol 26 0.99 9.1 ug/L 8270D 2,4-Dichlorophenol 520 E 1.2 9.1 ug/L 8270D 2-Chlorophenol 31 0.97 9.1 ug/		6300	D	40	1000	_	8260C	
Chloroform 270 DJ 48 1000 ug/L 8260C Toluene 21000 D 40 1000 ug/L 8260C Trichloroethene (TCE) 120 DJ 40 1000 ug/L 8260C trans-1,2-Dichloroethene 51 DJ 40 1000 ug/L 8260C 1,2,4-Trichlorobenzene 100 1.1 9.1 ug/L 8270D 1,2-Dichlorobenzene 34 1.1 9.1 ug/L 8270D 1,3-Dichlorobenzene 4.1 J 0.92 9.1 ug/L 8270D 1,4-Dichlorobenzene 97 1.1 9.1 ug/L 8270D 2,4,5-Trichlorophenol 26 0.99 9.1 ug/L 8270D 2,4-Dichlorophenol 520 E 1.2 9.1 ug/L 8270D 2,4-Dimethylphenol 31 0.97 9.1 ug/L 8270D 2-Methylphenol 32 0.91 9.1 ug/L	Chlorobenzene	2400	D	40	1000	_	8260C	
Toluene 21000 D 40 1000 ug/L 8260C Trichloroethene (TCE) 120 DJ 40 1000 ug/L 8260C trans-1,2-Dichloroethene 51 DJ 40 1000 ug/L 8260C 1,2,4-Trichlorobenzene 100 1.1 9.1 ug/L 8270D 1,2-Dichlorobenzene 34 1.1 9.1 ug/L 8270D 1,3-Dichlorobenzene 4.1 J 0.92 9.1 ug/L 8270D 1,4-Dichlorophenzene 97 1.1 9.1 ug/L 8270D 2,4,5-Trichlorophenol 26 0.99 9.1 ug/L 8270D 2,4-Dichlorophenol 520 E 1.2 9.1 ug/L 8270D 2,4-Dimethylphenol 6.9 J 1.3 9.1 ug/L 8270D 2-Methylphenol 32 0.91 9.1 ug/L 8270D 3- and 4-Methylphenol 40 0.98 9.1 ug/L <td>Chloroform</td> <td>270</td> <td>DJ</td> <td>48</td> <td>1000</td> <td>_</td> <td>8260C</td>	Chloroform	270	DJ	48	1000	_	8260C	
Trichloroethene (TCE) 120 DJ 40 1000 ug/L 8260C trans-1,2-Dichloroethene 51 DJ 40 1000 ug/L 8260C 1,2,4-Trichlorobenzene 100 1.1 9.1 ug/L 8270D 1,2-Dichlorobenzene 34 1.1 9.1 ug/L 8270D 1,3-Dichlorobenzene 4.1 J 0.92 9.1 ug/L 8270D 1,4-Dichlorobenzene 97 1.1 9.1 ug/L 8270D 2,4,5-Trichlorophenol 26 0.99 9.1 ug/L 8270D 2,4-Dichlorophenol 520 E 1.2 9.1 ug/L 8270D 2,4-Dimethylphenol 6.9 J 1.3 9.1 ug/L 8270D 2-Chlorophenol 31 0.97 9.1 ug/L 8270D 2-Methylphenol 32 0.91 9.1 ug/L 8270D 3- and 4-Methylphenol 40 0.98 9.1 ug/L	Toluene	21000			1000	_	8260C	
trans-1,2-Dichloroethene 51 DJ 40 1000 ug/L 8260C 1,2,4-Trichlorobenzene 100 1.1 9.1 ug/L 8270D 1,2-Dichlorobenzene 34 1.1 9.1 ug/L 8270D 1,3-Dichlorobenzene 4.1 J 0.92 9.1 ug/L 8270D 1,4-Dichlorobenzene 97 1.1 9.1 ug/L 8270D 2,4,5-Trichlorophenol 26 0.99 9.1 ug/L 8270D 2,4-Dichlorophenol 520 E 1.2 9.1 ug/L 8270D 2,4-Dimethylphenol 6.9 J 1.3 9.1 ug/L 8270D 2-Chlorophenol 31 0.97 9.1 ug/L 8270D 2-Methylphenol 32 0.91 9.1 ug/L 8270D 3- and 4-Methylphenol Coelution 71 1.1 9.1 ug/L 8270D 4-Chloro-3-methylphenol 40 0.98 9.1 ug/L 8270D Benzoic Acid 4100 E 33 91 ug/L 8270D Benzoic Acid 4100 E 33 91 ug/L 8270D Benzyl Alcohol 290 1.5 9.1 ug/L 8270D	Trichloroethene (TCE)	120	DJ	40	1000	_	8260C	
1,2,4-Trichlorobenzene 100 1.1 9.1 ug/L 8270D 1,2-Dichlorobenzene 34 1.1 9.1 ug/L 8270D 1,3-Dichlorobenzene 4.1 J 0.92 9.1 ug/L 8270D 1,4-Dichlorobenzene 97 1.1 9.1 ug/L 8270D 2,4,5-Trichlorophenol 26 0.99 9.1 ug/L 8270D 2,4-Dichlorophenol 520 E 1.2 9.1 ug/L 8270D 2,4-Dimethylphenol 6.9 J 1.3 9.1 ug/L 8270D 2-Chlorophenol 31 0.97 9.1 ug/L 8270D 2-Methylphenol 32 0.91 9.1 ug/L 8270D 3- and 4-Methylphenol Coelution 71 1.1 9.1 ug/L 8270D 4-Chloro-3-methylphenol 40 0.98 9.1 ug/L 8270D Benzoic Acid 4100 E 33 91 ug/L 8270D Benzyl Alcohol 290 1.5 9.1 ug/L 8270D </td <td>` ,</td> <td></td> <td></td> <td></td> <td></td> <td>_</td> <td></td>	` ,					_		
1,2-Dichlorobenzene 34 1.1 9.1 ug/L 8270D 1,3-Dichlorobenzene 4.1 J 0.92 9.1 ug/L 8270D 1,4-Dichlorobenzene 97 1.1 9.1 ug/L 8270D 2,4,5-Trichlorophenol 26 0.99 9.1 ug/L 8270D 2,4-Dichlorophenol 520 E 1.2 9.1 ug/L 8270D 2,4-Dimethylphenol 6.9 J 1.3 9.1 ug/L 8270D 2-Chlorophenol 31 0.97 9.1 ug/L 8270D 2-Methylphenol 32 0.91 9.1 ug/L 8270D 3- and 4-Methylphenol Coelution 71 1.1 9.1 ug/L 8270D 4-Chloro-3-methylphenol 40 0.98 9.1 ug/L 8270D Benzyl Alcohol 290 1.5 9.1 ug/L 8270D Beigle ablancethyl) Ether 34 1.2 0.4 0.4 0.4 0.4 0.4 0.4 0.4 0.4 0.4 0.4 0.4 0.4 <td>1,2,4-Trichlorobenzene</td> <td>100</td> <td></td> <td>1.1</td> <td>9.1</td> <td>_</td> <td>8270D</td>	1,2,4-Trichlorobenzene	100		1.1	9.1	_	8270D	
1,3-Dichlorobenzene 4.1 J 0.92 9.1 ug/L 8270D 1,4-Dichlorobenzene 97 1.1 9.1 ug/L 8270D 2,4,5-Trichlorophenol 26 0.99 9.1 ug/L 8270D 2,4-Dichlorophenol 520 E 1.2 9.1 ug/L 8270D 2,4-Dimethylphenol 6.9 J 1.3 9.1 ug/L 8270D 2-Chlorophenol 31 0.97 9.1 ug/L 8270D 2-Methylphenol 32 0.91 9.1 ug/L 8270D 3- and 4-Methylphenol Coelution 71 1.1 9.1 ug/L 8270D 4-Chloro-3-methylphenol 40 0.98 9.1 ug/L 8270D Benzoic Acid 4100 E 33 91 ug/L 8270D Benzyl Alcohol 290 1.5 9.1 ug/L 8270D				1.1	9.1	_		
1,4-Dichlorobenzene 97 1.1 9.1 ug/L 8270D 2,4,5-Trichlorophenol 26 0.99 9.1 ug/L 8270D 2,4-Dichlorophenol 520 E 1.2 9.1 ug/L 8270D 2,4-Dimethylphenol 6.9 J 1.3 9.1 ug/L 8270D 2-Chlorophenol 31 0.97 9.1 ug/L 8270D 2-Methylphenol 32 0.91 9.1 ug/L 8270D 3- and 4-Methylphenol Coelution 71 1.1 9.1 ug/L 8270D 4-Chloro-3-methylphenol 40 0.98 9.1 ug/L 8270D Benzoic Acid 4100 E 33 91 ug/L 8270D Benzyl Alcohol 290 1.5 9.1 ug/L 8270D	1,3-Dichlorobenzene		J			_		
2,4,5-Trichlorophenol 26 0.99 9.1 ug/L 8270D 2,4-Dichlorophenol 520 E 1.2 9.1 ug/L 8270D 2,4-Dimethylphenol 6.9 J 1.3 9.1 ug/L 8270D 2-Chlorophenol 31 0.97 9.1 ug/L 8270D 2-Methylphenol 32 0.91 9.1 ug/L 8270D 3- and 4-Methylphenol Coelution 71 1.1 9.1 ug/L 8270D 4-Chloro-3-methylphenol 40 0.98 9.1 ug/L 8270D Benzoic Acid 4100 E 33 91 ug/L 8270D Benzyl Alcohol 290 1.5 9.1 ug/L 8270D						_		
2,4-Dichlorophenol 520 E 1.2 9.1 ug/L 8270D 2,4-Dimethylphenol 6.9 J 1.3 9.1 ug/L 8270D 2-Chlorophenol 31 0.97 9.1 ug/L 8270D 2-Methylphenol 32 0.91 9.1 ug/L 8270D 3- and 4-Methylphenol Coelution 71 1.1 9.1 ug/L 8270D 4-Chloro-3-methylphenol 40 0.98 9.1 ug/L 8270D Benzoic Acid 4100 E 33 91 ug/L 8270D Benzyl Alcohol 290 1.5 9.1 ug/L 8270D Bic/2 ablescetbyl) Ether 34 1.2 9.4 1.2 9.2 1.2 9.2 <		26				_		
2,4-Dimethylphenol 6.9 J 1.3 9.1 ug/L 8270D 2-Chlorophenol 31 0.97 9.1 ug/L 8270D 2-Methylphenol 32 0.91 9.1 ug/L 8270D 3- and 4-Methylphenol Coelution 71 1.1 9.1 ug/L 8270D 4-Chloro-3-methylphenol 40 0.98 9.1 ug/L 8270D Benzoic Acid 4100 E 33 91 ug/L 8270D Benzyl Alcohol 290 1.5 9.1 ug/L 8270D Bio (2 ablass othyl) Ethos 34 1.2 0.4 ug/L 8270D	•		Е			_		
2-Chlorophenol 31 0.97 9.1 ug/L 8270D 2-Methylphenol 32 0.91 9.1 ug/L 8270D 3- and 4-Methylphenol Coelution 71 1.1 9.1 ug/L 8270D 4-Chloro-3-methylphenol 40 0.98 9.1 ug/L 8270D Benzoic Acid 4100 E 33 91 ug/L 8270D Benzyl Alcohol 290 1.5 9.1 ug/L 8270D Bio (2 ablescettyd) Ether 34 1.2 0.4 ug/L 8270D	•					_		
2-Methylphenol 32 0.91 9.1 ug/L 8270D 3- and 4-Methylphenol Coelution 71 1.1 9.1 ug/L 8270D 4-Chloro-3-methylphenol 40 0.98 9.1 ug/L 8270D Benzoic Acid 4100 E 33 91 ug/L 8270D Benzyl Alcohol 290 1.5 9.1 ug/L 8270D Bio (2 ablescetbyl) Ether 34 1.2 0.4 1.2 9.2 1.2 9.2<	• •					_		
3- and 4-Methylphenol Coelution 71 1.1 9.1 ug/L 8270D 4-Chloro-3-methylphenol 40 0.98 9.1 ug/L 8270D Benzoic Acid 4100 E 33 91 ug/L 8270D Benzyl Alcohol 290 1.5 9.1 ug/L 8270D Bio(3 ablass atbyl) Ether	•					_		
4-Chloro-3-methylphenol 40 0.98 9.1 ug/L 8270D Benzoic Acid 4100 E 33 91 ug/L 8270D Benzyl Alcohol 290 1.5 9.1 ug/L 8270D Bio(3 chloroothyl) Ethor 34 1.2 0.1 ug/L 8270D						_		
Benzoic Acid 4100 E 33 91 ug/L 8270D Benzyl Alcohol 290 1.5 9.1 ug/L 8270D Bio (3 oblescetbyl) Ether 34 4.3 0.4 ug/L 8370D	· ·					_		
Benzyl Alcohol 290 1.5 9.1 ug/L 8270D	* *		Е			_		
Dio/2 obless of bull Ether 24 4.2 0.4 us// 9270D			_			_		
	Bis(2-chloroethyl) Ether	24		1.2	9.1	ug/L	8270D	



LIENT ID: WG-9954-070120-SG-028	Lab ID: R2005701-006					
Analyte	Results	Flag	MDL	MRL	Units	Method
Hexachlorobutadiene	1.0	J	0.91	9.1	ug/L	8270D
Phenol	44		0.91	9.1	ug/L	8270D
1,2,4-Trichlorobenzene	76	J	26	230	ug/L	8270D
1,2-Dichlorobenzene	37	J	27	230	ug/L	8270D
1,4-Dichlorobenzene	100	J	26	230	ug/L	8270D
2,4,5-Trichlorophenol	34	J	25	230	ug/L	8270D
2,4-Dichlorophenol	350	D	29	230	ug/L	8270D
2-Chlorophenol	25	J	25	230	ug/L	8270D
2-Methylphenol	33	J	23	230	ug/L	8270D
3- and 4-Methylphenol Coelution	60	J	27	230	ug/L	8270D
4-Chloro-3-methylphenol	30	J	25	230	ug/L	8270D
Benzoic Acid	9200	E	820	2300	ug/L	8270D
Benzyl Alcohol	250		36	230	ug/L	8270D
Phenol	37	J	23	230	ug/L	8270D
1,2,4-Trichlorobenzene	72	J	51	450	ug/L	8270D
1,4-Dichlorobenzene	100	J	52	450	ug/L	8270D
2,4-Dichlorophenol	320	J	57	450	ug/L	8270D
3- and 4-Methylphenol Coelution	60	J	53	450	ug/L	8270D
Benzoic Acid	10000	D	1700	4500	ug/L	8270D
Benzyl Alcohol	220	J	71	450	ug/L	8270D
Aldrin	0.85		0.19	0.45	ug/L	8081B
alpha-BHC	25		0.19	0.45	ug/L	8081B
beta-BHC	6.9		0.19	0.45	ug/L	8081B
delta-BHC	8.7		0.19	0.45	ug/L	8081B
gamma-BHC (Lindane)	6.2		0.19	0.45	ug/L	8081B
alpha-BHC	18		0.19	0.45	ug/L	8081B
delta-BHC	6.3		0.19	0.45	ug/L	8081B
gamma-BHC (Lindane)	3.6		0.19	0.45	ug/L	8081B
gamma-Chlordane	1.5		0.19	0.45	ug/L	8081B
Aldrin	0.66		0.19	0.45	ug/L	8081B
Heptachlor	0.40	J	0.19	0.45	ug/L	8081B
beta-BHC	4.4		0.19	0.45	ug/L	8081B



Sample Receipt Information

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request:R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

SAMPLE CROSS-REFERENCE

SAMPLE #	CLIENT SAMPLE ID	<u>DATE</u>	<u>TIME</u>
R2005701-001	WG-9954-070120-SG-027	7/1/2020	1150
R2005701-002	WG-9954-070120-SG-026	7/1/2020	1045
R2005701-003	WG-9954-070120-SG-025	7/1/2020	0955
R2005701-004	WG-9954-070120-SG-024	7/1/2020	0845
R2005701-005	TB-9954-070120-SG-005	7/1/2020	0000
R2005701-006	WG-9954-070120-SG-028	7/1/2020	1150

CHAIN-OF-CUSTODY/Analytical Request Document The Chain-of-Custody is a LEGAL DOCUMENT. All relevant fields must be completed accurately.

Client Information					
GLEN SPRINGS HOLDINGS INC	Report To: Kathy Willy				
806 97TH STREET	Сору То:				
LOVE CANAL					
NIAGARA FALLS, NEW YORK 14304	Invoice To:				
Phone: 716-283-0111	PO:				
Fax: 716-283-2866	Project Name: LOVE CANAL ANNUAL GW				
Email: kathy.willy@ghd.com	Project Number: 9964				

	Information
Laboratory: ALS	
Laboratory Location: 15 BUILDING 300, SUITE 30	
ROCHESTER, NY 14623	
ROCHESTER, NY 14623	
Laboratory Contact: BR	ADY KALKMAN
Requested Due Date:	TAT: 10
QA/QC Requirements:	

Event Inform	nation
ID#: LC ANNUAL GW SA	MPLING 2020-05-1
SSOW Ref#: 292-	402-999-3100
ampler Name: S GARDN	ER. D. TVRAN

	Valid Matrix Code WG Groundwater WB Borehole Water								Sample Cond	lition
	WS Surface Water			Ţ.	<u>e</u>				Temp in C	
	SO Soil SE Sediment	oge	Collected	ected	s(None)	9		1	Received on ice	Y/N
		Ü	👸	Colle	l mi	ļ ē	ξ		Sealed Cooler	Y/N
		Matrix	Date (Time (estPC	SVOC(none)	VOA(HCI)		Samples Intact	Y/N
Sample Identification		Ž	ة	Ë	ď	6	>	Remarks		
WG-9954-070120-SG-027		WG	07/01/2020	11:50	2	2	3			
WG-9954-070120-SG-026		WG	07/01/2020	10:45	2	2	3			
WG-9954-070120-SG-025	<u></u>	WG	07/01/2020	09:55	2	2	3			
WG-9954-070120-SG-024		WG	07/01/2020	08:45	2	2	3			
TB-9954-070120-SG-005		WG Q	07/01/2020	00:00	-	-	3			
WG-9954-070120-SG-028		WG	07/01/2020	11:50	2	2	3			
Total Bottles				-	10	10	18	Grand Tota	ıl:38	····

SHIPMENT METHOD	NO. OF COOLERS	RELINQUISHED BY:	DATE	TIME	RECIEVED BY:	DATE	TIME
FedEx	1	Show Marcher	71/20	1400	Seem hat	7/2/1013	1050
AIRBILL#:					- 4 John	//	-

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ALS Enuir	ALS Environmental					1565			(585)	288-5		FAX (585) 2	ocheste 208-84	er, NY 14623 175	004 011	4, 005, 006, 007, 008, 009, 010, 1, 012, 013		T030477
Project Name: Love Canal:292-402-D02-3100						70		14D										,	
Project Number:	Report To Kathy Will	у] _	_			-	ļ	_		_	т —	ļ					
9954 Annual Long Term Monitoring Company / Address GHD Services Inc. 2055 Niagara Falls Blvd., Suite 3 Niagara Falls NY, 14304		1		CONTAINERS	8	:		e.											
Phone #	FAX#				est	ÇB	8	70C		1									
716-297-2160 Sampler Signature	716-297-2 Sampler F	rinted Name		1 8	=	/ P	S	-			l								•
				NUMBER OF	8081B / Pest	8082A / PCB	8270D / SVO	8260C			1				Domosli				
CLIENT SAMPLE ID	LABID	SAMPLING Date Time	Matrix	+	ĕ	ĕ	86	86	_	a	က	4	 "		Remark	(8			
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			Lkquid	1			Г	1			Ì		1	1					
2. 3.		<u> </u>	Liquid		1			-	\Box	Π				1					
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10.			Llquid					$oldsymbol{ol}}}}}}}}}}}}}}}}}}$		<u> </u>				1					
Special Instructions/Co	mments:							Tu					•		ents		eport Requirements		Invoice Information
,	•										SURC 1 (3 t			APP	PLY)	— II.	Results Only Results + QC Summaries (LCS, UP, MS/MSD as required)	P.	O.#
									•								Results + QC and Cilibration Summeries	В	ill To:
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									F	teque	sted	Rep	ort Da	ate			YesNo	=	
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Printed Name	Printed N	lame .	Prir	ited N	lame					7	Printe	ed N	ame				Printed Name		Printed Name
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Date/Time	Date/Time Date				ate/Time				70	Date/Time					Date/Time		Date/Time		



Cooler Receipt and Preservation Check Form

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<i>(</i>	R2005701 5	`
I	GHD Services (nc. Love Canal:292-402-D02-3100	
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ı		

2 Custody papers properly completed (ink, signed)? N 3 Did all bottles arrive in good condition (unbroken)? N 4 Circle: Wet Co Dry Ice Gel packs present? N 5 Where did the bottles originate? ALS/ROC CLIENT 7 Soil VOA received as: Bulk Encore 5035set NA 8. Temperature Readings Date: 1/2/102/20 Time: 1/2/20 1/20 1/2		ent_ <i>GH</i> 1	4-1			_Folder Nı	ımber	_			<u>.</u> .				
2 Custody papers properly completed (ink, signed)? N 3 Did all bottles arrive in good condition (unbroken)* N 4 Circle: (WeTC) Dry Ice Gel packs present? N 5 Did VOA received as: Bulk Encore 5035set NA 5. Temperature Readings Date:	Cooler receiv	ed on 7/2	zporo .	by: 12	100	CO	URIER:	ALS	UPS	PEDE	X) VEL	OCITY	CLIF	ENT	
3 Did all bottles arrive in good condition (unbroken)? N 4 Circle: Wet C Dry Ice Gel packs present? N 5 Soil VOA received as: Bulk Encore 5035sct NA 1 Temperature Readings Date: 1/2/22, Time: //LD ID: IR#7 RRFID From: Comp Blank Sample Bo Observed Temp (°C) / 3	1 Were Cu	istody seals or	outside of coole	r?	P) N 5a	Perch	lorate s	amples	have rec	quired he	adspac	e?	Y N	V (VA)
4 Circle: WetTee Dry Ice Gel packs present? N 7 Soil VOA received as: Bulk Encore 5035set NA Temperature Readings Date: 7/1/2020. Time: 1/LD ID: IR#7 IR#10 From: Comp Blank Sample Bo Observed Temp (°C) 1/3 0.9 Within 0-6°C? N N N N Y N Y N Y N Y N Y N Y N Y N Y	2 Custody	papers prope	rly completed (in	k, signed	1)? 😗) N 5b	Did	OA vial	s, Alk,	r Sulfid	e have si	g* bub	bles?	Ý) N	NA NA
Temperature Readings Date:	3 Did all b	ottles arrive in	good condition (unbroke	n)?(Y) N 6	Where	did the	bottles	originat	e?	ALS/	ROC>	CLIE	NT
Observed Temp (°C) Within 0-6°C? Within 0-6°C? Within 0-6°C? Within 0-6°C? Within 0-6°C? Within 0-6°C? Within 0-6°C, were samples frozen? Within 0-6°C? Within 0-6	4 Circle: (Wet Ice Dry	Ice Gel packs	prese	nt? (Y	N 7	Soil V	OA rec	eived as	; Bı	ılk E	ncore	5035	set 🔿	₹A>
Within 0-6°C? ON ON YN YN YN YN YN YN	. Temperatu	re Readings	Date: 7/2/	7 <u>07υ</u> Τ	ime:	1/12	ID:	IR#7	I R#10		From:	emp	Blank	Sam	ple Bottle
Within 0-6°C? Y N	Observed To	emp (°C)	1.3		0.9				1						
If out of Temperature, note packing/ice condition: Ice melted Poorly Packed (described below) Same Day Ru & Client Approval to Run Samples: Standing Approval Client aware at drop-off Client notified by:	Within 0-6°	C?	/ (Y) N	7		1 Y	N	Y	N	Y	N	Y	N	Y	N
All samples held in storage location: All samples held in storage location: 5035 samples placed in storage location: Description: Description: Description: All samples held in storage location: Description: Descriptio	If <0°C, wer	re samples froz	zen? Y N		Ý N	I Y	N	Y	N	Y	N	Y	N	Y	N
## Client Approval to Run Samples: Standing Approval Client aware at drop-off Client notified by: ### All samples held in storage location: But by ### on ### on ### at ### within 48 hours of sampling? Y N ### Cooler Breakdown/Preservation Check**: Date: ### D	If out of 7	Temperaturc.	note packing/ic	e conditi	ion:	•	Ice melt	ed P	oorly P	acked (d	lescribed	below)	Same I	Day Rule
All samples held in storage location: 5035 samples placed in storage location: by Solution S							_		-	· · · · · · · · · · · · · · · · · · ·					•
9. Were all bottle labels complete (i.e. analysis, preservation, etc.)? 10. Did all bottle labels and tags agree with custody papers? 11. Were correct containers used for the tests indicated? 12. Were 5035 vials acceptable (no extra labels, not leaking)? 13. Air Samples: Cassettes / Tubes Intact with MS? Canisters Pressurized pH	-			<u> </u>		<i></i>	1900	- -	_	within 4	8 hours	of sam	pling?	Y	N
Paper Yes No	9.	Were all bottle	labels complete (i a anal	ucic ni	eceruation e				-					
≥12	11. Y 12. Y	Were correct co Were 5035 vial	abels and tags agr ontainers used for Is acceptable (no	ee with of the tests extra lab	custody s indica els, no	y papers? ated? ot leaking)?	·	rized	Y Y Y	ES ES ES Cedlar®	NO NO NO	flated		X 7A (N/A)	
Section Sect	11. Y 12. Y 13. A	Were correct co Were 5035 vial Air Samples: O	abels and tags agr ontainers used for Is acceptable (no Cassettes / Tubes	ee with of the tests extra lab Intact w	custody s indica els, no rith MS	y papers? ated? ot leaking)? S? Caniste	ers Pressur	_	Samp	le ID	NO NO NO Bags In: Vol.	L	ot Adde	X7A N/A	Final
Second Second	11. Y 12. Y 13. P	Were correct co Were 5035 vial Air Samples: C Lot of test	abels and tags agrontainers used for ls acceptable (no assettes / Tubes Reagent	ee with of the tests extra lab Intact w	custody s indica els, no rith MS	y papers? ated? ot leaking)? S? Caniste	ers Pressur	_	Samp	le ID	NO NO NO Bags In: Vol.	L	ot Adde	N/A N/A	
S-9 For 608pest No=Notify for 3day	11. Y 12. Y 13. pH ≥12	Were correct co Were 5035 vial Air Samples: C Lot of test	abels and tags agrontainers used for its acceptable (no Cassettes / Tubes Reagent	ee with of the tests extra lab Intact w	custody s indica els, no rith MS	y papers? ated? ot leaking)? S? Caniste	ers Pressur	_	Samp	le ID	NO NO NO Bags In: Vol.	L	ot Adde	₹7A N/A ed	
S-9	11. Y 12. Y 13. pH ≥12 ≤2	Were correct co Were 5035 vial Air Samples: C Lot of test	abels and tags agrontainers used for ls acceptable (no Cassettes / Tubes Reagent NaOH HNO3	ee with of the tests extra lab Intact w	custody s indica els, no rith MS	y papers? ated? ot leaking)? S? Caniste	ers Pressur	_	Samp	le ID	NO NO NO Bags In: Vol.	L	ot Adde	N/A ed	
Residual Chlorine (-)	11. V 12. V 13. pH ≥12 ≤2 ≤2	Were correct co Were 5035 vial Air Samples: C Lot of test	abels and tags agrontainers used for ls acceptable (no Cassettes / Tubes Reagent NaOH HNO ₃ H ₂ SO ₄	ee with of the tests extra lab Intact w	custody s indica els, no rith MS	y papers? ated? ot leaking)? S? Caniste	ers Pressur	_	Samp	le ID	NO NO NO Bags In: Vol.	L	ot Adde	MA N/A ed	
Chlorine (-) Phenol, 625, 608pest, 522 Na ₂ S ₂ O ₃ ZnAcetate HCl ** ** CN), ascorbic (phenol). **VOAs and 1664 Not to be tested before analysis. Otherwise, all bottles of all samples with chemical preservative are checked (not just representatives). Bottle lot numbers: 2538, 05//20 - 18mc Explain all Discrepancies/ Other Comments:	11. V 12. V 13. pH ≥12 ≤2 ≤2 <4	Were correct co Were 5035 vial Air Samples: C Lot of test	abels and tags agrontainers used for ls acceptable (no Cassettes / Tubes Reagent NaOH HNO3 H ₂ SO ₄ NaHSO ₄	ee with of the tests extra lab Intact w	custody s indica els, no rith MS ed? No	y papers? ated? at leaking)? S? Caniste Lot Receive	ors Pressur d	_	Samp	le ID	NO NO NO Bags In: Vol.	L	ot Adde	NA NA a	
(-) 608pest, 522 CN), ascorbic (phenol). Na ₂ S ₂ O ₃ ZnAcetate Otherwise, all bottles of all samples with chemical preservative are checked (not just representatives). Bottle lot numbers: 2538, 05/120 - 18mc Explain all Discrepancies/ Other Comments:	11. Y 12. Y 13. pH ≥12 ≤2 ≤2 <4 5-9	Were correct co Were 5035 vial Air Samples: C Lot of test	hels and tags agrontainers used for ls acceptable (no cassettes / Tubes Reagent NaOH HNO3 H ₂ SO ₄ NaHSO ₄ For 608pest	ee with of the tests extra lab Intact w	custody s indica eels, no eith MS ed? No	y papers? ated? It leaking)? Caniste Lot Receive No=Notify fo	d d r 3day	_	Samp	le ID	NO NO NO Bags In: Vol.	L	ot Adde	NA NA ed	
Na ₂ S ₂ O ₃ ZnAcetate Otherwise, all bottles of all samples with chemical preservative are checked (not just representatives). Bottle lot numbers: 2538, 05/120 - 18mc Explain all Discrepancies/ Other Comments:	11. V 12. V 13. PH ≥12 ≤2 ≤2 <4 5-9 Residual	Were correct co Were 5035 vial Air Samples: C Lot of test	hels and tags agrontainers used for ls acceptable (no cassettes / Tubes Reagent NaOH HNO3 H ₂ SO ₄ NaHSO ₄ For 608pest For CN,	ee with of the tests extra lab Intact w	custody s indica sels, no rith MS ed? No	y papers? ated? tt leaking)? S? Caniste Lot Receive No=Notify fo If +, contact P	ors Pressur d or 3day M to add	_	Samp	le ID	NO NO NO Bags In: Vol.	L	ot Adde	ØA WA	
ZnAcetate - **VOAs and 1664 Not to be tested before analysis. Otherwise, all bottles of all samples with chemical preservative are checked (not just representatives). Bottle lot numbers: 2538, 05/120 - 18mc Explain all Discrepancies/ Other Comments:	11. N 12. N 13. P pH ≥12 ≤2 ≤2 <4 5-9 Residual Chlorine	Were correct co Were 5035 vial Air Samples: C Lot of test	hels and tags agrontainers used for ls acceptable (no cassettes / Tubes Reagent NaOH HNO3 H ₂ SO ₄ NaHSO ₄ For 608pest For CN, Phenol, 625,	ee with of the tests extra lab Intact w	custody s indica els, no rith MS ed? No	y papers? ated? at leaking)? S? Caniste Lot Receive No=Notify fo If +, contact P Na ₂ S ₂ O ₃ (625,	ors Pressur d or 3day M to add 608,	_	Samp	le ID	NO NO NO Bags In: Vol.	L	ot Adde	MA NA sd	
Bottle lot numbers: 2538, 05/120 - 18mc Explain all Discrepancies/ Other Comments:	11. V 12. V 13. PH ≥12 ≤2 ≤2 <4 5-9 Residual Chlorine	Were correct co Were 5035 vial Air Samples: C Lot of test	hels and tags agrontainers used for ls acceptable (no cassettes / Tubes Reagent NaOH HNO3 H ₂ SO ₄ NaHSO ₄ For 608pest For CN, Phenol, 625, 608pest, 522	ee with of the tests extra lab Intact w	custody s indica els, no rith MS ed? No	y papers? ated? at leaking)? S? Caniste Lot Receive No=Notify fo If +, contact P Na ₂ S ₂ O ₃ (625,	ors Pressur d or 3day M to add 608,	_	Samp	le ID	NO NO NO Bags In: Vol.	L	ot Adde	N/A N/A ed	
Bottle lot numbers: 2538, 05/130 - 18mc Explain all Discrepancies/ Other Comments:	11. N 12. N 13. P pH ≥12 ≤2 ≤2 <4 5-9 Residual Chlorine	Were correct co Were 5035 vial Air Samples: C Lot of test	hbels and tags agrontainers used for ls acceptable (no Cassettes / Tubes Reagent NaOH HNO ₃ H ₂ SO ₄ NaHSO ₄ For 608pest For CN, Phenol, 625, 608pest, 522 Na ₂ S ₂ O ₃	ee with of the tests extra lab Intact with Preserve Yes	custody s indica els, no rith MS ed? No	y papers? ated? at leaking)? S? Caniste Lot Receive No=Notify fo If +, contact P Na ₂ S ₂ O ₃ (625,	ors Pressur d or 3day M to add 608,	_	Samp Adjus	le ID ted	NO NO NO Bags In Vol. Added	I.			
	11. N 12. N 13. P pH ≥12 ≤2 ≤2 <4 5-9 Residual Chlorine	Were correct co Were 5035 vial Air Samples: C Lot of test	hels and tags agrontainers used for ls acceptable (no Cassettes / Tubes Reagent NaOH HNO3 H ₂ SO ₄ NaHSO ₄ For 608pest For CN, Phenol, 625, 608pest, 522 Na ₂ S ₂ O ₃ ZnAcetate	ee with of the tests extra lab Intact with Preserve Yes	custody s indica sels, no rith MS ed? No	y papers? ated? at leaking)? S? Caniste Lot Receive No=Notify fo If +, contact P Na ₂ S ₂ O ₃ (625,	ors Pressur d or 3day M to add 608,	_	Samp Adjus **VOA Otherw	ted as and 166 ise, all bo	NO NO NO Bags In Vol. Addec	Let let tested samples	before ar	nalysis.	pH
10 hadre 15 al	11. N 12. N 13. P pH ≥12 ≤2 ≤2 <4 5-9 Residual Chlorine (-)	Were correct of Were 5035 vial Air Samples: C Lot of test paper	hels and tags agrontainers used for ls acceptable (no lassettes / Tubes Reagent NaOH HNO3 H ₂ SO ₄ NaHSO ₄ For 608pest For CN, Phenol, 625, 608pest, 522 Na ₂ S ₂ O ₃ ZnAcetate HCl	ee with of the tests extra lab Intact were Yes	custody s indica sels, no rith MS ed? No	y papers? ated? tt leaking)? G? Caniste Lot Receive No=Notify fo If +, contact P Na ₂ S ₂ O ₃ (625, CN), ascorbic	ors Pressur d or 3day M to add 608,	_	Samp Adjus **VOA Otherw	ted as and 166 ise, all bo	NO NO NO Bags In Vol. Addec	Let let tested samples	before ar	nalysis.	pH

HPROD	BULK
HTR	FLDT
SUB	HGFB
ALS	LL3541

Labels secondary reviewed by:

PC Secondary Review:

*significant air bubbles: VOA > 5-6 mm : WC > 1 in. diameter Page 13 of 115



Miscellaneous Forms

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com



REPORT QUALIFIERS AND DEFINITIONS

- U Analyte was analyzed for but not detected. The sample quantitation limit has been corrected for dilution and for percent moisture, unless otherwise noted in the case narrative.
- J Estimated value due to either being a Tentatively Identified Compound (TIC) or that the concentration is between the MRL and the MDL. Concentrations are not verified within the linear range of the calibration. For DoD: concentration >40% difference between two GC columns (pesticides/Arclors).
- B Analyte was also detected in the associated method blank at a concentration that may have contributed to the sample result.
- E Inorganics- Concentration is estimated due to the serial dilution was outside control limits.
- E Organics- Concentration has exceeded the calibration range for that specific analysis.
- D Concentration is a result of a dilution, typically a secondary analysis of the sample due to exceeding the calibration range or that a surrogate has been diluted out of the sample and cannot be assessed.
- * Indicates that a quality control parameter has exceeded laboratory limits. Under the õNotesö column of the Form I, this qualifier denotes analysis was performed out of Holding Time.
- H Analysis was performed out of hold time for tests that have an õimmediateö hold time criteria.
- # Spike was diluted out.

- + Correlation coefficient for MSA is <0.995.
- N Inorganics- Matrix spike recovery was outside laboratory limits.
- N Organics- Presumptive evidence of a compound (reported as a TIC) based on the MS library search.
- S Concentration has been determined using Method of Standard Additions (MSA).
- W Post-Digestion Spike recovery is outside control limits and the sample absorbance is <50% of the spike absorbance.
- P Concentration >40% difference between the two GC columns.
- C Confirmed by GC/MS
- Q DoD reports: indicates a pesticide/Aroclor is not confirmed (×100% Difference between two GC columns).
- X See Case Narrative for discussion.
- MRL Method Reporting Limit. Also known as:
- LOQ Limit of Quantitation (LOQ)

 The lowest concentration at which the method analyte may be reliably quantified under the method conditions.
- MDL Method Detection Limit. A statistical value derived from a study designed to provide the lowest concentration that will be detected 99% of the time. Values between the MDL and MRL are estimated (see J qualifier).
- LOD Limit of Detection. A value at or above the MDL which has been verified to be detectable.
- ND Non-Detect. Analyte was not detected at the concentration listed. Same as U qualifier.



Rochester Lab ID # for State Certifications¹

Connecticut ID # PH0556	Maine ID #NY0032	Pennsylvania ID# 68-786
Delaware Approved	New Hampshire ID # 2941	Rhode Island ID # 158
DoD ELAP #65817	New York ID # 10145	Virginia #460167
Florida ID # E87674	North Carolina #676	

¹ Analyses were performed according to our laboratory

NELAP-approved quality assurance program and any applicable state or agency requirements. The test results meet requirements of the current NELAP/TNI standards or state or agency requirements, where applicable, except as noted in the case narrative. Since not all analyte/method/matrix combinations are offered for state/NELAC accreditation, this report may contain results which are not accredited. For a specific list of accredited analytes, contact the laboratory or go to https://www.alsglobal.com/locations/americas/north-america/usa/new-york/rochester-environmental

ALS Laboratory Group

Acronyms

ASTM American Society for Testing and Materials

A2LA American Association for Laboratory Accreditation

CARB California Air Resources Board

CAS Number Chemical Abstract Service registry Number

CFC Chlorofluorocarbon CFU Colony-Forming Unit

DEC Department of Environmental Conservation

DEQ Department of Environmental Quality

DHS Department of Health Services

DOE Department of Ecology DOH Department of Health

EPA U. S. Environmental Protection Agency

ELAP Environmental Laboratory Accreditation Program

GC Gas Chromatography

GC/MS Gas Chromatography/Mass Spectrometry

LUFT Leaking Underground Fuel Tank

M Modified

MCL Maximum Contaminant Level is the highest permissible concentration of a

substance allowed in drinking water as established by the USEPA.

MDL Method Detection Limit
MPN Most Probable Number
MRL Method Reporting Limit

NA Not Applicable NC Not Calculated

NCASI National Council of the Paper Industry for Air and Stream Improvement

ND Not Detected

NIOSH National Institute for Occupational Safety and Health

PQL Practical Quantitation Limit

RCRA Resource Conservation and Recovery Act

SIM Selected Ion Monitoring

TPH Total Petroleum Hydrocarbons

tr Trace level is the concentration of an analyte that is less than the PQL but

greater than or equal to the MDL.

Analyst Summary report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term **Project:**

Monitoring

Sample Name: WG-9954-070120-SG-027

Lab Code: R2005701-001

Sample Matrix: Water

Date Collected: 07/1/20

Date Received: 07/2/20

Service Request: R2005701

Analyzed By Extracted/Digested By Analysis Method

8081B **KSERCU JMISIUREWICZ** 8082A **KSERCU BALLGEIER** 8260C **KRUEST**

8270D **KSERCU JMISIUREWICZ**

Sample Name: WG-9954-070120-SG-027 **Date Collected:** 07/1/20

Lab Code: R2005701-001.R01 Date Received: 07/2/20

Sample Matrix: Water

Analyzed By Analysis Method Extracted/Digested By

8260C **KRUEST**

8270D **KSERCU JMISIUREWICZ**

Sample Name: WG-9954-070120-SG-027 **Date Collected:** 07/1/20

Lab Code: R2005701-001.R02 Date Received: 07/2/20

Sample Matrix: Water

Analyzed By Analysis Method Extracted/Digested By

8270D **KSERCU JMISIUREWICZ**

Sample Name: WG-9954-070120-SG-026 Date Collected: 07/1/20

Lab Code: R2005701-002 Date Received: 07/2/20

Sample Matrix: Water

Analyzed By Analysis Method Extracted/Digested By

8081B **KSERCU JMISIUREWICZ** 8082A **KSERCU BALLGEIER**

8260C **KRUEST**

8270D **KSERCU JMISIUREWICZ**

Analyst Summary report

Service Request: R2005701

Date Collected: 07/1/20

Date Received: 07/2/20

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Monitoring

Sample Name: WG-9954-070120-SG-025

Lab Code: R2005701-003

Sample Matrix: Water

Analysis Method Extracted/Digested By Analyzed By

8081B KSERCU JMISIUREWICZ 8082A KSERCU BALLGEIER 8260C KRUEST

8270D KSERCU JMISIUREWICZ

Sample Name: WG-9954-070120-SG-024 **Date Collected:** 07/1/20

Lab Code: R2005701-004 **Date Received:** 07/2/20

Sample Matrix: Water

Analysis Method Extracted/Digested By Analyzed By

8081BKSERCUJMISIUREWICZ8082AKSERCUBALLGEIER8260CKRUEST

8270D KSERCU JMISIUREWICZ

Sample Name: TB-9954-070120-SG-005 **Date Collected:** 07/1/20

Lab Code: R2005701-005 **Date Received:** 07/2/20

Sample Matrix: Water

Analysis Method Extracted/Digested By Analyzed By

8260C KRUEST

Sample Name: WG-9954-070120-SG-028 **Date Collected:** 07/1/20

Lab Code: R2005701-006 **Date Received:** 07/2/20

Sample Matrix: Water

Analysis Method Extracted/Digested By Analyzed By

8081B KSERCU BALLGEIER 8082A KSERCU BALLGEIER

8260C KRUEST

8270D KSERCU JMISIUREWICZ

Printed 7/30/2020 4:05:51 PM Superset Reference:20-0000555360 rev 00

Analyst Summary report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Monitoring

Sample Name: WG-9954-070120-SG-028

Lab Code: R2005701-006.R01

Sample Matrix: Water

Date Collected: 07/1/20

Date Received: 07/2/20

Service Request: R2005701

Analysis Method Extracted/Digested By Analyzed By

8081B KSERCU JMISIUREWICZ

8260C KRUEST

8270D KSERCU JMISIUREWICZ

Sample Name: WG-9954-070120-SG-028 **Date Collected:** 07/1/20

Lab Code: R2005701-006.R02 **Date Received:** 07/2/20

Sample Matrix: Water

Analysis Method Extracted/Digested By Analyzed By

8270D KSERCU JMISIUREWICZ



INORGANIC PREPARATION METHODS

The preparation methods associated with this report are found in these tables unless discussed in the case narrative.

Water/Liquid Matrix

Analytical Method	Preparation Method
200.7	200.2
200.8	200.2
6010C	3005A/3010A
6020A	ILM05.3
9034 Sulfide Acid Soluble	9030B
SM 4500-CN-E Residual Cyanide	SM 4500-CN-G
SM 4500-CN-E WAD Cyanide	SM 4500-CN-I

Solid/Soil/Non-Aqueous Matrix

Analytical Method	Preparation
	Method
6010C	3050B
6020A	3050B
6010C TCLP (1311)	3005A/3010A
extract	
6010 SPLP (1312) extract	3005A/3010A
7199	3060A
300.0 Anions/ 350.1/	DI extraction
353.2/ SM 2320B/ SM	
5210B/ 9056A Anions	
For analytical methods not listed,	
method is the same as the analytic	cal method
I ICICICILC.	



Sample Results

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com



Volatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

Analytical Report

Client: Service Request: R2005701 GHD (Formerly Conestoga-Rovers & Associates) **Date Collected:** 07/01/20 11:50

Love Canal:292-402-D02-3100/9954 Annual Long Term **Project:**

Monitoring

Date Received: 07/02/20 10:50 **Sample Matrix:** Water

Sample Name: WG-9954-070120-SG-027 Units: ug/L Lab Code: R2005701-001 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/13/20 21:35	
1,1,2,2-Tetrachloroethane	6.1	5.0	0.20	1	07/13/20 21:35	
1,1,2-Trichloroethane	12	5.0	0.20	1	07/13/20 21:35	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/13/20 21:35	
1,1-Dichloroethene (1,1-DCE)	0.63 J	5.0	0.20	1	07/13/20 21:35	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/13/20 21:35	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/13/20 21:35	
2-Butanone (MEK)	12	10	0.78	1	07/13/20 21:35	
2-Hexanone	10 U	10	0.20	1	07/13/20 21:35	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/13/20 21:35	
Acetone	79	10	5.0	1	07/13/20 21:35	
Benzene	1300 E	5.0	0.20	1	07/13/20 21:35	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/13/20 21:35	
Bromoform	5.0 U	5.0	0.25	1	07/13/20 21:35	
Bromomethane	5.0 U	5.0	0.70	1	07/13/20 21:35	
Carbon Disulfide	10 U	10	0.42	1	07/13/20 21:35	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/13/20 21:35	
Chlorobenzene	770 E	5.0	0.20	1	07/13/20 21:35	
Chloroethane	5.0 U	5.0	0.23	1	07/13/20 21:35	
Chloroform	140	5.0	0.24	1	07/13/20 21:35	
Chloromethane	5.0 U	5.0	0.28	1	07/13/20 21:35	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/13/20 21:35	
Dichloromethane	6.4	5.0	0.65	1	07/13/20 21:35	
Ethylbenzene	12	5.0	0.20	1	07/13/20 21:35	
Styrene	5.0 U	5.0	0.20	1	07/13/20 21:35	
Tetrachloroethene (PCE)	20	5.0	0.21	1	07/13/20 21:35	
Toluene	1700 E	5.0	0.20	1	07/13/20 21:35	
Trichloroethene (TCE)	120	5.0	0.20	1	07/13/20 21:35	
Vinyl Acetate	10 U	10	1.1	1	07/13/20 21:35	
Vinyl Chloride	10	5.0	0.20	1	07/13/20 21:35	
Xylenes, Total	58	5.0	0.23	1	07/13/20 21:35	
cis-1,2-Dichloroethene	26	5.0	0.23	1	07/13/20 21:35	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/13/20 21:35	
trans-1,2-Dichloroethene	29	5.0	0.20	1	07/13/20 21:35	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/13/20 21:35	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	114	85 - 122	07/13/20 21:35	
Dibromofluoromethane	89	89 - 119	07/13/20 21:35	
Toluene-d8	99	87 - 121	07/13/20 21:35	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 11:50

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-027
 Units: ug/L

 Lab Code:
 R2005701-001
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Tentatively Identified Compounds

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected

Analytical Report

Client: Service Request: R2005701 GHD (Formerly Conestoga-Rovers & Associates) Love Canal:292-402-D02-3100/9954 Annual Long Term **Project:**

Monitoring

Date Collected: 07/01/20 11:50

Sample Matrix: Water

Date Received: 07/02/20 10:50

Sample Name: WG-9954-070120-SG-027 Lab Code: R2005701-001

Units: ug/L Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	1000 U	1000	40	200	07/15/20 17:02	
1,1,2,2-Tetrachloroethane	1000 U	1000	40	200	07/15/20 17:02	
1,1,2-Trichloroethane	1000 U	1000	40	200	07/15/20 17:02	
1,1-Dichloroethane (1,1-DCA)	1000 U	1000	40	200	07/15/20 17:02	
1,1-Dichloroethene (1,1-DCE)	1000 U	1000	40	200	07/15/20 17:02	
1,2-Dichloroethane	1000 U	1000	40	200	07/15/20 17:02	
1,2-Dichloropropane	1000 U	1000	40	200	07/15/20 17:02	
2-Butanone (MEK)	2000 U	2000	160	200	07/15/20 17:02	
2-Hexanone	2000 U	2000	40	200	07/15/20 17:02	
4-Methyl-2-pentanone	2000 U	2000	40	200	07/15/20 17:02	
Acetone	2000 U	2000	1000	200	07/15/20 17:02	
Benzene	6500 D	1000	40	200	07/15/20 17:02	
Bromodichloromethane	60 DJ	1000	40	200	07/15/20 17:02	
Bromoform	1000 U	1000	50	200	07/15/20 17:02	
Bromomethane	1000 U	1000	140	200	07/15/20 17:02	
Carbon Disulfide	2000 U	2000	84	200	07/15/20 17:02	
Carbon Tetrachloride	1000 U	1000	68	200	07/15/20 17:02	
Chlorobenzene	2500 D	1000	40	200	07/15/20 17:02	
Chloroethane	1000 U	1000	46	200	07/15/20 17:02	
Chloroform	400 DJ	1000	48	200	07/15/20 17:02	
Chloromethane	1000 U	1000	56	200	07/15/20 17:02	
Dibromochloromethane	1000 U	1000	40	200	07/15/20 17:02	
Dichloromethane	1000 U	1000	130	200	07/15/20 17:02	
Ethylbenzene	1000 U	1000	40	200	07/15/20 17:02	
Styrene	1000 U	1000	40	200	07/15/20 17:02	
Tetrachloroethene (PCE)	1000 U	1000	42	200	07/15/20 17:02	
Toluene	22000 D	1000	40	200	07/15/20 17:02	
Trichloroethene (TCE)	100 DJ	1000	40	200	07/15/20 17:02	
Vinyl Acetate	2000 U	2000	220	200	07/15/20 17:02	
Vinyl Chloride	1000 U	1000	40	200	07/15/20 17:02	
Xylenes, Total	1000 U	1000	46	200	07/15/20 17:02	
cis-1,2-Dichloroethene	1000 U	1000	46	200	07/15/20 17:02	
cis-1,3-Dichloropropene	1000 U	1000	40	200	07/15/20 17:02	
trans-1,2-Dichloroethene	1000 U	1000	40	200	07/15/20 17:02	
trans-1,3-Dichloropropene	1000 U	1000	46	200	07/15/20 17:02	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	96	85 - 122	07/15/20 17:02	_
Dibromofluoromethane	96	89 - 119	07/15/20 17:02	
Toluene-d8	100	87 - 121	07/15/20 17:02	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 11:50

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-027
 Units: ug/L

 Lab Code:
 R2005701-001
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
000095-49-8	Benzene, 1-chloro-2-methyl-	11.16	16362.0	JN	
000108-41-8	Benzene, 1-chloro-3-methyl-	11.25	8134.0	JN	
032768-54-0	Benzene, 1,2-dichloro-3-methyl-	12.89	1534.0	JN	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 10:45

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-026
 Units: ug/L

 Lab Code:
 R2005701-002
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/14/20 17:31	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/14/20 17:31	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/14/20 17:31	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/14/20 17:31	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/14/20 17:31	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/14/20 17:31	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/14/20 17:31	
2-Butanone (MEK)	10 U	10	0.78	1	07/14/20 17:31	
2-Hexanone	10 U	10	0.20	1	07/14/20 17:31	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/14/20 17:31	
Acetone	10 U	10	5.0	1	07/14/20 17:31	
Benzene	5.0 U	5.0	0.20	1	07/14/20 17:31	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/14/20 17:31	
Bromoform	5.0 U	5.0	0.25	1	07/14/20 17:31	
Bromomethane	5.0 U	5.0	0.70	1	07/14/20 17:31	
Carbon Disulfide	3.3 J	10	0.42	1	07/14/20 17:31	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/14/20 17:31	
Chlorobenzene	0.89 J	5.0	0.20	1	07/14/20 17:31	
Chloroethane	5.0 U	5.0	0.23	1	07/14/20 17:31	
Chloroform	5.0 U	5.0	0.24	1	07/14/20 17:31	
Chloromethane	5.0 U	5.0	0.28	1	07/14/20 17:31	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/14/20 17:31	
Dichloromethane	5.0 U	5.0	0.65	1	07/14/20 17:31	
Ethylbenzene	5.0 U	5.0	0.20	1	07/14/20 17:31	
Styrene	5.0 U	5.0	0.20	1	07/14/20 17:31	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/14/20 17:31	
Toluene	0.26 BJ	5.0	0.20	1	07/14/20 17:31	
Trichloroethene (TCE)	6.5	5.0	0.20	1	07/14/20 17:31	
Vinyl Acetate	10 U	10	1.1	1	07/14/20 17:31	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/14/20 17:31	
Xylenes, Total	5.0 U	5.0	0.23	1	07/14/20 17:31	
cis-1,2-Dichloroethene	3.3 J	5.0	0.23	1	07/14/20 17:31	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/14/20 17:31	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/14/20 17:31	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/14/20 17:31	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 10:45

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-026
 Units: ug/L

 Lab Code:
 R2005701-002
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	95	85 - 122	07/14/20 17:31	
Dibromofluoromethane	105	89 - 119	07/14/20 17:31	
Toluene-d8	102	87 - 121	07/14/20 17:31	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 10:45

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-026
 Units: ug/L

 Lab Code:
 R2005701-002
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
007446-09-5	Sulfur dioxide	1.29	46.1	JN	
	unknown	1.63	20.0	J	
000098-15-7	Benzene, 1-chloro-3-(trifluorometh	10.66	6.1	JN	
000095-49-8	Benzene, 1-chloro-2-methyl-	11.16	8.1	JN	
000120-82-1	Benzene, 1,2,4-trichloro-	13.43	12.7	JN	
000066-25-1	Hexanal	9.25	6.7	JN	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 09:55

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-025
 Units: ug/L

 Lab Code:
 R2005701-003
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/14/20 17:53	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/14/20 17:53	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/14/20 17:53	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/14/20 17:53	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/14/20 17:53	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/14/20 17:53	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/14/20 17:53	
2-Butanone (MEK)	10 U	10	0.78	1	07/14/20 17:53	
2-Hexanone	10 U	10	0.20	1	07/14/20 17:53	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/14/20 17:53	
Acetone	10 U	10	5.0	1	07/14/20 17:53	
Benzene	5.0 U	5.0	0.20	1	07/14/20 17:53	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/14/20 17:53	
Bromoform	5.0 U	5.0	0.25	1	07/14/20 17:53	
Bromomethane	5.0 U	5.0	0.70	1	07/14/20 17:53	
Carbon Disulfide	10 U	10	0.42	1	07/14/20 17:53	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/14/20 17:53	
Chlorobenzene	5.0 U	5.0	0.20	1	07/14/20 17:53	
Chloroethane	5.0 U	5.0	0.23	1	07/14/20 17:53	
Chloroform	5.0 U	5.0	0.24	1	07/14/20 17:53	
Chloromethane	5.0 U	5.0	0.28	1	07/14/20 17:53	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/14/20 17:53	
Dichloromethane	5.0 U	5.0	0.65	1	07/14/20 17:53	
Ethylbenzene	5.0 U	5.0	0.20	1	07/14/20 17:53	
Styrene	5.0 U	5.0	0.20	1	07/14/20 17:53	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/14/20 17:53	
Toluene	0.27 BJ	5.0	0.20	1	07/14/20 17:53	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/14/20 17:53	
Vinyl Acetate	10 U	10	1.1	1	07/14/20 17:53	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/14/20 17:53	
Xylenes, Total	5.0 U	5.0	0.23	1	07/14/20 17:53	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/14/20 17:53	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/14/20 17:53	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/14/20 17:53	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/14/20 17:53	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 09:55

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-025
 Units: ug/L

 Lab Code:
 R2005701-003
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	98	85 - 122	07/14/20 17:53	
Dibromofluoromethane	100	89 - 119	07/14/20 17:53	
Toluene-d8	103	87 - 121	07/14/20 17:53	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 09:55

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-025
 Units: ug/L

 Lab Code:
 R2005701-003
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

			Result	
CAS#	Compound Identification	RT	ug/L	Q
007446-09-5	Sulfur dioxide	1.29	56.0	JN

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 08:45

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-024
 Units: ug/L

 Lab Code:
 R2005701-004
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/14/20 18:15	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/14/20 18:15	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/14/20 18:15	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/14/20 18:15	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/14/20 18:15	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/14/20 18:15	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/14/20 18:15	
2-Butanone (MEK)	10 U	10	0.78	1	07/14/20 18:15	
2-Hexanone	10 U	10	0.20	1	07/14/20 18:15	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/14/20 18:15	
Acetone	10 U	10	5.0	1	07/14/20 18:15	
Benzene	5.0 U	5.0	0.20	1	07/14/20 18:15	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/14/20 18:15	
Bromoform	5.0 U	5.0	0.25	1	07/14/20 18:15	
Bromomethane	5.0 U	5.0	0.70	1	07/14/20 18:15	
Carbon Disulfide	10 U	10	0.42	1	07/14/20 18:15	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/14/20 18:15	
Chlorobenzene	5.0 U	5.0	0.20	1	07/14/20 18:15	
Chloroethane	5.0 U	5.0	0.23	1	07/14/20 18:15	
Chloroform	5.0 U	5.0	0.24	1	07/14/20 18:15	
Chloromethane	5.0 U	5.0	0.28	1	07/14/20 18:15	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/14/20 18:15	
Dichloromethane	5.0 U	5.0	0.65	1	07/14/20 18:15	
Ethylbenzene	5.0 U	5.0	0.20	1	07/14/20 18:15	
Styrene	5.0 U	5.0	0.20	1	07/14/20 18:15	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/14/20 18:15	
Toluene	0.26 BJ	5.0	0.20	1	07/14/20 18:15	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/14/20 18:15	
Vinyl Acetate	10 U	10	1.1	1	07/14/20 18:15	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/14/20 18:15	
Xylenes, Total	5.0 U	5.0	0.23	1	07/14/20 18:15	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/14/20 18:15	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/14/20 18:15	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/14/20 18:15	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/14/20 18:15	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 08:45

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-024
 Units: ug/L

 Lab Code:
 R2005701-004
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	96	85 - 122	07/14/20 18:15	
Dibromofluoromethane	100	89 - 119	07/14/20 18:15	
Toluene-d8	102	87 - 121	07/14/20 18:15	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 08:45

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-024
 Units: ug/L

 Lab Code:
 R2005701-004
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Tentatively Identified Compounds

 CAS#
 Compound Identification
 RT
 ug/L
 Q

 007446-09-5
 Sulfur dioxide
 1.29
 66.0
 JN

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 00:00

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 TB-9954-070120-SG-005
 Units: ug/L

 Lab Code:
 R2005701-005
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/13/20 21:13	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/13/20 21:13	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/13/20 21:13	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/13/20 21:13	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/13/20 21:13	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/13/20 21:13	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/13/20 21:13	
2-Butanone (MEK)	10 U	10	0.78	1	07/13/20 21:13	
2-Hexanone	10 U	10	0.20	1	07/13/20 21:13	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/13/20 21:13	
Acetone	10 U	10	5.0	1	07/13/20 21:13	
Benzene	5.0 U	5.0	0.20	1	07/13/20 21:13	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/13/20 21:13	
Bromoform	5.0 U	5.0	0.25	1	07/13/20 21:13	
Bromomethane	5.0 U	5.0	0.70	1	07/13/20 21:13	
Carbon Disulfide	10 U	10	0.42	1	07/13/20 21:13	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/13/20 21:13	
Chlorobenzene	5.0 U	5.0	0.20	1	07/13/20 21:13	
Chloroethane	5.0 U	5.0	0.23	1	07/13/20 21:13	
Chloroform	5.0 U	5.0	0.24	1	07/13/20 21:13	
Chloromethane	5.0 U	5.0	0.28	1	07/13/20 21:13	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/13/20 21:13	
Dichloromethane	5.0 U	5.0	0.65	1	07/13/20 21:13	
Ethylbenzene	5.0 U	5.0	0.20	1	07/13/20 21:13	
Styrene	5.0 U	5.0	0.20	1	07/13/20 21:13	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/13/20 21:13	
Toluene	5.0 U	5.0	0.20	1	07/13/20 21:13	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/13/20 21:13	
Vinyl Acetate	10 U	10	1.1	1	07/13/20 21:13	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/13/20 21:13	
Xylenes, Total	5.0 U	5.0	0.23	1	07/13/20 21:13	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/13/20 21:13	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/13/20 21:13	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/13/20 21:13	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/13/20 21:13	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 00:00

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 TB-9954-070120-SG-005
 Units: ug/L

 Lab Code:
 R2005701-005
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	91	85 - 122	07/13/20 21:13	
Dibromofluoromethane	98	89 - 119	07/13/20 21:13	
Toluene-d8	100	87 - 121	07/13/20 21:13	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 00:00

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 TB-9954-070120-SG-005
 Units: ug/L

 Lab Code:
 R2005701-005
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Tentatively Identified Compounds

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected

Analytical Report

Client: Service Request: R2005701 GHD (Formerly Conestoga-Rovers & Associates) Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 11:50 **Project:**

Monitoring

Sample Name: WG-9954-070120-SG-028 Units: ug/L

Lab Code: R2005701-006 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Water

Sample Matrix:

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/13/20 23:02	
1,1,2,2-Tetrachloroethane	7.9	5.0	0.20	1	07/13/20 23:02	
1,1,2-Trichloroethane	14	5.0	0.20	1	07/13/20 23:02	
1,1-Dichloroethane (1,1-DCA)	0.27 J	5.0	0.20	1	07/13/20 23:02	
1,1-Dichloroethene (1,1-DCE)	0.78 J	5.0	0.20	1	07/13/20 23:02	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/13/20 23:02	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/13/20 23:02	
2-Butanone (MEK)	6.8 J	10	0.78	1	07/13/20 23:02	
2-Hexanone	2.1 J	10	0.20	1	07/13/20 23:02	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/13/20 23:02	
Acetone	71	10	5.0	1	07/13/20 23:02	
Benzene	1500 E	5.0	0.20	1	07/13/20 23:02	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/13/20 23:02	
Bromoform	5.0 U	5.0	0.25	1	07/13/20 23:02	
Bromomethane	5.0 U	5.0	0.70	1	07/13/20 23:02	
Carbon Disulfide	10 U	10	0.42	1	07/13/20 23:02	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/13/20 23:02	
Chlorobenzene	850 E	5.0	0.20	1	07/13/20 23:02	
Chloroethane	5.0 U	5.0	0.23	1	07/13/20 23:02	
Chloroform	140	5.0	0.24	1	07/13/20 23:02	
Chloromethane	5.0 U	5.0	0.28	1	07/13/20 23:02	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/13/20 23:02	
Dichloromethane	7.2	5.0	0.65	1	07/13/20 23:02	
Ethylbenzene	12	5.0	0.20	1	07/13/20 23:02	
Styrene	5.0 U	5.0	0.20	1	07/13/20 23:02	
Tetrachloroethene (PCE)	23	5.0	0.21	1	07/13/20 23:02	
Toluene	2000 E	5.0	0.20	1	07/13/20 23:02	
Trichloroethene (TCE)	130	5.0	0.20	1	07/13/20 23:02	
Vinyl Acetate	10 U	10	1.1	1	07/13/20 23:02	
Vinyl Chloride	11	5.0	0.20	1	07/13/20 23:02	
Xylenes, Total	58	5.0	0.23	1	07/13/20 23:02	
cis-1,2-Dichloroethene	28	5.0	0.23	1	07/13/20 23:02	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/13/20 23:02	
trans-1,2-Dichloroethene	35	5.0	0.20	1	07/13/20 23:02	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/13/20 23:02	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	113	85 - 122	07/13/20 23:02	
Dibromofluoromethane	99	89 - 119	07/13/20 23:02	
Toluene-d8	108	87 - 121	07/13/20 23:02	

Date Received: 07/02/20 10:50

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 11:50

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-028
 Units: ug/L

 Lab Code:
 R2005701-006
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Tentatively Identified Compounds

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected

Analytical Report

Client: Service Request: R2005701 GHD (Formerly Conestoga-Rovers & Associates) Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 11:50 **Project:**

Monitoring

Sample Matrix: Water

Date Received: 07/02/20 10:50

Sample Name: WG-9954-070120-SG-028 Units: ug/L Lab Code: R2005701-006 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	1000 U	1000	40	200	07/15/20 14:51	_
1,1,2,2-Tetrachloroethane	1000 U	1000	40	200	07/15/20 14:51	
1,1,2-Trichloroethane	1000 U	1000	40	200	07/15/20 14:51	
1,1-Dichloroethane (1,1-DCA)	1000 U	1000	40	200	07/15/20 14:51	
1,1-Dichloroethene (1,1-DCE)	1000 U	1000	40	200	07/15/20 14:51	
1,2-Dichloroethane	1000 U	1000	40	200	07/15/20 14:51	
1,2-Dichloropropane	1000 U	1000	40	200	07/15/20 14:51	
2-Butanone (MEK)	2000 U	2000	160	200	07/15/20 14:51	
2-Hexanone	2000 U	2000	40	200	07/15/20 14:51	
4-Methyl-2-pentanone	2000 U	2000	40	200	07/15/20 14:51	
Acetone	2000 U	2000	1000	200	07/15/20 14:51	
Benzene	6300 D	1000	40	200	07/15/20 14:51	
Bromodichloromethane	1000 U	1000	40	200	07/15/20 14:51	
Bromoform	1000 U	1000	50	200	07/15/20 14:51	
Bromomethane	1000 U	1000	140	200	07/15/20 14:51	
Carbon Disulfide	2000 U	2000	84	200	07/15/20 14:51	
Carbon Tetrachloride	1000 U	1000	68	200	07/15/20 14:51	
Chlorobenzene	2400 D	1000	40	200	07/15/20 14:51	
Chloroethane	1000 U	1000	46	200	07/15/20 14:51	
Chloroform	270 DJ	1000	48	200	07/15/20 14:51	
Chloromethane	1000 U	1000	56	200	07/15/20 14:51	
Dibromochloromethane	1000 U	1000	40	200	07/15/20 14:51	
Dichloromethane	1000 U	1000	130	200	07/15/20 14:51	
Ethylbenzene	1000 U	1000	40	200	07/15/20 14:51	
Styrene	1000 U	1000	40	200	07/15/20 14:51	
Tetrachloroethene (PCE)	1000 U	1000	42	200	07/15/20 14:51	
Toluene	21000 D	1000	40	200	07/15/20 14:51	
Trichloroethene (TCE)	120 DJ	1000	40	200	07/15/20 14:51	
Vinyl Acetate	2000 U	2000	220	200	07/15/20 14:51	
Vinyl Chloride	1000 U	1000	40	200	07/15/20 14:51	
Xylenes, Total	1000 U	1000	46	200	07/15/20 14:51	
cis-1,2-Dichloroethene	1000 U	1000	46	200	07/15/20 14:51	
cis-1,3-Dichloropropene	1000 U	1000	40	200	07/15/20 14:51	
trans-1,2-Dichloroethene	51 DJ	1000	40	200	07/15/20 14:51	
trans-1,3-Dichloropropene	1000 U	1000	46	200	07/15/20 14:51	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	96	85 - 122	07/15/20 14:51	_
Dibromofluoromethane	96	89 - 119	07/15/20 14:51	
Toluene-d8	101	87 - 121	07/15/20 14:51	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 11:50

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-028
 Units: ug/L

 Lab Code:
 R2005701-006
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
000095-49-8	Benzene, 1-chloro-2-methyl-	11.16	16222.0	JN	
000108-41-8	Benzene, 1-chloro-3-methyl-	11.25	7998.0	JN	
032768-54-0	Benzene, 1,2-dichloro-3-methyl-	12.89	1562.0	JN	



Semivolatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701 Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 11:50 **Project:**

Monitoring

Date Received: 07/02/20 10:50

Sample Matrix: Water

Units: ug/L

Sample Name: WG-9954-070120-SG-027 Basis: NA Lab Code: R2005701-001

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	93	9.1	1.1	1	07/09/20 12:04	7/7/20	
1,2-Dichlorobenzene	25	9.1	1.1	1	07/09/20 12:04	7/7/20	
1,3-Dichlorobenzene	3.6 J	9.1	0.92	1	07/09/20 12:04	7/7/20	
1,4-Dichlorobenzene	73	9.1	1.1	1	07/09/20 12:04	7/7/20	
2,4,5-Trichlorophenol	37	9.1	0.99	1	07/09/20 12:04	7/7/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/09/20 12:04	7/7/20	
2,4-Dichlorophenol	550 E	9.1	1.2	1	07/09/20 12:04	7/7/20	
2,4-Dimethylphenol	8.9 J	9.1	1.3	1	07/09/20 12:04	7/7/20	
2,4-Dinitrophenol	45 U	45	19	1	07/09/20 12:04	7/7/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/09/20 12:04	7/7/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/09/20 12:04	7/7/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/09/20 12:04	7/7/20	
2-Chlorophenol	27	9.1	0.97	1	07/09/20 12:04	7/7/20	
2-Methylnaphthalene	1.4 J	9.1	1.2	1	07/09/20 12:04	7/7/20	
2-Methylphenol	32	9.1	0.91	1	07/09/20 12:04	7/7/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/09/20 12:04	7/7/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/09/20 12:04	7/7/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/09/20 12:04	7/7/20	
3- and 4-Methylphenol Coelution	71	9.1	1.1	1	07/09/20 12:04	7/7/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/09/20 12:04	7/7/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/09/20 12:04	7/7/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/09/20 12:04	7/7/20	
4-Chloro-3-methylphenol	46	9.1	0.98	1	07/09/20 12:04	7/7/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/09/20 12:04	7/7/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/09/20 12:04	7/7/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/09/20 12:04	7/7/20	
4-Nitrophenol	45 U	45	5.8	1	07/09/20 12:04	7/7/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/09/20 12:04	7/7/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/09/20 12:04	7/7/20	
Anthracene	9.1 U	9.1	1.2	1	07/09/20 12:04	7/7/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/09/20 12:04	7/7/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/09/20 12:04	7/7/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/09/20 12:04	7/7/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/09/20 12:04	7/7/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/09/20 12:04	7/7/20	
Benzoic Acid	4500 E	91	33	1	07/09/20 12:04	7/7/20	
Benzyl Alcohol	290 290	9.1	1.5	1	07/09/20 12:04	7/7/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/09/20 12:04	7/7/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/09/20 12:04	7/7/20	
Bis(2-chloroethyl) Ether	20	9.1	1.3	1	07/09/20 12:04	7/7/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	07/09/20 12:04	7/7/20	
	9.1 U 9.1 U	9.1 9.1	1.3	1	07/09/20 12:04	7/7/20	
Butyl Benzyl Phthalate			1.3 1.1				
Chrysene	9.1 U	9.1	1.1	1	07/09/20 12:04	7/7/20	

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Superset Reference:20-0000555360 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 11:50

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-027
 Units: ug/L

 Lab Code:
 R2005701-001
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed Da	ate Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/09/20 12:04	7/7/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/09/20 12:04	7/7/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/09/20 12:04	7/7/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/09/20 12:04	7/7/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/09/20 12:04	7/7/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/09/20 12:04	7/7/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/09/20 12:04	7/7/20	
Fluorene	9.1 U	9.1	1.2	1	07/09/20 12:04	7/7/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/09/20 12:04	7/7/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/09/20 12:04	7/7/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/09/20 12:04	7/7/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/09/20 12:04	7/7/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/09/20 12:04	7/7/20	
Isophorone	9.1 U	9.1	1.3	1	07/09/20 12:04	7/7/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/09/20 12:04	7/7/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/09/20 12:04	7/7/20	
Naphthalene	9.1 U	9.1	1.1	1	07/09/20 12:04	7/7/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/09/20 12:04	7/7/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/09/20 12:04	7/7/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/09/20 12:04	7/7/20	
Phenol	40	9.1	0.91	1	07/09/20 12:04	7/7/20	
Pyrene	9.1 U	9.1	1.3	1	07/09/20 12:04	7/7/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	80	35 - 141	07/09/20 12:04	
2-Fluorobiphenyl	59	31 - 118	07/09/20 12:04	
2-Fluorophenol	29	10 - 105	07/09/20 12:04	
Nitrobenzene-d5	86	31 - 110	07/09/20 12:04	
Phenol-d6	24	10 - 107	07/09/20 12:04	
p-Terphenyl-d14	51	10 - 165	07/09/20 12:04	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
000611-95-0	Benzoic acid, 4-benzoyl-	10.20	780	JN	
000579-18-0	3-Benzoylbenzoic acid	10.22	290	JN	
	unknown	10.27	77	J	
000123-91-1	1,4-Dioxane	2.19	140	JN	
000108-88-3	Toluene	2.69	1800	JN	
000107-92-6	Butanoic acid	2.91	110	JN	
000108-90-7	Benzene, chloro-	3.32	350	JN	

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Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701 **Date Collected:** 07/01/20 11:50

Love Canal:292-402-D02-3100/9954 Annual Long Term **Project:**

Monitoring

Date Received: 07/02/20 10:50 **Sample Matrix:** Water

Sample Name: WG-9954-070120-SG-027 Units: ug/L Lab Code: R2005701-001 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
000108-93-0	Cyclohexanol	3.63	120	JN	
	unknown	3.70	42	J	
000591-23-1	Cyclohexanol, 3-methyl-	4.09	38	JN	
000095-49-8	Benzene, 1-chloro-2-methyl-	4.14	1800	JN	
000106-43-4	Benzene, 1-chloro-4-methyl-	4.18	1000	JN	
001195-79-5	Bicyclo[2.2.1]heptan-2-one, 1,3,3-	5.14	32	JN	
000106-48-9	Parachlorophenol	5.75	150	JN	
000074-11-3	Benzoic acid, 4-chloro-	6.68	1500	JN	
000535-80-8	Benzoic acid, 3-chloro-	6.90	9900	JN	
	unknown	6.93	360	J	
000591-35-5	Phenol, 3,5-dichloro-	7.00	370	JN	
	unknown	7.12	73	J	
000050-45-3	Benzoic acid, 2,3-dichloro-	7.60	28	JN	
000134-84-9	Methanone, (4-methylphenyl)phenyl-	8.73	31	JN	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701 **Date Collected:** 07/01/20 11:50

Love Canal:292-402-D02-3100/9954 Annual Long Term **Project:**

Monitoring

Date Received: 07/02/20 10:50 **Sample Matrix:** Water

Sample Name: Units: ug/L WG-9954-070120-SG-027 Basis: NA Lab Code: R2005701-001

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	59 J	230	26	25	07/13/20 14:46	7/7/20	
1,2-Dichlorobenzene	230 U	230	27	25	07/13/20 14:46	7/7/20	
1,3-Dichlorobenzene	230 U	230	23	25	07/13/20 14:46	7/7/20	
1,4-Dichlorobenzene	72 J	230	26	25	07/13/20 14:46	7/7/20	
2,4,5-Trichlorophenol	50 J	230	25	25	07/13/20 14:46	7/7/20	
2,4,6-Trichlorophenol	230 U	230	31	25	07/13/20 14:46	7/7/20	
2,4-Dichlorophenol	330 D	230	29	25	07/13/20 14:46	7/7/20	
2,4-Dimethylphenol	230 U	230	32	25	07/13/20 14:46	7/7/20	
2,4-Dinitrophenol	1100 U	1100	460	25	07/13/20 14:46	7/7/20	
2,4-Dinitrotoluene	230 U	230	54	25	07/13/20 14:46	7/7/20	
2,6-Dinitrotoluene	230 U	230	30	25	07/13/20 14:46	7/7/20	
2-Chloronaphthalene	230 U	230	31	25	07/13/20 14:46	7/7/20	
2-Chlorophenol	230 U	230	25	25	07/13/20 14:46	7/7/20	
2-Methylnaphthalene	230 U	230	30	25	07/13/20 14:46	7/7/20	
2-Methylphenol	32 J	230	23	25	07/13/20 14:46	7/7/20	
2-Nitroaniline	230 U	230	31	25	07/13/20 14:46	7/7/20	
2-Nitrophenol	230 U	230	34	25	07/13/20 14:46	7/7/20	
3,3'-Dichlorobenzidine	230 U	230	28	25	07/13/20 14:46	7/7/20	
3- and 4-Methylphenol Coelution	60 J	230	27	25	07/13/20 14:46	7/7/20	
3-Nitroaniline	230 U	230	56	25	07/13/20 14:46	7/7/20	
4,6-Dinitro-2-methylphenol	1100 U	1100	440	25	07/13/20 14:46	7/7/20	
4-Bromophenyl Phenyl Ether	230 U	230	38	25	07/13/20 14:46	7/7/20	
4-Chloro-3-methylphenol	31 J	230	25	25	07/13/20 14:46	7/7/20	
4-Chloroaniline	230 U	230	23	25	07/13/20 14:46	7/7/20	
4-Chlorophenyl Phenyl Ether	230 U	230	34	25	07/13/20 14:46	7/7/20	
4-Nitroaniline	230 U	230	62	25	07/13/20 14:46	7/7/20	
4-Nitrophenol	1100 U	1100	150	25	07/13/20 14:46	7/7/20	
Acenaphthene	230 U	230	31	25	07/13/20 14:46	7/7/20	
Acenaphthylene	230 U	230	31	25	07/13/20 14:46	7/7/20	
Anthracene	230 U	230	29	25	07/13/20 14:46	7/7/20	
Benz(a)anthracene	230 U	230	36	25	07/13/20 14:46	7/7/20	
Benzo(a)pyrene	230 U	230	26	25	07/13/20 14:46	7/7/20	
Benzo(b)fluoranthene	230 U	230	26	25	07/13/20 14:46	7/7/20	
Benzo(g,h,i)perylene	230 U	230	23	25	07/13/20 14:46	7/7/20	
Benzo(k)fluoranthene	230 U	230	28	25	07/13/20 14:46	7/7/20	
Benzoic Acid	9200 E	2300	820	25	07/13/20 14:46	7/7/20	
Benzyl Alcohol	240	230	36	25	07/13/20 14:46	7/7/20	
2,2'-Oxybis(1-chloropropane)	230 U	230	32	25	07/13/20 14:46	7/7/20	
Bis(2-chloroethoxy)methane	230 U	230	44	25	07/13/20 14:46	7/7/20	
Bis(2-chloroethyl) Ether	230 U	230	28	25	07/13/20 14:46	7/7/20	
Bis(2-ethylhexyl) Phthalate	230 U	230	23	25	07/13/20 14:46	7/7/20	
Butyl Benzyl Phthalate	230 U	230	23 32	25 25	07/13/20 14:46	7/7/20	
	230 U 230 U	230	32 27	25 25		7/7/20	
Chrysene	230 U	430	<i>∠1</i>	23	07/13/20 14:46	1/1/20	

