# RCRA Facility Investigation 2008 Soil Vapor Investigation Report Buildings 1, 2, 3, 4, and the Tank Storage Area Sampling Event Former Hampshire Chemical Corp. Facility Waterloo, New York NYD002234763

Prepared for

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**CH2MHILL** 

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# **Acronyms and Abbreviations**

μg/L micrograms per liter

μg/m³ micrograms per cubic meter

AF attenuation factor

AOC area of concern

COPI constituent of potential interest

CSM conceptual site model

DCA dichloroethane

DCE dichloroethene

DCP dichloropropane

GC gas chromatography

HCC Hampshire Chemical Corp.

MEK methyl ethyl ketone

MIBK methyl isobutyl ketone

MPA mercaptopropionic acid

MS mass spectrometry

NaSH sodium hydrosulfide

NYSDEC New York State Department of Environmental Conservation

NYSDOH New York State Department of Health

PCE tetrachloroethene

QA quality assurance

QC quality control

RCRA Resource Conservation and Recovery Act

RFI Resource Conservation and Recovery Act facility investigation

SOP standard operating procedure

SVI soil vapor investigation

SWMU solid waste management unit

T-acid thioglycolic acid

TCA trichloroethane

TCE trichloroethene

USEPA United States Environmental Protection Agency

VOC volatile organic compound

#### **SECTION 1**

# Introduction

On April 23 and 24, 2008, CH2M HILL conducted a subslab sampling event in support of evaluating the vapor intrusion pathway in Buildings 1, 2 (2-A and 2-B), 3, 4, and the Tank Storage Area at the Evans Chemetics manufacturing facility (former Hampshire Chemical Corp. [HCC]) in Waterloo, New York. Sampling activities were conducted at the request of the New York State Department of Environmental Conservation (NYSDEC) and New York State Department of Health (NYSDOH) in a correspondence dated November 8, 2006, that requested a soil vapor investigation work plan to address potential vapor intrusion pathways at the facility (NYSDEC and New York State Department of Health [NYSDOH] 2006). The *Resource Conservation and Recovery Act* [RCRA] *Facility Investigation* [RFI], *Soil Vapor Investigation Work Plan* was submitted to the NYSDEC and NYSDOH on February 26, 2007 (CH2M HILL 2007). In a letter dated April 25, 2007, the NYSDEC and NYSDOH approved the February 2007 work plan with the condition that a revised work plan be submitted to both agencies. The revised work plan was submitted to the NYSDEC and NYSDEC and NYSDOH in June 2007 (CH2M HILL 2007).

The purpose of the subslab sampling in Buildings 1 through 4 and the Tank Storage Area was to collect data to evaluate potential vapor intrusion pathways into the buildings located either above or in close proximity to volatile organic compounds (VOCs) identified in subsurface soil and/or groundwater. During this sampling event, subslab samples were collected from Buildings 1, 2, 3, 4, and the Tank Storage Area. In addition, two outdoor ambient air samples were collected within the facility property. These samples were collected in accordance with the *Guidance for Evaluating Soil Vapor Intrusion in the State of New York* (New York State Department of Health [NYSDOH] 2006) and followed the revised June 2007 RFI, *Soil Vapor Investigation Work Plan* (CH2M HILL 2007).

#### **SECTION 2**

# Site Background

The Evans Chemetics facility is located at 228 East Main Street in the Village of Waterloo, Seneca County, New York (Figure 1). The facility is bordered to the north by East Main Street, the east by Gorham Street, the west by East Water Street, and the south by the Seneca-Cayuga Canal.

RFI efforts have been performed since 1993 to determine the nature and extent of releases to the environment. The solid waste management units (SWMUs) and areas of concern (AOCs) that have been identified for further investigation and/or corrective measures are represented on Figure 2. The RFI data have been evaluated against the applicable NYSDEC screening levels for that time for each media (groundwater, soil, sediment), and various reports summarizing the data have been submitted to NYSDEC, including the RFI report by CH2M HILL in May 2006. Comments were provided by NYSDEC and NYSDOH in a letter dated November 8, 2006. In that letter, NYSDEC requested that an SVI work plan be submitted to address potential vapor intrusion pathways at the facility. That work plan was submitted to NYSDEC on February 26, 2007. During a site visit conducted on March 12, 2007, clarifications were obtained regarding the approach to the SVI evaluation, and a revised work plan was submitted to the NYSDEC in June 2007 (CH2M HILL 2007).

Buildings evaluated in this report where there is a potential for vapor intrusion are described in detail in the following sections and comprise:

- Building 1 is located adjacent to and south of Building 2. AOC C, "Gorham Street," is adjacent and east of Building 1.
- Buildings 2-A and 2-B are located in the northeast corner of the facility. AOC E is north of these buildings.
- Building 2 is located south of Buildings 2-A and 2-B, and includes the "Mercaptopropionic Acid (MPA) Process Area."
- Building 3 is located west of Building 1 along the Seneca-Cayuga Canal. AOC D is adjacent and southwest of Building 3.
- Building 4 is adjacent to and north of Building 3. AOC B underlies Building 4 and is the former Building 4 Pit, a below-grade sump that collected wastewater from Building 4.
- The Tank Storage Area is north of Building 4.

The locations of the buildings evaluated are shown on Figure 2.

Building surveys were completed during site walkthroughs in April 2006 and January 2007. The results of the building surveys are documented and provided in Section 2.3 of the SVI revised work plan (CH2M HILL 2007). During the April 2008 subslab sampling event, a list of constituents in each building was compiled and is presented as Appendix A.

# 2.1 Buildings 1 and 2

Buildings 2, 2-A, and 2-B are in the northeast corner of the site; Building 2-B is adjacent to AOC E (Figure 2). Buildings 2-A and 2-B contain the process for manufacturing thiodipropionate esters and 2-sulfoethyl methacrylate. Building 2 contains the process for manufacture of 3-MPA and low volume specialty constituents produced on a campaign basis. A small research and development laboratory and storage areas for intermediate process constituents are located in Building 2. Building 1, the site boiler room, is in the southeast area of the facility, south of Building 2.

Historically, a release of sodium hydrosulfide (NaSH) was reported in September 1995 from a vent located on the roof of Building 2, located within the property boundary of the former HCC facility. The release of less than 100 pounds of NaSH-contaminated site soil was later excavated (CH2M HILL 2006).

AOC C is a small area east of Building 1 and the MPA Process Area (Figure 2). Contaminated soil adjacent to Building 2 in AOC C was excavated in 1999. Although the bulk of the visibly stained soil was removed, several small veins of discolored soil reportedly remained in the excavation in the direction of the building foundation. Soil samples collected at the time of the excavation reported detected concentrations of acetone, benzene, methyl isobutyl ketone (4-methyl, 2-pentanone or MIBK), and methylene chloride above associated screening levels (CH2M HILL 2006). Soil samples were collected from AOC C in August 2007 as part of the RFI addendum field activities. No VOCs were detected in these samples.

Groundwater data associated with AOC C were collected from monitoring wells MW-7 and MW-8; MW-8 had a detection of chloroform at 1.08 micrograms per liter ( $\mu$ g/L) in 2005 (Figure 6). No other VOCs were detected in these wells.

AOC E is located to the north of Building 2-B (Figure 2). Several VOCs have been detected in soil at AOC E, including MIBK, acetone, 2-hexanone, methyl ethyl ketone (MEK or 2-butanone).