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Superset Reference:20-0000555360 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 11:50

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-027
 Units: ug/L

 Lab Code:
 R2005701-001
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	230 U	230	46	25	07/13/20 14:46	7/7/20	
Di-n-octyl Phthalate	230 U	230	74	25	07/13/20 14:46	7/7/20	
Dibenz(a,h)anthracene	230 U	230	24	25	07/13/20 14:46	7/7/20	
Dibenzofuran	230 U	230	31	25	07/13/20 14:46	7/7/20	
Diethyl Phthalate	230 U	230	25	25	07/13/20 14:46	7/7/20	
Dimethyl Phthalate	230 U	230	28	25	07/13/20 14:46	7/7/20	
Fluoranthene	230 U	230	34	25	07/13/20 14:46	7/7/20	
Fluorene	230 U	230	28	25	07/13/20 14:46	7/7/20	
Hexachlorobenzene	230 U	230	35	25	07/13/20 14:46	7/7/20	
Hexachlorobutadiene	230 U	230	23	25	07/13/20 14:46	7/7/20	
Hexachlorocyclopentadiene	230 U	230	49	25	07/13/20 14:46	7/7/20	
Hexachloroethane	230 U	230	24	25	07/13/20 14:46	7/7/20	
Indeno(1,2,3-cd)pyrene	230 U	230	40	25	07/13/20 14:46	7/7/20	
Isophorone	230 U	230	31	25	07/13/20 14:46	7/7/20	
N-Nitrosodi-n-propylamine	230 U	230	26	25	07/13/20 14:46	7/7/20	
N-Nitrosodiphenylamine	230 U	230	60	25	07/13/20 14:46	7/7/20	
Naphthalene	230 U	230	27	25	07/13/20 14:46	7/7/20	
Nitrobenzene	230 U	230	34	25	07/13/20 14:46	7/7/20	
Pentachlorophenol (PCP)	1100 U	1100	230	25	07/13/20 14:46	7/7/20	
Phenanthrene	230 U	230	31	25	07/13/20 14:46	7/7/20	
Phenol	32 J	230	23	25	07/13/20 14:46	7/7/20	
Pyrene	230 U	230	33	25	07/13/20 14:46	7/7/20	

Tentatively Identified Compounds

			Result	
CAS#	Compound Identification	RT	ug/L	Q

No Tentatively Identified Compounds Detected

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	0 *	35 - 141	07/13/20 14:46	D
2-Fluorobiphenyl	0 *	31 - 118	07/13/20 14:46	D
2-Fluorophenol	0 *	10 - 105	07/13/20 14:46	D
Nitrobenzene-d5	0 *	31 - 110	07/13/20 14:46	D
Phenol-d6	0 *	10 - 107	07/13/20 14:46	D
p-Terphenyl-d14	0 *	10 - 165	07/13/20 14:46	D

Analytical Report

Client:GHD (Formerly Conestoga-Rovers & Associates)Service Request:R2005701Project:Love Canal:292-402-D02-3100/9954 Annual Long TermDate Collected:07/01/20 11:50

Monitoring

2400 000000000

Sample Matrix: Water

Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-027
 Units: ug/L

 Lab Code:
 R2005701-001
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	67 J	450	51	50	07/13/20 15:42	7/7/20	
1,2-Dichlorobenzene	450 U	450	54	50	07/13/20 15:42	7/7/20	
1,3-Dichlorobenzene	450 U	450	46	50	07/13/20 15:42	7/7/20	
1,4-Dichlorobenzene	73 J	450	52	50	07/13/20 15:42	7/7/20	
2,4,5-Trichlorophenol	450 U	450	50	50	07/13/20 15:42	7/7/20	
2,4,6-Trichlorophenol	450 U	450	62	50	07/13/20 15:42	7/7/20	
2,4-Dichlorophenol	340 J	450	57	50	07/13/20 15:42	7/7/20	
2,4-Dimethylphenol	450 U	450	63	50	07/13/20 15:42	7/7/20	
2,4-Dinitrophenol	2300 U	2300	910	50	07/13/20 15:42	7/7/20	
2,4-Dinitrotoluene	450 U	450	110	50	07/13/20 15:42	7/7/20	
2,6-Dinitrotoluene	450 U	450	60	50	07/13/20 15:42	7/7/20	
2-Chloronaphthalene	450 U	450	62	50	07/13/20 15:42	7/7/20	
2-Chlorophenol	450 U	450	49	50	07/13/20 15:42	7/7/20	
2-Methylnaphthalene	450 U	450	59	50	07/13/20 15:42	7/7/20	
2-Methylphenol	450 U	450	46	50	07/13/20 15:42	7/7/20	
2-Nitroaniline	450 U	450	62	50	07/13/20 15:42	7/7/20	
2-Nitrophenol	450 U	450	68	50	07/13/20 15:42	7/7/20	
3,3'-Dichlorobenzidine	450 U	450	55	50	07/13/20 15:42	7/7/20	
3- and 4-Methylphenol Coelution	61 J	450	53	50	07/13/20 15:42	7/7/20	
3-Nitroaniline	450 U	450	120	50	07/13/20 15:42	7/7/20	
4,6-Dinitro-2-methylphenol	2300 U	2300	870	50	07/13/20 15:42	7/7/20	
4-Bromophenyl Phenyl Ether	450 U	450	75	50	07/13/20 15:42	7/7/20	
4-Chloro-3-methylphenol	450 U	450	49	50	07/13/20 15:42	7/7/20	
4-Chloroaniline	450 U	450	46	50	07/13/20 15:42	7/7/20	
4-Chlorophenyl Phenyl Ether	450 U	450	68	50	07/13/20 15:42	7/7/20	
4-Nitroaniline	450 U	450	130	50	07/13/20 15:42	7/7/20	
4-Nitrophenol	2300 U	2300	290	50	07/13/20 15:42	7/7/20	
Acenaphthene	450 U	450	61	50	07/13/20 15:42	7/7/20	
Acenaphthylene	450 U	450	61	50	07/13/20 15:42	7/7/20	
Anthracene	450 U	450	57	50	07/13/20 15:42	7/7/20	
Benz(a)anthracene	450 U	450	72	50	07/13/20 15:42	7/7/20	
Benzo(a)pyrene	450 U	450	51	50	07/13/20 15:42	7/7/20	
Benzo(b)fluoranthene	450 U	450	51	50	07/13/20 15:42	7/7/20	
Benzo(g,h,i)perylene	450 U	450	46	50	07/13/20 15:42	7/7/20	
Benzo(k)fluoranthene	450 U	450	55	50	07/13/20 15:42	7/7/20	
Benzoic Acid	11000 D	4500	1700	50	07/13/20 15:42	7/7/20	
Benzyl Alcohol	220 J	450	71	50	07/13/20 15:42	7/7/20	
2,2'-Oxybis(1-chloropropane)	450 U	450	64	50	07/13/20 15:42	7/7/20	
Bis(2-chloroethoxy)methane	450 U	450	87	50	07/13/20 15:42	7/7/20	
Bis(2-chloroethyl) Ether	450 U	450	56	50	07/13/20 15:42	7/7/20	
Bis(2-ethylhexyl) Phthalate	450 U	450	46	50	07/13/20 15:42	7/7/20	
Butyl Benzyl Phthalate	450 U	450	64	50	07/13/20 15:42	7/7/20	
Chrysene	450 U	450	53	50	07/13/20 15:42	7/7/20	
Chrysene	430 0	430	33	30	07/13/20 13.42	1/1/20	

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Superset Reference: 20-0000555360 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Love Canal:292-402-D02-3100/9954 Annual Long Term **Project:** Monitoring

Date Collected: 07/01/20 11:50

Sample Matrix: Water

Date Received: 07/02/20 10:50

Sample Name: WG-9954-070120-SG-027

Units: ug/L Lab Code: R2005701-001 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	450 U	450	91	50	07/13/20 15:42	7/7/20	
Di-n-octyl Phthalate	450 U	450	150	50	07/13/20 15:42	7/7/20	
Dibenz(a,h)anthracene	450 U	450	47	50	07/13/20 15:42	7/7/20	
Dibenzofuran	450 U	450	62	50	07/13/20 15:42	7/7/20	
Diethyl Phthalate	450 U	450	49	50	07/13/20 15:42	7/7/20	
Dimethyl Phthalate	450 U	450	56	50	07/13/20 15:42	7/7/20	
Fluoranthene	450 U	450	67	50	07/13/20 15:42	7/7/20	
Fluorene	450 U	450	56	50	07/13/20 15:42	7/7/20	
Hexachlorobenzene	450 U	450	70	50	07/13/20 15:42	7/7/20	
Hexachlorobutadiene	450 U	450	46	50	07/13/20 15:42	7/7/20	
Hexachlorocyclopentadiene	450 U	450	98	50	07/13/20 15:42	7/7/20	
Hexachloroethane	450 U	450	48	50	07/13/20 15:42	7/7/20	
Indeno(1,2,3-cd)pyrene	450 U	450	80	50	07/13/20 15:42	7/7/20	
Isophorone	450 U	450	62	50	07/13/20 15:42	7/7/20	
N-Nitrosodi-n-propylamine	450 U	450	52	50	07/13/20 15:42	7/7/20	
N-Nitrosodiphenylamine	450 U	450	120	50	07/13/20 15:42	7/7/20	
Naphthalene	450 U	450	54	50	07/13/20 15:42	7/7/20	
Nitrobenzene	450 U	450	67	50	07/13/20 15:42	7/7/20	
Pentachlorophenol (PCP)	2300 U	2300	450	50	07/13/20 15:42	7/7/20	
Phenanthrene	450 U	450	61	50	07/13/20 15:42	7/7/20	
Phenol	450 U	450	46	50	07/13/20 15:42	7/7/20	
Pyrene	450 U	450	65	50	07/13/20 15:42	7/7/20	

Tentatively Identified Compounds

			Result	
CAS#	Compound Identification	RT	ug/L	Q

No Tentatively Identified Compounds Detected

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q	
2,4,6-Tribromophenol	0 *	35 - 141	07/13/20 15:42	D	
2-Fluorobiphenyl	0 *	31 - 118	07/13/20 15:42	D	
2-Fluorophenol	0 *	10 - 105	07/13/20 15:42	D	
Nitrobenzene-d5	0 *	31 - 110	07/13/20 15:42	D	
Phenol-d6	0 *	10 - 107	07/13/20 15:42	D	
p-Terphenyl-d14	0 *	10 - 165	07/13/20 15:42	D	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 10:45

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-026
 Units: ug/L

 Lab Code:
 R2005701-002
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	7.7 J	9.1	1.1	1	07/09/20 13:32	7/7/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/09/20 13:32	7/7/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/09/20 13:32	7/7/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/09/20 13:32	7/7/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/09/20 13:32	7/7/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/09/20 13:32	7/7/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/09/20 13:32	7/7/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/09/20 13:32	7/7/20	
2,4-Dinitrophenol	45 U	45	19	1	07/09/20 13:32	7/7/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/09/20 13:32	7/7/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/09/20 13:32	7/7/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/09/20 13:32	7/7/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/09/20 13:32	7/7/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/09/20 13:32	7/7/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/09/20 13:32	7/7/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/09/20 13:32	7/7/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/09/20 13:32	7/7/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/09/20 13:32	7/7/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/09/20 13:32	7/7/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/09/20 13:32	7/7/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/09/20 13:32	7/7/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/09/20 13:32	7/7/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/09/20 13:32	7/7/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/09/20 13:32	7/7/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/09/20 13:32	7/7/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/09/20 13:32	7/7/20	
4-Nitrophenol	45 U	45	5.8	1	07/09/20 13:32	7/7/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/09/20 13:32	7/7/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/09/20 13:32	7/7/20	
Anthracene	9.1 U	9.1	1.2	1	07/09/20 13:32	7/7/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/09/20 13:32	7/7/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/09/20 13:32	7/7/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/09/20 13:32	7/7/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/09/20 13:32	7/7/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/09/20 13:32	7/7/20	
Benzoic Acid	91 U	91	33	1	07/09/20 13:32	7/7/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/09/20 13:32	7/7/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/09/20 13:32	7/7/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/09/20 13:32	7/7/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	07/09/20 13:32	7/7/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	07/09/20 13:32	7/7/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	07/09/20 13:32	7/7/20	
Chrysene	9.1 U	9.1	1.1	1	07/09/20 13:32	7/7/20	
	-			-			

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Superset Reference:20-0000555360 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 10:45

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-026
 Units: ug/L

 Lab Code:
 R2005701-002
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/09/20 13:32	7/7/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/09/20 13:32	7/7/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/09/20 13:32	7/7/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/09/20 13:32	7/7/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/09/20 13:32	7/7/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/09/20 13:32	7/7/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/09/20 13:32	7/7/20	
Fluorene	9.1 U	9.1	1.2	1	07/09/20 13:32	7/7/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/09/20 13:32	7/7/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/09/20 13:32	7/7/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/09/20 13:32	7/7/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/09/20 13:32	7/7/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/09/20 13:32	7/7/20	
Isophorone	9.1 U	9.1	1.3	1	07/09/20 13:32	7/7/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/09/20 13:32	7/7/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/09/20 13:32	7/7/20	
Naphthalene	9.1 U	9.1	1.1	1	07/09/20 13:32	7/7/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/09/20 13:32	7/7/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/09/20 13:32	7/7/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/09/20 13:32	7/7/20	
Phenol	9.1 U	9.1	0.91	1	07/09/20 13:32	7/7/20	
Pyrene	9.1 U	9.1	1.3	1	07/09/20 13:32	7/7/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	111	35 - 141	07/09/20 13:32	
2-Fluorobiphenyl	75	31 - 118	07/09/20 13:32	
2-Fluorophenol	45	10 - 105	07/09/20 13:32	
Nitrobenzene-d5	70	31 - 110	07/09/20 13:32	
Phenol-d6	33	10 - 107	07/09/20 13:32	
p-Terphenyl-d14	47	10 - 165	07/09/20 13:32	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown hydrocarbon	13.12	5.4	J	
	unknown hydrocarbon	14.57	4.9	J	
	unknown hydrocarbon	16.11	4.8	J	
000095-49-8	Benzene, 1-chloro-2-methyl-	4.13	7.9	JN	
000143-07-7	Dodecanoic acid	7.67	6.1	JN	
013798-23-7	Sulfur	7.78	14	JN	
	unknown	9.45	11	J	

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Analytical Report

Client: Service Request: R2005701 GHD (Formerly Conestoga-Rovers & Associates) **Date Collected:** 07/01/20 09:55

Love Canal:292-402-D02-3100/9954 Annual Long Term **Project:**

Monitoring

Date Received: 07/02/20 10:50 **Sample Matrix:** Water

Sample Name: WG-9954-070120-SG-025 Units: ug/L Basis: NA Lab Code: R2005701-003

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/09/20 14:02	7/7/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/09/20 14:02	7/7/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/09/20 14:02	7/7/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/09/20 14:02	7/7/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/09/20 14:02	7/7/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/09/20 14:02	7/7/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/09/20 14:02	7/7/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/09/20 14:02	7/7/20	
2,4-Dinitrophenol	45 U	45	19	1	07/09/20 14:02	7/7/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/09/20 14:02	7/7/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/09/20 14:02	7/7/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/09/20 14:02	7/7/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/09/20 14:02	7/7/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/09/20 14:02	7/7/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/09/20 14:02	7/7/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/09/20 14:02	7/7/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/09/20 14:02	7/7/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/09/20 14:02	7/7/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/09/20 14:02	7/7/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/09/20 14:02	7/7/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/09/20 14:02	7/7/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/09/20 14:02	7/7/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/09/20 14:02	7/7/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/09/20 14:02	7/7/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/09/20 14:02	7/7/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/09/20 14:02	7/7/20	
4-Nitrophenol	45 U	45	5.8	1	07/09/20 14:02	7/7/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/09/20 14:02	7/7/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/09/20 14:02	7/7/20	
Anthracene	9.1 U	9.1	1.2	1	07/09/20 14:02	7/7/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/09/20 14:02	7/7/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/09/20 14:02	7/7/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/09/20 14:02	7/7/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/09/20 14:02	7/7/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/09/20 14:02	7/7/20	
Benzoic Acid	91 U	91	33	1	07/09/20 14:02	7/7/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/09/20 14:02	7/7/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/09/20 14:02	7/7/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/09/20 14:02	7/7/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	07/09/20 14:02	7/7/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	07/09/20 14:02	7/7/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	07/09/20 14:02	7/7/20	
Chrysene	9.1 U	9.1	1.1	1	07/09/20 14:02	7/7/20	

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Superset Reference:20-0000555360 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 09:55

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-025
 Units: ug/L

 Lab Code:
 R2005701-003
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/09/20 14:02	7/7/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/09/20 14:02	7/7/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/09/20 14:02	7/7/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/09/20 14:02	7/7/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/09/20 14:02	7/7/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/09/20 14:02	7/7/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/09/20 14:02	7/7/20	
Fluorene	9.1 U	9.1	1.2	1	07/09/20 14:02	7/7/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/09/20 14:02	7/7/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/09/20 14:02	7/7/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/09/20 14:02	7/7/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/09/20 14:02	7/7/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/09/20 14:02	7/7/20	
Isophorone	9.1 U	9.1	1.3	1	07/09/20 14:02	7/7/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/09/20 14:02	7/7/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/09/20 14:02	7/7/20	
Naphthalene	9.1 U	9.1	1.1	1	07/09/20 14:02	7/7/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/09/20 14:02	7/7/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/09/20 14:02	7/7/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/09/20 14:02	7/7/20	
Phenol	9.1 U	9.1	0.91	1	07/09/20 14:02	7/7/20	
Pyrene	9.1 U	9.1	1.3	1	07/09/20 14:02	7/7/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	93	35 - 141	07/09/20 14:02	
2-Fluorobiphenyl	63	31 - 118	07/09/20 14:02	
2-Fluorophenol	40	10 - 105	07/09/20 14:02	
Nitrobenzene-d5	62	31 - 110	07/09/20 14:02	
Phenol-d6	25	10 - 107	07/09/20 14:02	
p-Terphenyl-d14	51	10 - 165	07/09/20 14:02	

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
000095-49-8	Benzene, 1-chloro-2-methyl-	4.13	7.6	JN	
000108-41-8	Benzene, 1-chloro-3-methyl-	4.18	3.8	JN	
013798-23-7	Sulfur	7.79	4.8	JN	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Co

Date Collected: 07/01/20 08:45

Monitoring
Sample Matrix: Water

Water **Date Received:** 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-024
 Units: ug/L

 Lab Code:
 R2005701-004
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/09/20 14:33	7/7/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/09/20 14:33	7/7/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/09/20 14:33	7/7/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/09/20 14:33	7/7/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/09/20 14:33	7/7/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/09/20 14:33	7/7/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/09/20 14:33	7/7/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/09/20 14:33	7/7/20	
2,4-Dinitrophenol	45 U	45	19	1	07/09/20 14:33	7/7/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/09/20 14:33	7/7/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/09/20 14:33	7/7/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/09/20 14:33	7/7/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/09/20 14:33	7/7/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/09/20 14:33	7/7/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/09/20 14:33	7/7/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/09/20 14:33	7/7/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/09/20 14:33	7/7/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/09/20 14:33	7/7/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/09/20 14:33	7/7/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/09/20 14:33	7/7/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/09/20 14:33	7/7/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/09/20 14:33	7/7/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/09/20 14:33	7/7/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/09/20 14:33	7/7/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/09/20 14:33	7/7/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/09/20 14:33	7/7/20	
4-Nitrophenol	45 U	45	5.8	1	07/09/20 14:33	7/7/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/09/20 14:33	7/7/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/09/20 14:33	7/7/20	
Anthracene	9.1 U	9.1	1.2	1	07/09/20 14:33	7/7/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/09/20 14:33	7/7/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/09/20 14:33	7/7/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/09/20 14:33	7/7/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/09/20 14:33	7/7/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/09/20 14:33	7/7/20	
Benzoic Acid	9.1 U	9.1	33	1	07/09/20 14:33	7/7/20	
	9.1 U	9.1	33 1.5	1	07/09/20 14:33	7/7/20	
Benzyl Alcohol 2,2'-Oxybis(1-chloropropane)		9.1 9.1	1.3	1	07/09/20 14:33	7/7/20	
	9.1 U	9.1 9.1	1.3			7/7/20 7/7/20	
Bis(2-chloroethoxy)methane	9.1 U			1	07/09/20 14:33		
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	07/09/20 14:33	7/7/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	07/09/20 14:33	7/7/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	07/09/20 14:33	7/7/20	
Chrysene	9.1 U	9.1	1.1	1	07/09/20 14:33	7/7/20	

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Superset Reference: 20-0000555360 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 08:45

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-024
 Units: ug/L

 Lab Code:
 R2005701-004
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/09/20 14:33	7/7/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/09/20 14:33	7/7/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/09/20 14:33	7/7/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/09/20 14:33	7/7/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/09/20 14:33	7/7/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/09/20 14:33	7/7/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/09/20 14:33	7/7/20	
Fluorene	9.1 U	9.1	1.2	1	07/09/20 14:33	7/7/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/09/20 14:33	7/7/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/09/20 14:33	7/7/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/09/20 14:33	7/7/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/09/20 14:33	7/7/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/09/20 14:33	7/7/20	
Isophorone	9.1 U	9.1	1.3	1	07/09/20 14:33	7/7/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/09/20 14:33	7/7/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/09/20 14:33	7/7/20	
Naphthalene	9.1 U	9.1	1.1	1	07/09/20 14:33	7/7/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/09/20 14:33	7/7/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/09/20 14:33	7/7/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/09/20 14:33	7/7/20	
Phenol	9.1 U	9.1	0.91	1	07/09/20 14:33	7/7/20	
Pyrene	9.1 U	9.1	1.3	1	07/09/20 14:33	7/7/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	80	35 - 141	07/09/20 14:33	
2-Fluorobiphenyl	59	31 - 118	07/09/20 14:33	
2-Fluorophenol	35	10 - 105	07/09/20 14:33	
Nitrobenzene-d5	56	31 - 110	07/09/20 14:33	
Phenol-d6	26	10 - 107	07/09/20 14:33	
p-Terphenyl-d14	65	10 - 165	07/09/20 14:33	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown	11.35	5.7	J	
	unknown hydrocarbon	11.89	8.4	J	
	unknown	12.48	11	J	
	unknown hydrocarbon	13.13	11	J	
	unknown hydrocarbon	13.83	10	J	
	unknown hydrocarbon	14.58	8.2	J	
1000309-12-4	Sulfurous acid, 2-propyl tridecyl	15.37	7.0	JN	

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Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 08:45

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-024
 Units: ug/L

 Lab Code:
 R2005701-004
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown	16.11	5.8	J	
	unknown	4.13	5.1	J	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Date Collected: 07/01/20 11:50

Monitoring Sample Matrix: Water

Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-028
 Units: ug/L

 Lab Code:
 R2005701-006
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

1.2.4-Trichlorobenzene	Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1.3-Dichlorobenzene	1,2,4-Trichlorobenzene	100	9.1	1.1	1	07/09/20 15:04	7/7/20	
JDichlorobenzene 97 9.1 1.1 1 07/09/20 15:04 77/20 24.5-Trichlorophenol 26 9.1 0.99 1 07/09/20 15:04 77/20 24.5-Trichlorophenol 91 U 9.1 1.3 1 07/09/20 15:04 77/20 2.4-Dinthylphenol 6.9 J 9.1 1.3 1 07/09/20 15:04 77/20 2.4-Dinitrophenol 45 U 45 19 1 07/09/20 15:04 77/20 2.4-Dinitrophenol 45 U 45 19 1 07/09/20 15:04 77/20 2.4-Dinitrophenol 9.1 U 9.1 1.2 1 07/09/20 15:04 77/20 2.4-Dinitrophenol 9.1 U 9.1 1.2 1 07/09/20 15:04 77/20 2.6-Dinitrotoluene 9.1 U 9.1 1.2 1 07/09/20 15:04 77/20 2.6-Dinitrotoluene 9.1 U 9.1 1.3 1 07/09/20 15:04 77/20 2.6-Dinitrotoluene 9.1 U 9.1 1.3 1 07/09/20 15:04 77/20 2.6-Dinitrotoluene 9.1 U 9.1 1.3 1 07/09/20 15:04 77/20 2.6-Dinitrotoluene 9.1 U 9.1 1.2 1 07/09/20 15:04 77/20 2.6-Dinitrotoluene 9.1 U 9.1 1.2 1 07/09/20 15:04 77/20 2.6-Dinitrotoluene 9.1 U 9.1 1.2 1 07/09/20 15:04 77/20 2.6-Dinitrotoluene 9.1 U 9.1 1.3 1 07/09/20 15:04 77/20 2.6-Dinitrotoluene 9.1 U 9.1 1.3 1 07/09/20 15:04 77/20 2.6-Dinitrotoluene 9.1 U 9.1 1.3 1 07/09/20 15:04 77/20 2.6-Dinitrotoluene 9.1 U 9.1 1.3 1 07/09/20 15:04 77/20 2.6-Dinitrotoluene 9.1 U 9.1 1.1 1 07/09/20 15:04 77/20 2.6-Dinitrotoluene 9.1 U 9.1 1.1 1 07/09/20 15:04 77/20 2.6-Dinitrotoluene 9.1 U 9.1 1.1 1 07/09/20 15:04 77/20 2.6-Dinitrotoluene 9.1 U 9.1 1.1 1 07/09/20 15:04 77/20 2.6-Dinitrotoluene 9.1 U 9.1 1.1 1 07/09/20 15:04 77/20 2.6-Dinitrotoluene 9.1 U 9.1 1.1 1 07/09/20 15:04 77/20 2.6-Dinitrotoluene 9.1 U 9.1 1.3 1 07/09/20 15:04 77/20 3.6-Dinitrotoluene 9.1 U 9.1 1.3 1 07/09/20 15:04 77/20 3.6-Dinit	1,2-Dichlorobenzene	34	9.1	1.1	1	07/09/20 15:04	7/7/20	
2.45-Trichlorophenol 26	1,3-Dichlorobenzene	4.1 J	9.1	0.92	1	07/09/20 15:04	7/7/20	
2.4.6-Trichlorophenol	1,4-Dichlorobenzene	97	9.1	1.1	1	07/09/20 15:04	7/7/20	
2.4-Dinterhylphenol 6.9 J 9.1 1.2 1 07/09/20 15:04 77/20	2,4,5-Trichlorophenol	26	9.1	0.99	1	07/09/20 15:04	7/7/20	
2.4-Dimitryphenol	2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/09/20 15:04	7/7/20	
2.4-Dimitrophenol	2,4-Dichlorophenol	520 E	9.1	1.2	1	07/09/20 15:04	7/7/20	
2.4-Dinitrotoluene		6.9 J	9.1	1.3	1	07/09/20 15:04	7/7/20	
2.4-Dinitrotoluene	2,4-Dinitrophenol	45 U	45	19	1	07/09/20 15:04	7/7/20	
2-Chloronaphthalene		9.1 U	9.1	2.2	1	07/09/20 15:04	7/7/20	
2-Chlorophenol 31 9.1 0.97 1 07/09/20 15:04 77/20		9.1 U	9.1	1.2	1	07/09/20 15:04	7/7/20	
2-Chlorophenol 31 9.1 0.97 1 07/09/20 15:04 77/20	2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/09/20 15:04	7/7/20	
2-Methylnaphthalene		31	9.1	0.97	1	07/09/20 15:04	7/7/20	
2-Methylphenol 32 9.1 0.91 1 07/09/20 15:04 77/20		9.1 U	9.1		1	07/09/20 15:04	7/7/20	
2-Nitroaniline								
2-Nitrophenol	2-Nitroaniline							
3,3'-Dichlorobenzidine								
3- and 4-Methylphenol Coelution 9.1 U 9.1 2.3 1 07/09/20 15:04 77/20								
3-Nitroaniline								
4,6-Dinitro-2-methylphenol 45 U 45 I8 1 07/09/20 15:04 7/7/20 4-Bromophenyl Phenyl Ether 9.1 U 9.1 1.5 1 07/09/20 15:04 7/7/20 4-Chloro-3-methylphenol 40 9.1 0.98 1 07/09/20 15:04 7/7/20 4-Chloroaniline 9.1 U 9.1 0.91 1 07/09/20 15:04 7/7/20 4-Chlorophenyl Phenyl Ether 9.1 U 9.1 1.4 1 07/09/20 15:04 7/7/20 4-Nitroaniline 9.1 U 9.1 U 9.1 2.5 1 07/09/20 15:04 7/7/20 4-Nitrophenol 45 U 45 S.8 1 07/09/20 15:04 7/7/20 4-Nitrophenol 45 U 45 S.8 1 07/09/20 15:04 7/7/20 Acenaphthene 9.1 U 9.1 I.3 1 07/09/20 15:04 7/7/20 Acenaphthylene 9.1 U 9.1 I.3 1 07/09/20 15:04 7/7/20 Achtracene 9.1 U 9.1 I.2 1 07/09/20 15:04 7/7/20 Benzo(a)pyrene 9.1 U								
4-Bromophenyl Phenyl Ether 9.1 U 9.1 1.5 1 07/09/20 15:04 7/7/20 4-Chloro-3-methylphenol 40 9.1 0.98 1 07/09/20 15:04 7/7/20 4-Chloroaniline 9.1 U 9.1 0.91 1 07/09/20 15:04 7/7/20 4-Chlorophenyl Phenyl Ether 9.1 U 9.1 1.4 1 07/09/20 15:04 7/7/20 4-Nitrophenol 45 U 45 5.8 1 07/09/20 15:04 7/7/20 4-Nitrophenol 45 U 45 5.8 1 07/09/20 15:04 7/7/20 Acenaphthene 9.1 U 9.1 U 9.1 1.3 1 07/09/20 15:04 7/7/20 Acenaphthylene 9.1 U 9.1 1.3 1 07/09/20 15:04 7/7/20 Anthracene 9.1 U 9.1 1.3 1 07/09/20 15:04 7/7/20 Benz(a)phtracene 9.1 U 9.1 1.5 1 07/09/20 15:04 7/7/20 Benzo(b)fluoranthene 9.1 U 9.1 1.1 1 07/09/20 15:04 7/7/20 Benzo(b)fluoranthene 9.1 U 9.1 1.1 1 07/09/20 15:04 7/7/20 Benzo(k)fluoranthene 9.1 U 9.1 1.1 1 07/09/20 15:04 7/7/20 Benzo(k)fluoranthene 9.1 U								
4-Chloro-3-methylphenol 40 9.1 0.98 1 07/09/20 15:04 7/7/20 4-Chloroaniline 9.1 U 9.1 0.91 1 07/09/20 15:04 7/7/20 4-Chlorophenyl Phenyl Ether 9.1 U 9.1 1.4 1 07/09/20 15:04 7/7/20 4-Nitroaniline 9.1 U 9.1 2.5 1 07/09/20 15:04 7/7/20 4-Nitrophenol 45 U 45 5.8 1 07/09/20 15:04 7/7/20 Acenaphthene 9.1 U 9.1 1.3 1 07/09/20 15:04 7/7/20 Acenaphthylene 9.1 U 9.1 1.3 1 07/09/20 15:04 7/7/20 Acenaphthylene 9.1 U 9.1 1.3 1 07/09/20 15:04 7/7/20 Acenaphthylene 9.1 U 9.1 1.5 1 07/09/20 15:04 7/7/20 Acenaphthylene 9.1 U 9.1 1.5 1 07/09/20 15:04 7/7/20 Benzo(a)pyrene 9.1 U 9.1 U 9.1 <								
4-Chloroaniline 9.1 U 9.1 U 9.1 U 9.1 U 0.91 U 1 07/09/20 15:04 7/7/20 4-Chlorophenyl Phenyl Ether 9.1 U 9.1 U 9.1 U.4 U 1.4 U 1 07/09/20 15:04 7/7/20 4-Nitroaniline 9.1 U 9.1 U 9.1 D.5 1 07/09/20 15:04 7/7/20 4-Nitrophenol 45 U 45 D.5.8 D.5.8 D.5.8 D.5.04 7/7/20 1 07/09/20 15:04 7/7/20 Acenaphthene 9.1 U D.5 D.1 U D.5 D.1 D.3 D.5.04 7/7/20 1.3 D.7/09/20 15:04 7/7/20 Acenaphthylene 9.1 U D.5 D.1 D.5 D.1 D.5 D.4 7/7/20 Anthracene 9.1 U D.5 D.1 D.5 D.1 D.5 D.4 7/7/20 Benz(a)anthracene 9.1 U D.5 D.1 D.5 D.1 D.5 D.4 7/7/20 Benzo(a)pyrene 9.1 U D.5 D.1 D.5 D.1 D.5 D.4 7/7/20 Benzo(b)fluoranthene 9.1 U D.5 D.1 D.5 D.1 D.5 D.4 7/7/20 Benzo(g,h,i)perylene 9.1 U D.5 D.1 D.1 D.5 D.1 D.7/09/20 15:04 7/7/20 Benzo(k)fluoranthene 9.1 U D.5 D.1 D.1 D.5 D.1 D.7/09/20 15:04 7/7/20 Benzoic Acid 4100 E. D.1 D.1 D.1 D.7/09/20 15:04 7/7/20 Benzyl Alcohol 290 D.1 D.1 D.1 D.1 D.5 D.1 D.7/09/20 15:04 7/7/20 Bis(2-chloroethoxy)methane 9.1 U D.1 D.1 D.1 D.1 D.7/09/20 15:04 7/7/20 Bis(
4-Chlorophenyl Phenyl Ether 9.1 U 9.1 1.4 1 07/09/20 15:04 7/7/20 4-Nitroaniline 9.1 U 9.1 2.5 1 07/09/20 15:04 7/7/20 4-Nitrophenol 45 U 45 5.8 1 07/09/20 15:04 7/7/20 Acenaphthene 9.1 U 9.1 1.3 1 07/09/20 15:04 7/7/20 Acenaphthylene 9.1 U 9.1 1.3 1 07/09/20 15:04 7/7/20 Anthracene 9.1 U 9.1 1.2 1 07/09/20 15:04 7/7/20 Benza(a)anthracene 9.1 U 9.1 1.5 1 07/09/20 15:04 7/7/20 Benzo(a)pyrene 9.1 U 9.1 1.1 1 07/09/20 15:04 7/7/20 Benzo(b)fluoranthene 9.1 U 9.1 1.1 1 07/09/20 15:04 7/7/20 Benzo(c,h,i)perylene 9.1 U 9.1 0.91 1.1 07/09/20 15:04 7/7/20 Benzo(k)fluoranthene 9.1 U 9.1 1.1 1 07/09/20 15:04 7/7/20 Benzo(c Acid 4100 E 91 33 1 07/09/20 15:04 7/7/20 Benzyl Alcohol 290 9.1 1.5 1 07/09/20 15:04 7								
4-Nitroaniline 9.1 U 9.1 Z.5 1 07/09/20 15:04 7/7/20 4-Nitrophenol 45 U 45 S.8 1 07/09/20 15:04 7/7/20 Acenaphthene 9.1 U 9.1 I.3 1 07/09/20 15:04 7/7/20 Acenaphthylene 9.1 U 9.1 I.3 1 07/09/20 15:04 7/7/20 Anthracene 9.1 U 9.1 I.2 1 07/09/20 15:04 7/7/20 Benz(a)anthracene 9.1 U 9.1 I.5 1 07/09/20 15:04 7/7/20 Benzo(a)pyrene 9.1 U 9.1 I.1 1 07/09/20 15:04 7/7/20 Benzo(b)fluoranthene 9.1 U 9.1 I.1 1 07/09/20 15:04 7/7/20 Benzo(b)fluoranthene 9.1 U 9.1 U.9.1 1.1 1 07/09/20 15:04 7/7/20 Benzo(k)fluoranthene 9.1 U.9.1 1.1 1 07/09/20 15:04 7/7/20 Benzoic Acid 4100 E.91 91 33 1 07/09/20 15:04 7/7/20 Benzyl Alcohol 290 9.1 I.5 1 07/09/20 15:04 7/7/20 2,2'-Oxybis(1-chloropropane) 9.1 U.9.1 1.3								
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Butyl Benzyl Phthalate 9.1 U 9.1 1.3 1 07/09/20 15:04 7/7/20								
	Chrysene	9.1 U	9.1	1.1	1	07/09/20 15:04	7/7/20	

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Superset Reference: 20-0000555360 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 11:50

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-028
 Units: ug/L

 Lab Code:
 R2005701-006
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/09/20 15:04	7/7/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/09/20 15:04	7/7/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/09/20 15:04	7/7/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/09/20 15:04	7/7/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/09/20 15:04	7/7/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/09/20 15:04	7/7/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/09/20 15:04	7/7/20	
Fluorene	9.1 U	9.1	1.2	1	07/09/20 15:04	7/7/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/09/20 15:04	7/7/20	
Hexachlorobutadiene	1.0 J	9.1	0.91	1	07/09/20 15:04	7/7/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/09/20 15:04	7/7/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/09/20 15:04	7/7/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/09/20 15:04	7/7/20	
Isophorone	9.1 U	9.1	1.3	1	07/09/20 15:04	7/7/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/09/20 15:04	7/7/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/09/20 15:04	7/7/20	
Naphthalene	9.1 U	9.1	1.1	1	07/09/20 15:04	7/7/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/09/20 15:04	7/7/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/09/20 15:04	7/7/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/09/20 15:04	7/7/20	
Phenol	44	9.1	0.91	1	07/09/20 15:04	7/7/20	
Pyrene	9.1 U	9.1	1.3	1	07/09/20 15:04	7/7/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	66	35 - 141	07/09/20 15:04	
2-Fluorobiphenyl	57	31 - 118	07/09/20 15:04	
2-Fluorophenol	32	10 - 105	07/09/20 15:04	
Nitrobenzene-d5	95	31 - 110	07/09/20 15:04	
Phenol-d6	27	10 - 107	07/09/20 15:04	
p-Terphenyl-d14	46	10 - 165	07/09/20 15:04	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
000611-95-0	Benzoic acid, 4-benzoyl-	10.19	630	JN	
000579-18-0	3-Benzoylbenzoic acid	10.22	170	JN	
000123-91-1	1,4-Dioxane	2.19	140	JN	
000107-92-6	Butanoic acid	2.92	100	JN	
000108-90-7	Benzene, chloro-	3.32	400	JN	
000108-93-0	Cyclohexanol	3.64	130	JN	
	unknown	3.70	38	J	

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Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 11:50

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-028
 Units: ug/L

 Lab Code:
 R2005701-006
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

		Result			
CAS#	Compound Identification	RT	ug/L	Q	
007443-52-9	Cyclohexanol, 2-methyl-, trans-	4.05	18	JN	
000591-23-1	Cyclohexanol, 3-methyl-	4.09	29	JN	
000095-49-8	Benzene, 1-chloro-2-methyl-	4.14	2000	JN	
000106-43-4	Benzene, 1-chloro-4-methyl-	4.19	1200	JN	
001195-79-5	Bicyclo[2.2.1]heptan-2-one, 1,3,3-	5.15	32	JN	
000106-48-9	Parachlorophenol	5.75	150	JN	
000535-80-8	Benzoic acid, 3-chloro-	6.62	180	JN	
000074-11-3	Benzoic acid, 4-chloro-	6.89	6200	JN	
000095-77-2	Phenol, 3,4-dichloro-	7.01	260	JN	
	unknown	7.12	63	J	
000050-79-3	Benzoic acid, 2,5-dichloro-	7.60	30	JN	
000713-36-0	Benzene, 1-methyl-2-(phenylmethyl)	7.64	17	JN	
000134-84-9	Methanone, (4-methylphenyl)phenyl-	8.73	27	JN	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701 **Date Collected:** 07/01/20 11:50

Love Canal:292-402-D02-3100/9954 Annual Long Term **Project:**

Monitoring

Date Received: 07/02/20 10:50 **Sample Matrix:** Water

Sample Name: Units: ug/L WG-9954-070120-SG-028 Basis: NA Lab Code: R2005701-006

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	76 J	230	26	25	07/13/20 15:14	7/7/20	
1,2-Dichlorobenzene	37 J	230	27	25	07/13/20 15:14	7/7/20	
1,3-Dichlorobenzene	230 U	230	23	25	07/13/20 15:14	7/7/20	
1,4-Dichlorobenzene	100 J	230	26	25	07/13/20 15:14	7/7/20	
2,4,5-Trichlorophenol	34 J	230	25	25	07/13/20 15:14	7/7/20	
2,4,6-Trichlorophenol	230 U	230	31	25	07/13/20 15:14	7/7/20	
2,4-Dichlorophenol	350 D	230	29	25	07/13/20 15:14	7/7/20	
2,4-Dimethylphenol	230 U	230	32	25	07/13/20 15:14	7/7/20	
2,4-Dinitrophenol	1100 U	1100	460	25	07/13/20 15:14	7/7/20	
2,4-Dinitrotoluene	230 U	230	54	25	07/13/20 15:14	7/7/20	
2,6-Dinitrotoluene	230 U	230	30	25	07/13/20 15:14	7/7/20	
2-Chloronaphthalene	230 U	230	31	25	07/13/20 15:14	7/7/20	
2-Chlorophenol	25 J	230	25	25	07/13/20 15:14	7/7/20	
2-Methylnaphthalene	230 U	230	30	25	07/13/20 15:14	7/7/20	
2-Methylphenol	33 J	230	23	25	07/13/20 15:14	7/7/20	
2-Nitroaniline	230 U	230	31	25	07/13/20 15:14	7/7/20	
2-Nitrophenol	230 U	230	34	25	07/13/20 15:14	7/7/20	
3,3'-Dichlorobenzidine	230 U	230	28	25	07/13/20 15:14	7/7/20	
3- and 4-Methylphenol Coelution	60 J	230	27	25	07/13/20 15:14	7/7/20	
3-Nitroaniline	230 U	230	56	25	07/13/20 15:14	7/7/20	
4,6-Dinitro-2-methylphenol	1100 U	1100	440	25	07/13/20 15:14	7/7/20	
4-Bromophenyl Phenyl Ether	230 U	230	38	25	07/13/20 15:14	7/7/20	
4-Chloro-3-methylphenol	30 J	230	25	25	07/13/20 15:14	7/7/20	
4-Chloroaniline	230 U	230	23	25	07/13/20 15:14	7/7/20	
4-Chlorophenyl Phenyl Ether	230 U	230	34	25	07/13/20 15:14	7/7/20	
4-Nitroaniline	230 U	230	62	25	07/13/20 15:14	7/7/20	
4-Nitrophenol	1100 U	1100	150	25	07/13/20 15:14	7/7/20	
Acenaphthene	230 U	230	31	25	07/13/20 15:14	7/7/20	
Acenaphthylene	230 U	230	31	25	07/13/20 15:14	7/7/20	
Anthracene	230 U	230	29	25	07/13/20 15:14	7/7/20	
Benz(a)anthracene	230 U	230	36	25	07/13/20 15:14	7/7/20	
Benzo(a)pyrene	230 U	230	26	25	07/13/20 15:14	7/7/20	
Benzo(b)fluoranthene	230 U	230	26	25	07/13/20 15:14	7/7/20	
Benzo(g,h,i)perylene	230 U	230	23	25	07/13/20 15:14	7/7/20	
Benzo(k)fluoranthene	230 U	230	28	25	07/13/20 15:14	7/7/20	
Benzoic Acid	9200 E	2300	820	25	07/13/20 15:14	7/7/20	
Benzyl Alcohol	250	230	36	25	07/13/20 15:14	7/7/20	
2,2'-Oxybis(1-chloropropane)	230 U	230	32	25	07/13/20 15:14	7/7/20	
Bis(2-chloroethoxy)methane	230 U	230	44	25	07/13/20 15:14	7/7/20	
Bis(2-chloroethyl) Ether	230 U	230	28	25	07/13/20 15:14	7/7/20	
Bis(2-ethylhexyl) Phthalate	230 U	230	23	25	07/13/20 15:14	7/7/20	
Butyl Benzyl Phthalate	230 U	230	32	25	07/13/20 15:14	7/7/20	
Chrysene	230 U	230	27	25	07/13/20 15:14	7/7/20	

Printed 7/30/2020 4:06:15 PM

Superset Reference:20-0000555360 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701 **Date Collected:** 07/01/20 11:50

Love Canal:292-402-D02-3100/9954 Annual Long Term **Project:**

Monitoring

Date Received: 07/02/20 10:50 **Sample Matrix:** Water

Sample Name: WG-9954-070120-SG-028 Units: ug/L Lab Code: R2005701-006 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	230 U	230	46	25	07/13/20 15:14	7/7/20	
Di-n-octyl Phthalate	230 U	230	74	25	07/13/20 15:14	7/7/20	
Dibenz(a,h)anthracene	230 U	230	24	25	07/13/20 15:14	7/7/20	
Dibenzofuran	230 U	230	31	25	07/13/20 15:14	7/7/20	
Diethyl Phthalate	230 U	230	25	25	07/13/20 15:14	7/7/20	
Dimethyl Phthalate	230 U	230	28	25	07/13/20 15:14	7/7/20	
Fluoranthene	230 U	230	34	25	07/13/20 15:14	7/7/20	
Fluorene	230 U	230	28	25	07/13/20 15:14	7/7/20	
Hexachlorobenzene	230 U	230	35	25	07/13/20 15:14	7/7/20	
Hexachlorobutadiene	230 U	230	23	25	07/13/20 15:14	7/7/20	
Hexachlorocyclopentadiene	230 U	230	49	25	07/13/20 15:14	7/7/20	
Hexachloroethane	230 U	230	24	25	07/13/20 15:14	7/7/20	
Indeno(1,2,3-cd)pyrene	230 U	230	40	25	07/13/20 15:14	7/7/20	
Isophorone	230 U	230	31	25	07/13/20 15:14	7/7/20	
N-Nitrosodi-n-propylamine	230 U	230	26	25	07/13/20 15:14	7/7/20	
N-Nitrosodiphenylamine	230 U	230	60	25	07/13/20 15:14	7/7/20	
Naphthalene	230 U	230	27	25	07/13/20 15:14	7/7/20	
Nitrobenzene	230 U	230	34	25	07/13/20 15:14	7/7/20	
Pentachlorophenol (PCP)	1100 U	1100	230	25	07/13/20 15:14	7/7/20	
Phenanthrene	230 U	230	31	25	07/13/20 15:14	7/7/20	
Phenol	37 J	230	23	25	07/13/20 15:14	7/7/20	
Pyrene	230 U	230	33	25	07/13/20 15:14	7/7/20	

Tentatively Identified Compounds

			Result	
CAS#	Compound Identification	RT	ug/L	Q

No Tentatively Identified Compounds Detected

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	0 *	35 - 141	07/13/20 15:14	D
2-Fluorobiphenyl	0 *	31 - 118	07/13/20 15:14	D
2-Fluorophenol	0 *	10 - 105	07/13/20 15:14	D
Nitrobenzene-d5	0 *	31 - 110	07/13/20 15:14	D
Phenol-d6	0 *	10 - 107	07/13/20 15:14	D
p-Terphenyl-d14	0 *	10 - 165	07/13/20 15:14	D

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701 Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 11:50 **Project:**

Monitoring

Date Received: 07/02/20 10:50 **Sample Matrix:** Water

Sample Name: WG-9954-070120-SG-028 Units: ug/L Lab Code: R2005701-006 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

1,2,4-Trichlorobenzene	Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1.3-Dichlorobenzene 450 U 450 50 07/13/20 16:10 7/7/20	1,2,4-Trichlorobenzene	72 J	450	51	50	07/13/20 16:10	7/7/20	
1.4-Dicklorobenzene 100 J 450 S2 50 O7/13/20 16:10 77/20 2.4.5-Tricklorophenol 450 U 450 S0 50 O7/13/20 16:10 77/20 2.4.5-Tricklorophenol 450 U 450 S0 50 O7/13/20 16:10 77/20 2.4-Dichlorophenol 320 J 450 S7 50 O7/13/20 16:10 77/20 2.4-Dimitylphenol 2300 U 2300 P10 S0 50 O7/13/20 16:10 77/20 2.4-Dinitrotoluene 450 U 450 H0 50 O7/13/20 16:10 77/20 2.4-Dinitrotoluene 450 U 450 H0 50 O7/13/20 16:10 77/20 2.6-Dinitrotoluene 450 U 450 H0 60 S0 O7/13/20 16:10 77/20 2.6-Dinitrotoluene 450 U 450 H0 60 S0 O7/13/20 16:10 77/20 2.6-Dinitrotoluene 450 U 450 H0 50 O7/13/20 16:10 77/20 2.Chlorophenol 450 U 450 H0 50 O7/13/20 16:10 77/20 2.Methylphenol 450 U 450 H0 50 O7/13/20 16:10 77/20 2.Methylphenol 450 U 450 B0 S0 S0 S0 O7/13/20 16:10	1,2-Dichlorobenzene	450 U	450	54	50	07/13/20 16:10	7/7/20	
2.4.5-Trichlorophenol	1,3-Dichlorobenzene	450 U	450	46	50	07/13/20 16:10	7/7/20	
2,4-Dichlorophenol 450 U 450 62 50 07/13/20 16:10 77/20 2,4-Dichlorophenol 450 U 450 63 50 07/13/20 16:10 77/20 2,4-Dimitrophenol 2300 U 2300 910 50 07/13/20 16:10 77/20 2,4-Dimitrophenol 2300 U 2300 910 50 07/13/20 16:10 77/20 2,4-Dimitrophenol 450 U 450 10 50 07/13/20 16:10 77/20 2,4-Dimitroplenol 450 U 450 10 50 07/13/20 16:10 77/20 2,4-Dimitroplenol 450 U 450 60 50 07/13/20 16:10 77/20 2,4-Dimitroplenol 450 U 450 62 50 07/13/20 16:10 77/20 2,4-Dimitroplenol 450 U 450 62 50 07/13/20 16:10 77/20 2,4-Dimitroplenol 450 U 450 62 50 07/13/20 16:10 77/20 2,4-Dimitroplenol 450 U 450 49 50 07/13/20 16:10 77/20 2,4-Dimitroplenol 450 U 450 49 50 07/13/20 16:10 77/20 2,4-Dimitroplenol 450 U 450 46 50 07/13/20 16:10 77/20 2,4-Dimitroplenol 450 U 450 46 50 07/13/20 16:10 77/20 2,4-Dimitroplenol 450 U 450 62 50 07/13/20 16:10 77/20 2,4-Dimitroplenol 450 U 450 68 50 07/13/20 16:10 77/20 2,4-Dimitroplenol 450 U 450 68 50 07/13/20 16:10 77/20 3,4-Dimitroplenol 450 U 450 55 50 07/13/20 16:10 77/20 3,4-Dimitroplenol 450 U 450 55 50 07/13/20 16:10 77/20 3,-Nitroanline 450 U 450 53 50 07/13/20 16:10 77/20 4,-Dimitro-2-methylphenol 2300 U 2300 870 50 07/13/20 16:10 77/20 4,-Dimitro-2-methylphenol 450 U 450 450 50 07/13/20 16:10 77/20 4,-Dimitro-2-methylphenol 450 U 450 450 50 07/13/20 16:10 77/20 4,-Dimitro-2-methylphenol 450 U 450 450 50 07/13/20 16:10 77/20 4,-Dimitro-2-methylphenol 450 U 450 68 50 07/13/20 16:10 77/20 4,-Dimitro-2-methylphenol 450 U 450 68 50 07/13/20 16:10 77/20 4,-Dimitro-2-methylphenol 450 U 450 68 50 07/13/20 16:10 77/20 4,-Dimitro-2-methylphenol 450 U 450 68 50 07/13/20 16:10 77/20 4,-Dimitro-2-methylphenol 450 U 4	1,4-Dichlorobenzene	100 J	450	52	50	07/13/20 16:10	7/7/20	
2,4-Dinchlorophenol 320 J 450 57 50 07/13/20 16:10 77/20 2,4-Dinitrophenol 2300 U 2300 910 50 07/13/20 16:10 77/20 2,4-Dinitrotoluene 450 U 450 110 50 07/13/20 16:10 77/20 2,4-Dinitrotoluene 450 U 450 110 50 07/13/20 16:10 77/20 2,6-Dinitrotoluene 450 U 450 60 50 07/13/20 16:10 77/20 2-Chlorophenol 450 U 450 62 50 07/13/20 16:10 77/20 2-Methylaphthalene 450 U 450 49 50 07/13/20 16:10 77/20 2-Methylaphthalene 450 U 450 46 50 07/13/20 16:10 77/20 2-Methylaphthalene 450 U 450 46 50 07/13/20 16:10 77/20 2-Methylaphthalene 450 U 450 68 50 07/13/20 16:10 77/20 2-Nitrophenol 450 U 450 U 55 <td>2,4,5-Trichlorophenol</td> <td>450 U</td> <td>450</td> <td>50</td> <td>50</td> <td>07/13/20 16:10</td> <td>7/7/20</td> <td></td>	2,4,5-Trichlorophenol	450 U	450	50	50	07/13/20 16:10	7/7/20	
2.4-Dimitrlyphenol 450 U 450 U 350 07/13/20 16:10 77/20 2.4-Dimitrophenol 2300 U 2300 910 50 07/13/20 16:10 77/20 2.4-Dimitrotoluene 450 U 450 III 50 07/13/20 16:10 77/20 2.6-Dimitrotoluene 450 U 450 60 50 07/13/20 16:10 77/20 2Chloronaphthalene 450 U 450 62 50 07/13/20 16:10 77/20 2Chlorophenol 450 U 450 49 50 07/13/20 16:10 77/20 2Methylphenol 450 U 450 46 50 07/13/20 16:10 77/20 2Methylphenol 450 U 450 46 50 07/13/20 16:10 77/20 2Mitrophenol 450 U 450 62 50 07/13/20 16:10 77/20 2Nitrophenol 450 U 450 68 50 07/13/20 16:10 77/20 2Nitrophenol 450 U 450 68 50 07/13/20 16:10 77/20 3and 4-Methylphenol Coelution 460 J 450 55 50 07/13/20 16:10 77/20 3bitroaniline 450 U 450 53 50 07/13/20 16:10 77/20 4Bromophenyl Phenyl Ether 450 U 450 120 50 07/13/20 16:10 77/20 4Bromophenyl Phenyl Ether 450 U 450 46 50 07/13/20 16:10 77/20 4-Chloroaniline 4	2,4,6-Trichlorophenol	450 U	450	62	50	07/13/20 16:10	7/7/20	
2.4-Dinitrophenol 2300 U 2300 D 910 S0 07/13/20 16:10 7/7/20 2.4-Dinitrotoluene 450 U 450 U 450 D 110 S0 07/13/20 16:10 7/7/20 2.6-Dinitrotoluene 450 U 450 G2 50 07/13/20 16:10 7/7/20 2-Chlorophenol 450 U 450 G2 50 07/13/20 16:10 7/7/20 2-Methylaphthalene 450 U 450 S9 S0 07/13/20 16:10 7/7/20 2-Methylaphthalene 450 U 450 S9 S0 07/13/20 16:10 7/7/20 2-Methylaphthalene 450 U 450 G2 50 07/13/20 16:10 7/7/20 2-Mitrophenol 450 U 450 G2 50 07/13/20 16:10 7/7/20 2-Nitrophenol 450 U 450 G8 S0 07/13/20 16:10 7/7/20 2-Nitrophenol 450 U 450 S5 S0 07/13/20 16:10 7/7/20 3-3r-Dichlorobenzidine 450 U 450 S5 S0 07/13/20 16:10 7/7/20 3-Nitroaniline 450 U 450 S0 07/13/20 16:10 7/7/20 4-Bromophenyl Phenyl Ether 450 U 450 U 50 07/13/20 16:10	2,4-Dichlorophenol	320 J	450	57	50	07/13/20 16:10	7/7/20	
2,4-Dinitrotoluene 450 U 450 U 450 do 50 07/13/20 16:10 7/7/20 2,6-Dinitrotoluene 450 U 450 do 50 07/13/20 16:10 7/7/20 2-Chloronaphthalene 450 U 450 do 450 do 9 50 07/13/20 16:10 7/7/20 2-Methylnaphthalene 450 U 450 do 450 do 59 50 07/13/20 16:10 7/7/20 2-Methylphenol 450 U 450 do 450 do 59 50 07/13/20 16:10 7/7/20 2-Methylphenol 450 U 450 do 62 50 07/13/20 16:10 7/7/20 2-Nitrophenol 450 U 450 do 68 50 07/13/20 16:10 7/7/20 3,3'-Dichlorobenzidine 450 U 450 55 50 07/13/20 16:10 7/7/20 3,3'-Dichlorobenzidine 450 U 450 55 50 07/13/20 16:10 7/7/20 3,3'-Dichlorobenzidine 450 U 450 55 50 07/13/20 16:10 7/7/20 4,6'Dinitro-2-nethylphenol 2300	2,4-Dimethylphenol	450 U	450	63	50	07/13/20 16:10	7/7/20	
2,6-Dinitrotoluene 450 U 450 60 50 07/13/20 16:10 7/7/20 2-Chloronaphthalene 450 U 450 62 50 07/13/20 16:10 7/7/20 2-Chlorophenol 450 U 450 49 50 07/13/20 16:10 7/7/20 2-Methylaphthalene 450 U 450 59 50 07/13/20 16:10 7/7/20 2-Methylphenol 450 U 450 46 50 07/13/20 16:10 7/7/20 2-Nitroaniline 450 U 450 62 50 07/13/20 16:10 7/7/20 2-Nitrophenol 450 U 450 68 50 07/13/20 16:10 7/7/20 3,3-Dichlorobenzidine 450 U 450 55 50 07/13/20 16:10 7/7/20 3,3-Dichlorobenzidine 450 U 450 53 50 07/13/20 16:10 7/7/20 3,3-Dichlorobenzidine 450 U 450 53 50 07/13/20 16:10 7/7/20 3,3-Dichlorobenzidine 450 U 450 120 50	2,4-Dinitrophenol	2300 U	2300	910	50	07/13/20 16:10	7/7/20	
2-Chloronaphthalene 450 U 450 decided by the company of the company o		450 U	450	110	50	07/13/20 16:10	7/7/20	
2-Chlorophenol 450 U 450 U 450 S9 50 07/13/20 16:10 77/20 2-Methylnphthalene 450 U 450 S9 50 07/13/20 16:10 77/20 2-Methylphenol 450 U 450 G2 50 07/13/20 16:10 77/20 2-Nitrophenol 450 U 450 G8 50 07/13/20 16:10 77/20 2-Nitrophenol 450 U 450 G8 50 07/13/20 16:10 77/20 3-3-Dichlorobenzidine 450 U 450 S5 55 00 07/13/20 16:10 77/20 3-and 4-Methylphenol Coclution 60 J 450 U 450 U 50 07/13/20 16:10 77/20 3-Nitroaniline 450 U 450 U 2300 U 2300 U 200 S0 870 S0 07/13/20 16:10 77/20 4-Chloro-3-methylphenol 450 U 450 T5 50 07/13/20 16:10 77/20 4-Chlorophenyl Phenyl Ether 450 U 450 G8 50 07/13/20 16:10 77/20 4-Nitrophenol 2300 U 450 G8 50 07/13/20 16:10	2,6-Dinitrotoluene	450 U	450	60	50	07/13/20 16:10	7/7/20	
2-Chlorophenol 450 U 450 U 450 S9 50 07/13/20 16:10 77/20 2-Methylnphthalene 450 U 450 S9 50 07/13/20 16:10 77/20 2-Methylphenol 450 U 450 G2 50 07/13/20 16:10 77/20 2-Nitrophenol 450 U 450 G8 50 07/13/20 16:10 77/20 2-Nitrophenol 450 U 450 G8 50 07/13/20 16:10 77/20 3-3-Dichlorobenzidine 450 U 450 S5 55 00 07/13/20 16:10 77/20 3-and 4-Methylphenol Coclution 60 J 450 U 450 U 50 07/13/20 16:10 77/20 3-Nitroaniline 450 U 450 U 2300 U 2300 U 200 S0 870 S0 07/13/20 16:10 77/20 4-Chloro-3-methylphenol 450 U 450 T5 50 07/13/20 16:10 77/20 4-Chlorophenyl Phenyl Ether 450 U 450 G8 50 07/13/20 16:10 77/20 4-Nitrophenol 2300 U 450 G8 50 07/13/20 16:10		450 U	450	62	50	07/13/20 16:10	7/7/20	
2-Methylnaphthalene 450 U 450 59 50 07/13/20 16:10 77/720 2-Methylphenol 450 U 450 46 50 07/13/20 16:10 77/720 2-Nitrophenol 450 U 450 68 50 07/13/20 16:10 77/720 3,3-Dichlorobenzidine 450 U 450 55 50 07/13/20 16:10 77/720 3,3-Dichlorobenzidine 450 U 450 55 50 07/13/20 16:10 77/720 3-and 4-Methylphenol Coelution 66 J 450 L 50 07/13/20 16:10 77/720 3-Nitroaniline 450 U 450 120 50 07/13/20 16:10 77/720 4-Bromophenyl Phenyl Ether 450 U 450 75 50 07/13/20 16:10 77/720 4-Chloroaniline 450 U 450 45 46 50 07/13/20 16:10 77/720 4-Chloroaphenyl Phenyl Ether 450 U <td< td=""><td></td><td>450 U</td><td>450</td><td>49</td><td></td><td>07/13/20 16:10</td><td>7/7/20</td><td></td></td<>		450 U	450	49		07/13/20 16:10	7/7/20	
2-Methylphenol		450 U	450	59	50	07/13/20 16:10	7/7/20	
2-Nitroaniline 450 U 450 Gel 50 07/13/20 16:10 7/7/20 2-Nitrophenol 450 U 450 Gel 50 07/13/20 16:10 7/7/20 3,3"-Dichlorobenzidine 450 U 450 Sel 55 50 07/13/20 16:10 7/7/20 3- and 4-Methylphenol Coelution 60 J 450 U 250 Sel 50 07/13/20 16:10 7/7/20 3-Nitroaniline 450 U 450 U 200 Sel 50 07/13/20 16:10 7/7/20 4,6-Dinitro-2-methylphenol 2300 U 2300 Sel 870 Sel 50 07/13/20 16:10 7/7/20 4-Bromophenyl Phenyl Ether 450 U 450 Tel 50 07/13/20 16:10 7/7/20 4-Chloro-3-methylphenol 450 U 450 del 50 07/13/20 16:10 7/7/20 4-Chloro-alinen 450 U 450 del 50 07/13/20 16:10 7/7/20 4-Chlorophenyl Phenyl Ether 450 U 450 del 50 07/13/20 16:10 7/7/20 4-Nitroaniline 450 U 450 del 50 07/13/20 16:10 7/7/20 4-Nitroaniline 450 U 450 del 50 07/13/20 16:10 7/7/20 4-Nitroaniline 450 U 450 del 50 07/13/20 16:10 7/7/20 4-Nitroaniline 450 U		450 U		46	50	07/13/20 16:10	7/7/20	
2-Nitrophenol 450 U 450 68 50 07/13/20 16:10 7/7/20 3,3'-Dichlorobenzidine 450 U 450 55 50 07/13/20 16:10 77/20 3-And 4-Methylphenol Coelution 60 J 450 U 55 50 07/13/20 16:10 77/20 3-Nitroaniline 450 U 450 120 50 07/13/20 16:10 77/20 4.6-Dinitro-2-methylphenol 2300 U 2300 870 50 07/13/20 16:10 77/20 4-Bromophenyl Phenyl Ether 450 U 450 75 50 07/13/20 16:10 77/20 4-Chloro-3-methylphenol 450 U 450 450 49 50 07/13/20 16:10 77/20 4-Chloro-3-methylphenol 450 U 450 46 50 07/13/20 16:10 77/20 4-Chloro-alline 450 U 450 46 50 07/13/20 16:10 77/20 4-Chlorophenyl Phenyl Ether 450 U 450 46 50 07/13/20 16:10 77/20 4-Chlorophenyl Phenyl Ether 450 U 450 68 50 07/13/20 16:10 77/20 4-Chlorophenyl Phenyl Ether 450 U 450 68 50 07/13/20 16:10 77/20 4-Chlorophenyl Phenyl Ether 450 U 450 68 50 07/13/20 16:10 77/20 4-Nitroaniline 450 U 450 68 50 07/13/20 16:10 77/20 4-Nitroaniline 450 U 450 61 50 07/13/20 16:10 77/20 4-Chaphthene 450 U 450 61 50 07/13/20 16:10 77/20 Acenaphthene 450 U 450 61 50 07/13/20 16:10 77/20 Acenaphthylene 450 U 450 61 50 07/13/20 16:10 77/20 Acenaphthylene 450 U 450 61 50 07/13/20 16:10 77/20 Benz(a)anthracene 450 U 450 57 50 07/13/20 16:10 77/20 Benz(a)pyrene 450 U 450 51 50 07/13/20 16:10 77/20 Benz(a)pyrene 450 U 450 51 50 07/13/20 16:10 77/20 Benz(b)fluoranthene 450 U 450 51 50 07/13/20 16:10 77/20 Benz(b)fluoranthene 450 U 450 51 50 07/13/20 16:10 77/20 Benz(b)fluoranthene 450 U 450 55 50 07/13/20 16:10 77/20 Benz(b)fluoranthene 450 U 450 51 50 07/13/20 16:10 77/20 Benz(b)fluoranthene 450 U 450 55 50 07/13/20 16:10 77/20 Benz(b)fluoranthene 450 U 450 64 50 07/13/20 16:10 77/20 Benz(b)fluoranthene 450 U 450 64 50 07/13/20 16:10 77/20 Benz(b)fluoranthene 450 U 450 64 50 07/13/20 16:10 77/20 Benz(b)fluoranthene 450 U 450 64 50 07/13/20 16:10 77/20 Benz(b)fluoranthene 450 U 450 64 50 07/13/20 16:10 77/20 Benz(b)fluoranthene 450 U 450 64 50 07/13/20 16:10 77/20 Benz(b)fluoranthene 450 U 450 64 50 07/13/20 16:10 77/20 Benz(b)fluoranthene 450 U 450 64 50 07/13/20 16:10 77/20 Benz(b)fluoranthene 450 U 450 64 50								
3,3'-Dichlorobenzidine 450 U 450 S5 50 07/13/20 16:10 77/20 3- and 4-Methylphenol Coelution 60 J 450 U 53 S0 50 07/13/20 16:10 77/20 3-Nitroaniline 450 U 450 I20 50 07/13/20 16:10 77/20 4-G-Dinitro-2-methylphenol 2300 U 2300 R70 50 07/13/20 16:10 77/20 4-Bromophenyl Phenyl Ether 450 U 450 V 450 A9 50 07/13/20 16:10 77/20 4-Chloro-3-methylphenol 450 U 450 V 450 A9 50 07/13/20 16:10 77/20 4-Chloro-almine 450 U 450 A9 50 07/13/20 16:10 77/20 4-Chlorophenyl Phenyl Ether 450 U 450 A6 50 07/13/20 16:10 77/20 4-Nitroaniline 450 U 450 A6 50 07/13/20 16:10 77/20 4-Nitroaniline 450 U 450 A6 50 07/13/20 16:10 77/20 4-Nitroaniline 450 U 450 A6 50 07/13/20 16:10 77/20 4-Nitroaniline 450 U 450 B6 50 07/13/20 16:10 77/20 Accnapht	2-Nitrophenol	450 U	450	68	50	07/13/20 16:10	7/7/20	
3- and 4-Methylphenol Coelution 60 J 450 U 450 D 53 50 O7/13/20 16:10 O7/12/0 7/7/20 3-Nitroaniline 450 U 450 U 2300 U 2300 B70 S0 O7/13/20 16:10 O7/12/0 7/7/20 4-G-Dinitro-2-methylphenol 2300 U 450 U 450 U 50 O7/13/20 16:10 O7/12/0 7/7/20 4-Bromophenyl Phenyl Ether 450 U 450 U 49 S0 O7/13/20 16:10 O7/12/0 7/7/20 4-Chloroaniline 450 U 450 U 450 U 46 S0 O7/13/20 16:10 O7/12/0 7/7/20 4-Chlorophenyl Phenyl Ether 450 U 450 G8 S0 O7/13/20 16:10 O7/12/0 7/7/20 4-Nitroaniline 450 U 450 G8 S0 O7/13/20 16:10 O7/12/0 7/7/20 4-Nitroaniline 450 U 450 G8 S0 O7/13/20 16:10 O7/12/0 7/7/20 4-Nitroaniline 450 U 450 G1 S0 O7/13/20 16:10 O7/12/0 7/7/20 4-Nitroaniline 450 U 450 G1 S0 O7/13/20 16:10 O7/13/20 16:10 O7/12/0 7/7/20 4-Nitrophenol 2300 U 2300 U 2300 O7/13/20 16:10 O7/13/20 16:10 O7/13/20 16:10 O7/12/0 7/7/20 Acenaphthylene 450 U 450 G1 S0 O7/								
3-Nitroaniline		60 J						
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Butyl Benzyl Phthalate 450 U 450 64 50 07/13/20 16:10 7/7/20								
	Chrysene	450 U	450	53	50	07/13/20 16:10	7/7/20	

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Superset Reference:20-0000555360 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 11:50

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-028
 Units: ug/L

 Lab Code:
 R2005701-006
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	450 U	450	91	50	07/13/20 16:10	7/7/20	
Di-n-octyl Phthalate	450 U	450	150	50	07/13/20 16:10	7/7/20	
Dibenz(a,h)anthracene	450 U	450	47	50	07/13/20 16:10	7/7/20	
Dibenzofuran	450 U	450	62	50	07/13/20 16:10	7/7/20	
Diethyl Phthalate	450 U	450	49	50	07/13/20 16:10	7/7/20	
Dimethyl Phthalate	450 U	450	56	50	07/13/20 16:10	7/7/20	
Fluoranthene	450 U	450	67	50	07/13/20 16:10	7/7/20	
Fluorene	450 U	450	56	50	07/13/20 16:10	7/7/20	
Hexachlorobenzene	450 U	450	70	50	07/13/20 16:10	7/7/20	
Hexachlorobutadiene	450 U	450	46	50	07/13/20 16:10	7/7/20	
Hexachlorocyclopentadiene	450 U	450	98	50	07/13/20 16:10	7/7/20	
Hexachloroethane	450 U	450	48	50	07/13/20 16:10	7/7/20	
Indeno(1,2,3-cd)pyrene	450 U	450	80	50	07/13/20 16:10	7/7/20	
Isophorone	450 U	450	62	50	07/13/20 16:10	7/7/20	
N-Nitrosodi-n-propylamine	450 U	450	52	50	07/13/20 16:10	7/7/20	
N-Nitrosodiphenylamine	450 U	450	120	50	07/13/20 16:10	7/7/20	
Naphthalene	450 U	450	54	50	07/13/20 16:10	7/7/20	
Nitrobenzene	450 U	450	67	50	07/13/20 16:10	7/7/20	
Pentachlorophenol (PCP)	2300 U	2300	450	50	07/13/20 16:10	7/7/20	
Phenanthrene	450 U	450	61	50	07/13/20 16:10	7/7/20	
Phenol	450 U	450	46	50	07/13/20 16:10	7/7/20	
Pyrene	450 U	450	65	50	07/13/20 16:10	7/7/20	

Tentatively Identified Compounds

			Result	
CAS#	Compound Identification	RT	ug/L	Q

No Tentatively Identified Compounds Detected

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	0 *	35 - 141	07/13/20 16:10	D
2-Fluorobiphenyl	0 *	31 - 118	07/13/20 16:10	D
2-Fluorophenol	0 *	10 - 105	07/13/20 16:10	D
Nitrobenzene-d5	0 *	31 - 110	07/13/20 16:10	D
Phenol-d6	0 *	10 - 107	07/13/20 16:10	D
p-Terphenyl-d14	0 *	10 - 165	07/13/20 16:10	D



Semivolatile Organic Compounds by GC

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701 Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 11:50 **Project:**

Monitoring

Date Received: 07/02/20 10:50 **Sample Matrix:** Water

Sample Name: WG-9954-070120-SG-027 Units: ug/L Lab Code: R2005701-001 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.45 U	0.45	0.19	10	07/10/20 15:00	7/7/20	_
4,4'-DDE	0.45 U	0.45	0.19	10	07/10/20 15:00	7/7/20	
4,4'-DDT	0.45 U	0.45	0.19	10	07/10/20 15:00	7/7/20	
Aldrin	1.0 P	0.45	0.19	10	07/10/20 15:00	7/7/20	
Dieldrin	0.45 U	0.45	0.19	10	07/10/20 15:00	7/7/20	
Endosulfan I	0.45 U	0.45	0.19	10	07/10/20 15:00	7/7/20	
Endosulfan II	0.45 U	0.45	0.19	10	07/10/20 15:00	7/7/20	
Endosulfan Sulfate	0.45 U	0.45	0.19	10	07/10/20 15:00	7/7/20	
Endrin	0.45 U	0.45	0.19	10	07/10/20 15:00	7/7/20	
Endrin Ketone	0.45 U	0.45	0.19	10	07/10/20 15:00	7/7/20	
Heptachlor	0.45 U	0.45	0.19	10	07/10/20 15:00	7/7/20	
Heptachlor Epoxide	0.45 U	0.45	0.19	10	07/10/20 15:00	7/7/20	
Methoxychlor	0.45 U	0.45	0.19	10	07/10/20 15:00	7/7/20	
Toxaphene	4.6 U	4.6	4.6	10	07/10/20 15:00	7/7/20	
alpha-BHC	27	0.45	0.19	10	07/10/20 15:00	7/7/20	
alpha-Chlordane	0.45 U	0.45	0.19	10	07/10/20 15:00	7/7/20	
beta-BHC	6.8 P	0.45	0.19	10	07/10/20 15:00	7/7/20	
delta-BHC	9.2	0.45	0.19	10	07/10/20 15:00	7/7/20	
gamma-BHC (Lindane)	6.5	0.45	0.19	10	07/10/20 15:00	7/7/20	
gamma-Chlordane	0.45 U	0.45	0.19	10	07/10/20 15:00	7/7/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q	
Decachlorobiphenyl	14	10 - 164	07/10/20 15:00		
Tetrachloro-m-xylene	171 *	10 - 147	07/10/20 15:00	*	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 10:45

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-026
 Units: ug/L

 Lab Code:
 R2005701-002
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/08/20 15:51	7/7/20	_
4,4'-DDE	0.045 U	0.045	0.019	1	07/08/20 15:51	7/7/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/08/20 15:51	7/7/20	
Aldrin	0.045 U	0.045	0.019	1	07/08/20 15:51	7/7/20	
Dieldrin	0.045 U	0.045	0.019	1	07/08/20 15:51	7/7/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/08/20 15:51	7/7/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/08/20 15:51	7/7/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/08/20 15:51	7/7/20	
Endrin	0.045 U	0.045	0.019	1	07/08/20 15:51	7/7/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/08/20 15:51	7/7/20	
Heptachlor	0.045 U	0.045	0.019	1	07/08/20 15:51	7/7/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/08/20 15:51	7/7/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/08/20 15:51	7/7/20	
Toxaphene	0.46 U	0.46	0.46	1	07/08/20 15:51	7/7/20	
alpha-BHC	0.27	0.045	0.019	1	07/08/20 15:51	7/7/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/08/20 15:51	7/7/20	
beta-BHC	0.040 J	0.045	0.019	1	07/08/20 15:51	7/7/20	
delta-BHC	0.17	0.045	0.019	1	07/08/20 15:51	7/7/20	
gamma-BHC (Lindane)	0.21	0.045	0.019	1	07/08/20 15:51	7/7/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/08/20 15:51	7/7/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	23	10 - 164	07/08/20 15:51	
Tetrachloro-m-xylene	62	10 - 147	07/08/20 15:51	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701 **Date Collected:** 07/01/20 09:55

Love Canal:292-402-D02-3100/9954 Annual Long Term **Project:**

Monitoring

Date Received: 07/02/20 10:50 **Sample Matrix:** Water

Sample Name: WG-9954-070120-SG-025 Units: ug/L Lab Code: R2005701-003 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/08/20 16:10	7/7/20	
4,4'-DDE	0.045 U	0.045	0.019	1	07/08/20 16:10	7/7/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/08/20 16:10	7/7/20	
Aldrin	0.045 U	0.045	0.019	1	07/08/20 16:10	7/7/20	
Dieldrin	0.045 U	0.045	0.019	1	07/08/20 16:10	7/7/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/08/20 16:10	7/7/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/08/20 16:10	7/7/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/08/20 16:10	7/7/20	
Endrin	0.045 U	0.045	0.019	1	07/08/20 16:10	7/7/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/08/20 16:10	7/7/20	
Heptachlor	0.045 U	0.045	0.019	1	07/08/20 16:10	7/7/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/08/20 16:10	7/7/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/08/20 16:10	7/7/20	
Toxaphene	0.46 U	0.46	0.46	1	07/08/20 16:10	7/7/20	
alpha-BHC	0.13	0.045	0.019	1	07/08/20 16:10	7/7/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/08/20 16:10	7/7/20	
beta-BHC	0.045 U	0.045	0.019	1	07/08/20 16:10	7/7/20	
delta-BHC	0.029 J	0.045	0.019	1	07/08/20 16:10	7/7/20	
gamma-BHC (Lindane)	0.076	0.045	0.019	1	07/08/20 16:10	7/7/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/08/20 16:10	7/7/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	17	10 - 164	07/08/20 16:10	
Tetrachloro-m-xylene	56	10 - 147	07/08/20 16:10	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701 **Date Collected:** 07/01/20 08:45

Love Canal:292-402-D02-3100/9954 Annual Long Term **Project:**

Monitoring

Date Received: 07/02/20 10:50 **Sample Matrix:** Water

Sample Name: WG-9954-070120-SG-024 Units: ug/L Lab Code: R2005701-004 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/08/20 16:29	7/7/20	
4,4'-DDE	0.045 U	0.045	0.019	1	07/08/20 16:29	7/7/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/08/20 16:29	7/7/20	
Aldrin	0.045 U	0.045	0.019	1	07/08/20 16:29	7/7/20	
Dieldrin	0.045 U	0.045	0.019	1	07/08/20 16:29	7/7/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/08/20 16:29	7/7/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/08/20 16:29	7/7/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/08/20 16:29	7/7/20	
Endrin	0.045 U	0.045	0.019	1	07/08/20 16:29	7/7/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/08/20 16:29	7/7/20	
Heptachlor	0.045 U	0.045	0.019	1	07/08/20 16:29	7/7/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/08/20 16:29	7/7/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/08/20 16:29	7/7/20	
Toxaphene	0.46 U	0.46	0.46	1	07/08/20 16:29	7/7/20	
alpha-BHC	0.045 U	0.045	0.019	1	07/08/20 16:29	7/7/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/08/20 16:29	7/7/20	
beta-BHC	0.045 U	0.045	0.019	1	07/08/20 16:29	7/7/20	
delta-BHC	0.045 U	0.045	0.019	1	07/08/20 16:29	7/7/20	
gamma-BHC (Lindane)	0.045 U	0.045	0.019	1	07/08/20 16:29	7/7/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/08/20 16:29	7/7/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	34	10 - 164	07/08/20 16:29	
Tetrachloro-m-xylene	52	10 - 147	07/08/20 16:29	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701 **Project:**

Love Canal:292-402-D02-3100/9954 Annual Long Term

Date Collected: 07/01/20 11:50

Monitoring **Sample Matrix:** Water

Lab Code:

Date Received: 07/02/20 10:50

Sample Name: WG-9954-070120-SG-028

Units: ug/L R2005701-006 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.45 U	0.45	0.19	10	07/10/20 15:19	7/7/20	_
4,4'-DDE	0.45 U	0.45	0.19	10	07/10/20 15:19	7/7/20	
4,4'-DDT	0.45 U	0.45	0.19	10	07/10/20 15:19	7/7/20	
Aldrin	0.85	0.45	0.19	10	07/10/20 15:19	7/7/20	
Dieldrin	0.45 U	0.45	0.19	10	07/10/20 15:19	7/7/20	
Endosulfan I	0.45 U	0.45	0.19	10	07/10/20 15:19	7/7/20	
Endosulfan II	0.45 U	0.45	0.19	10	07/10/20 15:19	7/7/20	
Endosulfan Sulfate	0.45 U	0.45	0.19	10	07/10/20 15:19	7/7/20	
Endrin	0.45 U	0.45	0.19	10	07/10/20 15:19	7/7/20	
Endrin Ketone	0.45 U	0.45	0.19	10	07/10/20 15:19	7/7/20	
Heptachlor	0.45 U	0.45	0.19	10	07/10/20 15:19	7/7/20	
Heptachlor Epoxide	0.45 U	0.45	0.19	10	07/10/20 15:19	7/7/20	
Methoxychlor	0.45 U	0.45	0.19	10	07/10/20 15:19	7/7/20	
Toxaphene	4.6 U	4.6	4.6	10	07/10/20 15:19	7/7/20	
alpha-BHC	25	0.45	0.19	10	07/10/20 15:19	7/7/20	
alpha-Chlordane	0.45 U	0.45	0.19	10	07/10/20 15:19	7/7/20	
beta-BHC	6.9	0.45	0.19	10	07/10/20 15:19	7/7/20	
delta-BHC	8.7	0.45	0.19	10	07/10/20 15:19	7/7/20	
gamma-BHC (Lindane)	6.2	0.45	0.19	10	07/10/20 15:19	7/7/20	
gamma-Chlordane	0.45 U	0.45	0.19	10	07/10/20 15:19	7/7/20	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 11:50

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-028
 Units: ug/L

 Lab Code:
 R2005701-006
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analysis Method: 8081B **Prep Method:** EPA 3510C

 Surrogate Name
 % Rec
 Control Limits
 Date Analyzed
 Q

 Decachlorobiphenyl
 23
 10 - 164
 07/10/20 15:19

 Tetrachloro-m-xylene
 154 *
 10 - 147
 07/10/20 15:19
 *

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 11:50

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-028
 Units: ug/L

 Lab Code:
 R2005701-006
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.45 U	0.45	0.19	10	07/24/20 23:43	7/14/20	*
4,4'-DDE	0.45 U	0.45	0.19	10	07/24/20 23:43	7/14/20	*
4,4'-DDT	0.45 U	0.45	0.19	10	07/24/20 23:43	7/14/20	*
Aldrin	0.66	0.45	0.19	10	07/24/20 23:43	7/14/20	*
Dieldrin	0.45 U	0.45	0.19	10	07/24/20 23:43	7/14/20	*
Endosulfan I	0.45 U	0.45	0.19	10	07/24/20 23:43	7/14/20	*
Endosulfan II	0.45 U	0.45	0.19	10	07/24/20 23:43	7/14/20	*
Endosulfan Sulfate	0.45 U	0.45	0.19	10	07/24/20 23:43	7/14/20	*
Endrin	0.45 U	0.45	0.19	10	07/24/20 23:43	7/14/20	*
Endrin Ketone	0.45 U	0.45	0.19	10	07/24/20 23:43	7/14/20	*
Heptachlor	0.40 J	0.45	0.19	10	07/24/20 23:43	7/14/20	*
Heptachlor Epoxide	0.45 U	0.45	0.19	10	07/24/20 23:43	7/14/20	*
Methoxychlor	0.45 U	0.45	0.19	10	07/24/20 23:43	7/14/20	*
Toxaphene	4.6 U	4.6	4.6	10	07/24/20 23:43	7/14/20	*
alpha-BHC	18	0.45	0.19	10	07/24/20 23:43	7/14/20	*
alpha-Chlordane	0.45 U	0.45	0.19	10	07/24/20 23:43	7/14/20	*
beta-BHC	4.4	0.45	0.19	10	07/24/20 23:43	7/14/20	*
delta-BHC	6.3	0.45	0.19	10	07/24/20 23:43	7/14/20	*
gamma-BHC (Lindane)	3.6	0.45	0.19	10	07/24/20 23:43	7/14/20	*
gamma-Chlordane	1.5	0.45	0.19	10	07/24/20 23:43	7/14/20	*

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 11:50

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-028
 Units: ug/L

 Lab Code:
 R2005701-006
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analysis Method: 8081B

Prep Method: EPA 3510C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	36	10 - 164	07/24/20 23:43	
Tetrachloro-m-xylene	317 *	10 - 147	07/24/20 23:43	*

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 11:50

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-027
 Units: ug/L

 Lab Code:
 R2005701-001
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/08/20 15:36	7/7/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/08/20 15:36	7/7/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/08/20 15:36	7/7/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/08/20 15:36	7/7/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/08/20 15:36	7/7/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/08/20 15:36	7/7/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/08/20 15:36	7/7/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	22	10 - 152	07/08/20 15:36	
Tetrachloro-m-xvlene	50	14 - 129	07/08/20 15:36	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 10:45

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-026
 Units: ug/L

 Lab Code:
 R2005701-002
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/10/20 11:50	7/7/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/10/20 11:50	7/7/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/10/20 11:50	7/7/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/10/20 11:50	7/7/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/10/20 11:50	7/7/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/10/20 11:50	7/7/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/10/20 11:50	7/7/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	43	10 - 152	07/10/20 11:50	
Tetrachloro-m-xylene	42	14 - 129	07/10/20 11:50	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 09:55

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-025
 Units: ug/L

 Lab Code:
 R2005701-003
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/10/20 12:10	7/7/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/10/20 12:10	7/7/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/10/20 12:10	7/7/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/10/20 12:10	7/7/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/10/20 12:10	7/7/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/10/20 12:10	7/7/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/10/20 12:10	7/7/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	18	10 - 152	07/10/20 12:10	
Tetrachloro-m-xylene	39	14 - 129	07/10/20 12:10	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 08:45

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-024
 Units: ug/L

 Lab Code:
 R2005701-004
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/10/20 13:09	7/7/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/10/20 13:09	7/7/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/10/20 13:09	7/7/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/10/20 13:09	7/7/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/10/20 13:09	7/7/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/10/20 13:09	7/7/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/10/20 13:09	7/7/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	32	10 - 152	07/10/20 13:09	
Tetrachloro-m-xylene	34	14 - 129	07/10/20 13:09	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/01/20 11:50

Monitoring

Sample Matrix: Water Date Received: 07/02/20 10:50

 Sample Name:
 WG-9954-070120-SG-028
 Units: ug/L

 Lab Code:
 R2005701-006
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/08/20 17:35	7/7/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/08/20 17:35	7/7/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/08/20 17:35	7/7/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/08/20 17:35	7/7/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/08/20 17:35	7/7/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/08/20 17:35	7/7/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/08/20 17:35	7/7/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	24	10 - 152	07/08/20 17:35	
Tetrachloro-m-xvlene	57	14 - 129	07/08/20 17:35	



QC Summary Forms

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com



Volatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request: R2005701 Love Canal:292-402-D02-3100/9954 Annual Long Term

Project:

Sample Matrix: Water

> SURROGATE RECOVERY SUMMARY Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

		4-Bromofluorobenzene	Dibromofluoromethane	Toluene-d8
Sample Name	Lab Code	85-122	89-119	87-121
WG-9954-070120-SG-027 DL	R2005701-001	96	96	100
WG-9954-070120-SG-026	R2005701-002	95	105	102
WG-9954-070120-SG-025	R2005701-003	98	100	103
WG-9954-070120-SG-024	R2005701-004	96	100	102
WG-9954-070120-SG-028 DL	R2005701-006	96	96	101
Method Blank	RQ2007501-04	95	99	101
Method Blank	RQ2007565-04	94	99	100
Lab Control Sample	RQ2007501-03	103	103	105
Lab Control Sample	RQ2007565-03	98	99	101
WG-9954-070120-SG-027	R2005701-001	114	89	99
TB-9954-070120-SG-005	R2005701-005	91	98	100
WG-9954-070120-SG-028	R2005701-006	113	99	108
Method Blank	RQ2007451-04	95	99	102
Lab Control Sample	RQ2007451-03	99	100	101

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007451-04
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/13/20 19:02	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/13/20 19:02	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/13/20 19:02	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/13/20 19:02	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/13/20 19:02	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/13/20 19:02	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/13/20 19:02	
2-Butanone (MEK)	10 U	10	0.78	1	07/13/20 19:02	
2-Hexanone	10 U	10	0.20	1	07/13/20 19:02	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/13/20 19:02	
Acetone	10 U	10	5.0	1	07/13/20 19:02	
Benzene	5.0 U	5.0	0.20	1	07/13/20 19:02	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/13/20 19:02	
Bromoform	5.0 U	5.0	0.25	1	07/13/20 19:02	
Bromomethane	5.0 U	5.0	0.70	1	07/13/20 19:02	
Carbon Disulfide	10 U	10	0.42	1	07/13/20 19:02	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/13/20 19:02	
Chlorobenzene	5.0 U	5.0	0.20	1	07/13/20 19:02	
Chloroethane	5.0 U	5.0	0.23	1	07/13/20 19:02	
Chloroform	5.0 U	5.0	0.24	1	07/13/20 19:02	
Chloromethane	0.32 J	5.0	0.28	1	07/13/20 19:02	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/13/20 19:02	
Dichloromethane	5.0 U	5.0	0.65	1	07/13/20 19:02	
Ethylbenzene	5.0 U	5.0	0.20	1	07/13/20 19:02	
Styrene	5.0 U	5.0	0.20	1	07/13/20 19:02	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/13/20 19:02	
Toluene	5.0 U	5.0	0.20	1	07/13/20 19:02	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/13/20 19:02	
Vinyl Acetate	10 U	10	1.1	1	07/13/20 19:02	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/13/20 19:02	
Xylenes, Total	5.0 U	5.0	0.23	1	07/13/20 19:02	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/13/20 19:02	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/13/20 19:02	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/13/20 19:02	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/13/20 19:02	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA **Project:**

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name: Method Blank Units: ug/L Basis: NA Lab Code: RQ2007451-04

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	95	85 - 122	07/13/20 19:02	
Dibromofluoromethane	99	89 - 119	07/13/20 19:02	
Toluene-d8	102	87 - 121	07/13/20 19:02	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007451-04Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Tentatively Identified Compounds

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007501-04
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/14/20 14:25	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/14/20 14:25	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/14/20 14:25	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/14/20 14:25	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/14/20 14:25	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/14/20 14:25	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/14/20 14:25	
2-Butanone (MEK)	10 U	10	0.78	1	07/14/20 14:25	
2-Hexanone	10 U	10	0.20	1	07/14/20 14:25	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/14/20 14:25	
Acetone	10 U	10	5.0	1	07/14/20 14:25	
Benzene	5.0 U	5.0	0.20	1	07/14/20 14:25	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/14/20 14:25	
Bromoform	5.0 U	5.0	0.25	1	07/14/20 14:25	
Bromomethane	5.0 U	5.0	0.70	1	07/14/20 14:25	
Carbon Disulfide	10 U	10	0.42	1	07/14/20 14:25	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/14/20 14:25	
Chlorobenzene	5.0 U	5.0	0.20	1	07/14/20 14:25	
Chloroethane	5.0 U	5.0	0.23	1	07/14/20 14:25	
Chloroform	5.0 U	5.0	0.24	1	07/14/20 14:25	
Chloromethane	5.0 U	5.0	0.28	1	07/14/20 14:25	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/14/20 14:25	
Dichloromethane	5.0 U	5.0	0.65	1	07/14/20 14:25	
Ethylbenzene	5.0 U	5.0	0.20	1	07/14/20 14:25	
Styrene	5.0 U	5.0	0.20	1	07/14/20 14:25	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/14/20 14:25	
Toluene	0.41 J	5.0	0.20	1	07/14/20 14:25	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/14/20 14:25	
Vinyl Acetate	10 U	10	1.1	1	07/14/20 14:25	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/14/20 14:25	
Xylenes, Total	5.0 U	5.0	0.23	1	07/14/20 14:25	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/14/20 14:25	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/14/20 14:25	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/14/20 14:25	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/14/20 14:25	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007501-04Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	95	85 - 122	07/14/20 14:25	
Dibromofluoromethane	99	89 - 119	07/14/20 14:25	
Toluene-d8	101	87 - 121	07/14/20 14:25	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007501-04Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown	1.61	8.5	J	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007565-04
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/15/20 13:34	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/15/20 13:34	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/15/20 13:34	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/15/20 13:34	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/15/20 13:34	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/15/20 13:34	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/15/20 13:34	
2-Butanone (MEK)	10 U	10	0.78	1	07/15/20 13:34	
2-Hexanone	10 U	10	0.20	1	07/15/20 13:34	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/15/20 13:34	
Acetone	10 U	10	5.0	1	07/15/20 13:34	
Benzene	5.0 U	5.0	0.20	1	07/15/20 13:34	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/15/20 13:34	
Bromoform	5.0 U	5.0	0.25	1	07/15/20 13:34	
Bromomethane	5.0 U	5.0	0.70	1	07/15/20 13:34	
Carbon Disulfide	10 U	10	0.42	1	07/15/20 13:34	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/15/20 13:34	
Chlorobenzene	5.0 U	5.0	0.20	1	07/15/20 13:34	
Chloroethane	5.0 U	5.0	0.23	1	07/15/20 13:34	
Chloroform	5.0 U	5.0	0.24	1	07/15/20 13:34	
Chloromethane	5.0 U	5.0	0.28	1	07/15/20 13:34	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/15/20 13:34	
Dichloromethane	5.0 U	5.0	0.65	1	07/15/20 13:34	
Ethylbenzene	5.0 U	5.0	0.20	1	07/15/20 13:34	
Styrene	5.0 U	5.0	0.20	1	07/15/20 13:34	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/15/20 13:34	
Toluene	5.0 U	5.0	0.20	1	07/15/20 13:34	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/15/20 13:34	
Vinyl Acetate	10 U	10	1.1	1	07/15/20 13:34	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/15/20 13:34	
Xylenes, Total	5.0 U	5.0	0.23	1	07/15/20 13:34	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/15/20 13:34	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/15/20 13:34	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/15/20 13:34	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/15/20 13:34	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007565-04Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	94	85 - 122	07/15/20 13:34	
Dibromofluoromethane	99	89 - 119	07/15/20 13:34	
Toluene-d8	100	87 - 121	07/15/20 13:34	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name: Method Blank Units: ug/L

Lab Code: RQ2007565-04 **Basis:** NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Project:

Tentatively Identified Compounds

Result

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/13/20

Sample Matrix: Water

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005701

Lab Control Sample

RQ2007451-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
1,1,1-Trichloroethane (TCA)	8260C	21.0	20.0	105	75-125
1,1,2,2-Tetrachloroethane	8260C	23.6	20.0	118	78-126
1,1,2-Trichloroethane	8260C	21.4	20.0	107	82-121
1,1-Dichloroethane (1,1-DCA)	8260C	20.6	20.0	103	80-124
1,1-Dichloroethene (1,1-DCE)	8260C	24.0	20.0	120 *	71-118
1,2-Dichloroethane	8260C	20.8	20.0	104	71-127
1,2-Dichloropropane	8260C	21.3	20.0	106	80-119
2-Butanone (MEK)	8260C	23.3	20.0	116	61-137
2-Hexanone	8260C	20.5	20.0	102	63-124
4-Methyl-2-pentanone	8260C	21.0	20.0	105	66-124
Acetone	8260C	23.4	20.0	117	40-161
Benzene	8260C	21.2	20.0	106	79-119
Bromodichloromethane	8260C	21.1	20.0	105	81-123
Bromoform	8260C	21.4	20.0	107	65-146
Bromomethane	8260C	23.4	20.0	117	42-166
Carbon Disulfide	8260C	21.0	20.0	105	66-128
Carbon Tetrachloride	8260C	21.5	20.0	108	70-127
Chlorobenzene	8260C	21.1	20.0	106	80-121
Chloroethane	8260C	18.6	20.0	93	62-131
Chloroform	8260C	20.5	20.0	103	79-120
Chloromethane	8260C	22.9	20.0	115	65-135
Dibromochloromethane	8260C	22.6	20.0	113	72-128
Dichloromethane	8260C	20.3	20.0	102	73-122
Ethylbenzene	8260C	20.7	20.0	103	76-120
Styrene	8260C	21.1	20.0	106	80-124
Tetrachloroethene (PCE)	8260C	19.1	20.0	95	72-125
Toluene	8260C	21.4	20.0	107	79-119
Trichloroethene (TCE)	8260C	19.6	20.0	98	74-122
Vinyl Acetate	8260C	30.4	20.0	152	52-174
Vinyl Chloride	8260C	22.9	20.0	115	74-159
cis-1,2-Dichloroethene	8260C	21.2	20.0	106	80-121
cis-1,3-Dichloropropene	8260C	21.0	20.0	105	77-122
trans-1,2-Dichloroethene	8260C	23.2	20.0	116	73-118
Printed 7/30/2020 4:05:55 PM			Supers	et Reference:20-000	00555360 rev 00

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix:

Project:

Water

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

> Units:ug/L Basis:NA

Service Request: R2005701

Date Analyzed: 07/13/20

Lab Control Sample

RQ2007451-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
trans-1 3-Dichloropropene	8260C	20.5	20.0	103	71-133

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring **Date Analyzed:** 07/14/20

Sample Matrix: Water

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005701

Lab Control Sample

RQ2007501-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
1,1,1-Trichloroethane (TCA)	8260C	20.1	20.0	100	75-125
1,1,2,2-Tetrachloroethane	8260C	22.4	20.0	112	78-126
1,1,2-Trichloroethane	8260C	21.9	20.0	110	82-121
1,1-Dichloroethane (1,1-DCA)	8260C	20.0	20.0	100	80-124
1,1-Dichloroethene (1,1-DCE)	8260C	22.5	20.0	112	71-118
1,2-Dichloroethane	8260C	19.9	20.0	99	71-127
1,2-Dichloropropane	8260C	19.8	20.0	99	80-119
2-Butanone (MEK)	8260C	22.7	20.0	114	61-137
2-Hexanone	8260C	19.9	20.0	99	63-124
4-Methyl-2-pentanone	8260C	20.8	20.0	104	66-124
Acetone	8260C	21.8	20.0	109	40-161
Benzene	8260C	20.3	20.0	101	79-119
Bromodichloromethane	8260C	21.0	20.0	105	81-123
Bromoform	8260C	19.8	20.0	99	65-146
Bromomethane	8260C	19.2	20.0	96	42-166
Carbon Disulfide	8260C	21.2	20.0	106	66-128
Carbon Tetrachloride	8260C	21.0	20.0	105	70-127
Chlorobenzene	8260C	20.7	20.0	104	80-121
Chloroethane	8260C	18.2	20.0	91	62-131
Chloroform	8260C	20.1	20.0	101	79-120
Chloromethane	8260C	19.9	20.0	100	65-135
Dibromochloromethane	8260C	22.1	20.0	111	72-128
Dichloromethane	8260C	19.2	20.0	96	73-122
Ethylbenzene	8260C	21.4	20.0	107	76-120
Styrene	8260C	20.7	20.0	103	80-124
Tetrachloroethene (PCE)	8260C	19.9	20.0	100	72-125
Toluene	8260C	21.2	20.0	106	79-119
Trichloroethene (TCE)	8260C	18.0	20.0	90	74-122
Vinyl Acetate	8260C	35.0	20.0	175 *	52-174
Vinyl Chloride	8260C	18.8	20.0	94	74-159
cis-1,2-Dichloroethene	8260C	19.8	20.0	99	80-121
cis-1,3-Dichloropropene	8260C	20.6	20.0	103	77-122
trans-1,2-Dichloroethene	8260C	22.7	20.0	114	73-118
Printed 7/30/2020 4:05:55 PM			Supers	et Reference:20-000	00555360 rev 00

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix:

Project:

Water

Service Request: R2005701

Date Analyzed: 07/14/20

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

> Units:ug/L Basis:NA

Lab Control Sample

RQ2007501-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
trans-1 3-Dichloropropene	8260C	19 9	20.0	100	71-133

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring **Date Analyzed:** 07/15/20

Sample Matrix: Water

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005701

Lab Control Sample

RQ2007565-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
1,1,1-Trichloroethane (TCA)	8260C	20.1	20.0	101	75-125
1,1,2,2-Tetrachloroethane	8260C	20.4	20.0	102	78-126
1,1,2-Trichloroethane	8260C	19.5	20.0	98	82-121
1,1-Dichloroethane (1,1-DCA)	8260C	20.4	20.0	102	80-124
1,1-Dichloroethene (1,1-DCE)	8260C	23.2	20.0	116	71-118
1,2-Dichloroethane	8260C	18.9	20.0	95	71-127
1,2-Dichloropropane	8260C	20.6	20.0	103	80-119
2-Butanone (MEK)	8260C	21.0	20.0	105	61-137
2-Hexanone	8260C	18.7	20.0	93	63-124
4-Methyl-2-pentanone	8260C	20.0	20.0	100	66-124
Acetone	8260C	20.0	20.0	100	40-161
Benzene	8260C	20.7	20.0	103	79-119
Bromodichloromethane	8260C	19.6	20.0	98	81-123
Bromoform	8260C	19.9	20.0	99	65-146
Bromomethane	8260C	15.4	20.0	77	42-166
Carbon Disulfide	8260C	21.2	20.0	106	66-128
Carbon Tetrachloride	8260C	21.1	20.0	105	70-127
Chlorobenzene	8260C	20.0	20.0	100	80-121
Chloroethane	8260C	21.4	20.0	107	62-131
Chloroform	8260C	20.0	20.0	100	79-120
Chloromethane	8260C	24.7	20.0	124	65-135
Dibromochloromethane	8260C	21.2	20.0	106	72-128
Dichloromethane	8260C	19.8	20.0	99	73-122
Ethylbenzene	8260C	20.5	20.0	103	76-120
Styrene	8260C	20.5	20.0	102	80-124
Tetrachloroethene (PCE)	8260C	20.0	20.0	100	72-125
Toluene	8260C	20.6	20.0	103	79-119
Trichloroethene (TCE)	8260C	18.7	20.0	93	74-122
Vinyl Acetate	8260C	29.2	20.0	146	52-174
Vinyl Chloride	8260C	23.2	20.0	116	74-159
cis-1,2-Dichloroethene	8260C	19.5	20.0	98	80-121
cis-1,3-Dichloropropene	8260C	19.4	20.0	97	77-122
trans-1,2-Dichloroethene	8260C	22.4	20.0	112	73-118
Printed 7/30/2020 4:05:56 PM			Supers	et Reference:20-000	00555360 rev 00

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix:

Project:

Water

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

> Units:ug/L Basis:NA

Service Request: R2005701

Date Analyzed: 07/15/20

Lab Control Sample

RQ2007565-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
trans-1.3-Dichloropropene	8260C	19 4	20.0	97	71-133



Semivolatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

QA/QC Report

Service Request: R2005701

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

Project:

SURROGATE RECOVERY SUMMARY Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Extraction Method:** EPA 3510C

		2,4,6-Tribromophenol	2-Fluorobiphenyl	2-Fluorophenol
Sample Name	Lab Code	35-141	31-118	10-105
WG-9954-070120-SG-027	R2005701-001	80	59	29
WG-9954-070120-SG-027 DL	R2005701-001	0*	0*	0*
WG-9954-070120-SG-027 DL	R2005701-001	0*	0*	0*
WG-9954-070120-SG-026	R2005701-002	111	75	45
WG-9954-070120-SG-025	R2005701-003	93	63	40
WG-9954-070120-SG-024	R2005701-004	80	59	35
WG-9954-070120-SG-028	R2005701-006	66	57	32
WG-9954-070120-SG-028 DL	R2005701-006	0*	0*	0*
WG-9954-070120-SG-028 DL	R2005701-006	0*	0*	0*
Method Blank	RQ2007188-01	84	68	48
Lab Control Sample	RQ2007188-02	103	79	52
Duplicate Lab Control Sample	RQ2007188-03	97	86	45
WG-9954-070120-SG-027 MS	RQ2007188-04	67	51	35
WG-9954-070120-SG-027 DMS	RQ2007188-05	71	49	30

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

SURROGATE RECOVERY SUMMARY Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Extraction Method:** EPA 3510C

		Nitrobenzene-d5	Phenol-d6	p-Terphenyl-d14
Sample Name	Lab Code	31-110	10-107	10-165
WG-9954-070120-SG-027	R2005701-001	86	24	51
WG-9954-070120-SG-027 DL	R2005701-001	0*	0*	0*
WG-9954-070120-SG-027 DL	R2005701-001	0*	0*	0*
WG-9954-070120-SG-026	R2005701-002	70	33	47
WG-9954-070120-SG-025	R2005701-003	62	25	51
WG-9954-070120-SG-024	R2005701-004	56	26	65
WG-9954-070120-SG-028	R2005701-006	95	27	46
WG-9954-070120-SG-028 DL	R2005701-006	0*	0*	0*
WG-9954-070120-SG-028 DL	R2005701-006	0*	0*	0*
Method Blank	RQ2007188-01	75	35	67
Lab Control Sample	RQ2007188-02	83	39	71
Duplicate Lab Control Sample	RQ2007188-03	82	35	73
WG-9954-070120-SG-027 MS	RQ2007188-04	94	27	39
WG-9954-070120-SG-027 DMS	RQ2007188-05	82	23	37

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Date Collected: 07/01/20 **Date Received:** 07/02/20

Sample Matrix: Water

07/02/20 07/9/20

Date Analyzed:
Date Extracted:

Duplicate Matrix Spike

72.7

72.7

72.7

50

54

47

46-126

44-114

41-127

<1

2

<1

30

30

30

07/7/20

Duplicate Matrix Spike Summary Semivolatile Organic Compounds by GC/MS

Matrix Spike

Sample Name: WG-9954-070120-SG-027

Units: ug/L

Lab Code: R2005701-001

Basis: NA

Analysis Method: 8270D **Prep Method:** EPA 3510C

		RQ2	2007188-04		R	Q2007188-0)5			
	Sample		Spike			Spike		% Rec		RPD
Analyte Name	Result	Result	Amount	% Rec	Result	Amount	% Rec	Limits	RPD	Limit
1,2,4-Trichlorobenzene	93	148	72.7	75	141	72.7	66	10-127	13	30
1,2-Dichlorobenzene	25	66.6	72.7	57	63.4	72.7	52	17-105	9	30
1,3-Dichlorobenzene	3.6 J	36.5	72.7	45	33.6	72.7	41	21-99	9	30
1,4-Dichlorobenzene	73	130	72.7	77	125	72.7	71	10-124	8	30
2,4,5-Trichlorophenol	37	65.3	72.7	39 *	73.6	72.7	51	48-134	27	30
2,4,6-Trichlorophenol	9.1 U	42.2	72.7	58	41.5	72.7	57	44-135	2	30
2,4-Dichlorophenol	550 E	504 E	72.7	-62 #	514 E	72.7	-48 #	40-130	NC	30
2,4-Dimethylphenol	8.9 J	66.5	72.7	79	64.6	72.7	77	42-121	3	30
2,4-Dinitrophenol	45 U	67.1	72.7	92	66.8	72.7	92	21-168	<1	30
2,4-Dinitrotoluene	9.1 U	48.6	72.7	67	51.7	72.7	71	37-143	6	30
2,6-Dinitrotoluene	9.1 U	42.5	72.7	58	46.4	72.7	64	39-136	10	30
2-Chloronaphthalene	9.1 U	29.6	72.7	41	28.6	72.7	39 *	40-108	5	30
2-Chlorophenol	27	61.5	72.7	48	58.8	72.7	44	37-112	9	30
2-Methylnaphthalene	1.4 J	56.3	72.7	75	54.3	72.7	73	34-102	3	30
2-Methylphenol	32	63.8	72.7	44	60.1	72.7	38	37-102	15	30
2-Nitroaniline	9.1 U	57.2	72.7	79	57.8	72.7	79	40-136	<1	30
2-Nitrophenol	9.1 U	65.1	72.7	90	60.9	72.7	84	27-143	7	30
3,3'-Dichlorobenzidine	9.1 U	2.48 J	72.7	3 *	3.08 J	72.7	4 *	11-131	29	30
3- and 4-Methylphenol Coelution	71	95.5	72.7	33	89.7	72.7	25 *	30-95	28	30
3-Nitroaniline	9.1 U	28.4	72.7	39	30.1	72.7	41	19-117	5	30
4,6-Dinitro-2-methylphenol	45 U	47.2	72.7	65	48.6	72.7	67	25-154	3	30
4-Bromophenyl Phenyl Ether	9.1 U	34.6	72.7	48	36.9	72.7	51	39-115	6	30
4-Chloro-3-methylphenol	46	101	72.7	77	100	72.7	75	41-126	3	30
4-Chloroaniline	9.1 U	48.1	72.7	66	50.2	72.7	69	19-111	4	30
4-Chlorophenyl Phenyl Ether	9.1 U	31.6	72.7	43	31.6	72.7	43	41-111	<1	30
4-Nitroaniline	9.1 U	26.1	72.7	36	29.6	72.7	41	18-143	13	30
4-Nitrophenol	45 U	31.9 J	72.7	44	30.3 J	72.7	42	10-126	5	30
Acenaphthene	9.1 U	42.2	72.7	58	40.9	72.7	56	43-117	4	30
Acenaphthylene	9.1 U	45.8	72.7	63	45.5	72.7	63	45-119	<1	30
Anthracene	9.1 U	44.0	72.7	60	44.8	72.7	62	45-127	3	30

Results flagged with an asterisk (*) indicate values outside control criteria.

Benz(a)anthracene

Benzo(b)fluoranthene

Benzo(a)pyrene

Results flagged with a pound (#) indicate the control criteria is not applicable.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

36.1

38.6

34.4

9.1 U

9.1 U

9.1 U

Matrix Spike and Matrix Spike Duplicate Data is presented for information purposes only. The matrix may or may not be relevant to samples reported in this report. The laboratory evaluates system performance based on the LCS and LCSD control limits.

72.7

72.7

72.7

50

53

47

36.2

39.1

34.0

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request:

R2005701

Project: Lo

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Date Collected:

07/01/20

Sample Matrix: Water

Date Received:

07/02/20

Date Analyzed: Date Extracted: 07/9/20 07/7/20

Duplicate Matrix Spike Summary Semivolatile Organic Compounds by GC/MS

Sample Name: WG-9954-070120-SG-027

Units:

ug/L

Lab Code:

R2005701-001

8270D

Basis: NA

Analysis Method: Prep Method:

EPA 3510C

Matrix Spike

Duplicate Matrix Spike

RO2007188-05

		RQ2	007188-04		R	Q2007188-0)5			
	Sample		Spike			Spike		% Rec		RPD
Analyte Name	Result	Result	Amount	% Rec	Result	Amount	% Rec	Limits	RPD	Limit
Benzo(g,h,i)perylene	9.1 U	40.2	72.7	55	39.3	72.7	54	50-143	2	30
Benzo(k)fluoranthene	9.1 U	36.6	72.7	50	37.4	72.7	51	46-139	2	30
Benzoic Acid	4500 E	4010 E	109	-482 #	91 U	109	-4155#	10-94	NC	30
Benzyl Alcohol	290	272	72.7	-30 *	282	72.7	-17 *	31-109	NC	30
2,2'-Oxybis(1-chloropropane)	9.1 U	38.4	72.7	53	34.5	72.7	47	21-126	12	30
Bis(2-chloroethoxy)methane	9.1 U	62.4	72.7	86	61.3	72.7	84	41-118	2	30
Bis(2-chloroethyl) Ether	20	59.0	72.7	54	55.0	72.7	49	33-108	10	30
Bis(2-ethylhexyl) Phthalate	9.1 U	38.0	72.7	52	36.4	72.7	50	41-132	4	30
Butyl Benzyl Phthalate	9.1 U	44.1	72.7	61	42.9	72.7	59	41-148	3	30
Chrysene	9.1 U	38.0	72.7	52	38.1	72.7	52	47-126	<1	30
Di-n-butyl Phthalate	9.1 U	46.5	72.7	64	47.5	72.7	65	43-130	2	30
Di-n-octyl Phthalate	9.1 U	41.7	72.7	57	41.2	72.7	57	40-139	<1	30
Dibenz(a,h)anthracene	9.1 U	43.3	72.7	60	43.0	72.7	59	43-136	2	30
Dibenzofuran	9.1 U	43.3	72.7	60	42.5	72.7	58	46-119	3	30
Diethyl Phthalate	9.1 U	42.8	72.7	59	43.1	72.7	59	36-122	<1	30
Dimethyl Phthalate	9.1 U	41.3	72.7	57	42.6	72.7	59	33-123	3	30
Fluoranthene	9.1 U	44.6	72.7	61	46.1	72.7	63	43-135	3	30
Fluorene	9.1 U	39.0	72.7	54	39.4	72.7	54	43-113	<1	30
Hexachlorobenzene	9.1 U	35.4	72.7	49	37.5	72.7	51	42-125	4	30
Hexachlorobutadiene	9.1 U	55.2	72.7	76	50.7	72.7	70	10-111	8	30
Hexachlorocyclopentadiene	9.1 U	20.9	72.7	29	20.5	72.7	28	10-103	4	30
Hexachloroethane	9.1 U	32.1	72.7	44	29.8	72.7	41	12-101	7	30
Indeno(1,2,3-cd)pyrene	9.1 U	37.9	72.7	52	36.6	72.7	50	49-140	4	30
Isophorone	9.1 U	4.33 J	72.7	6 *	5.01 J	72.7	7 *	40-111	15	30
N-Nitrosodi-n-propylamine	9.1 U	42.4	72.7	58	38.6	72.7	53	35-108	9	30
N-Nitrosodiphenylamine	9.1 U	49.0	72.7	67	53.0	72.7	73	43-127	9	30
Naphthalene	9.1 U	2040 E	72.7	2804 *	2060 E	72.7	2836 *	37-108	1	30
Nitrobenzene	9.1 U	61.0	72.7	84	57.2	72.7	79	35-112	6	30
Pentachlorophenol (PCP)	45 U	76.5	72.7	105	77.3	72.7	106	29-164	<1	30
Phenanthrene	9.1 U	43.3	72.7	60	44.3	72.7	61	46-123	2	30
Phenol	40	60.4	72.7	28	56.5	72.7	23	10-113	20	30
Pyrene	9.1 U	43.2	72.7	59	44.9	72.7	62	44-129	5	30

Results flagged with an asterisk (*) indicate values outside control criteria.

Results flagged with a pound (#) indicate the control criteria is not applicable.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

Matrix Spike and Matrix Spike Duplicate Data is presented for information purposes only. The matrix may or may not be relevant to samples reported in this report. The laboratory evaluates system performance based on the LCS and LCSD control limits.

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007188-01
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	10 U	10	1.2	1	07/09/20 10:38	7/7/20	
1,2-Dichlorobenzene	10 U	10	1.2	1	07/09/20 10:38	7/7/20	
1,3-Dichlorobenzene	10 U	10	1.1	1	07/09/20 10:38	7/7/20	
1,4-Dichlorobenzene	10 U	10	1.2	1	07/09/20 10:38	7/7/20	
2,4,5-Trichlorophenol	10 U	10	1.1	1	07/09/20 10:38	7/7/20	
2,4,6-Trichlorophenol	10 U	10	1.4	1	07/09/20 10:38	7/7/20	
2,4-Dichlorophenol	10 U	10	1.3	1	07/09/20 10:38	7/7/20	
2,4-Dimethylphenol	10 U	10	1.4	1	07/09/20 10:38	7/7/20	
2,4-Dinitrophenol	50 U	50	20	1	07/09/20 10:38	7/7/20	
2,4-Dinitrotoluene	10 U	10	2.4	1	07/09/20 10:38	7/7/20	
2,6-Dinitrotoluene	10 U	10	1.4	1	07/09/20 10:38	7/7/20	
2-Chloronaphthalene	10 U	10	1.4	1	07/09/20 10:38	7/7/20	
2-Chlorophenol	10 U	10	1.1	1	07/09/20 10:38	7/7/20	
2-Methylnaphthalene	10 U	10	1.3	1	07/09/20 10:38	7/7/20	
2-Methylphenol	10 U	10	1.0	1	07/09/20 10:38	7/7/20	
2-Nitroaniline	10 U	10	1.4	1	07/09/20 10:38	7/7/20	
2-Nitrophenol	10 U	10	1.5	1	07/09/20 10:38	7/7/20	
3,3'-Dichlorobenzidine	10 U	10	1.2	1	07/09/20 10:38	7/7/20	
3- and 4-Methylphenol Coelution	10 U	10	1.2	1	07/09/20 10:38	7/7/20	
3-Nitroaniline	10 U	10	2.5	1	07/09/20 10:38	7/7/20	
4,6-Dinitro-2-methylphenol	50 U	50	20	1	07/09/20 10:38	7/7/20	
4-Bromophenyl Phenyl Ether	10 U	10	1.7	1	07/09/20 10:38	7/7/20	
4-Chloro-3-methylphenol	10 U	10	1.1	1	07/09/20 10:38	7/7/20	
4-Chloroaniline	10 U	10	1.0	1	07/09/20 10:38	7/7/20	
4-Chlorophenyl Phenyl Ether	10 U	10	1.5	1	07/09/20 10:38	7/7/20	
4-Nitroaniline	10 U	10	2.7	1	07/09/20 10:38	7/7/20	
4-Nitrophenol	50 U	50	6.4	1	07/09/20 10:38	7/7/20	
Acenaphthene	10 U	10	1.4	1	07/09/20 10:38	7/7/20	
Acenaphthylene	10 U	10	1.4	1	07/09/20 10:38	7/7/20	
Anthracene	10 U	10	1.3	1	07/09/20 10:38	7/7/20	
Benz(a)anthracene	10 U	10	1.6	1	07/09/20 10:38	7/7/20	
Benzo(a)pyrene	10 U	10	1.2	1	07/09/20 10:38	7/7/20	
Benzo(b)fluoranthene	10 U	10	1.2	1	07/09/20 10:38	7/7/20	
Benzo(g,h,i)perylene	10 U	10	1.0	1	07/09/20 10:38	7/7/20	
Benzo(k)fluoranthene	10 U	10	1.3	1	07/09/20 10:38	7/7/20	
Benzoic Acid	100 U	100	36	1	07/09/20 10:38	7/7/20	
Benzyl Alcohol	10 U	10	1.6	1	07/09/20 10:38	7/7/20	
2,2'-Oxybis(1-chloropropane)	10 U	10	1.4	1	07/09/20 10:38	7/7/20	
Bis(2-chloroethoxy)methane	10 U	10	1.9	1	07/09/20 10:38	7/7/20	
Bis(2-chloroethyl) Ether	10 U	10	1.3	1	07/09/20 10:38	7/7/20	
Bis(2-ethylhexyl) Phthalate	10 U	10	1.0	1	07/09/20 10:38	7/7/20	
Butyl Benzyl Phthalate	10 U	10	1.4	1	07/09/20 10:38	7/7/20	
Chrysene	10 U	10	1.2	1	07/09/20 10:38	7/7/20	

Printed 7/30/2020 4:06:16 PM

Superset Reference: 20-0000555360 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007188-01
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	10 U	10	2.0	1	07/09/20 10:38	7/7/20	
Di-n-octyl Phthalate	10 U	10	3.3	1	07/09/20 10:38	7/7/20	
Dibenz(a,h)anthracene	10 U	10	1.1	1	07/09/20 10:38	7/7/20	
Dibenzofuran	10 U	10	1.4	1	07/09/20 10:38	7/7/20	
Diethyl Phthalate	10 U	10	1.1	1	07/09/20 10:38	7/7/20	
Dimethyl Phthalate	10 U	10	1.3	1	07/09/20 10:38	7/7/20	
Fluoranthene	10 U	10	1.5	1	07/09/20 10:38	7/7/20	
Fluorene	10 U	10	1.3	1	07/09/20 10:38	7/7/20	
Hexachlorobenzene	10 U	10	1.6	1	07/09/20 10:38	7/7/20	
Hexachlorobutadiene	10 U	10	1.0	1	07/09/20 10:38	7/7/20	
Hexachlorocyclopentadiene	10 U	10	2.2	1	07/09/20 10:38	7/7/20	
Hexachloroethane	10 U	10	1.1	1	07/09/20 10:38	7/7/20	
Indeno(1,2,3-cd)pyrene	10 U	10	1.8	1	07/09/20 10:38	7/7/20	
Isophorone	10 U	10	1.4	1	07/09/20 10:38	7/7/20	
N-Nitrosodi-n-propylamine	10 U	10	1.2	1	07/09/20 10:38	7/7/20	
N-Nitrosodiphenylamine	10 U	10	2.7	1	07/09/20 10:38	7/7/20	
Naphthalene	10 U	10	1.2	1	07/09/20 10:38	7/7/20	
Nitrobenzene	10 U	10	1.5	1	07/09/20 10:38	7/7/20	
Pentachlorophenol (PCP)	50 U	50	9.8	1	07/09/20 10:38	7/7/20	
Phenanthrene	10 U	10	1.4	1	07/09/20 10:38	7/7/20	
Phenol	10 U	10	1.0	1	07/09/20 10:38	7/7/20	
Pyrene	10 U	10	1.5	1	07/09/20 10:38	7/7/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	84	35 - 141	07/09/20 10:38	
2-Fluorobiphenyl	68	31 - 118	07/09/20 10:38	
2-Fluorophenol	48	10 - 105	07/09/20 10:38	
Nitrobenzene-d5	75	31 - 110	07/09/20 10:38	
Phenol-d6	35	10 - 107	07/09/20 10:38	
p-Terphenyl-d14	67	10 - 165	07/09/20 10:38	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown hydrocarbon	11.89	4.3	J	
	unknown hydrocarbon	12.48	6.3	J	
	unknown hydrocarbon	13.13	6.7	J	
	unknown	13.83	5.3	J	
	unknown hydrocarbon	14.58	4.1	J	
000544-25-2	1,3,5-Cycloheptatriene	2.68	9.2	JN	
000095-49-8	Benzene, 1-chloro-2-methyl-	4.12	4.2	JN	

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QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/09/20

Sample Matrix: Water

Duplicate Lab Control Sample Summary Semivolatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005701

Lab Control Sample

Duplicate Lab Control Sample

RQ2007188-02

RQ2007188-03

Analyte Name	Analytica l Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
1,2,4-Trichlorobenzene	8270D	52.8	80.0	66	51.4	80.0	64	10-127	3	30
1,2-Dichlorobenzene	8270D	52.1	80.0	65	46.3	80.0	58	23-130	11	30
1,3-Dichlorobenzene	8270D	50.7	80.0	63	44.6	80.0	56	21-90	12	30
1,4-Dichlorobenzene	8270D	49.3	80.0	62	45.1	80.0	56	10-124	10	30
2,4,5-Trichlorophenol	8270D	67.5	80.0	84	67.2	80.0	84	48-134	<1	30
2,4,6-Trichlorophenol	8270D	64.9	80.0	81	64.6	80.0	81	44-135	<1	30
2,4-Dichlorophenol	8270D	56.7	80.0	71	55.8	80.0	70	48-127	1	30
2,4-Dimethylphenol	8270D	63.5	80.0	79	66.6	80.0	83	59-113	5	30
2,4-Dinitrophenol	8270D	57.6	80.0	72	58.1	80.0	73	21-154	1	30
2,4-Dinitrotoluene	8270D	74.5	80.0	93	75.5	80.0	94	54-130	1	30
2,6-Dinitrotoluene	8270D	79.0	80.0	99	80.7	80.0	101	51-127	2	30
2-Chloronaphthalene	8270D	64.2	80.0	80	68.1	80.0	85	40-108	6	30
2-Chlorophenol	8270D	51.3	80.0	64	45.4	80.0	57	42-112	12	30
2-Methylnaphthalene	8270D	57.8	80.0	72	61.4	80.0	77	34-102	7	30
2-Methylphenol	8270D	56.4	80.0	70	51.6	80.0	65	47-100	7	30
2-Nitroaniline	8270D	76.1	80.0	95	76.9	80.0	96	52-133	1	30
2-Nitrophenol	8270D	60.3	80.0	75	61.4	80.0	77	43-131	3	30
3,3'-Dichlorobenzidine	8270D	73.6	80.0	92	72.6	80.0	91	43-126	1	30
3- and 4-Methylphenol Coelution	8270D	51.8	80.0	65	47.2	80.0	59	40-92	10	30
3-Nitroaniline	8270D	73.4	80.0	92	68.6	80.0	86	42-111	7	30
4,6-Dinitro-2-methylphenol	8270D	63.1	80.0	79	58.2	80.0	73	36-152	8	30
4-Bromophenyl Phenyl Ether	8270D	69.7	80.0	87	70.4	80.0	88	48-114	1	30
4-Chloro-3-methylphenol	8270D	64.8	80.0	81	63.6	80.0	80	52-113	1	30
4-Chloroaniline	8270D	66.8	80.0	84	62.7	80.0	78	44-109	7	30
4-Chlorophenyl Phenyl Ether	8270D	63.3	80.0	79	64.8	80.0	81	51-107	3	30
4-Nitroaniline	8270D	71.8	80.0	90	70.0	80.0	88	54-133	2	30
4-Nitrophenol	8270D	27.1 J	80.0	34	24.5 J	80.0	31	10-126	9	30
Acenaphthene	8270D	67.8	80.0	85	71.4	80.0	89	52-107	5	30
Acenaphthylene	8270D	74.2	80.0	93	77.1	80.0	96	55-109	3	30
Anthracene	8270D	77.5	80.0	97	78.0	80.0	98	55-116	1	30
Benz(a)anthracene	8270D	75.0	80.0	94	74.7	80.0	93	61-121	1	30
Benzo(a)pyrene	8270D	85.1	80.0	106	81.7	80.0	102	44-114	4	30
Benzo(b)fluoranthene	8270D	72.5	80.0	91	72.6	80.0	91	62-115	<1	30
D: 17/20/2020 106 16 D) 5							. D. C	20 00005	55060	00

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QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/09/20

Sample Matrix: Water

Duplicate Lab Control Sample Summary Semivolatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005701

Lab Control Sample

Duplicate Lab Control Sample

RQ2007188-02

RQ2007188-03

Analyte Name	Analytica l Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
Benzo(g,h,i)perylene	8270D	88.4	80.0	111	88.6	80.0	111	63-136	<1	30
Benzo(k)fluoranthene	8270D	80.1	80.0	100	82.4	80.0	103	49-133	3	30
Benzoic Acid	8270D	87.5 J	120	73	87.8 J	120	73	10-94	<1	30
Benzyl Alcohol	8270D	66.2	80.0	83	64.2	80.0	80	31-109	4	30
2,2'-Oxybis(1-chloropropane)	8270D	58.8	80.0	74	56.4	80.0	70	32-122	6	30
Bis(2-chloroethoxy)methane	8270D	59.7	80.0	75	69.6	80.0	87	55-110	15	30
Bis(2-chloroethyl) Ether	8270D	55.3	80.0	69	50.2	80.0	63	46-102	9	30
Bis(2-ethylhexyl) Phthalate	8270D	80.6	80.0	101	81.1	80.0	101	51-132	<1	30
Butyl Benzyl Phthalate	8270D	77.6	80.0	97	79.1	80.0	99	41-148	2	30
Chrysene	8270D	79.3	80.0	99	80.4	80.0	100	57-118	1	30
Di-n-butyl Phthalate	8270D	84.9	80.0	106	84.3	80.0	105	57-128	<1	30
Di-n-octyl Phthalate	8270D	82.0	80.0	102	80.4	80.0	101	62-124	<1	30
Dibenz(a,h)anthracene	8270D	97.4	80.0	122	96.1	80.0	120	54-135	2	30
Dibenzofuran	8270D	72.8	80.0	91	76.1	80.0	95	55-110	4	30
Diethyl Phthalate	8270D	70.7	80.0	88	73.7	80.0	92	53-113	4	30
Dimethyl Phthalate	8270D	77.1	80.0	96	78.7	80.0	98	51-112	2	30
Fluoranthene	8270D	83.7	80.0	105	82.5	80.0	103	66-127	2	30
Fluorene	8270D	72.1	80.0	90	75.3	80.0	94	54-106	4	30
Hexachlorobenzene	8270D	82.6	80.0	103	83.6	80.0	104	53-123	<1	30
Hexachlorobutadiene	8270D	53.9	80.0	67	56.1	80.0	70	16-95	4	30
Hexachlorocyclopentadiene	8270D	30.9	80.0	39	33.7	80.0	42	10-99	7	30
Hexachloroethane	8270D	49.3	80.0	62	43.0	80.0	54	15-92	14	30
Indeno(1,2,3-cd)pyrene	8270D	81.7	80.0	102	78.1	80.0	98	62-137	4	30
Isophorone	8270D	56.8	80.0	71	59.4	80.0	74	50-116	4	30
N-Nitrosodi-n-propylamine	8270D	63.4	80.0	79	65.7	80.0	82	49-115	4	30
N-Nitrosodiphenylamine	8270D	89.2	80.0	112	89.6	80.0	112	45-123	<1	30
Naphthalene	8270D	59.0	80.0	74	59.6	80.0	74	38-99	<1	30
Nitrobenzene	8270D	62.3	80.0	78	63.7	80.0	80	46-108	3	30
Pentachlorophenol (PCP)	8270D	77.6	80.0	97	75.7	80.0	95	29-164	2	30
Phenanthrene	8270D	75.0	80.0	94	75.8	80.0	95	58-118	1	30
Phenol	8270D	33.1	80.0	41	29.5	80.0	37	10-113	10	30
Pyrene	8270D	82.2	80.0	103	83.4	80.0	104	61-122	<1	30

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Semivolatile Organic Compounds by GC

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

QA/QC Report

Service Request: R2005701

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

SURROGATE RECOVERY SUMMARY

Organochlorine Pesticides by Gas Chromatography

Analysis Method: 8081B **Extraction Method:** EPA 3510C

		Decachlorobiphenyl	Tetrachloro-m-xylene	
Sample Name	Lab Code	10-164	10-147	
WG-9954-070120-SG-027	R2005701-001	14	171*	
WG-9954-070120-SG-026	R2005701-002	23	62	
WG-9954-070120-SG-025	R2005701-003	17	56	
WG-9954-070120-SG-024	R2005701-004	34	52	
WG-9954-070120-SG-028	R2005701-006	23	154*	
WG-9954-070120-SG-028 RE	R2005701-006	36	317*	
Method Blank	RQ2007190-01	70	67	
Method Blank	RQ2007457-01	67	60	
Lab Control Sample	RQ2007190-02	61	63	
Duplicate Lab Control Sample	RQ2007190-03	61	63	
Lab Control Sample	RQ2007457-02	66	67	
Duplicate Lab Control Sample	RQ2007457-03	64	63	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007190-01
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analysis Method: 8081B **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.050 U	0.050	0.020	1	07/08/20 14:15	7/7/20	_
4,4'-DDE	0.050 U	0.050	0.020	1	07/08/20 14:15	7/7/20	
4,4'-DDT	0.050 U	0.050	0.020	1	07/08/20 14:15	7/7/20	
Aldrin	0.050 U	0.050	0.020	1	07/08/20 14:15	7/7/20	
Dieldrin	0.050 U	0.050	0.020	1	07/08/20 14:15	7/7/20	
Endosulfan I	0.050 U	0.050	0.020	1	07/08/20 14:15	7/7/20	
Endosulfan II	0.050 U	0.050	0.020	1	07/08/20 14:15	7/7/20	
Endosulfan Sulfate	0.050 U	0.050	0.020	1	07/08/20 14:15	7/7/20	
Endrin	0.050 U	0.050	0.020	1	07/08/20 14:15	7/7/20	
Endrin Ketone	0.050 U	0.050	0.020	1	07/08/20 14:15	7/7/20	
Heptachlor	0.050 U	0.050	0.020	1	07/08/20 14:15	7/7/20	
Heptachlor Epoxide	0.050 U	0.050	0.020	1	07/08/20 14:15	7/7/20	
Methoxychlor	0.050 U	0.050	0.020	1	07/08/20 14:15	7/7/20	
Toxaphene	0.50 U	0.50	0.50	1	07/08/20 14:15	7/7/20	
alpha-BHC	0.050 U	0.050	0.020	1	07/08/20 14:15	7/7/20	
alpha-Chlordane	0.050 U	0.050	0.020	1	07/08/20 14:15	7/7/20	
beta-BHC	0.050 U	0.050	0.020	1	07/08/20 14:15	7/7/20	
delta-BHC	0.050 U	0.050	0.020	1	07/08/20 14:15	7/7/20	
gamma-BHC (Lindane)	0.050 U	0.050	0.020	1	07/08/20 14:15	7/7/20	
gamma-Chlordane	0.050 U	0.050	0.020	1	07/08/20 14:15	7/7/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	70	10 - 164	07/08/20 14:15	
Tetrachloro-m-xylene	67	10 - 147	07/08/20 14:15	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007457-01
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analysis Method: 8081B **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
4,4'-DDE	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
4,4'-DDT	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
Aldrin	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
Dieldrin	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
Endosulfan I	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
Endosulfan II	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
Endosulfan Sulfate	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
Endrin	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
Endrin Ketone	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
Heptachlor	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
Heptachlor Epoxide	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
Methoxychlor	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
Toxaphene	0.50 U	0.50	0.50	1	07/15/20 14:09	7/14/20	
alpha-BHC	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
alpha-Chlordane	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
beta-BHC	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
delta-BHC	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
gamma-BHC (Lindane)	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
gamma-Chlordane	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	67	10 - 164	07/15/20 14:09	
Tetrachloro-m-xylene	60	10 - 147	07/15/20 14:09	

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/08/20

Sample Matrix: Water

Duplicate Lab Control Sample Summary Organochlorine Pesticides by Gas Chromatography

Units:ug/L Basis:NA

Service Request: R2005701

Lab Control Sample

Duplicate Lab Control Sample

RQ2007190-02

RQ2007190-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
4,4'-DDD	8081B	0.323	0.400	81	0.333	0.400	83	42-159	3	30
4,4'-DDE	8081B	0.301	0.400	75	0.312	0.400	78	47-147	3	30
4,4'-DDT	8081B	0.319	0.400	80	0.321	0.400	80	41-149	<1	30
Aldrin	8081B	0.250	0.400	62	0.260	0.400	65	22-137	4	30
Dieldrin	8081B	0.331	0.400	83	0.345	0.400	86	52-144	4	30
Endosulfan I	8081B	0.324	0.400	81	0.338	0.400	84	52-136	4	30
Endosulfan II	8081B	0.340	0.400	85	0.350	0.400	88	57-138	3	30
Endosulfan Sulfate	8081B	0.309	0.400	77	0.314	0.400	78	34-156	2	30
Endrin	8081B	0.331	0.400	83	0.341	0.400	85	56-143	3	30
Endrin Ketone	8081B	0.341	0.400	85	0.348	0.400	87	59-143	2	30
Heptachlor	8081B	0.232	0.400	58	0.235	0.400	59	32-141	1	30
Heptachlor Epoxide	8081B	0.325	0.400	81	0.338	0.400	85	51-143	4	30
Methoxychlor	8081B	0.330	0.400	82	0.333	0.400	83	56-149	<1	30
alpha-BHC	8081B	0.312	0.400	78	0.330	0.400	83	36-151	6	30
alpha-Chlordane	8081B	0.316	0.400	79	0.330	0.400	82	50-139	4	30
beta-BHC	8081B	0.335	0.400	84	0.349	0.400	87	55-149	4	30
delta-BHC	8081B	0.317	0.400	79	0.325	0.400	81	29-159	3	30
gamma-BHC (Lindane)	8081B	0.317	0.400	79	0.332	0.400	83	41-149	5	30
gamma-Chlordane	8081B	0.321	0.400	80	0.237	0.400	59	50-140	30	30

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request: R2005701 **Project:** Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/15/20

Sample Matrix: Water

Duplicate Lab Control Sample Summary Organochlorine Pesticides by Gas Chromatography

Units:ug/L Basis:NA

Lab Control Sample

Duplicate Lab Control Sample

RQ2007457-02

RQ2007457-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
4,4'-DDD	8081B	0.295	0.400	74	0.278	0.400	69	42-159	6	30
4,4'-DDE	8081B	0.296	0.400	74	0.281	0.400	70	47-147	5	30
4,4'-DDT	8081B	0.317	0.400	79	0.308	0.400	77	41-149	3	30
Aldrin	8081B	0.248	0.400	62	0.240	0.400	60	22-137	3	30
Dieldrin	8081B	0.319	0.400	80	0.314	0.400	79	52-144	1	30
Endosulfan I	8081B	0.312	0.400	78	0.307	0.400	77	52-136	2	30
Endosulfan II	8081B	0.231	0.400	58	0.288	0.400	72	57-138	22	30
Endosulfan Sulfate	8081B	0.228	0.400	57	0.267	0.400	67	34-156	16	30
Endrin	8081B	0.325	0.400	81	0.318	0.400	80	56-143	2	30
Endrin Ketone	8081B	0.288	0.400	72	0.305	0.400	76	59-143	6	30
Heptachlor	8081B	0.246	0.400	62	0.250	0.400	63	32-141	2	30
Heptachlor Epoxide	8081B	0.316	0.400	79	0.309	0.400	77	51-143	2	30
Methoxychlor	8081B	0.311	0.400	78	0.315	0.400	79	56-149	1	30
alpha-BHC	8081B	0.311	0.400	78	0.297	0.400	74	36-151	5	30
alpha-Chlordane	8081B	0.309	0.400	77	0.301	0.400	75	50-139	3	30
beta-BHC	8081B	0.312	0.400	78	0.310	0.400	77	55-149	<1	30
delta-BHC	8081B	0.252	0.400	63	0.279	0.400	70	29-159	10	30
gamma-BHC (Lindane)	8081B	0.315	0.400	79	0.301	0.400	75	41-149	4	30
gamma-Chlordane	8081B	0.304	0.400	76	0.297	0.400	74	50-140	2	30

QA/QC Report

Service Request: R2005701

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

Project:

SURROGATE RECOVERY SUMMARY Polychlorinated Biphenyls (PCBs) by GC

Analysis Method: 8082A **Extraction Method:** EPA 3510C

		Decachlorobiphenyl	Tetrachloro-m-xylene	
Sample Name	Lab Code	10-152	14-129	
WG-9954-070120-SG-027	R2005701-001	22	50	
WG-9954-070120-SG-026	R2005701-002	43	42	
WG-9954-070120-SG-025	R2005701-003	18	39	
WG-9954-070120-SG-024	R2005701-004	32	34	
WG-9954-070120-SG-028	R2005701-006	24	57	
Method Blank	RQ2007190-01	67	55	
Lab Control Sample	RQ2007190-02	62	47	
Duplicate Lab Control Sample	RQ2007190-03	66	52	
WG-9954-070120-SG-025 MS	RQ2007190-04	24	31	
WG-9954-070120-SG-025 DMS	RQ2007190-05	25	31	

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates) **Service Request:** R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring **Date Collected:** 07/01/20 **Date Received:**

Sample Matrix: Water

07/02/20

07/10/20

Date Extracted:

Date Analyzed:

07/7/20

Duplicate Matrix Spike Summary Polychlorinated Biphenyls (PCBs) by GC

Sample Name: WG-9954-070120-SG-025 **Units:** ug/L

Lab Code: R2005701-003 **Basis:** NA

Analysis Method: 8082A **Prep Method:** EPA 3510C

Matrix Spike

Duplicate Matrix Spike

RQ2007190-05

	Sample		Spike			Spike		% Rec		RPD
Analyte Name	Result	Result	Amount	% Rec	Result	Amount	% Rec	Limits	RPD	Limit
Aroclor 1016	0.91 U	1.70	3.64	47	1.93	3.64	53	32-142	13	30
Aroclor 1260	0.91 U	1.77	3.64	49	1.89	3.64	52	28-142	3	30

RQ2007190-04

Results flagged with an asterisk (*) indicate values outside control criteria.

Results flagged with a pound (#) indicate the control criteria is not applicable.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

Matrix Spike and Matrix Spike Duplicate Data is presented for information purposes only. The matrix may or may not be relevant to samples reported in this report. The laboratory evaluates system performance based on the LCS and LCSD control limits.

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005701

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007190-01Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analysis Method: 8082A **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	1.0 U	1.0	0.50	1	07/08/20 14:37	7/7/20	
Aroclor 1221	2.0 U	2.0	1.0	1	07/08/20 14:37	7/7/20	
Aroclor 1232	1.0 U	1.0	0.50	1	07/08/20 14:37	7/7/20	
Aroclor 1242	1.0 U	1.0	0.50	1	07/08/20 14:37	7/7/20	
Aroclor 1248	1.0 U	1.0	0.50	1	07/08/20 14:37	7/7/20	
Aroclor 1254	1.0 U	1.0	0.50	1	07/08/20 14:37	7/7/20	
Aroclor 1260	1.0 U	1.0	0.50	1	07/08/20 14:37	7/7/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	67	10 - 152	07/08/20 14:37	
Tetrachloro-m-xvlene	55	14 - 129	07/08/20 14:37	

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix:

Water

Service Request: R2005701 Date Analyzed: 07/08/20

Duplicate Lab Control Sample Summary Polychlorinated Biphenyls (PCBs) by GC

Units:ug/L Basis:NA

Lab Control Sample

Duplicate Lab Control Sample

RQ2007190-02

RQ2007190-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
Aroclor 1016	8082A	2.88	4.00	72	3.30	4.00	82	49-123	13	30
Aroclor 1260	8082A	3.13	4.00	78	3.62	4.00	91	30-120	15	30



Service Request No:R2005820

Ms. Kathy Willy GHD Services Inc. 2055 Niagara Falls Blvd., Niagara Falls, NY 14304

Laboratory Results for: Love Canal:292-402-D02-3100

Dear Ms. Willy,

Enclosed are the results of the sample(s) submitted to our laboratory July 08, 2020 For your reference, these analyses have been assigned our service request number **R2005820**.

All testing was performed according to our laboratory's quality assurance program and met the requirements of the TNI standards except as noted in the case narrative report. Any testing not included in the lab's accreditation is identified on a Non-Certified Analytes report. All results are intended to be considered in their entirety. ALS Environmental is not responsible for use of less than the complete report. Results apply only to the individual samples submitted to the lab for analysis, as listed in the report. The measurement uncertainty of the results included in this report is within that expected when using the prescribed method(s), and represented by Laboratory Control Sample control limits. Any events, such as QC failures or Holding Time exceedances, which may add to the uncertainty are explained in the report narrative or are flagged with qualifiers. The flags are explained in the Report Qualifiers and Definitions page of this report.

Please contact me if you have any questions. My extension is 7471. You may also contact me via email at Brady.Kalkman@alsglobal.com.

Respectfully submitted,

Goody Kullen

ALS Group USA, Corp. dba ALS Environmental

Brady Kalkman Project Manager

ADDRESS



Narrative Documents

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com



Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100 Date Received: 07/08/2020

Sample Matrix: Water

CASE NARRATIVE

All analyses were performed consistent with the quality assurance program of ALS Environmental. This report contains analytical results for samples for the Tier level IV requested by the client.

Sample Receipt:

Five water samples were received for analysis at ALS Environmental on 07/08/2020. Any discrepancies upon initial sample inspection are annotated on the sample receipt and preservation form included within this report. The samples were stored at minimum in accordance with the analytical method requirements.

Semivolatiles by GC/MS:

Method 8270D, 07/14/2020: The lower control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). Since there were no detections of the analyte(s) above the MRL in the associated field samples, the quantitation is not affected. The data quality was not significantly affected and no further corrective action was taken.

Method 8270D, 07/14/2020: The Method Blank contained a low level of one or more analytes at concentrations above the Method Detection Limit (MDL). Since there were no detections of the analyte(s) in the associated field samples, the data quality was not significantly affected and no further corrective action was taken.

Method 8270D, 07/14/2020: The lower control limit for the spike recovery of the Matrix Spike Duplicate (MSD) was exceeded for one or more analyte. Precision is also outside limits. The LCS/LCSD/MS were within limits for all analytes. The analytes affected are flagged in the MS Summary.

Semivoa GC:

Method 8081B, 07/14/2020: The upper control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). The field samples analyzed in this sequence did not contain the analyte(s) in question above the Method Reporting Limit (MRL). Since the exceedance equates to a potential high bias, the data quality was not significantly affected and no further corrective action was taken.

Method 8081B, 07/13/2020: The lower control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). Since there were no detections of the analyte(s) above the MRL in the associated field samples, the quantitation is not affected. The data quality was not significantly affected and no further corrective action was taken. Method 8082A, 07/16/2020: The upper control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). The field samples analyzed in this sequence did not contain the analyte(s) in question above the Method Reporting Limit (MRL). Since the exceedance equates to a potential high bias, the data quality was not significantly affected and no further corrective action was taken.

Method 8082A, 07/13/2020: The upper control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). The field samples analyzed in this sequence did not contain the analyte(s) in question above the Method Reporting Limit (MRL). Since the exceedance equates to a potential high bias, the data quality was not significantly affected and no further corrective action was taken.

Volatiles by GC/MS:

Method 8260C, 07/20/2020: The lower control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). Since there were no detections of the analyte(s) above the MRL in the associated field samples, the quantitation is not affected. The data quality was not significantly affected and no further corrective action was taken.

Method 8260C, 07/20/2020: The upper control criterion was exceeded for one or more analytes in the Laboratory Control Sample (LCS). There were no detections of the analyte(s) above the MRL in the associated field samples. The error associated with elevated recovery equates to a high bias. The sample data is not significantly affected. No further corrective action was appropriate.



Method 8260C, 07/19/2020: The lower control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). Since there were no detections of the analyte(s) above the MRL in the associated field samples, the quantitation is not affected. The data quality was not significantly affected and no further corrective action was taken. Method 8260C, 07/19/2020: The upper control criterion was exceeded for one or more analytes in the Laboratory Control Sample (LCS). There were no detections of the analyte(s) above the MRL in the associated field samples. The error associated with elevated recovery equates to a high bias. The sample data is not significantly affected. No further corrective action was appropriate.

Approved by _____

Date 07/27/2020



SAMPLE DETECTION SUMMARY

Lab ID: R2005820-002

CLIENT ID: WG-9954-070620-SG-029

Analyte	Results	Flag	MDL	MRL	Units	Method		
Carbon Disulfide	6.4	J	0.42	10	ug/L	8260C		
alpha-BHC	0.47		0.019	0.045	ug/L	8081B		
delta-BHC	0.13		0.019	0.045	ug/L	8081B		
gamma-BHC (Lindane)	0.48		0.019	0.045	ug/L	8081B		
CLIENT ID: WG-9954-070720-SG-032	Lab ID: R2005820-004							
Analyte	Results	Flag	MDL	MRL	Units	Method		
Carbon Disulfide	3.7	J	0.42	10	ug/L	8260C		
alpha-BHC	0.070		0.019	0.045	ug/L	8081B		
delta-BHC	0.067		0.019	0.045	ug/L	8081B		
gamma-BHC (Lindane)	0.098		0.019	0.045	ug/L	8081B		
CLIENT ID: WG-9954-070720-SG-030		Lak	ID: R2005	820-005				
Analyte	Results	Flag	MDL	MRL	Units	Method		
Carbon Disulfide	3.4	J	0.42	10	ug/L	8260C		
alpha-BHC	0.24		0.019	0.045	ug/L	8081B		
delta-BHC	0.084		0.019	0.045	ug/L	8081B		
gamma-BHC (Lindane)	0.26		0.019	0.045	ug/L	8081B		



Sample Receipt Information

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request:R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

SAMPLE CROSS-REFERENCE

SAMPLE #	CLIENT SAMPLE ID	<u>DATE</u>	<u>TIME</u>
R2005820-001	TB-9954-070620-SG-006	7/6/2020	0000
R2005820-002	WG-9954-070620-SG-029	7/6/2020	1355
R2005820-003	WG-9954-070720-SG-031	7/7/2020	1030
R2005820-004	WG-9954-070720-SG-032	7/7/2020	1300
R2005820-005	WG-9954-070720-SG-030	7/7/2020	1215

*CHAIN-OF-CUSTODY/Analytical Request Document The Cheln-of-Custody is a LEGAL DOCUMENT. All relevant fields must be completed accurately.