Groundwater data associated with AOC E were collected from monitoring well MW-10 and temporary piezometers installed around MW-10 during previous investigations at the site (Figure 6). Monitoring well MW-10 is located at the northern end of Gorham Street in the northeast area of the facility; five VOCs have been detected in groundwater from this area (MIBK, acetone, benzene, carbon disulfide, and toluene).

Groundwater data collected from this area are summarized in the RFI addendum report, which was submitted to the NYSDEC in November 2008 (CH2M HILL 2008).

The primary vapor intrusion pathway for the MPA Process Area of Building 2 and Building 1 is VOCs migrating from soil at AOC C through the soil vadose zone and into the subslab vapor space. Another potential pathway is VOCs migrating from groundwater at AOCs C and E through the vadose zone into the subslab vapor space. For Building 2, including the MPA Process Area, the vapor intrusion pathway is probably weak because of the slightly negative pressurization of the building air space, and limited or absent preferential pathways (Section 2.3.1, CH2M HILL 2007).

# 2.2 Building 3

Building 3 is directly west of Building 1 (Figure 2). Building 3 contains the process for manufacture of thioglycolic acid (T-acid). T-acid manufacturing processes occupy half of the ground and first floors of the building. The remainder of the ground floor is used for raw material storage and in-process chemical storage. The remainder of the first floor is maintenance storage area, including the maintenance stockroom.

AOC D is a small area to the southwest of Building 3 (Figure 2). In 2004, MEK was detected in soil at AOC D. No other VOCs have been detected in soil at AOC D.

Groundwater data associated with AOC D were collected from monitoring wells MW-11S, MW-21, MW-23, and MW-24 (Figure 6). Eight VOCs have been detected in groundwater in these wells (MIBK, acetone, benzene, carbon disulfide, chloromethane, ethylbenzene, toluene, and xylenes [total]).

Groundwater data collected from this area are summarized in the RFI addendum report, which was submitted to the NYSDEC in November 2008 (CH2M HILL 2008).

The primary vapor intrusion pathway for Building 3 is likely to be VOCs conveyed into the building from soil or groundwater at AOC D. A vapor intrusion pathway may exist through the Building 3 control room floor drains (Section 2.3.2, CH2M HILL 2007).

# 2.3 Building 4

Building 4/4A is adjacent to and northeast of Building 3 (Figure 2). Building 4/4A contains the process for purification of T-acid and the manufacture of thioglycolates.

AOC B, the former Building 4 Pit, was a below-grade sump located along the southern end of Building 4, extending from the west wall nearly the length of the building. The sump was a collection point for wash water in Building 4. Liquid that collected in the sump was pumped to the facility wastewater treatment plant for treatment and discharge to the Seneca-Cayuga Canal via a permitted outfall.

In July 1994, discolored soil was encountered in the foundation excavation for a new T-acid tower in Building 4. The discolored soil triggered an investigation of the area and found the source of impacts to be the below-grade sump later designated as AOC B.

A number of VOCs have been detected in soil at AOC B, including MIBK, acetone, carbon disulfide, chloroform, toluene, and xylenes (total).

Groundwater data associated with AOC B were collected from monitoring wells, piezometers, and a sampling port: MW-01, MW-02, MW-03, MW-22, MW-23, PZ-1, PZ-4, PZ-5, PZ-6, PZ-7, BLDG4-FD, BLDG4-PW, Bldg 4 Pit Sump (Figure 6). VOCs have been detected in groundwater in these wells, including 1,1-dichloroethane (1,1-DCA); 1,2-dichloropropane (1,2-DCP); MEK; MIBK; acetone; benzene; carbon disulfide; chlorobenzene; chloroform; cis-1,2-dichloroethene (cis-1,2-DCE); ethylbenzene; methylene chloride; toluene; trans-1,2-DCE; trichloroethene (TCE); vinyl chloride; and xylenes (total). Because of the detection of VOCs in site groundwater, soil vapor sampling was conducted for this building.

Soil and groundwater data collected from this area are summarized in the RFI addendum report, which was submitted to the NYSDEC in November 2008 (CH2M HILL 2008).

The primary potential vapor intrusion pathway for Building 4 includes VOCs potentially migrating from groundwater and soil at AOC B through the soil vadose zone and into the subslab vapor space. A potential vapor intrusion pathway exists because of the proximity of the former source to the buildings and the depressurized conditions indoors (Section 2.3.3, CH2M HILL 2007).

# 2.4 Tank Storage Area

The Tank Storage Area is an enclosed area north of Building 4 (Figure 2) and to the west and below the relative grade of Building 9. Sodium hydrosulfide, glycerin, glycerol monothioglycolate, and ammonium thioglycolate tanks are within this area. In addition, the Tank Storage Area is adjacent to monitoring well MW-01 (Figure 6).

AOC B is south of the Tank Storage Area. VOCs detected in soil and groundwater at AOC B are discussed in Section 2.3.

As part of the RFI addendum field activities, PZ-3 was installed inside the Tank Storage Area (Figure 6). In December 2007, six VOCs were detected in groundwater from PZ-3 (1,2-DCA; 1,2-DCP; MIBK; acetone; carbon disulfide; and toluene) (CH2M HILL 2008). Because of the detection of VOCs in groundwater, soil vapor sampling was conducted for this area.

Groundwater data collected from this area are summarized in the RFI addendum report, which was submitted to the NYSDEC in November 2008 (CH2M HILL 2008).

A potential vapor intrusion pathway exists because of VOCs in groundwater in the vicinity and the negative pressurization of the Tank Storage Area relative to the outside air (Section 2.3.4, CH2M HILL 2007).

#### **SECTION 3**

# Vapor Intrusion Conceptual Site Model

The vapor intrusion conceptual site model (CSM) presented as Figure 3 depicts sources of constituents, migration pathways, potential receptor populations, and potential exposure pathways to the receptors.

Potential indoor air exposures in the buildings may result from VOCs in subsurface soil and/or shallow groundwater volatilizing, migrating vertically (and horizontally to a limited extent) through the soil column and entering the buildings through cracks. The VOCs may then be inhaled by building occupants. Sources of constituents potentially contributing to vapor intrusion are the VOCs detected in soil and in groundwater underneath or in close proximity to the buildings that were discussed in Section 2.

# Sampling Procedures and Methods

On April 23 and 24, 2008, 12 subslab and two ambient air samples were collected in accordance with the *Guidance for Evaluating Soil Vapor Intrusion in the State of New York* (NYSDOH 2006) and the *RFI Soil Vapor Investigation Work Plan*, revised in June 2007 (CH2M HILL 2007). The sample locations within each building and area are shown on Figures 4 and 5. The rationale for placement of these subslab probes were adjusted and agreed upon by NYSDEC during a March 2007 site visit. The placement of the subslab probes are presented for each building in the *RFI Soil Vapor Investigation Work Plan* (Section 4.1, CH2M HILL 2007).

Two ambient air samples were collected within the property boundary, and the locations are shown on Figures 4 and 5. Sampling and analysis of the subslab and ambient air samples were conducted in accordance with the work plan (CH2M HILL 2007). The field sampling log sheets are provided as Appendix B.

# 4.1 Subslab Vapor Sampling

Twelve subslab samples (SG-1 through SG-9, SG-3A, SG-5A, and SG-7A) were collected over an 8-hour period from inside Buildings 1 through 4 and the Tank Storage Area (Figure 2) as described in the standard operating procedure (SOP) (CH2M HILL 2007).