. Client In	formation
GLEN SPRINGS HOLDINGS INC	Report To: Kathy Willy
806 95TH STREET	Сору То:
LOVE CANAL	
NIAGARA FALLS, NEW YORK 14304	Involce To:
Phone: 716-283-0111	PO:
Fax: 716-283-2856	Project Neme: LOVE CANAL ANNUAL GW
Email: kathy.willy@ghd.com	Project Number: 9954

Laboratory: ALS				
Laboratory Locatic BUILDING 300, SU ROCHESTER, NY	TE 360	EFFERSON	IRD	
ROCHESTER, NY	4623			
Laboratory Contac	t: BRADY	KALKMAN		
Requested Due Da	te:	TAT: 10		

	Event Information
ID#: L	.C ANNUAL GW SAMPLING 2020-06-1
	SSOW Ref#: 292-402-999-3100
Sampler Na	IME:S GARONFR D TURAN

Sample identification	Valld Matrix Code WG Groundwater WB Borehole Water WS Surface Water SO Soil SE Sediment	Matrix Code	Date Collected	Time Collected	PestPCBs(None)	SVOC(none)	VOA(HCI)	Remarks	Sample Cond Temp in C Received on ice Sealed Cooler Samples Intact	Y/N Y/N Y/N	
TB-9954-070620-SG-006		WG Q	07/06/2020	00:00	0	0	3 .				
WG-9954-070620-SG-029	•	WG	07/06/2020	13:55	2	2	3				
WG-9954-070720-SG-031		WG	07/07/2020	10;30	2	2	3		,		
WG-9954-070720-SG-032		WG	07/07/2020	13:00	2	2	3				
WG-9954-070720-SG-030		WG	07/07/2020	12:15	2	2	3				
Total Bottles					8	8	16	Grand Tot	al:31		

SHIPMENT METHOD	NO. OF COOLERS	RELINQUISHED BY:	DATE	TIME	RECIEVED BY:			DATE	TIME
FedEx		Shapen Mardon	77202	0143	11110	1111111	A3	782021	061010
AIRBILL#:				,	1000				

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ALS Enuir	onme	ental				1565			(5 65)	286-5	5380 /		585)	•	ester, NY 1462 -8475	3 QC	04, 005, 006, 007, 008, 009, 010, 11, 012, 013		T030477
Project Name: Love Canal:292-402-D02-3100	· .				Ι	2		14D	_		,			1				,	
Project Number: 9954 Annual Long Term Monitoring	Report To Kathy Will	v		1		<u>~</u>	_	-	1	_	,		_	4					
Company / Address GHD Services Inc. 2055 Niagara Falts Blvd., Suite 3				CONTAINERS	,			,								•			
Niagara Falis NY, 14304			-	Ş	8	ĺ	٦	G FP		1				١					•
Phone # 716-297-2160	FAX # 716-297-2			P	l &	Ω̈	[≳	9									•		,
Sampler Signature	Sampler F	rinted Name		NUMBER	8081B / Pest	8082A / PCB	3270D / SVO	3260C / VOC					١		Rema	rke			
CLIENT SAMPLE ID	LABID	SAMPLING Date Time	Matrix	_	<u> </u>	ž	180	86	-	10	6	4	-	+	Kema	11/2	1 .		
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R2005820	5
Love Canal:292-402-D02-3100	

Date/Time

Date/Time

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Cooler Receipt and Preservation Check Form

	R2005820		5
	GHD Services Inc. Love Canal:292-402-D02-310	Nima kana kan	
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roject/Clie	nt GH	D		/	_Folde	r Numbe	er				T I BEIRIN	40 nan	— — — — — — — — — — — — — — — — — — —		
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1 Were Cu	stody seals on	outside of cooler	?	(Y) (Y)	5a Pe	erchlora	te sa	amples	have req	uired he	adspa	ice?	YN	(NA)
2 Custody	papers proper	ly completed (ink	, signe	d)? (N	5b D	d VOA	yial	Alk,c	r Sulfide	have si	g* bu	bbles?	YA	NA
		good condition (1				6 W	here did	the	bottles	originat	e?	(AL	S/ROC	CLIE	NT
4 Circle:	Wet Ice Dry	Ice Gel packs	prese	nt?	N	7 S	oil VOA	rece	ived as	: Bu	lk E	псоге	503	Sset N	TA)
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Observed Te	mp (°C)	25							_					<u> </u>	
Within 0-6°C	7?	(Y N		Y	N	Y N		Y	N		N	Y		<u> Y</u>	N
If <0°C, wer	e samples froz	en? Y N		Y	N	Y N		Y	N	Y	N	Y	<u>N</u> _	Y	N
If out of T	emperature,	note packing/ice	condi	tion:		Ice	melted	P	oorly P	acked (d	escribed	l belo	w)	Same I	ay Rule
	=	un Samples:			ding App	roval C	lient aw	are a	at drop-	off Cl	ient not	ified l	oy:		
	held in storages placed in storage	e location: orage location:	R.60.		y <u>NE</u> y —	on		at]; at		within 4	8 hours	of sa	mpling?	Y	N
9. V 10. I 11. V 12. V	Vere all bottle Did all bottle la Vere correct co Vere 5035 vial	ervation Check** labels complete (bels and tags agreementainers used for s acceptable (no classettes / Tubes)	i.e. ana ee with the tes extra la	lysis, j custo ts indi bels, r	preservat dy paper icated? not leakin	non, etc.)? s?		đ		ES ES ES Tedlar®	NO NO NO NO Bags In			N/A N/A))
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<u>≥12</u> ≤2	·	HNO ₃			· · · · · · · · · · · · · · · · · · ·		-				<u> </u>				
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Bottle lot Explain a	numbers: N	ies/ Other Comm	Jo John	Infa	for l	10A V1	als, O	511	20-JE	anc .			<u> </u>		
*11	p Blank:	A113 VA	al3												,

HPROD	BULK
HTR	FLDT
SUB	HGFB
ALS	LL3541

Labels secondary reviewed by: _______PC Secondary Review: ______

*significant air bubbles: VOA > 5-6 mm : WC > 1 in. diameter

Internal Chain of Custody Report

Client: GHD Services Inc. Service Request: R2005820

Bottle ID	Methods	Date	Time	Sample Location / User	Disposed On
R2005820-001.01					
	8260C				
		7/8/2020	1649	SMO / DWARD	
		7/8/2020	1651	R-001 / DWARD	
		7/19/2020	1230	In Lab / KRUEST	
		7/19/2020	1251	R-001-S12 / KRUEST	
R2005820-001.02					
		7/8/2020	1640	SMO / DWARD	
			1649		
		7/8/2020	1651	R-001 / DWARD	
R2005820-001.03					
		7/8/2020	1649	SMO / DWARD	
		7/8/2020	1651	R-001 / DWARD	
R2005820-002.01		1, 3, 2020	1001	R out / D WIND	
N2UU302U-UU2.U1	8081B,8082A				
	000111,000211	7/8/2020	1649	SMO / DWARD	
		7/8/2020	1651	R-002 / DWARD	
		7/10/2020	1200	In Lab / VSTAUFFER	
R2005820-002.02					
12003020-002.02					
		7/8/2020	1649	SMO / DWARD	
		7/8/2020	1651	R-002 / DWARD	
R2005820-002.03					
		7/8/2020	1649	SMO / DWARD	
		7/8/2020	1651	R-001 / DWARD	
		7/19/2020	1230	In Lab / KRUEST	
		7/19/2020	1250	R-001-S12 / KRUEST	
R2005820-002.04	00.00				
	8260C	7/9/2020	1.640	CMO / DWARD	
		7/8/2020	1649	SMO / DWARD	
		7/8/2020	1651	R-001 / DWARD	
		7/20/2020	1209	In Lab / KRUEST	
		7/20/2020	1359	R-001-S12 / KRUEST	
R2005820-002.05					
		7/8/2020	1649	SMO / DWARD	
		7/8/2020	/	R-001 / DWARD	

Internal Chain of Custody Report

Client: GHD Services Inc. Service Request: R2005820

Bottle ID	Methods	Date	Time	Sample Location / User	Disposed On
		7/8/2020	1649	SMO / DWARD	
		7/8/2020	1651	R-002 / DWARD	
R2005820-002.07					
	8270D				
		7/8/2020	1649	SMO / DWARD	
		7/8/2020	1651	R-002 / DWARD	
		7/13/2020	0808	In Lab / VSTAUFFER	
R2005820-003.01					
		7/8/2020	1649	SMO / DWARD	
		7/8/2020	1651	R-002 / DWARD	
R2005820-003.02	00017 00004				
	8081B,8082A	7/8/2020	1649	SMO / DWARD	
		7/8/2020	1651	R-002 / DWARD	
		7/10/2020	1200	In Lab / VSTAUFFER	
R2005820-003.03					
	8260C				
		7/8/2020	1649	SMO / DWARD	
		7/8/2020	1651	R-001 / DWARD	
		7/19/2020	1230	In Lab / KRUEST	
		7/19/2020	1250	R-001-S12 / KRUEST	
R2005820-003.04					
		7/8/2020	1649	SMO / DWARD	
		7/8/2020	1651	R-001 / DWARD	
R2005820-003.05					
		7/8/2020	1649	SMO / DWARD	
		7/8/2020	1651	R-001 / DWARD	
R2005820-003.06					
	8270D				
		7/8/2020	1649	SMO / DWARD	
		7/8/2020	1651	R-002 / DWARD	
		7/13/2020	0808	In Lab / VSTAUFFER	
R2005820-003.07					
		7/8/2020	1649	SMO / DWARD	
		7/8/2020	1651	R-002 / DWARD	

Internal Chain of Custody Report

Client: GHD Services Inc. Service Request: R2005820

Bottle ID	Methods	Date	Time	Sample Location / User	Disposed On
R2005820-004.01					
	8081B,8082A	= (0 (= 0 = 0		21.50 / P.W. / P.	
		7/8/2020	1649	SMO / DWARD	
		7/8/2020	1651	R-002 / DWARD	
		7/10/2020	1200	In Lab / VSTAUFFER	
R2005820-004.02					
		7/8/2020	1649	SMO / DWARD	
		7/8/2020	1651	R-002 / DWARD	
R2005820-004.03					
		7/8/2020	1649	SMO / DWARD	
		7/8/2020	1651	R-001 / DWARD	
		7/19/2020	1230	In Lab / KRUEST	
R2005820-004.04					
	8260C	7/8/2020	1649	SMO / DWARD	
		7/8/2020	1651	R-001 / DWARD	
		7/20/2020	1209	In Lab / KRUEST	
		7/20/2020	1359	R-001-S12 / KRUEST	
R2005820-004.05					
		7/8/2020	1649	SMO / DWARD	
		7/8/2020	1651	R-001 / DWARD	
R2005820-004.06					
	8270D	7/8/2020	1640	SMO / DWARD	
		7/8/2020	1649 1651	R-002 / DWARD	
		7/13/2020	0808	In Lab / VSTAUFFER	
R2005820-004.07		1/13/2020		III Dao / VOIMOII DA	
		7/8/2020	1649	SMO / DWARD	
		7/8/2020	1651	R-002 / DWARD	
R2005820-005.01					
	8081B,8082A				
		7/8/2020	1649	SMO / DWARD	
		7/8/2020	1651	R-002 / DWARD	
		7/10/2020	1201	In Lab / VSTAUFFER	
R2005820-005.02					
		7/8/2020	1649	SMO / DWARD	
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Internal Chain of Custody Report

Client: GHD Services Inc. Service Request: R2005820

Bottle ID	Methods	Date	Time	Sample Location / User	Disposed On
		7/8/2020	1651	R-002 / DWARD	
		7/13/2020	0808	In Lab / VSTAUFFER	
R2005820-005.03					
	8260C				
		7/8/2020	1649	SMO / DWARD	
		7/8/2020	1651	R-001 / DWARD	
		7/19/2020	1230	In Lab / KRUEST	
		7/19/2020	1250	R-001-S12 / KRUEST	
		7/20/2020	1209	In Lab / KRUEST	
		7/20/2020	1359	R-001-S12 / KRUEST	
R2005820-005.04					
		7/8/2020	1649	SMO / DWARD	
		7/8/2020	1651	R-001 / DWARD	
R2005820-005.05					
		7/0/2020	1640		
		7/8/2020	1649	SMO / DWARD	
		7/8/2020	1651	R-001 / DWARD	
R2005820-005.06					
	8270D				
		7/8/2020	1649	SMO / DWARD	
		7/8/2020	1651	R-002 / DWARD	
		7/13/2020	0808	In Lab / VSTAUFFER	
R2005820-005.07					
		7/8/2020	1649	SMO / DWARD	
		7/8/2020	1651	R-002 / DWARD	
		7/13/2020	0808	In Lab / VSTAUFFER	
		1/13/2020	0000	III Lau / VSIAUITEK	



Miscellaneous Forms

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com



REPORT QUALIFIERS AND DEFINITIONS

- U Analyte was analyzed for but not detected. The sample quantitation limit has been corrected for dilution and for percent moisture, unless otherwise noted in the case narrative.
- J Estimated value due to either being a Tentatively Identified Compound (TIC) or that the concentration is between the MRL and the MDL. Concentrations are not verified within the linear range of the calibration. For DoD: concentration >40% difference between two GC columns (pesticides/Arclors).
- B Analyte was also detected in the associated method blank at a concentration that may have contributed to the sample result.
- E Inorganics- Concentration is estimated due to the serial dilution was outside control limits.
- E Organics- Concentration has exceeded the calibration range for that specific analysis.
- D Concentration is a result of a dilution, typically a secondary analysis of the sample due to exceeding the calibration range or that a surrogate has been diluted out of the sample and cannot be assessed.
- * Indicates that a quality control parameter has exceeded laboratory limits. Under the õNotesö column of the Form I, this qualifier denotes analysis was performed out of Holding Time.
- H Analysis was performed out of hold time for tests that have an õimmediateö hold time criteria.
- # Spike was diluted out.

- + Correlation coefficient for MSA is <0.995.
- N Inorganics- Matrix spike recovery was outside laboratory limits.
- N Organics- Presumptive evidence of a compound (reported as a TIC) based on the MS library search.
- S Concentration has been determined using Method of Standard Additions (MSA).
- W Post-Digestion Spike recovery is outside control limits and the sample absorbance is <50% of the spike absorbance.
- P Concentration >40% difference between the two GC columns.
- C Confirmed by GC/MS
- Q DoD reports: indicates a pesticide/Aroclor is not confirmed (×100% Difference between two GC columns).
- X See Case Narrative for discussion.
- MRL Method Reporting Limit. Also known as:
- LOQ Limit of Quantitation (LOQ)

 The lowest concentration at which the method analyte may be reliably quantified under the method conditions.
- MDL Method Detection Limit. A statistical value derived from a study designed to provide the lowest concentration that will be detected 99% of the time. Values between the MDL and MRL are estimated (see J qualifier).
- LOD Limit of Detection. A value at or above the MDL which has been verified to be detectable.
- ND Non-Detect. Analyte was not detected at the concentration listed. Same as U qualifier.



Rochester Lab ID # for State Certifications¹

Connecticut ID # PH0556	Maine ID #NY0032	Pennsylvania ID# 68-786	
Delaware Approved	New Hampshire ID # 2941	Rhode Island ID # 158	
DoD ELAP #65817	New York ID # 10145	Virginia #460167	
Florida ID # E87674	North Carolina #676		

¹ Analyses were performed according to our laboratory

NELAP-approved quality assurance program and any applicable state or agency requirements. The test results meet requirements of the current NELAP/TNI standards or state or agency requirements, where applicable, except as noted in the case narrative. Since not all analyte/method/matrix combinations are offered for state/NELAC accreditation, this report may contain results which are not accredited. For a specific list of accredited analytes, contact the laboratory or go to https://www.alsglobal.com/locations/americas/north-america/usa/new-york/rochester-environmental

9/28/18

ALS Laboratory Group

Acronyms

ASTM American Society for Testing and Materials

A2LA American Association for Laboratory Accreditation

CARB California Air Resources Board

CAS Number Chemical Abstract Service registry Number

CFC Chlorofluorocarbon CFU Colony-Forming Unit

DEC Department of Environmental Conservation

DEQ Department of Environmental Quality

DHS Department of Health Services

DOE Department of Ecology DOH Department of Health

EPA U. S. Environmental Protection Agency

ELAP Environmental Laboratory Accreditation Program

GC Gas Chromatography

GC/MS Gas Chromatography/Mass Spectrometry

LUFT Leaking Underground Fuel Tank

M Modified

MCL Maximum Contaminant Level is the highest permissible concentration of a

substance allowed in drinking water as established by the USEPA.

MDL Method Detection Limit
MPN Most Probable Number
MRL Method Reporting Limit

NA Not Applicable NC Not Calculated

NCASI National Council of the Paper Industry for Air and Stream Improvement

ND Not Detected

NIOSH National Institute for Occupational Safety and Health

PQL Practical Quantitation Limit

RCRA Resource Conservation and Recovery Act

SIM Selected Ion Monitoring

TPH Total Petroleum Hydrocarbons

tr Trace level is the concentration of an analyte that is less than the PQL but

greater than or equal to the MDL.

Analyst Summary report

Service Request: R2005820

Date Collected: 07/6/20

Date Received: 07/8/20

Analyzed By

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term **Project:**

Monitoring

Sample Name: TB-9954-070620-SG-006

Lab Code: R2005820-001

Sample Matrix: Water

Analyzed By Analysis Method Extracted/Digested By

8260C **KRUEST**

Sample Name: WG-9954-070620-SG-029 **Date Collected:** 07/6/20

Lab Code: R2005820-002 Date Received: 07/8/20

Sample Matrix: Water

Sample Matrix:

Water

Extracted/Digested By Analyzed By Analysis Method

8081B **KSERCU JMISIUREWICZ** 8082A **KSERCU BALLGEIER**

KRUEST 8260C

8270D **KSERCU JMISIUREWICZ**

Sample Name: WG-9954-070720-SG-031 **Date Collected:** 07/7/20

Lab Code: R2005820-003 **Date Received:** 07/8/20

Analysis Method Extracted/Digested By 8081B **KSERCU JMISIUREWICZ KSERCU BALLGEIER** 8082A

8260C **KRUEST**

8270D **KSERCU JMISIUREWICZ**

Sample Name: WG-9954-070720-SG-032 **Date Collected:** 07/7/20

Lab Code: R2005820-004 Date Received: 07/8/20 **Sample Matrix:** Water

Analyzed By Extracted/Digested By Analysis Method

KSERCU 8081B **JMISIUREWICZ** 8082A **KSERCU BALLGEIER**

8260C **KRUEST**

8270D **JMISIUREWICZ KSERCU**

Printed 7/27/2020 10:42:09 AM Superset Reference:20-0000555771 rev 00

Analyst Summary report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Monitoring

Sample Name: WG-9954-070720-SG-030

Lab Code: R2005820-005

Sample Matrix: Water

Service Request: R2005820

Date Collected: 07/7/20

Date Received: 07/8/20

Analysis Method	Extracted/Digested By	Analyzed By
8081B	KSERCU	JMISIUREWICZ
8082A	KSERCU	BALLGEIER
8260C		KRUEST

8270D KSERCU JMISIUREWICZ



INORGANIC PREPARATION METHODS

The preparation methods associated with this report are found in these tables unless discussed in the case narrative.

Water/Liquid Matrix

Analytical Method	Preparation Method
200.7	200.2
200.8	200.2
6010C	3005A/3010A
6020A	ILM05.3
9034 Sulfide Acid Soluble	9030B
SM 4500-CN-E Residual	SM 4500-CN-G
Cyanide	
SM 4500-CN-E WAD	SM 4500-CN-I
Cyanide	

Solid/Soil/Non-Aqueous Matrix

Analytical Method	Preparation Method			
6010C	3050B			
6020A	3050B			
6010C TCLP (1311)	3005A/3010A			
extract				
6010 SPLP (1312) extract	3005A/3010A			
7199	3060A			
300.0 Anions/ 350.1/	DI extraction			
353.2/ SM 2320B/ SM				
5210B/ 9056A Anions				
For analytical methods not listed, the preparation method is the same as the analytical method reference.				



Sample Results

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com



Volatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/06/20 00:00

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 TB-9954-070620-SG-006
 Units: ug/L

 Lab Code:
 R2005820-001
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/19/20 18:54	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/19/20 18:54	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/19/20 18:54	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/19/20 18:54	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/19/20 18:54	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/19/20 18:54	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/19/20 18:54	
2-Butanone (MEK)	10 U	10	0.78	1	07/19/20 18:54	
2-Hexanone	10 U	10	0.20	1	07/19/20 18:54	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/19/20 18:54	
Acetone	10 U	10	5.0	1	07/19/20 18:54	
Benzene	5.0 U	5.0	0.20	1	07/19/20 18:54	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/19/20 18:54	
Bromoform	5.0 U	5.0	0.25	1	07/19/20 18:54	
Bromomethane	5.0 U	5.0	0.70	1	07/19/20 18:54	
Carbon Disulfide	10 U	10	0.42	1	07/19/20 18:54	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/19/20 18:54	
Chlorobenzene	5.0 U	5.0	0.20	1	07/19/20 18:54	
Chloroethane	5.0 U	5.0	0.23	1	07/19/20 18:54	
Chloroform	5.0 U	5.0	0.24	1	07/19/20 18:54	
Chloromethane	5.0 U	5.0	0.28	1	07/19/20 18:54	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/19/20 18:54	
Dichloromethane	5.0 U	5.0	0.65	1	07/19/20 18:54	
Ethylbenzene	5.0 U	5.0	0.20	1	07/19/20 18:54	
Styrene	5.0 U	5.0	0.20	1	07/19/20 18:54	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/19/20 18:54	
Toluene	5.0 U	5.0	0.20	1	07/19/20 18:54	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/19/20 18:54	
Vinyl Acetate	10 U	10	1.1	1	07/19/20 18:54	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/19/20 18:54	
Xylenes, Total	5.0 U	5.0	0.23	1	07/19/20 18:54	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/19/20 18:54	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/19/20 18:54	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/19/20 18:54	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/19/20 18:54	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/06/20 00:00

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 TB-9954-070620-SG-006
 Units: ug/L

 Lab Code:
 R2005820-001
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	94	85 - 122	07/19/20 18:54	
Dibromofluoromethane	98	89 - 119	07/19/20 18:54	
Toluene-d8	101	87 - 121	07/19/20 18:54	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/06/20 00:00

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 TB-9954-070620-SG-006
 Units: ug/L

 Lab Code:
 R2005820-001
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/06/20 13:55

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070620-SG-029
 Units: ug/L

 Lab Code:
 R2005820-002
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/20/20 15:11	_
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/20/20 15:11	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/20/20 15:11	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/20/20 15:11	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/20/20 15:11	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/20/20 15:11	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/20/20 15:11	
2-Butanone (MEK)	10 U	10	0.78	1	07/20/20 15:11	
2-Hexanone	10 U	10	0.20	1	07/20/20 15:11	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/20/20 15:11	
Acetone	10 U	10	5.0	1	07/20/20 15:11	
Benzene	5.0 U	5.0	0.20	1	07/20/20 15:11	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/20/20 15:11	
Bromoform	5.0 U	5.0	0.25	1	07/20/20 15:11	
Bromomethane	5.0 U	5.0	0.70	1	07/20/20 15:11	
Carbon Disulfide	6.4 J	10	0.42	1	07/20/20 15:11	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/20/20 15:11	
Chlorobenzene	5.0 U	5.0	0.20	1	07/20/20 15:11	
Chloroethane	5.0 U	5.0	0.23	1	07/20/20 15:11	
Chloroform	5.0 U	5.0	0.24	1	07/20/20 15:11	
Chloromethane	5.0 U	5.0	0.28	1	07/20/20 15:11	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/20/20 15:11	
Dichloromethane	5.0 U	5.0	0.65	1	07/20/20 15:11	
Ethylbenzene	5.0 U	5.0	0.20	1	07/20/20 15:11	
Styrene	5.0 U	5.0	0.20	1	07/20/20 15:11	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/20/20 15:11	
Toluene	5.0 U	5.0	0.20	1	07/20/20 15:11	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/20/20 15:11	
Vinyl Acetate	10 U	10	1.1	1	07/20/20 15:11	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/20/20 15:11	
Xylenes, Total	5.0 U	5.0	0.23	1	07/20/20 15:11	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/20/20 15:11	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/20/20 15:11	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/20/20 15:11	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/20/20 15:11	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/06/20 13:55

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070620-SG-029
 Units: ug/L

 Lab Code:
 R2005820-002
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	97	85 - 122	07/20/20 15:11	
Dibromofluoromethane	98	89 - 119	07/20/20 15:11	
Toluene-d8	101	87 - 121	07/20/20 15:11	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/06/20 13:55

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070620-SG-029
 Units: ug/L

 Lab Code:
 R2005820-002
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown	1.21	6.8	J	
007446-09-5	Sulfur dioxide	1.32	303.7	JN	
000066-25-1	Hexanal	9.24	7.7	JN	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/07/20 10:30

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070720-SG-031
 Units: ug/L

 Lab Code:
 R2005820-003
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/19/20 19:16	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/19/20 19:16	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/19/20 19:16	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/19/20 19:16	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/19/20 19:16	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/19/20 19:16	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/19/20 19:16	
2-Butanone (MEK)	10 U	10	0.78	1	07/19/20 19:16	
2-Hexanone	10 U	10	0.20	1	07/19/20 19:16	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/19/20 19:16	
Acetone	10 U	10	5.0	1	07/19/20 19:16	
Benzene	5.0 U	5.0	0.20	1	07/19/20 19:16	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/19/20 19:16	
Bromoform	5.0 U	5.0	0.25	1	07/19/20 19:16	
Bromomethane	5.0 U	5.0	0.70	1	07/19/20 19:16	
Carbon Disulfide	10 U	10	0.42	1	07/19/20 19:16	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/19/20 19:16	
Chlorobenzene	5.0 U	5.0	0.20	1	07/19/20 19:16	
Chloroethane	5.0 U	5.0	0.23	1	07/19/20 19:16	
Chloroform	5.0 U	5.0	0.24	1	07/19/20 19:16	
Chloromethane	5.0 U	5.0	0.28	1	07/19/20 19:16	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/19/20 19:16	
Dichloromethane	5.0 U	5.0	0.65	1	07/19/20 19:16	
Ethylbenzene	5.0 U	5.0	0.20	1	07/19/20 19:16	
Styrene	5.0 U	5.0	0.20	1	07/19/20 19:16	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/19/20 19:16	
Toluene	5.0 U	5.0	0.20	1	07/19/20 19:16	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/19/20 19:16	
Vinyl Acetate	10 U	10	1.1	1	07/19/20 19:16	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/19/20 19:16	
Xylenes, Total	5.0 U	5.0	0.23	1	07/19/20 19:16	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/19/20 19:16	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/19/20 19:16	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/19/20 19:16	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/19/20 19:16	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/07/20 10:30

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070720-SG-031
 Units: ug/L

 Lab Code:
 R2005820-003
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	93	85 - 122	07/19/20 19:16	
Dibromofluoromethane	98	89 - 119	07/19/20 19:16	
Toluene-d8	99	87 - 121	07/19/20 19:16	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/07/20 10:30

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070720-SG-031
 Units: ug/L

 Lab Code:
 R2005820-003
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

Result

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/07/20 13:00

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070720-SG-032
 Units: ug/L

 Lab Code:
 R2005820-004
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/20/20 15:33	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/20/20 15:33	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/20/20 15:33	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/20/20 15:33	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/20/20 15:33	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/20/20 15:33	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/20/20 15:33	
2-Butanone (MEK)	10 U	10	0.78	1	07/20/20 15:33	
2-Hexanone	10 U	10	0.20	1	07/20/20 15:33	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/20/20 15:33	
Acetone	10 U	10	5.0	1	07/20/20 15:33	
Benzene	5.0 U	5.0	0.20	1	07/20/20 15:33	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/20/20 15:33	
Bromoform	5.0 U	5.0	0.25	1	07/20/20 15:33	
Bromomethane	5.0 U	5.0	0.70	1	07/20/20 15:33	
Carbon Disulfide	3.7 J	10	0.42	1	07/20/20 15:33	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/20/20 15:33	
Chlorobenzene	5.0 U	5.0	0.20	1	07/20/20 15:33	
Chloroethane	5.0 U	5.0	0.23	1	07/20/20 15:33	
Chloroform	5.0 U	5.0	0.24	1	07/20/20 15:33	
Chloromethane	5.0 U	5.0	0.28	1	07/20/20 15:33	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/20/20 15:33	
Dichloromethane	5.0 U	5.0	0.65	1	07/20/20 15:33	
Ethylbenzene	5.0 U	5.0	0.20	1	07/20/20 15:33	
Styrene	5.0 U	5.0	0.20	1	07/20/20 15:33	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/20/20 15:33	
Toluene	5.0 U	5.0	0.20	1	07/20/20 15:33	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/20/20 15:33	
Vinyl Acetate	10 U	10	1.1	1	07/20/20 15:33	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/20/20 15:33	
Xylenes, Total	5.0 U	5.0	0.23	1	07/20/20 15:33	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/20/20 15:33	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/20/20 15:33	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/20/20 15:33	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/20/20 15:33	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/07/20 13:00

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070720-SG-032
 Units: ug/L

 Lab Code:
 R2005820-004
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	98	85 - 122	07/20/20 15:33	
Dibromofluoromethane	104	89 - 119	07/20/20 15:33	
Toluene-d8	104	87 - 121	07/20/20 15:33	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/07/20 13:00

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070720-SG-032
 Units: ug/L

 Lab Code:
 R2005820-004
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Tentatively Identified Compounds

 CAS#
 Compound Identification
 RT
 ug/L
 Q

 007446-09-5
 Sulfur dioxide
 1.21
 1.4
 JN

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/07/20 12:15

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070720-SG-030
 Units: ug/L

 Lab Code:
 R2005820-005
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/20/20 15:55	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/20/20 15:55	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/20/20 15:55	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/20/20 15:55	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/20/20 15:55	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/20/20 15:55	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/20/20 15:55	
2-Butanone (MEK)	10 U	10	0.78	1	07/20/20 15:55	
2-Hexanone	10 U	10	0.20	1	07/20/20 15:55	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/20/20 15:55	
Acetone	10 U	10	5.0	1	07/20/20 15:55	
Benzene	5.0 U	5.0	0.20	1	07/20/20 15:55	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/20/20 15:55	
Bromoform	5.0 U	5.0	0.25	1	07/20/20 15:55	
Bromomethane	5.0 U	5.0	0.70	1	07/20/20 15:55	
Carbon Disulfide	3.4 J	10	0.42	1	07/20/20 15:55	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/20/20 15:55	
Chlorobenzene	5.0 U	5.0	0.20	1	07/20/20 15:55	
Chloroethane	5.0 U	5.0	0.23	1	07/20/20 15:55	
Chloroform	5.0 U	5.0	0.24	1	07/20/20 15:55	
Chloromethane	5.0 U	5.0	0.28	1	07/20/20 15:55	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/20/20 15:55	
Dichloromethane	5.0 U	5.0	0.65	1	07/20/20 15:55	
Ethylbenzene	5.0 U	5.0	0.20	1	07/20/20 15:55	
Styrene	5.0 U	5.0	0.20	1	07/20/20 15:55	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/20/20 15:55	
Toluene	5.0 U	5.0	0.20	1	07/20/20 15:55	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/20/20 15:55	
Vinyl Acetate	10 U	10	1.1	1	07/20/20 15:55	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/20/20 15:55	
Xylenes, Total	5.0 U	5.0	0.23	1	07/20/20 15:55	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/20/20 15:55	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/20/20 15:55	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/20/20 15:55	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/20/20 15:55	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/07/20 12:15

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070720-SG-030
 Units: ug/L

 Lab Code:
 R2005820-005
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	95	85 - 122	07/20/20 15:55	
Dibromofluoromethane	100	89 - 119	07/20/20 15:55	
Toluene-d8	101	87 - 121	07/20/20 15:55	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/07/20 12:15

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070720-SG-030
 Units: ug/L

 Lab Code:
 R2005820-005
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Tentatively Identified Compounds



Semivolatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/06/20 13:55

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070620-SG-029
 Units: ug/L

 Lab Code:
 R2005820-002
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/14/20 18:03 7/13/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/14/20 18:03 7/13/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/14/20 18:03 7/13/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/14/20 18:03 7/13/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/14/20 18:03 7/13/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/14/20 18:03 7/13/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/14/20 18:03 7/13/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/14/20 18:03 7/13/20	
2,4-Dinitrophenol	45 U	45	19	1	07/14/20 18:03 7/13/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/14/20 18:03 7/13/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/14/20 18:03 7/13/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/14/20 18:03 7/13/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/14/20 18:03 7/13/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/14/20 18:03 7/13/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/14/20 18:03 7/13/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/14/20 18:03 7/13/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/14/20 18:03 7/13/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/14/20 18:03 7/13/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/14/20 18:03 7/13/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/14/20 18:03 7/13/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/14/20 18:03 7/13/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/14/20 18:03 7/13/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/14/20 18:03 7/13/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/14/20 18:03 7/13/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/14/20 18:03 7/13/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/14/20 18:03 7/13/20	
4-Nitrophenol	45 U	45	5.8	1	07/14/20 18:03 7/13/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/14/20 18:03 7/13/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/14/20 18:03 7/13/20	
Anthracene	9.1 U	9.1	1.2	1	07/14/20 18:03 7/13/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/14/20 18:03 7/13/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/14/20 18:03 7/13/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/14/20 18:03 7/13/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/14/20 18:03 7/13/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/14/20 18:03 7/13/20	
Benzoic Acid	91 U	91	33	1	07/14/20 18:03 7/13/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/14/20 18:03 7/13/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/14/20 18:03 7/13/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/14/20 18:03 7/13/20 07/14/20 18:03 7/13/20	
Bis(2-chloroethyl) Ether	9.1 U 9.1 U	9.1	1.3	1	07/14/20 18:03	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	07/14/20 18:03	
	9.1 U 9.1 U	9.1 9.1	1.3	1	07/14/20 18:03	
Butyl Benzyl Phthalate						
Chrysene	9.1 U	9.1	1.1	1	07/14/20 18:03 7/13/20	

Printed 7/27/2020 10:42:24 AM

Superset Reference: 20-0000555771 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820 **Date Collected:** 07/06/20 13:55

Love Canal:292-402-D02-3100/9954 Annual Long Term **Project:**

Monitoring

Date Received: 07/08/20 10:20 **Sample Matrix:** Water

Sample Name: WG-9954-070620-SG-029 Units: ug/L Lab Code: R2005820-002 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/14/20 18:03	7/13/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/14/20 18:03	7/13/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/14/20 18:03	7/13/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/14/20 18:03	7/13/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/14/20 18:03	7/13/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/14/20 18:03	7/13/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/14/20 18:03	7/13/20	
Fluorene	9.1 U	9.1	1.2	1	07/14/20 18:03	7/13/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/14/20 18:03	7/13/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/14/20 18:03	7/13/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/14/20 18:03	7/13/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/14/20 18:03	7/13/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/14/20 18:03	7/13/20	
Isophorone	9.1 U	9.1	1.3	1	07/14/20 18:03	7/13/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/14/20 18:03	7/13/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/14/20 18:03	7/13/20	
Naphthalene	9.1 U	9.1	1.1	1	07/14/20 18:03	7/13/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/14/20 18:03	7/13/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/14/20 18:03	7/13/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/14/20 18:03	7/13/20	
Phenol	9.1 U	9.1	0.91	1	07/14/20 18:03	7/13/20	
Pvrene	9.1 U	9.1	1.3	1	07/14/20 18:03	7/13/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	86	35 - 141	07/14/20 18:03	
2-Fluorobiphenyl	48	31 - 118	07/14/20 18:03	
2-Fluorophenol	36	10 - 105	07/14/20 18:03	
Nitrobenzene-d5	46	31 - 110	07/14/20 18:03	
Phenol-d6	30	10 - 107	07/14/20 18:03	
p-Terphenyl-d14	41	10 - 165	07/14/20 18:03	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
013798-23-7	Sulfur	7.82	13	JN	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/07/20 10:30

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070720-SG-031
 Units: ug/L

 Lab Code:
 R2005820-003
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/14/20 18:32	7/13/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/14/20 18:32	7/13/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/14/20 18:32	7/13/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/14/20 18:32	7/13/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/14/20 18:32	7/13/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/14/20 18:32	7/13/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/14/20 18:32	7/13/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/14/20 18:32	7/13/20	
2,4-Dinitrophenol	45 U	45	19	1	07/14/20 18:32	7/13/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/14/20 18:32	7/13/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/14/20 18:32	7/13/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/14/20 18:32	7/13/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/14/20 18:32	7/13/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/14/20 18:32	7/13/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/14/20 18:32	7/13/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/14/20 18:32	7/13/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/14/20 18:32	7/13/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/14/20 18:32	7/13/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/14/20 18:32	7/13/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/14/20 18:32	7/13/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/14/20 18:32	7/13/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/14/20 18:32	7/13/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/14/20 18:32	7/13/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/14/20 18:32	7/13/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/14/20 18:32	7/13/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/14/20 18:32	7/13/20	
4-Nitrophenol	45 U	45	5.8	1	07/14/20 18:32	7/13/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/14/20 18:32	7/13/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/14/20 18:32	7/13/20	
Anthracene	9.1 U	9.1	1.2	1	07/14/20 18:32	7/13/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/14/20 18:32	7/13/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/14/20 18:32	7/13/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/14/20 18:32	7/13/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/14/20 18:32	7/13/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/14/20 18:32	7/13/20	
Benzoic Acid	9.1 U	9.1	33	1	07/14/20 18:32	7/13/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/14/20 18:32	7/13/20	
2,2'-Oxybis(1-chloropropane)	9.1 U 9.1 U	9.1 9.1	1.3	1	07/14/20 18:32	7/13/20	
Bis(2-chloroethoxy)methane	9.1 U 9.1 U	9.1	1.3	1	07/14/20 18:32	7/13/20	
			1.8				
Bis(2-chloroethyl) Ether	9.1 U 9.1 U	9.1 9.1		1	07/14/20 18:32	7/13/20	
Bis(2-ethylhexyl) Phthalate			0.91	1	07/14/20 18:32	7/13/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	07/14/20 18:32	7/13/20	
Chrysene	9.1 U	9.1	1.1	1	07/14/20 18:32	7/13/20	

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Superset Reference: 20-0000555771 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Date Collected: 07/07/20 10:30

Monitoring Sample Matrix: Water

Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070720-SG-031
 Units: ug/L

 Lab Code:
 R2005820-003
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed Date Extra	acted Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/14/20 18:32 7/13/2	.0
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/14/20 18:32 7/13/2	.0
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/14/20 18:32 7/13/2	0
Dibenzofuran	9.1 U	9.1	1.3	1	07/14/20 18:32 7/13/2	0
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/14/20 18:32 7/13/2	0
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/14/20 18:32 7/13/2	.0
Fluoranthene	9.1 U	9.1	1.4	1	07/14/20 18:32 7/13/2	0
Fluorene	9.1 U	9.1	1.2	1	07/14/20 18:32 7/13/2	.0
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/14/20 18:32 7/13/2	.0
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/14/20 18:32 7/13/2	0
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/14/20 18:32 7/13/2	0
Hexachloroethane	9.1 U	9.1	0.96	1	07/14/20 18:32 7/13/2	0
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/14/20 18:32 7/13/2	0
Isophorone	9.1 U	9.1	1.3	1	07/14/20 18:32 7/13/2	0
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/14/20 18:32 7/13/2	0
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/14/20 18:32 7/13/2	0
Naphthalene	9.1 U	9.1	1.1	1	07/14/20 18:32 7/13/2	0
Nitrobenzene	9.1 U	9.1	1.4	1	07/14/20 18:32 7/13/2	0
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/14/20 18:32 7/13/2	0
Phenanthrene	9.1 U	9.1	1.3	1	07/14/20 18:32 7/13/2	0
Phenol	9.1 U	9.1	0.91	1	07/14/20 18:32 7/13/2	0
Pyrene	9.1 U	9.1	1.3	1	07/14/20 18:32 7/13/2	0

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	63	35 - 141	07/14/20 18:32	
2-Fluorobiphenyl	52	31 - 118	07/14/20 18:32	
2-Fluorophenol	36	10 - 105	07/14/20 18:32	
Nitrobenzene-d5	45	31 - 110	07/14/20 18:32	
Phenol-d6	26	10 - 107	07/14/20 18:32	
p-Terphenyl-d14	54	10 - 165	07/14/20 18:32	

Tentatively Identified Compounds

			Result	
CAS#	Compound Identification	RT	ug/L	Q

No Tentatively Identified Compounds

Detected

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/07/20 13:00

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070720-SG-032
 Units: ug/L

 Lab Code:
 R2005820-004
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/14/20 19:00	7/13/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/14/20 19:00	7/13/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/14/20 19:00	7/13/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/14/20 19:00	7/13/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/14/20 19:00	7/13/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/14/20 19:00	7/13/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/14/20 19:00	7/13/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/14/20 19:00	7/13/20	
2,4-Dinitrophenol	45 U	45	19	1	07/14/20 19:00	7/13/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/14/20 19:00	7/13/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/14/20 19:00	7/13/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/14/20 19:00	7/13/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/14/20 19:00	7/13/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/14/20 19:00	7/13/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/14/20 19:00	7/13/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/14/20 19:00	7/13/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/14/20 19:00	7/13/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/14/20 19:00	7/13/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/14/20 19:00	7/13/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/14/20 19:00	7/13/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/14/20 19:00	7/13/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/14/20 19:00	7/13/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/14/20 19:00	7/13/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/14/20 19:00	7/13/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/14/20 19:00	7/13/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/14/20 19:00	7/13/20	
4-Nitrophenol	45 U	45	5.8	1	07/14/20 19:00	7/13/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/14/20 19:00	7/13/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/14/20 19:00	7/13/20	
Anthracene	9.1 U	9.1	1.3	1	07/14/20 19:00	7/13/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/14/20 19:00	7/13/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/14/20 19:00	7/13/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/14/20 19:00	7/13/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/14/20 19:00	7/13/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/14/20 19:00	7/13/20	
Benzoic Acid	9.1 U	9.1	33	1	07/14/20 19:00	7/13/20	
	9.1 U	9.1	33 1.5	1	07/14/20 19:00	7/13/20	
Benzyl Alcohol	9.1 U 9.1 U	9.1 9.1	1.3	1	07/14/20 19:00	7/13/20	
2,2'-Oxybis(1-chloropropane)			1.3				
Bis(2-chloroethoxy)methane	9.1 U	9.1		1	07/14/20 19:00	7/13/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	07/14/20 19:00	7/13/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	07/14/20 19:00	7/13/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	07/14/20 19:00	7/13/20	
Chrysene	9.1 U	9.1	1.1	1	07/14/20 19:00	7/13/20	

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Superset Reference: 20-0000555771 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/07/20 13:00

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070720-SG-032
 Units: ug/L

 Lab Code:
 R2005820-004
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/14/20 19:00	7/13/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/14/20 19:00	7/13/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/14/20 19:00	7/13/20	,
Dibenzofuran	9.1 U	9.1	1.3	1	07/14/20 19:00	7/13/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/14/20 19:00	7/13/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/14/20 19:00	7/13/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/14/20 19:00	7/13/20	
Fluorene	9.1 U	9.1	1.2	1	07/14/20 19:00	7/13/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/14/20 19:00	7/13/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/14/20 19:00	7/13/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/14/20 19:00	7/13/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/14/20 19:00	7/13/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/14/20 19:00	7/13/20	
Isophorone	9.1 U	9.1	1.3	1	07/14/20 19:00	7/13/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/14/20 19:00	7/13/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/14/20 19:00	7/13/20	
Naphthalene	9.1 U	9.1	1.1	1	07/14/20 19:00	7/13/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/14/20 19:00	7/13/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/14/20 19:00	7/13/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/14/20 19:00	7/13/20	
Phenol	9.1 U	9.1	0.91	1	07/14/20 19:00	7/13/20	
Pvrene	9.1 U	9.1	1.3	1	07/14/20 19:00	7/13/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	80	35 - 141	07/14/20 19:00	
2-Fluorobiphenyl	58	31 - 118	07/14/20 19:00	
2-Fluorophenol	39	10 - 105	07/14/20 19:00	
Nitrobenzene-d5	55	31 - 110	07/14/20 19:00	
Phenol-d6	27	10 - 107	07/14/20 19:00	
p-Terphenyl-d14	39	10 - 165	07/14/20 19:00	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown hydrocarbon	11.88	4.4	J	
	unknown hydrocarbon	12.48	5.7	J	
	unknown hydrocarbon	13.12	5.6	J	
	unknown hydrocarbon	13.82	4.9	J	
	unknown hydrocarbon	14.57	5.6	J	
	unknown hydrocarbon	16.10	4.4	J	
013798-23-7	Sulfur	7.78	4.2	JN	

Printed 7/27/2020 10:42:25 AM

Superset Reference:20-0000555771 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/07/20 12:15

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070720-SG-030
 Units: ug/L

 Lab Code:
 R2005820-005
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed D	ate Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/14/20 19:28	7/13/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/14/20 19:28	7/13/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/14/20 19:28	7/13/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/14/20 19:28	7/13/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/14/20 19:28	7/13/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/14/20 19:28	7/13/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/14/20 19:28	7/13/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/14/20 19:28	7/13/20	
2,4-Dinitrophenol	45 U	45	19	1	07/14/20 19:28	7/13/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/14/20 19:28	7/13/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/14/20 19:28	7/13/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/14/20 19:28	7/13/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/14/20 19:28	7/13/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/14/20 19:28	7/13/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/14/20 19:28	7/13/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/14/20 19:28	7/13/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/14/20 19:28	7/13/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/14/20 19:28	7/13/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/14/20 19:28	7/13/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/14/20 19:28	7/13/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/14/20 19:28	7/13/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/14/20 19:28	7/13/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/14/20 19:28	7/13/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/14/20 19:28	7/13/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/14/20 19:28	7/13/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/14/20 19:28	7/13/20	
4-Nitrophenol	45 U	45	5.8	1	07/14/20 19:28	7/13/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/14/20 19:28	7/13/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/14/20 19:28	7/13/20	
Anthracene	9.1 U	9.1	1.2	1	07/14/20 19:28	7/13/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/14/20 19:28	7/13/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/14/20 19:28	7/13/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/14/20 19:28	7/13/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/14/20 19:28	7/13/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/14/20 19:28	7/13/20	
Benzoic Acid	91 U	91	33	1	07/14/20 19:28	7/13/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/14/20 19:28	7/13/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/14/20 19:28	7/13/20	
Bis(2-chloroethoxy)methane	9.1 U 9.1 U	9.1 9.1	1.5	1	07/14/20 19:28	7/13/20	
Bis(2-chloroethyl) Ether	9.1 U 9.1 U	9.1 9.1	1.8	1	07/14/20 19:28	7/13/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	07/14/20 19:28	7/13/20	
	9.1 U 9.1 U	9.1 9.1	1.3	1	07/14/20 19:28	7/13/20	
Butyl Benzyl Phthalate							
Chrysene	9.1 U	9.1	1.1	1	07/14/20 19:28	7/13/20	

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Superset Reference: 20-0000555771 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/07/20 12:15

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070720-SG-030
 Units: ug/L

 Lab Code:
 R2005820-005
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/14/20 19:28	7/13/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/14/20 19:28	7/13/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/14/20 19:28	7/13/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/14/20 19:28	7/13/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/14/20 19:28	7/13/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/14/20 19:28	7/13/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/14/20 19:28	7/13/20	
Fluorene	9.1 U	9.1	1.2	1	07/14/20 19:28	7/13/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/14/20 19:28	7/13/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/14/20 19:28	7/13/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/14/20 19:28	7/13/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/14/20 19:28	7/13/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/14/20 19:28	7/13/20	
Isophorone	9.1 U	9.1	1.3	1	07/14/20 19:28	7/13/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/14/20 19:28	7/13/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/14/20 19:28	7/13/20	
Naphthalene	9.1 U	9.1	1.1	1	07/14/20 19:28	7/13/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/14/20 19:28	7/13/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/14/20 19:28	7/13/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/14/20 19:28	7/13/20	
Phenol	9.1 U	9.1	0.91	1	07/14/20 19:28	7/13/20	
Pyrene	9.1 U	9.1	1.3	1	07/14/20 19:28	7/13/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	79	35 - 141	07/14/20 19:28	
2-Fluorobiphenyl	52	31 - 118	07/14/20 19:28	
2-Fluorophenol	35	10 - 105	07/14/20 19:28	
Nitrobenzene-d5	52	31 - 110	07/14/20 19:28	
Phenol-d6	26	10 - 107	07/14/20 19:28	
p-Terphenyl-d14	50	10 - 165	07/14/20 19:28	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown hydrocarbon	11.88	6.0	J	
	unknown hydrocarbon	12.47	7.3	J	
	unknown hydrocarbon	13.12	8.3	J	
	unknown hydrocarbon	13.82	6.8	J	
	unknown hydrocarbon	14.57	6.0	J	
	unknown hydrocarbon	15.35	4.6	J	
	unknown	2.87	11	J	

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Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/07/20 12:15

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070720-SG-030
 Units: ug/L

 Lab Code:
 R2005820-005
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Tentatively Identified Compounds



Semivolatile Organic Compounds by GC

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/06/20 13:55

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070620-SG-029
 Units: ug/L

 Lab Code:
 R2005820-002
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/14/20 09:49	7/10/20	_
4,4'-DDE	0.045 U	0.045	0.019	1	07/14/20 09:49	7/10/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/14/20 09:49	7/10/20	
Aldrin	0.045 U	0.045	0.019	1	07/14/20 09:49	7/10/20	
Dieldrin	0.045 U	0.045	0.019	1	07/14/20 09:49	7/10/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/14/20 09:49	7/10/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/14/20 09:49	7/10/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/14/20 09:49	7/10/20	
Endrin	0.045 U	0.045	0.019	1	07/14/20 09:49	7/10/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/14/20 09:49	7/10/20	
Heptachlor	0.045 U	0.045	0.019	1	07/14/20 09:49	7/10/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/14/20 09:49	7/10/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/14/20 09:49	7/10/20	
Toxaphene	0.46 U	0.46	0.46	1	07/14/20 09:49	7/10/20	
alpha-BHC	0.47	0.045	0.019	1	07/14/20 09:49	7/10/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/14/20 09:49	7/10/20	
beta-BHC	0.045 U	0.045	0.019	1	07/14/20 09:49	7/10/20	
delta-BHC	0.13	0.045	0.019	1	07/14/20 09:49	7/10/20	
gamma-BHC (Lindane)	0.48	0.045	0.019	1	07/14/20 09:49	7/10/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/14/20 09:49	7/10/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	51	10 - 164	07/14/20 09:49	
Tetrachloro-m-xylene	63	10 - 147	07/14/20 09:49	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/07/20 10:30

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070720-SG-031
 Units: ug/L

 Lab Code:
 R2005820-003
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/13/20 21:05	7/10/20	
4,4'-DDE	0.045 U	0.045	0.019	1	07/13/20 21:05	7/10/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/13/20 21:05	7/10/20	
Aldrin	0.045 U	0.045	0.019	1	07/13/20 21:05	7/10/20	
Dieldrin	0.045 U	0.045	0.019	1	07/13/20 21:05	7/10/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/13/20 21:05	7/10/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/13/20 21:05	7/10/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/13/20 21:05	7/10/20	
Endrin	0.045 U	0.045	0.019	1	07/13/20 21:05	7/10/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/13/20 21:05	7/10/20	
Heptachlor	0.045 U	0.045	0.019	1	07/13/20 21:05	7/10/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/13/20 21:05	7/10/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/13/20 21:05	7/10/20	
Toxaphene	0.46 U	0.46	0.46	1	07/13/20 21:05	7/10/20	
alpha-BHC	0.045 U	0.045	0.019	1	07/13/20 21:05	7/10/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/13/20 21:05	7/10/20	
beta-BHC	0.045 U	0.045	0.019	1	07/13/20 21:05	7/10/20	
delta-BHC	0.045 U	0.045	0.019	1	07/13/20 21:05	7/10/20	
gamma-BHC (Lindane)	0.045 U	0.045	0.019	1	07/13/20 21:05	7/10/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/13/20 21:05	7/10/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	61	10 - 164	07/13/20 21:05	
Tetrachloro-m-xylene	56	10 - 147	07/13/20 21:05	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/07/20 13:00

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070720-SG-032
 Units: ug/L

 Lab Code:
 R2005820-004
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/14/20 10:08	7/10/20	
4,4'-DDE	0.045 U	0.045	0.019	1	07/14/20 10:08	7/10/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/14/20 10:08	7/10/20	
Aldrin	0.045 U	0.045	0.019	1	07/14/20 10:08	7/10/20	
Dieldrin	0.045 U	0.045	0.019	1	07/14/20 10:08	7/10/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/14/20 10:08	7/10/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/14/20 10:08	7/10/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/14/20 10:08	7/10/20	
Endrin	0.045 U	0.045	0.019	1	07/14/20 10:08	7/10/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/14/20 10:08	7/10/20	
Heptachlor	0.045 U	0.045	0.019	1	07/14/20 10:08	7/10/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/14/20 10:08	7/10/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/14/20 10:08	7/10/20	
Toxaphene	0.46 U	0.46	0.46	1	07/14/20 10:08	7/10/20	
alpha-BHC	0.070	0.045	0.019	1	07/14/20 10:08	7/10/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/14/20 10:08	7/10/20	
beta-BHC	0.045 U	0.045	0.019	1	07/14/20 10:08	7/10/20	
delta-BHC	0.067	0.045	0.019	1	07/14/20 10:08	7/10/20	
gamma-BHC (Lindane)	0.098	0.045	0.019	1	07/14/20 10:08	7/10/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/14/20 10:08	7/10/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	66	10 - 164	07/14/20 10:08	
Tetrachloro-m-xylene	63	10 - 147	07/14/20 10:08	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/07/20 12:15

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070720-SG-030
 Units: ug/L

 Lab Code:
 R2005820-005
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/14/20 10:27	7/10/20	
4,4'-DDE	0.045 U	0.045	0.019	1	07/14/20 10:27	7/10/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/14/20 10:27	7/10/20	
Aldrin	0.045 U	0.045	0.019	1	07/14/20 10:27	7/10/20	
Dieldrin	0.045 U	0.045	0.019	1	07/14/20 10:27	7/10/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/14/20 10:27	7/10/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/14/20 10:27	7/10/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/14/20 10:27	7/10/20	
Endrin	0.045 U	0.045	0.019	1	07/14/20 10:27	7/10/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/14/20 10:27	7/10/20	
Heptachlor	0.045 U	0.045	0.019	1	07/14/20 10:27	7/10/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/14/20 10:27	7/10/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/14/20 10:27	7/10/20	
Toxaphene	0.46 U	0.46	0.46	1	07/14/20 10:27	7/10/20	
alpha-BHC	0.24	0.045	0.019	1	07/14/20 10:27	7/10/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/14/20 10:27	7/10/20	
beta-BHC	0.045 U	0.045	0.019	1	07/14/20 10:27	7/10/20	
delta-BHC	0.084	0.045	0.019	1	07/14/20 10:27	7/10/20	
gamma-BHC (Lindane)	0.26	0.045	0.019	1	07/14/20 10:27	7/10/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/14/20 10:27	7/10/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	77	10 - 164	07/14/20 10:27	
Tetrachloro-m-xylene	73	10 - 147	07/14/20 10:27	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/06/20 13:55

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070620-SG-029
 Units: ug/L

 Lab Code:
 R2005820-002
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/14/20 17:02	7/10/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/14/20 17:02	7/10/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/14/20 17:02	7/10/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/14/20 17:02	7/10/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/14/20 17:02	7/10/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/14/20 17:02	7/10/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/14/20 17:02	7/10/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q	
Decachlorobiphenyl	50	10 - 152	07/14/20 17:02		
Tetrachloro-m-xvlene	54	14 - 129	07/14/20 17:02		

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/07/20 10:30

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070720-SG-031
 Units: ug/L

 Lab Code:
 R2005820-003
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/13/20 19:13	7/10/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/13/20 19:13	7/10/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/13/20 19:13	7/10/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/13/20 19:13	7/10/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/13/20 19:13	7/10/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/13/20 19:13	7/10/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/13/20 19:13	7/10/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	59	10 - 152	07/13/20 19:13	
Tetrachloro-m-xylene	47	14 - 129	07/13/20 19:13	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/07/20 13:00

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070720-SG-032
 Units: ug/L

 Lab Code:
 R2005820-004
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/14/20 17:22	7/10/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/14/20 17:22	7/10/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/14/20 17:22	7/10/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/14/20 17:22	7/10/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/14/20 17:22	7/10/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/14/20 17:22	7/10/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/14/20 17:22	7/10/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	63	10 - 152	07/14/20 17:22	
Tetrachloro-m-xvlene	52	14 - 129	07/14/20 17:22	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/07/20 12:15

Monitoring

Sample Matrix: Water Date Received: 07/08/20 10:20

 Sample Name:
 WG-9954-070720-SG-030
 Units: ug/L

 Lab Code:
 R2005820-005
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/16/20 14:40	7/10/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/16/20 14:40	7/10/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/16/20 14:40	7/10/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/16/20 14:40	7/10/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/16/20 14:40	7/10/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/16/20 14:40	7/10/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/16/20 14:40	7/10/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	68	10 - 152	07/16/20 14:40	
Tetrachloro-m-xylene	55	14 - 129	07/16/20 14:40	



QC Summary Forms

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com



Volatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

SURROGATE RECOVERY SUMMARY Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Extraction Method: EPA 5030C

		4-Bromofluorobenzene	Dibromofluoromethane	Toluene-d8
Sample Name	Lab Code	85-122	89-119	87-121
TB-9954-070620-SG-006	R2005820-001	94	98	101
WG-9954-070620-SG-029	R2005820-002	97	98	101
WG-9954-070720-SG-031	R2005820-003	93	98	99
WG-9954-070720-SG-032	R2005820-004	98	104	104
WG-9954-070720-SG-030	R2005820-005	95	100	101
Method Blank	RQ2007740-04	95	99	98
Method Blank	RQ2007775-04	94	98	100
Lab Control Sample	RQ2007740-03	99	100	102
Lab Control Sample	RQ2007775-03	101	100	105

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007740-04
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/19/20 13:26	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/19/20 13:26	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/19/20 13:26	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/19/20 13:26	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/19/20 13:26	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/19/20 13:26	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/19/20 13:26	
2-Butanone (MEK)	10 U	10	0.78	1	07/19/20 13:26	
2-Hexanone	10 U	10	0.20	1	07/19/20 13:26	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/19/20 13:26	
Acetone	10 U	10	5.0	1	07/19/20 13:26	
Benzene	5.0 U	5.0	0.20	1	07/19/20 13:26	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/19/20 13:26	
Bromoform	5.0 U	5.0	0.25	1	07/19/20 13:26	
Bromomethane	5.0 U	5.0	0.70	1	07/19/20 13:26	
Carbon Disulfide	10 U	10	0.42	1	07/19/20 13:26	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/19/20 13:26	
Chlorobenzene	5.0 U	5.0	0.20	1	07/19/20 13:26	
Chloroethane	5.0 U	5.0	0.23	1	07/19/20 13:26	
Chloroform	5.0 U	5.0	0.24	1	07/19/20 13:26	
Chloromethane	5.0 U	5.0	0.28	1	07/19/20 13:26	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/19/20 13:26	
Dichloromethane	5.0 U	5.0	0.65	1	07/19/20 13:26	
Ethylbenzene	5.0 U	5.0	0.20	1	07/19/20 13:26	
Styrene	5.0 U	5.0	0.20	1	07/19/20 13:26	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/19/20 13:26	
Toluene	5.0 U	5.0	0.20	1	07/19/20 13:26	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/19/20 13:26	
Vinyl Acetate	10 U	10	1.1	1	07/19/20 13:26	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/19/20 13:26	
Xylenes, Total	5.0 U	5.0	0.23	1	07/19/20 13:26	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/19/20 13:26	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/19/20 13:26	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/19/20 13:26	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/19/20 13:26	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007740-04Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	95	85 - 122	07/19/20 13:26	
Dibromofluoromethane	99	89 - 119	07/19/20 13:26	
Toluene-d8	98	87 - 121	07/19/20 13:26	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007740-04Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Tentatively Identified Compounds

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007775-04
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/20/20 12:31	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/20/20 12:31	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/20/20 12:31	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/20/20 12:31	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/20/20 12:31	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/20/20 12:31	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/20/20 12:31	
2-Butanone (MEK)	10 U	10	0.78	1	07/20/20 12:31	
2-Hexanone	10 U	10	0.20	1	07/20/20 12:31	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/20/20 12:31	
Acetone	10 U	10	5.0	1	07/20/20 12:31	
Benzene	5.0 U	5.0	0.20	1	07/20/20 12:31	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/20/20 12:31	
Bromoform	5.0 U	5.0	0.25	1	07/20/20 12:31	
Bromomethane	5.0 U	5.0	0.70	1	07/20/20 12:31	
Carbon Disulfide	10 U	10	0.42	1	07/20/20 12:31	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/20/20 12:31	
Chlorobenzene	5.0 U	5.0	0.20	1	07/20/20 12:31	
Chloroethane	5.0 U	5.0	0.23	1	07/20/20 12:31	
Chloroform	5.0 U	5.0	0.24	1	07/20/20 12:31	
Chloromethane	5.0 U	5.0	0.28	1	07/20/20 12:31	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/20/20 12:31	
Dichloromethane	5.0 U	5.0	0.65	1	07/20/20 12:31	
Ethylbenzene	5.0 U	5.0	0.20	1	07/20/20 12:31	
Styrene	5.0 U	5.0	0.20	1	07/20/20 12:31	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/20/20 12:31	
Toluene	5.0 U	5.0	0.20	1	07/20/20 12:31	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/20/20 12:31	
Vinyl Acetate	10 U	10	1.1	1	07/20/20 12:31	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/20/20 12:31	
Xylenes, Total	5.0 U	5.0	0.23	1	07/20/20 12:31	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/20/20 12:31	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/20/20 12:31	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/20/20 12:31	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/20/20 12:31	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007775-04Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	94	85 - 122	07/20/20 12:31	
Dibromofluoromethane	98	89 - 119	07/20/20 12:31	
Toluene-d8	100	87 - 121	07/20/20 12:31	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007775-04Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Tentatively Identified Compounds

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/19/20

Sample Matrix: Water

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005820

Lab Control Sample

RQ2007740-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits		
1,1,1-Trichloroethane (TCA)	8260C	21.2	20.0	106	75-125		
1,1,2,2-Tetrachloroethane	8260C	22.4	20.0	112	78-126		
1,1,2-Trichloroethane	8260C	21.0	20.0	105	82-121		
1,1-Dichloroethane (1,1-DCA)	8260C	20.8	20.0	104	80-124		
1,1-Dichloroethene (1,1-DCE)	8260C	25.1	20.0	125 *	71-118		
1,2-Dichloroethane	8260C	20.0	20.0	100	71-127		
1,2-Dichloropropane	8260C	21.2	20.0	106	80-119		
2-Butanone (MEK)	8260C	21.6	20.0	108	61-137		
2-Hexanone	8260C	19.1	20.0	96	63-124		
4-Methyl-2-pentanone	8260C	19.7	20.0	99	66-124		
Acetone	8260C	23.0	20.0	115	40-161		
Benzene	8260C	21.1	20.0	106	79-119		
Bromodichloromethane	8260C	20.0	20.0	100	81-123		
Bromoform	8260C	20.4	20.0	102	65-146		
Bromomethane	8260C	18.0	20.0	90	42-166		
Carbon Disulfide	8260C	21.3	20.0	106	66-128		
Carbon Tetrachloride	8260C	21.7	20.0	109	70-127		
Chlorobenzene	8260C	20.8	20.0	104	80-121		
Chloroethane	8260C	20.8	20.0	104	62-131		
Chloroform	8260C	19.9	20.0	100	79-120		
Chloromethane	8260C	24.0	20.0	120	65-135		
Dibromochloromethane	8260C	21.9	20.0	109	72-128		
Dichloromethane	8260C	20.8	20.0	104	73-122		
Ethylbenzene	8260C	21.5	20.0	108	76-120		
Styrene	8260C	21.3	20.0	106	80-124		
Tetrachloroethene (PCE)	8260C	21.2	20.0	106	72-125		
Toluene	8260C	21.2	20.0	106	79-119		
Trichloroethene (TCE)	8260C	19.4	20.0	97	74-122		
Vinyl Acetate	8260C	30.1	20.0	151	52-174		
Vinyl Chloride	8260C	24.3	20.0	121	74-159		
cis-1,2-Dichloroethene	8260C	20.2	20.0	101	80-121		
cis-1,3-Dichloropropene	8260C	20.3	20.0	101	77-122		
trans-1,2-Dichloroethene	8260C	22.8	20.0	114	73-118		
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QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix:

Project:

Water

Service Request: R2005820 **Date Analyzed:** 07/19/20

Date Analyzeu: 07/1

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

> Units:ug/L Basis:NA

Lab Control Sample

RQ2007740-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
trans-1.3-Dichloropropene	8260C	19.9	20.0	99	71-133

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/20/20

Sample Matrix: Water

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005820

Lab Control Sample

RQ2007775-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
1,1,1-Trichloroethane (TCA)	8260C	20.2	20.0	101	75-125
1,1,2,2-Tetrachloroethane	8260C	22.7	20.0	114	78-126
1,1,2-Trichloroethane	8260C	21.1	20.0	105	82-121
1,1-Dichloroethane (1,1-DCA)	8260C	21.1	20.0	105	80-124
1,1-Dichloroethene (1,1-DCE)	8260C	23.9	20.0	119 *	71-118
1,2-Dichloroethane	8260C	19.6	20.0	98	71-127
1,2-Dichloropropane	8260C	20.8	20.0	104	80-119
2-Butanone (MEK)	8260C	21.8	20.0	109	61-137
2-Hexanone	8260C	19.5	20.0	98	63-124
4-Methyl-2-pentanone	8260C	20.3	20.0	101	66-124
Acetone	8260C	23.4	20.0	117	40-161
Benzene	8260C	20.8	20.0	104	79-119
Bromodichloromethane	8260C	19.8	20.0	99	81-123
Bromoform	8260C	19.4	20.0	97	65-146
Bromomethane	8260C	18.2	20.0	91	42-166
Carbon Disulfide	8260C	21.6	20.0	108	66-128
Carbon Tetrachloride	8260C	19.8	20.0	99	70-127
Chlorobenzene	8260C	20.5	20.0	103	80-121
Chloroethane	8260C	22.3	20.0	112	62-131
Chloroform	8260C	20.3	20.0	101	79-120
Chloromethane	8260C	25.0	20.0	125	65-135
Dibromochloromethane	8260C	21.8	20.0	109	72-128
Dichloromethane	8260C	20.8	20.0	104	73-122
Ethylbenzene	8260C	20.3	20.0	101	76-120
Styrene	8260C	20.6	20.0	103	80-124
Tetrachloroethene (PCE)	8260C	19.5	20.0	98	72-125
Toluene	8260C	20.6	20.0	103	79-119
Trichloroethene (TCE)	8260C	18.1	20.0	90	74-122
Vinyl Acetate	8260C	31.4	20.0	157	52-174
Vinyl Chloride	8260C	25.3	20.0	126	74-159
cis-1,2-Dichloroethene	8260C	20.9	20.0	105	80-121
cis-1,3-Dichloropropene	8260C	19.6	20.0	98	77-122
trans-1,2-Dichloroethene	8260C	23.8	20.0	119 *	73-118
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QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix: Water

Project:

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

> Units:ug/L Basis:NA

Service Request: R2005820

Date Analyzed: 07/20/20

Lab Control Sample

RQ2007775-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
trans-1.3-Dichloropropene	8260C	19.3	20.0	97	71-133



Semivolatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

QA/QC Report

Service Request: R2005820

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

Project:

SURROGATE RECOVERY SUMMARY Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Extraction Method:** EPA 3510C

		2,4,6-Tribromophenol	2-Fluorobiphenyl	2-Fluorophenol
Sample Name	Lab Code	35-141	31-118	10-105
WG-9954-070620-SG-029	R2005820-002	86	48	36
WG-9954-070720-SG-031	R2005820-003	63	52	36
WG-9954-070720-SG-032	R2005820-004	80	58	39
WG-9954-070720-SG-030	R2005820-005	79	52	35
Method Blank	RQ2007408-01	67	54	45
Lab Control Sample	RQ2007408-02	82	64	40
Duplicate Lab Control Sample	RQ2007408-03	89	70	42
WG-9954-070720-SG-030 MS	RQ2007408-04	94	72	41
WG-9954-070720-SG-030 DMS	RQ2007408-05	65	52	29

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

SURROGATE RECOVERY SUMMARY Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Extraction Method:** EPA 3510C

		Nitrobenzene-d5	Phenol-d6	p-Terphenyl-d14	
Sample Name	Lab Code	31-110	10-107	10-165	
WG-9954-070620-SG-029	R2005820-002	46	30	41	
WG-9954-070720-SG-031	R2005820-003	45	26	54	
WG-9954-070720-SG-032	R2005820-004	55	27	39	
WG-9954-070720-SG-030	R2005820-005	52	26	50	
Method Blank	RQ2007408-01	55	32	60	
Lab Control Sample	RQ2007408-02	65	32	58	
Duplicate Lab Control Sample	RQ2007408-03	63	32	62	
WG-9954-070720-SG-030 MS	RQ2007408-04	72	36	47	
WG-9954-070720-SG-030 DMS	RQ2007408-05	54	21	29	

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates) **Service Request:**

R2005820

Project:

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Date Collected:

07/07/20

Sample Matrix: Water

Date Received:

07/08/20

Date Analyzed: **Date Extracted:**

07/14/20 07/13/20

Duplicate Matrix Spike Summary Semivolatile Organic Compounds by GC/MS

Sample Name: WG-9954-070720-SG-030 **Units:**

ug/L

Lab Code:

R2005820-005

Basis:

NA

Analysis Method: Prep Method:

8270D

EPA 3510C

Matrix Spike

Duplicate Matrix Spike

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Results flagged with an asterisk (*) indicate values outside control criteria.

Results flagged with a pound (#) indicate the control criteria is not applicable.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

Matrix Spike and Matrix Spike Duplicate Data is presented for information purposes only. The matrix may or may not be relevant to samples reported in this report. The laboratory evaluates system performance based on the LCS and LCSD control limits.

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates) **Service Request:**

R2005820 07/07/20

Project: Sample Matrix:

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Date Received: 07/08/20

Date Collected: Date Analyzed:

07/14/20

Date Extracted:

07/13/20

NA

Duplicate Matrix Spike Summary Semivolatile Organic Compounds by GC/MS

Sample Name: WG-9954-070720-SG-030

Water

Units: ug/L

Basis:

Lab Code:

R2005820-005

8270D

Analysis Method: Prep Method:

EPA 3510C

Matrix Spike Duplicate Matrix Spike RQ2007408-04 RQ2007408-05

		1102	007 100 01			Q2007 100 C	,,,			
	Sample		Spike			Spike		% Rec		RPD
Analyte Name	Result	Result	Amount	% Rec	Result	Amount	% Rec	Limits	RPD	Limit
Benzo(g,h,i)perylene	9.1 U	46.0	72.7	63	26.5	72.7	36 *	50-143	55*	30
Benzo(k)fluoranthene	9.1 U	44.2	72.7	61	25.7	72.7	35 *	46-139	54*	30
Benzoic Acid	91 U	79.9 J	109	73	56.3 J	109	52	10-94	34*	30
Benzyl Alcohol	9.1 U	52.0	72.7	72	35.1	72.7	48	31-109	40*	30
2,2'-Oxybis(1-chloropropane)	9.1 U	49.6	72.7	68	37.5	72.7	52	21-126	27	30
Bis(2-chloroethoxy)methane	9.1 U	55.7	72.7	77	40.7	72.7	56	41-118	32*	30
Bis(2-chloroethyl) Ether	9.1 U	46.8	72.7	64	31.6	72.7	43	33-108	39*	30
Bis(2-ethylhexyl) Phthalate	9.1 U	42.8	72.7	59	25.9	72.7	36 *	41-132	48*	30
Butyl Benzyl Phthalate	9.1 U	54.8	72.7	75	37.1	72.7	51	41-148	38*	30
Chrysene	9.1 U	46.3	72.7	64	27.6	72.7	38 *	47-126	51*	30
Di-n-butyl Phthalate	9.1 U	63.6	72.7	87	42.2	72.7	58	43-130	40*	30
Di-n-octyl Phthalate	9.1 U	42.0	72.7	58	27.2	72.7	37 *	40-139	44*	30
Dibenz(a,h)anthracene	9.1 U	49.5	72.7	68	28.4	72.7	39 *	43-136	54*	30
Dibenzofuran	9.1 U	60.2	72.7	83	42.1	72.7	58	46-119	35*	30
Diethyl Phthalate	9.1 U	58.4	72.7	80	45.1	72.7	62	36-122	25	30
Dimethyl Phthalate	9.1 U	63.4	72.7	87	44.9	72.7	62	33-123	34*	30
Fluoranthene	9.1 U	62.2	72.7	85	37.7	72.7	52	43-135	48*	30
Fluorene	9.1 U	60.3	72.7	83	41.9	72.7	58	43-113	35*	30
Hexachlorobenzene	9.1 U	54.3	72.7	75	34.7	72.7	48	42-125	44*	30
Hexachlorobutadiene	9.1 U	46.0	72.7	63	30.7	72.7	42	10-111	40*	30
Hexachlorocyclopentadiene	9.1 U	4.26 J	72.7	6 *	3.77 J	72.7	5 *	10-103	18	30
Hexachloroethane	9.1 U	41.8	72.7	57	30.8	72.7	42	12-101	30	30
Indeno(1,2,3-cd)pyrene	9.1 U	39.9	72.7	55	22.7	72.7	31 *	49-140	56*	30
Isophorone	9.1 U	46.4	72.7	64	34.7	72.7	48	40-111	29	30
N-Nitrosodi-n-propylamine	9.1 U	57.6	72.7	79	42.0	72.7	58	35-108	31*	30
N-Nitrosodiphenylamine	9.1 U	72.9	72.7	100	53.9	72.7	74	43-127	30	30
Naphthalene	9.1 U	50.7	72.7	70	36.4	72.7	50	37-108	33*	30
Nitrobenzene	9.1 U	50.2	72.7	69	38.0	72.7	52	35-112	28	30
Pentachlorophenol (PCP)	45 U	61.4	72.7	84	40.3 J	72.7	55	29-164	42*	30
Phenanthrene	9.1 U	59.0	72.7	81	39.9	72.7	55	46-123	38*	30
Phenol	9.1 U	27.4	72.7	38	17.5	72.7	24	10-113	45*	30
Pyrene	9.1 U	60.7	72.7	83	39.6	72.7	55	44-129	41*	30
-										

Results flagged with an asterisk (*) indicate values outside control criteria.

Results flagged with a pound (#) indicate the control criteria is not applicable.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

Matrix Spike and Matrix Spike Duplicate Data is presented for information purposes only. The matrix may or may not be relevant to samples reported in this report. The laboratory evaluates system performance based on the LCS and LCSD control limits.

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007408-01
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed Date Extracted	Q
1,2,4-Trichlorobenzene	10 U	10	1.2	1	07/14/20 16:10 7/13/20	
1,2-Dichlorobenzene	10 U	10	1.2	1	07/14/20 16:10 7/13/20	
1,3-Dichlorobenzene	10 U	10	1.1	1	07/14/20 16:10 7/13/20	
1,4-Dichlorobenzene	10 U	10	1.2	1	07/14/20 16:10 7/13/20	
2,4,5-Trichlorophenol	10 U	10	1.1	1	07/14/20 16:10 7/13/20	
2,4,6-Trichlorophenol	10 U	10	1.4	1	07/14/20 16:10 7/13/20	
2,4-Dichlorophenol	10 U	10	1.3	1	07/14/20 16:10 7/13/20	
2,4-Dimethylphenol	10 U	10	1.4	1	07/14/20 16:10 7/13/20	
2,4-Dinitrophenol	50 U	50	20	1	07/14/20 16:10 7/13/20	
2,4-Dinitrotoluene	10 U	10	2.4	1	07/14/20 16:10 7/13/20	
2,6-Dinitrotoluene	10 U	10	1.4	1	07/14/20 16:10 7/13/20	
2-Chloronaphthalene	10 U	10	1.4	1	07/14/20 16:10 7/13/20	
2-Chlorophenol	10 U	10	1.1	1	07/14/20 16:10 7/13/20	
2-Methylnaphthalene	10 U	10	1.3	1	07/14/20 16:10 7/13/20	
2-Methylphenol	10 U	10	1.0	1	07/14/20 16:10 7/13/20	
2-Nitroaniline	10 U	10	1.4	1	07/14/20 16:10 7/13/20	
2-Nitrophenol	10 U	10	1.5	1	07/14/20 16:10 7/13/20	
3,3'-Dichlorobenzidine	10 U	10	1.2	1	07/14/20 16:10 7/13/20	
3- and 4-Methylphenol Coelution	10 U	10	1.2	1	07/14/20 16:10 7/13/20	
3-Nitroaniline	10 U	10	2.5	1	07/14/20 16:10 7/13/20	
4,6-Dinitro-2-methylphenol	50 U	50	20	1	07/14/20 16:10 7/13/20	
4-Bromophenyl Phenyl Ether	10 U	10	1.7	1	07/14/20 16:10 7/13/20	
4-Chloro-3-methylphenol	10 U	10	1.1	1	07/14/20 16:10 7/13/20	
4-Chloroaniline	10 U	10	1.0	1	07/14/20 16:10 7/13/20	
4-Chlorophenyl Phenyl Ether	10 U	10	1.5	1	07/14/20 16:10 7/13/20	
4-Nitroaniline	10 U	10	2.7	1	07/14/20 16:10 7/13/20	
4-Nitrophenol	50 U	50	6.4	1	07/14/20 16:10 7/13/20	
Acenaphthene	10 U	10	1.4	1	07/14/20 16:10 7/13/20	
Acenaphthylene	10 U	10	1.4	1	07/14/20 16:10 7/13/20	
Anthracene	10 U	10	1.3	1	07/14/20 16:10 7/13/20	
Benz(a)anthracene	10 U	10	1.6	1	07/14/20 16:10 7/13/20	
Benzo(a)pyrene	10 U	10	1.2	1	07/14/20 16:10 7/13/20	
Benzo(b)fluoranthene	10 U	10	1.2	1	07/14/20 16:10 7/13/20	
Benzo(g,h,i)perylene	10 U	10	1.0	1	07/14/20 16:10 7/13/20	
Benzo(k)fluoranthene	10 U	10	1.3	1	07/14/20 16:10 7/13/20	
Benzoic Acid	100 U	100	36	1	07/14/20 16:10 7/13/20	
Benzyl Alcohol	10 U	100	1.6	1	07/14/20 16:10 7/13/20	
2,2'-Oxybis(1-chloropropane)	10 U	10	1.4	1	07/14/20 16:10 7/13/20	
Bis(2-chloroethoxy)methane	10 U	10	1.4	1	07/14/20 16:10 7/13/20 07/14/20 16:10 7/13/20	
Bis(2-chloroethyl) Ether	10 U	10	1.3	1	07/14/20 16:10 7/13/20 07/14/20 16:10 7/13/20	
Bis(2-ethylhexyl) Phthalate	10 U	10	1.0	1	07/14/20 16:10	
Butyl Benzyl Phthalate	10 U	10	1.0	1	07/14/20 16:10	
	10 U	10	1.4	1		
Chrysene	10 U	10	1.2	1	07/14/20 16:10 7/13/20	

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Superset Reference: 20-0000555771 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007408-01
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed D	Date Extracted	Q
Di-n-butyl Phthalate	2.4 J	10	2.0	1	07/14/20 16:10	7/13/20	
Di-n-octyl Phthalate	10 U	10	3.3	1	07/14/20 16:10	7/13/20	
Dibenz(a,h)anthracene	10 U	10	1.1	1	07/14/20 16:10	7/13/20	
Dibenzofuran	10 U	10	1.4	1	07/14/20 16:10	7/13/20	
Diethyl Phthalate	10 U	10	1.1	1	07/14/20 16:10	7/13/20	
Dimethyl Phthalate	10 U	10	1.3	1	07/14/20 16:10	7/13/20	
Fluoranthene	10 U	10	1.5	1	07/14/20 16:10	7/13/20	
Fluorene	10 U	10	1.3	1	07/14/20 16:10	7/13/20	
Hexachlorobenzene	10 U	10	1.6	1	07/14/20 16:10	7/13/20	
Hexachlorobutadiene	10 U	10	1.0	1	07/14/20 16:10	7/13/20	
Hexachlorocyclopentadiene	10 U	10	2.2	1	07/14/20 16:10	7/13/20	
Hexachloroethane	10 U	10	1.1	1	07/14/20 16:10	7/13/20	
Indeno(1,2,3-cd)pyrene	10 U	10	1.8	1	07/14/20 16:10	7/13/20	
Isophorone	10 U	10	1.4	1	07/14/20 16:10	7/13/20	
N-Nitrosodi-n-propylamine	10 U	10	1.2	1	07/14/20 16:10	7/13/20	
N-Nitrosodiphenylamine	10 U	10	2.7	1	07/14/20 16:10	7/13/20	
Naphthalene	10 U	10	1.2	1	07/14/20 16:10	7/13/20	
Nitrobenzene	10 U	10	1.5	1	07/14/20 16:10	7/13/20	
Pentachlorophenol (PCP)	50 U	50	9.8	1	07/14/20 16:10	7/13/20	
Phenanthrene	10 U	10	1.4	1	07/14/20 16:10	7/13/20	
Phenol	10 U	10	1.0	1	07/14/20 16:10	7/13/20	
Pyrene	10 U	10	1.5	1	07/14/20 16:10	7/13/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	67	35 - 141	07/14/20 16:10	
2-Fluorobiphenyl	54	31 - 118	07/14/20 16:10	
2-Fluorophenol	45	10 - 105	07/14/20 16:10	
Nitrobenzene-d5	55	31 - 110	07/14/20 16:10	
Phenol-d6	32	10 - 107	07/14/20 16:10	
p-Terphenyl-d14	60	10 - 165	07/14/20 16:10	

Tentatively Identified Compounds

			Result	
CAS#	Compound Identification	RT	ug/L	Q

No Tentatively Identified Compounds

Detected

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/14/20

Sample Matrix: Water

Duplicate Lab Control Sample Summary Semivolatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005820

Lab Control Sample

Duplicate Lab Control Sample

RQ2007408-02

RQ2007408-03

Analyte Name	Analytica l Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
1,2,4-Trichlorobenzene	8270D	41.0	80.0	51	45.3	80.0	57	10-127	11	30
1,2-Dichlorobenzene	8270D	40.1	80.0	50	43.3	80.0	54	23-130	8	30
1,3-Dichlorobenzene	8270D	38.8	80.0	48	41.6	80.0	52	21-90	8	30
1,4-Dichlorobenzene	8270D	38.9	80.0	49	41.5	80.0	52	10-124	6	30
2,4,5-Trichlorophenol	8270D	54.6	80.0	68	62.1	80.0	78	48-134	14	30
2,4,6-Trichlorophenol	8270D	53.8	80.0	67	57.6	80.0	72	44-135	7	30
2,4-Dichlorophenol	8270D	44.9	80.0	56	49.6	80.0	62	48-127	10	30
2,4-Dimethylphenol	8270D	51.1	80.0	64	55.3	80.0	69	59-113	8	30
2,4-Dinitrophenol	8270D	49.7 J	80.0	62	47.6 J	80.0	60	21-154	3	30
2,4-Dinitrotoluene	8270D	60.9	80.0	76	66.3	80.0	83	54-130	9	30
2,6-Dinitrotoluene	8270D	64.2	80.0	80	71.6	80.0	90	51-127	12	30
2-Chloronaphthalene	8270D	51.3	80.0	64	55.8	80.0	70	40-108	9	30
2-Chlorophenol	8270D	40.5	80.0	51	43.4	80.0	54	42-112	6	30
2-Methylnaphthalene	8270D	47.8	80.0	60	48.8	80.0	61	34-102	2	30
2-Methylphenol	8270D	44.5	80.0	56	48.2	80.0	60	47-100	7	30
2-Nitroaniline	8270D	60.0	80.0	75	64.4	80.0	80	52-133	6	30
2-Nitrophenol	8270D	47.5	80.0	59	49.6	80.0	62	43-131	5	30
3,3'-Dichlorobenzidine	8270D	59.6	80.0	75	64.0	80.0	80	43-126	6	30
3- and 4-Methylphenol Coelution	8270D	41.8	80.0	52	43.3	80.0	54	40-92	4	30
3-Nitroaniline	8270D	57.1	80.0	71	61.3	80.0	77	42-111	8	30
4,6-Dinitro-2-methylphenol	8270D	53.3	80.0	67	58.2	80.0	73	36-152	9	30
4-Bromophenyl Phenyl Ether	8270D	61.6	80.0	77	65.6	80.0	82	48-114	6	30
4-Chloro-3-methylphenol	8270D	54.4	80.0	68	58.7	80.0	73	52-113	7	30
4-Chloroaniline	8270D	54.2	80.0	68	56.9	80.0	71	44-109	4	30
4-Chlorophenyl Phenyl Ether	8270D	53.2	80.0	67	56.3	80.0	70	51-107	4	30
4-Nitroaniline	8270D	58.4	80.0	73	64.4	80.0	80	54-133	9	30
4-Nitrophenol	8270D	19.5 J	80.0	24	19.5 J	80.0	24	10-126	<1	30
Acenaphthene	8270D	55.9	80.0	70	59.6	80.0	74	52-107	6	30
Acenaphthylene	8270D	59.5	80.0	74	64.7	80.0	81	55-109	9	30
Anthracene	8270D	63.3	80.0	79	68.2	80.0	85	55-116	7	30
Benz(a)anthracene	8270D	59.3	80.0	74	63.9	80.0	80	61-121	8	30
Benzo(a)pyrene	8270D	61.3	80.0	77	69.0	80.0	86	44-114	11	30
Benzo(b)fluoranthene	8270D	58.2	80.0	73	65.0	80.0	81	62-115	10	30
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QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/14/20

Sample Matrix: Water

Duplicate Lab Control Sample Summary Semivolatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005820

Lab Control Sample

Duplicate Lab Control Sample

RQ2007408-02

RQ2007408-03

Analyte Name	Analytica l Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
Benzo(g,h,i)perylene	8270D	69.3	80.0	87	76.1	80.0	95	63-136	9	30
Benzo(k)fluoranthene	8270D	65.8	80.0	82	72.3	80.0	90	49-133	9	30
Benzoic Acid	8270D	67.3 J	120	56	69.8 J	120	58	10-94	4	30
Benzyl Alcohol	8270D	52.5	80.0	66	52.8	80.0	66	31-109	<1	30
2,2'-Oxybis(1-chloropropane)	8270D	45.4	80.0	57	48.4	80.0	61	32-122	7	30
Bis(2-chloroethoxy)methane	8270D	51.7	80.0	65	55.3	80.0	69	55-110	6	30
Bis(2-chloroethyl) Ether	8270D	43.2	80.0	54	47.0	80.0	59	46-102	9	30
Bis(2-ethylhexyl) Phthalate	8270D	62.9	80.0	79	66.6	80.0	83	51-132	5	30
Butyl Benzyl Phthalate	8270D	60.3	80.0	75	66.3	80.0	83	41-148	10	30
Chrysene	8270D	64.6	80.0	81	70.0	80.0	88	57-118	8	30
Di-n-butyl Phthalate	8270D	68.0	80.0	85	75.5	80.0	94	57-128	10	30
Di-n-octyl Phthalate	8270D	61.7	80.0	77	67.7	80.0	85	62-124	10	30
Dibenz(a,h)anthracene	8270D	74.6	80.0	93	83.6	80.0	105	54-135	12	30
Dibenzofuran	8270D	61.0	80.0	76	65.1	80.0	81	55-110	6	30
Diethyl Phthalate	8270D	60.3	80.0	75	63.5	80.0	79	53-113	5	30
Dimethyl Phthalate	8270D	66.0	80.0	83	70.2	80.0	88	51-112	6	30
Fluoranthene	8270D	70.6	80.0	88	74.4	80.0	93	66-127	6	30
Fluorene	8270D	60.8	80.0	76	65.1	80.0	81	54-106	6	30
Hexachlorobenzene	8270D	69.2	80.0	86	73.2	80.0	91	53-123	6	30
Hexachlorobutadiene	8270D	43.3	80.0	54	47.6	80.0	60	16-95	11	30
Hexachlorocyclopentadiene	8270D	14.5	80.0	18	18.2	80.0	23	10-99	24	30
Hexachloroethane	8270D	39.0	80.0	49	42.8	80.0	54	15-92	10	30
Indeno(1,2,3-cd)pyrene	8270D	62.9	80.0	79	68.6	80.0	86	62-137	8	30
Isophorone	8270D	46.0	80.0	57	47.6	80.0	60	50-116	5	30
N-Nitrosodi-n-propylamine	8270D	55.5	80.0	69	56.9	80.0	71	49-115	3	30
N-Nitrosodiphenylamine	8270D	74.5	80.0	93	82.0	80.0	102	45-123	9	30
Naphthalene	8270D	45.5	80.0	57	49.9	80.0	62	38-99	8	30
Nitrobenzene	8270D	49.9	80.0	62	49.9	80.0	62	46-108	<1	30
Pentachlorophenol (PCP)	8270D	58.4	80.0	73	61.3	80.0	77	29-164	5	30
Phenanthrene	8270D	62.0	80.0	78	66.7	80.0	83	58-118	6	30
Phenol	8270D	27.8	80.0	35	28.7	80.0	36	10-113	3	30
Pyrene	8270D	66.1	80.0	83	71.5	80.0	89	61-122	7	30

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Superset Reference:20-0000555771 rev 00



Semivolatile Organic Compounds by GC

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

SURROGATE RECOVERY SUMMARY

Organochlorine Pesticides by Gas Chromatography

Analysis Method: 8081B **Extraction Method:** EPA 3510C

		Decachlorobiphenyl	Tetrachloro-m-xylene	
Sample Name	Lab Code	10-164	10-147	
WG-9954-070620-SG-029	R2005820-002	51	63	
WG-9954-070720-SG-031	R2005820-003	61	56	
WG-9954-070720-SG-032	R2005820-004	66	63	
WG-9954-070720-SG-030	R2005820-005	77	73	
Method Blank	RQ2007347-01	63	65	
Lab Control Sample	RQ2007347-02	56	58	
Duplicate Lab Control Sample	RQ2007347-03	62	55	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007347-01Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
4,4'-DDE	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
4,4'-DDT	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
Aldrin	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
Dieldrin	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
Endosulfan I	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
Endosulfan II	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
Endosulfan Sulfate	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
Endrin	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
Endrin Ketone	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
Heptachlor	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
Heptachlor Epoxide	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
Methoxychlor	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
Toxaphene	0.50 U	0.50	0.50	1	07/13/20 14:42	7/10/20	
alpha-BHC	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
alpha-Chlordane	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
beta-BHC	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
delta-BHC	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
gamma-BHC (Lindane)	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	
gamma-Chlordane	0.050 U	0.050	0.020	1	07/13/20 14:42	7/10/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	63	10 - 164	07/13/20 14:42	
Tetrachloro-m-xylene	65	10 - 147	07/13/20 14:42	

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/13/20

Sample Matrix: Water

Duplicate Lab Control Sample Summary Organochlorine Pesticides by Gas Chromatography

Units:ug/L Basis:NA

Service Request: R2005820

Lab Control Sample

Duplicate Lab Control Sample

RQ2007347-02

RQ2007347-03

	Analytical		Spike			Spike		% Rec		RPD
Analyte Name	Method	Result	Amount	% Rec	Result	Amount	% Rec	Limits	RPD	Limit
4,4'-DDD	8081B	0.303	0.400	76	0.342	0.400	86	42-159	12	30
4,4'-DDE	8081B	0.280	0.400	70	0.322	0.400	81	47-147	14	30
4,4'-DDT	8081B	0.309	0.400	77	0.353	0.400	88	41-149	13	30
Aldrin	8081B	0.236	0.400	59	0.234	0.400	59	22-137	<1	30
Dieldrin	8081B	0.307	0.400	77	0.346	0.400	86	52-144	12	30
Endosulfan I	8081B	0.296	0.400	74	0.329	0.400	82	52-136	11	30
Endosulfan II	8081B	0.236	0.400	59	0.283	0.400	71	57-138	18	30
Endosulfan Sulfate	8081B	0.267	0.400	67	0.310	0.400	78	34-156	15	30
Endrin	8081B	0.319	0.400	80	0.357	0.400	89	56-143	11	30
Endrin Ketone	8081B	0.301	0.400	75	0.346	0.400	87	59-143	14	30
Heptachlor	8081B	0.274	0.400	68	0.283	0.400	71	32-141	3	30
Heptachlor Epoxide	8081B	0.307	0.400	77	0.338	0.400	84	51-143	10	30
Methoxychlor	8081B	0.318	0.400	80	0.359	0.400	90	56-149	12	30
alpha-BHC	8081B	0.297	0.400	74	0.313	0.400	78	36-151	5	30
alpha-Chlordane	8081B	0.293	0.400	73	0.322	0.400	80	50-139	9	30
beta-BHC	8081B	0.315	0.400	79	0.341	0.400	85	55-149	8	30
delta-BHC	8081B	0.292	0.400	73	0.328	0.400	82	29-159	12	30
gamma-BHC (Lindane)	8081B	0.305	0.400	76	0.326	0.400	82	41-149	7	30
gamma-Chlordane	8081B	0.296	0.400	74	0.323	0.400	81	50-140	9	30

QA/QC Report

Service Request: R2005820

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

Project:

SURROGATE RECOVERY SUMMARY Polychlorinated Biphenyls (PCBs) by GC

Analysis Method: 8082A **Extraction Method:** EPA 3510C

		Decachlorobiphenyl	Tetrachloro-m-xylene
Sample Name	Lab Code	10-152	14-129
WG-9954-070620-SG-029	R2005820-002	50	54
WG-9954-070720-SG-031	R2005820-003	59	47
WG-9954-070720-SG-032	R2005820-004	63	52
WG-9954-070720-SG-030	R2005820-005	68	55
Method Blank	RQ2007347-01	69	59
Lab Control Sample	RQ2007347-02	70	48
Duplicate Lab Control Sample	RQ2007347-03	68	53

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005820

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007347-01Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	1.0 U	1.0	0.50	1	07/13/20 14:13	7/10/20	
Aroclor 1221	2.0 U	2.0	1.0	1	07/13/20 14:13	7/10/20	
Aroclor 1232	1.0 U	1.0	0.50	1	07/13/20 14:13	7/10/20	
Aroclor 1242	1.0 U	1.0	0.50	1	07/13/20 14:13	7/10/20	
Aroclor 1248	1.0 U	1.0	0.50	1	07/13/20 14:13	7/10/20	
Aroclor 1254	1.0 U	1.0	0.50	1	07/13/20 14:13	7/10/20	
Aroclor 1260	1.0 U	1.0	0.50	1	07/13/20 14:13	7/10/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	69	10 - 152	07/13/20 14:13	
Tetrachloro-m-xylene	59	14 - 129	07/13/20 14:13	

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix:

Project:

Water

Duplicate Lab Control Sample Summary Polychlorinated Biphenyls (PCBs) by GC

Units:ug/L Basis:NA

Service Request: R2005820

Date Analyzed: 07/13/20

Lab Control Sample

Duplicate Lab Control Sample

RQ2007347-02

RQ2007347-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
Aroclor 1016	8082A	3.35	4.00	84	3.40	4.00	85	49-123	2	30
Aroclor 1260	8082A	3.98	4.00	100	3.79	4.00	95	30-120	5	30



Service Request No:R2005980

Ms. Kathy Willy GHD Services Inc. 2055 Niagara Falls Blvd., Niagara Falls, NY 14304

Laboratory Results for: Love Canal:292-402-D02-3100

Dear Ms. Willy,

Enclosed are the results of the sample(s) submitted to our laboratory July 11, 2020 For your reference, these analyses have been assigned our service request number R2005980.

All testing was performed according to our laboratory's quality assurance program and met the requirements of the TNI standards except as noted in the case narrative report. Any testing not included in the lab's accreditation is identified on a Non-Certified Analytes report. All results are intended to be considered in their entirety. ALS Environmental is not responsible for use of less than the complete report. Results apply only to the individual samples submitted to the lab for analysis, as listed in the report. The measurement uncertainty of the results included in this report is within that expected when using the prescribed method(s), and represented by Laboratory Control Sample control limits. Any events, such as QC failures or Holding Time exceedances, which may add to the uncertainty are explained in the report narrative or are flagged with qualifiers. The flags are explained in the Report Qualifiers and Definitions page of this report.

Please contact me if you have any questions. My extension is 7471. You may also contact me via email at Brady.Kalkman@alsglobal.com.

Respectfully submitted,

Goody Kullen

ALS Group USA, Corp. dba ALS Environmental

Brady Kalkman

Project Manager



Narrative Documents

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com



Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100

Service Request: R2005980

Date Received: 07/11/2020

Sample Matrix: Water

CASE NARRATIVE

All analyses were performed consistent with the quality assurance program of ALS Environmental. This report contains analytical results for samples for the Tier level IV requested by the client.

Sample Receipt:

Five water samples were received for analysis at ALS Environmental on 07/11/2020. Any discrepancies upon initial sample inspection are annotated on the sample receipt and preservation form included within this report. The samples were stored at minimum in accordance with the analytical method requirements.

Semivolatiles by GC/MS:

Method 8270D, 07/20/2020: The upper control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). The field samples analyzed in this sequence did not contain the analyte(s) in question above the Method Reporting Limit (MRL). Since the exceedance equates to a potential high bias, the data quality was not significantly affected and no further corrective action was taken.

Method 8270D, 07/20/2020: The lower control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). Since there were no detections of the analyte(s) above the MRL in the associated field samples, the quantitation is not affected. The data quality was not significantly affected and no further corrective action was taken.

Semivoa GC:

Method 8081B, 07/15/2020: The lower control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). Since there were no detections of the analyte(s) above the MRL in the associated field samples, the quantitation is not affected. The data quality was not significantly affected and no further corrective action was taken. Method 8082A, 07/16/2020: The upper control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). The field samples analyzed in this sequence did not contain the analyte(s) in question above the Method Reporting Limit (MRL). Since the exceedance equates to a potential high bias, the data quality was not significantly affected and no further corrective action was taken.

Method 8082A, 07/16/2020: The upper control criterion was exceeded for one or more analytes in the Laboratory Control Sample (LCS). There were no detections of the analyte(s) above the MRL in the associated field samples. 1260 went out high on front column of the lcsd. The error associated with elevated recovery equates to a high bias. The sample data is not significantly affected. No further corrective action was appropriate.

Method 8082A, 07/22/2020: The upper control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). The field samples analyzed in this sequence did not contain the analyte(s) in question above the Method Reporting Limit (MRL). Since the exceedance equates to a potential high bias, the data quality was not significantly affected and no further corrective action was taken.

Volatiles by GC/MS:

Method 8260C, 07/21/2020: The lower control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). Since there were no detections of the analyte(s) above the MRL in the associated field samples, the quantitation is not affected. The data quality was not significantly affected and no further corrective action was taken.

Method 8260C, 687805: Sample(s) required dilution due to the foaming nature of the matrix. The reporting limits are adjusted to reflect the dilution.

Method 8260C, 688008: Sample(s) required dilution due to the foaming nature of the matrix. The reporting limits are adjusted to reflect the dilution.



Method 8260C, 07/22/2020: The upper control criterion was exceeded for one or more analytes in the Laboratory Control Sample (LCS). There were no detections of the analyte(s) above the MRL in the associated field samples. The error associated with elevated recovery equates to a high bias. The sample data is not significantly affected. No further corrective action was appropriate.

Method 8260C, 07/22/2020: The lower control limit was exceeded for one or more analytes in the Continuing Calibration Verification (CCV). Since there were no detections of the analyte(s) above the MRL in the associated field samples, the quantitation is not affected. The data quality was not significantly affected and no further corrective action was taken.

Approved by _____

Date 07/30/2020



SAMPLE DETECTION SUMMARY

CLIENT ID: WG-9954-070920-SG-033								
Analyte	Results	Flag	MDL	MRL	Units	Method		
Carbon Disulfide	9.2	J	0.42	10	ug/L	8260C		
Bis(2-ethylhexyl) Phthalate	3.9	J	0.91	9.1	ug/L	8270D		
Di-n-butyl Phthalate	2.3	J	1.9	9.1	ug/L	8270D		
alpha-BHC	0.12		0.019	0.045	ug/L	8081B		
delta-BHC	0.088		0.019	0.045	ug/L	8081B		
gamma-BHC (Lindane)	0.15		0.019	0.045	ug/L	8081B		
LIENT ID: WG-9954-070920-SG-034		Lat	D: R2005	980-002				
Analyte	Results	Flag	MDL	MRL	Units	Method		
Carbon Disulfide	160		2.1	50	ug/L	8260C		
Naphthalene	1.2	J	1.1	9.1	ug/L	8270D		
alpha-BHC	0.069		0.019	0.045	ug/L	8081B		
delta-BHC	0.065		0.019	0.045	ug/L	8081B		
gamma-BHC (Lindane)	0.077		0.019	0.045	ug/L	8081B		
LIENT ID: WG-9954-070920-SG-035	Lab ID: R2005980-003							
Analyte	Results	Flag	MDL	MRL	Units	Method		
Carbon Disulfide	25	J	2.1	50	ug/L	8260C		
alpha-BHC	0.072		0.019	0.045	ug/L	8081B		
beta-BHC	0.019	J	0.019	0.045	ug/L	8081B		
delta-BHC	0.19		0.019	0.045	ug/L	8081B		
gamma-BHC (Lindane)	0.11		0.019	0.045	ug/L	8081B		
LIENT ID: WG-9954-070920-SG-036		Lab ID: R2005980-005						
Analyte	Results	Flag	MDL	MRL	Units	Method		
Bis(2-ethylhexyl) Phthalate	1.3	J	0.91	9.1	ug/L	8270D		
Methoxychlor	0.041	J	0.019	0.045	ug/L	8081B		
alpha-BHC	0.069		0.019	0.045	ug/L	8081B		
delta-BHC	0.021	J	0.019	0.045	ug/L	8081B		
gamma-BHC (Lindane)	0.058		0.019	0.045	ug/L	8081B		



Sample Receipt Information

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request:R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

SAMPLE CROSS-REFERENCE

SAMPLE #	CLIENT SAMPLE ID	<u>DATE</u>	<u>TIME</u>
R2005980-001	WG-9954-070920-SG-033	7/9/2020	1225
R2005980-002	WG-9954-070920-SG-034	7/9/2020	1255
R2005980-003	WG-9954-070920-SG-035	7/9/2020	1315
R2005980-004	TB-9954-070920-SG-007	7/9/2020	0000
R2005980-005	WG-9954-070920-SG-036	7/9/2020	0815

Page 7 of 78

EVENT COMPLETE

CHAIN-OF-CUSTODY/Analytical Request Document The Chain-of-Custody is a LEGAL DOCUMENT, All relevant fields must be completed accurately.

Report To: Kathy Willy
Сору То:
Involce To:
PO:
Project Name: LOVE CANAL ANNUAL GW
Project Number: 9954

Lai	Information .
Laboratory: ALS	
Laboratory Location: 11 BUILDING 300, SUITE 3 ROCHESTER, NY 14623	60
ROCHESTER, NY 14623	3
Laboratory Contact: BR	RADY KALKMAN
Requested Due Date:	TAT: 10
QA/QC Regulrements:	

Event Information ID#: LC ANNUAL GW SAMPLING 2020-07-1 SSOW Ref#: 292-402-999-3100 Sampler Name: S GARDNER, D

	Valid Matrix Code WG Groundwater								Sample Condition		
	WB Borehole Water WS Surface Water		_		<u>ē</u>				Temp in C		
	SO Soil SE Sediment	a a	Collected	ollected	PestPCBs(None)	<u> </u>			Received on ice	Y/N	
		Ö) je	1 8	8	Ē	ᅙ		Sealed Cooler	Y/N	
		Matrix	Date (Time	£	SVOC(none)	VOA(HCI)		Samples Intact	Y/N	
Sample Identification		₹	ä	Ē	ے	is	>	Remarks			
WG-9954-070920-SG-03	3	WG	07/09/2020	12:25	2	2	3				
WG-9954-070920-SG-03	4	WG	07/09/2020	12:55	2	2	3				
WG-9954-070920-SG-03	5	WG	07/09/2020	13:15	2	2	3	-			
TB-9954-070920-SG-007		WG Q	07/09/2020	00:00	0	0	3				
WG-9954-071020-SG-03	6	WG	07/10/2020	08:15	2	2	3			·	
Total Bottles					8	В	15	Grand Tot	al:31		

SHIPMENT METHOD	NO. OF COOLERS	RELINQUISHED BY:	DATE	TIME	RECIEVED BY:	DATÉ	TIME
FedEx	1	Shoun Haidnin	7/10/20	2013	os alutulus	7/11/2	12/CA05
AIRBILL#:							****

CHAIN OF CUSTODY / LABORATORY ANALYSIS REQUEST FORM SR# 004, 005, 006, 007, 008, 009, 010, 1565 Jefferson Road, Bldg 300, Suite 360, Rochester, NY 14623 T030477 011, 012, 013 Phone (585) 288-5380 / FAX (585) 288-8475 Environmental www,alsglobal.com 14D 2 Love Canal:292-402-D02-3100 Project Number: 9954 Annual Long Term Monitoring Report To Kathy Willy NUMBER OF CONTAINERS Company / Address GHD Services Inc. 2055 Niagara Falls Blvd., Suite 3 Niagara Falls NY, 14304 081B / Pest OC 8082A / PCB 3270D / SVO 3260C / VOC Phone # 716-297-2265 716-297-2160 Sampler Printed Name Sampler Signature Remarks SAMPLING Matrix **CLIENT SAMPLE ID** LABID Date Time Liquid Liquid Liquid Liquid Llauld Liquid Llauid Liquid Liquid Liquid 10. Report Requirements Invoice Information **Turnaround Requirements** Special Instructions/Comments: RUSH (SURCHARGES APPLY) I. Results Only P.O.# II. Results + QC Summaries (LCS, DUP, MS/MSD as required) Standard (3 weeks) III. Results + QC and Cilibration Bill To:____ Summaries X . IV. Data Validation Report REQUESTED FAX DATE with Raw Data EDate Yes No Requested Report Date Received By: Relinquished By: Received By: Received By: Relinquished By: Relinquished By: Signature Signature Signature Signature Signature Signature

Printed Name

Firm

Date/Time

Printed Name

Firm

Date/Time

Printed Name

Firm

Date/Time

Printed Name

Firm

Date/Time

Printed Name

Date/Time

Firm

Printed Name

Date/Time

Firm



Cooler Receipt and Preservation Check Form

R2005980 GHD Services Inc.	5
Love Canal:292-402-D02-3100	

Project/Clie	ent GHO)			_Fok	der Num	ber								<u> </u>
	ed on 7/11/7		bу: <u>М</u>	n		COUR	ŒR:	ALS	UPS	PEDEX	(VEI	OCITY	CLIE	NT	
		outside of cooler	r?		Y A	5a	Perch	lorate s	amples	have req	uired he	adspace	?	Y N	(AIA)
2 Custody	papers prope	rly completed (in	k, sign		S N		Did Y	OA vial	s, Alk,	or Sulfide	have s	ig* bub	bles?	18 N	NA _t
3 Did all b	ottles arrive in	good condition ((unbrok	cen)?	Y; N	6	Where	did the	bottles	originate	?	A&S/	ROC	CLIE	NT
4 Circle:	Wet Ice Dry	Ice Gel packs	pres	ent?	N &	7.	Soil V	OA rec	eived a	s: Bu	lk F	псоге	5035s	et 🔿	IA)
3. Temperatur	re Readings	Date: 7/11/70	70	Time:	097	30	ID:	(R#)	IR#10)	From	: Temp	Blank	Sarraj	ole Bottle
Observed Te	emp (°C)	1,30		·											
Within 0-6°	C?	Ø N		Y	N	YN	J	Y	N	Y	N	Y	N	Y	N
If <0°C, wer	e samples froz	en? Y N		Y	N	Y	1	Y	N	Y	N	Y	N	Y	N
If out of 7		note packing/ic	e cond			To	e melt	ed P	oorly P	acked (d	escribe	l below)		Same D	ay Rule
	-	un Samples:							•	-					
- CHOICH P	xpprovar to x	di Sampies		_		 									
	held in storag		K~		y M			at C				ē.			
5035 sample	es placed in st	orage location:		t	у	on		_ at _		within 4	8 hours	of samp	oling?	Y	N
									-			-			
9. V 10. I 11. V 12. V	Were all bottle Did all bottle la Were correct co Were 5035 vial	ervation Check** labels complete (abels and tags agr ontainers used for ls acceptable (no	(i.e. and ree with r the tea extra la	alysis, custo sts indi abels, r	preserving preserving	ration, etc.) ers? ring)?)?		-	PES (ES)	NO NO NO NO	a . 1		₩ <u>A</u>	
		Cassettes / Tubes	,				Pressu			Tedlar®				<u>(YA)</u> _	Final
pH	Lot of test paper	Reagent	Yes Yes	No No	Lot R	Received		Ехр	Samı Adju	ole ID sted	Vol. Adde		ot Adde	u .	pH
≥12		NaOH													
≤2		HNO ₃									<u> </u>				
≤2		H ₂ SO ₄						<u> </u>	<u> </u>		ļ		_		
<4		NaHSO ₄							ļ	·	<u> </u>				
5-9		For 608pest	<u> </u>			lotify for 3		 			 	- -			
Residual		For CN,				ontact PM t O ₃ (625, 60					1	1			1
Chlorine		Phenol, 625, 608pest, 522	<u> </u>			ascorbic (ph									
(-)	·	Na ₂ S ₂ O ₃						· ·	 	·	 				
L	_L	ZnAcetate	+	+			··	+	**VO	As and 16	4 Not to	be tested	before an	alysis.	<u> </u>
		HCl	**	**	-	 		+	1	•		-		mical pr	eservatives
		L	1						are ch	ecked (not	just repre	senianve	s).		
Explain a	11 Discrepanc	lus Corus ies/ Other Comm 4-070170-86-	nents:	05	1120	- /BM	nc		•	·		 .		<u>-</u> .	
31/20	3 vials; Triv Blank														

HPROD	BULK
HTR	FLDT
SUB	HGFB
ALS	LL3541

Labels second	lary reviewed	by:
PC Secondary	Review:	

*significant air bubbles: VOA > 5-6 mm: WC > 1 in. diameter Page 10 of 78



Miscellaneous Forms

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com



REPORT QUALIFIERS AND DEFINITIONS

- U Analyte was analyzed for but not detected. The sample quantitation limit has been corrected for dilution and for percent moisture, unless otherwise noted in the case narrative.
- J Estimated value due to either being a Tentatively Identified Compound (TIC) or that the concentration is between the MRL and the MDL. Concentrations are not verified within the linear range of the calibration. For DoD: concentration >40% difference between two GC columns (pesticides/Arclors).
- B Analyte was also detected in the associated method blank at a concentration that may have contributed to the sample result.
- E Inorganics- Concentration is estimated due to the serial dilution was outside control limits.
- E Organics- Concentration has exceeded the calibration range for that specific analysis.
- D Concentration is a result of a dilution, typically a secondary analysis of the sample due to exceeding the calibration range or that a surrogate has been diluted out of the sample and cannot be assessed.
- * Indicates that a quality control parameter has exceeded laboratory limits. Under the õNotesö column of the Form I, this qualifier denotes analysis was performed out of Holding Time.
- H Analysis was performed out of hold time for tests that have an õimmediateö hold time criteria.
- # Spike was diluted out.

- + Correlation coefficient for MSA is <0.995.
- N Inorganics- Matrix spike recovery was outside laboratory limits.
- N Organics- Presumptive evidence of a compound (reported as a TIC) based on the MS library search.
- S Concentration has been determined using Method of Standard Additions (MSA).
- W Post-Digestion Spike recovery is outside control limits and the sample absorbance is <50% of the spike absorbance.
- P Concentration >40% difference between the two GC columns.
- C Confirmed by GC/MS
- Q DoD reports: indicates a pesticide/Aroclor is not confirmed (×100% Difference between two GC columns).
- X See Case Narrative for discussion.
- MRL Method Reporting Limit. Also known as:
- LOQ Limit of Quantitation (LOQ)

 The lowest concentration at which the method analyte may be reliably quantified under the method conditions.
- MDL Method Detection Limit. A statistical value derived from a study designed to provide the lowest concentration that will be detected 99% of the time. Values between the MDL and MRL are estimated (see J qualifier).
- LOD Limit of Detection. A value at or above the MDL which has been verified to be detectable.
- ND Non-Detect. Analyte was not detected at the concentration listed. Same as U qualifier.



Rochester Lab ID # for State Certifications¹

Connecticut ID # PH0556	Maine ID #NY0032	Pennsylvania ID# 68-786
Delaware Approved	New Hampshire ID # 2941	Rhode Island ID # 158
DoD ELAP #65817	New York ID # 10145	Virginia #460167
Florida ID # E87674	North Carolina #676	

¹ Analyses were performed according to our laboratory

NELAP-approved quality assurance program and any applicable state or agency requirements. The test results meet requirements of the current NELAP/TNI standards or state or agency requirements, where applicable, except as noted in the case narrative. Since not all analyte/method/matrix combinations are offered for state/NELAC accreditation, this report may contain results which are not accredited. For a specific list of accredited analytes, contact the laboratory or go to https://www.alsglobal.com/locations/americas/north-america/usa/new-york/rochester-environmental

ALS Laboratory Group

Acronyms

ASTM American Society for Testing and Materials

A2LA American Association for Laboratory Accreditation

CARB California Air Resources Board

CAS Number Chemical Abstract Service registry Number

CFC Chlorofluorocarbon CFU Colony-Forming Unit

DEC Department of Environmental Conservation

DEQ Department of Environmental Quality

DHS Department of Health Services

DOE Department of Ecology DOH Department of Health

EPA U. S. Environmental Protection Agency

ELAP Environmental Laboratory Accreditation Program

GC Gas Chromatography

GC/MS Gas Chromatography/Mass Spectrometry

LUFT Leaking Underground Fuel Tank

M Modified

MCL Maximum Contaminant Level is the highest permissible concentration of a

substance allowed in drinking water as established by the USEPA.

MDL Method Detection Limit
MPN Most Probable Number
MRL Method Reporting Limit

NA Not Applicable NC Not Calculated

NCASI National Council of the Paper Industry for Air and Stream Improvement

ND Not Detected

NIOSH National Institute for Occupational Safety and Health

PQL Practical Quantitation Limit

RCRA Resource Conservation and Recovery Act

SIM Selected Ion Monitoring

TPH Total Petroleum Hydrocarbons

tr Trace level is the concentration of an analyte that is less than the PQL but

greater than or equal to the MDL.

Analyst Summary report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Monitoring

Sample Name: WG-9954-070920-SG-033

Lab Code: R2005980-001

Sample Matrix: Water

Date Received: 07/11/20

Service Request: R2005980

Date Collected: 07/9/20

Analysis Method Extracted/Digested By Analyzed By

8081B KSERCU BALLGEIER 8082A KSERCU JMISIUREWICZ

8260C KRUEST

8270D KSERCU JMISIUREWICZ

Sample Name: WG-9954-070920-SG-034 Date Collected: 07/9/20

Lab Code: R2005980-002 **Date Received:** 07/11/20

Sample Matrix: Water

Analysis Method Extracted/Digested By Analyzed By

8081B KSERCU BALLGEIER 8082A KSERCU JMISIUREWICZ

8260C KRUEST

8270D KSERCU JMISIUREWICZ

Sample Name: WG-9954-070920-SG-035 **Date Collected:** 07/9/20

Lab Code: R2005980-003 **Date Received:** 07/11/20

Sample Matrix: Water

Analysis Method Extracted/Digested By Analyzed By

8081B KSERCU BALLGEIER 8082A KSERCU JMISIUREWICZ

8260C KRUEST

8270D KSERCU JMISIUREWICZ

Sample Name: TB-9954-070920-SG-007 **Date Collected:** 07/9/20

Lab Code: R2005980-004 **Date Received:** 07/11/20

Sample Matrix: Water

Analysis Method Extracted/Digested By Analyzed By

8260C KRUEST

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Analyst Summary report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Monitoring

Sample Name: WG-9954-070920-SG-036

Lab Code: R2005980-005 **Date Received:** 07/11/20

Sample Matrix: Water

Analysis MethodExtracted/Digested ByAnalyzed By8081BKSERCUBALLGEIER8082AKSERCUBALLGEIER8260CKRUEST

8270D KSERCU JMISIUREWICZ

Service Request: R2005980

Date Collected: 07/9/20



INORGANIC PREPARATION METHODS

The preparation methods associated with this report are found in these tables unless discussed in the case narrative.

Water/Liquid Matrix

Analytical Method	Preparation Method
200.7	200.2
200.8	200.2
6010C	3005A/3010A
6020A	ILM05.3
9034 Sulfide Acid Soluble	9030B
SM 4500-CN-E Residual Cyanide	SM 4500-CN-G
SM 4500-CN-E WAD Cyanide	SM 4500-CN-I

Solid/Soil/Non-Aqueous Matrix

Analytical Method	Preparation				
	Method				
6010C	3050B				
6020A	3050B				
6010C TCLP (1311)	3005A/3010A				
extract					
6010 SPLP (1312) extract	3005A/3010A				
7199	3060A				
300.0 Anions/ 350.1/	DI extraction				
353.2/ SM 2320B/ SM					
5210B/ 9056A Anions					
For analytical methods not listed,					
method is the same as the analytical method					
I ICICICILC.					



Sample Results

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com



Volatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 12:25

Monitoring

Sample Matrix: Water Date Received: 07/11/20 09:05

 Sample Name:
 WG-9954-070920-SG-033
 Units: ug/L

 Lab Code:
 R2005980-001
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/22/20 13:03	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/22/20 13:03	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/22/20 13:03	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/22/20 13:03	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/22/20 13:03	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/22/20 13:03	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/22/20 13:03	
2-Butanone (MEK)	10 U	10	0.78	1	07/22/20 13:03	
2-Hexanone	10 U	10	0.20	1	07/22/20 13:03	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/22/20 13:03	
Acetone	10 U	10	5.0	1	07/22/20 13:03	
Benzene	5.0 U	5.0	0.20	1	07/22/20 13:03	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/22/20 13:03	
Bromoform	5.0 U	5.0	0.25	1	07/22/20 13:03	
Bromomethane	5.0 U	5.0	0.70	1	07/22/20 13:03	
Carbon Disulfide	9.2 J	10	0.42	1	07/22/20 13:03	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/22/20 13:03	
Chlorobenzene	5.0 U	5.0	0.20	1	07/22/20 13:03	
Chloroethane	5.0 U	5.0	0.23	1	07/22/20 13:03	
Chloroform	5.0 U	5.0	0.24	1	07/22/20 13:03	
Chloromethane	5.0 U	5.0	0.28	1	07/22/20 13:03	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/22/20 13:03	
Dichloromethane	5.0 U	5.0	0.65	1	07/22/20 13:03	
Ethylbenzene	5.0 U	5.0	0.20	1	07/22/20 13:03	
Styrene	5.0 U	5.0	0.20	1	07/22/20 13:03	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/22/20 13:03	
Toluene	5.0 U	5.0	0.20	1	07/22/20 13:03	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/22/20 13:03	
Vinyl Acetate	10 U	10	1.1	1	07/22/20 13:03	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/22/20 13:03	
Xylenes, Total	5.0 U	5.0	0.23	1	07/22/20 13:03	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/22/20 13:03	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/22/20 13:03	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/22/20 13:03	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/22/20 13:03	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	96	85 - 122	07/22/20 13:03	
Dibromofluoromethane	101	89 - 119	07/22/20 13:03	
Toluene-d8	104	87 - 121	07/22/20 13:03	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 12:25

Monitoring

Sample Matrix: Water Date Received: 07/11/20 09:05

 Sample Name:
 WG-9954-070920-SG-033
 Units: ug/L

 Lab Code:
 R2005980-001
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
000074-98-6	Propane	1.19	7.7	JN	
000075-28-5	Isobutane	1.31	11.4	JN	
	unknown	1.41	10.4	J	
000074-93-1	Methanethiol	1.59	5.8	JN	
000078-78-4	Butane, 2-methyl-	1.77	5.4	JN	
003658-80-8	Trisulfide, dimethyl	11.47	9.9	JN	
000124-19-6	Nonanal	12.69	7.0	JN	
000075-18-3	Dimethyl sulfide	2.45	325.6	JN	
000624-89-5	Ethane, (methylthio)-	4.32	25.8	JN	
000624-92-0	Disulfide, dimethyl	8.09	33.4	JN	
000066-25-1	Hexanal	9.24	10.3	JN	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 12:55

Monitoring

Sample Matrix: Water Date Received: 07/11/20 09:05

 Sample Name:
 WG-9954-070920-SG-034
 Units: ug/L

 Lab Code:
 R2005980-002
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	25 U	25	1.0	5	07/22/20 13:24	
1,1,2,2-Tetrachloroethane	25 U	25	1.0	5	07/22/20 13:24	
1,1,2-Trichloroethane	25 U	25	1.0	5	07/22/20 13:24	
1,1-Dichloroethane (1,1-DCA)	25 U	25	1.0	5	07/22/20 13:24	
1,1-Dichloroethene (1,1-DCE)	25 U	25	1.0	5	07/22/20 13:24	
1,2-Dichloroethane	25 U	25	1.0	5	07/22/20 13:24	
1,2-Dichloropropane	25 U	25	1.0	5	07/22/20 13:24	
2-Butanone (MEK)	50 U	50	3.9	5	07/22/20 13:24	
2-Hexanone	50 U	50	1.0	5	07/22/20 13:24	
4-Methyl-2-pentanone	50 U	50	1.0	5	07/22/20 13:24	
Acetone	50 U	50	25	5	07/22/20 13:24	
Benzene	25 U	25	1.0	5	07/22/20 13:24	
Bromodichloromethane	25 U	25	1.0	5	07/22/20 13:24	
Bromoform	25 U	25	1.3	5	07/22/20 13:24	
Bromomethane	25 U	25	3.5	5	07/22/20 13:24	
Carbon Disulfide	160	50	2.1	5	07/22/20 13:24	
Carbon Tetrachloride	25 U	25	1.7	5	07/22/20 13:24	
Chlorobenzene	25 U	25	1.0	5	07/22/20 13:24	
Chloroethane	25 U	25	1.2	5	07/22/20 13:24	
Chloroform	25 U	25	1.2	5	07/22/20 13:24	
Chloromethane	25 U	25	1.4	5	07/22/20 13:24	
Dibromochloromethane	25 U	25	1.0	5	07/22/20 13:24	
Dichloromethane	25 U	25	3.3	5	07/22/20 13:24	
Ethylbenzene	25 U	25	1.0	5	07/22/20 13:24	
Styrene	25 U	25	1.0	5	07/22/20 13:24	
Tetrachloroethene (PCE)	25 U	25	1.1	5	07/22/20 13:24	
Toluene	25 U	25	1.0	5	07/22/20 13:24	
Trichloroethene (TCE)	25 U	25	1.0	5	07/22/20 13:24	
Vinyl Acetate	50 U	50	5.5	5	07/22/20 13:24	
Vinyl Chloride	25 U	25	1.0	5	07/22/20 13:24	
Xylenes, Total	25 U	25	1.2	5	07/22/20 13:24	
cis-1,2-Dichloroethene	25 U	25	1.2	5	07/22/20 13:24	
cis-1,3-Dichloropropene	25 U	25	1.0	5	07/22/20 13:24	
trans-1,2-Dichloroethene	25 U	25	1.0	5	07/22/20 13:24	
trans-1,3-Dichloropropene	25 U	25	1.2	5	07/22/20 13:24	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 12:55

Monitoring

Sample Matrix: Water Date Received: 07/11/20 09:05

 Sample Name:
 WG-9954-070920-SG-034
 Units: ug/L

 Lab Code:
 R2005980-002
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	96	85 - 122	07/22/20 13:24	
Dibromofluoromethane	102	89 - 119	07/22/20 13:24	
Toluene-d8	105	87 - 121	07/22/20 13:24	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 12:55

Monitoring

Sample Matrix: Water Date Received: 07/11/20 09:05

 Sample Name:
 WG-9954-070920-SG-034
 Units: ug/L

 Lab Code:
 R2005980-002
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
007446-09-5	Sulfur dioxide	1.29	213.2	JN	
000074-93-1	Methanethiol	1.59	78.0	JN	
000075-18-3	Dimethyl sulfide	2.45	594.4	JN	
000624-89-5	Ethane, (methylthio)-	4.32	50.8	JN	
000624-92-0	Disulfide, dimethyl	8.09	45.7	JN	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 13:15

Monitoring

Sample Matrix: Water Date Received: 07/11/20 09:05

 Sample Name:
 WG-9954-070920-SG-035
 Units: ug/L

 Lab Code:
 R2005980-003
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	25 U	25	1.0	5	07/21/20 16:17	
1,1,2,2-Tetrachloroethane	25 U	25	1.0	5	07/21/20 16:17	
1,1,2-Trichloroethane	25 U	25	1.0	5	07/21/20 16:17	
1,1-Dichloroethane (1,1-DCA)	25 U	25	1.0	5	07/21/20 16:17	
1,1-Dichloroethene (1,1-DCE)	25 U	25	1.0	5	07/21/20 16:17	
1,2-Dichloroethane	25 U	25	1.0	5	07/21/20 16:17	
1,2-Dichloropropane	25 U	25	1.0	5	07/21/20 16:17	
2-Butanone (MEK)	50 U	50	3.9	5	07/21/20 16:17	
2-Hexanone	50 U	50	1.0	5	07/21/20 16:17	
4-Methyl-2-pentanone	50 U	50	1.0	5	07/21/20 16:17	
Acetone	50 U	50	25	5	07/21/20 16:17	
Benzene	25 U	25	1.0	5	07/21/20 16:17	
Bromodichloromethane	25 U	25	1.0	5	07/21/20 16:17	
Bromoform	25 U	25	1.3	5	07/21/20 16:17	
Bromomethane	25 U	25	3.5	5	07/21/20 16:17	
Carbon Disulfide	25 J	50	2.1	5	07/21/20 16:17	
Carbon Tetrachloride	25 U	25	1.7	5	07/21/20 16:17	
Chlorobenzene	25 U	25	1.0	5	07/21/20 16:17	
Chloroethane	25 U	25	1.2	5	07/21/20 16:17	
Chloroform	25 U	25	1.2	5	07/21/20 16:17	
Chloromethane	25 U	25	1.4	5	07/21/20 16:17	
Dibromochloromethane	25 U	25	1.0	5	07/21/20 16:17	
Dichloromethane	25 U	25	3.3	5	07/21/20 16:17	
Ethylbenzene	25 U	25	1.0	5	07/21/20 16:17	
Styrene	25 U	25	1.0	5	07/21/20 16:17	
Tetrachloroethene (PCE)	25 U	25	1.1	5	07/21/20 16:17	
Toluene	25 U	25	1.0	5	07/21/20 16:17	
Trichloroethene (TCE)	25 U	25	1.0	5	07/21/20 16:17	
Vinyl Acetate	50 U	50	5.5	5	07/21/20 16:17	
Vinyl Chloride	25 U	25	1.0	5	07/21/20 16:17	
Xylenes, Total	25 U	25	1.2	5	07/21/20 16:17	
cis-1,2-Dichloroethene	25 U	25	1.2	5	07/21/20 16:17	
cis-1,3-Dichloropropene	25 U	25	1.0	5	07/21/20 16:17	
trans-1,2-Dichloroethene	25 U	25	1.0	5	07/21/20 16:17	
trans-1,3-Dichloropropene	25 U	25	1.2	5	07/21/20 16:17	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 13:15

Monitoring

Sample Matrix: Water Date Received: 07/11/20 09:05

 Sample Name:
 WG-9954-070920-SG-035
 Units: ug/L

 Lab Code:
 R2005980-003
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	97	85 - 122	07/21/20 16:17	
Dibromofluoromethane	103	89 - 119	07/21/20 16:17	
Toluene-d8	103	87 - 121	07/21/20 16:17	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 13:15

Monitoring

Sample Matrix: Water Date Received: 07/11/20 09:05

 Sample Name:
 WG-9954-070920-SG-035
 Units: ug/L

 Lab Code:
 R2005980-003
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
007446-09-5	Sulfur dioxide	1.32	1552.5	JN	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 00:00

Monitoring

Sample Matrix: Water Date Received: 07/11/20 09:05

 Sample Name:
 TB-9954-070920-SG-007
 Units: ug/L

 Lab Code:
 R2005980-004
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/21/20 15:34	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/21/20 15:34	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/21/20 15:34	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/21/20 15:34	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/21/20 15:34	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/21/20 15:34	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/21/20 15:34	
2-Butanone (MEK)	10 U	10	0.78	1	07/21/20 15:34	
2-Hexanone	10 U	10	0.20	1	07/21/20 15:34	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/21/20 15:34	
Acetone	10 U	10	5.0	1	07/21/20 15:34	
Benzene	5.0 U	5.0	0.20	1	07/21/20 15:34	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/21/20 15:34	
Bromoform	5.0 U	5.0	0.25	1	07/21/20 15:34	
Bromomethane	5.0 U	5.0	0.70	1	07/21/20 15:34	
Carbon Disulfide	10 U	10	0.42	1	07/21/20 15:34	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/21/20 15:34	
Chlorobenzene	5.0 U	5.0	0.20	1	07/21/20 15:34	
Chloroethane	5.0 U	5.0	0.23	1	07/21/20 15:34	
Chloroform	5.0 U	5.0	0.24	1	07/21/20 15:34	
Chloromethane	5.0 U	5.0	0.28	1	07/21/20 15:34	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/21/20 15:34	
Dichloromethane	5.0 U	5.0	0.65	1	07/21/20 15:34	
Ethylbenzene	5.0 U	5.0	0.20	1	07/21/20 15:34	
Styrene	5.0 U	5.0	0.20	1	07/21/20 15:34	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/21/20 15:34	
Toluene	5.0 U	5.0	0.20	1	07/21/20 15:34	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/21/20 15:34	
Vinyl Acetate	10 U	10	1.1	1	07/21/20 15:34	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/21/20 15:34	
Xylenes, Total	5.0 U	5.0	0.23	1	07/21/20 15:34	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/21/20 15:34	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/21/20 15:34	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/21/20 15:34	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/21/20 15:34	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 00:00

Monitoring

Sample Matrix: Water Date Received: 07/11/20 09:05

 Sample Name:
 TB-9954-070920-SG-007
 Units: ug/L

 Lab Code:
 R2005980-004
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	96	85 - 122	07/21/20 15:34	
Dibromofluoromethane	97	89 - 119	07/21/20 15:34	
Toluene-d8	102	87 - 121	07/21/20 15:34	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 00:00

Monitoring

Sample Matrix: Water Date Received: 07/11/20 09:05

 Sample Name:
 TB-9954-070920-SG-007
 Units: ug/L

 Lab Code:
 R2005980-004
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Tentatively Identified Compounds

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 08:15

Monitoring

Sample Matrix: Water Date Received: 07/11/20 13:09

 Sample Name:
 WG-9954-070920-SG-036
 Units: ug/L

 Lab Code:
 R2005980-005
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/21/20 15:55	_
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/21/20 15:55	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/21/20 15:55	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/21/20 15:55	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/21/20 15:55	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/21/20 15:55	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/21/20 15:55	
2-Butanone (MEK)	10 U	10	0.78	1	07/21/20 15:55	
2-Hexanone	10 U	10	0.20	1	07/21/20 15:55	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/21/20 15:55	
Acetone	10 U	10	5.0	1	07/21/20 15:55	
Benzene	5.0 U	5.0	0.20	1	07/21/20 15:55	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/21/20 15:55	
Bromoform	5.0 U	5.0	0.25	1	07/21/20 15:55	
Bromomethane	5.0 U	5.0	0.70	1	07/21/20 15:55	
Carbon Disulfide	10 U	10	0.42	1	07/21/20 15:55	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/21/20 15:55	
Chlorobenzene	5.0 U	5.0	0.20	1	07/21/20 15:55	
Chloroethane	5.0 U	5.0	0.23	1	07/21/20 15:55	
Chloroform	5.0 U	5.0	0.24	1	07/21/20 15:55	
Chloromethane	5.0 U	5.0	0.28	1	07/21/20 15:55	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/21/20 15:55	
Dichloromethane	5.0 U	5.0	0.65	1	07/21/20 15:55	
Ethylbenzene	5.0 U	5.0	0.20	1	07/21/20 15:55	
Styrene	5.0 U	5.0	0.20	1	07/21/20 15:55	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/21/20 15:55	
Toluene	5.0 U	5.0	0.20	1	07/21/20 15:55	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/21/20 15:55	
Vinyl Acetate	10 U	10	1.1	1	07/21/20 15:55	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/21/20 15:55	
Xylenes, Total	5.0 U	5.0	0.23	1	07/21/20 15:55	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/21/20 15:55	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/21/20 15:55	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/21/20 15:55	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/21/20 15:55	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 08:15

Monitoring

Sample Matrix: Water Date Received: 07/11/20 13:09

 Sample Name:
 WG-9954-070920-SG-036
 Units: ug/L

 Lab Code:
 R2005980-005
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	93	85 - 122	07/21/20 15:55	
Dibromofluoromethane	99	89 - 119	07/21/20 15:55	
Toluene-d8	100	87 - 121	07/21/20 15:55	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 08:15

Monitoring

Sample Matrix: Water Date Received: 07/11/20 13:09

 Sample Name:
 WG-9954-070920-SG-036
 Units: ug/L

 Lab Code:
 R2005980-005
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected



Semivolatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 12:25

Monitoring

Sample Matrix: Water Date Received: 07/11/20 09:05

 Sample Name:
 WG-9954-070920-SG-033
 Units: ug/L

 Lab Code:
 R2005980-001
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/20/20 15:12	7/15/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/20/20 15:12	7/15/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/20/20 15:12	7/15/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/20/20 15:12	7/15/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/20/20 15:12	7/15/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/20/20 15:12	7/15/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/20/20 15:12	7/15/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/20/20 15:12	7/15/20	
2,4-Dinitrophenol	45 U	45	19	1	07/20/20 15:12	7/15/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/20/20 15:12	7/15/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/20/20 15:12	7/15/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/20/20 15:12	7/15/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/20/20 15:12	7/15/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/20/20 15:12	7/15/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/20/20 15:12	7/15/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/20/20 15:12	7/15/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/20/20 15:12	7/15/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/20/20 15:12	7/15/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/20/20 15:12	7/15/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/20/20 15:12	7/15/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/20/20 15:12	7/15/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/20/20 15:12	7/15/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/20/20 15:12	7/15/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/20/20 15:12	7/15/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/20/20 15:12	7/15/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/20/20 15:12	7/15/20	
4-Nitrophenol	45 U	45	5.8	1	07/20/20 15:12	7/15/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/20/20 15:12	7/15/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/20/20 15:12	7/15/20	
Anthracene	9.1 U	9.1	1.2	1	07/20/20 15:12	7/15/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/20/20 15:12	7/15/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/20/20 15:12	7/15/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/20/20 15:12	7/15/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/20/20 15:12	7/15/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/20/20 15:12	7/15/20	
Benzoic Acid	91 U	91	33	1	07/20/20 15:12	7/15/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/20/20 15:12	7/15/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/20/20 15:12	7/15/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/20/20 15:12	7/15/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	07/20/20 15:12	7/15/20	
Bis(2-ethylhexyl) Phthalate	3.9 J	9.1	0.91	1	07/20/20 15:12	7/15/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	07/20/20 15:12	7/15/20	
Chrysene	9.1 U	9.1	1.1	1	07/20/20 15:12	7/15/20	

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Superset Reference: 20-0000556447 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 12:25

Monitoring

Sample Matrix: Water Date Received: 07/11/20 09:05

 Sample Name:
 WG-9954-070920-SG-033
 Units: ug/L

 Lab Code:
 R2005980-001
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	2.3 Ј	9.1	1.9	1	07/20/20 15:12	7/15/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/20/20 15:12	7/15/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/20/20 15:12	7/15/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/20/20 15:12	7/15/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/20/20 15:12	7/15/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/20/20 15:12	7/15/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/20/20 15:12	7/15/20	
Fluorene	9.1 U	9.1	1.2	1	07/20/20 15:12	7/15/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/20/20 15:12	7/15/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/20/20 15:12	7/15/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/20/20 15:12	7/15/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/20/20 15:12	7/15/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/20/20 15:12	7/15/20	
Isophorone	9.1 U	9.1	1.3	1	07/20/20 15:12	7/15/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/20/20 15:12	7/15/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/20/20 15:12	7/15/20	
Naphthalene	9.1 U	9.1	1.1	1	07/20/20 15:12	7/15/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/20/20 15:12	7/15/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/20/20 15:12	7/15/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/20/20 15:12	7/15/20	
Phenol	9.1 U	9.1	0.91	1	07/20/20 15:12	7/15/20	
Pyrene	9.1 U	9.1	1.3	1	07/20/20 15:12	7/15/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	106	35 - 141	07/20/20 15:12	
2-Fluorobiphenyl	66	31 - 118	07/20/20 15:12	
2-Fluorophenol	53	10 - 105	07/20/20 15:12	
Nitrobenzene-d5	68	31 - 110	07/20/20 15:12	
Phenol-d6	47	10 - 107	07/20/20 15:12	
p-Terphenyl-d14	55	10 - 165	07/20/20 15:12	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown hydrocarbon	11.34	4.3	J	
	unknown	11.87	7.2	J	
	unknown hydrocarbon	12.47	9.5	J	
	unknown hydrocarbon	13.11	10	J	
	unknown hydrocarbon	13.81	8.6	J	
	unknown hydrocarbon	14.56	10	J	
	unknown hydrocarbon	15.34	7.1	J	

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Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 12:25

Monitoring

Sample Matrix: Water Date Received: 07/11/20 09:05

 Sample Name:
 WG-9954-070920-SG-033
 Units: ug/L

 Lab Code:
 R2005980-001
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown hydrocarbon	16.10	11	J	
000624-92-0	Disulfide, dimethyl	2.47	21	JN	
	unknown	6.08	4.0	J	
013798-23-7	Sulfur	7.74	8.7	JN	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 12:55

Monitoring

Sample Matrix: Water Date Received: 07/11/20 09:05

 Sample Name:
 WG-9954-070920-SG-034
 Units: ug/L

 Lab Code:
 R2005980-002
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/20/20 15:40	7/15/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/20/20 15:40	7/15/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/20/20 15:40	7/15/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/20/20 15:40	7/15/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/20/20 15:40	7/15/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/20/20 15:40	7/15/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/20/20 15:40	7/15/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/20/20 15:40	7/15/20	
2,4-Dinitrophenol	45 U	45	19	1	07/20/20 15:40	7/15/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/20/20 15:40	7/15/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/20/20 15:40	7/15/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/20/20 15:40	7/15/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/20/20 15:40	7/15/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/20/20 15:40	7/15/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/20/20 15:40	7/15/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/20/20 15:40	7/15/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/20/20 15:40	7/15/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/20/20 15:40	7/15/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/20/20 15:40	7/15/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/20/20 15:40	7/15/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/20/20 15:40	7/15/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/20/20 15:40	7/15/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/20/20 15:40	7/15/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/20/20 15:40	7/15/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/20/20 15:40	7/15/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/20/20 15:40	7/15/20	
4-Nitrophenol	45 U	45	5.8	1	07/20/20 15:40	7/15/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/20/20 15:40	7/15/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/20/20 15:40	7/15/20	
Anthracene	9.1 U	9.1	1.2	1	07/20/20 15:40	7/15/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/20/20 15:40	7/15/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/20/20 15:40	7/15/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/20/20 15:40	7/15/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/20/20 15:40	7/15/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/20/20 15:40	7/15/20	
Benzoic Acid	91 U	91	33	1	07/20/20 15:40	7/15/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/20/20 15:40	7/15/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/20/20 15:40	7/15/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/20/20 15:40	7/15/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	07/20/20 15:40	7/15/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	07/20/20 15:40	7/15/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	07/20/20 15:40	7/15/20	
Chrysene	9.1 U	9.1	1.1	1	07/20/20 15:40	7/15/20	

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Superset Reference: 20-0000556447 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 12:55

Monitoring

Sample Matrix: Water Date Received: 07/11/20 09:05

 Sample Name:
 WG-9954-070920-SG-034
 Units: ug/L

 Lab Code:
 R2005980-002
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/20/20 15:40	7/15/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/20/20 15:40	7/15/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/20/20 15:40	7/15/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/20/20 15:40	7/15/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/20/20 15:40	7/15/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/20/20 15:40	7/15/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/20/20 15:40	7/15/20	
Fluorene	9.1 U	9.1	1.2	1	07/20/20 15:40	7/15/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/20/20 15:40	7/15/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/20/20 15:40	7/15/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/20/20 15:40	7/15/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/20/20 15:40	7/15/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/20/20 15:40	7/15/20	
Isophorone	9.1 U	9.1	1.3	1	07/20/20 15:40	7/15/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/20/20 15:40	7/15/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/20/20 15:40	7/15/20	
Naphthalene	1.2 J	9.1	1.1	1	07/20/20 15:40	7/15/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/20/20 15:40	7/15/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/20/20 15:40	7/15/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/20/20 15:40	7/15/20	
Phenol	9.1 U	9.1	0.91	1	07/20/20 15:40	7/15/20	
Pyrene	9.1 U	9.1	1.3	1	07/20/20 15:40	7/15/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	106	35 - 141	07/20/20 15:40	
2-Fluorobiphenyl	69	31 - 118	07/20/20 15:40	
2-Fluorophenol	56	10 - 105	07/20/20 15:40	
Nitrobenzene-d5	67	31 - 110	07/20/20 15:40	
Phenol-d6	60	10 - 107	07/20/20 15:40	
p-Terphenyl-d14	56	10 - 165	07/20/20 15:40	

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown hydrocarbon	11.87	3.7	J	
	unknown hydrocarbon	12.46	4.5	J	
	unknown hydrocarbon	13.11	5.0	J	
	unknown hydrocarbon	13.81	3.7	J	
000624-92-0	Disulfide, dimethyl	2.47	24	JN	
	unknown	3.81	3.8	J	
005756-24-1	Tetrasulfide, dimethyl	5.95	190	JN	

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Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 13:15

Monitoring

Sample Matrix: Water Date Received: 07/11/20 09:05

 Sample Name:
 WG-9954-070920-SG-035
 Units: ug/L

 Lab Code:
 R2005980-003
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/20/20 16:09	7/15/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/20/20 16:09	7/15/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/20/20 16:09	7/15/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/20/20 16:09	7/15/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/20/20 16:09	7/15/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/20/20 16:09	7/15/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/20/20 16:09	7/15/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/20/20 16:09	7/15/20	
2,4-Dinitrophenol	45 U	45	19	1	07/20/20 16:09	7/15/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/20/20 16:09	7/15/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/20/20 16:09	7/15/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/20/20 16:09	7/15/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/20/20 16:09	7/15/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/20/20 16:09	7/15/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/20/20 16:09	7/15/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/20/20 16:09	7/15/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/20/20 16:09	7/15/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/20/20 16:09	7/15/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/20/20 16:09	7/15/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/20/20 16:09	7/15/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/20/20 16:09	7/15/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/20/20 16:09	7/15/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/20/20 16:09	7/15/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/20/20 16:09	7/15/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/20/20 16:09	7/15/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/20/20 16:09	7/15/20	
4-Nitrophenol	45 U	45	5.8	1	07/20/20 16:09	7/15/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/20/20 16:09	7/15/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/20/20 16:09	7/15/20	
Anthracene	9.1 U	9.1	1.2	1	07/20/20 16:09	7/15/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/20/20 16:09	7/15/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/20/20 16:09	7/15/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/20/20 16:09	7/15/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/20/20 16:09	7/15/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/20/20 16:09	7/15/20	
Benzoic Acid	91 U	91	33	1	07/20/20 16:09	7/15/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/20/20 16:09	7/15/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/20/20 16:09	7/15/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/20/20 16:09	7/15/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	07/20/20 16:09	7/15/20	
Bis(2-ethylhexyl) Phthalate	9.1 U	9.1	0.91	1	07/20/20 16:09	7/15/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	07/20/20 16:09	7/15/20	
Chrysene	9.1 U	9.1	1.1	1	07/20/20 16:09	7/15/20	

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Superset Reference: 20-0000556447 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 13:15

Monitoring

Sample Matrix: Water Date Received: 07/11/20 09:05

 Sample Name:
 WG-9954-070920-SG-035
 Units: ug/L

 Lab Code:
 R2005980-003
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/20/20 16:09	7/15/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/20/20 16:09	7/15/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/20/20 16:09	7/15/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/20/20 16:09	7/15/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/20/20 16:09	7/15/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/20/20 16:09	7/15/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/20/20 16:09	7/15/20	
Fluorene	9.1 U	9.1	1.2	1	07/20/20 16:09	7/15/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/20/20 16:09	7/15/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/20/20 16:09	7/15/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/20/20 16:09	7/15/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/20/20 16:09	7/15/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/20/20 16:09	7/15/20	
Isophorone	9.1 U	9.1	1.3	1	07/20/20 16:09	7/15/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/20/20 16:09	7/15/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/20/20 16:09	7/15/20	
Naphthalene	9.1 U	9.1	1.1	1	07/20/20 16:09	7/15/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/20/20 16:09	7/15/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/20/20 16:09	7/15/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/20/20 16:09	7/15/20	
Phenol	9.1 U	9.1	0.91	1	07/20/20 16:09	7/15/20	
Pyrene	9.1 U	9.1	1.3	1	07/20/20 16:09	7/15/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	91	35 - 141	07/20/20 16:09	
2-Fluorobiphenyl	51	31 - 118	07/20/20 16:09	
2-Fluorophenol	33	10 - 105	07/20/20 16:09	
Nitrobenzene-d5	55	31 - 110	07/20/20 16:09	
Phenol-d6	23	10 - 107	07/20/20 16:09	
p-Terphenyl-d14	46	10 - 165	07/20/20 16:09	

Tentatively Identified Compounds

			Result	Q	
CAS#	Compound Identification	RT	ug/L		
	unknown hydrocarbon	11.87	4.6	J	
	unknown hydrocarbon	12.47	6.3	J	
	unknown	13.11	7.0	J	
	unknown hydrocarbon	13.81	5.9	J	
	unknown hydrocarbon	14.56	5.6	J	
	unknown hydrocarbon	15.34	4.1	J	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 08:15

Monitoring

Sample Matrix: Water Date Received: 07/11/20 13:09

 Sample Name:
 WG-9954-070920-SG-036
 Units: ug/L

 Lab Code:
 R2005980-005
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	9.1 U	9.1	1.1	1	07/20/20 16:37	7/15/20	
1,2-Dichlorobenzene	9.1 U	9.1	1.1	1	07/20/20 16:37	7/15/20	
1,3-Dichlorobenzene	9.1 U	9.1	0.92	1	07/20/20 16:37	7/15/20	
1,4-Dichlorobenzene	9.1 U	9.1	1.1	1	07/20/20 16:37	7/15/20	
2,4,5-Trichlorophenol	9.1 U	9.1	0.99	1	07/20/20 16:37	7/15/20	
2,4,6-Trichlorophenol	9.1 U	9.1	1.3	1	07/20/20 16:37	7/15/20	
2,4-Dichlorophenol	9.1 U	9.1	1.2	1	07/20/20 16:37	7/15/20	
2,4-Dimethylphenol	9.1 U	9.1	1.3	1	07/20/20 16:37	7/15/20	
2,4-Dinitrophenol	45 U	45	19	1	07/20/20 16:37	7/15/20	
2,4-Dinitrotoluene	9.1 U	9.1	2.2	1	07/20/20 16:37	7/15/20	
2,6-Dinitrotoluene	9.1 U	9.1	1.2	1	07/20/20 16:37	7/15/20	
2-Chloronaphthalene	9.1 U	9.1	1.3	1	07/20/20 16:37	7/15/20	
2-Chlorophenol	9.1 U	9.1	0.97	1	07/20/20 16:37	7/15/20	
2-Methylnaphthalene	9.1 U	9.1	1.2	1	07/20/20 16:37	7/15/20	
2-Methylphenol	9.1 U	9.1	0.91	1	07/20/20 16:37	7/15/20	
2-Nitroaniline	9.1 U	9.1	1.3	1	07/20/20 16:37	7/15/20	
2-Nitrophenol	9.1 U	9.1	1.4	1	07/20/20 16:37	7/15/20	
3,3'-Dichlorobenzidine	9.1 U	9.1	1.1	1	07/20/20 16:37	7/15/20	
3- and 4-Methylphenol Coelution	9.1 U	9.1	1.1	1	07/20/20 16:37	7/15/20	
3-Nitroaniline	9.1 U	9.1	2.3	1	07/20/20 16:37	7/15/20	
4,6-Dinitro-2-methylphenol	45 U	45	18	1	07/20/20 16:37	7/15/20	
4-Bromophenyl Phenyl Ether	9.1 U	9.1	1.5	1	07/20/20 16:37	7/15/20	
4-Chloro-3-methylphenol	9.1 U	9.1	0.98	1	07/20/20 16:37	7/15/20	
4-Chloroaniline	9.1 U	9.1	0.91	1	07/20/20 16:37	7/15/20	
4-Chlorophenyl Phenyl Ether	9.1 U	9.1	1.4	1	07/20/20 16:37	7/15/20	
4-Nitroaniline	9.1 U	9.1	2.5	1	07/20/20 16:37	7/15/20	
4-Nitrophenol	45 U	45	5.8	1	07/20/20 16:37	7/15/20	
Acenaphthene	9.1 U	9.1	1.3	1	07/20/20 16:37	7/15/20	
Acenaphthylene	9.1 U	9.1	1.3	1	07/20/20 16:37	7/15/20	
Anthracene	9.1 U	9.1	1.2	1	07/20/20 16:37	7/15/20	
Benz(a)anthracene	9.1 U	9.1	1.5	1	07/20/20 16:37	7/15/20	
Benzo(a)pyrene	9.1 U	9.1	1.1	1	07/20/20 16:37	7/15/20	
Benzo(b)fluoranthene	9.1 U	9.1	1.1	1	07/20/20 16:37	7/15/20	
Benzo(g,h,i)perylene	9.1 U	9.1	0.91	1	07/20/20 16:37	7/15/20	
Benzo(k)fluoranthene	9.1 U	9.1	1.1	1	07/20/20 16:37	7/15/20	
Benzoic Acid	91 U	91	33	1	07/20/20 16:37	7/15/20	
Benzyl Alcohol	9.1 U	9.1	1.5	1	07/20/20 16:37	7/15/20	
2,2'-Oxybis(1-chloropropane)	9.1 U	9.1	1.3	1	07/20/20 16:37	7/15/20	
Bis(2-chloroethoxy)methane	9.1 U	9.1	1.8	1	07/20/20 16:37	7/15/20	
Bis(2-chloroethyl) Ether	9.1 U	9.1	1.2	1	07/20/20 16:37	7/15/20	
Bis(2-ethylhexyl) Phthalate	1.3 J	9.1	0.91	1	07/20/20 16:37	7/15/20	
Butyl Benzyl Phthalate	9.1 U	9.1	1.3	1	07/20/20 16:37	7/15/20	
Chrysene	9.1 U	9.1	1.1	1	07/20/20 16:37	7/15/20	

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Superset Reference: 20-0000556447 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 08:15

Monitoring

Sample Matrix: Water Date Received: 07/11/20 13:09

 Sample Name:
 WG-9954-070920-SG-036
 Units: ug/L

 Lab Code:
 R2005980-005
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	9.1 U	9.1	1.9	1	07/20/20 16:37	7/15/20	
Di-n-octyl Phthalate	9.1 U	9.1	3.0	1	07/20/20 16:37	7/15/20	
Dibenz(a,h)anthracene	9.1 U	9.1	0.93	1	07/20/20 16:37	7/15/20	
Dibenzofuran	9.1 U	9.1	1.3	1	07/20/20 16:37	7/15/20	
Diethyl Phthalate	9.1 U	9.1	0.98	1	07/20/20 16:37	7/15/20	
Dimethyl Phthalate	9.1 U	9.1	1.2	1	07/20/20 16:37	7/15/20	
Fluoranthene	9.1 U	9.1	1.4	1	07/20/20 16:37	7/15/20	
Fluorene	9.1 U	9.1	1.2	1	07/20/20 16:37	7/15/20	
Hexachlorobenzene	9.1 U	9.1	1.4	1	07/20/20 16:37	7/15/20	
Hexachlorobutadiene	9.1 U	9.1	0.91	1	07/20/20 16:37	7/15/20	
Hexachlorocyclopentadiene	9.1 U	9.1	2.0	1	07/20/20 16:37	7/15/20	
Hexachloroethane	9.1 U	9.1	0.96	1	07/20/20 16:37	7/15/20	
Indeno(1,2,3-cd)pyrene	9.1 U	9.1	1.6	1	07/20/20 16:37	7/15/20	
Isophorone	9.1 U	9.1	1.3	1	07/20/20 16:37	7/15/20	
N-Nitrosodi-n-propylamine	9.1 U	9.1	1.1	1	07/20/20 16:37	7/15/20	
N-Nitrosodiphenylamine	9.1 U	9.1	2.4	1	07/20/20 16:37	7/15/20	
Naphthalene	9.1 U	9.1	1.1	1	07/20/20 16:37	7/15/20	
Nitrobenzene	9.1 U	9.1	1.4	1	07/20/20 16:37	7/15/20	
Pentachlorophenol (PCP)	45 U	45	8.9	1	07/20/20 16:37	7/15/20	
Phenanthrene	9.1 U	9.1	1.3	1	07/20/20 16:37	7/15/20	
Phenol	9.1 U	9.1	0.91	1	07/20/20 16:37	7/15/20	
Pyrene	9.1 U	9.1	1.3	1	07/20/20 16:37	7/15/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	69	35 - 141	07/20/20 16:37	
2-Fluorobiphenyl	56	31 - 118	07/20/20 16:37	
2-Fluorophenol	40	10 - 105	07/20/20 16:37	
Nitrobenzene-d5	59	31 - 110	07/20/20 16:37	
Phenol-d6	28	10 - 107	07/20/20 16:37	
p-Terphenyl-d14	53	10 - 165	07/20/20 16:37	

Tentatively Identified Compounds

			Result	
CAS#	Compound Identification	RT	ug/L	Q

No Tentatively Identified Compounds

Detected



Semivolatile Organic Compounds by GC

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 12:25

Monitoring

Sample Matrix: Water Date Received: 07/11/20 09:05

 Sample Name:
 WG-9954-070920-SG-033
 Units: ug/L

 Lab Code:
 R2005980-001
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/15/20 15:26	7/14/20	
4,4'-DDE	0.045 U	0.045	0.019	1	07/15/20 15:26	7/14/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/15/20 15:26	7/14/20	
Aldrin	0.045 U	0.045	0.019	1	07/15/20 15:26	7/14/20	
Dieldrin	0.045 U	0.045	0.019	1	07/15/20 15:26	7/14/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/15/20 15:26	7/14/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/15/20 15:26	7/14/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/15/20 15:26	7/14/20	
Endrin	0.045 U	0.045	0.019	1	07/15/20 15:26	7/14/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/15/20 15:26	7/14/20	
Heptachlor	0.045 U	0.045	0.019	1	07/15/20 15:26	7/14/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/15/20 15:26	7/14/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/15/20 15:26	7/14/20	
Toxaphene	0.46 U	0.46	0.46	1	07/15/20 15:26	7/14/20	
alpha-BHC	0.12	0.045	0.019	1	07/15/20 15:26	7/14/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/15/20 15:26	7/14/20	
beta-BHC	0.045 U	0.045	0.019	1	07/15/20 15:26	7/14/20	
delta-BHC	0.088	0.045	0.019	1	07/15/20 15:26	7/14/20	
gamma-BHC (Lindane)	0.15	0.045	0.019	1	07/15/20 15:26	7/14/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/15/20 15:26	7/14/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	27	10 - 164	07/15/20 15:26	
Tetrachloro-m-xylene	59	10 - 147	07/15/20 15:26	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 12:55

Monitoring

Sample Matrix: Water Date Received: 07/11/20 09:05

 Sample Name:
 WG-9954-070920-SG-034
 Units: ug/L

 Lab Code:
 R2005980-002
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/25/20 00:02	7/14/20	_
4,4'-DDE	0.045 U	0.045	0.019	1	07/25/20 00:02	7/14/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/25/20 00:02	7/14/20	
Aldrin	0.045 U	0.045	0.019	1	07/25/20 00:02	7/14/20	
Dieldrin	0.045 U	0.045	0.019	1	07/25/20 00:02	7/14/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/25/20 00:02	7/14/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/25/20 00:02	7/14/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/25/20 00:02	7/14/20	
Endrin	0.045 U	0.045	0.019	1	07/25/20 00:02	7/14/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/25/20 00:02	7/14/20	
Heptachlor	0.045 U	0.045	0.019	1	07/25/20 00:02	7/14/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/25/20 00:02	7/14/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/25/20 00:02	7/14/20	
Toxaphene	0.46 U	0.46	0.46	1	07/25/20 00:02	7/14/20	
alpha-BHC	0.069	0.045	0.019	1	07/25/20 00:02	7/14/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/25/20 00:02	7/14/20	
beta-BHC	0.045 U	0.045	0.019	1	07/25/20 00:02	7/14/20	
delta-BHC	0.065	0.045	0.019	1	07/25/20 00:02	7/14/20	
gamma-BHC (Lindane)	0.077	0.045	0.019	1	07/25/20 00:02	7/14/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/25/20 00:02	7/14/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	69	10 - 164	07/25/20 00:02	
Tetrachloro-m-xylene	62	10 - 147	07/25/20 00:02	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 13:15

Monitoring

Sample Matrix: Water Date Received: 07/11/20 09:05

 Sample Name:
 WG-9954-070920-SG-035
 Units: ug/L

 Lab Code:
 R2005980-003
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/25/20 00:21	7/14/20	
4,4'-DDE	0.045 U	0.045	0.019	1	07/25/20 00:21	7/14/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/25/20 00:21	7/14/20	
Aldrin	0.045 U	0.045	0.019	1	07/25/20 00:21	7/14/20	
Dieldrin	0.045 U	0.045	0.019	1	07/25/20 00:21	7/14/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/25/20 00:21	7/14/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/25/20 00:21	7/14/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/25/20 00:21	7/14/20	
Endrin	0.045 U	0.045	0.019	1	07/25/20 00:21	7/14/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/25/20 00:21	7/14/20	
Heptachlor	0.045 U	0.045	0.019	1	07/25/20 00:21	7/14/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/25/20 00:21	7/14/20	
Methoxychlor	0.045 U	0.045	0.019	1	07/25/20 00:21	7/14/20	
Toxaphene	0.46 U	0.46	0.46	1	07/25/20 00:21	7/14/20	
alpha-BHC	0.072	0.045	0.019	1	07/25/20 00:21	7/14/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/25/20 00:21	7/14/20	
beta-BHC	0.019 J	0.045	0.019	1	07/25/20 00:21	7/14/20	
delta-BHC	0.19	0.045	0.019	1	07/25/20 00:21	7/14/20	
gamma-BHC (Lindane)	0.11	0.045	0.019	1	07/25/20 00:21	7/14/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/25/20 00:21	7/14/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	42	10 - 164	07/25/20 00:21	
Tetrachloro-m-xylene	52	10 - 147	07/25/20 00:21	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 08:15

Monitoring

Sample Matrix: Water Date Received: 07/11/20 13:09

 Sample Name:
 WG-9954-070920-SG-036
 Units: ug/L

 Lab Code:
 R2005980-005
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.045 U	0.045	0.019	1	07/15/20 16:24	7/14/20	
4,4'-DDE	0.045 U	0.045	0.019	1	07/15/20 16:24	7/14/20	
4,4'-DDT	0.045 U	0.045	0.019	1	07/15/20 16:24	7/14/20	
Aldrin	0.045 U	0.045	0.019	1	07/15/20 16:24	7/14/20	
Dieldrin	0.045 U	0.045	0.019	1	07/15/20 16:24	7/14/20	
Endosulfan I	0.045 U	0.045	0.019	1	07/15/20 16:24	7/14/20	
Endosulfan II	0.045 U	0.045	0.019	1	07/15/20 16:24	7/14/20	
Endosulfan Sulfate	0.045 U	0.045	0.019	1	07/15/20 16:24	7/14/20	
Endrin	0.045 U	0.045	0.019	1	07/15/20 16:24	7/14/20	
Endrin Ketone	0.045 U	0.045	0.019	1	07/15/20 16:24	7/14/20	
Heptachlor	0.045 U	0.045	0.019	1	07/15/20 16:24	7/14/20	
Heptachlor Epoxide	0.045 U	0.045	0.019	1	07/15/20 16:24	7/14/20	
Methoxychlor	0.041 J	0.045	0.019	1	07/15/20 16:24	7/14/20	
Toxaphene	0.46 U	0.46	0.46	1	07/15/20 16:24	7/14/20	
alpha-BHC	0.069	0.045	0.019	1	07/15/20 16:24	7/14/20	
alpha-Chlordane	0.045 U	0.045	0.019	1	07/15/20 16:24	7/14/20	
beta-BHC	0.045 U	0.045	0.019	1	07/15/20 16:24	7/14/20	
delta-BHC	0.021 J	0.045	0.019	1	07/15/20 16:24	7/14/20	
gamma-BHC (Lindane)	0.058	0.045	0.019	1	07/15/20 16:24	7/14/20	
gamma-Chlordane	0.045 U	0.045	0.019	1	07/15/20 16:24	7/14/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	44	10 - 164	07/15/20 16:24	
Tetrachloro-m-xylene	61	10 - 147	07/15/20 16:24	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 12:25