The final locations of the subslab vapor probes were based on site conditions and operations and are identified in Figures 4 and 5. Subslab vapor sampling was not conducted for at least 24 hours after probe installation was completed and the samples were collected in the manner stated in the NYSDOH guidance to obtain representative samples and to minimize possible discrepancies due to outdoor air infiltration (NYSDOH 2006). Helium was used as a tracer gas to verify the soil vapor sample was not affected by indoor air. Tracer gas readings were collected before and after sample collection.

## 4.2 Outdoor Air Sampling

Two outdoor (i.e., ambient) air samples (SG-B2 and SG-B4) were collected over an 8-hour period during daytime working hours. The ambient air samples were collected concurrently with the subslab soil vapor samples (Tables 1 and 2). Sample SG-B2 was collected outside Building 2 on the east side; sample SG-B4 was collected outside Building 11 on the east side. The locations of samples SG-B2 and SG-B4 are shown on Figures 4 and 5, respectively.

# 4.3 Quality Assurance/Quality Control

Quality assurance/quality control (QA/QC) samples for this sampling event included a co-located sample (field duplicate), which was collected by placing two canisters side by side and opening the valves simultaneously. The tubing from the field duplicate subslab probe was connected to each of the canisters utilizing a "T" fixture so the samples drew the

same air from the subslab probe. One field blank was collected by placing two canisters side by side and not opening one of the valves. This field blank was a control sample, and the SUMMA<sup>TM</sup> canisters were cleaned to the reporting limits by the laboratory. Zero grade humidified air was used by the laboratory to pressurize the canisters.

# 4.4 Laboratory Analysis and Validation

Columbia Analytical Services of Simi Valley, California (NYSDOH Environmental Laboratory Approval Program, New York Laboratory Identification Number 11221) supplied certified clean SUMMA<sup>TM</sup> canisters with individual tracking numbers and calibrated flow regulators. Air samples (subslab and outdoor air) were collected using evacuated stainless steel SUMMA<sup>TM</sup> canisters and analyzed in accordance with *Compendium Method TO-15: Determination of VOCs in Air Collected in Specially Prepared Canisters and Analyzed by Gas Chromatography (GC)/ Mass Spectrometry (MS)* (U.S. Environmental Protection Agency [USEPA] 1999). The reporting limits for Method TO-15 presented in the work plan (CH2M HILL 2007) were achieved (Appendix D).

The data were validated using applicable quality criteria in the *National Functional Guidelines* for Organic Data Review (USEPA 1994) and USEPA Region 2 data validation procedures (USEPA 2007). Appendix D contains the validated laboratory data package for the samples collected during this investigation.

### 4.5 Field Observations

Because of the thickness of the concrete slab in several of the buildings, selected vapor probe locations were moved from 1 to 3 feet away from the proposed locations to areas where the probe installation could be completed, with the exception of two locations (SG-5 and SG-5a), which were moved 5 to 10 feet away from the proposed locations. The locations that were moved 1 to 3 feet comprised SG-2, SG-3, SG-5, SG-5a, SG-6, SG-7, and SG-8. The final locations of the samples collected are presented on Figures 4 and 5. During installation of the vapor probe at location SG-5, an unknown liquid was observed on the drill bit, and an odor was noted during one of several attempts to install the vapor probe. Once the subslab probe at SG-5 was installed, the unsuccessful attempts were filled with concrete.

Prior to sampling, the field team noted activities in or near the sampling areas such as constituent use or storage and operation of forklifts or other vehicles that may influence the ambient and/or indoor air quality. A list of chemicals in each building is included as Appendix A.

#### **SECTION 5**

# Subslab Soil Vapor Data Evaluation

The subslab sample results from each building were evaluated for the detected constituents, and are presented in Tables 1 through 9. The following lines of evidence were evaluated in this report to determine the potential significance of the vapor intrusion pathways in the various buildings:

- Comparison between types of sampling results subslab soil vapor and ambient outdoor air, and available soil and groundwater data
- Comparison of the subslab soil vapor results to background levels of volatile constituents in indoor air – 90th percentile indoor air background levels from NYSDOH (2006)
- Comparison of the subslab soil vapor results to the NYSDOH air guideline values (2006)
- Comparison of the subslab soil vapor results to the USEPA (2002) generic risk-based target shallow soil gas-to-indoor air screening level at attenuation factors (AFs) of 0.1 and 0.01

The concentration levels of the constituents were compared to the criteria described in Section 5.1. The subslab soil vapor data evaluation approach is described in Section 5.2. The data assessment results are summarized in Section 5.3.

## 5.1 Criteria Used for Comparison

## 5.1.1 Site-Specific Ambient Outdoor Air Concentrations

As stated in NYSDOH guidance (2006), "New York State currently does not have any standards, criteria or guidance values for concentrations of compounds measured in subslab soil vapor samples. Additionally, there are currently no databases available of background levels of volatile constituents in soil vapor." NYSDOH (2006) guidance suggests that a comparison of subslab soil vapor results with background outdoor air concentrations will provide some perspective on the need for additional investigations. Therefore, as a first step, the subslab vapor sample results from the buildings/areas were compared with the site-specific ambient air concentrations.

Ambient air sample SG-B2 was collected near Building 2-B, Building 2, MPA Process Area (Building 2), and Building 1, and thus represents outdoor air conditions for these buildings. Sample SG-B4 was collected near Building 3, Building 4, and the Tank Storage Area, and thus represents outdoor air conditions for these buildings. The site-specific ambient air analytical results are listed in Tables 1 and 2. Subslab vapor data were compared to site-specific ambient air data to provide one line of evidence to determine if vapor intrusion is a potential concern.

#### 5.1.2 90th Percentile Indoor Air Background Levels (NYSDOH)

The 90th percentile indoor air background levels are provided in Table C1: *NYSDOH* 2003 *Study of Volatile Organic Chemicals in Air of Fuel Oil Heated Homes of Appendix C* (NYSDOH 2006). Note that background indoor air concentrations are not risk-based, and an exceedance only indicates if the subslab soil vapor concentration is different from background aboveground indoor air concentrations, which provides one line of evidence in determining if the vapor detections are potentially site-related and/or if additional investigations are needed to further assess the soil vapor intrusion pathway. Additionally, some of the detected constituents do not have background indoor air concentrations listed in the NYSDOH guidance document (2006). The comparison is provided in Tables 3 through 9.

#### 5.1.3 Air Guideline Values (NYSDOH)

NYSDOH (2006) provides air guideline values for five constituents (methylene chloride, polychlorinated biphenyls, tetrachlorodibenzo-*p*-dioxin equivalents, tetrachloroethene [PCE], and TCE). Detected concentrations in subslab soil vapor were compared against the available NYSDOH air guideline values. As noted above, NYSDOH currently does not have guidance values for subslab soil vapor. However, NYSDOH (2006) guidance suggests that a comparison of subslab soil vapor results with indoor air guideline values may provide some perspective on the concentrations measured in subslab soil vapor. The comparison of subslab soil vapor concentrations to air guideline values is provided in Tables 3 through 9.

# 5.1.4 Target Shallow Soil Gas-to-Indoor Air Screening Level at AFs of 0.1and 0.01 from USEPA (2002)

Target shallow soil gas concentrations were provided from USEPA (2002), *Draft Guidance for Evaluating the Vapor Intrusion to Indoor Air Pathway from Groundwater and Soils (Subsurface Vapor Intrusion Guidance)*. USEPA (2002) generic risk-based soil gas-to-indoor air screening levels based on AFs of 0.1 and 0.01 were used. Values based on a range of AFs (0.1 to 0.01) were used given the significant uncertainties in the actual attenuation and given statements from USEPA at various vapor intrusion workshops the last few years (http://iavi.rti.org/WorkshopsAndConferences.cfm) that it has been considering revising the default shallow soil gas-to-indoor air AF of 0.1 to a value closer to 0.01. Tables 3 to 9 include results of the comparisons against these criteria.

# 5.2 Subslab Soil Vapor Data Evaluation Approach

This soil vapor data evaluation generally was conducted in accordance with the NYSDOH *Guidance for Evaluating Soil Vapor Intrusion in the State of New York* (NYSDOH 2006). Consistent with NYSDOH SVI guidance (2006), multiple lines of evidence were considered in this SVI evaluation. The criteria used for comparison during this data evaluation are presented in Section 5.1.

Subslab soil vapor results were compared with site-specific outdoor air data collected in April 2008. Constituents of potential interest (COPIs) (i.e., constituents with detected concentrations above the site-specific outdoor air concentrations) were identified for each building. COPIs at each sample location exceeding the NYSDOH 90th percentile background indoor air concentrations are presented on Figures 4 and 5.

The subslab soil vapor results also were considered relative to historical site soil and groundwater data, if available, for the area where each subslab soil vapor sample was collected. As discussed in Section 5.1, subslab soil vapor concentrations also were compared with (1) the range of New York background indoor air levels listed in NYSDOH (2006), (2) NYSDOH air guideline values, and (3) USEPA (2002) generic soil gas-to-indoor air screening values in order to help focus on the VOCs that are of a vapor intrusion concerns.

Where applicable, groundwater data from co-located monitoring wells or piezometers (i.e., wells located in close proximity to the subslab soil vapor samples) were compared to generic groundwater screening levels. The generic groundwater screening levels were the target groundwater concentrations from USEPA (2002), *Draft Guidance for Evaluating the Vapor Intrusion to Indoor Air Pathway from Groundwater and Soils (Subsurface Vapor Intrusion Guidance)*, which are based on a default groundwater-to-indoor air AF of 0.001 and the assumption that partitioning across the water table obeys Henry's Law. These comparisons are presented in Tables 10 through 15. Co-located monitoring wells and peizometers are identified on Figure 6.

The subslab data collected in April 2008 are discussed in detail below for each of the buildings and areas (Building 2-A/2-B, Building 2, MPA Process Area [Building 2], Building 1, Building 3, Building 4, and Tank Storage Area). The COPIs identified are discussed individually under each site along with lines of evidence for their occurrence and vapor intrusion potential. While all exceedances of applicable criteria were evaluated, not all of the COPIs identified correlate to materials that were managed in the respective SWMUs and/or AOCs as identified in the RFI report (CH2M HILL 2008).

As listed in Tables 3 through 9, soil vapor concentration exceedances of indoor air background levels occur at each building. Therefore, additional samples (i.e., concurrent subslab and indoor air samples) were collected in accordance with the 2007 work plan (CH2M HILL 2007) in November 2008; the results of which were presented in the *Soil Vapor Investigation Report* submitted to NYSDEC on February 8, 2010 (CH2M HILL 2010).

## 5.3 Data Assessment Results

Each COPI is discussed below for its potential to be of vapor intrusion concern based on a review of multiple lines of evidence where applicable, including a comparison of soil vapor concentrations against the NYSDOH established air guideline values, 90th percentile background levels and the maximum background levels from NYSDOH guidance, and the USEPA (2002) generic risk-based soil gas-to-indoor air screening levels (Tables 3 to 9) based on an AF of 0.1 and an AF of 0.01. Groundwater comparisons are based on analytical results from sampling periods indicated in Tables 10 through 15.

Marginal levels of acetone, carbon disulfide, MIBK, MEK, and toluene were detected in the field blank. Appendix C contains the QA/QC results table.

## 5.3.1 Building 2-A/2-B

One subslab sample (SG-1) was collected at Building 2-B (Figure 4) to represent both Buildings 2-A and 2-B. VOCs detected in subslab soil vapor above the site-specific outdoor air concentrations at Building 2-B were identified as COPIs in this vapor intrusion evaluation (Table 1). In addition, groundwater data for COPIs in MW-10, which is located

just north of Building 2-B, were compared to generic groundwater screening levels (Table 10 and Figure 6). Note that Buildings 2-A and 2-B are considered the same air space and contain the process location for manufacturing thiodipropionate esters and 2-sulfoethyl methacrylate.

The following observations were made for COPIs detected in the soil vapor sample associated with Buildings 2-A and 2-B.

#### Acrylonitrile

Acrylonitrile was detected in the subslab sample associated with Buildings 2-A and 2-B, and results of comparison against various criteria are discussed below.

- The subslab soil vapor concentration exceeded the corresponding site-specific outdoor air concentration (Table 1).
- No NYSDOH guidance based background levels are available for comparison for this
  constituent.
- The subslab soil vapor concentration in the sample (5.9 micrograms per cubic meter  $[\mu g/m^3]$ ) exceeded the USEPA (2002) generic risk-based soil gas-to-indoor air screening level of 3.6  $\mu g/m^3$  at an AF of 0.01 (Table 3).
- Acrylonitrile was not detected in groundwater in monitoring well MW-10, which is identified as AOC E, and located outside the northern side of Building 2-B (Table 10).
- Buildings 2-A and 2-B contain the process location for manufacturing thiodipropionate esters and 2-sulfoethyl methacrylate, which includes the use of acrylonitrile.

#### Chloroform and TCE

Chloroform and TCE were detected in subslab samples from Buildings 2A and 2B, and results of comparison against various criteria are discussed below.

- Subslab soil vapor concentrations exceeded the corresponding site-specific outdoor air concentrations (Table 1).
- Subslab soil vapor concentrations exceeded the NYSDOH 90th percentile background indoor air concentrations.
- TCE in subslab soil vapor does not exceed the established NYSDOH air guideline value of  $5 \mu g/m^3$  (Table 3).
- Subslab soil vapor concentrations were above the generic risk-based soil gas-to-indoor air screening level based on an AF of 0.1 criteria, but not above the screening level based on an AF of 0.01 criteria from USEPA (2002) (Table 3).
- Chloroform and TCE were not detected in groundwater in monitoring well MW-10 (Table 10).

#### **MIBK**

MIBK was detected in sub-slab sample from Buildings 2A and 2B, and results of comparison against various criteria are discussed below.

- The subslab soil vapor concentration exceeded the corresponding site-specific outdoor air concentration (Table 1).
- The subslab soil vapor concentration exceeded the NYSDOH 90th percentile and maximum background indoor air concentrations (Table 3).
- The subslab soil vapor concentration did not exceed the generic risk-based soil gas-to-indoor air screening level based on an AF of 0.1 criteria from USEPA (2002) (Table 3).
- MIBK was detected in groundwater in monitoring well MW-10 (Table 10).
- MIBK is a constituent of Building 2A operations and was detected in ambient air samples, albeit at approximately 10 times lower levels.

#### Other Detected Constituents

All other results of detected constituents in subslab soil vapor samples, comprising carbon tetrachloride, chloromethane, methylene chloride, and PCE are discussed below:

- The subslab soil vapor concentrations exceeded the corresponding site-specific outdoor air concentrations (Table 1).
- The subslab soil vapor concentrations did not exceed the NYSDOH 90th percentile or the maximum background indoor air concentrations (Table 3).
- PCE in the subslab soil vapor does not exceed the established NYSDOH air guideline value of  $100 \,\mu\text{g/m}^3$  (Table 3).
- Subslab soil vapor concentrations were below the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at a AF of 0.1 and 0.01 (Table 3).
- None of these COPIs was detected in groundwater in well MW-10 located outside the northern site of Building 2-B (Table 10).