Monitoring

Sample Matrix: Water Date Received: 07/11/20 09:05

 Sample Name:
 WG-9954-070920-SG-033
 Units: ug/L

 Lab Code:
 R2005980-001
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/22/20 16:41	7/14/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/22/20 16:41	7/14/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/22/20 16:41	7/14/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/22/20 16:41	7/14/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/22/20 16:41	7/14/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/22/20 16:41	7/14/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/22/20 16:41	7/14/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	22	10 - 152	07/22/20 16:41	
Tetrachloro-m-xylene	45	14 - 129	07/22/20 16:41	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 12:55

Monitoring

Sample Matrix: Water Date Received: 07/11/20 09:05

 Sample Name:
 WG-9954-070920-SG-034
 Units: ug/L

 Lab Code:
 R2005980-002
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/22/20 17:01	7/14/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/22/20 17:01	7/14/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/22/20 17:01	7/14/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/22/20 17:01	7/14/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/22/20 17:01	7/14/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/22/20 17:01	7/14/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/22/20 17:01	7/14/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	55	10 - 152	07/22/20 17:01	
Tetrachloro-m-xylene	45	14 - 129	07/22/20 17:01	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 13:15

Monitoring

Sample Matrix: Water Date Received: 07/11/20 09:05

 Sample Name:
 WG-9954-070920-SG-035
 Units: ug/L

 Lab Code:
 R2005980-003
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/22/20 17:21	7/14/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/22/20 17:21	7/14/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/22/20 17:21	7/14/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/22/20 17:21	7/14/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/22/20 17:21	7/14/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/22/20 17:21	7/14/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/22/20 17:21	7/14/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	41	10 - 152	07/22/20 17:21	
Tetrachloro-m-xylene	47	14 - 129	07/22/20 17:21	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term **Date Collected:** 07/09/20 08:15

Monitoring

Sample Matrix: Water Date Received: 07/11/20 13:09

 Sample Name:
 WG-9954-070920-SG-036
 Units: ug/L

 Lab Code:
 R2005980-005
 Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	0.91 U	0.91	0.46	1	07/16/20 17:02	7/14/20	
Aroclor 1221	1.8 U	1.8	0.91	1	07/16/20 17:02	7/14/20	
Aroclor 1232	0.91 U	0.91	0.46	1	07/16/20 17:02	7/14/20	
Aroclor 1242	0.91 U	0.91	0.46	1	07/16/20 17:02	7/14/20	
Aroclor 1248	0.91 U	0.91	0.46	1	07/16/20 17:02	7/14/20	
Aroclor 1254	0.91 U	0.91	0.46	1	07/16/20 17:02	7/14/20	
Aroclor 1260	0.91 U	0.91	0.46	1	07/16/20 17:02	7/14/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	47	10 - 152	07/16/20 17:02	
Tetrachloro-m-xylene	49	14 - 129	07/16/20 17:02	



QC Summary Forms

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com



Volatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

SURROGATE RECOVERY SUMMARYVolatile Organic Compounds by GC/MS

Analysis Method: 8260C

Extraction Method: EPA 5030C

		4-Bromofluorobenzene	Dibromofluoromethane	Toluene-d8
Sample Name	Lab Code	85-122	89-119	87-121
WG-9954-070920-SG-033	R2005980-001	96	101	104
WG-9954-070920-SG-034	R2005980-002	96	102	105
WG-9954-070920-SG-035	R2005980-003	97	103	103
TB-9954-070920-SG-007	R2005980-004	96	97	102
WG-9954-070920-SG-036	R2005980-005	93	99	100
Method Blank	RQ2007864-04	93	101	103
Method Blank	RQ2007931-04	93	99	102
Lab Control Sample	RQ2007864-03	99	101	103
Lab Control Sample	RQ2007931-03	98	101	102

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007864-04Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/21/20 12:31	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/21/20 12:31	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/21/20 12:31	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/21/20 12:31	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/21/20 12:31	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/21/20 12:31	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/21/20 12:31	
2-Butanone (MEK)	10 U	10	0.78	1	07/21/20 12:31	
2-Hexanone	10 U	10	0.20	1	07/21/20 12:31	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/21/20 12:31	
Acetone	10 U	10	5.0	1	07/21/20 12:31	
Benzene	5.0 U	5.0	0.20	1	07/21/20 12:31	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/21/20 12:31	
Bromoform	5.0 U	5.0	0.25	1	07/21/20 12:31	
Bromomethane	5.0 U	5.0	0.70	1	07/21/20 12:31	
Carbon Disulfide	10 U	10	0.42	1	07/21/20 12:31	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/21/20 12:31	
Chlorobenzene	5.0 U	5.0	0.20	1	07/21/20 12:31	
Chloroethane	5.0 U	5.0	0.23	1	07/21/20 12:31	
Chloroform	5.0 U	5.0	0.24	1	07/21/20 12:31	
Chloromethane	5.0 U	5.0	0.28	1	07/21/20 12:31	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/21/20 12:31	
Dichloromethane	5.0 U	5.0	0.65	1	07/21/20 12:31	
Ethylbenzene	5.0 U	5.0	0.20	1	07/21/20 12:31	
Styrene	5.0 U	5.0	0.20	1	07/21/20 12:31	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/21/20 12:31	
Toluene	5.0 U	5.0	0.20	1	07/21/20 12:31	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/21/20 12:31	
Vinyl Acetate	10 U	10	1.1	1	07/21/20 12:31	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/21/20 12:31	
Xylenes, Total	5.0 U	5.0	0.23	1	07/21/20 12:31	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/21/20 12:31	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/21/20 12:31	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/21/20 12:31	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/21/20 12:31	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007864-04Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	93	85 - 122	07/21/20 12:31	
Dibromofluoromethane	101	89 - 119	07/21/20 12:31	
Toluene-d8	103	87 - 121	07/21/20 12:31	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007864-04Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Tentatively Identified Compounds

Result

CAS# Compound Identification RT ug/L Q

No Tentatively Identified Compounds

Detected

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007931-04
 Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C **Prep Method:** EPA 5030C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Q
1,1,1-Trichloroethane (TCA)	5.0 U	5.0	0.20	1	07/22/20 12:26	
1,1,2,2-Tetrachloroethane	5.0 U	5.0	0.20	1	07/22/20 12:26	
1,1,2-Trichloroethane	5.0 U	5.0	0.20	1	07/22/20 12:26	
1,1-Dichloroethane (1,1-DCA)	5.0 U	5.0	0.20	1	07/22/20 12:26	
1,1-Dichloroethene (1,1-DCE)	5.0 U	5.0	0.20	1	07/22/20 12:26	
1,2-Dichloroethane	5.0 U	5.0	0.20	1	07/22/20 12:26	
1,2-Dichloropropane	5.0 U	5.0	0.20	1	07/22/20 12:26	
2-Butanone (MEK)	10 U	10	0.78	1	07/22/20 12:26	
2-Hexanone	10 U	10	0.20	1	07/22/20 12:26	
4-Methyl-2-pentanone	10 U	10	0.20	1	07/22/20 12:26	
Acetone	10 U	10	5.0	1	07/22/20 12:26	
Benzene	5.0 U	5.0	0.20	1	07/22/20 12:26	
Bromodichloromethane	5.0 U	5.0	0.20	1	07/22/20 12:26	
Bromoform	5.0 U	5.0	0.25	1	07/22/20 12:26	
Bromomethane	5.0 U	5.0	0.70	1	07/22/20 12:26	
Carbon Disulfide	10 U	10	0.42	1	07/22/20 12:26	
Carbon Tetrachloride	5.0 U	5.0	0.34	1	07/22/20 12:26	
Chlorobenzene	5.0 U	5.0	0.20	1	07/22/20 12:26	
Chloroethane	5.0 U	5.0	0.23	1	07/22/20 12:26	
Chloroform	5.0 U	5.0	0.24	1	07/22/20 12:26	
Chloromethane	5.0 U	5.0	0.28	1	07/22/20 12:26	
Dibromochloromethane	5.0 U	5.0	0.20	1	07/22/20 12:26	
Dichloromethane	5.0 U	5.0	0.65	1	07/22/20 12:26	
Ethylbenzene	5.0 U	5.0	0.20	1	07/22/20 12:26	
Styrene	5.0 U	5.0	0.20	1	07/22/20 12:26	
Tetrachloroethene (PCE)	5.0 U	5.0	0.21	1	07/22/20 12:26	
Toluene	5.0 U	5.0	0.20	1	07/22/20 12:26	
Trichloroethene (TCE)	5.0 U	5.0	0.20	1	07/22/20 12:26	
Vinyl Acetate	10 U	10	1.1	1	07/22/20 12:26	
Vinyl Chloride	5.0 U	5.0	0.20	1	07/22/20 12:26	
Xylenes, Total	5.0 U	5.0	0.23	1	07/22/20 12:26	
cis-1,2-Dichloroethene	5.0 U	5.0	0.23	1	07/22/20 12:26	
cis-1,3-Dichloropropene	5.0 U	5.0	0.20	1	07/22/20 12:26	
trans-1,2-Dichloroethene	5.0 U	5.0	0.20	1	07/22/20 12:26	
trans-1,3-Dichloropropene	5.0 U	5.0	0.23	1	07/22/20 12:26	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007931-04Basis: NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
4-Bromofluorobenzene	93	85 - 122	07/22/20 12:26	
Dibromofluoromethane	99	89 - 119	07/22/20 12:26	
Toluene-d8	102	87 - 121	07/22/20 12:26	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name: Method Blank Units: ug/L

Lab Code: RQ2007931-04 **Basis:** NA

Volatile Organic Compounds by GC/MS

Analysis Method: 8260C

Prep Method: EPA 5030C

Tentatively Identified Compounds

			Result		
CAS#	Compound Identification	RT	ug/L	Q	
	unknown	1.62	5.7	J	

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/21/20

Sample Matrix: Water

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005980

Lab Control Sample

RQ2007864-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
1,1,1-Trichloroethane (TCA)	8260C	18.4	20.0	92	75-125
1,1,2,2-Tetrachloroethane	8260C	20.8	20.0	104	78-126
1,1,2-Trichloroethane	8260C	19.7	20.0	98	82-121
1,1-Dichloroethane (1,1-DCA)	8260C	18.8	20.0	94	80-124
1,1-Dichloroethene (1,1-DCE)	8260C	21.3	20.0	106	71-118
1,2-Dichloroethane	8260C	19.4	20.0	97	71-127
1,2-Dichloropropane	8260C	20.3	20.0	102	80-119
2-Butanone (MEK)	8260C	19.7	20.0	99	61-137
2-Hexanone	8260C	17.0	20.0	85	63-124
4-Methyl-2-pentanone	8260C	18.2	20.0	91	66-124
Acetone	8260C	20.3	20.0	101	40-161
Benzene	8260C	20.0	20.0	100	79-119
Bromodichloromethane	8260C	19.6	20.0	98	81-123
Bromoform	8260C	18.4	20.0	92	65-146
Bromomethane	8260C	15.6	20.0	78	42-166
Carbon Disulfide	8260C	20.8	20.0	104	66-128
Carbon Tetrachloride	8260C	18.4	20.0	92	70-127
Chlorobenzene	8260C	18.9	20.0	94	80-121
Chloroethane	8260C	19.5	20.0	98	62-131
Chloroform	8260C	19.3	20.0	97	79-120
Chloromethane	8260C	21.8	20.0	109	65-135
Dibromochloromethane	8260C	19.5	20.0	97	72-128
Dichloromethane	8260C	19.0	20.0	95	73-122
Ethylbenzene	8260C	18.8	20.0	94	76-120
Styrene	8260C	19.3	20.0	97	80-124
Tetrachloroethene (PCE)	8260C	17.9	20.0	89	72-125
Toluene	8260C	19.7	20.0	99	79-119
Trichloroethene (TCE)	8260C	17.6	20.0	88	74-122
Vinyl Acetate	8260C	27.8	20.0	139	52-174
Vinyl Chloride	8260C	21.3	20.0	106	74-159
cis-1,2-Dichloroethene	8260C	19.8	20.0	99	80-121
cis-1,3-Dichloropropene	8260C	18.5	20.0	92	77-122
trans-1,2-Dichloroethene	8260C	20.6	20.0	103	73-118
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QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix:

Project:

Water

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

> Units:ug/L Basis:NA

Service Request: R2005980

Date Analyzed: 07/21/20

Lab Control Sample

RQ2007864-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
trans-1 3-Dichloropropene	8260C	18.6	20.0	93	71-133

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/22/20

Sample Matrix: Water

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005980

Lab Control Sample

RQ2007931-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
1,1,1-Trichloroethane (TCA)	8260C	21.2	20.0	106	75-125
1,1,2,2-Tetrachloroethane	8260C	21.9	20.0	110	78-126
1,1,2-Trichloroethane	8260C	20.6	20.0	103	82-121
1,1-Dichloroethane (1,1-DCA)	8260C	21.6	20.0	108	80-124
1,1-Dichloroethene (1,1-DCE)	8260C	24.6	20.0	123 *	71-118
1,2-Dichloroethane	8260C	19.5	20.0	98	71-127
1,2-Dichloropropane	8260C	21.1	20.0	105	80-119
2-Butanone (MEK)	8260C	20.6	20.0	103	61-137
2-Hexanone	8260C	18.9	20.0	94	63-124
4-Methyl-2-pentanone	8260C	19.6	20.0	98	66-124
Acetone	8260C	21.9	20.0	109	40-161
Benzene	8260C	21.1	20.0	106	79-119
Bromodichloromethane	8260C	19.8	20.0	99	81-123
Bromoform	8260C	19.1	20.0	96	65-146
Bromomethane	8260C	17.2	20.0	86	42-166
Carbon Disulfide	8260C	22.2	20.0	111	66-128
Carbon Tetrachloride	8260C	21.2	20.0	106	70-127
Chlorobenzene	8260C	20.8	20.0	104	80-121
Chloroethane	8260C	23.7	20.0	119	62-131
Chloroform	8260C	20.3	20.0	102	79-120
Chloromethane	8260C	24.7	20.0	124	65-135
Dibromochloromethane	8260C	20.5	20.0	102	72-128
Dichloromethane	8260C	21.2	20.0	106	73-122
Ethylbenzene	8260C	21.2	20.0	106	76-120
Styrene	8260C	20.9	20.0	105	80-124
Tetrachloroethene (PCE)	8260C	20.7	20.0	104	72-125
Toluene	8260C	21.2	20.0	106	79-119
Trichloroethene (TCE)	8260C	19.8	20.0	99	74-122
Vinyl Acetate	8260C	32.1	20.0	160	52-174
Vinyl Chloride	8260C	25.4	20.0	127	74-159
cis-1,2-Dichloroethene	8260C	20.9	20.0	104	80-121
cis-1,3-Dichloropropene	8260C	19.4	20.0	97	77-122
trans-1,2-Dichloroethene	8260C	23.9	20.0	119 *	73-118
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QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix:

Project:

Water

Lab Control Sample Summary Volatile Organic Compounds by GC/MS

> Units:ug/L Basis:NA

Service Request: R2005980

Date Analyzed: 07/22/20

Lab Control Sample

RQ2007931-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	% Rec Limits
trans-1,3-Dichloropropene	8260C	19.9	20.0	100	71-133



Semivolatile Organic Compounds by GC/MS

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request: R2005980

Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

Project:

SURROGATE RECOVERY SUMMARY Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Extraction Method:** EPA 3510C

		2,4,6-Tribromophenol	2-Fluorobiphenyl	2-Fluorophenol
Sample Name	Lab Code	35-141	31-118	10-105
WG-9954-070920-SG-033	R2005980-001	106	66	53
WG-9954-070920-SG-034	R2005980-002	106	69	56
WG-9954-070920-SG-035	R2005980-003	91	51	33
WG-9954-070920-SG-036	R2005980-005	69	56	40
Method Blank	RQ2007538-01	90	67	49
Lab Control Sample	RQ2007538-02	94	66	43
Duplicate Lab Control Sample	RQ2007538-03	98	70	43

QA/QC Report

Service Request: R2005980

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

Project:

SURROGATE RECOVERY SUMMARY Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Extraction Method:** EPA 3510C

		Nitrobenzene-d5	Phenol-d6	p-Terphenyl-d14
Sample Name	Lab Code	31-110	10-107	10-165
WG-9954-070920-SG-033	R2005980-001	68	47	55
WG-9954-070920-SG-034	R2005980-002	67	60	56
WG-9954-070920-SG-035	R2005980-003	55	23	46
WG-9954-070920-SG-036	R2005980-005	59	28	53
Method Blank	RQ2007538-01	70	35	70
Lab Control Sample	RQ2007538-02	57	32	64
Duplicate Lab Control Sample	RQ2007538-03	66	32	70

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007538-01
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
1,2,4-Trichlorobenzene	10 U	10	1.2	1	07/20/20 13:13	7/15/20	
1,2-Dichlorobenzene	10 U	10	1.2	1	07/20/20 13:13	7/15/20	
1,3-Dichlorobenzene	10 U	10	1.1	1	07/20/20 13:13	7/15/20	
1,4-Dichlorobenzene	10 U	10	1.2	1	07/20/20 13:13	7/15/20	
2,4,5-Trichlorophenol	10 U	10	1.1	1	07/20/20 13:13	7/15/20	
2,4,6-Trichlorophenol	10 U	10	1.4	1	07/20/20 13:13	7/15/20	
2,4-Dichlorophenol	10 U	10	1.3	1	07/20/20 13:13	7/15/20	
2,4-Dimethylphenol	10 U	10	1.4	1	07/20/20 13:13	7/15/20	
2,4-Dinitrophenol	50 U	50	20	1	07/20/20 13:13	7/15/20	
2,4-Dinitrotoluene	10 U	10	2.4	1	07/20/20 13:13	7/15/20	
2,6-Dinitrotoluene	10 U	10	1.4	1	07/20/20 13:13	7/15/20	
2-Chloronaphthalene	10 U	10	1.4	1	07/20/20 13:13	7/15/20	
2-Chlorophenol	10 U	10	1.1	1	07/20/20 13:13	7/15/20	
2-Methylnaphthalene	10 U	10	1.3	1	07/20/20 13:13	7/15/20	
2-Methylphenol	10 U	10	1.0	1	07/20/20 13:13	7/15/20	
2-Nitroaniline	10 U	10	1.4	1	07/20/20 13:13	7/15/20	
2-Nitrophenol	10 U	10	1.5	1	07/20/20 13:13	7/15/20	
3,3'-Dichlorobenzidine	10 U	10	1.2	1	07/20/20 13:13	7/15/20	
3- and 4-Methylphenol Coelution	10 U	10	1.2	1	07/20/20 13:13	7/15/20	
3-Nitroaniline	10 U	10	2.5	1	07/20/20 13:13	7/15/20	
4,6-Dinitro-2-methylphenol	50 U	50	20	1	07/20/20 13:13	7/15/20	
4-Bromophenyl Phenyl Ether	10 U	10	1.7	1	07/20/20 13:13	7/15/20	
4-Chloro-3-methylphenol	10 U	10	1.1	1	07/20/20 13:13	7/15/20	
4-Chloroaniline	10 U	10	1.0	1	07/20/20 13:13	7/15/20	
4-Chlorophenyl Phenyl Ether	10 U	10	1.5	1	07/20/20 13:13	7/15/20	
4-Nitroaniline	10 U	10	2.7	1	07/20/20 13:13	7/15/20	
4-Nitrophenol	50 U	50	6.4	1	07/20/20 13:13	7/15/20	
Acenaphthene	10 U	10	1.4	1	07/20/20 13:13	7/15/20	
Acenaphthylene	10 U	10	1.4	1	07/20/20 13:13	7/15/20	
Anthracene	10 U	10	1.3	1	07/20/20 13:13	7/15/20	
Benz(a)anthracene	10 U	10	1.6	1	07/20/20 13:13	7/15/20	
Benzo(a)pyrene	10 U	10	1.2	1	07/20/20 13:13	7/15/20	
Benzo(b)fluoranthene	10 U	10	1.2	1	07/20/20 13:13	7/15/20	
Benzo(g,h,i)perylene	10 U	10	1.0	1	07/20/20 13:13	7/15/20	
Benzo(k)fluoranthene	10 U	10	1.3	1	07/20/20 13:13	7/15/20	
Benzoic Acid	100 U	100	36	1	07/20/20 13:13	7/15/20	
Benzyl Alcohol	10 U	10	1.6	1	07/20/20 13:13	7/15/20	
2,2'-Oxybis(1-chloropropane)	10 U	10	1.4	1	07/20/20 13:13	7/15/20	
Bis(2-chloroethoxy)methane	10 U	10	1.9	1	07/20/20 13:13	7/15/20	
Bis(2-chloroethyl) Ether	10 U	10	1.3	1	07/20/20 13:13	7/15/20	
Bis(2-ethylhexyl) Phthalate	10 U	10	1.0	1	07/20/20 13:13	7/15/20	
Butyl Benzyl Phthalate	10 U	10	1.4	1	07/20/20 13:13	7/15/20	
Chrysene	10 U	10	1.2	1	07/20/20 13:13	7/15/20	
Cinybone	10 0	10	1.2	1	01120120 13.13	1/13/40	

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Superset Reference: 20-0000556447 rev 00

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007538-01
 Basis: NA

Semivolatile Organic Compounds by GC/MS

Analysis Method: 8270D **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Di-n-butyl Phthalate	10 U	10	2.0	1	07/20/20 13:13	7/15/20	
Di-n-octyl Phthalate	10 U	10	3.3	1	07/20/20 13:13	7/15/20	
Dibenz(a,h)anthracene	10 U	10	1.1	1	07/20/20 13:13	7/15/20	
Dibenzofuran	10 U	10	1.4	1	07/20/20 13:13	7/15/20	
Diethyl Phthalate	10 U	10	1.1	1	07/20/20 13:13	7/15/20	
Dimethyl Phthalate	10 U	10	1.3	1	07/20/20 13:13	7/15/20	
Fluoranthene	10 U	10	1.5	1	07/20/20 13:13	7/15/20	
Fluorene	10 U	10	1.3	1	07/20/20 13:13	7/15/20	
Hexachlorobenzene	10 U	10	1.6	1	07/20/20 13:13	7/15/20	
Hexachlorobutadiene	10 U	10	1.0	1	07/20/20 13:13	7/15/20	
Hexachlorocyclopentadiene	10 U	10	2.2	1	07/20/20 13:13	7/15/20	
Hexachloroethane	10 U	10	1.1	1	07/20/20 13:13	7/15/20	
Indeno(1,2,3-cd)pyrene	10 U	10	1.8	1	07/20/20 13:13	7/15/20	
Isophorone	10 U	10	1.4	1	07/20/20 13:13	7/15/20	
N-Nitrosodi-n-propylamine	10 U	10	1.2	1	07/20/20 13:13	7/15/20	
N-Nitrosodiphenylamine	10 U	10	2.7	1	07/20/20 13:13	7/15/20	
Naphthalene	10 U	10	1.2	1	07/20/20 13:13	7/15/20	
Nitrobenzene	10 U	10	1.5	1	07/20/20 13:13	7/15/20	
Pentachlorophenol (PCP)	50 U	50	9.8	1	07/20/20 13:13	7/15/20	
Phenanthrene	10 U	10	1.4	1	07/20/20 13:13	7/15/20	
Phenol	10 U	10	1.0	1	07/20/20 13:13	7/15/20	
Pyrene	10 U	10	1.5	1	07/20/20 13:13	7/15/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
2,4,6-Tribromophenol	90	35 - 141	07/20/20 13:13	
2-Fluorobiphenyl	67	31 - 118	07/20/20 13:13	
2-Fluorophenol	49	10 - 105	07/20/20 13:13	
Nitrobenzene-d5	70	31 - 110	07/20/20 13:13	
Phenol-d6	35	10 - 107	07/20/20 13:13	
p-Terphenyl-d14	70	10 - 165	07/20/20 13:13	

Tentatively Identified Compounds

			Result	
CAS#	Compound Identification	RT	ug/L	Q

No Tentatively Identified Compounds

Detected

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/20/20

Sample Matrix: Water

Duplicate Lab Control Sample Summary Semivolatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005980

Lab Control Sample

Duplicate Lab Control Sample

RQ2007538-02

RQ2007538-03

Analyte Name	Analytica l Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
1,2,4-Trichlorobenzene	8270D	42.2	80.0	53	44.0	80.0	55	10-127	4	30
1,2-Dichlorobenzene	8270D	39.6	80.0	50	40.3	80.0	50	23-130	<1	30
1,3-Dichlorobenzene	8270D	40.7	80.0	51	39.4	80.0	49	21-90	4	30
1,4-Dichlorobenzene	8270D	39.0	80.0	49	39.2	80.0	49	10-124	<1	30
2,4,5-Trichlorophenol	8270D	58.4	80.0	73	63.8	80.0	80	48-134	9	30
2,4,6-Trichlorophenol	8270D	53.6	80.0	67	58.8	80.0	73	44-135	9	30
2,4-Dichlorophenol	8270D	46.5	80.0	58	48.0	80.0	60	48-127	3	30
2,4-Dimethylphenol	8270D	51.7	80.0	65	55.8	80.0	70	59-113	7	30
2,4-Dinitrophenol	8270D	54.5	80.0	68	64.5	80.0	81	21-154	17	30
2,4-Dinitrotoluene	8270D	69.0	80.0	86	74.4	80.0	93	54-130	8	30
2,6-Dinitrotoluene	8270D	72.6	80.0	91	81.8	80.0	102	51-127	11	30
2-Chloronaphthalene	8270D	51.7	80.0	65	55.2	80.0	69	40-108	6	30
2-Chlorophenol	8270D	42.7	80.0	53	41.7	80.0	52	42-112	2	30
2-Methylnaphthalene	8270D	50.7	80.0	63	51.9	80.0	65	34-102	3	30
2-Methylphenol	8270D	44.5	80.0	56	48.3	80.0	60	47-100	7	30
2-Nitroaniline	8270D	66.2	80.0	83	74.0	80.0	92	52-133	10	30
2-Nitrophenol	8270D	46.8	80.0	59	50.8	80.0	64	43-131	8	30
3,3'-Dichlorobenzidine	8270D	69.5	80.0	87	74.0	80.0	93	43-126	7	30
3- and 4-Methylphenol Coelution	8270D	40.5	80.0	51	44.5	80.0	56	40-92	9	30
3-Nitroaniline	8270D	63.3	80.0	79	68.5	80.0	86	42-111	8	30
4,6-Dinitro-2-methylphenol	8270D	62.5	80.0	78	67.6	80.0	85	36-152	9	30
4-Bromophenyl Phenyl Ether	8270D	65.6	80.0	82	71.2	80.0	89	48-114	8	30
4-Chloro-3-methylphenol	8270D	60.1	80.0	75	63.4	80.0	79	52-113	5	30
4-Chloroaniline	8270D	54.7	80.0	68	55.8	80.0	70	44-109	3	30
4-Chlorophenyl Phenyl Ether	8270D	56.1	80.0	70	61.0	80.0	76	51-107	8	30
4-Nitroaniline	8270D	64.4	80.0	81	70.5	80.0	88	54-133	8	30
4-Nitrophenol	8270D	24.4 J	80.0	31	26.8 J	80.0	34	10-126	9	30
Acenaphthene	8270D	58.4	80.0	73	62.8	80.0	78	52-107	7	30
Acenaphthylene	8270D	62.4	80.0	78	67.2	80.0	84	55-109	7	30
Anthracene	8270D	66.7	80.0	83	74.4	80.0	93	55-116	11	30
Benz(a)anthracene	8270D	65.3	80.0	82	71.9	80.0	90	61-121	9	30
Benzo(a)pyrene	8270D	69.6	80.0	87	79.9	80.0	100	44-114	14	30
Benzo(b)fluoranthene	8270D	63.6	80.0	80	74.5	80.0	93	62-115	15	30
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QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/20/20

Sample Matrix: Water

Duplicate Lab Control Sample Summary Semivolatile Organic Compounds by GC/MS

Units:ug/L Basis:NA

Service Request: R2005980

Lab Control Sample

Duplicate Lab Control Sample

RQ2007538-02

RQ2007538-03

Analyte Name	Analytica l Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
Benzo(g,h,i)perylene	8270D	71.7	80.0	90	82.0	80.0	102	63-136	13	30
Benzo(k)fluoranthene	8270D	70.1	80.0	88	77.4	80.0	97	49-133	10	30
Benzoic Acid	8270D	53.4 J	120	44	63.7 J	120	53	10-94	19	30
Benzyl Alcohol	8270D	52.8	80.0	66	57.7	80.0	72	31-109	9	30
2,2'-Oxybis(1-chloropropane)	8270D	42.6	80.0	53	49.3	80.0	62	32-122	16	30
Bis(2-chloroethoxy)methane	8270D	52.0	80.0	65	55.9	80.0	70	55-110	7	30
Bis(2-chloroethyl) Ether	8270D	44.1	80.0	55	44.2	80.0	55	46-102	<1	30
Bis(2-ethylhexyl) Phthalate	8270D	80.4	80.0	100	86.5	80.0	108	51-132	8	30
Butyl Benzyl Phthalate	8270D	74.6	80.0	93	79.7	80.0	100	41-148	7	30
Chrysene	8270D	67.1	80.0	84	76.7	80.0	96	57-118	13	30
Di-n-butyl Phthalate	8270D	82.2	80.0	103	87.9	80.0	110	57-128	7	30
Di-n-octyl Phthalate	8270D	80.9	80.0	101	92.3	80.0	115	62-124	13	30
Dibenz(a,h)anthracene	8270D	81.4	80.0	102	91.7	80.0	115	54-135	12	30
Dibenzofuran	8270D	62.2	80.0	78	68.1	80.0	85	55-110	9	30
Diethyl Phthalate	8270D	65.9	80.0	82	68.3	80.0	85	53-113	4	30
Dimethyl Phthalate	8270D	71.9	80.0	90	76.2	80.0	95	51-112	5	30
Fluoranthene	8270D	75.3	80.0	94	82.3	80.0	103	66-127	9	30
Fluorene	8270D	64.1	80.0	80	69.4	80.0	87	54-106	8	30
Hexachlorobenzene	8270D	72.0	80.0	90	82.7	80.0	103	53-123	13	30
Hexachlorobutadiene	8270D	46.9	80.0	59	47.4	80.0	59	16-95	<1	30
Hexachlorocyclopentadiene	8270D	13.6	80.0	17	15.6	80.0	20	10-99	16	30
Hexachloroethane	8270D	39.4	80.0	49	38.8	80.0	49	15-92	<1	30
Indeno(1,2,3-cd)pyrene	8270D	69.2	80.0	87	82.3	80.0	103	62-137	17	30
Isophorone	8270D	45.7	80.0	57	51.7	80.0	65	50-116	13	30
N-Nitrosodi-n-propylamine	8270D	53.9	80.0	67	60.4	80.0	76	49-115	13	30
N-Nitrosodiphenylamine	8270D	78.4	80.0	98	85.6	80.0	107	45-123	9	30
Naphthalene	8270D	45.4	80.0	57	48.1	80.0	60	38-99	5	30
Nitrobenzene	8270D	43.3	80.0	54	49.6	80.0	62	46-108	14	30
Pentachlorophenol (PCP)	8270D	76.9	80.0	96	84.2	80.0	105	29-164	9	30
Phenanthrene	8270D	65.1	80.0	81	72.5	80.0	91	58-118	12	30
Phenol	8270D	28.0	80.0	35	27.6	80.0	35	10-113	<1	30
Pyrene	8270D	73.5	80.0	92	77.6	80.0	97	61-122	5	30

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Superset Reference: 20-0000556447 rev 00



Semivolatile Organic Compounds by GC

ALS Environmental—Rochester Laboratory 1565 Jefferson Road, Building 300, Suite 360, Rochester, NY 14623 Phone (585) 288-5380 Fax (585) 288-8475 www.alsglobal.com

QA/QC Report

Service Request: R2005980

Client: GHD (Formerly Conestoga-Rovers & Associates)

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

SURROGATE RECOVERY SUMMARY

Organochlorine Pesticides by Gas Chromatography

Analysis Method: 8081B **Extraction Method:** EPA 3510C

		Decachlorobiphenyl	Tetrachloro-m-xylene	
Sample Name	Lab Code	10-164	10-147	
WG-9954-070920-SG-033	R2005980-001	27	59	
WG-9954-070920-SG-034	R2005980-002	69	62	
WG-9954-070920-SG-035	R2005980-003	42	52	
WG-9954-070920-SG-036	R2005980-005	44	61	
Method Blank	RQ2007457-01	67	60	
Lab Control Sample	RQ2007457-02	66	67	
Duplicate Lab Control Sample	RQ2007457-03	64	63	

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

 Sample Name:
 Method Blank
 Units: ug/L

 Lab Code:
 RQ2007457-01
 Basis: NA

Organochlorine Pesticides by Gas Chromatography

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
4,4'-DDD	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
4,4'-DDE	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
4,4'-DDT	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
Aldrin	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
Dieldrin	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
Endosulfan I	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
Endosulfan II	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
Endosulfan Sulfate	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
Endrin	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
Endrin Ketone	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
Heptachlor	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
Heptachlor Epoxide	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
Methoxychlor	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
Toxaphene	0.50 U	0.50	0.50	1	07/15/20 14:09	7/14/20	
alpha-BHC	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
alpha-Chlordane	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
beta-BHC	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
delta-BHC	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
gamma-BHC (Lindane)	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	
gamma-Chlordane	0.050 U	0.050	0.020	1	07/15/20 14:09	7/14/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	67	10 - 164	07/15/20 14:09	
Tetrachloro-m-xylene	60	10 - 147	07/15/20 14:09	

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring Date Analyzed: 07/15/20

Sample Matrix: Water

Project:

Duplicate Lab Control Sample Summary Organochlorine Pesticides by Gas Chromatography

Units:ug/L Basis:NA

Service Request: R2005980

Lab Control Sample

Duplicate Lab Control Sample

RQ2007457-02

RQ2007457-03

	Analytical		Spike			Spike		% Rec		RPD
Analyte Name	Method	Result	Amount	% Rec	Result	Amount	% Rec	Limits	RPD	Limit
4,4'-DDD	8081B	0.295	0.400	74	0.278	0.400	69	42-159	6	30
4,4'-DDE	8081B	0.296	0.400	74	0.281	0.400	70	47-147	5	30
4,4'-DDT	8081B	0.317	0.400	79	0.308	0.400	77	41-149	3	30
Aldrin	8081B	0.248	0.400	62	0.240	0.400	60	22-137	3	30
Dieldrin	8081B	0.319	0.400	80	0.314	0.400	79	52-144	1	30
Endosulfan I	8081B	0.312	0.400	78	0.307	0.400	77	52-136	2	30
Endosulfan II	8081B	0.231	0.400	58	0.288	0.400	72	57-138	22	30
Endosulfan Sulfate	8081B	0.228	0.400	57	0.267	0.400	67	34-156	16	30
Endrin	8081B	0.325	0.400	81	0.318	0.400	80	56-143	2	30
Endrin Ketone	8081B	0.288	0.400	72	0.305	0.400	76	59-143	6	30
Heptachlor	8081B	0.246	0.400	62	0.250	0.400	63	32-141	2	30
Heptachlor Epoxide	8081B	0.316	0.400	79	0.309	0.400	77	51-143	2	30
Methoxychlor	8081B	0.311	0.400	78	0.315	0.400	79	56-149	1	30
alpha-BHC	8081B	0.311	0.400	78	0.297	0.400	74	36-151	5	30
alpha-Chlordane	8081B	0.309	0.400	77	0.301	0.400	75	50-139	3	30
beta-BHC	8081B	0.312	0.400	78	0.310	0.400	77	55-149	<1	30
delta-BHC	8081B	0.252	0.400	63	0.279	0.400	70	29-159	10	30
gamma-BHC (Lindane)	8081B	0.315	0.400	79	0.301	0.400	75	41-149	4	30
gamma-Chlordane	8081B	0.304	0.400	76	0.297	0.400	74	50-140	2	30

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Service Request: R2005980 Love Canal:292-402-D02-3100/9954 Annual Long Term

Sample Matrix: Water

Project:

SURROGATE RECOVERY SUMMARY Polychlorinated Biphenyls (PCBs) by GC

Analysis Method: 8082A **Extraction Method:** EPA 3510C

		Decachlorobiphenyl	Tetrachloro-m-xylene	
Sample Name	Lab Code	10-152	14-129	
WG-9954-070920-SG-033	R2005980-001	22	45	
WG-9954-070920-SG-034	R2005980-002	55	45	
WG-9954-070920-SG-035	R2005980-003	41	47	
WG-9954-070920-SG-036	R2005980-005	47	49	
Method Blank	RQ2007457-01	82	59	
Lab Control Sample	RQ2007457-02	76	54	
Duplicate Lab Control Sample	RQ2007457-03	79	59	

ALS Group USA, Corp. dba ALS Environmental

Analytical Report

Client: GHD (Formerly Conestoga-Rovers & Associates) Service Request: R2005980

Project: Love Canal:292-402-D02-3100/9954 Annual Long Term Date Collected: NA

Monitoring

Sample Matrix: Water Date Received: NA

Sample Name:Method BlankUnits: ug/LLab Code:RQ2007457-01Basis: NA

Polychlorinated Biphenyls (PCBs) by GC

Analysis Method: 8082A **Prep Method:** EPA 3510C

Analyte Name	Result	MRL	MDL	Dil.	Date Analyzed	Date Extracted	Q
Aroclor 1016	1.0 U	1.0	0.50	1	07/16/20 15:01	7/14/20	
Aroclor 1221	2.0 U	2.0	1.0	1	07/16/20 15:01	7/14/20	
Aroclor 1232	1.0 U	1.0	0.50	1	07/16/20 15:01	7/14/20	
Aroclor 1242	1.0 U	1.0	0.50	1	07/16/20 15:01	7/14/20	
Aroclor 1248	1.0 U	1.0	0.50	1	07/16/20 15:01	7/14/20	
Aroclor 1254	1.0 U	1.0	0.50	1	07/16/20 15:01	7/14/20	
Aroclor 1260	1.0 U	1.0	0.50	1	07/16/20 15:01	7/14/20	

Surrogate Name	% Rec	Control Limits	Date Analyzed	Q
Decachlorobiphenyl	82	10 - 152	07/16/20 15:01	
Tetrachloro-m-xylene	59	14 - 129	07/16/20 15:01	

ALS Group USA, Corp. dba ALS Environmental

QA/QC Report

Client: GHD (Formerly Conestoga-Rovers & Associates)

Love Canal:292-402-D02-3100/9954 Annual Long Term Monitoring

Sample Matrix:

Project:

Water

Duplicate Lab Control Sample Summary Polychlorinated Biphenyls (PCBs) by GC

> Units:ug/L Basis:NA

Service Request: R2005980

Date Analyzed: 07/16/20

Lab Control Sample

Duplicate Lab Control Sample

RQ2007457-02

RQ2007457-03

Analyte Name	Analytical Method	Result	Spike Amount	% Rec	Result	Spike Amount	% Rec	% Rec Limits	RPD	RPD Limit
Aroclor 1016	8082A	3.23	4.00	81	3.79	4.00	95	49-123	16	30
Aroclor 1260	8082A	3.92	4.00	98	4.38	4.00	110	30-120	11	30

Appendix F Data Validation Memorandum



Memorandum

August 26, 2020

To: Joe Branch [joseph_branch@oxy.com] Ref. No.: 009954

YW

From: Kathy Willy/adh/65 Tel: 716-205-1942

cc: John Pentilchuk, Dennis Hoyt, Maggie Popek

Subject: Analytical Results and Full Validation

Love Canal Annual Long-Term Monitoring Program

Glenn Springs Holdings, Inc. Niagara Falls, New York

June-July 2020

Introduction

This document details a validation of analytical results for water samples collected in support of the Annual Long-Term Monitoring Program at the Love Canal Site during June and July 2020. Samples were submitted to ALS Laboratories located in Rochester, New York. A sample collection and analysis summary is presented in Table 1. The validated analytical results are summarized in Table 2. A summary of the analytical methodology is presented in Table 3.

Full Contract Laboratory Program (CLP) equivalent raw data deliverables were provided by the laboratory. Evaluation of the data was based on information obtained from the finished data sheets, raw data, chain of custody forms, calibration data, blank data, recovery data from surrogate spikes/laboratory control samples (LCS)/matrix spike (MS) samples, and field quality assurance/quality control (QA/QC) samples. The assessment of analytical and in-house data included checks for: data consistency (by observing comparability of duplicate analyses), adherence to accuracy and precision criteria, and transmittal errors.

The QA/QC criteria by which these data have been assessed are outlined in the analytical methods referenced in Table 3 and applicable guidance from the documents entitled:

- "Quality Assurance Project Plan", Appendix B of "Sampling Manual Long-Term Groundwater Monitoring Program", June 2013
- "USEPA Contract Laboratory Program National Functional Guidelines for Organic Data Review",
 United States Environmental Protection Agency (USEPA) 540-R 2016-002, September 2016

Item ii) will subsequently be referred to as the "Guidelines" in this Memorandum.





2. Sample Holding Time and Preservation

The sample holding time criteria for the analyses are summarized in Table 3. Sample chain of custody documents and analytical reports were used to determine sample holding times. All samples were prepared and analyzed within the required holding times.

All samples were properly preserved, delivered on ice, and stored by the laboratory at the required temperature (0-6°C).

 Gas Chromatography/Mass Spectrometry (GC/MS) – Tuning and Mass Calibration (Instrument Performance Check)

3.1 Organic Analyses

Prior to volatile organic compound (VOC) and semi-volatile organic compound (SVOC) analysis, GC/MS instrumentation is tuned to ensure optimization over the mass range of interest. To evaluate instrument tuning, methods require the analysis of specific tuning compounds bromofluorobenzene (BFB) and decafluorotriphenylphosphine (DFTPP), respectively. The resulting spectra must meet the criteria cited in the methods before analysis is initiated. Analysis of the tuning compound must then be repeated every 12 hours throughout sample analysis to ensure the continued optimization of the instrument.

Tuning compounds were analyzed at the required frequency throughout VOC and SVOC analysis periods. All tuning criteria were met indicating that proper optimization of the instrumentation was achieved.

4. Initial Calibration - Organic Analyses

4.1 GC/MS

To quantify VOCs and SVOCs of interest in samples, calibration of the GC/MS over a specific concentration range must be performed. Initially, a five-point calibration curve containing all compounds of interest is analyzed to characterize instrument response for each analyte over a specific concentration range. Linearity of the calibration curve and instrument sensitivity are evaluated against the following criteria:

- i) All relative response factors (RRFs) must be greater than or equal to 0.05 (0.01 for poor responders).
- ii) The percent relative standard deviation (RSD) values must not exceed 20.0 percent (40 percent for poor responders) or a minimum correlation coefficient (R) of 0.995 and minimum coefficient of determination (R²) of 0.99 if linear and quadratic equation calibration curves, respectively, are used.

The initial calibration data for VOCs and SVOCs were reviewed. All compounds met the above criteria for linearity and sensitivity.

4.2 GC

To quantify pesticides, the performance evaluation mixture (PEM) is analyzed at the beginning and end of the initial calibration sequence and throughout the analytical sequence. The results of these analyses are



used to evaluate dichlorodiphenyltrichloroethane (DDT)/endrin breakdown, using the method degradation criteria of <15 percent. PEM standards were analyzed at the required frequency throughout sample analysis, and all method performance criteria were met.

In order to quantify organic compounds of interest by GC, calibration of the gas chromatograph over a specific concentration range must be performed. Initially, a calibration curve consisting of a minimum of five concentration levels is analyzed for all single component compounds of interest and for polychlorinated biphenyls (PCBs) (Aroclors 1016 and 1260). A single calibration standard is analyzed for all other multi-response compounds. Linearity of the calibration curve is acceptable if all RSD values are less than or equal to 20.0 percent or if the correlation coefficient (R) is 0.995 or greater for linear regression curves.

Retention time windows are also calculated from the initial calibration analyses. These windows are then used to identify all compounds of interest in subsequent analyses.

All initial calibration standards were analyzed at the required frequencies. All retention time, peak resolution, and linearity criteria were satisfied as specified in the methods.

5. Continuing Calibration - Organic Analyses

5.1 GC/MS

To ensure that instrument calibration for VOC and SVOC analyses is acceptable throughout the sample analysis period, continuing calibration standards must be analyzed and compared to the initial calibration curve every 12 hours.

The following criteria were employed to evaluate continuing calibration data:

- i) All RRF values must be greater than or equal to 0.05 (0.01 for poor responders).
- ii) Percent difference (%D) values must not exceed 25 percent (40 percent for poor responders).

Calibration standards were analyzed at the required frequency, and the results met the above criteria for instrument sensitivity. Both vinyl acetate and bis(2-chloroethyl)ether demonstrated some variability. A summary of qualified results is presented in Table 4.

5.2 GC

To ensure that the calibration of the instrument for organic analyses by GC is valid throughout the sample analysis period, continuing calibration standards are analyzed and evaluated on a regular basis. To evaluate the continued linearity of the calibration, %D values are calculated for each compound. As specified in the methods, all %D values should not exceed 15 percent. To ensure that compound retention times do not vary over the analysis period, all retention times for continuing calibration compounds must fall within the established retention time windows.

Continuing calibration standards were analyzed at the required frequency. All %D values and compound retention times met the above criteria, indicating acceptable instrument calibration throughout the analysis period.



6. Laboratory Blank Analyses

Method blanks are prepared from a purified matrix and analyzed with investigative samples to determine the existence and magnitude of sample contamination introduced during the analytical procedures.

For this study, laboratory method blanks were analyzed at a minimum frequency of 1 per 20 investigative samples and/or 1 per analytical batch.

Most method blank results were non-detect, indicating that laboratory contamination was not a factor for this investigation. Low concentrations of toluene, chloromethane, and di-n-butylphthalate were reported. All associated sample results with concentrations similar to that found in the blanks were considered to be a reflection of laboratory contamination and were qualified as non-detect. Sample results that were either non-detect or significantly greater in concentration than that found in the blanks would not have been impacted by the potential contamination, and no qualification of the data was required. A summary of qualified results is presented in Table 5.

7. Surrogate Spike Recoveries

In accordance with the methods employed, all samples, blanks, and QC samples analyzed for organics are spiked with surrogate compounds prior to sample extraction and/or analysis. Surrogate recoveries provide a means to evaluate the effects of laboratory performance on individual sample matrices.

All samples submitted for VOC, SVOC, pesticides, and PCB determinations were spiked with the appropriate number of surrogate compounds prior to sample extraction and/or analysis.

Each individual surrogate compound is expected to meet the laboratory control limits with the exception of SVOC analyses. According to the "Guidelines" for SVOC analyses, up to one outlying surrogate in the base/neutral or acid fractions is acceptable as long as the recovery is at least 10 percent.

Surrogate recoveries were assessed against laboratory control limits. Some surrogate recoveries could not be assessed due to necessary secondary dilutions performed on the samples. All other surrogate recoveries were within acceptable limits, with the exception of a severely low (<10 percent) pesticide surrogate recovery and low SVOC surrogate recoveries in one sample. The associated samples results were qualified as follows:

- i) Non-detect pesticide results were rejected due to the demonstrated poor analytical efficiency.
- ii) Positive pesticide results were qualified as estimated to reflect the implied low bias.
- iii) Associated SVOC results were qualified as estimated to reflect the implied low bias.

A summary of qualified results are presented in Table 6.



8. Internal Standards (IS) Analyses

IS data were evaluated for all VOC and SVOC sample analyses.

To ensure that changes in the GC/MS sensitivity and response do not affect sample analysis results, IS compounds are added to each sample prior to analysis. All results are then calculated as a ratio of the IS responses.

The sample IS results were evaluated against the following criteria:

- i) The retention time of the IS must not vary more than ±30 seconds from the associated calibration standard.
- ii) IS area counts must not vary by more than a factor of two (-50 percent to +100 percent) from the associated calibration standard.

All organic IS recoveries and retention times met the above criteria.

9. Laboratory Control Sample Analyses

LCS and/or laboratory control sample duplicates (LCSD) are prepared and analyzed as samples to assess the analytical efficiencies of the methods employed, independent of sample matrix effects. The relative percent difference (RPD) of the LCS/LCSD recoveries is used to evaluate analytical precision.

For this study, LCS/LCSD were analyzed at a minimum frequency of 1 per 20 investigative samples and/or 1 per analytical batch.

The LCS/LCSD contained all compounds of interest. All LCS recoveries and RPDs were within the laboratory control limits, demonstrating acceptable analytical accuracy and precision with the exception of some low SVOC recoveries and outlying RPD values. All associated sample results were qualified as estimated to reflect the potential low bias and indicated variability. A summary of qualified results is presented in Table 7.

10. Matrix Spike/Matrix Spike Duplicate (MS/MSD) Analyses

To evaluate the effects of sample matrices on the distillation process, measurement procedures, and accuracy of a particular analysis, samples are spiked with a known concentration of the analyte of concern and analyzed as MS/MSD samples. The RPD between the MS and MSD is used to assess analytical precision.

MS/MSD analyses were performed as specified in Table 1.

The MS/MSD samples were spiked with all compounds of interest. All percent recoveries and RPD values were within the laboratory control limits, demonstrating acceptable analytical accuracy and precision with the exception of some severely low (<10 percent) SVOC recoveries. The associated non-detect sample results



were rejected due to the demonstrated poor analytical efficiency. A summary of qualified results is presented in Table 8.

11. Field QA/QC Samples

The field QA/QC consisted of seven trip blank samples, two rinse blank samples, and three field duplicate sample sets.

11.1 Trip Blank Sample Analysis

To evaluate contamination from sample collection, transportation, storage, and analytical activities, seven trip blanks were submitted to the laboratory for VOC analysis. All results were non-detect for the compounds of interest with the exception of a low concentration of toluene reported in one blank. All associated sample results with concentrations similar to that found in the trip blank had been previously qualified as non-detect due to potential method blank contamination, and no further qualification of the data was required.

11.2 Rinse Blank Sample Analysis

To assess field decontamination procedures, ambient conditions at the site, and cleanliness of sample containers, two rinse blanks were submitted for analysis, as identified in Table 1. Low concentrations of several target analytes were detected. Associated sample results with concentrations similar to that found in the blanks were qualified as non-detect. Sample results that were either non-detect or significantly greater in concentration than the blanks would not have been impacted, and no qualification of the data was required. A summary of qualified results is presented in Table 9.

11.3 Field Duplicate Sample Analysis

To assess the analytical and sampling protocol precision, three field duplicate samples were collected and submitted "blind" to the laboratory, as specified in Table 1. The RPDs associated with these duplicate samples must be less than 50 percent for water samples. If the reported concentration in either the investigative sample or its duplicate is less than five times the reporting limit (RL), the evaluation criterion is one times the RL value for water samples.

All field duplicate results were within agreement, demonstrating acceptable sampling and analytical precision with the exception of delta-BHC, which showed some variability in results between one sample and its field duplicate. A summary of qualified results is presented in Table 10.

12. Tentatively Identified Compounds (TICS)

Chromatographic peaks recorded during VOC and SVOC sample GC/MS analyses that are not target compounds, surrogates, or IS, are potential TICs.



A summary of the TICs reported by the laboratory is presented in Table 2. Per the "Guidelines", TICs that were present in the method blanks or identified as solvent preservatives/aldol reaction products were rejected and are not included in the table.

13. Dual Column Analysis

Pesticide analyses were performed using dual column analysis. All pesticide results showed good correlation between the two columns (<40 percent).

14. Analyte Reporting

The laboratory reported detected results down to the laboratory's Method Detection Limit (MDL) for each analyte. Positive analyte detections less than the Practical Quantitation Limit (PQL) but greater than the MDL were qualified as estimated (J) in Table 2 unless qualified otherwise in this memorandum. Non-detect results were presented as non-detect at the RL in Table 2.

15. Target Compound Identification

To minimize erroneous compound identification during organic analyses, qualitative criteria including compound retention time and mass spectra (if applicable) were evaluated according to the identification criteria established by the methods. The samples identified in Table 1 were reviewed. The organic compounds reported adhered to the specified identification criteria.

16. Conclusion

Based on the assessment detailed in the foregoing, the data summarized in Table 2 are acceptable with the specific exceptions and qualifications noted herein.

Table 1

Sample Collection and Analysis Summary Love Canal Annual Long-Term Monitoring Program Glenn Springs Holdings, Inc. Niagara Falls, New York June-July 2020

Analysis/Parameters

					Analysis/Parameter			eters	<u>; </u>	
Sample Identification	Location	Matrix	Collection Date (mm/dd/yyyy)	Collection Time (hr:min)	VOCs	SVOCs	Pesticides	PCBs	Comments	
RB-9954-062520-SG-001	_	Water	06/25/2020	13:50	Х	Х	Х	Χ	Rinse Blank	
RB-9954-063020-SG-002	_	Water	06/30/2020	13:00	X	X	X	X	Rinse Blank	
TB-9954-062420-SG-001	_	Water	06/24/2020	-	X	^	^	^	Trip Blank	
TB-9954-062520-SG-002	_	Water	06/25/2020	_	X				Trip Blank	
TB-9954-062620-RM-003	_	Water	06/26/2020	_	X				Trip Blank	
TB-9954-063020-SG-004	_	Water	06/30/2020	_	X				Trip Blank	
TB-9954-070120-SG-005	_	Water	07/01/2020	_	X				Trip Blank	
TB-9954-070620-SG-006		Water	07/06/2020	_	X				Trip Blank	
TB-9954-070920-SG-007	_	Water	07/09/2020	_	X				Trip Blank	
WG-9954-062420-SG-001	7115	Groundwater	06/24/2020	09:10	X	Х	Χ	X	The blank	
WG-9954-062420-SG-002	7125	Groundwater	06/24/2020	09:55	X	X	X	X		
WG-9954-062420-SG-003	7130	Groundwater	06/24/2020	10:30	X	X	X	X		
WG-9954-062420-SG-004	7132	Groundwater	06/24/2020	11:05	X	X	X	X		
WG-9954-062420-SG-005	8115	Groundwater	06/24/2020	11:50	X	X	X	X		
WG-9954-062420-SG-006	8125	Groundwater	06/24/2020	12:45	X	X	X	X		
WG-9954-062420-SG-007	9105	Groundwater	06/24/2020	13:35	X	X	X	X		
WG-9954-062520-SG-008	9113	Groundwater	06/25/2020	09:15	X	X	X	X		
WG-9954-062520-SG-009	9118	Groundwater	06/25/2020	10:05	X	X	X	X		
WG-9954-062520-SG-010	9118	Groundwater	06/25/2020	10:05	X	X	X	X	Field duplicate of sample WG-9954-062520-SG-009	
WG-9954-062520-SG-011	6209	Groundwater	06/25/2020	11:10	X	X	X	X	riola daplicate of cample 11 0 000 1 002020 00 000	
WG-9954-062520-SG-012	9210	Groundwater	06/25/2020	12:25	X	X	X	X	Matrix Spike/Matrix Spike Duplicate	
WG-9954-062520-SG-013	10205	Groundwater	06/25/2020	13:15	X	X	X	X	man woping man woping 2 aprioate	
WG-9954-062620-RM-014	7205	Groundwater	06/26/2020	08:50	X	X	X	X		
WG-9954-062620-RM-015	8210	Groundwater	06/26/2020	10:00	X	Χ	Χ	X		
WG-9954-062620-RM-016	9205	Groundwater	06/26/2020	11:05	X	X	X	X		
WG-9954-062620-RM-017	8106	Groundwater	06/26/2020	11:55	Х	X	X	X		
WG-9954-062620-RM-018	3257	Groundwater	06/26/2020	12:40	X	Χ	Χ	X		
WG-9954-063020-RM-023	MW-01	Groundwater	06/30/2020	12:15	X	X	X	X	Matrix Spike/Matrix Spike Duplicate	
WG-9954-063020-SG-019	10210B	Groundwater	06/30/2020	09:05	Х	Х	X	Χ	The state of the s	
WG-9954-063020-SG-020	10210C	Groundwater	06/30/2020	10:10	X	Х	X	Χ		
WG-9954-063020-SG-021	10210C	Groundwater	06/30/2020	10:10	X	Х	X	X	Field duplicate of sample WG-9954-063020-SG-020	
WG-9954-063020-SG-022	10215	Groundwater	06/30/2020	11:00	X	Х	Χ	Χ		
WG-9954-070120-SG-024	MW-02	Groundwater	07/01/2020	08:45	X	Х	X	Χ		
WG-9954-070120-SG-025	5221	Groundwater	07/01/2020	09:55	X	Х	X	Χ		
WG-9954-070120-SG-026	10225C	Groundwater	07/01/2020	10:45	X	Χ	X	X		
WG-9954-070120-SG-027	10135	Groundwater	07/01/2020	11:50	Χ	Χ	Χ	Χ		

Table 1

Sample Collection and Analysis Summary Love Canal Annual Long-Term Monitoring Program Glenn Springs Holdings, Inc. Niagara Falls, New York June-July 2020

					Analysis/Parameters				
Sample Identification	Location	Matrix	Collection Date (mm/dd/yyyy)	Collection Time (hr:min)	VOCs	SVOCs	Pesticides	PCBs	Comments
WG-9954-070120-SG-028	10135	Groundwater	07/01/2020	11:50	Χ	Х	Χ	X	Field duplicate of sample WG-9954-070120-SG-027
WG-9954-070620-SG-029	10278	Groundwater	07/06/2020	13:55	X	X	X	Χ	
WG-9954-070720-SG-030	10272	Groundwater	07/07/2020	12:15	Χ	Χ	Χ	Χ	
WG-9954-070720-SG-031	10178A	Groundwater	07/07/2020	10:30	X	X	X	Χ	
WG-9954-070720-SG-032	10270	Groundwater	07/07/2020	13:00	Χ	Χ	Χ	Χ	
WG-9954-070920-SG-033	10210A	Groundwater	07/09/2020	12:25	X	X	X	Χ	
WG-9954-070920-SG-034	10225A	Groundwater	07/09/2020	12:55	X	X	X	Χ	
WG-9954-070920-SG-035	10225B	Groundwater	07/09/2020	13:15	X	X	X	Χ	
WG-9954-071020-SG-036	10178B	Groundwater	07/10/2020	08:15	X	X	X	Χ	

Notes:

VOCs - Volatile Organic Compounds SVOCs - Semi-volatile Organic Compounds PCBs - Polychlorinated Biphenyls - Not applicable

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	Location ID: Sample Name: Sample Date:	3257 WG-9954-062620-RM-018 06/26/2020	5221 WG-9954-070120-SG-025 07/01/2020	6209 WG-9954-062520-SG-011 06/25/2020	7115 WG-9954-062420-SG-001 06/24/2020
Parameters	Unit				
Volatile Organic Compounds (VOCs)					
1,1,1-Trichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1,2,2-Tetrachloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1,2-Trichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1-Dichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1-Dichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,2-Dichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,2-Dichloropropane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
2-Butanone (Methyl ethyl ketone) (MEK)	μg/L	10 U	10 U	10 U	10 U
2-Hexanone	μg/L	10 U	10 U	10 U	10 U
4-Methyl-2-pentanone (Methyl isobutyl ketone)		10 U	10 U	10 U	10 U
Acetone	μg/L	10 U	10 U	10 U	10 U
Benzene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Bromodichloromethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Bromoform	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Bromomethane (Methyl bromide)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Carbon disulfide	μg/L	1.2 J	10 U	2.0 J	10 U
Carbon tetrachloride	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chlorobenzene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chloroform (Trichloromethane)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chloromethane (Methyl chloride)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
cis-1,2-Dichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
cis-1,3-Dichloropropene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Dibromochloromethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Ethylbenzene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Methylene chloride	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Styrene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Tetrachloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Toluene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
trans-1,2-Dichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
trans-1,3-Dichloropropene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Trichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Vinyl acetate	μg/L	10 U	10 U	10 U	10 UJ
Vinyl chloride	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Xylenes (total)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U

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Locat Sample Sampl		3257 WG-9954-062620-RM-018 06/26/2020	5221 WG-9954-070120-SG-025 07/01/2020	6209 WG-9954-062520-SG-011 06/25/2020	7115 WG-9954-062420-SG-001 06/24/2020
Parameters	Unit				
VOCs, Tentatively Identified Compounds (TICs)					
1,2,4-Trichlorobenzene	μg/L	-	-	-	-
1-Chloro-2-methyl-benzene	μg/L	-	-	-	-
1-Chloro-3-methylbenzene	μg/L	-	-	-	-
2,6-Dichlorotoluene	μg/L	-	-	-	-
2-Methylbutane	μg/L	-	-	-	-
6-Methyl-5-hepten-2-one	μg/L	-	-	-	-
Dimethyl disulfide	μg/L	-	-	-	-
Dimethyl sulfide	μg/L	-	-	-	-
Dimethyl trisulfide	μg/L	-	-	-	-
Hexanal	μg/L	-	-	-	-
Isobutane	μg/L	-	-	-	-
m-Monochlorobenzotrifluoride	μg/L	-	-	-	-
Methanethiol	μg/L	-	-	-	-
Methoxytrimethyl-silane	μg/L	-	-	-	-
Methylthioethane	μg/L	-	-	-	-
Nonanal	μg/L	-	-	-	-
Propane	μg/L	-	-	-	-
Sulfur dioxide (SO2)	μg/L	35.9 J	56.0 J	-	-
Trimethylfluorosilane	μg/L	-	-	-	-
Unknown	μg/L	-	-	82.8 J	-
Semi-volatile Organic Compounds (SVOCs)					
1,2,4-Trichlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
1,2-Dichlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
1,3-Dichlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
1,4-Dichlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,2'-Oxybis(1-chloropropane) (bis(2-Chloroisopropyl) ether		9.1 U	9.1 U	9.1 U	9.1 U
2,4,5-Trichlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4,6-Trichlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4-Dichlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4-Dimethylphenol	μg/L	9.1 U	9.1 U	9.1 UJ	9.1 U
2,4-Dinitrophenol	μg/L	45 U	45 U	45 U	45 U
2,4-Dinitrotoluene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,6-Dinitrotoluene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Chloronaphthalene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Chlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Methylnaphthalene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Methylphenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U

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	Location ID: Sample Name: Sample Date:	3257 WG-9954-062620-RM-018 06/26/2020	5221 WG-9954-070120-SG-025 07/01/2020	6209 WG-9954-062520-SG-011 06/25/2020	7115 WG-9954-062420-SG-001 06/24/2020
Parameters	Unit				
SVOCs-Continued					
2-Nitroaniline	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Nitrophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
3&4-Methylphenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
3,3'-Dichlorobenzidine	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
3-Nitroaniline	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4,6-Dinitro-2-methylphenol	μg/L	45 U	45 U	45 U	45 U
4-Bromophenyl phenyl ether	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Chloro-3-methylphenol	μg/L	9.1 U	9.1 U	9.1 UJ	9.1 U
4-Chloroaniline	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Chlorophenyl phenyl ether	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Nitroaniline	μg/L	9.1 U	9.1 U	9.1 UJ	9.1 U
4-Nitrophenol	μg/L	45 U	45 U	45 U	45 U
Acenaphthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Acenaphthylene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Anthracene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(a)anthracene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(a)pyrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(b)fluoranthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(g,h,i)perylene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(k)fluoranthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzoic acid	μg/L	91 U	91 U	91 U	91 U
Benzyl alcohol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
bis(2-Chloroethoxy)methane	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
bis(2-Chloroethyl)ether	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
bis(2-Ethylhexyl)phthalate (DEHP)	μg/L	9.1 U	9.1 U	9.1 U	1.7 J
Butyl benzylphthalate (BBP)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Chrysene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Di-n-butylphthalate (DBP)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Di-n-octyl phthalate (DnOP)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Dibenz(a,h)anthracene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Dibenzofuran	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Diethyl phthalate	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Dimethyl phthalate	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Fluoranthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Fluorene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Hexachlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Hexachlorobutadiene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Hexachlorocyclopentadiene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U

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	Location ID: Sample Name: Sample Date:	3257 WG-9954-062620-RM-018 06/26/2020	5221 WG-9954-070120-SG-025 07/01/2020	6209 WG-9954-062520-SG-011 06/25/2020	7115 WG-9954-062420-SG-001 06/24/2020
Parameters	Unit				
SVOCs-Continued					
Hexachloroethane	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Indeno(1,2,3-cd)pyrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Isophorone	μg/L	9.1 U	9.1 U	9.1 UJ	9.1 U
N-Nitrosodi-n-propylamine	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
N-Nitrosodiphenylamine	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Naphthalene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Nitrobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Pentachlorophenol	μg/L	45 U	45 U	45 U	45 U
Phenanthrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Phenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Pyrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
SVOC, TICs					
(4-Methylphenyl)phenylmethanone	μg/L	-	-	-	-
1,4-Dioxane	μg/L	-	-	-	-
1-Chloro-2-methyl-benzene	μg/L	-	7.6 JN	-	-
1-Chloro-3-methylbenzene	μg/L	-	3.8 JN	-	-
1-Methyl-2-(phenylmethyl)-benzene	μg/L	-	-	-	-
2,3-Dichlorobenzoic acid	μg/L	-	-	-	-
2,5-Dichlorobenzoic acid	μg/L	-	-	-	-
2-Ethyl-ethanoic acid	μg/L	-	-	-	-
2-Propyl tridecyl-sulfurous acid	μg/L	-	-	-	-
3,4-Dichlorophenol	μg/L	-	-	-	-
3,5-Dichlorophenol	μg/L	-	-	-	-
3-Benzoylbenzoic acid	μg/L	-	-	-	-
3-Chlorobenzoic acid	μg/L	-	-	-	-
3-Methyl-cyclohexanol	μg/L	-	-	-	-
4-Benzoyl-(rel)-benzoic acid	μg/L	-	-	-	-
4-Chlorobenzoic acid	μg/L	-	-	-	-
4-Chlorophenol	μg/L	-	-	-	-
4-Chlorotoluene	μg/L	-	-	-	-
6-Octadecenoic acid	μg/L	-	-	-	-
Bicyclo[2.2.1]heptan-2-one, 1,3,3-trimethyl	μg/L	-	-	-	-
Chlorobenzene	μg/L	-	-	-	-
Cyclohexanol	μg/L	-	-	-	-
Dimethyl disulfide	μg/L	-	-	-	-
Dimethyl tetrasulfide	μg/L	-	-	-	-
Dodecanoic acid	μg/L	-	-	-	-

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	Location ID: Sample Name: Sample Date:	3257 WG-9954-062620-RM-018 06/26/2020	5221 WG-9954-070120-SG-025 07/01/2020	6209 WG-9954-062520-SG-011 06/25/2020	7115 WG-9954-062420-SG-001 06/24/2020
Parameters	Unit				
SVOCs, TICs-Continued					
Hexadecanoic acid	μg/L	-	-	-	-
Homomenthyl salicylate	μg/L	-	-	-	-
Nonanoic acid	μg/L	-	-	-	-
Sulfur	μg/L	9.1 JN	4.8 JN	8.0 JN	-
Toluene	μg/L	-	-	-	-
trans-2-Methylcyclohexanol	μg/L	-	-	-	-
Unknown	μg/L	52.6 J	-	-	5.1 J
Polychlorinated Biphenyls (PCBs)					
Aroclor-1016 (PCB-1016)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1221 (PCB-1221)	μg/L	1.8 U	1.8 U	1.8 U	1.8 U
Aroclor-1232 (PCB-1232)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1242 (PCB-1242)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1248 (PCB-1248)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1254 (PCB-1254)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1260 (PCB-1260)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Pesticides					
4,4'-DDD	μg/L	R	0.045 U	0.045 U	0.045 U
4,4'-DDE	μg/L	0.029 J	0.045 U	0.045 U	0.045 U
4,4'-DDT	μg/L	R	0.045 U	0.045 U	0.045 U
Aldrin	μg/L	R	0.045 U	0.045 U	0.045 U
alpha-BHC	μg/L	R	0.13	0.31	0.045 U
alpha-Chlordane	μg/L	R	0.045 U	0.045 U	0.045 U
beta-BHC	μg/L	0.045 J	0.045 U	0.045 U	0.045 U
delta-BHC	μg/L	R	0.029 J	0.13	0.045 U
Dieldrin	μg/L	R	0.045 U	0.045 U	0.045 U
Endosulfan I	μg/L	0.037 J	0.045 U	0.045 U	0.045 U
Endosulfan II	μg/L	R	0.045 U	0.045 U	0.045 U
Endosulfan sulfate	μg/L	R	0.045 U	0.045 U	0.045 U
Endrin	μg/L	R	0.045 U	0.045 U	0.045 U
Endrin ketone	μg/L	R	0.045 U	0.045 U	0.045 U
gamma-BHC (lindane)	μg/L	0.026 J	0.076	0.25	0.045 U
gamma-Chlordane	μg/L	0.020 NJ	0.045 U	0.045 U	0.045 U
Heptachlor	μg/L	R	0.045 U	0.045 U	0.045 U
Heptachlor epoxide	μg/L	R	0.045 U	0.045 U	0.045 U
Methoxychlor	μg/L	R	0.045 U	0.045 U	0.045 U
Toxaphene	μg/L	R	0.46 U	0.46 U	0.46 U

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	Location ID: Sample Name: Sample Date:	3257 WG-9954-062620-RM-018 06/26/2020	5221 WG-9954-070120-SG-025 07/01/2020	6209 WG-9954-062520-SG-011 06/25/2020	7115 WG-9954-062420-SG-001 06/24/2020
Parameters	Unit				
Field Parameters					
Temperature, field	Deg C	12.9	13.7	12.8	11.7
Conductivity, field	mS/cm	3.82	3.31	5.19	1.60
Turbidity, field	NTU	13.9	7.22	5.80	369
pH, field	s.u.	6.93	7.56	7.35	7.31

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	Location ID: Sample Name: Sample Date:	7125 WG-9954-062420-SG-002 06/24/2020	7130 WG-9954-062420-SG-003 06/24/2020	7132 WG-9954-062420-SG-004 06/24/2020	7205 WG-9954-062620-RM-014 06/26/2020
Parameters	Unit				
Volatile Organic Compounds (VOCs)					
1,1,1-Trichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1,2,2-Tetrachloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1,2-Trichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1-Dichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1-Dichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,2-Dichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,2-Dichloropropane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
2-Butanone (Methyl ethyl ketone) (MEK)	μg/L	10 U	10 U	10 U	10 U
2-Hexanone	μg/L	10 U	10 U	10 U	10 U
4-Methyl-2-pentanone (Methyl isobutyl ketone) (MIBK) µg/L	10 U	10 U	10 U	2.4 J
Acetone	μg/L	10 U	10 U	10 U	10 U
Benzene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Bromodichloromethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Bromoform	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Bromomethane (Methyl bromide)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Carbon disulfide	μg/L	10 U	10 U	0.83 J	4.0 J
Carbon tetrachloride	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chlorobenzene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chloroform (Trichloromethane)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chloromethane (Methyl chloride)	μg/L	5.0 U	5.0 U	5.0 U	0.30 J
cis-1,2-Dichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
cis-1,3-Dichloropropene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Dibromochloromethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Ethylbenzene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Methylene chloride	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Styrene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Tetrachloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Toluene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
trans-1,2-Dichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
trans-1,3-Dichloropropene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Trichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Vinyl acetate	μg/L	10 U	10 U	10 U	10 UJ
Vinyl chloride	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Xylenes (total)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U

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	Location ID: Sample Name: Sample Date:	7125 WG-9954-062420-SG-002 06/24/2020	7130 WG-9954-062420-SG-003 06/24/2020	7132 WG-9954-062420-SG-004 06/24/2020	7205 WG-9954-062620-RM-014 06/26/2020
Parameters	Unit				
VOCs, Tentatively Identified Compounds (TIC	s)				
1,2,4-Trichlorobenzene	μg/L	-	-	-	-
1-Chloro-2-methyl-benzene	μg/L	-	-	-	-
1-Chloro-3-methylbenzene	μg/L	-	-	-	-
2,6-Dichlorotoluene	μg/L	-	-	-	-
2-Methylbutane	μg/L	-	-	-	-
6-Methyl-5-hepten-2-one	μg/L	-	-	-	10.2 J
Dimethyl disulfide	μg/L	-	-	-	-
Dimethyl sulfide	μg/L	-	-	-	-
Dimethyl trisulfide	μg/L	-	-	-	-
Hexanal	μg/L	-	-	-	-
Isobutane	μg/L	-	-	-	-
m-Monochlorobenzotrifluoride	μg/L	-	-	-	-
Methanethiol	μg/L	-	-	-	-
Methoxytrimethyl-silane	μg/L	-	-	-	-
Methylthioethane	μg/L	-	-	-	-
Nonanal	μg/L	-	-	-	-
Propane	μg/L	-	-	-	-
Sulfur dioxide (SO2)	μg/L	-	-	-	-
Trimethylfluorosilane	μg/L	-	-	-	-
Unknown	μg/L	-	-	-	8.5 J
Semi-volatile Organic Compounds (SVOCs)					
1,2,4-Trichlorobenzene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
1,2-Dichlorobenzene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
1,3-Dichlorobenzene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
1,4-Dichlorobenzene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
2,2'-Oxybis(1-chloropropane) (bis(2-Chloroisopro		9.1 UJ	9.1 U	9.1 U	9.1 U
2,4,5-Trichlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4,6-Trichlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4-Dichlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4-Dimethylphenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4-Dinitrophenol	μg/L	45 U	45 U	45 U	45 U
2,4-Dinitrotoluene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
2,6-Dinitrotoluene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
2-Chloronaphthalene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
2-Chlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Methylnaphthalene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
2-Methylphenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U

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	Location ID: Sample Name: Sample Date:	7125 WG-9954-062420-SG-002 06/24/2020	7130 WG-9954-062420-SG-003 06/24/2020	7132 WG-9954-062420-SG-004 06/24/2020	7205 WG-9954-062620-RM-014 06/26/2020
Parameters	Unit				
SVOCs-Continued					
2-Nitroaniline	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
2-Nitrophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
3&4-Methylphenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
3,3'-Dichlorobenzidine	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
3-Nitroaniline	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
4,6-Dinitro-2-methylphenol	μg/L	45 U	45 U	45 U	45 U
4-Bromophenyl phenyl ether	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
4-Chloro-3-methylphenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Chloroaniline	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
4-Chlorophenyl phenyl ether	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
4-Nitroaniline	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
4-Nitrophenol	μg/L	45 U	45 U	45 U	45 U
Acenaphthene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
Acenaphthylene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
Anthracene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
Benzo(a)anthracene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
Benzo(a)pyrene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
Benzo(b)fluoranthene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
Benzo(g,h,i)perylene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
Benzo(k)fluoranthene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
Benzoic acid	μg/L	91 UJ	91 U	91 U	91 U
Benzyl alcohol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
bis(2-Chloroethoxy)methane	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
bis(2-Chloroethyl)ether	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 UJ
bis(2-Ethylhexyl)phthalate (DEHP)	μg/L	9.1 UJ	9.1 U	9.1 U	2.2 J
Butyl benzylphthalate (BBP)	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
Chrysene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
Di-n-butylphthalate (DBP)	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
Di-n-octyl phthalate (DnOP)	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
Dibenz(a,h)anthracene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
Dibenzofuran	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
Diethyl phthalate	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
Dimethyl phthalate	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
Fluoranthene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
Fluorene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
Hexachlorobenzene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
Hexachlorobutadiene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
Hexachlorocyclopentadiene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U

	Location ID: Sample Name: Sample Date:	7125 WG-9954-062420-SG-002 06/24/2020	7130 WG-9954-062420-SG-003 06/24/2020	7132 WG-9954-062420-SG-004 06/24/2020	7205 WG-9954-062620-RM-014 06/26/2020
Parameters	Unit				
SVOCs-Continued					
Hexachloroethane	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
Indeno(1,2,3-cd)pyrene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
Isophorone	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
N-Nitrosodi-n-propylamine	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
N-Nitrosodiphenylamine	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
Naphthalene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
Nitrobenzene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
Pentachlorophenol	μg/L	45 U	45 U	45 U	45 U
Phenanthrene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
Phenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Pyrene	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 U
SVOC, TICs					
(4-Methylphenyl)phenylmethanone	μg/L	-	-	-	-
1,4-Dioxane	μg/L	-	-	-	-
1-Chloro-2-methyl-benzene	μg/L	-	-	-	-
1-Chloro-3-methylbenzene	μg/L	-	-	-	-
1-Methyl-2-(phenylmethyl)-benzene	μg/L	-	-	-	-
2,3-Dichlorobenzoic acid	μg/L	-	-	-	-
2,5-Dichlorobenzoic acid	μg/L	-	-	-	-
2-Ethyl-ethanoic acid	μg/L	-	-	-	-
2-Propyl tridecyl-sulfurous acid	μg/L	-	-	-	-
3,4-Dichlorophenol	μg/L	-	-	-	-
3,5-Dichlorophenol	μg/L	-	-	-	-
3-Benzoylbenzoic acid	μg/L	-	-	-	-
3-Chlorobenzoic acid	μg/L	-	-	-	-
3-Methyl-cyclohexanol	μg/L	-	-	-	-
4-Benzoyl-(rel)-benzoic acid	μg/L	-	-	-	-
4-Chlorobenzoic acid	μg/L	-	-	-	-
4-Chlorophenol	μg/L	-	-	-	-
4-Chlorotoluene	μg/L	-	-	-	-
6-Octadecenoic acid	μg/L	-	-	-	-
Bicyclo[2.2.1]heptan-2-one, 1,3,3-trimethyl	μg/L	-	-	-	-
Chlorobenzene	μg/L	-	-	-	-
Cyclohexanol	μg/L	-	-	-	-
Dimethyl disulfide	μg/L	-	-	-	-
Dimethyl tetrasulfide	μg/L	-	-	-	-
Dodecanoic acid	μg/L	-	-	-	-