Overall, several VOCs are present in subslab soil vapor under Buildings 2-A and 2-B. Of specific interest is acrylonitrile, as the subslab soil vapor concentration exceeded the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.01. However, only one of the detected constituents was in common with those detected in AOC E groundwater, which is MIBK.

## 5.3.2 Building 2

One subslab sample (SG-2) was collected at Building 2 (Figure 4). A duplicate sample was collected for the same duration and was a split sample of SG-2. There are no known VOC impacts to subsurface soil or groundwater in the area of Building 2. VOCs detected in subslab soil vapor above the site-specific outdoor air concentrations at Building 2 were identified as COPIs in this vapor intrusion evaluation (Table 1). Note that Building 2 contains indoor sources for constituents because of the process location for manufacture of 3-MPA and low volume specialty constituents produced on a campaign basis.

The following observations were made for COPIs detected in soil vapor at Building 2.

#### Chloroform and TCE

Chloroform and TCE were detected in the subslab sample from Building 2, and results of comparison against various criteria are discussed below.

- Subslab soil vapor concentrations exceeded the corresponding site-specific outdoor air concentrations (Table 1).
- Subslab soil vapor concentrations exceeded the NYSDOH 90th percentile and/or maximum background indoor air concentrations.
- TCE does not exceed the established NYSDOH air guideline value of 5 μg/m³ (Table 4).
- Subslab soil vapor concentrations were above the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.01 (Table 4).

#### 1,2-DCA, Carbon Tetrachloride, and PCE

1,2-DCA, carbon tetrachloride, and PCE were detected in the subslab sample from Building 2, and results of comparison against various criteria are discussed below.

- Subslab soil vapor concentrations exceeded the corresponding site-specific outdoor air concentrations (Table 1).
- Subslab soil vapor concentrations exceeded the NYSDOH 90th percentile background indoor air concentrations; however, they are below the maximum background level (Table 4).
- PCE does not exceed the established NYSDOH air guideline value of 100 μg/m<sup>3</sup> (Table 4).
- Subslab soil vapor concentration were above the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.1 criteria, but below an AF of 0.01 (Table 4).

#### 1,1-DCE, 1,2-DCP, and Acetone

1,1-DCE, 1,2-DCP, and acetone also were detected in subslab sample from Building 2, and results of comparison against various criteria are discussed below.

- The subslab soil vapor concentrations exceeded the corresponding site-specific outdoor air concentrations (Table 1).
- The subslab soil vapor concentrations exceeded the NYSDOH 90th percentile; however, they are below the maximum background level (Table 4).
- Subslab soil vapor concentrations were below the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.1 (Table 4).

At Building 2, the maximum detected concentration of all other constituents was below the most conservative screening criteria. Overall, no VOCs were detected in subsurface soil or groundwater in the area of Building 2. Only chloroform and TCE indicated levels in the soil vapor sample exceeding the USEPA (2002) generic risk-based soil gas-to-indoor air

screening level at an AF of 0.01. All other VOCs were generally below the health-protection based conservative screening values from USEPA with an AF of 0.01 and the NYSDOH background maximum levels.

#### 5.3.3 MPA Process Area (Building 2)

Two subslab samples (SG-3 and SG-3a) were collected in the MPA Process Area (Figure 4). VOCs detected in subslab soil vapor above the site-specific outdoor air concentrations at the MPA Process Area were identified as COPIs (Table 1). This area is the process location for manufacture of 3-MPA within Building 2.

The following observations were made for COPIs detected in soil vapor at the MPA Process Area (Building 2).

#### Acrylonitrile

Acrylonitrile was detected in subslab samples from MPA Process Area of Building 2, and results of comparison against various criteria are discussed below.

- Subslab soil vapor concentrations exceeded the corresponding site-specific outdoor air concentrations (Table 1).
- No NYSDOH guidance based background levels are available for comparison for this constituent (Table 5).
- Subslab soil vapor concentrations exceeded the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.1 and 0.01 (Table 5).
- Acrylonitrile is a constituent of Building 2 operations.

#### Benzene, Chloroform, and TCE

Benzene, chloroform, and TCE were detected in subslab samples from MPA process area of Building 2 and results of comparison against various criteria are discussed below.

- Subslab soil vapor concentrations exceeded the corresponding site-specific outdoor air concentrations (Table 1).
- Subslab soil vapor concentrations exceeded the corresponding site-specific outdoor air concentrations and the NYSDOH 90th percentile background indoor air concentrations, but all three constituents detected in subslab soil vapor samples were below the NYSDOH maximum background levels (Table 5).
- One of the two subslab sample concentrations exceeded the NYSDOH established air guideline value of 5  $\mu$ g/m³ for TCE (Table 5).
- Subslab soil vapor concentrations exceeded the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.01 (Table 5).

#### Chloroethane; cis-1,2-DCE; MIBK; and PCE

Chloroethane; cis-1,2-DCE; MIBK; and PCE also were detected in subslab samples from the MPA Process Area of Building 2, and results of comparison against various criteria are discussed below.

- The subslab soil vapor concentrations exceeded the corresponding site-specific outdoor air concentrations (Table 1).
- The subslab soil vapor concentrations exceeded the NYSDOH 90th percentile; however, it was below the maximum background level (Table 5).
- Subslab soil vapor concentrations were below the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.1 and 0.01 (Table 5).
- PCE does not exceed the established NYSDOH air guideline value of  $100 \,\mu g/m^3$  (Table 5).
- MIBK is related to production in Building 2.

At the MPA Process Area of Building 2, the maximum detected concentrations of all other constituents were below the most conservative screening criteria.

Overall, in the MPA Process Area, there are no known VOC impacts to subsurface soil or groundwater. Acrylonitrile, benzene, chloroform, and TCE was detected at levels in soil vapor samples exceeding the USEPA (2002) generic risk-based soil gas-to-indoor air screening levels at an AF of 0.01. In addition, TCE exceeded the NYSDOH indoor air screening level of 5  $\mu$ g/m³. All other VOCs were generally below the health protection-based conservative screening values from USEPA with an AF of 0.01 and the NYSDOH background maximum levels.

## 5.3.4 **Building 1**

One subslab sample (SG-4) was collected in Building 1 (Figure 4). VOCs detected in subslab soil vapor above the site-specific outdoor air concentrations at Building 1 were identified as COPIs (Table 1). Note that Building 1 is the site boiler room, and VOC sources have not been identified in this area.

The following observations were made for COPIs detected in soil gas at Building 1.

#### Chloroform and TCE

Chloroform and TCE were detected in the subslab sample from Building 1, and results of comparison against various criteria are discussed below.

- Subslab soil vapor concentrations exceeded the corresponding site-specific outdoor air concentrations (Table 1).
- Subslab soil vapor concentrations exceeded the NYSDOH 90th percentile and/or maximum background indoor air concentrations.
- TCE does not exceed the established NYSDOH air guideline value of 5 μg/m³ (Table 6).

• Subslab soil vapor concentrations were above the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.01 (Table 6).

#### **PCE**

PCE was detected in subslab sample from Building 1, and results of comparison against various criteria are discussed below.

- The subslab soil vapor concentration exceeded the corresponding site-specific outdoor air concentration (Table 1).
- The subslab soil vapor concentration exceeded the NYSDOH 90th percentile background indoor air concentration, but was below the maximum background level.
- PCE did not exceed the established NYSDOH air guideline value of  $100 \mu g/m^3$  (Table 6).
- Subslab soil vapor concentration was above the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.1, but not an AF of 0.01 (Table 6).

#### 1,1,1-Trichloroethane; 1,1,-DCA; and 1,2-DCP

1,1,1-Trichloroethane (TCA); 1,1,-DCA; and 1,2-DCP were detected in subslab sample from Building 1, and results of comparison against various criteria are discussed below.

- The subslab soil vapor concentrations exceeded the corresponding site-specific outdoor air concentrations (Table 1).
- The subslab soil vapor concentrations exceeded the NYSDOH 90th percentile background indoor air concentrations, but concentrations were below the NYSDOH maximum background indoor air concentrations (Table 6).
- Subslab soil vapor concentrations were below the USEPA (2002) generic risk-based soil gas-to-indoor air screening levels at an AF of 0.1 (Table 6).

Benzene, Chloromethane, Ethylbenzene, Methylene Chloride, Toluene, and Xylenes Benzene, chloromethane, ethylbenzene, methylene chloride, toluene, and xylenes were detected in subslab sample from Building 1, and results of comparison against various criteria are discussed below.

- The subslab soil vapor concentrations exceeded the corresponding site-specific outdoor air concentration (Table 1).
- The subslab soil vapor concentrations did not exceed the NYSDOH 90th percentile or the maximum background indoor air concentrations, if available.
- Methylene chloride concentrations are below the established NYSDOH air guideline value (Table 6).
- The subslab soil vapor concentrations were below the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.1 (Table 6).

Overall, there are no known VOC impacts to subsurface soil or groundwater in the area of Building 1. Only chloroform and TCE indicated levels in the soil vapor sample exceeding the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.01.

All other VOCs were generally below the health protection-based conservative screening values from USEPA with an AF of 0.01 and the NYSDOH background maximum levels.

#### **5.3.5** Building 3

Three subslab samples (SG-5, SG-5a, and SG-6) were collected within Building 3 (Figure 5). VOCs detected in subslab soil vapor above the site-specific outdoor air concentrations at Building 3 were identified as COPIs in this vapor intrusion evaluation (Table 2). In addition, groundwater data for COPIs in MW-11S, which is located southwest of Building 3 (downgradient), were compared to generic groundwater screening levels (Table 11 and Figure 6). Note that Building 3 is the process location for manufacture of T-acid.

#### Chloroform and TCE

Chloroform was detected at  $1,100 \,\mu\text{g/m}^3$ , and TCE was detected at  $20 \,\mu\text{g/m}^3$  at Building 3. These concentrations were compared against various criteria, and results are discussed below.

- Subslab soil vapor concentrations exceeded the corresponding site-specific outdoor air concentrations (Table 2).
- Subslab soil vapor concentrations exceeded the NYSDOH 90th percentile and/or the maximum background indoor air concentrations (Table 7).
- One of the three subslab sample concentrations exceeded the NYSDOH established air guideline value of 5  $\mu$ g/m³ for TCE (Table 7).
- Subslab soil vapor concentrations exceeded the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.01 (Table 7).
- Chloroform was not detected in groundwater in well MW-11S located southwest of Building 3 (Table 11).

#### **MIBK**

MIBK concentrations from Building 3 were compared against criteria as presented below.

- Subslab soil vapor concentrations exceeded the corresponding site-specific outdoor air concentrations (Table 2).
- Subslab soil vapor concentrations exceeded the NYSDOH 90th percentile and maximum background indoor air concentrations (Table 7).
- MIBK subslab soil vapor concentrations were below the conservative USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.1 (Table 7).
- MIBK was detected in groundwater in well MW-11S located southwest of Building 3; however, concentrations were below the risk-based target groundwater concentrations (Table 11).

#### 1,1,-DCA; cis-1,2-DCE; and PCE

1,1,-DCA; cis-1,2-DCE; and PCE concentrations from Building 3 subslab soil vapor samples were compared against criteria as discussed below.

- Subslab soil vapor concentrations exceeded the corresponding site-specific outdoor air concentrations (Table 2).
- Subslab soil vapor concentrations exceeded the NYSDOH 90th percentile background indoor air concentrations, but concentrations were below the NYSDOH maximum background indoor air concentrations (Table 7).
- PCE concentrations are below the established NYSDOH air guideline value (Table 7).
- Subslab soil vapor concentrations were below the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.1 (Table 7).
- 1,1,-DCA; cis-1,2-DCE; and PCE were not detected in groundwater at MW-11S (Table 11).

# Bromodichloromethane, Carbon Disulfide, Carbon Tetrachloride, Methylene Chloride, and Xylenes

Bromodichloromethane, carbon disulfide, carbon tetrachloride, methylene chloride, and xylenes concentrations from Building 3 subslab soil vapor samples were compared against criteria as discussed below.

- Subslab soil vapor concentrations exceeded the corresponding site-specific outdoor air concentrations (Table 2).
- Subslab soil vapor concentrations did not exceed the NYSDOH 90th percentile indoor air concentrations, if a value is available in NYSDOH guidance.
- Methylene chloride concentrations are below the established NYSDOH air guideline value (Table 7).
- Subslab soil vapor concentrations were below the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.1 (Table 7).
- Carbon disulfide and xylenes were detected in groundwater at MW-11S; however, their concentrations were below the risk-based target groundwater concentrations (Table 11).
- Bromodichloromethane, carbon tetrachloride, and methylene chloride were not detected in groundwater at MW-11S (Table 11).

Overall, Building 3 chloroform and TCE subslab soil vapor levels were well above screening levels and exceeded the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.01.

## 5.3.6 Building 4

Three subslab samples (SG-7, SG-7a, and SG-8) were collected at Building 4 (Figure 5). VOCs detected in subslab soil vapor above the site-specific outdoor air concentrations at Building 4 were identified as COPIs in this vapor intrusion evaluation (Table 2). In

addition, potential source area in groundwater as monitored by groundwater monitoring wells for COPIs in MW-03, PZ-1, and BLDG4-PIT-SSP, which are located inside Building 4 (Figure 6). The VOC data from these wells were compared to generic groundwater screening levels (Tables 12, 13, and 14). Note that Building 4 (and 4A) is the process location for purification of T-acid and the manufacture of thioglycolates.

The following observations were made for constituents detected in soil vapor at Building 4.

#### Chloroform, ethylbenzene, MIBK, and TCE

Chloroform, ethylbenzene, MIBK, and TCE concentrations from Building 4 subslab soil vapor samples were compared against criteria as discussed below.

- Subslab soil vapor concentrations exceeded the corresponding site-specific outdoor air concentrations (Table 2).
- Subslab soil vapor concentrations exceeded the NYSDOH 90th percentile and maximum background indoor air concentrations (Table 8).
- TCE exceeded the established NYSDOH air guideline value in one of three samples collected (Table 8).
- Subslab soil vapor concentrations were above the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.01 (Table 8).
- Chloroform was detected in groundwater in PZ-1 and BLDG4-PIT-SSP, both located inside Building 4, above the risk-based target groundwater concentration (Tables 12 and 14).
- MIBK was detected in groundwater in PZ-1, MW-03, and BLDG4-PIT-SSP, all located inside Building 4; however, the concentrations were below the risk-based target groundwater concentration at PZ-1 and MW-03. At least one of the analyses at BLDG4-PIT-SSP exceeded the risk-based target groundwater concentration (Tables 12, 13, and 14).
- Ethylbenzene and TCE were detected in groundwater in MW-03, located inside Building 4; however, the concentrations were below the risk-based target groundwater concentration (Table 13).
- MIBK is a constituent of Building 4 operations and was detected in ambient air samples.