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	Location ID: Sample Name: Sample Date:	7125 WG-9954-062420-SG-002 06/24/2020	7130 WG-9954-062420-SG-003 06/24/2020	7132 WG-9954-062420-SG-004 06/24/2020	7205 WG-9954-062620-RM-014 06/26/2020
Parameters	Unit				
SVOCs, TICs-Continued					
Hexadecanoic acid	μg/L	-	-	-	-
Homomenthyl salicylate	μg/L	-	-	-	-
Nonanoic acid	μg/L	-	-	-	-
Sulfur	μg/L	-	-	-	14 JN
Toluene	μg/L	-	-	-	-
trans-2-Methylcyclohexanol	μg/L	-	-	-	-
Unknown	μg/L	-	3.7 J	-	29.2 J
Polychlorinated Biphenyls (PCBs)					
Aroclor-1016 (PCB-1016)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1221 (PCB-1221)	μg/L	1.8 U	1.8 U	1.8 U	1.8 U
Aroclor-1232 (PCB-1232)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1242 (PCB-1242)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1248 (PCB-1248)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1254 (PCB-1254)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1260 (PCB-1260)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Pesticides	,,	0.04511	0.045.11	0.04511	0.04511
4,4'-DDD	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
4,4'-DDE	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
4,4'-DDT	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Aldrin	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
alpha-BHC	μg/L	0.045 U	0.045 U	0.045 U	0.088
alpha-Chlordane	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
beta-BHC	μg/L	0.045 U	0.045 U	0.045 U	0.036 J
delta-BHC Dieldrin	μg/L	0.045 U 0.045 U	0.045 U 0.045 U	0.045 U	0.32
	μg/L	0.045 U 0.045 U	0.045 U	0.045 U 0.045 U	0.045 U 0.045 U
Endosulfan I Endosulfan II	μg/L	0.045 U 0.045 U	0.045 U 0.045 U	0.045 U 0.045 U	0.045 U 0.045 U
Endosulfan fil Endosulfan sulfate	μg/L	0.045 U 0.045 U	0.045 U	0.045 U 0.045 U	0.045 U 0.045 U
Endosulian sullate Endrin	μg/L				
	μg/L	0.045 U 0.045 U	0.045 U 0.045 U	0.045 U 0.045 U	0.045 U
Endrin ketone	μg/L	0.045 U 0.045 U	0.045 U 0.045 U	0.045 U 0.045 U	0.045 U 0.11
gamma-BHC (lindane)	μg/L	0.045 U 0.045 U	0.045 U 0.045 U	0.045 U 0.045 U	0.11 0.045 U
gamma-Chlordane	μg/L			0.045 U 0.045 U	
Heptachlor analysida	μg/L	0.045 U	0.045 U		0.045 U
Heptachlor epoxide	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Methoxychlor	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Toxaphene	μg/L	0.46 U	0.46 U	0.46 U	0.46 U

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	Location ID: Sample Name: Sample Date:	7125 WG-9954-062420-SG-002 06/24/2020	7130 WG-9954-062420-SG-003 06/24/2020	7132 WG-9954-062420-SG-004 06/24/2020	7205 WG-9954-062620-RM-014 06/26/2020
Parameters	Unit				
Field Parameters					
Temperature, field	Deg C	10.2	11.1	11.5	11.6
Conductivity, field	mS/cm	1.25	1.04	1.17	2.91
Turbidity, field	NTU	47.2	12.7	4.00	4.78
pH, field	s.u.	7.50	7.58	7.78	7.35

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	Location ID: ample Name: Sample Date:	8106 WG-9954-062620-RM-017 06/26/2020	8115 WG-9954-062420-SG-005 06/24/2020	8125 WG-9954-062420-SG-006 06/24/2020	8210 WG-9954-062620-RM-015 06/26/2020
Parameters	Unit				
Volatile Organic Compounds (VOCs)					
1,1,1-Trichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1,2,2-Tetrachloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1,2-Trichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1-Dichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1-Dichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,2-Dichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,2-Dichloropropane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
2-Butanone (Methyl ethyl ketone) (MEK)	μg/L	10 U	10 U	10 U	10 U
2-Hexanone	μg/L	10 U	10 U	10 U	10 U
4-Methyl-2-pentanone (Methyl isobutyl ketone) (MII	3K) μg/L	10 U	10 U	10 U	10 U
Acetone	μg/L	10 U	10 U	10 U	10 U
Benzene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Bromodichloromethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Bromoform	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Bromomethane (Methyl bromide)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Carbon disulfide	μg/L	10 U	10 U	10 U	2.7 J
Carbon tetrachloride	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chlorobenzene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chloroform (Trichloromethane)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chloromethane (Methyl chloride)	μg/L	0.38 J	5.0 U	5.0 U	0.35 J
cis-1,2-Dichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
cis-1,3-Dichloropropene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Dibromochloromethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Ethylbenzene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Methylene chloride	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Styrene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Tetrachloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Toluene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
trans-1,2-Dichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
trans-1,3-Dichloropropene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Trichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Vinyl acetate	μg/L	10 UJ	10 UJ	10 UJ	10 UJ
Vinyl chloride	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Xylenes (total)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U

Locatio Sample N Sample	ame:	8106 WG-9954-062620-RM-017 06/26/2020	8115 WG-9954-062420-SG-005 06/24/2020	8125 WG-9954-062420-SG-006 06/24/2020	8210 WG-9954-062620-RM-015 06/26/2020
Parameters	Unit				
VOCs, Tentatively Identified Compounds (TICs)					
1,2,4-Trichlorobenzene	μg/L	-	-	-	-
1-Chloro-2-methyl-benzene	μg/L	-	-	-	-
1-Chloro-3-methylbenzene	μg/L	-	-	-	-
2,6-Dichlorotoluene	μg/L	-	-	-	-
2-Methylbutane	μg/L	-	-	-	-
6-Methyl-5-hepten-2-one	μg/L	-	-	-	-
Dimethyl disulfide	μg/L	-	-	-	-
Dimethyl sulfide	μg/L	-	-	-	-
Dimethyl trisulfide	μg/L	-	-	-	-
Hexanal	μg/L	-	-	-	-
Isobutane	μg/L	-	-	-	-
m-Monochlorobenzotrifluoride	μg/L	-	-	-	-
Methanethiol	μg/L	-	-	-	-
Methoxytrimethyl-silane	μg/L	-	-	-	-
Methylthioethane	μg/L	-	-	-	-
Nonanal	μg/L	-	-	-	-
Propane	μg/L	-	-	-	-
Sulfur dioxide (SO2)	μg/L	-	-	-	-
Trimethylfluorosilane	μg/L	-	-	-	10.2 J
Unknown	μg/L	-	-	-	-
Semi-volatile Organic Compounds (SVOCs)					
1,2,4-Trichlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
1,2-Dichlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
1,3-Dichlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
1,4-Dichlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,2'-Oxybis(1-chloropropane) (bis(2-Chloroisopropyl) ether)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4,5-Trichlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4,6-Trichlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4-Dichlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4-Dimethylphenol	μg/L	9.1 U	9.1 U	9.1 UJ	9.1 U
2,4-Dinitrophenol	μg/L	45 U	45 U	45 U	45 U
2,4-Dinitrotoluene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,6-Dinitrotoluene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Chloronaphthalene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Chlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Methylnaphthalene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Methylphenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U

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	Location ID: Sample Name: Sample Date:	8106 WG-9954-062620-RM-017 06/26/2020	8115 WG-9954-062420-SG-005 06/24/2020	8125 WG-9954-062420-SG-006 06/24/2020	8210 WG-9954-062620-RM-015 06/26/2020
Parameters	Unit				
SVOCs-Continued					
2-Nitroaniline	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Nitrophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
3&4-Methylphenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
3,3'-Dichlorobenzidine	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
3-Nitroaniline	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4,6-Dinitro-2-methylphenol	μg/L	45 U	45 U	45 U	45 U
4-Bromophenyl phenyl ether	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Chloro-3-methylphenol	μg/L	9.1 U	9.1 U	9.1 UJ	9.1 U
4-Chloroaniline	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Chlorophenyl phenyl ether	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Nitroaniline	μg/L	9.1 U	9.1 U	9.1 UJ	9.1 U
4-Nitrophenol	μg/L	45 U	45 U	45 U	45 U
Acenaphthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Acenaphthylene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Anthracene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(a)anthracene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(a)pyrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(b)fluoranthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(g,h,i)perylene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(k)fluoranthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzoic acid	μg/L	91 U	91 U	91 U	91 U
Benzyl alcohol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
bis(2-Chloroethoxy)methane	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
bis(2-Chloroethyl)ether	μg/L	9.1 UJ	9.1 U	9.1 U	9.1 UJ
bis(2-Ethylhexyl)phthalate (DEHP)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Butyl benzylphthalate (BBP)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Chrysene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Di-n-butylphthalate (DBP)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Di-n-octyl phthalate (DnOP)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Dibenz(a,h)anthracene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Dibenzofuran	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Diethyl phthalate	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Dimethyl phthalate	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Fluoranthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Fluorene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Hexachlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Hexachlorobutadiene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Hexachlorocyclopentadiene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U

	Location ID: Sample Name: Sample Date:	8106 WG-9954-062620-RM-017 06/26/2020	8115 WG-9954-062420-SG-005 06/24/2020	8125 WG-9954-062420-SG-006 06/24/2020	8210 WG-9954-062620-RM-015 06/26/2020
Parameters	Unit				
SVOCs-Continued					
Hexachloroethane	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Indeno(1,2,3-cd)pyrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Isophorone	μg/L	9.1 U	9.1 U	9.1 UJ	9.1 U
N-Nitrosodi-n-propylamine	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
N-Nitrosodiphenylamine	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Naphthalene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Nitrobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Pentachlorophenol	μg/L	45 U	45 U	45 U	45 U
Phenanthrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Phenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Pyrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
SVOC, TICs					
(4-Methylphenyl)phenylmethanone	μg/L	-	-	-	-
1,4-Dioxane	μg/L	-	-	-	-
1-Chloro-2-methyl-benzene	μg/L	-	-	-	-
1-Chloro-3-methylbenzene	μg/L	-	-	-	-
1-Methyl-2-(phenylmethyl)-benzene	μg/L	-	-	-	-
2,3-Dichlorobenzoic acid	μg/L	-	-	-	-
2,5-Dichlorobenzoic acid	μg/L	-	-	-	-
2-Ethyl-ethanoic acid	μg/L	-	-	-	-
2-Propyl tridecyl-sulfurous acid	μg/L	-	-	-	-
3,4-Dichlorophenol	μg/L	-	-	-	-
3,5-Dichlorophenol	μg/L	-	-	-	-
3-Benzoylbenzoic acid	μg/L	-	-	-	-
3-Chlorobenzoic acid	μg/L	-	-	-	-
3-Methyl-cyclohexanol	μg/L	-	-	-	-
4-Benzoyl-(rel)-benzoic acid	μg/L	-	-	-	-
4-Chlorobenzoic acid	μg/L	-	-	-	-
4-Chlorophenol	μg/L	-	-	-	-
4-Chlorotoluene	μg/L	-	-	-	-
6-Octadecenoic acid	μg/L 	-	-	-	-
Bicyclo[2.2.1]heptan-2-one, 1,3,3-trimethyl	μg/L 	-	-	-	-
Chlorobenzene	μg/L 	-	-	-	-
Cyclohexanol	μg/L 	-	-	-	-
Dimethyl disulfide	μg/L	-	-	-	-
Dimethyl tetrasulfide	μg/L	-	-	-	-
Dodecanoic acid	μg/L	-	-	-	-

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Parameters Unit
Hexadecanoic acid
Homomenthyl salicylate
Nonancic acid μg/L - - - - - - - - -
Sulfur Foliable
trans-2-Methylcyclohexanol Unknown
Polychlorinated Biphenyls (PCBs) Color-1016 (PCB-1016) μg/L 0.91 U
Polychlorinated Biphenyls (PCBs) Aroclor-1016 (PCB-1016) μg/L 0.91 U
Aroclor-1016 (PCB-1016) μg/L 0.91 U
Aroclor-1016 (PCB-1016) μg/L 0.91 U
Aroclor-1221 (PCB-1221) μg/L 1.8 U 1.8
Aroclor-1232 (PCB-1232) μg/L 0.91 U <
Aroclor-1242 (PCB-1242) μg/L 0.91 U 0.94 U 0.91 U 0.94 U 0.94 U
Aroclor-1248 (PCB-1248) μg/L 0.91 U 0.945 U 0.945 U 0.945 U 0.945 U 0.945 U 0.945 U 0.945 U 0.945 U 0.945 U
Aroclor-1254 (PCB-1254) μg/L 0.91 U 0.945 U 0.045 U
Aroclor-1260 (PCB-1260) μg/L 0.91 U 0.91 U 0.91 U 0.91 U Pesticides 4,4*-DDD μg/L 0.045 U 0.045 U 0.045 U 0.045 U 4,4*-DDE μg/L 0.045 U 0.045 U 0.045 U 0.045 U 4,4*-DDT μg/L 0.045 U 0.045 U 0.045 U 0.045 U Aldrin μg/L 0.045 U 0.045 U 0.045 U 0.045 U alpha-BHC μg/L 0.045 U 0.045 U 0.045 U 0.045 U alpha-Chlordane μg/L 0.045 U 0.045 U 0.045 U 0.045 U beta-BHC μg/L 0.045 U 0.045 U 0.045 U 0.045 U delta-BHC μg/L 0.045 U 0.045 U 0.045 U 0.045 U
4,4*-DDD µg/L 0.045 U
4,4*-DDD µg/L 0.045 U
4,4'-DDE µg/L 0.045 U
4,4'-DDT µg/L 0.045 U 0.064 0.064 0.045 U <
Aldrin µg/L 0.045 U 0.042 J
alpha-BHC µg/L 0.045 U 0.045 U 0.045 U 0.064 alpha-Chlordane µg/L 0.045 U 0.045 U 0.045 U 0.045 U beta-BHC µg/L 0.045 U 0.045 U 0.045 U 0.045 U 0.045 U delta-BHC µg/L 0.045 U 0.045 U 0.045 U 0.045 U 0.042 J
alpha-Chlordane µg/L 0.045 U 0.045 U </td
beta-BHC μg/L 0.045 U 0.042 J
delta-BHC
Dieldrin μg/L 0.045 U 0.045 U 0.045 U 0.045 U 0.045 U
Endosulfan I μg/L 0.045 U 0.045 U 0.045 U 0.045 U
Endosulfan II μg/L 0.045 U 0.045 U 0.045 U 0.045 U
Endosulfan sulfate µg/L 0.045 U 0.045 U 0.045 U 0.045 U
Endrin μg/L 0.045 U 0.045 U 0.045 U 0.045 U
Endrin ketone μg/L 0.045 U 0.045 U 0.045 U 0.045 U
gamma-BHC (lindane) µg/L 0.045 U 0.045 U 0.045 U 0.060
gamma-Chlordane µg/L 0.045 U 0.045 U 0.045 U 0.045 U
Heptachlor µg/L 0.045 U 0.045 U 0.045 U 0.045 U
Heptachlor epoxide μg/L 0.045 U 0.045 U 0.045 U 0.045 U
Methoxychlor μg/L 0.045 U 0.045 U 0.045 U 0.045 U
Toxaphene µg/L 0.46 U 0.46 U 0.46 U 0.46 U

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	Location ID: Sample Name: Sample Date:	8106 WG-9954-062620-RM-017 06/26/2020	8115 WG-9954-062420-SG-005 06/24/2020	8125 WG-9954-062420-SG-006 06/24/2020	8210 WG-9954-062620-RM-015 06/26/2020
Parameters	Unit				
Field Parameters					
Temperature, field	Deg C	11.7	10.6	10.4	12.3
Conductivity, field	mS/cm	2.28	1.44	1.62	3.92
Turbidity, field	NTU	206	456	335	2.05
pH, field	s.u.	7.20	7.49	7.33	7.23

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	Location ID: Sample Name: Sample Date:	9105 WG-9954-062420-SG-007 06/24/2020	9113 WG-9954-062520-SG-008 06/25/2020	9118 WG-9954-062520-SG-009 06/25/2020	9118 WG-9954-062520-SG-010 06/25/2020 Duplicate
Parameters	Unit				
Volatile Organic Compounds (VOCs)					
1,1,1-Trichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1,2,2-Tetrachloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1,2-Trichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1-Dichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1-Dichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,2-Dichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,2-Dichloropropane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
2-Butanone (Methyl ethyl ketone) (MEK)	μg/L	10 U	10 U	10 U	10 U
2-Hexanone	μg/L	10 U	10 U	10 U	10 U
4-Methyl-2-pentanone (Methyl isobutyl ketone) (M	IBK) μg/L	10 U	0.55 J	10 U	10 U
Acetone	μg/L	10 U	10 U	10 U	10 U
Benzene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Bromodichloromethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Bromoform	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Bromomethane (Methyl bromide)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Carbon disulfide	μg/L	10 U	10 U	0.48 J	0.75 J
Carbon tetrachloride	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chlorobenzene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chloroform (Trichloromethane)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chloromethane (Methyl chloride)	μg/L	5.0 U	0.32 J	0.38 J	5.0 U
cis-1,2-Dichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
cis-1,3-Dichloropropene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Dibromochloromethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Ethylbenzene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Methylene chloride	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Styrene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Tetrachloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Toluene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
trans-1,2-Dichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
trans-1,3-Dichloropropene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Trichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Vinyl acetate	μg/L	10 UJ	10 UJ	10 U	10 U
Vinyl chloride	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Xylenes (total)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U

Location Sample Nar Sample Da	ne:	9105 WG-9954-062420-SG-007 06/24/2020	9113 WG-9954-062520-SG-008 06/25/2020	9118 WG-9954-062520-SG-009 06/25/2020	9118 WG-9954-062520-SG-010 06/25/2020 Duplicate
Parameters	Unit				
VOCs, Tentatively Identified Compounds (TICs)					
1,2,4-Trichlorobenzene	μg/L	-	-	-	-
1-Chloro-2-methyl-benzene	μg/L	-	-	-	-
1-Chloro-3-methylbenzene	μg/L	-	-	-	-
2,6-Dichlorotoluene	μg/L	-	-	-	-
2-Methylbutane	μg/L	-	-	-	-
6-Methyl-5-hepten-2-one	μg/L	-	-	-	-
Dimethyl disulfide	μg/L	-	-	-	-
Dimethyl sulfide	μg/L	-	-	-	-
Dimethyl trisulfide	μg/L	-	-	-	-
Hexanal	μg/L	-	-	-	-
Isobutane	μg/L	-	-	-	-
m-Monochlorobenzotrifluoride	μg/L	-	-	-	-
Methanethiol	μg/L	-	-	-	-
Methoxytrimethyl-silane	μg/L	-	-	-	-
Methylthioethane	μg/L	-	-	-	-
Nonanal	μg/L	-	-	-	-
Propane	μg/L	-	-	-	<u>-</u>
Sulfur dioxide (SO2)	μg/L	-	-	-	28.9 JN
Trimethylfluorosilane	μg/L	-	-	-	-
Unknown	μg/L	-	-	-	-
Semi-volatile Organic Compounds (SVOCs)					
1,2,4-Trichlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
1,2-Dichlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
1,3-Dichlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
1,4-Dichlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,2'-Oxybis(1-chloropropane) (bis(2-Chloroisopropyl) ether)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4,5-Trichlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4,6-Trichlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4-Dichlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4-Dimethylphenol	μg/L	9.1 UJ	9.1 UJ	9.1 UJ	9.1 UJ
2,4-Dinitrophenol	μg/L	45 U	45 U	45 U	45 U
2,4-Dinitrotoluene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,6-Dinitrotoluene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Chloronaphthalene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Chlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Methylnaphthalene	μg/L	9.1 U	9.1 U	9.1 U 9.1 U	9.1 U
2-Methylphenol	μg/L	9.1 U	9.1 U	9.10	9.1 U

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	Location ID: Sample Name: Sample Date:	9105 WG-9954-062420-SG-007 06/24/2020	9113 WG-9954-062520-SG-008 06/25/2020	9118 WG-9954-062520-SG-009 06/25/2020	9118 WG-9954-062520-SG-010 06/25/2020 Duplicate
Parameters	Unit				
SVOCs-Continued					
2-Nitroaniline	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Nitrophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
3&4-Methylphenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
3,3'-Dichlorobenzidine	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
3-Nitroaniline	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4,6-Dinitro-2-methylphenol	μg/L	45 U	45 U	45 U	45 U
4-Bromophenyl phenyl ether	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Chloro-3-methylphenol	μg/L	9.1 UJ	9.1 UJ	9.1 UJ	9.1 UJ
4-Chloroaniline	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Chlorophenyl phenyl ether	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Nitroaniline	μg/L	9.1 UJ	9.1 UJ	9.1 UJ	9.1 UJ
4-Nitrophenol	μg/L	45 U	45 U	45 U	45 U
Acenaphthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Acenaphthylene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Anthracene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(a)anthracene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(a)pyrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(b)fluoranthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(g,h,i)perylene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(k)fluoranthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzoic acid	μg/L	91 U	91 U	91 U	91 U
Benzyl alcohol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
bis(2-Chloroethoxy)methane	μg/L 	9.1 U	9.1 U	9.1 U	9.1 U
bis(2-Chloroethyl)ether	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
bis(2-Ethylhexyl)phthalate (DEHP)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Butyl benzylphthalate (BBP)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Chrysene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Di-n-butylphthalate (DBP)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Di-n-octyl phthalate (DnOP)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Dibenz(a,h)anthracene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Dibenzofuran	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Diethyl phthalate	μg/L	9.1 U 9.1 U	9.1 U 9.1 U	9.1 U 9.1 U	9.1 U 9.1 U
Dimethyl phthalate	μg/L				
Fluoranthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Fluorene	μg/L	9.1 U	9.1 U 9.1 U	9.1 U	9.1 U
Hexachlorobenzene	μg/L	9.1 U		9.1 U	9.1 U
Hexachloroputadiene	μg/L	9.1 U 9.1 U	9.1 U 9.1 U	9.1 U 9.1 U	9.1 U 9.1 U
Hexachlorocyclopentadiene	μg/L	9.10	9.10	9.10	9.10

Sample	tion ID: Name: e Date:	9105 WG-9954-062420-SG-007 06/24/2020	9113 WG-9954-062520-SG-008 06/25/2020	9118 WG-9954-062520-SG-009 06/25/2020	9118 WG-9954-062520-SG-010 06/25/2020 Duplicate
Parameters	Unit				
SVOCs-Continued					
Hexachloroethane	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Indeno(1,2,3-cd)pyrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Isophorone	μg/L	9.1 UJ	9.1 UJ	9.1 UJ	9.1 UJ
N-Nitrosodi-n-propylamine	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
N-Nitrosodiphenylamine	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Naphthalene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Nitrobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Pentachlorophenol	μg/L	45 U	45 U	45 U	45 U
Phenanthrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Phenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Pyrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
SVOC, TICs					
(4-Methylphenyl)phenylmethanone	μg/L	-	-	-	-
1,4-Dioxane	μg/L	-	-	-	-
1-Chloro-2-methyl-benzene	μg/L	-	-	-	-
1-Chloro-3-methylbenzene	μg/L	-	-	-	-
1-Methyl-2-(phenylmethyl)-benzene	μg/L	-	-	-	-
2,3-Dichlorobenzoic acid	μg/L	-	-	-	-
2,5-Dichlorobenzoic acid	μg/L	-	-	-	-
2-Ethyl-ethanoic acid	μg/L	-	-	-	-
2-Propyl tridecyl-sulfurous acid	μg/L	-	-	-	-
3,4-Dichlorophenol	μg/L	-	-	-	-
3,5-Dichlorophenol	μg/L	-	-	-	-
3-Benzoylbenzoic acid	μg/L	-	-	-	-
3-Chlorobenzoic acid	μg/L	-	-	-	-
3-Methyl-cyclohexanol	μg/L	-	-	-	-
4-Benzoyl-(rel)-benzoic acid	μg/L	-	-	-	-
4-Chlorobenzoic acid	μg/L	-	-	-	-
4-Chlorophenol	μg/L	-	-	-	-
4-Chlorotoluene	μg/L	-	-	-	-
6-Octadecenoic acid	μg/L	-	11 JN	-	-
Bicyclo[2.2.1]heptan-2-one, 1,3,3-trimethyl	μg/L	-	-	-	-
Chlorobenzene	μg/L	-	-	-	-
Cyclohexanol	μg/L	-	-	-	-
Dimethyl disulfide	μg/L	-	-	-	-
Dimethyl tetrasulfide	μg/L	-	-	-	-
Dodecanoic acid	μg/L	-	-	-	-

	Location ID: Sample Name: Sample Date:	9105 WG-9954-062420-SG-007 06/24/2020	9113 WG-9954-062520-SG-008 06/25/2020	9118 WG-9954-062520-SG-009 06/25/2020	9118 WG-9954-062520-SG-010 06/25/2020 Duplicate
Parameters	Unit				
SVOCs, TICs-Continued					
Hexadecanoic acid	μg/L	-	5.8 JN	-	-
Homomenthyl salicylate	μg/L	-	-	-	-
Nonanoic acid	μg/L	-	-	-	-
Sulfur	μg/L	-	-	-	-
Toluene	μg/L	-	-	-	-
trans-2-Methylcyclohexanol	μg/L	-	-	-	-
Unknown	μg/L	83.2 J	4.6 J	59.2 J	-
Polychlorinated Biphenyls (PCBs)					
Aroclor-1016 (PCB-1016)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1221 (PCB-1221)	μg/L	1.8 U	1.8 U	1.8 U	1.8 U
Aroclor-1232 (PCB-1232)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1242 (PCB-1242)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1248 (PCB-1248)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1254 (PCB-1254)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1260 (PCB-1260)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Pesticides					
4,4'-DDD	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
4,4'-DDE	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
4,4'-DDT	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Aldrin	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
alpha-BHC	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
alpha-Chlordane	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
beta-BHC	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
delta-BHC	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Dieldrin	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Endosulfan I	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Endosulfan II	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Endosulfan sulfate	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Endrin	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Endrin ketone	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
gamma-BHC (lindane)	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
gamma-Chlordane	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Heptachlor	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Heptachlor epoxide	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Methoxychlor	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Toxaphene	μg/L	0.46 U	0.46 U	0.46 U	0.46 U

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	Location ID: Sample Name: Sample Date:	9105 WG-9954-062420-SG-007 06/24/2020	9113 WG-9954-062520-SG-008 06/25/2020	9118 WG-9954-062520-SG-009 06/25/2020	9118 WG-9954-062520-SG-010 06/25/2020 Duplicate
Parameters	Unit				
Field Parameters					
Temperature, field	Deg C	11.0	11.1	11.3	11.3
Conductivity, field	mS/cm	1.72	0.79	0.96	0.96
Turbidity, field	NTU	1000	79.8	202	202
pH, field	s.u.	7.48	10.46	7.47	7.47

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	Location ID: ample Name: Sample Date:	9205 WG-9954-062620-RM-016 06/26/2020	9210 WG-9954-062520-SG-012 06/25/2020	10135 WG-9954-070120-SG-027 07/01/2020	10135 WG-9954-070120-SG-028 07/01/2020 Duplicate
Parameters	Unit				
Volatile Organic Compounds (VOCs)					
1,1,1-Trichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1,2,2-Tetrachloroethane	μg/L	5.0 U	5.0 U	6.1	7.9
1,1,2-Trichloroethane	μg/L	5.0 U	5.0 U	12	14
1,1-Dichloroethane	μg/L	5.0 U	5.0 U	5.0 U	0.27 J
1,1-Dichloroethene	μg/L	5.0 U	5.0 U	0.63 J	0.78 J
1,2-Dichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,2-Dichloropropane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
2-Butanone (Methyl ethyl ketone) (MEK)	μg/L	10 U	10 U	12	6.8 J
2-Hexanone	μg/L	10 U	10 U	10 U	2.1 J
4-Methyl-2-pentanone (Methyl isobutyl ketone) (MIE		10 U	10 U	10 U	10 U
Acetone	μg/L	10 U	10 U	79	71
Benzene	μg/L	5.0 U	5.0 U	6500	6300
Bromodichloromethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Bromoform	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Bromomethane (Methyl bromide)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Carbon disulfide	μg/L	1.9 J	2.6 J	10 U	10 U
Carbon tetrachloride	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chlorobenzene	μg/L	5.0 U	5.0 U	2500	2400
Chloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chloroform (Trichloromethane)	μg/L	5.0 U	5.0 U	140	140
Chloromethane (Methyl chloride)	μg/L	5.0 U	0.30 J	5.0 U	5.0 U
cis-1,2-Dichloroethene	μg/L	5.0 U	5.0 U	26	28
cis-1,3-Dichloropropene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Dibromochloromethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Ethylbenzene	μg/L	5.0 U	5.0 U	12	12
Methylene chloride	μg/L	5.0 U	5.0 U	6.4	7.2
Styrene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Tetrachloroethene	μg/L	5.0 U	5.0 U	20	23
Toluene	μg/L	5.0 U	5.0 U	22000	21000
trans-1,2-Dichloroethene	μg/L	5.0 U	5.0 U	29	35
trans-1,3-Dichloropropene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Trichloroethene	μg/L	5.0 U	5.0 U	120	130
Vinyl acetate	μg/L	10 U	10 U	10 U	10 U
Vinyl chloride	μg/L	5.0 U	5.0 U	10	11
Xylenes (total)	μg/L	5.0 U	5.0 U	58	58

Location ID Sample Name Sample Date	•	9205 WG-9954-062620-RM-016 06/26/2020	9210 WG-9954-062520-SG-012 06/25/2020	10135 WG-9954-070120-SG-027 07/01/2020	10135 WG-9954-070120-SG-028 07/01/2020 Duplicate
Parameters	Unit				
VOCs, Tentatively Identified Compounds (TICs)					
1,2,4-Trichlorobenzene	μg/L	-	-	-	-
1-Chloro-2-methyl-benzene	μg/L	-	-	16362.0 J	-
1-Chloro-3-methylbenzene	μg/L	-	-	8134.0 J	-
2,6-Dichlorotoluene	μg/L	-	-	1534.0 J	-
2-Methylbutane	μg/L	-	-	-	-
6-Methyl-5-hepten-2-one	μg/L	-	-	-	-
Dimethyl disulfide	μg/L	-	-	-	-
Dimethyl sulfide	μg/L	-	-	-	-
Dimethyl trisulfide	μg/L	-	-	-	-
Hexanal	μg/L	-	-	-	-
Isobutane	μg/L	-	-	-	-
m-Monochlorobenzotrifluoride	μg/L	-	-	-	-
Methanethiol	μg/L	-	-	-	-
Methoxytrimethyl-silane	μg/L	-	-	-	-
Methylthioethane	μg/L	-	-	-	-
Nonanal	μg/L	-	-	-	-
Propane	μg/L	-	-	-	-
Sulfur dioxide (SO2)	μg/L	71.1 J	-	-	-
Trimethylfluorosilane	μg/L	-	-	-	-
Unknown	μg/L	-	-	-	-
Semi-volatile Organic Compounds (SVOCs)	_				
1,2,4-Trichlorobenzene	μg/L	9.1 U	9.1 U	93	100
1,2-Dichlorobenzene	μg/L	9.1 U	9.1 U	25	34
1,3-Dichlorobenzene	μg/L	9.1 U	9.1 U	3.6 J	4.1 J
1,4-Dichlorobenzene	μg/L	9.1 U	9.1 U	73	97
2,2'-Oxybis(1-chloropropane) (bis(2-Chloroisopropyl) ether)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4,5-Trichlorophenol	μg/L	9.1 U	9.1 U	37	26
2,4,6-Trichlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4-Dichlorophenol	μg/L	9.1 U	9.1 U	330	350
2,4-Dimethylphenol	μg/L	9.1 U	9.1 UJ	8.9 J	6.9 J
2,4-Dinitrophenol	μg/L	45 U	45 U	45 U	45 U
2,4-Dinitrotoluene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,6-Dinitrotoluene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Chloronaphthalene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Chlorophenol	μg/L	9.1 U	9.1 U	27	31
2-Methylnaphthalene	μg/L	9.1 U	9.1 U	1.4 J	9.1 U
2-Methylphenol	μg/L	9.1 U	9.1 U	32	32

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	Location ID: Sample Name: Sample Date:	9205 WG-9954-062620-RM-016 06/26/2020	9210 WG-9954-062520-SG-012 06/25/2020	10135 WG-9954-070120-SG-027 07/01/2020	10135 WG-9954-070120-SG-028 07/01/2020 Duplicate
Parameters	Unit				
SVOCs-Continued					
2-Nitroaniline	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Nitrophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
3&4-Methylphenol	μg/L	9.1 U	9.1 U	71	71
3,3'-Dichlorobenzidine	μg/L	9.1 U	9.1 U	R	9.1 U
3-Nitroaniline	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4,6-Dinitro-2-methylphenol	μg/L	45 U	45 U	45 U	45 U
4-Bromophenyl phenyl ether	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Chloro-3-methylphenol	μg/L	9.1 U	9.1 UJ	46	40
4-Chloroaniline	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Chlorophenyl phenyl ether	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Nitroaniline	μg/L	9.1 U	9.1 UJ	9.1 U	9.1 U
4-Nitrophenol	μg/L	45 U	45 U	45 U	45 U
Acenaphthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Acenaphthylene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Anthracene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(a)anthracene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(a)pyrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(b)fluoranthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(g,h,i)perylene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(k)fluoranthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzoic acid	μg/L	91 U	91 U	11000	10000
Benzyl alcohol	μg/L	9.1 U	9.1 U	290	290
bis(2-Chloroethoxy)methane	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
bis(2-Chloroethyl)ether	μg/L	9.1 UJ	9.1 U	20	24
bis(2-Ethylhexyl)phthalate (DEHP)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Butyl benzylphthalate (BBP)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Chrysene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Di-n-butylphthalate (DBP)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Di-n-octyl phthalate (DnOP)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Dibenz(a,h)anthracene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Dibenzofuran	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Diethyl phthalate	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Dimethyl phthalate	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Fluoranthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Fluorene	μ <mark>g</mark> /L	9.1 U	9.1 U	9.1 U	9.1 U
Hexachlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Hexachlorobutadiene	μg/L	9.1 U	9.1 U	9.1 U	1.0 J
Hexachlorocyclopentadiene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U

	Location ID: Sample Name: Sample Date:	9205 WG-9954-062620-RM-016 06/26/2020	9210 WG-9954-062520-SG-012 06/25/2020	10135 WG-9954-070120-SG-027 07/01/2020	10135 WG-9954-070120-SG-028 07/01/2020 Duplicate
Parameters	Unit				
SVOCs-Continued					
Hexachloroethane	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Indeno(1,2,3-cd)pyrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Isophorone	μg/L	9.1 U	9.1 UJ	R	9.1 U
N-Nitrosodi-n-propylamine	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
N-Nitrosodiphenylamine	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Naphthalene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Nitrobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Pentachlorophenol	μg/L	45 U	45 U	45 U	45 U
Phenanthrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Phenol	μg/L	9.1 U	9.1 U	40	44
Pyrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
SVOC, TICs					
(4-Methylphenyl)phenylmethanone	μg/L	-	-	31 JN	27 JN
1,4-Dioxane	μg/L	-	-	140 JN	140 JN
1-Chloro-2-methyl-benzene	μg/L	-	-	1800 JN	2000 JN
1-Chloro-3-methylbenzene	μg/L	-	-	-	-
1-Methyl-2-(phenylmethyl)-benzene	μg/L	-	-	-	17 JN
2,3-Dichlorobenzoic acid	μg/L	-	-	28 JN	-
2,5-Dichlorobenzoic acid	μg/L	-	-	-	30 JN
2-Ethyl-ethanoic acid	μg/L	-	-	110 JN	100 JN
2-Propyl tridecyl-sulfurous acid	μg/L	-	-	-	-
3,4-Dichlorophenol	μg/L	-	-	-	260 JN
3,5-Dichlorophenol	μg/L	-	-	370 JN	-
3-Benzoylbenzoic acid	μg/L	-	-	290 JN	170 JN
3-Chlorobenzoic acid	μg/L	-	-	9900 JN	180 JN
3-Methyl-cyclohexanol	μg/L	-	-	38 JN	29 JN
4-Benzoyl-(rel)-benzoic acid	μg/L	-	-	780 JN	630 JN
4-Chlorobenzoic acid	μg/L	-	-	1500 JN	6200 JN
4-Chlorophenol	μg/L	-	-	150 JN	150 JN
4-Chlorotoluene	μg/L	-	-	1000 JN	1200 JN
6-Octadecenoic acid	μg/L	-	-	-	-
Bicyclo[2.2.1]heptan-2-one, 1,3,3-trimethyl	μg/L	-	-	32 JN	32 JN
Chlorobenzene	μg/L	-	-	350 JN	400 JN
Cyclohexanol	μg/L	-	-	120 JN	130 JN
Dimethyl disulfide	μg/L	-	-	-	-
Dimethyl tetrasulfide	μg/L	-	-	-	-
Dodecanoic acid	μg/L	-	-	-	-

	Location ID: Sample Name: Sample Date:	9205 WG-9954-062620-RM-016 06/26/2020	9210 WG-9954-062520-SG-012 06/25/2020	10135 WG-9954-070120-SG-027 07/01/2020	10135 WG-9954-070120-SG-028 07/01/2020 Duplicate
Parameters	Unit				
SVOCs, TICs-Continued					
Hexadecanoic acid	μg/L	-	-	-	-
Homomenthyl salicylate	μg/L	-	-	-	-
Nonanoic acid	μg/L	-	-	-	-
Sulfur	μg/L	4.0 JN	9.6 JN	-	-
Toluene	μg/L	-	-	1800 JN	-
trans-2-Methylcyclohexanol	μg/L	-	-	-	18 JN
Unknown	μg/L	86.4 J	-	552 J	101 J
Polychlorinated Biphenyls (PCBs)					
Aroclor-1016 (PCB-1016)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1221 (PCB-1221)	μg/L	1.8 U	1.8 U	1.8 U	1.8 U
Aroclor-1232 (PCB-1232)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1242 (PCB-1242)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1248 (PCB-1248)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1254 (PCB-1254)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1260 (PCB-1260)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Pesticides					
4,4'-DDD	μg/L	0.045 U	0.045 U	0.45 U	0.45 U
4,4'-DDE	μg/L	0.045 U	0.045 U	0.45 U	0.45 U
4,4'-DDT	μg/L	0.045 U	0.045 U	0.45 U	0.45 U
Aldrin	μg/L	0.045 U	0.045 U	1.0 J	0.85
alpha-BHC	μg/L	0.097	0.033 J	27	25
alpha-Chlordane	μg/L	0.045 U	0.045 U	0.45 U	0.45 U
beta-BHC	μg/L	0.045 U	0.045 U	6.8 J	6.9
delta-BHC	μg/L	0.10	0.068	9.2	8.7
Dieldrin	μg/L	0.045 U	0.045 U	0.45 U	0.45 U
Endosulfan I	μg/L	0.045 U	0.045 U	0.45 U	0.45 U
Endosulfan II	μg/L	0.045 U	0.045 U	0.45 U	0.45 U
Endosulfan sulfate	μg/L	0.045 U	0.045 U	0.45 U	0.45 U
Endrin	μg/L	0.045 U	0.045 U	0.45 U	0.45 U
Endrin ketone	μg/L	0.045 U	0.045 U	0.45 U	0.45 U
gamma-BHC (lindane)	μg/L	0.096	0.037 J	6.5	6.2
gamma-Chlordane	μg/L	0.045 U	0.028 J	0.45 U	0.45 U
Heptachlor	μg/L	0.045 U	0.045 U	0.45 U	0.45 U
Heptachlor epoxide	μg/L	0.045 U	0.045 U	0.45 U	0.45 U
Methoxychlor	μg/L	0.045 U	0.045 U	0.45 U	0.45 U
Toxaphene	μg/L	0.46 U	0.46 U	4.6 U	4.6 U

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	Location ID: Sample Name: Sample Date:	9205 WG-9954-062620-RM-016 06/26/2020	9210 WG-9954-062520-SG-012 06/25/2020	10135 WG-9954-070120-SG-027 07/01/2020	10135 WG-9954-070120-SG-028 07/01/2020 Duplicate
Parameters	Unit				
Field Parameters					
Temperature, field	Deg C	13.1	13.8	11.3	11.3
Conductivity, field	mS/cm	2.34	5.51	2.97	2.97
Turbidity, field	NTU	14.9	3.69	869	869
pH, field	s.u.	7.32	7.02	6.98	6.98

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	Location ID: ample Name: Sample Date:	10178A WG-9954-070720-SG-031 07/07/2020	10178B WG-9954-071020-SG-036 07/10/2020	10205 WG-9954-062520-SG-013 06/25/2020	10210A WG-9954-070920-SG-033 07/09/2020
Parameters	Unit				
Volatile Organic Compounds (VOCs)					
1,1,1-Trichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1,2,2-Tetrachloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1,2-Trichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1-Dichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1-Dichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,2-Dichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,2-Dichloropropane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
2-Butanone (Methyl ethyl ketone) (MEK)	μg/L	10 U	10 U	10 U	10 U
2-Hexanone	μg/L	10 U	10 U	10 U	10 U
4-Methyl-2-pentanone (Methyl isobutyl ketone) (MIE	βK) μg/L	10 U	10 U	10 U	10 U
Acetone	μg/L	10 U	10 U	10 U	10 U
Benzene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Bromodichloromethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Bromoform	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Bromomethane (Methyl bromide)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Carbon disulfide	μg/L	10 U	10 U	5.4 J	9.2 J
Carbon tetrachloride	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chlorobenzene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chloroform (Trichloromethane)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chloromethane (Methyl chloride)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
cis-1,2-Dichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
cis-1,3-Dichloropropene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Dibromochloromethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Ethylbenzene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Methylene chloride	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Styrene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Tetrachloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Toluene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
trans-1,2-Dichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
trans-1,3-Dichloropropene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Trichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Vinyl acetate	μg/L	10 U	10 U	10 U	10 U
Vinyl chloride	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Xylenes (total)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U

Location ID Sample Name Sample Date		10178A WG-9954-070720-SG-031 07/07/2020	10178B WG-9954-071020-SG-036 07/10/2020	10205 WG-9954-062520-SG-013 06/25/2020	10210A WG-9954-070920-SG-033 07/09/2020
Parameters	Unit				
VOCs, Tentatively Identified Compounds (TICs)					
1,2,4-Trichlorobenzene	μg/L	-	-	-	-
1-Chloro-2-methyl-benzene	μg/L	-	-	-	-
1-Chloro-3-methylbenzene	μg/L	-	-	-	-
2,6-Dichlorotoluene	μg/L	-	-	-	-
2-Methylbutane	μg/L	-	-	-	5.4 J
6-Methyl-5-hepten-2-one	μg/L	-	-	-	-
Dimethyl disulfide	μg/L	-	-	-	33.4 J
Dimethyl sulfide	μg/L	-	-	-	325.6 J
Dimethyl trisulfide	μg/L	-	-	-	9.9 J
Hexanal	μg/L	-	-	-	10.3 J
Isobutane	μg/L	-	-	-	11.4 J
m-Monochlorobenzotrifluoride	μg/L	-	-	-	-
Methanethiol	μg/L	-	-	-	5.8 J
Methoxytrimethyl-silane	μg/L	-	-	-	-
Methylthioethane	μg/L	-	-	-	25.8 J
Nonanal	μg/L	-	-	-	7.0 J
Propane	μg/L	-	-	-	7.7 J
Sulfur dioxide (SO2)	μg/L	-	-	-	-
Trimethylfluorosilane	μg/L	-	-	-	-
Unknown	μg/L	-	-	-	-
Semi-volatile Organic Compounds (SVOCs)					
1,2,4-Trichlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
1,2-Dichlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
1,3-Dichlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
1,4-Dichlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,2'-Oxybis(1-chloropropane) (bis(2-Chloroisopropyl) ether)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4,5-Trichlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4,6-Trichlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4-Dichlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4-Dimethylphenol	μg/L	9.1 U	9.1 U	9.1 UJ	9.1 U
2,4-Dinitrophenol	μg/L	45 U	45 U	45 U	45 U
2,4-Dinitrotoluene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,6-Dinitrotoluene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Chloronaphthalene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Chlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Methylnaphthalene	μg/L	9.1 U 9.1 U	9.1 U	9.1 U	9.1 U
2-Methylphenol	μg/L	9.10	9.1 U	9.1 U	9.1 U

	Location ID: Sample Name: Sample Date:	10178A WG-9954-070720-SG-031 07/07/2020	10178B WG-9954-071020-SG-036 07/10/2020	10205 WG-9954-062520-SG-013 06/25/2020	10210A WG-9954-070920-SG-033 07/09/2020
Parameters	Unit				
SVOCs-Continued					
2-Nitroaniline	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Nitrophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
3&4-Methylphenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
3,3'-Dichlorobenzidine	μ <u>σ</u> /L	9.1 U	9.1 U	9.1 U	9.1 U
3-Nitroaniline	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4,6-Dinitro-2-methylphenol	μg/L	45 U	45 U	45 U	45 U
4-Bromophenyl phenyl ether	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Chloro-3-methylphenol	μg/L	9.1 U	9.1 U	9.1 UJ	9.1 U
4-Chloroaniline	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Chlorophenyl phenyl ether	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Nitroaniline	μg/L	9.1 U	9.1 U	9.1 UJ	9.1 U
4-Nitrophenol	μg/L	45 U	45 U	45 U	45 U
Acenaphthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Acenaphthylene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Anthracene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(a)anthracene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(a)pyrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(b)fluoranthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(g,h,i)perylene	μ <mark>g</mark> /L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(k)fluoranthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzoic acid	μg/L	91 U	91 U	91 U	91 U
Benzyl alcohol	μ <mark>g</mark> /L	9.1 U	9.1 U	9.1 U	9.1 U
bis(2-Chloroethoxy)methane	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
bis(2-Chloroethyl)ether	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
bis(2-Ethylhexyl)phthalate (DEHP)	μg/L	9.1 U	1.3 J	3.7 J	3.9 J
Butyl benzylphthalate (BBP)	μ <u>σ</u> /L	9.1 U	9.1 U	9.1 U	9.1 U
Chrysene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Di-n-butylphthalate (DBP)	μg/L	9.1 U	9.1 U	9.1 U	2.3 J
Di-n-octyl phthalate (DnOP)	μ <u>σ</u> /L	9.1 U	9.1 U	9.1 U	9.1 U
Dibenz(a,h)anthracene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Dibenzofuran	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Diethyl phthalate	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Dimethyl phthalate	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Fluoranthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Fluorene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Hexachlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Hexachlorobutadiene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Hexachlorocyclopentadiene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U

	Location ID: Sample Name: Sample Date:	10178A WG-9954-070720-SG-031 07/07/2020	10178B WG-9954-071020-SG-036 07/10/2020	10205 WG-9954-062520-SG-013 06/25/2020	10210A WG-9954-070920-SG-033 07/09/2020
Parameters	Unit				
SVOCs-Continued					
Hexachloroethane	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Indeno(1,2,3-cd)pyrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Isophorone	μg/L	9.1 U	9.1 U	9.1 UJ	9.1 U
N-Nitrosodi-n-propylamine	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
N-Nitrosodiphenylamine	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Naphthalene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Nitrobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Pentachlorophenol	μg/L	45 U	45 U	45 U	45 U
Phenanthrene	μg/L 	9.1 U	9.1 U	9.1 U	9.1 U
Phenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Pyrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
SVOC, TICs					
(4-Methylphenyl)phenylmethanone	μg/L	-	-	-	-
1,4-Dioxane	μg/L	-	-	-	-
1-Chloro-2-methyl-benzene	μg/L	-	-	-	-
1-Chloro-3-methylbenzene	μg/L	-	-	-	-
1-Methyl-2-(phenylmethyl)-benzene	μg/L	-	-	-	-
2,3-Dichlorobenzoic acid	μg/L	-	-	-	-
2,5-Dichlorobenzoic acid	μg/L	-	-	-	-
2-Ethyl-ethanoic acid	μg/L	-	-	-	-
2-Propyl tridecyl-sulfurous acid	μg/L	-	-	-	-
3,4-Dichlorophenol	μg/L	-	-	-	-
3,5-Dichlorophenol	μg/L	-	-	-	-
3-Benzoylbenzoic acid	μg/L	-	-	-	-
3-Chlorobenzoic acid	μg/L	-	-	-	-
3-Methyl-cyclohexanol	μg/L 	-	-	-	-
4-Benzoyl-(rel)-benzoic acid	μg/L 	-	-	-	-
4-Chlorobenzoic acid	μg/L	-	-	-	-
4-Chlorophenol	μg/L	-	-	-	-
4-Chlorotoluene	μg/L	-	-	-	-
6-Octadecenoic acid	μg/L	-	-	-	-
Bicyclo[2.2.1]heptan-2-one, 1,3,3-trimethyl	μg/L	-	-	-	-
Chlorobenzene	μg/L	-	-	-	-
Cyclohexanol	μg/L	-	-	-	- 04 IN
Dimethyl disulfide	μg/L	-	-	-	21 JN
Dimethyl tetrasulfide	μg/L	-	-	-	-
Dodecanoic acid	μg/L	-	-	-	-

	Location ID: Sample Name: Sample Date:	10178A WG-9954-070720-SG-031 07/07/2020	10178B WG-9954-071020-SG-036 07/10/2020	10205 WG-9954-062520-SG-013 06/25/2020	10210A WG-9954-070920-SG-033 07/09/2020
Parameters	Unit				
SVOCs, TICs-Continued					
Hexadecanoic acid	μg/L	-	-	-	-
Homomenthyl salicylate	μg/L	-	-	-	-
Nonanoic acid	μg/L	-	-	4.5 JN	-
Sulfur	μg/L	-	-	42 JN	8.7 JN
Toluene	μg/L	-	-	-	-
trans-2-Methylcyclohexanol	μg/L	-	-	-	-
Unknown	μg/L	-	-	81.1 J	71.7 J
Polychlorinated Biphenyls (PCBs)					
Aroclor-1016 (PCB-1016)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1221 (PCB-1221)	μg/L	1.8 U	1.8 U	1.8 U	1.8 U
Aroclor-1232 (PCB-1232)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1242 (PCB-1242)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1248 (PCB-1248)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1254 (PCB-1254)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1260 (PCB-1260)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Pesticides					
4,4'-DDD	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
4,4'-DDE	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
4.4'-DDT	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Aldrin	μg/L	0.045 U	0.045 U	0.019 J	0.045 U
alpha-BHC	μg/L	0.045 U	0.069	0.18	0.12
alpha-Chlordane	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
beta-BHC	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
delta-BHC	μg/L	0.045 U	0.021 J	0.28	0.088
Dieldrin	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Endosulfan I	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Endosulfan II	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Endosulfan sulfate	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Endrin	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Endrin ketone	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
gamma-BHC (lindane)	μg/L	0.045 U	0.058	0.21	0.15
gamma-Chlordane	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Heptachlor	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Heptachlor epoxide	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Methoxychlor	μg/L	0.045 U	0.041 J	0.045 U	0.045 UJ
Toxaphene	μg/L	0.46 U	0.46 U	0.46 U	0.46 U

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	Location ID:	10178A	10178B	10205	10210A
	Sample Name:	WG-9954-070720-SG-031	WG-9954-071020-SG-036	WG-9954-062520-SG-013	WG-9954-070920-SG-033
	Sample Date:	07/07/2020	07/10/2020	06/25/2020	07/09/2020
Parameters	Unit				
Field Parameters Temperature, field Conductivity, field Turbidity, field	Deg C	12.7	13.0	13.4	12.9
	mS/cm	1.70	2.22	5.74	58.56
	NTU	72.7	111	28.2	620
pH, field	S.U.	7.29	8.16	6.99	6.96

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Sai	Location ID: mple Name: ample Date:	10210B WG-9954-063020-SG-019 06/30/2020	10210C WG-9954-063020-SG-020 06/30/2020	10210C WG-9954-063020-SG-021 06/30/2020 Duplicate	10215 WG-9954-063020-SG-022 06/30/2020
Parameters	Unit				
Volatile Organic Compounds (VOCs)					
1,1,1-Trichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1,2,2-Tetrachloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1,2-Trichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1-Dichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1-Dichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,2-Dichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,2-Dichloropropane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
2-Butanone (Methyl ethyl ketone) (MEK)	μg/L	10 U	10 U	10 U	10 U
2-Hexanone	μg/L	10 U	10 U	10 U	10 U
4-Methyl-2-pentanone (Methyl isobutyl ketone) (MIBH		10 U	10 U	10 U	10 U
Acetone	μg/L	10 U	10 U	10 U	10 U
Benzene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Bromodichloromethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Bromoform	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Bromomethane (Methyl bromide)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Carbon disulfide	μg/L	4.3 J	4.2 J	3.1 J	3.5 J
Carbon tetrachloride	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chlorobenzene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chloroform (Trichloromethane)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chloromethane (Methyl chloride)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
cis-1,2-Dichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
cis-1,3-Dichloropropene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Dibromochloromethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Ethylbenzene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Methylene chloride	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Styrene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Tetrachloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Toluene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
trans-1,2-Dichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
trans-1,3-Dichloropropene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Trichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Vinyl acetate	μg/L	10 U	10 U	10 U	10 U
Vinyl chloride	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Xylenes (total)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U

Location Sample Nan Sample Da	ie:	10210B WG-9954-063020-SG-019 06/30/2020	10210C WG-9954-063020-SG-020 06/30/2020	10210C WG-9954-063020-SG-021 06/30/2020 Duplicate	10215 WG-9954-063020-SG-022 06/30/2020
Parameters	Unit				
VOCs, Tentatively Identified Compounds (TICs)					
1,2,4-Trichlorobenzene	μg/L	-	-	-	-
1-Chloro-2-methyl-benzene	μg/L	-	-	-	-
1-Chloro-3-methylbenzene	μg/L	-	-	-	-
2,6-Dichlorotoluene	μg/L	-	-	-	-
2-Methylbutane	μg/L	-	-	-	-
6-Methyl-5-hepten-2-one	μg/L	-	-	-	-
Dimethyl disulfide	μg/L	-	-	-	-
Dimethyl sulfide	μg/L	-	-	-	-
Dimethyl trisulfide	μg/L	-	-	-	-
Hexanal	μg/L	-	-	-	-
Isobutane	μg/L	-	-	-	-
m-Monochlorobenzotrifluoride	μg/L	-	-	-	-
Methanethiol	μg/L	-	-	-	-
Methoxytrimethyl-silane	μg/L	-	-	-	-
Methylthioethane	μg/L	-	-	-	-
Nonanal	μg/L	-	-	-	-
Propane	μg/L	-	-	-	-
Sulfur dioxide (SO2)	μg/L	78.4 J	41.3 J	46.7 J	51.5 J
Trimethylfluorosilane	μg/L	- 206 J	- 47.4 J	-	- 24.7 J
Unknown	μg/L	206 J	47.4 J	24.0 J	24.7 J
Semi-volatile Organic Compounds (SVOCs)					
1,2,4-Trichlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
1,2-Dichlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
1,3-Dichlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
1,4-Dichlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,2'-Oxybis(1-chloropropane) (bis(2-Chloroisopropyl) ether)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4,5-Trichlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4,6-Trichlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4-Dichlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4-Dimethylphenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4-Dinitrophenol	μg/L	45 U	45 U	45 U	45 U
2,4-Dinitrotoluene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,6-Dinitrotoluene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Chloronaphthalene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Chlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Methylnaphthalene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Methylphenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U

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	Location ID: Sample Name: Sample Date:	10210B WG-9954-063020-SG-019 06/30/2020	10210C WG-9954-063020-SG-020 06/30/2020	10210C WG-9954-063020-SG-021 06/30/2020 Duplicate	10215 WG-9954-063020-SG-022 06/30/2020
Parameters	Unit				
SVOCs-Continued					
2-Nitroaniline	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Nitrophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
3&4-Methylphenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
3,3'-Dichlorobenzidine	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
3-Nitroaniline	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4,6-Dinitro-2-methylphenol	μg/L	45 U	45 U	45 U	45 U
4-Bromophenyl phenyl ether	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Chloro-3-methylphenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Chloroaniline	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Chlorophenyl phenyl ether	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Nitroaniline	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Nitrophenol	μg/L	45 U	45 U	45 U	45 U
Acenaphthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Acenaphthylene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Anthracene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(a)anthracene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(a)pyrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(b)fluoranthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(g,h,i)perylene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(k)fluoranthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzoic acid	μg/L	91 U	91 U	91 U	91 U
Benzyl alcohol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
bis(2-Chloroethoxy)methane	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
bis(2-Chloroethyl)ether	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
bis(2-Ethylhexyl)phthalate (DEHP)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Butyl benzylphthalate (BBP)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Chrysene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Di-n-butylphthalate (DBP)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Di-n-octyl phthalate (DnOP)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Dibenz(a,h)anthracene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Dibenzofuran	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Diethyl phthalate	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Dimethyl phthalate	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Fluoranthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Fluorene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Hexachlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Hexachlorobutadiene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Hexachlorocyclopentadiene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U

	Location ID: Sample Name: Sample Date:	10210B WG-9954-063020-SG-019 06/30/2020	10210C WG-9954-063020-SG-020 06/30/2020	10210C WG-9954-063020-SG-021 06/30/2020 Duplicate	10215 WG-9954-063020-SG-022 06/30/2020
Parameters	Unit				
SVOCs-Continued					
Hexachloroethane	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Indeno(1,2,3-cd)pyrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Isophorone	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
N-Nitrosodi-n-propylamine	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
N-Nitrosodiphenylamine	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Naphthalene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Nitrobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Pentachlorophenol	μg/L	45 U	45 U	45 U	45 U
Phenanthrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Phenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Pyrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
SVOC, TICs					
(4-Methylphenyl)phenylmethanone	μg/L	-	-	-	-
1,4-Dioxane	μg/L	-	-	-	-
1-Chloro-2-methyl-benzene	μg/L	-	-	-	-
1-Chloro-3-methylbenzene	μg/L	-	-	-	-
1-Methyl-2-(phenylmethyl)-benzene	μg/L	_	-	<u>-</u>	_
2,3-Dichlorobenzoic acid	μg/L	_	-	<u>-</u>	_
2,5-Dichlorobenzoic acid	μg/L	_	_	_	_
2-Ethyl-ethanoic acid	μg/L	_	_	_	_
2-Propyl tridecyl-sulfurous acid	μg/L	_	_	_	_
3,4-Dichlorophenol	μg/L	_	_	_	_
3,5-Dichlorophenol	μg/L	_	_	_	_
3-Benzoylbenzoic acid	μg/L		_		_
3-Chlorobenzoic acid	μg/L μg/L		- -	<u>-</u>	- -
3-Methyl-cyclohexanol	μg/L μg/L		- -	<u>-</u>	- -
4-Benzoyl-(rel)-benzoic acid	μg/L μg/L	-	<u>-</u>	<u>-</u>	-
4-Chlorobenzoic acid	μg/L	-	-	-	-
4-Chlorophenol	μg/L	-	-	-	-
•	μg/L	-	-	-	-
4-Chlorotoluene	μg/L	-	-	-	-
6-Octadecenoic acid	μg/L	-	-	-	-
Bicyclo[2.2.1]heptan-2-one, 1,3,3-trimethyl	μg/L	-	-	-	-
Chlorobenzene	μg/L	-	-	-	-
Cyclohexanol	μg/L	-	-	-	-
Dimethyl disulfide	μg/L 	-	-	-	-
Dimethyl tetrasulfide	μg/L 	-	-	-	-
Dodecanoic acid	μg/L	-	-	-	-

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	Location ID: Sample Name: Sample Date:	10210B WG-9954-063020-SG-019 06/30/2020	10210C WG-9954-063020-SG-020 06/30/2020	10210C WG-9954-063020-SG-021 06/30/2020 Duplicate	10215 WG-9954-063020-SG-022 06/30/2020
Parameters	Unit				
SVOCs, TICs-Continued					
Hexadecanoic acid	μg/L	-	-	-	-
Homomenthyl salicylate	μg/L	-	-	-	4.5 JN
Nonanoic acid	μg/L 	-		-	-
Sulfur	μg/L	-	6.4 JN	-	-
Toluene	μg/L	-	-	-	-
trans-2-Methylcyclohexanol	μg/L 		-	-	
Unknown	μg/L	31.9 J	36.9 J	20.6 J	64.2 J
Polychlorinated Biphenyls (PCBs)					
Aroclor-1016 (PCB-1016)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1221 (PCB-1221)	μg/L	1.8 U	1.8 U	1.8 U	1.8 U
Aroclor-1232 (PCB-1232)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1242 (PCB-1242)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1248 (PCB-1248)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1254 (PCB-1254)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1260 (PCB-1260)	μ <mark>g</mark> /L	0.91 U	0.91 U	0.91 U	0.91 U
Pesticides					
4,4'-DDD	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
4,4'-DDE	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
4,4'-DDT	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Aldrin	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
alpha-BHC	μg/L	0.12	0.081	0.10	0.064
alpha-Chlordane	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
beta-BHC	μg/L	0.045 U	0.045 U	0.045 U	0.027 J
delta-BHC	μg/L	0.076	0.068 J	0.12 J	0.29
Dieldrin	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Endosulfan I	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Endosulfan II	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Endosulfan sulfate	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Endrin	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Endrin ketone	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
gamma-BHC (lindane)	μg/L	0.10	0.089	0.11	0.074
gamma-Chlordane	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Heptachlor	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Heptachlor epoxide	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Methoxychlor	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Toxaphene	μg/L	0.46 U	0.46 U	0.46 U	0.46 U

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	Location ID: Sample Name: Sample Date:	10210B WG-9954-063020-SG-019 06/30/2020	10210C WG-9954-063020-SG-020 06/30/2020	10210C WG-9954-063020-SG-021 06/30/2020 Duplicate	10215 WG-9954-063020-SG-022 06/30/2020
Parameters	Unit				
Field Parameters					
Temperature, field	Deg C	13.5	14.0	14.0	14.4
Conductivity, field	mS/cm	11.63	5.91	5.91	5.44
Turbidity, field	NTU	19.0	3.96	3.96	6.25
pH, field	s.u.	7.06	7.07	7.07	7.16

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	Location ID: Sample Name: Sample Date:	10225A WG-9954-070920-SG-034 07/09/2020	10225B WG-9954-070920-SG-035 07/09/2020	10225C WG-9954-070120-SG-026 07/01/2020	10270 WG-9954-070720-SG-032 07/07/2020
Parameters	Unit				
Volatile Organic Compounds (VOCs)					
1,1,1-Trichloroethane	μg/L	25 U	25 U	5.0 U	5.0 U
1,1,2,2-Tetrachloroethane	μg/L	25 U	25 U	5.0 U	5.0 U
1,1,2-Trichloroethane	μg/L	25 U	25 U	5.0 U	5.0 U
1,1-Dichloroethane	μg/L	25 U	25 U	5.0 U	5.0 U
1,1-Dichloroethene	μg/L	25 U	25 U	5.0 U	5.0 U
1,2-Dichloroethane	μg/L	25 U	25 U	5.0 U	5.0 U
1,2-Dichloropropane	μg/L	25 U	25 U	5.0 U	5.0 U
2-Butanone (Methyl ethyl ketone) (MEK)	μg/L	50 U	50 U	10 U	10 U
2-Hexanone	μg/L	50 U	50 U	10 U	10 U
4-Methyl-2-pentanone (Methyl isobutyl ketone) (50 U	50 U	10 U	10 U
Acetone	μg/L	50 U	50 U	10 U	10 U
Benzene	μg/L	25 U	25 U	5.0 U	5.0 U
Bromodichloromethane	μg/L	25 U	25 U	5.0 U	5.0 U
Bromoform	μg/L	25 U	25 U	5.0 U	5.0 U
Bromomethane (Methyl bromide)	μg/L	25 U	25 U	5.0 U	5.0 U
Carbon disulfide	μg/L	160	25 J	3.3 J	3.7 J
Carbon tetrachloride	μg/L	25 U	25 U	5.0 U	5.0 U
Chlorobenzene	μg/L	25 U	25 U	0.89 J	5.0 U
Chloroethane	μg/L	25 U	25 U	5.0 U	5.0 U
Chloroform (Trichloromethane)	μg/L	25 U	25 U	5.0 U	5.0 U
Chloromethane (Methyl chloride)	μg/L	25 U	25 U	5.0 U	5.0 U
cis-1,2-Dichloroethene	μg/L	25 U	25 U	3.3 J	5.0 U
cis-1,3-Dichloropropene	μg/L	25 U	25 U	5.0 U	5.0 U
Dibromochloromethane	μg/L	25 U	25 U	5.0 U	5.0 U
Ethylbenzene	μg/L	25 U	25 U	5.0 U	5.0 U
Methylene chloride	μg/L	25 U	25 U	5.0 U	5.0 U
Styrene	μg/L	25 U	25 U	5.0 U	5.0 U
Tetrachloroethene	μg/L	25 U	25 U	5.0 U	5.0 U
Toluene	μg/L	25 U	25 U	5.0 U	5.0 U
trans-1,2-Dichloroethene	μg/L	25 U	25 U	5.0 U	5.0 U
trans-1,3-Dichloropropene	μg/L	25 U	25 U	5.0 U	5.0 U
Trichloroethene	μg/L	25 U	25 U	6.5	5.0 U
Vinyl acetate	μg/L	50 U	50 U	10 U	10 U
Vinyl chloride	μg/L	25 U	25 U	5.0 U	5.0 U
Xylenes (total)	μg/L	25 U	25 U	5.0 U	5.0 U

Location IE Sample Name Sample Date):	10225A WG-9954-070920-SG-034 07/09/2020	10225B WG-9954-070920-SG-035 07/09/2020	10225C WG-9954-070120-SG-026 07/01/2020	10270 WG-9954-070720-SG-032 07/07/2020
Parameters	Unit				
VOCs, Tentatively Identified Compounds (TICs)					
1,2,4-Trichlorobenzene	μg/L	-	-	12.7 J	-
1-Chloro-2-methyl-benzene	μg/L	-	-	8.1 J	-
1-Chloro-3-methylbenzene	μg/L	-	-	-	-
2,6-Dichlorotoluene	μg/L	-	-	-	-
2-Methylbutane	μg/L	-	-	-	-
6-Methyl-5-hepten-2-one	μg/L	-	-	-	-
Dimethyl disulfide	μg/L	45.7 J	-	-	-
Dimethyl sulfide	μg/L	594.4 J	-	-	-
Dimethyl trisulfide	μg/L	-	-		-
Hexanal	μg/L	-	-	6.7 J	-
Isobutane	μg/L	-	-	-	-
m-Monochlorobenzotrifluoride	μg/L	-	-	6.1 J	-
Methanethiol	μg/L	78.0 J	-	=	-
Methoxytrimethyl-silane	μg/L	-	-	-	-
Methylthioethane	μg/L	50.8 J	-	-	-
Nonanal	μg/L	-	-	-	-
Propane	μg/L	-	- 4550 5 1	-	-
Sulfur dioxide (SO2)	μg/L	213.2 J	1552.5 J	46.1 J	1.4 J
Trimethylfluorosilane	μg/L	-	-	-	-
Unknown	μg/L	-	-	=	-
Semi-volatile Organic Compounds (SVOCs)					
1,2,4-Trichlorobenzene	μg/L	9.1 U	9.1 U	7.7 J	9.1 U
1,2-Dichlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
1,3-Dichlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
1,4-Dichlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,2'-Oxybis(1-chloropropane) (bis(2-Chloroisopropyl) ether)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4,5-Trichlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4,6-Trichlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4-Dichlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4-Dimethylphenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4-Dinitrophenol	μg/L	45 U	45 U	45 U	45 U
2,4-Dinitrotoluene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,6-Dinitrotoluene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Chloronaphthalene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Chlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Methylnaphthalene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Methylphenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U

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	Location ID: Sample Name: Sample Date:	10225A WG-9954-070920-SG-034 07/09/2020	10225B WG-9954-070920-SG-035 07/09/2020	10225C WG-9954-070120-SG-026 07/01/2020	10270 WG-9954-070720-SG-032 07/07/2020
Parameters	Unit				
SVOCs-Continued					
2-Nitroaniline	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Nitrophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
3&4-Methylphenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
3,3'-Dichlorobenzidine	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
3-Nitroaniline	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4,6-Dinitro-2-methylphenol	μg/L	45 U	45 U	45 U	45 U
4-Bromophenyl phenyl ether	μ <mark>g</mark> /L	9.1 U	9.1 U	9.1 U	9.1 U
4-Chloro-3-methylphenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Chloroaniline	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Chlorophenyl phenyl ether	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Nitroaniline	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Nitrophenol	μg/L	45 U	45 U	45 U	45 U
Acenaphthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Acenaphthylene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Anthracene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(a)anthracene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(a)pyrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(b)fluoranthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(g,h,i)perylene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(k)fluoranthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzoic acid	μg/L	91 U	91 U	91 U	91 U
Benzyl alcohol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
bis(2-Chloroethoxy)methane	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
bis(2-Chloroethyl)ether	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
bis(2-Ethylhexyl)phthalate (DEHP)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Butyl benzylphthalate (BBP)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Chrysene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Di-n-butylphthalate (DBP)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Di-n-octyl phthalate (DnOP)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Dibenz(a,h)anthracene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Dibenzofuran	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Diethyl phthalate	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Dimethyl phthalate	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Fluoranthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Fluorene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Hexachlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Hexachlorobutadiene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Hexachlorocyclopentadiene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U

	Location ID: Sample Name: Sample Date:	10225A WG-9954-070920-SG-034 07/09/2020	10225B WG-9954-070920-SG-035 07/09/2020	10225C WG-9954-070120-SG-026 07/01/2020	10270 WG-9954-070720-SG-032 07/07/2020
Parameters	Unit				
SVOCs-Continued					
Hexachloroethane	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Indeno(1,2,3-cd)pyrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Isophorone	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
N-Nitrosodi-n-propylamine	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
N-Nitrosodiphenylamine	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Naphthalene	μg/L	1.2 J	9.1 U	9.1 U	9.1 U
Nitrobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Pentachlorophenol	μg/L	45 U	45 U	45 U	45 U
Phenanthrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Phenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Pyrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
SVOC, TICs					
(4-Methylphenyl)phenylmethanone	μg/L	-	-	-	-
1,4-Dioxane	μg/L	-	-	-	-
1-Chloro-2-methyl-benzene	μg/L	-	-	7.9 JN	-
1-Chloro-3-methylbenzene	μg/L	-	-	-	-
1-Methyl-2-(phenylmethyl)-benzene	μg/L	-	-	-	-
2,3-Dichlorobenzoic acid	μg/L	-	-	-	-
2,5-Dichlorobenzoic acid	μg/L	-	-	-	-
2-Ethyl-ethanoic acid	μg/L	-	-	-	-
2-Propyl tridecyl-sulfurous acid	μg/L	-	-	-	-
3,4-Dichlorophenol	μg/L	-	-	-	-
3,5-Dichlorophenol	μg/L	-	-	-	-
3-Benzoylbenzoic acid	μg/L	-	-	-	-
3-Chlorobenzoic acid	μg/L	-	-	-	-
3-Methyl-cyclohexanol	μg/L	-	-	-	-
4-Benzoyl-(rel)-benzoic acid	μg/L	-	-	-	-
4-Chlorobenzoic acid	μg/L	-	-	-	-
4-Chlorophenol	μg/L	-	-	-	-
4-Chlorotoluene	μg/L	-	-	-	-
6-Octadecenoic acid	μg/L	-	-	-	-
Bicyclo[2.2.1]heptan-2-one, 1,3,3-trimethyl	μg/L	-	-	-	-
Chlorobenzene	μg/L	-	-	-	-
Cyclohexanol	μg/L		-	-	-
Dimethyl disulfide	μg/L	24 JN	-	-	-
Dimethyl tetrasulfide	μg/L	190 JN	-		-
Dodecanoic acid	μg/L	-	-	6.1 JN	-

	Location ID: Sample Name: Sample Date:	10225A WG-9954-070920-SG-034 07/09/2020	10225B WG-9954-070920-SG-035 07/09/2020	10225C WG-9954-070120-SG-026 07/01/2020	10270 WG-9954-070720-SG-032 07/07/2020
Parameters	Unit				
SVOCs, TICs-Continued					
Hexadecanoic acid	μg/L	-	-	-	-
Homomenthyl salicylate	μg/L	-	-	-	-
Nonanoic acid	μg/L	-	-	-	-
Sulfur	μg/L	-	-	14 JN	4.2 JN
Toluene	μg/L	-	-	-	-
trans-2-Methylcyclohexanol	μg/L	-	-	-	-
Unknown	μg/L		20.7 J	26.1 J	30.6 J
Polychlorinated Biphenyls (PCBs)					
Aroclor-1016 (PCB-1016)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1221 (PCB-1221)	μg/L	1.8 U	1.8 U	1.8 U	1.8 U
Aroclor-1232 (PCB-1232)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1242 (PCB-1242)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1248 (PCB-1248)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1254 (PCB-1254)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1260 (PCB-1260)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Pesticides					
4,4'-DDD	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
4,4'-DDE	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
4,4'-DDT	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Aldrin	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
alpha-BHC	μg/L	0.069	0.072	0.27	0.070
alpha-Chlordane	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
beta-BHC	μg/L	0.045 U	0.019 J	0.040 J	0.045 U
delta-BHC	μg/L	0.065	0.19	0.17	0.067
Dieldrin	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Endosulfan I	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Endosulfan II	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Endosulfan sulfate	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Endrin	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Endrin ketone	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
gamma-BHC (lindane)	μg/L	0.077	0.11	0.21	0.098
gamma-Chlordane	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Heptachlor	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Heptachlor epoxide	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Methoxychlor	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Toxaphene	μg/L	0.46 U	0.46 U	0.46 U	0.46 U

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	Location ID: Sample Name: Sample Date:	10225A WG-9954-070920-SG-034 07/09/2020	10225B WG-9954-070920-SG-035 07/09/2020	10225C WG-9954-070120-SG-026 07/01/2020	10270 WG-9954-070720-SG-032 07/07/2020
Parameters	Unit				
Field Parameters					
Temperature, field	Deg C	12.8	12.6	13.1	14.4
Conductivity, field	mS/cm	128.8	6.50	3.68	3.56
Turbidity, field	NTU	67.8	59.9	1.97	6.40
pH, field	s.u.	7.60	8.13	7.31	7.18

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	Location ID: cample Name: Sample Date:	10272 WG-9954-070720-SG-030 07/07/2020	10278 WG-9954-070620-SG-029 07/06/2020	MW-01 WG-9954-063020-RM-023 06/30/2020	MW-02 WG-9954-070120-SG-024 07/01/2020
Parameters	Unit				
Volatile Organic Compounds (VOCs)					
1,1,1-Trichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1,2,2-Tetrachloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1,2-Trichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1-Dichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1-Dichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,2-Dichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,2-Dichloropropane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
2-Butanone (Methyl ethyl ketone) (MEK)	μg/L	10 U	10 U	10 U	10 U
2-Hexanone	μg/L	10 U	10 U	10 U	10 U
4-Methyl-2-pentanone (Methyl isobutyl ketone) (MI		10 U	10 U	10 U	10 U
Acetone	μg/L	10 U	10 U	10 U	10 U
Benzene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Bromodichloromethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Bromoform	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Bromomethane (Methyl bromide)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Carbon disulfide	μg/L	3.4 J	6.4 J	2.9 J	10 U
Carbon tetrachloride	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chlorobenzene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chloroform (Trichloromethane)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chloromethane (Methyl chloride)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
cis-1,2-Dichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
cis-1,3-Dichloropropene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Dibromochloromethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Ethylbenzene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Methylene chloride	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Styrene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Tetrachloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Toluene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
trans-1,2-Dichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
trans-1,3-Dichloropropene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Trichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Vinyl acetate	μg/L	10 U	10 U	10 U	10 U
Vinyl chloride	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Xylenes (total)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U