#### Methylene chloride and PCE

Methylene chloride and PCE concentrations from Building 4 subslab soil vapor samples were compared against criteria as discussed below.

- Subslab soil vapor concentrations exceeded the corresponding site-specific outdoor air concentrations (Table 2).
- Subslab soil vapor concentrations exceeded the NYSDOH 90th percentile background indoor air concentrations, but were below the maximum background levels (Table 8).

- Methylene chloride exceeded the established NYSDOH air guideline value in one of three samples collected; however, PCE concentrations did not exceed the air guideline value (Table 8).
- Subslab soil vapor concentrations were above the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.1 but below the AF of 0.01 (Table 8).
- Methyl chloride was detected in groundwater at PZ-1 and BLDG4-PIT-SSP; however, the concentrations at PZ-1 were below the risk-based target groundwater concentration, while the concentrations at PZ-1 exceeded the risk-based target groundwater concentration (Tables 12 and 14).
- PCE was not detected in groundwater at PZ-1, MW-03, and BLDG4-PIT-SSP (Tables 12, 13, and 14).

#### Acrylonitrile and Benzene

Acrylonitrile and benzene concentrations from Building 4 subslab soil vapor samples were compared against criteria as discussed below.

- Subslab soil vapor concentrations exceeded the corresponding site-specific outdoor air concentrations (Table 2).
- Benzene was below the NYSDOH 90th percentile and maximum background indoor air concentrations. Acrylonitrile does not have NYSDOH background values (Table 8).
- Subslab soil vapor concentrations for both constituents were above the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.01, but below the AF of 0.1 (Table 8).
- Benzene was detected in groundwater at MW-03 and BLDG4-PIT-SSP; however, the concentrations were below the risk-based target groundwater concentration (Tables 13 and 14).
- Acrylonitrile was not detected in groundwater at PZ-1, MW-03 and BLDG4-PIT-SSP (Tables 12, 13, and 14).

#### Cis-1,2-DCE and Xylenes

Cis-1,2-DCE and xylenes concentrations from Building 4 subslab soil vapor samples were compared against criteria as discussed below.

- Subslab soil vapor concentrations exceeded the corresponding site-specific outdoor air concentrations (Table 2).
- Subslab soil vapor concentrations exceeded the NYSDOH 90th percentile and maximum background indoor air concentrations (Table 8).
- Cis-1,2-DCE and xylenes in subslab soil vapor samples were below the USEPA (2002) generic risk-based soil gas-to-indoor air screening levels at an AF of 0.1 (Table 8).
- Cis-1,2-DCE was detected in PZ-1, MW-03, and BLDG4-PIT-SSP, and xylenes (total) were detected in MW-01 and BLDG4-PIT-SSP; however, the concentrations were below the risk-based target groundwater concentration (Tables 12, 13, and 14).

#### 1,1-DCA; 1,2-DCP; Chlorobenzene; MEK; and Toluene

1,1-DCA; 1,2-DCP; chlorobenzene; MEK; and toluene concentrations from Building 4 subslab soil vapor samples were compared against criteria as discussed below.

- Subslab soil vapor concentrations exceeded the corresponding site-specific outdoor air concentrations (Table 2).
- Subslab soil vapor concentrations for all five of the compounds exceeded the NYSDOH 90th percentile background indoor air concentrations, while subslab soil vapor concentrations for 1,1-DCA; 1,2-DCP; and chlorobenzene exceeded the NYSDOH maximum background indoor air concentrations (Table 8).
- Subslab soil vapor concentrations were below the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.1 (Table 8).
- 1,1-DCA and MEK were detected in groundwater at PZ-1, MW-03, and BLDG4-PIT-SSP; however, the concentrations were below the risk-based target groundwater concentrations (Tables 12, 13, and 14).
- Chlorobenzene and toluene were detected in groundwater at MW-03 and BLDG4-PIT-SSP; however, the concentrations were below the risk-based target groundwater concentrations (Tables 13 and 14).
- 1,2-DCP was detected in groundwater at MW-03; however, the concentrations were below the risk-based target groundwater concentration (Table 13).

#### Carbon Disulfide and Chloromethane

Carbon disulfide and chloromethane concentrations from Building 4 subslab soil vapor sample were compared against criteria as discussed below.

- Subslab soil vapor concentrations exceeded the corresponding site-specific outdoor air concentrations (Table 2).
- Subslab soil vapor concentrations did not exceed the NYSDOH 90th percentile or the maximum background indoor air concentrations, if available (Table 8).
- Subslab soil vapor concentrations were below the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.1 (Table 8).
- Carbon disulfide was detected in groundwater at PZ-1, MW-03, and BLDG4-PIT-SSP; however, the concentrations were below the risk-based target groundwater concentration (Tables 12, 13, and 14).
- Chloromethane was not detected in groundwater at PZ-1, MW-03, and BLDG4-PIT-SSP (Tables 12, 13, and 14).

Building 4 has a potential groundwater source area for several VOCs. Most of the VOCs detected in groundwater also were detected in the subslab soil vapor samples. However, some of the constituents detected also are part of the ongoing manufacturing processes in the building. Chloroform, ethylbenzene, MIBK, and TCE exceeded the USEPA (2002)

generic risk-based soil gas-to-indoor air screening level at an AF of 0.01. TCE and methylene chloride also exceeded the NYSDOH indoor air screening value.

Based on the soil sample results reported in the RFI addendum report (CH2M HILL 2008), the VOCs in soil sample location PZ-1 were low concentrations or not detected.

#### 5.3.7 Tank Storage Area

One subslab sample (SG-9) was collected at the Tank Storage Area building (Figure 5). VOCs detected in subslab soil vapor above the site-specific outdoor air concentrations at the Tank Storage Area building were identified as COPIs in this vapor intrusion evaluation (Table 2). In addition, groundwater data for COPIs in PZ-3, which is located inside the Tank Storage Area building (Figure 6), were compared to generic groundwater screening levels (Table 15).

#### Chloroform and TCE

Chloroform and TCE concentrations from Tank Storage Area subslab soil vapor sample were compared against criteria as discussed below.

- Subslab soil vapor concentrations exceeded the corresponding site-specific outdoor air concentrations (Table 2).
- Subslab soil vapor concentrations exceeded the NYSDOH 90th percentile and maximum background indoor air concentrations (Table 9).
- The concentration of TCE exceeds the established NYSDOH air guideline value (Table 9).
- Subslab soil vapor concentrations were above the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.1 and 0.01 (Table 9).
- Chloroform and TCE were not detected in groundwater at PZ-3 (Table 15).

#### 1,2-DCP and PCE

1,2-DCP and PCE concentrations from Tank Storage Area subslab soil vapor sample were compared against criteria as discussed below.

- Subslab soil vapor concentrations exceeded the corresponding site-specific outdoor air concentrations (Table 2).
- Subslab soil vapor concentrations exceeded the NYSDOH 90th percentile background indoor air concentrations, but concentrations were below the NYSDOH maximum background indoor air concentrations.
- PCE concentrations are below the established NYSDOH air guideline value (Table 9).
- Subslab soil vapor concentrations were below the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at a AF of 0.1 (Table 9).
- 1,2-Dichloropropane was detected in groundwater at PZ-3; however, the concentration was below the risk-based target groundwater concentration (Table 15).
- PCE was not detected in groundwater at PZ-3 (Table 15).