Location II Sample Name Sample Date) :	10272 WG-9954-070720-SG-030 07/07/2020	10278 WG-9954-070620-SG-029 07/06/2020	MW-01 WG-9954-063020-RM-023 06/30/2020	MW-02 WG-9954-070120-SG-024 07/01/2020
Parameters	Unit				
VOCs, Tentatively Identified Compounds (TICs)					
1,2,4-Trichlorobenzene	μg/L	-	-	-	-
1-Chloro-2-methyl-benzene	μg/L	-	-	-	-
1-Chloro-3-methylbenzene	μg/L	-	-	-	-
2,6-Dichlorotoluene	μg/L	-	-	-	-
2-Methylbutane	μg/L	-	-	-	-
6-Methyl-5-hepten-2-one	μg/L	-	-	-	-
Dimethyl disulfide	μg/L	-	-	-	-
Dimethyl sulfide	μg/L	-	-	-	-
Dimethyl trisulfide	μg/L	-	-	-	-
Hexanal	μg/L	-	7.7 J	-	-
Isobutane	μg/L	-	-	-	-
m-Monochlorobenzotrifluoride	μg/L	-	-	-	-
Methanethiol	μg/L	-	-	-	-
Methoxytrimethyl-silane	μg/L	-	-	-	-
Methylthioethane	μg/L	-	-	-	-
Nonanal	μg/L	-	-	-	-
Propane	μg/L	- -		- 	<u>-</u>
Sulfur dioxide (SO2)	μg/L	48.8 J	303.7 J	47.6 J	66.0 J
Trimethylfluorosilane	μg/L	-	-	-	-
Unknown	μg/L	-	-	-	-
Semi-volatile Organic Compounds (SVOCs)					
1,2,4-Trichlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
1,2-Dichlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
1,3-Dichlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
1,4-Dichlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,2'-Oxybis(1-chloropropane) (bis(2-Chloroisopropyl) ether)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4,5-Trichlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4,6-Trichlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4-Dichlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4-Dimethylphenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,4-Dinitrophenol	μg/L	45 U	45 U	45 U	45 U
2,4-Dinitrotoluene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2,6-Dinitrotoluene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Chloronaphthalene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Chlorophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Methylnaphthalene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Methylphenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U

	Location ID: Sample Name: Sample Date:	10272 WG-9954-070720-SG-030 07/07/2020	10278 WG-9954-070620-SG-029 07/06/2020	MW-01 WG-9954-063020-RM-023 06/30/2020	MW-02 WG-9954-070120-SG-024 07/01/2020
Parameters	Unit				
SVOCs-Continued					
2-Nitroaniline	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
2-Nitrophenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
3&4-Methylphenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
3,3'-Dichlorobenzidine	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
3-Nitroaniline	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4,6-Dinitro-2-methylphenol	μg/L	45 U	45 U	45 U	45 U
4-Bromophenyl phenyl ether	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Chloro-3-methylphenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Chloroaniline	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Chlorophenyl phenyl ether	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Nitroaniline	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
4-Nitrophenol	μg/L	45 U	45 U	45 U	45 U
Acenaphthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Acenaphthylene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Anthracene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(a)anthracene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(a)pyrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(b)fluoranthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(g,h,i)perylene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzo(k)fluoranthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Benzoic acid	μg/L	91 U	91 U	91 U	91 U
Benzyl alcohol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
bis(2-Chloroethoxy)methane	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
bis(2-Chloroethyl)ether	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
bis(2-Ethylhexyl)phthalate (DEHP)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Butyl benzylphthalate (BBP)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Chrysene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Di-n-butylphthalate (DBP)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Di-n-octyl phthalate (DnOP)	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Dibenz(a,h)anthracene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Dibenzofuran	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Diethyl phthalate	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Dimethyl phthalate	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Fluoranthene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Fluorene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Hexachlorobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Hexachlorobutadiene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Hexachlorocyclopentadiene	μg/L	R	9.1 U	9.1 U	9.1 U

	Location ID: Sample Name: Sample Date:	10272 WG-9954-070720-SG-030 07/07/2020	10278 WG-9954-070620-SG-029 07/06/2020	MW-01 WG-9954-063020-RM-023 06/30/2020	MW-02 WG-9954-070120-SG-024 07/01/2020
Parameters	Unit				
SVOCs-Continued					
Hexachloroethane	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Indeno(1,2,3-cd)pyrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Isophorone	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
N-Nitrosodi-n-propylamine	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
N-Nitrosodiphenylamine	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Naphthalene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Nitrobenzene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Pentachlorophenol	μg/L	45 U	45 U	45 U	45 U
Phenanthrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Phenol	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
Pyrene	μg/L	9.1 U	9.1 U	9.1 U	9.1 U
SVOC, TICs					
(4-Methylphenyl)phenylmethanone	μg/L	-	-	-	-
1,4-Dioxane	μg/L	-	-	-	-
1-Chloro-2-methyl-benzene	μg/L	-	-	-	-
1-Chloro-3-methylbenzene	μg/L	-	-	-	-
1-Methyl-2-(phenylmethyl)-benzene	μg/L	-	-	-	-
2,3-Dichlorobenzoic acid	μg/L	-	-	-	-
2,5-Dichlorobenzoic acid	μg/L	-	-	-	-
2-Ethyl-ethanoic acid	μg/L	-	-	-	-
2-Propyl tridecyl-sulfurous acid	μg/L	-	-	-	7.0 JN
3,4-Dichlorophenol	μg/L	-	-	-	-
3,5-Dichlorophenol	μg/L	-	-	-	-
3-Benzoylbenzoic acid	μg/L	-	-	-	-
3-Chlorobenzoic acid	μg/L	-	-	-	-
3-Methyl-cyclohexanol	μg/L	-	-	-	-
4-Benzoyl-(rel)-benzoic acid	μg/L	-	-	-	-
4-Chlorobenzoic acid	μg/L	-	-	-	-
4-Chlorophenol	μg/L	-	-	-	-
4-Chlorotoluene	μg/L	-	-	-	-
6-Octadecenoic acid	μg/L	-	-	-	-
Bicyclo[2.2.1]heptan-2-one, 1,3,3-trimethyl	μg/L	-	-	-	-
Chlorobenzene	μg/L	-	-	-	-
Cyclohexanol	μg/L	-	-	-	-
Dimethyl disulfide	μg/L 	-	-	-	-
Dimethyl tetrasulfide	μg/L	-	-	-	-
Dodecanoic acid	μg/L	-	-	-	-

	Location ID: Sample Name: Sample Date:	10272 WG-9954-070720-SG-030 07/07/2020	10278 WG-9954-070620-SG-029 07/06/2020	MW-01 WG-9954-063020-RM-023 06/30/2020	MW-02 WG-9954-070120-SG-024 07/01/2020
Parameters	Unit				
SVOCs, TICs-Continued					
Hexadecanoic acid	μg/L	-	-	-	-
Homomenthyl salicylate	μg/L	-	-	-	-
Nonanoic acid	μg/L	-	-	-	-
Sulfur	μg/L	-	13 JN	-	-
Toluene	μg/L	-	-	-	-
trans-2-Methylcyclohexanol	μg/L	-	-	-	-
Unknown	μg/L	57.6 J	-	59.1 J	65.2 J
Polychlorinated Biphenyls (PCBs)					
Aroclor-1016 (PCB-1016)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1221 (PCB-1221)	μg/L	1.8 U	1.8 U	1.8 U	1.8 U
Aroclor-1232 (PCB-1232)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1242 (PCB-1242)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1248 (PCB-1248)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1254 (PCB-1254)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Aroclor-1260 (PCB-1260)	μg/L	0.91 U	0.91 U	0.91 U	0.91 U
Pesticides					
4,4'-DDD	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
4,4'-DDE	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
4,4'-DDT	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Aldrin	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
alpha-BHC	μg/L	0.24	0.47	0.045 U	0.045 U
alpha-Chlordane	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
beta-BHC	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
delta-BHC	μg/L	0.084	0.13	0.045 U	0.045 U
Dieldrin	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Endosulfan I	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Endosulfan II	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Endosulfan sulfate	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Endrin	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Endrin ketone	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
gamma-BHC (lindane)	μg/L	0.26	0.48	0.045 U	0.045 U
gamma-Chlordane	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Heptachlor	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Heptachlor epoxide	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Methoxychlor	μg/L	0.045 U	0.045 U	0.045 U	0.045 U
Toxaphene	μg/L	0.46 U	0.46 U	0.46 U	0.46 U

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	Location ID:	10272	10278	MW-01	MW-02
	Sample Name:	WG-9954-070720-SG-030	WG-9954-070620-SG-029	WG-9954-063020-RM-023	WG-9954-070120-SG-024
	Sample Date:	07/07/2020	07/06/2020	06/30/2020	07/01/2020
Parameters	Unit				
Field Parameters Temperature, field Conductivity, field Turbidity, field pH, field	Deg C	14.5	13.4	13.3	13.0
	mS/cm	5.20	5.13	3.75	3.40
	NTU	13.0	3.31	59.0	20.5
	s.u.	6.97	7.33	7.18	7.19

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	Location ID: Sample Name: Sample Date:	RINSE BLANK RB-9954-062520-SG-001 06/25/2020	RINSE BLANK RB-9954-063020-SG-002 06/30/2020	Trip Blank TB-9954-062420-SG-001 06/24/2020	Trip Blank TB-9954-062520-SG-002 06/25/2020
Parameters	Unit				
Volatile Organic Compounds (VOCs)					
1,1,1-Trichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1,2,2-Tetrachloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1,2-Trichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1-Dichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,1-Dichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,2-Dichloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
1,2-Dichloropropane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
2-Butanone (Methyl ethyl ketone) (MEK)	μg/L	10 U	10 U	10 U	10 U
2-Hexanone	μg/L	10 U	10 U	10 U	10 U
4-Methyl-2-pentanone (Methyl isobutyl ketone)	(MIBK) μg/L	10 U	10 U	10 U	10 U
Acetone	μg/L	10 U	10 U	10 U	10 U
Benzene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Bromodichloromethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Bromoform	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Bromomethane (Methyl bromide)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Carbon disulfide	μg/L	10 U	10 U	10 U	10 U
Carbon tetrachloride	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chlorobenzene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chloroethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chloroform (Trichloromethane)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Chloromethane (Methyl chloride)	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
cis-1,2-Dichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
cis-1,3-Dichloropropene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Dibromochloromethane	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Ethylbenzene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Methylene chloride	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Styrene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Tetrachloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Toluene	μg/L	5.0 U	0.25 J	5.0 U	5.0 U
trans-1,2-Dichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
trans-1,3-Dichloropropene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Trichloroethene	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Vinyl acetate	μg/L	10 U	10 U	10 U	10 U
Vinyl chloride	μg/L	5.0 U	5.0 U	5.0 U	5.0 U
Xylenes (total)	μ <mark>g</mark> /L	5.0 U	5.0 U	5.0 U	5.0 U

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Sample	tion ID: Name: e Date:	RINSE BLANK RB-9954-062520-SG-001 06/25/2020	RINSE BLANK RB-9954-063020-SG-002 06/30/2020	Trip Blank TB-9954-062420-SG-001 06/24/2020	Trip Blank TB-9954-062520-SG-002 06/25/2020
Parameters	Unit				
VOCs, Tentatively Identified Compounds (TICs)					
1,2,4-Trichlorobenzene	μg/L	-	-	-	-
1-Chloro-2-methyl-benzene	μg/L	-	-	-	-
1-Chloro-3-methylbenzene	μg/L	-	-	-	-
2,6-Dichlorotoluene	μg/L	-	-	-	-
2-Methylbutane	μg/L	-	-	-	-
6-Methyl-5-hepten-2-one	μg/L	-	-	-	-
Dimethyl disulfide	μg/L	-	-	-	-
Dimethyl sulfide	μg/L	-	-	-	-
Dimethyl trisulfide	μg/L	-	-	-	-
Hexanal	μg/L	-	-	-	-
Isobutane	μg/L	-	-	-	-
m-Monochlorobenzotrifluoride	μg/L	-	-	-	-
Methanethiol	μg/L	-	-	-	-
Methoxytrimethyl-silane	μg/L	-	-	-	-
Methylthioethane	μg/L	-	-	-	-
Nonanal	μg/L	-	-	-	-
Propane	μg/L	-	-	-	-
Sulfur dioxide (SO2)	μg/L	-	84.1 J	-	-
Trimethylfluorosilane	μg/L	-	-	-	-
Unknown	μg/L	-	-	-	-
Semi-volatile Organic Compounds (SVOCs)					
1,2,4-Trichlorobenzene	μg/L	9.1 U	9.1 U	-	-
1,2-Dichlorobenzene	μg/L	9.1 U	9.1 U	-	-
1,3-Dichlorobenzene	μg/L	9.1 U	9.1 U	-	-
1,4-Dichlorobenzene	μg/L	9.1 U	9.1 U	-	-
2,2'-Oxybis(1-chloropropane) (bis(2-Chloroisopropyl) ethe	r) µg/L	9.1 U	9.1 U	-	-
2,4,5-Trichlorophenol	μg/L	9.1 U	9.1 U	-	-
2,4,6-Trichlorophenol	μg/L	9.1 U	9.1 U	-	-
2,4-Dichlorophenol	μg/L	9.1 U	9.1 U	-	-
2,4-Dimethylphenol	μg/L	9.1 U	9.1 U	-	-
2,4-Dinitrophenol	μg/L	45 U	45 U	-	-
2,4-Dinitrotoluene	μg/L	9.1 U	9.1 U	-	-
2,6-Dinitrotoluene	μg/L	9.1 U	9.1 U	-	-
2-Chloronaphthalene	μg/L	9.1 U	9.1 U	-	-
2-Chlorophenol	μg/L	9.1 U	9.1 U	-	-
2-Methylnaphthalene	μg/L	9.1 U	9.1 U	-	-
2-Methylphenol	μg/L	9.1 U	9.1 U	-	-

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Parameters		Location ID: Sample Name: Sample Date:	RINSE BLANK RB-9954-062520-SG-001 06/25/2020	RINSE BLANK RB-9954-063020-SG-002 06/30/2020	Trip Blank TB-9954-062420-SG-001 06/24/2020	Trip Blank TB-9954-062520-SG-002 06/25/2020
2-Nitrophenol	Parameters	Unit				
2-Nitrophenol	SVOCs-Continued					
33-Helhylphenol	2-Nitroaniline	μg/L			-	-
3.3-Dichlorobenzdine	2-Nitrophenol	μg/L	9.1 U		-	-
3.3-Dichlorobenzidine	3&4-Methylphenol	μg/L	9.1 U	9.1 U	-	-
4.6-Dinitro-Z-methylphenol μg/L 9.1 U 9		μg/L			-	-
4-Bromophenyl phenyl ether μg/L 9.1 U 9.1 U -	3-Nitroaniline	μg/L			-	-
4-Chlora-methylphenol μg/L 9.1 U 9.1		μg/L	45 U		-	-
4-Chloraniline μg/L 9.1 U					-	-
4-Chlorophenyl phenyl ether μg/L 9.1 U		μg/L			-	-
4-Nitrophenol μg/L 45 U 45 U		μg/L			-	-
4-Nitrophenol μg/L 45 U 45 U -		μg/L			-	-
Acenaphthene μg/L 9.1 U 9.1 U - - Acenaphthylene μg/L 9.1 U 9.1 U - - Anthracene μg/L 9.1 U 9.1 U - - Benzo(a)anthracene μg/L 9.1 U 9.1 U - - Benzo(b)fluoranthene μg/L 9.1 U 9.1 U - - Benzo(k)fluoranthene μg/L 9.1 U 9.1 U - - Benzo(k)fluoranthene μg/L 9.1 U 9.1 U - - Benzolc acid μg/L 9.1 U 9.1 U - - Benzylachol μg/L 9.1 U 9.1 U - - Benzylachol μg/L 9.1 U 9.1 U - - bis(2-Chlorethoxy)methane μg/L 9.1 U 9.1 U - - bis(2-Chlorethoxy)methane μg/L 9.1 U 9.1 U - - bis(2-Chlorethoxy)methane μg/L 9.1 U 9.1 U <	4-Nitroaniline				-	-
Acenaphthylene Anthracene µg/L Anthracene µg/L Benzo(a)anthracene µg/L Benzo(a)pyrene µg/L Benzo(a)pyrene µg/L Benzo(a)fluoranthene µg/L Benzo(b)fluoranthene µg/L Benzo(b)fluoranthene µg/L Benzo(b)fluoranthene µg/L Benzo(b,fluoranthene µg/L Benzo(b,fluoranthene) µg/L Benzo(b,fluoranthene) µg/L Benzo(acid µg/L Benzo(acid µg/L Benzo(acid µg/L Benzo(acid µg/L Benzo(acid µg/L Benzo(acid µg/L Benzo(acid µg/L Benzo(acid µg/L Benzo(acid µg/L Benzo(acid µg/L Benzo(acid µg/L Benzo(acid Benzo(acid µg/L Benzo(acid µg/L Benzo(acid Benzo(acid µg/L Benzo(acid Benzo(acid µg/L Benzo(acid Ben	4-Nitrophenol	μg/L			-	-
Anthracene µg/L 9.1 U		μg/L			-	-
Benzo(a)anthracene µg/L 9.1 U 9.1 U -	Acenaphthylene	μg/L			-	-
Benzo(a)pyrene μg/L 9.1 U 9.1 U - - Benzo(b)fluoranthene μg/L 9.1 U 9.1 U - - Benzo(k)fluoranthene μg/L 9.1 U 9.1 U - - Benzol acid μg/L 9.1 U 9.1 U - - Benzyl alcohol μg/L 9.1 U 9.1 U - - bis(2-Chloroethoxy)methane μg/L 9.1 U 9.1 U - - bis(2-Chloroethy)ether μg/L 9.1 U 9.1 U - - bis(2-Chlybexyl)phthalate (DEHP) μg/L 9.1 U 9.1 U - - Butyl benzylphthalate (BBP) μg/L 9.1 U 9.1 U - - Butyl benzylphthalate (BBP) μg/L 9.1 U 9.1 U - - Butyl benzylphthalate (BBP) μg/L 9.1 U 9.1 U - - Butyl benzylphthalate (BBP) μg/L 9.1 U 9.1 U - - Di-n-butylphthalate (BBP)					-	-
Benzo(b)fluoranthene µg/L 9.1 U 9.1 U - - Benzo(g,h.i)perylene µg/L 9.1 U 9.1 U - - Benzo(s,h.i)perylene µg/L 9.1 U 9.1 U - - Benzoic acid µg/L 91 U 91 U - - Benzyl alcohol µg/L 9.1 U 9.1 U - - bis(2-Chloroethoxy)methane µg/L 9.1 U 9.1 U - - bis(2-Chloroethoxy)ghthralete (DEHP) µg/L 9.1 U 9.1 U - - bis(2-Ethylhexy)phthalate (DEHP) µg/L 9.1 U 9.1 U - - Butyl benzylphthalate (DEHP) µg/L 9.1 U 9.1 U - - Butyl benzylphthalate (BBP) µg/L 9.1 U 9.1 U - - Chrysene µg/L 9.1 U 9.1 U - - Di-n-butylphthalate (BBP) µg/L 9.1 U 9.1 U - - Di-n-butylphthalate (DnOP) <					-	-
Benzo(g,h,i)perylene µg/L 9.1 U 9.1 U - - Benzo(k)fluoranthene µg/L 9.1 U 9.1 U - - Benzol cacid µg/L 91 U 91 U - - Benzyl alcohol µg/L 9.1 U 9.1 U - - bis(2-Chloroethylymether µg/L 9.1 U 9.1 U - - bis(2-Ethylhexyl)phthalate (DEHP) µg/L 9.1 U 9.1 U - - Butyl benzylphthalate (BBP) µg/L 9.1 U 9.1 U - - Chrysene µg/L 9.1 U 9.1 U - - - Di-n-butylphthalate (DBP) µg/L 9.1 U 9.1 U - - - Di-n-butylphthalate (DnOP) µg/L 9.1 U 9.1 U - - - Di-n-butylphthalate (DnOP) µg/L 9.1 U 9.1 U - - - Di-n-butylphthalate (DnOP) µg/L 9.1 U 9.1 U - -	Benzo(a)pyrene				-	-
Benzo(k)fluoranthene µg/L 9.1 U 9.1 U - - Benzoic acid µg/L 91 U 91 U - - Benzyl alcohol µg/L 9.1 U 9.1 U - - bis(2-Chloroethoxy)methane µg/L 9.1 U 9.1 U - - bis(2-Chloroethyl)ether µg/L 9.1 U 9.1 U - - bis(2-Ethylpextyl)phthalate (DEHP) µg/L 9.1 U 9.1 U - - Butyl benzylphthalate (BBP) µg/L 9.1 U 9.1 U - - Chrysene µg/L 9.1 U 9.1 U - - - Di-n-butylphthalate (DBP) µg/L 9.1 U 9.1 U - - - Di-n-octyl phthalate (DnOP) µg/L 9.1 U 9.1 U - - - Dibenz(a,h)anthracene µg/L 9.1 U 9.1 U - - - Dibenz(bright) phthalate µg/L 9.1 U 9.1 U - -	Benzo(b)fluoranthene				-	-
Benzoic acid µg/L 91 U 91 U - - Benzyl alcohol µg/L 9.1 U 9.1 U - - bis(2-Chloreethoxy)methane µg/L 9.1 U 9.1 U - - bis(2-Chloreethy)lether µg/L 9.1 U 9.1 U - - bis(2-Ethylhexyl)phthalate (DEHP) µg/L 9.1 U 9.1 U - - Butyl benzylphthalate (BBP) µg/L 9.1 U 9.1 U - - Chrysene µg/L 9.1 U 9.1 U - - - Chrysene µg/L 9.1 U 9.1 U - - - Chrysene µg/L 9.1 U 9.1 U - - - Chrysene µg/L 9.1 U 9.1 U - - - Di-n-octyl phthalate (DBP) µg/L 9.1 U 9.1 U - - - Dibenz(a, h)anthracene µg/L 9.1 U 9.1 U - - -	Benzo(g,h,i)perylene	μg/L			-	-
Benzyl alcohol pg/L 9.1 U 9.1	Benzo(k)fluoranthene	μg/L	9.1 U		-	-
bis(2-Chloroethoxy)methane µg/L 9.1 U 9.1 U -	Benzoic acid	μg/L			-	-
bis(2-Chloroethyl)ether µg/L 9.1 U 9.1 U -		μg/L			-	-
bis(2-Ethylhexyl)phthalate (DEHP) µg/L 9.1 U 9.1 U - <td></td> <td>μg/L</td> <td></td> <td></td> <td>-</td> <td>-</td>		μg/L			-	-
Butyl benzylphthalate (BBP) μg/L 9.1 U 9.1 U - - Chrysene μg/L 9.1 U 9.1 U - - Di-n-butylphthalate (DBP) μg/L 9.1 U 9.1 U - - Di-n-octyl phthalate (DnOP) μg/L 9.1 U 9.1 U - - Dibenz(a,h)anthracene μg/L 9.1 U 9.1 U - - Dibenzofuran μg/L 9.1 U 9.1 U - - Diethyl phthalate μg/L 9.1 U 9.1 U - - Dimethyl phthalate μg/L 9.1 U 9.1 U - - Fluoranthene μg/L 9.1 U 9.1 U - - Fluorene μg/L 9.1 U 9.1 U - - Hexachlorobenzene μg/L 9.1 U 9.1 U - - Hexachlorobutadiene μg/L 9.1 U 9.1 U - -		μg/L			-	-
Chrysene μg/L 9.1 U 9.1 U - - Di-n-butylphthalate (DBP) μg/L 9.1 U 9.1 U - - Di-n-octyl phthalate (DnOP) μg/L 9.1 U 9.1 U - - Dibenz(a,h)anthracene μg/L 9.1 U 9.1 U - - Dibenzofuran μg/L 9.1 U 9.1 U - - Diethyl phthalate μg/L 9.1 U 9.1 U - - Dimethyl phthalate μg/L 9.1 U 9.1 U - - Fluoranthene μg/L 9.1 U 9.1 U - - Fluorene μg/L 9.1 U 9.1 U - - Hexachlorobenzene μg/L 9.1 U 9.1 U - - Hexachlorobutadiene μg/L 9.1 U 9.1 U - -		μg/L			-	-
Di-n-butylphthalate (DBP) μg/L 9.1 U 9.1 U -	Butyl benzylphthalate (BBP)	μg/L			-	-
Di-n-octyl phthalate (DnOP) μg/L 9.1 U 9.1 U -	Chrysene				-	-
Dibenz(a,h)anthracene $\mu g/L$ 9.1 U 9.1 U	Di-n-butylphthalate (DBP)	μg/L	9.1 U		-	-
Dibenzofuran $\mu g/L$ 9.1 U 9.1 U	Di-n-octyl phthalate (DnOP)	μg/L			-	-
Diethyl phthalate $\begin{array}{cccccccccccccccccccccccccccccccccccc$	Dibenz(a,h)anthracene		9.1 U		-	-
Dimethyl phthalate μg/L 9.1 U 9.1 U - <th< td=""><td>Dibenzofuran</td><td>μg/L</td><td>9.1 U</td><td>9.1 U</td><td>-</td><td>-</td></th<>	Dibenzofuran	μg/L	9.1 U	9.1 U	-	-
Fluoranthene $\mu g/L$ 9.1 U 9.1 U Fluorene $\mu g/L$ 9.1 U 9.1 U	Diethyl phthalate	μg/L	9.1 U	9.1 U	-	-
Fluoranthene $\mu g/L$ 9.1 U 9.1 U Fluorene $\mu g/L$ 9.1 U 9.1 U		μg/L			-	-
Fluorene $\mu g/L$ 9.1 U 9.1 U Hexachlorobenzene $\mu g/L$ 9.1 U 9.1 U		μg/L	9.1 U		-	-
Hexachlorobenzene $\mu g/L$ 9.1 U 9.1 U Hexachlorobutadiene $\mu g/L$ 9.1 U 9.1 U	Fluorene	μg/L			-	-
Hexachlorobutadiene μg/L 9.1 U 9.1 U	Hexachlorobenzene	μg/L	9.1 U	9.1 U	-	-
	Hexachlorobutadiene	μg/L	9.1 U	9.1 U	-	-
	Hexachlorocyclopentadiene		9.1 U	9.1 U	-	-

	Location ID: Sample Name: Sample Date:	RINSE BLANK RB-9954-062520-SG-001 06/25/2020	RINSE BLANK RB-9954-063020-SG-002 06/30/2020	Trip Blank TB-9954-062420-SG-001 06/24/2020	Trip Blank TB-9954-062520-SG-002 06/25/2020
Parameters	Unit				
SVOCs-Continued					
Hexachloroethane	μg/L	9.1 U	9.1 U	-	-
Indeno(1,2,3-cd)pyrene	μg/L	9.1 U	9.1 U	-	-
Isophorone	μg/L	9.1 U	9.1 U	-	-
N-Nitrosodi-n-propylamine	μg/L	9.1 U	9.1 U	-	-
N-Nitrosodiphenylamine	μg/L	9.1 U	9.1 U	-	-
Naphthalene	μg/L	9.1 U	9.1 U	-	-
Nitrobenzene	μg/L	9.1 U	9.1 U	-	-
Pentachlorophenol	μg/L	45 U	45 U	-	-
Phenanthrene	μg/L	9.1 U	9.1 U	-	-
Phenol	μg/L	9.1 U	0.94 J	-	-
Pyrene	μg/L	9.1 U	9.1 U	-	-
SVOC, TICs					
(4-Methylphenyl)phenylmethanone	μg/L	-	-	-	-
1,4-Dioxane	μg/L	-	-	-	-
1-Chloro-2-methyl-benzene	μg/L	-	-	-	-
1-Chloro-3-methylbenzene	μg/L	-	-	-	-
1-Methyl-2-(phenylmethyl)-benzene	μg/L	-	-	-	-
2,3-Dichlorobenzoic acid	μg/L	-	-	-	-
2,5-Dichlorobenzoic acid	μg/L	-	-	-	-
2-Ethyl-ethanoic acid	μg/L	-	-	-	-
2-Propyl tridecyl-sulfurous acid	μg/L	-	-	-	-
3,4-Dichlorophenol	μg/L	-	-	-	-
3,5-Dichlorophenol	μg/L	-	-	-	-
3-Benzoylbenzoic acid	μg/L	-	-	-	-
3-Chlorobenzoic acid	μg/L	-	-	-	-
3-Methyl-cyclohexanol	μg/L	-	-	-	-
4-Benzoyl-(rel)-benzoic acid	μg/L	-	-	-	-
4-Chlorobenzoic acid	μg/L	-	-	-	-
4-Chlorophenol	μg/L	-	-	-	-
4-Chlorotoluene	μg/L	-	-	-	-
6-Octadecenoic acid	μg/L 	-	-	-	-
Bicyclo[2.2.1]heptan-2-one, 1,3,3-trimethyl	μg/L	-	-	-	-
Chlorobenzene	μg/L	-	-	-	-
Cyclohexanol	μg/L	-	-	-	-
Dimethyl disulfide	μg/L	-	-	-	-
Dimethyl tetrasulfide	μg/L	-	-	-	-
Dodecanoic acid	μg/L	-	-	-	-

	Location ID: Sample Name: Sample Date:	RINSE BLANK RB-9954-062520-SG-001 06/25/2020	RINSE BLANK RB-9954-063020-SG-002 06/30/2020	Trip Blank TB-9954-062420-SG-001 06/24/2020	Trip Blank TB-9954-062520-SG-002 06/25/2020
Parameters	Unit				
SVOCs, TICs-Continued					
Hexadecanoic acid	μg/L	-	-	-	-
Homomenthyl salicylate	μg/L	-	-	-	-
Nonanoic acid	μg/L	-	-	-	-
Sulfur	μg/L 	-	-	-	-
Toluene	μg/L	-	-	-	-
trans-2-Methylcyclohexanol	μg/L 		-	-	-
Unknown	μg/L	17.9 J	25.2 J	-	-
Polychlorinated Biphenyls (PCBs)					
Aroclor-1016 (PCB-1016)	μg/L	0.91 U	0.91 U	_	_
Aroclor-1221 (PCB-1221)	μg/L	1.8 U	1.8 U	_	_
Aroclor-1232 (PCB-1232)	μg/L	0.91 U	0.91 U	_	_
Aroclor-1242 (PCB-1242)	μg/L	0.91 U	0.91 U	_	_
Aroclor-1248 (PCB-1248)	μg/L	0.91 U	0.91 U	-	-
Aroclor-1254 (PCB-1254)	μg/L	0.91 U	R	-	-
Aroclor-1260 (PCB-1260)	μg/L	0.91 U	R	-	-
Pesticides		0.04511	_		
4,4'-DDD	μg/L	0.045 U	R	-	-
4,4'-DDE	μg/L	0.045 U	R	-	-
4,4'-DDT	μg/L	0.045 U	R	-	-
Aldrin	μg/L	0.045 U	R	-	-
alpha-BHC	μg/L	0.079 0.045 U	R R	-	-
alpha-Chlordane beta-BHC	μg/L	0.045 U 0.045 U	R R	-	-
delta-BHC	μg/L	0.045 U 0.044 J	0.037 J	-	-
Dieldrin	μg/L μg/L	0.044 J 0.045 U	0.037 J R	-	-
Endosulfan I	μg/L μg/L	0.045 U	R	-	-
Endosulfan II	μg/L μg/L	0.045 U	R	-	-
Endosulfan sulfate	μg/L μg/L	0.045 U	R	- -	<u> </u>
Endrin	μg/L	0.045 U	R	-	_
Endrin ketone	μg/L μg/L	0.045 U	R	-	<u>-</u>
gamma-BHC (lindane)	μg/L	0.075	0.058 J	_	_
gamma-Chlordane	μg/L	0.045 U	R	_	_
Heptachlor	μg/L	0.045 U	R	-	-
Heptachlor epoxide	μg/L	0.045 U	R	-	-
Methoxychlor	μg/L	0.045 U	R	_	-
Toxaphene	μg/L	0.46 U	R	-	-
•	r-9· -				

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	Location ID: Sample Name: Sample Date:	RINSE BLANK RB-9954-062520-SG-001 06/25/2020	RINSE BLANK RB-9954-063020-SG-002 06/30/2020	Trip Blank TB-9954-062420-SG-001 06/24/2020	Trip Blank TB-9954-062520-SG-002 06/25/2020
Parameters	Unit				
Field Parameters	Dog C				
Temperature, field Conductivity, field	Deg C mS/cm	- -	- -	- -	- -
Turbidity, field	NTU	-	-	-	-
pH, field	s.u.	-	-	-	-

Table 2 Page 61 of 66

	Location ID: Sample Name: Sample Date:	Trip Blank TB-9954-062620-RM-003 06/26/2020	Trip Blank TB-9954-063020-SG-004 06/30/2020	Trip Blank TB-9954-070120-SG-005 07/01/2020	Trip Blank TB-9954-070620-SG-006 07/06/2020	Trip Blank TB-9954-070920-SG-007 07/09/2020
Parameters	Unit					
Volatile Organic Compounds (VOCs)						
1,1,1-Trichloroethane	μg/L	5.0 U				
1,1,2,2-Tetrachloroethane	μg/L	5.0 U				
1,1,2-Trichloroethane	μg/L	5.0 U				
1,1-Dichloroethane	μg/L	5.0 U				
1,1-Dichloroethene	μg/L	5.0 U				
1,2-Dichloroethane	μg/L	5.0 U				
1,2-Dichloropropane	μg/L	5.0 U				
2-Butanone (Methyl ethyl ketone) (MEK)	μg/L	10 U				
2-Hexanone	μg/L	10 U				
4-Methyl-2-pentanone (Methyl isobutyl ketone) (Methyl isobutyl ketone)		10 U				
Acetone	μg/L	10 U				
Benzene	μg/L	5.0 U				
Bromodichloromethane	μg/L	5.0 U				
Bromoform	μg/L	5.0 U				
Bromomethane (Methyl bromide)	μg/L	5.0 U				
Carbon disulfide	μg/L	10 U				
Carbon tetrachloride	μg/L	5.0 U				
Chlorobenzene	μg/L	5.0 U				
Chloroethane	μg/L	5.0 U				
Chloroform (Trichloromethane)	μg/L	5.0 U				
Chloromethane (Methyl chloride)	μg/L	5.0 U				
cis-1,2-Dichloroethene	μg/L	5.0 U				
cis-1,3-Dichloropropene	μg/L	5.0 U				
Dibromochloromethane	μg/L	5.0 U				
Ethylbenzene	μg/L	5.0 U				
Methylene chloride	μg/L	5.0 U				
Styrene	μg/L	5.0 U				
Tetrachloroethene	μg/L	5.0 U				
Toluene	μg/L	5.0 U	0.30 J	5.0 U	5.0 U	5.0 U
trans-1,2-Dichloroethene	μg/L	5.0 U				
trans-1,3-Dichloropropene	μg/L	5.0 U				
Trichloroethene	μg/L	5.0 U				
Vinyl acetate	μg/L	10 U				
Vinyl chloride	μg/L	5.0 U				
Xylenes (total)	μg/L	5.0 U				

Sample Na	Location ID: Sample Name: Sample Date:		Trip Blank TB-9954-063020-SG-004 06/30/2020	Trip Blank TB-9954-070120-SG-005 07/01/2020	Trip Blank TB-9954-070620-SG-006 07/06/2020	Trip Blank TB-9954-070920-SG-007 07/09/2020
Parameters	Unit					
VOCs, Tentatively Identified Compounds (TICs)						
1,2,4-Trichlorobenzene	μg/L	-	-	-	-	-
1-Chloro-2-methyl-benzene	μg/L	-	-	-	-	-
1-Chloro-3-methylbenzene	μg/L	-	-	-	-	-
2,6-Dichlorotoluene	μg/L	-	-	-	-	-
2-Methylbutane	μg/L	-	-	-	-	-
6-Methyl-5-hepten-2-one	μg/L	-	-	-	-	-
Dimethyl disulfide	μg/L	-	-	-	-	-
Dimethyl sulfide	μg/L	-	-	-	-	-
Dimethyl trisulfide	μg/L	-	-	-	-	-
Hexanal	μg/L	-	-	-	-	-
Isobutane	μg/L	-	-	-	-	-
m-Monochlorobenzotrifluoride	μg/L	-	-	-	-	-
Methanethiol	μg/L	-	.	-	-	-
Methoxytrimethyl-silane	μg/L	-	11.9 J	-	-	-
Methylthioethane	μg/L	-	-	-	-	-
Nonanal	μg/L	-	-	-	-	-
Propane	μg/L	-	-	-	-	-
Sulfur dioxide (SO2)	μg/L	-	-	-	-	-
Trimethylfluorosilane	μg/L	-	-	-	-	-
Unknown	μg/L	-	17.3 J	-	-	-
Semi-volatile Organic Compounds (SVOCs)						
1,2,4-Trichlorobenzene	μg/L	-	_	-	-	-
1,2-Dichlorobenzene	μg/L	-	-	-	-	-
1,3-Dichlorobenzene	μg/L	_	_	-	-	-
1,4-Dichlorobenzene	μg/L	-	-	-	-	-
2,2'-Oxybis(1-chloropropane) (bis(2-Chloroisopropyl) ether)	μg/L	-	-	-	-	-
2,4,5-Trichlorophenol	μg/L	-	-	-	-	-
2,4,6-Trichlorophenol	μg/L	-	-	-	-	-
2,4-Dichlorophenol	μg/L	-	-	-	-	-
2,4-Dimethylphenol	μg/L	-	-	-	-	-
2,4-Dinitrophenol	μg/L	-	-	-	-	-
2,4-Dinitrotoluene	μg/L	-	-	-	-	-
2,6-Dinitrotoluene	μg/L	-	-	-	-	-
2-Chloronaphthalene	μg/L	-	-	-	-	-
2-Chlorophenol	μg/L	-	-	-	-	-
2-Methylnaphthalene	μg/L	-	-	-	-	-
2-Methylphenol	μg/L	-	-	-	-	-

	Location ID: Sample Name: Sample Date:	Trip Blank TB-9954-062620-RM-003 06/26/2020	Trip Blank TB-9954-063020-SG-004 06/30/2020	Trip Blank TB-9954-070120-SG-005 07/01/2020	Trip Blank TB-9954-070620-SG-006 07/06/2020	Trip Blank TB-9954-070920-SG-007 07/09/2020
Parameters	Unit					
SVOCs-Continued						
2-Nitroaniline	μg/L	-	-	-	-	-
2-Nitrophenol	μg/L	-	-	-	-	-
3&4-Methylphenol	μg/L	-	-	-	-	-
3,3'-Dichlorobenzidine	μg/L	-	-	-	-	-
3-Nitroaniline	μg/L	-	-	-	-	-
4,6-Dinitro-2-methylphenol	μg/L	-	-	-	-	-
4-Bromophenyl phenyl ether	μg/L	-	-	-	-	-
4-Chloro-3-methylphenol	μg/L	-	-	-	-	-
4-Chloroaniline	μg/L	-	-	-	-	-
4-Chlorophenyl phenyl ether	μg/L	-	-	-	-	-
4-Nitroaniline	μg/L	-	-	-	-	-
4-Nitrophenol	μg/L	-	-	-	-	-
Acenaphthene	μg/L	-	-	-	-	-
Acenaphthylene	μg/L	-	-	-	-	-
Anthracene	μg/L	-	-	-	-	-
Benzo(a)anthracene	μg/L	-	-	-	-	-
Benzo(a)pyrene	μg/L	-	-	-	-	-
Benzo(b)fluoranthene	μg/L	-	-	-	-	-
Benzo(g,h,i)perylene	μg/L 	-	-	-	-	-
Benzo(k)fluoranthene	μg/L 	-	-	-	-	-
Benzoic acid	μg/L	-	-	-	-	-
Benzyl alcohol	μg/L	-	-	-	-	-
bis(2-Chloroethoxy)methane	μg/L	-	-	-	-	-
bis(2-Chloroethyl)ether	μg/L	-	-	-	-	-
bis(2-Ethylhexyl)phthalate (DEHP)	μg/L	-	-	-	-	-
Butyl benzylphthalate (BBP)	μg/L	-	-	-	-	-
Chrysene	μg/L	-	-	-	-	-
Di-n-butylphthalate (DBP)	μg/L	-	-	-	-	-
Di-n-octyl phthalate (DnOP) Dibenz(a,h)anthracene	μg/L	-	-	-	-	-
Dibenzofuran	μg/L	-	-	-	-	-
	μg/L	-	-	-	-	-
Diethyl phthalate Dimethyl phthalate	μg/L	-	-	-	-	-
Fluoranthene	μg/L	-	-	-	-	-
Fluoranthene Fluorene	μg/L	-	-	-	-	-
Hexachlorobenzene	μg/L	-	-	-	-	-
Hexachlorobutadiene	μg/L	-	-	-	-	-
Hexachlorocyclopentadiene	μg/L	- -	- -	-	<u>-</u>	-
r iezacinorocyclopentaulene	μg/L	-	-	-	-	-

	Location ID: Sample Name: Sample Date:	Trip Blank TB-9954-062620-RM-003 06/26/2020	Trip Blank TB-9954-063020-SG-004 06/30/2020	Trip Blank TB-9954-070120-SG-005 07/01/2020	Trip Blank TB-9954-070620-SG-006 07/06/2020	Trip Blank TB-9954-070920-SG-007 07/09/2020
Parameters	Unit					
SVOCs-Continued						
Hexachloroethane	μg/L	-	-	-	-	-
Indeno(1,2,3-cd)pyrene	μg/L	-	-	-	-	-
Isophorone	μg/L	-	-	-	-	-
N-Nitrosodi-n-propylamine	μg/L	-	-	-	-	-
N-Nitrosodiphenylamine	μg/L	-	-	-	-	-
Naphthalene	μg/L	-	-	-	-	-
Nitrobenzene	μg/L	-	-	-	-	-
Pentachlorophenol	μg/L	-	-	-	-	-
Phenanthrene	μg/L	-	-	-	-	-
Phenol	μg/L	-	-	-	-	-
Pyrene	μg/L	-	-	-	-	-
SVOC, TICs						
(4-Methylphenyl)phenylmethanone	μg/L	_	-	_	-	_
1,4-Dioxane	μg/L	_	-	_	-	_
1-Chloro-2-methyl-benzene	μg/L	-	-	-	-	_
1-Chloro-3-methylbenzene	µg/L	-	-	-	-	-
1-Methyl-2-(phenylmethyl)-benzene	μg/L	-	-	-	-	-
2,3-Dichlorobenzoic acid	μg/L	-	-	-	-	-
2,5-Dichlorobenzoic acid	μg/L	-	-	-	-	-
2-Ethyl-ethanoic acid	μg/L	-	-	-	-	-
2-Propyl tridecyl-sulfurous acid	μg/L	-	-	-	-	-
3,4-Dichlorophenol	μg/L	-	-	-	-	-
3,5-Dichlorophenol	μg/L	-	-	-	-	-
3-Benzoylbenzoic acid	μg/L	-	-	-	-	-
3-Chlorobenzoic acid	μg/L	-	-	-	-	-
3-Methyl-cyclohexanol	μg/L	-	-	-	-	-
4-Benzoyl-(rel)-benzoic acid	μg/L	-	-	-	-	-
4-Chlorobenzoic acid	μg/L	-	-	-	-	-
4-Chlorophenol	μg/L	-	-	-	-	-
4-Chlorotoluene	μg/L	-	-	-	-	-
6-Octadecenoic acid	μg/L	-	-	-	-	-
Bicyclo[2.2.1]heptan-2-one, 1,3,3-trimethyl	μg/L	-	-	-	-	-
Chlorobenzene	μg/L	-	-	-	-	-
Cyclohexanol	μg/L	-	-	-	-	-
Dimethyl disulfide	μg/L	-	-	-	-	-
Dimethyl tetrasulfide	μg/L	-	-	-	-	-
Dodecanoic acid	μg/L	-	-	-	-	-

	Location ID: Sample Name: Sample Date:	Trip Blank TB-9954-062620-RM-003 06/26/2020	Trip Blank TB-9954-063020-SG-004 06/30/2020	Trip Blank TB-9954-070120-SG-005 07/01/2020	Trip Blank TB-9954-070620-SG-006 07/06/2020	Trip Blank TB-9954-070920-SG-007 07/09/2020
Parameters	Unit					
SVOCs, TICs-Continued						
Hexadecanoic acid	μg/L	-	-	-	-	-
Homomenthyl salicylate	μg/L	-	-	-	-	-
Nonanoic acid	μg/L	-	-	-	-	-
Sulfur	μg/L	-	-	-	-	-
Toluene	μg/L	-	-	-	-	-
trans-2-Methylcyclohexanol	μg/L	-	-	-	-	-
Unknown	μg/L	-	-	-	-	-
Debughteringtod Binhanda (BODs)						
Polychlorinated Biphenyls (PCBs) Aroclor-1016 (PCB-1016)						
Aroclor-1016 (PCB-1016) Aroclor-1221 (PCB-1221)	μg/L	-	-	-	-	-
Aroclor-1221 (FCB-1221) Aroclor-1232 (PCB-1232)	μg/L	-	-	-	-	-
Aroclor-1232 (PCB-1232) Aroclor-1242 (PCB-1242)	μg/L μg/L	-	-	-	-	-
Aroclor-1242 (FCB-1242) Aroclor-1248 (PCB-1248)	μg/L μg/L	<u>-</u>	<u>-</u>	<u>-</u>	<u>-</u>	-
Aroclor-1246 (FCB-1246) Aroclor-1254 (PCB-1254)	μg/L	- -	- -		- -	- -
Aroclor-1254 (FGB-1254) Aroclor-1260 (PCB-1260)	μg/L	_	<u> </u>		_	<u>-</u>
A10Cl01-1200 (1 CB-1200)	µg/L	<u>-</u>	_	<u>-</u>	-	-
Pesticides						
4,4'-DDD	μg/L	-	-	-	-	-
4,4'-DDE	μg/L	-	-	-	-	-
4,4'-DDT	μg/L	-	-	-	-	-
Aldrin	μg/L	-	-	-	-	-
alpha-BHC	μg/L	-	-	-	-	-
alpha-Chlordane	μg/L	-	-	-	-	-
beta-BHC	μg/L	-	-	-	-	-
delta-BHC	μg/L	-	-	-	-	-
Dieldrin	μg/L	-	-	-	-	-
Endosulfan I	μg/L	-	-	-	-	-
Endosulfan II	μg/L	-	-	-	-	-
Endosulfan sulfate	μg/L	-	-	-	-	-
Endrin	μg/L	-	-	-	-	-
Endrin ketone	μg/L	-	-	-	-	-
gamma-BHC (lindane)	μg/L	-	-	-	-	-
gamma-Chlordane	μg/L	-	-	-	-	-
Heptachlor	μg/L	-	-	-	-	-
Heptachlor epoxide	μg/L	-	-	-	-	-
Methoxychlor	μg/L	-	-	-	-	-
Toxaphene	μg/L	-	-	-	-	-

Table 2 Page 66 of 66

Analytical Results Summary Love Canal Annual Long-Term Monitoring Program Glenn Springs Holdings, Inc. Niagara Falls, New York June-July 2020

	Location ID: Sample Name: Sample Date:	Trip Blank TB-9954-062620-RM-003 06/26/2020	Trip Blank TB-9954-063020-SG-004 06/30/2020	Trip Blank TB-9954-070120-SG-005 07/01/2020	Trip Blank TB-9954-070620-SG-006 07/06/2020	Trip Blank TB-9954-070920-SG-007 07/09/2020
Parameters	Unit					
Field Parameters						
Temperature, field	Deg C	-	_	-	-	-
Conductivity, field	mS/cm	-	-	-	-	-
Turbidity, field	NTU	<u>-</u>	_	-	-	-
pH, field	s.u.	-	-	-	-	-

Notes:

NJ

J - Estimated concentration

- Presumptively present at estimated value JN Not detected at the associated reporting limit
 Not detected; associated reporting limit is estimated U ŨJ

- Tentatively identified compound, estimated concentration

R - Rejected

- Not applicable

Table 3

Analytical Methods Love Canal Annual Long-Term Monitoring Program Glenn Springs Holdings, Inc. Niagara Falls, New York June-July 2020

			Holding Time		
B			Collection to Extraction	Collection or Extraction to Analysis	
Parameter	Method	Matrix	(Days)	(Days)	
Volatile Organic Compounds (VOCs)	SW-846 8260B	Water	-	14	
Semi-volatile Organic Compounds (SVOCs)	SW-846 8270C	Water	7	40	
Polychlorinated Biphenyls (PCBs)	SW-846 8082	Water	7	40	
Organochlorine Pesticides	SW-846 8081A	Water	7	40	

Notes:

- Not applicable

Method Reference:

SW-846 - "Test Methods for Evaluating Solid Waste, Physical/Chemical Methods", SW-846, Third Edition, 1986, with subsequent revisions

Table 4

Qualified Sample Results Due to Outlying Continuing Calibration Results Love Canal Annual Long-Term Monitoring Program Glenn Springs Holdings, Inc. Niagara Falls, New York June-July 2020

Parameter	Analyte	Calibration Date	%D	Associated Sample ID	Qualified Result	Units
		(mm/dd/yyyy)				
VOCs	Vinyl acetate	06/30/2020	44	WG-9954-062420-SG-001	10 UJ	μg/L
				WG-9954-062420-SG-005	10 UJ	μg/L
				WG-9954-062420-SG-006	10 UJ	μg/L
				WG-9954-062420-SG-007	10 UJ	μg/L
				WG-9954-062620-RM-014	10 UJ	μg/L
				WG-9954-062620-RM-015	10 UJ	μg/L
				WG-9954-062620-RM-017	10 UJ	μg/L
				WG-9954-062520-SG-008	10 UJ	μg/L
SVOCs	bis(2-Chloroethyl)ether	07/06/2020	35.7	WG-9954-062620-RM-014	9.1 UJ	μg/L
				WG-9954-062620-RM-015	9.1 UJ	μg/L
				WG-9954-062620-RM-016	9.1 UJ	μg/L
				WG-9954-062620-RM-017	9.1 UJ	μg/L
				WG-9954-062620-RM-018	9.1 UJ	μg/L

Notes:

%D - Percent difference

UJ - Not detected; associated reporting limit is estimated

VOCs - Volatile Organic Compounds SVOCs - Semi-volatile Organic Compounds

Table 5

Qualified Sample Results Due to Analyte Concentrations in the Method Blanks Love Canal Annual Long-Term Monitoring Program Glenn Springs Holdings, Inc. Niagara Falls, New York June-July 2020

Parameter	Analyte	Analysis Date (mm/dd/yyyy)	Blank Result *	Sample ID	Original Result	Qualified Result	Units
VOCs	Toluene	07/14/2020	0.41 J	WG-9954-063020-RM-023 WG-9954-063020-SG-019 WG-9954-063020-SG-020 WG-9954-063020-SG-021 WG-9954-063020-SG-022 WG-9954-070120-SG-024 WG-9954-070120-SG-025	0.31 J 0.34 J 0.35 J 0.31 J 0.38 J 0.26 J 0.27 J	5.0 U 5.0 U 5.0 U 5.0 U 5.0 U 5.0 U	µg/L µg/L µg/L µg/L µg/L
				WG-9954-070120-SG-025 WG-9954-070120-SG-026	0.27 J 0.26 J	5.0 U 5.0 U	μg/L μg/L

Notes:

* - Blank result adjusted for sample factors where applicable

U - Not detected at the associated reporting limit

J - Estimated concentration

VOCs - Volatile Organic Compounds

9.1 UJ

μg/L

Hexachlorocyclopentadiene

Table 6

Qualified Sample Data Due to Outlying of Surrogate Recoveries Love Canal Annual Long-Term Monitoring Program Glenn Springs Holdings, Inc. Niagara Falls, New York June-July 2020

Parameter	Sample ID	Surrogate	Surrogate % Recovery	Control Limits % Recovery	Analyte	Qualified Result	Units
SVOCs	WG-9954-062420-SG-002	2,4,6-Tribromophenol	27	35 - 141	1,2,4-Trichlorobenzene	9.1 UJ	μg/L
		2-Fluorobiphenyl	26	31 - 118	1,2-Dichlorobenzene	9.1 UJ	μg/L
		Nitrobenzene-D5	22	31 - 110	1,3-Dichlorobenzene	9.1 UJ	μg/L
					1,4-Dichlorobenzene	9.1 UJ	μg/L
					2,2'-Oxybis(1-chloropropane)	9.1 UJ	μg/L
					2,4-Dinitrotoluene	9.1 UJ	μg/L
					2,6-Dinitrotoluene	9.1 UJ	μg/L
					2-Chloronaphthalene	9.1 UJ	μg/L
					2-Methylnaphthalene	9.1 UJ	μg/L
					2-Nitroaniline	9.1 UJ	μg/L
					3,3'-Dichlorobenzidine	9.1 UJ	μg/L
					3-Nitroaniline	9.1 UJ	μg/L
					4-Bromophenyl phenyl ether	9.1 UJ	μg/L
					4-Chloroaniline	9.1 UJ	μg/L
					4-Chlorophenyl phenyl ether	9.1 UJ	μg/L
					4-Nitroaniline	9.1 UJ	μg/L
					Acenaphthene	9.1 UJ	μg/L
					Acenaphthylene	9.1 UJ	μg/L
					Anthracene	9.1 UJ	μg/L
					Benzo(a)anthracene	9.1 UJ	μg/L
					Benzo(a)pyrene	9.1 UJ	μg/L
					Benzo(b)fluoranthene	9.1 UJ	μg/L
					Benzo(g,h,i)perylene	9.1 UJ	μg/L
					Benzo(k)fluoranthene	9.1 UJ	μg/L
					Benzoic acid	91 UJ	μg/L
					bis(2-Chloroethoxy)methane	9.1 UJ	μg/L
					bis(2-Chloroethyl)ether	9.1 UJ	μg/L
					bis(2-Ethylhexyl)phthalate (DEHP)	9.1 UJ	μg/L
					Butyl benzylphthalate (BBP)	9.1 UJ	μg/L
					Chrysene	9.1 UJ	μg/L
					Dibenz(a,h)anthracene	9.1 UJ	μg/L
					Dibenzofuran	9.1 UJ	μg/L
					Diethyl phthalate	9.1 UJ	μg/L
					Dimethyl phthalate	9.1 UJ	μg/L
					Di-n-butylphthalate (DBP)	9.1 UJ	μg/L
					Di-n-octyl phthalate (DnOP)	9.1 UJ	μg/L
					Fluoranthene	9.1 UJ	μg/L
					Fluorene	9.1 UJ	μg/L
					Hexachlorobenzene	9.1 UJ	μg/L
					Hexachlorobutadiene	9.1 UJ	μg/L
					Hayaahlaraayalanantadiana	0.1.111	ug/l

Table 6

Qualified Sample Data Due to Outlying of Surrogate Recoveries Love Canal Annual Long-Term Monitoring Program Glenn Springs Holdings, Inc. Niagara Falls, New York June-July 2020

Parameter	Sample ID	Surrogate	Surrogate % Recovery	Control Limits % Recovery	Analyte	Qualified Result	Units
SVOCs	WG-9954-062420-SG-002	2,4,6-Tribromophenol	27	35 - 141	Hexachloroethane	9.1 UJ	μg/L
		2-Fluorobiphenyl	26	31 - 118	Indeno(1,2,3-cd)pyrene	9.1 UJ	μg/L
		Nitrobenzene-D5	22	31 - 110	Isophorone	9.1 UJ	μg/L
					Naphthalene	9.1 UJ	μg/L
					Nitrobenzene	9.1 UJ	μg/L
					N-Nitrosodi-n-propylamine	9.1 UJ	μg/L
					N-Nitrosodiphenylamine	9.1 UJ	μg/L
					Phenanthrene	9.1 UJ	μg/L
					Pyrene	9.1 UJ	μg/L
Pesticides	WG-9954-062620-RM-018	(PCB 209) Decachlorobiphenyl	3	10 - 164	4,4'-DDD	R	μg/L
		, , ,			4,4'-DDE	0.029 J	μg/L
					4,4'-DDT	R	μg/L
					Aldrin	R	μg/L
					alpha-BHC	R	μg/L
					alpha-Chlordane	R	μg/L
					beta-BHC	0.045 J	μg/L
					delta-BHC	R	μg/L
					Dieldrin	R	μg/L
					Endosulfan I	0.037 J	μg/L
					Endosulfan II	R	μg/L
					Endosulfan sulfate	R	μg/L
					Endrin	R	μg/L
					Endrin ketone	R	μg/L
					gamma-BHC (lindane)	0.026 J	μg/L
					gamma-Chlordane	0.020 NJ	μg/L
					Heptachlor	R	μg/L
					Heptachlor epoxide	R	μg/L
					Methoxychlor	R	μg/L
					Toxaphene	R	μg/L

Notes:

Estimated concentration

UJ - Not detected; associated reporting limit is estimated NJ - Tentatively identified compound, estimated concentration

R - Rejected

SVOCs - Semi-volatile Organic Compounds

Table 7

Qualified Sample Results Due to Outlying LCS/LCSD Results Love Canal Annual Long-Term Monitoring Program Glenn Springs Holdings, Inc. Niagara Falls, New York June-July 2020

Parameter	Analyte	LCS Date (mm/dd/yyyy)	LCS % Recovery	LCSD % Recovery	RPD (percent)	Control Lin % Recovery	nits RPD	Associated Sample ID	Qualified Result	Units
SVOCs	2,4-Dimethylphenol	07/01/2020	82	56	38	59 - 113	30	WG-9954-062420-SG-006 WG-9954-062420-SG-007 WG-9954-062520-SG-008 WG-9954-062520-SG-009 WG-9954-062520-SG-010 WG-9954-062520-SG-011 WG-9954-062520-SG-012 WG-9954-062520-SG-013	9.1 UJ 9.1 UJ 9.1 UJ 9.1 UJ 9.1 UJ 9.1 UJ 9.1 UJ 9.1 UJ	μg/L μg/L μg/L μg/L μg/L μg/L μg/L
SVOCs	4-Chloro-3-methylphenol	07/01/2020	83	51	48	52 - 113	30	WG-9954-062420-SG-006 WG-9954-062420-SG-007 WG-9954-062520-SG-008 WG-9954-062520-SG-009 WG-9954-062520-SG-010 WG-9954-062520-SG-011 WG-9954-062520-SG-012 WG-9954-062520-SG-013	9.1 UJ 9.1 UJ 9.1 UJ 9.1 UJ 9.1 UJ 9.1 UJ 9.1 UJ 9.1 UJ	μg/L μg/L μg/L μg/L μg/L μg/L μg/L
SVOCs	4-Nitroaniline	07/01/2020	79	53	39	54 - 133	30	WG-9954-062420-SG-006 WG-9954-062420-SG-007 WG-9954-062520-SG-008 WG-9954-062520-SG-009 WG-9954-062520-SG-010 WG-9954-062520-SG-011 WG-9954-062520-SG-012 WG-9954-062520-SG-013	9.1 UJ 9.1 UJ 9.1 UJ 9.1 UJ 9.1 UJ 9.1 UJ 9.1 UJ 9.1 UJ	μg/L μg/L μg/L μg/L μg/L μg/L μg/L
SVOCs	Isophorone	07/01/2020	76	46	49	50 - 116	30	WG-9954-062420-SG-006 WG-9954-062420-SG-007 WG-9954-062520-SG-008 WG-9954-062520-SG-009 WG-9954-062520-SG-010 WG-9954-062520-SG-011 WG-9954-062520-SG-012 WG-9954-062520-SG-013	9.1 UJ 9.1 UJ 9.1 UJ 9.1 UJ 9.1 UJ 9.1 UJ 9.1 UJ 9.1 UJ	μg/L μg/L μg/L μg/L μg/L μg/L μg/L μg/L

Notes:

LCS - Laboratory Control Sample
LCSD - Laboratory Control Sample Duplicate
RPD - Relative Percent Difference
- Not detected: associated

UJ - Not detected; associated reporting limit is estimated SVOCs - Semi-volatile Organic Compounds

Table 8

Qualified Sample Results Due to Outlying MS/MSD Results Love Canal Annual Long-Term Monitoring Program Glenn Springs Holdings, Inc. Niagara Falls, New York June-July 2020

			MS	MSD		Control Li				
Parameter	Sample ID	Analyte	% Recovery	% Recovery	RPD (percent)	% Recovery	RPD	Result	Units	
SVOCs	WG-9954-070120-SG-027	3,3'-Dichlorobenzidine	3	4	29	11 - 131	30	R	μg/L	
SVOCs	WG-9954-070120-SG-027	Isophorone	6	7	15	40 - 111	30	R	μg/L	
SVOCs	WG-9954-070720-SG-030	Hexachlorocyclopentadiene	6	5	18	10 - 103	30	R	μg/L	

Notes:

MS - Matrix Spike

MSD - Matrix Spike Duplicate RPD - Relative Percent Difference

R - Rejected

SVOCs - Semi-volatile Organic Compounds

Table 9

Qualified Sample Data Due to Analyte Concentrations in the Rinse Blanks Love Canal Annual Long-Term Monitoring Program Glenn Springs Holdings, Inc. Niagara Falls, New York June-July 2020

Parameter	Rinse Blank ID	Blank Date (dd/mm/yyyy)	Analyte	Blank Result	Associated Sample ID	Original Result	Qualified Result	Units
SVOCs	RB-9954-063020-SG-002	06/30/2020	Phenol	0.94 J	WG-9954-063020-SG-021	1.1 J	9.1 U	μg/L

Notes:

U - Not detected at the associated reporting limit

J - Estimated concentration

SVOCs - Semi-volatile Organic Compounds

Table 10

Qualified Sample Data Due to Variability in Field Duplicate Results Love Canal Annual Long-Term Monitoring Program Glenn Springs Holdings, Inc. Niagara Falls, New York June-July 2020

Parameter	Analyte	RPD	Sample ID	Qualified Result	Field Duplicate Sample ID	Qualified Result	Units
Pesticides	delta-BHC	55	WG-9954-063020-SG-020	0.068 J	WG-9954-063020-SG-021	0.12 J	μg/L

Notes:

RPD - Relative Percent Difference
J - Estimated concentration

Appendix G Niagara Falls Water Board Annual Inspection Letter



NIAGARA FALLS WATER BOARD

ANNUAL PLANT INSPECTION

INDUSTRIAL PRETREATMENT PROGRAM

PAGE 1 OF 6

Name and Address of SIU

Glenn Springs Holding Inc. Love Canal Leachate Treatment Facility 805 – 97th Street

Niagara Falls, NY 14304

Permit Number: 44 SIC Codes: 4951

Date of Last Inspection: 4/11/2019

CATEGORICAL IU? NO

Day/Date and Time of Inspection: Friday September 4th, 2020 @ 11:00AM

SIU Representative:

Darrell Crockett

Inspectors Name:

Stephen Stewart

Contact Phone No.: 998-5804

PART I FLOW RECORDING AND SAMPLING INSTRUMENTATION

calibration): **Quarterly – August 31**st, **2020**

a)	Flow measure	ement instrument m	eets permit requirements?	YES
b)	Primary flow r	measurement device	e properly installed?	YES
c)	Type of flow r	measurement device	e	
	Weir []	Flume(s)[]	Water meter []	Mag meter [X]
d)	Does device	measure flow adequ	uately?	YES
e)	Is primary me maintained? [easuring device prop [1960.6]	perly operated and	YES
f)		ry instruments (recor ated and maintained	,	YES
g)	Calibration fre	equency adequate?	(date of last	YES

PART II SAMPLE COLLECTION

a)		permit require SIU to submit Periodic Nonitoring Reports? [40 CFR403.12h]	YES
b)	-	s", does the sample collection frequency and ant type conform with permit requirements?	YES
c)	the pe	ne sample collection locations as described in ermit adequate for representative sample stion? [1960.6 (a)]	YES
d)	permi	the method of sample collection conform with t requirements, Sewer Use Ordinance and Federal dards?	YES
	i)	Sample refrigerated throughout collection and storage?	YES
	ii)	Are samples properly preserved?	YES
	iii)	Are samples collected using flow proportion composite or grab sampling where appropriate ? [40 CFR12(b)(5)(iii)].	YES
	iv)	Sample holding times appropriate? [40 CFR136.3]	YES
PART	III LA	ABORATORY FACILITIES	
a)	ls a c	ommercial laboratory used?	YES
	i) ii)	Name of laboratory: ALS Environmental	
	ii)	Address: 1565 Jefferson Rd. Rochester, NY 14623	
	iii)	Is laboratory State certified?	YES
b)	Does	SIU perform its own analysis?	NO
	i)	Is the SIU's laboratory State certified?	N/A
	ii)	Are your laboratory wastes properly disposed of?	N/A
c)	Are E	PA approved testing methods used?	YES

PAGE

OF

6

3

e)	Has this facility been evaluated <i>OR</i> re-evaluated for its' potential to experience a slug discharge? 9/4/2020 1) Is a Slug Control Plan required for this facility [40 CFR403.8(f)(2)(vi)]?	YES NO
	A. Has the facility <u>Developed</u> and <u>Implemented</u> a Slug Control Plan?	N/A
	 a. The date of the plan's last update: N/A b. Is the latest update on file at the NFWB? c. Does it contain the correct Water Board 	N/A
	phone numbers and extensions.	N/A
	2) Has the facility experienced a slug discharge since The last inspection?	NO
f)	Have there been any significant manufacturing or process changes? [1960.5 (c)] List: <i>None</i>	NO
	Who was contacted prior to implementation of these changes? **N/A** Date: N/A**	
g)	Describe your hazardous waste storage area(s). Double contained in the decontamination/storage facility.	
	Do they meet DEC & EPA containment requirements?	YES
	Are all containers correctly labeled and time limits adhered to?	YES
	Describe your method of disposal:	
	Incineration as needed through Veolia Services. at their Port Arthur Texas. Approximately once per quarter.	
h)	Regarding the blueprints that you submitted with your last permit application, have there been any significant changes made to your process or sewer lines?	NO
	Have revised blueprints been sent to the WWTP?	N/A

PART VI PRETREATMENT

a)	•		equired pretrea Carbon treati		/WTP	•	
b)	Are all p	retreatment f	acilities proper	ly maintai	ned?		YES
c)		ny pH probes contain? <i>0</i>	does your pH	monitorir	ıg		
	List the f	requency for	calibration.				
d)	un-permi	tted waste or	is anyone disc waste not prop ? [40 CFR 403	erly pretr		d	NO
e)	Were WW	/TP personal	notified?				N/A
	Prior to	discharge to	sewer?				N/A
	During	or after disch	arge?				N/A
	Who?	N/A	Date:	N/	Ά	Time:	N/A
	SIU persor	nal who conta	cted WWTP:	N/A			
			given to the W		nin		
	[40 CFR 40						N/A

f) List any pretreatment changes that were made in the past 12 months. **None**

Sent to: N/A From: N/A Date: N/A

Who was contacted ${\bf prior}$ to implementation of these pretreatment changes? ${\it N/A}$

PART VII COMPLIANCE AND ENFORCEMENT

a)	Has the SIU had any violations since the last	
-	inspection? List: None	NO

b) If numeric violations were noted by SIU, was a repeat sample collection and analysis performed within 30 days and the results submitted to the POTW [40 CFR403.12(g)]?

N/A

c) Is SIU currently on an administrative order and/or compliance schedule?

d) If yes, have milestone dates on schedule been met? N/A

e) Was escalating enforcement action required to achieve compliance? Describe: **None Required** NO

PART VIII RECOMMENDATIONS, REQUIREMENTS AND COMMENTS:

Hazardous waste from Love Canal consists of Debris from wells, Clarifier solids, Filters, and PPE.



January 25, 2021

Mr. Darrell Crockett – Facility Manager Glenn Springs Holding Inc. Love Canal Leachate Treatment Facility 805 – 97th Street Niagara Falls, New York 14304

Dear Mr. Crockett:

Enclosed please find the results of the Plant Inspection that was conducted at your facility on Friday, September 4^{th} , 2020 @ 11:00AM. These inspections are the annual compliance inspections performed by NFWB as required by 40 CFR 403.8(f)(2)(v).

If you have any questions, I can be reached at 283-9770 ext 1701.

Sincerely,

NIAGARA FALLS WATER BOARD WASTEWATER FACILITIES

Stephen C. Stewart

Stephen C. Stewart Senior Industrial Waste Inspector

Cc: J. Paradise \rightarrow S. Stewart \rightarrow File: - I-44 Emailed to SIU

Appendix H
Test and Backflow Prevention
Device Reports

NEW YORK STATE DEPARTMENT OF HEALTH Bureau of Public Water Supply Protection Empire State Plaza - Corning Tower Room 1110 Albany, NY 12237

ГА	Please us	se a separ	ate form for each de	vice.		For the year Initia	il test - Con	nplete e	entire form Part A only
c Water Supply	y of f NI	2000	Account No.	,	Count	ragara	Block		Lot
	City- 01 14	arjanta,		Location	of Device	,			
ility Name	City of Ni Glean Spring 5-95th St.	s Rea	a Falls	<u></u>	ainte	nance	BD6		
Street	Manufacturer	i	Type RPZ RCV	Model	09M2G	Size (in	inches)		Serial Number 02-9613
rmation	Check Valve		Check Valve N		Differenti	al Pressure R	telief	Line P	ressure 73 psi
	CIRCK VALVE			<u> </u>				ate	
st fore pair	Leaked Closed tight		Leaked Closed tight] .	Opened at	. <u>2. </u> psid		O S	2 25 20 D Y
	Pressure drop across f	irst check va	1146						
								Name_	Repaired by
escribe epairs and eaterials								Lic #	
sed						•			epaired:
inal test	Closed tight		Closed tight		Opened	atps	id	Date	M D Y
	Pressure drop acros	ss first psid							
Water Meter	7,507	315	Meter Reading	1	Don	Service: (chechestic • Fi	re •	Other_	ents)
Remarks (D	escribe deficiencies; bypasse	es, cutiets befo	re the device, connections bet	veen the dev	ice and point	of entry, missing	or inadequat	e sirgaps	, 5.6.7
Aust	This device me I hereby certify the forego	eets,* Ding data to b	does NOT meet, the require correct. (3) 42 Certified Tester No.	rements of	an accepta	ble containne	ent device a	00	ne of testing 1 / 30 / 27. ation Date
Print Name Property ow		certification	that test was performed:			mature		الد	798 5864 Telephone
Print Name	Certification that instr		accordance with the appro			(To be compl supplier.)	eted by the c	iesign en	gineer or architect or water
l hereby co	ertify that this installation	is in accord	ance with the approved pla	ins.	Date				NYS DOH Log #
Name			Title		-	m	d	у	
License N			Phone ()	Des	cribe minor	installation ch		·	
Represen	ung								
Address		State	Zip		•	•			
City		عاجات ا	1						



NEW YORK STATE DEPARTMENT OF HEALTH NEW YORK STATE DEPARTMENT OF HEALTH Bureau of Public Water Supply Protection Empire State Plaza - Corning Tower Room 1110 Albany, NY 12237

RTA	Please use a	separate fo	rm for each de	vice.			l test - Co	2020 omplete ent Complete Pa	ire form
blic Water Suppl	у С А	C 11.	Account No.		County	10000	Block		Lot
	City of Ninga	ara radis		Location					
	(1. Socions	Remod	intion	Location	J. Device				
cility Name _ idress <u>&O</u>	City of Ninge Glen Springs 5-05" St N	lagala	<u>Folls</u>	Tre	street	BDG	· (^	lech	Room)
Street	Manufacturer	Туре	RPZ DCV	Model 90	19	Size (in i	nches)		rial Number
formation	Watts Check Valve No.	1	Check Valve No		Differentia	i Pressure R	elief	Line Pre	ssure 7/ psi
	CHECK VALVE IVE							Date	
est efore	Leaked X		Leaked Closed tight		Opened at	2 <u>()</u> psid		<u>0</u> 2	25 20 D Y
epair	Pressure drop across first 7.1 psid	check valve						IVI	
escribe epairs and naterials								Name	Repaired by
used								Date repa	aired:
Final test	Closed tight		Closed tight]	Opened	atpsi	ď	Date	
*	Pressure drop across fi	rst I							
Water Meter	Number 7 19 2-33	7 A		イクラー	· Dom		e •	Other	4
Remarks (C	escribe deficiencies: bypasses, co								
1	meets. I hereby certify the foregoing	data to be corre	OT meet, the requirect.	ements of a	in acceptab	le containme	nt device	at the time Ogli Expiration	301 Lile
Print Name	vners (or owners agent) ber	tification that te	st was performed:		M	Ctt		716	PB5864
Dal	1011 Usitell		Fitte		Sign	ature		<u></u>	Telephone
Print Name	Certification that installat	ion is in accorda	ance with the appro	ved plans.		(To be complising supplier.)	eted by the	design engin	eer or architect or water
l hereby c	ertify that this installation is it	accordance wi	th the approved pla	ns.	1	1			NYS DOH Log #
Name		Title			Date				N19 DON LOG #
License N	lumber	Phone	e()			m 3.	d	у	
Represen				Desc	ibe minor i	nstallation ch	anges		
Address						,			
City		State	Zip						
Signature			tinent representative a			upplier within 3	n days of th	ne testing dev	ice.

NEW YORK STATE DEPARTMENT OF HEALTH Bureau of Public Water Supply Protection Empire State Plaza - Corning Tower Room 1110 Albany, NY 12237

RT A	Please use a separa	te form for each de	vice.	ini ini	tial test - Comp	olete entire form
		A manufat NO		County	Block	Lot
olic Water Suppl	City of Niagara Fal Glenn Springs Ro G-ast St. Niagara City	Account No.		Miagara		
	My of Magard to		Location of I	Device '		
cility Name	Glenn Springs Ro	emediation				
dress 1201	5- 95+ St. Niagara	Falls 710	_Loc	ner Room		
Street	City	Type RPZ	Model	Size (i	n inches)	Serial Number
vice ormation	Manufacturer WATT 3	DCV	909	ifferential Pressure		ine Pressurepsi
	Check Valve No. 1	Check Valve N	0.2	Valve	Da	*0
		Leaked] 0,	ened at 2.1 ps	sid [022520
est efore	Leaked Closed tight	Closed tight				M D Y
epair	Pressure drop across first check val	ve				IVI 5
					١,	Repaired by
Describe repairs and					Ì	ame
materials						IC #
used					1	Date repaired:
				_		M D Y
Final test	Closed tight	Closed tight		Opened at		Date V
-	Pressure drop across first					
	check valvepsid			Type of Service: (ch	eck one)	
Water Meter	Number 31671117	Meter Reading	L L	· Domestic	Fire • C	Other
	Describe deficiencies: bypasses, outlets befor	, , , , , , , , , , , , , , , , , , ,	ween the device a	ind point of entry, missi	ng or inadequate	airgaps, etc.)
Remarks (D						
	×	ioes NOT meet, the requi	irements of an	acceptable contains	gent device at	the time of testing
Certification	n: This device meets.* I hereby certify the foregoing data to be	e correct. 13192		12	2-1	<u>09 30 10</u> Expiration Date
Print Name	10 Crossley	Certified Tester No.		Ignature /		LADITATION DE LA
1	wners of owners agent) certification t	hat test was performed:		Mont	•	711 9985864
Dam	Il Godott	Icon		Signature		Telephone
Print Name		Title	wad plans	(To be con	pleted by the de	sign engineer or architect or water
PART B	Certification that installation is in a	ccordance with the appro	oved plans.	supplier.)		
l hereby c	ertify that this installation is in accorda	nce with the approved pla	ans.	·		NYS DOH Log #
Name		Title		Date		
License N	lumber	Phone ()				у
Represen			Describ	e minor installation	changes	
Address						
City	State	Zip				
Signature	2	n department representative a			30 days of the t	esting device. DOH- 1013(9/91)

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exice Macormation	Check Valve No. 1 Leaked Closed tight	Check Valve No	Model 009	Device Incom	BDC Size (in ir			erial Number
exice Macormation	Check Valve No. 1 Leaked Closed tight	Check Valve No	Model 009	Device Incom	Size (in ir	nches)		erial Number
exice Macormation	Check Valve No. 1 Leaked Closed tight	Check Valve No	Model 009	dness m3Q	Size (in ir	nches)		erial Number
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est C pair P	Check Valve No. 1 Leaked	Check Valve No		ifferentia				61327
efore C	closed tight		i		l Pressure Re Valve		Line Property	ressure 72 psi
. P	والمستنينا بالمسا	Closed tight	1	pened at j	2,0 psid		Q 2 M	25 20 D Y
\ ≥	ressure drop across first check valves. Opposite	/e		<u></u>				
escribe epairs and naterials						1	Lic #	Repaired by
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Print Name Property owne	irs (opowners agent) certification th	Certified Tester No.	(M	ature		716	98564 Telephone
Print Name	Certification that installation is in a	Title	red plans.	O.g.		ted by the o	design eng	ineer or architect or water
	fy that this installation is in accordan		is.					NYS DOH Log #
Name		Title		Date		<u> </u>		
License Num	nber	Phone ()			m .	d 	У	
Representing			Describ	e minor i	nstallation ch	anges		
Address					,			
City	State	Zip						

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ГА	Please use a sep	evice.	For the year			orm only		
c Water Supply	() () () () () ()	Falls Account No.		County	ggara	Block		Lot
	City of Niagara	1 / (27.73	Location o	f Device				
ility Name	City of Niagara Cilena Springs ast St. Niaga	Remediation rea Falls			at B	06.	ße)	
Street	City	Zip Type X RPZ	Model		Size (in i	nches)		Number
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escribe	7,0 psid					1	lame	paired by
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Print Name	Certification that installation is	in accordance with the app			(To be comp supplier.)	leted by the d	esign engineer	or architect or water
! hereby c	ertify that this installation is in acc		lans.	Date			NYS	S DOH Log #
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	**			,55 ,,,,,,,		-		
Represer	nting		}					

2020 Cap Regra	ading Photo	Appendix I graphic Log



Photo 1 - View east of depression in access road before regrading



Photo 2 - View northeast of depression in cap before regrading



Cap Regrading Photographs – October 5-13, 2020



Photo 3 - View north of depression in cap before regrading



Photo 4 - View east of depression in access road before regrading





Photo 5 - View southeast of topsoil brought in to regrade cap



Photo 6 - View northeast of topsoil unloaded to regrade cap



Cap Regrading Photographs – October 5-13, 2020



Photo 7 - View east of access road and equipment used during regrading



Photo 8 - View north of access road after regrading





Photo 9 - View southeast of access road and cap after regrading



Photo 10 - View southeast of cap after regrading





about GHD

GHD is one of the world's leading professional services companies operating in the global markets of water, energy and resources, environment, property and buildings, and transportation. We provide engineering, environmental, and construction services to private and public sector clients.

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