#### Benzene and Xylenes

Benzene and xylenes concentrations from Tank Storage Area subslab soil vapor sample were compared against criteria as discussed below.

- Subslab soil vapor concentrations exceeded the corresponding site-specific outdoor air concentrations (Table 2).
- Subslab soil vapor concentrations did not exceed the NYSDOH 90th percentile or the maximum background indoor air concentrations (Table 9).
- Subslab soil vapor concentrations were below the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.1 (Table 9).
- Benzene and xylenes were not detected in groundwater at PZ-3 (Table 15).

The Tank Storage Area has a potential groundwater source area for 1,2-DCP. This compound also was detected in the subslab soil vapor sample, and while it exceeded the NYSDOH 90th percentile background indoor air concentrations, it did not have any other exceedances.

Chloroform and TCE exceeded the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.01. TCE also exceeded the NYSDOH indoor air screening value. The adjacent Building 4 has a potential groundwater source area for VOCs.

Based on the soil sample results reported in the RFI addendum report (CH2M HILL 2008), the VOCs in soil sample location PZ-3 were low concentrations or not detected.

#### **SECTION 6**

# **Summary and Recommendations**

Subslab soil vapor sampling has been conducted to evaluate the potential subsurface-to-indoor air vapor intrusion pathway at Buildings 2-A, 2-B, 2, 1, 3, 4, MPA Process Area, and Tank Storage Area. A number of VOCs were detected in the subslab soil vapor samples at each building. A review of these detections indicated several constituents were above site-specific background levels at each building. In addition, some soil vapor concentrations were shown to exceed statewide background indoor air concentrations and USEPA (2002) generic risk-based soil gas-to-indoor air screening levels at an AF of 0.01 or 0.1.

Some VOCs were detected beneath all of the buildings and included chloroform, PCE, and TCE. Chloroform and MIBK also were detected in the soil vapor samples at Building 4 (Table 2).

- Buildings 2-A and 2-B: Several VOCs are present in subslab soil vapor under Buildings 2-A and 2-B. Of specific interest is acrylonitrile, as the subslab soil vapor concentration exceeded the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.01.
- Building 2: Chloroform and TCE subslab soil vapor concentrations were above the USEPA (2002) generic risk-based soil gas-to-indoor air screening levels at an AF of 0.01.
- MPA Process Area: Acrylonitrile, benzene, chloroform, and TCE had levels in subslab soil gas samples exceeding the USEPA (2002) generic risk-based soil gas-to-indoor air screening levels at an AF of 0.01. In addition, TCE in subslab soil vapor exceeded the NYSDOH indoor air screening level of  $5 \mu g/m^3$ .
- Building 1: Chloroform and TCE subslab soil vapor concentrations were above the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.01.
- Building 3: Chloroform and TCE subslab soil vapor levels were above the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.01.
- Building 4: Chloroform, ethylbenzene, MIBK, and TCE exceeded the USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.01. TCE and methylene chloride in subslab soil vapor also exceeded the NYSDOH indoor air screening value.
- Tank Storage Area: 1,2-DCP was detected in the subslab soil vapor sample, and while it
  exceeded the NYSDOH 90th percentile background indoor air concentrations, it did not
  have any other exceedances of risk-based criteria. Chloroform and TCE exceeded the
  USEPA (2002) generic risk-based soil gas-to-indoor air screening level at an AF of 0.01.
  TCE in subslab soil vapor also exceeded the NYSDOH indoor air screening value and
  may have migrated.

As discussed in NYSDOH (2006) and other state and federal vapor intrusion guidance documents, elevated subslab soil gas concentrations do not necessarily mean vapors are

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intruding into a building, but rather suggest the potential exists. Based on the results presented in this report and following the NYSDOH (2006) vapor intrusion evaluation guidance and the work plan previously submitted to NYSDEC (CH2M HILL 2007), additional indoor air sampling was performed for the buildings detailed in this report. This sampling included concurrent indoor air, outdoor air, and subslab vapor sampling, and the results were presented in the *Soil Vapor Investigation Report* submitted to NYSDEC on February 8, 2010 (CH2M HILL 2010). Additional investigation activities also have been performed in other media, including soil and groundwater, after developing the 2007 work plan. Based on this additional information, a revised work plan for continued vapor intrusion investigation will be prepared and submitted to NYSDEC for review.

It should be noted that comments were received from NYSDEC in June 2010 on the *Soil Vapor Investigation Report* (CH2M HILL 2010) referenced above, after a draft of this report had been prepared. The NYSDEC comments stated that use of USEPA RSLs, attenuation factors, and other USEPA-recommended risk-based screening levels is not part of the NYSDOH VI guidance (NYSDEC 2010). NYSDEC requested that USEPA methodologies not be included in future VI reports for this facility. Because this report was already in preparation, changes were not made to delete the USEPA risk-based results. However, it is important to note that the conclusions of this report would be the same if the USEPA risk-based criteria were not used in the evaluation. The conclusions from this evaluation are that all of the buildings included in this report would have additional sampling done and no buildings were excluded from further evaluation based on the USEPA risk-based screening levels.

#### **SECTION 7**

# References

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TABLE 1 Subslab Soil Vapor Sample Results - Buildings 1 and 2 Soil Vapor Investigation - Buildings 1, 2, 3, 4, and Tank Storage Area Former Hampshire Chemical Corp. Facility, Waterloo, New York

Field Sample ID		Outside	Building 2-B	Bui	ilding 2	MPA Process	Area (Building 2)	Building 1
. Ioia campic is		WAT-SG-B2-042308	WAT-SG-1-042308	WAT-SG-2-042308	WAT-SG-DUP-042308	WAT-SG-3-042308	WAT-SG-3a-042308	WAT-SG-4-042308
Location ID		SG-B2	SG-1	SG-2	SG-2	SG-3	SG-3A	SG-4
Sample Date		4/23/2008	4/23/2008	4/23/2008	4/23/2008	4/23/2008	4/23/2008	4/23/2008
Sample Type		Site-Specific Ambient Air	Normal	Normal	Duplicate	Normal	Normal	Normal
Lab Sample ID		P0801215-001	P0801215-002	P0801215-003	P0801215-007	P0801215-004	P0801215-005	P0801215-006
Dilution		1.52	4.13	3.4	12.5	12.3	2.92	4.13
Matrix		AIR	AIR	AIR	AIR	AIR	AIR	AIR
Parameter	CAS#							
Volatile Organics, by Method TO15 (UG/M3)								
1,1,1-TRICHLOROETHANE	71-55-6	0.15 U	0.41 U	0.34 U	1.3 U	1.2 U	0.29 U	12
1,1,2,2-TETRACHLOROETHANE	79-34-5	0.15 U	0.41 U	0.34 U	1.3 U	1.2 U	0.29 U	0.41 U
1,1,2-TRICHLOROETHANE	79-00-5	0.15 U	0.41 U	0.34 U	1.3 U	1.2 U	0.29 U	0.41 U
1,1-DICHLOROETHANE	75-34-3	0.15 U	0.41 U	0.34 U	1.3 U	1.2 U	0.29 U	2.1
1,1-DICHLOROETHENE	75-35-4	0.15 U	0.41 U	4.2	4	1.2 U	0.29 U	0.41 U
1,2-DICHLOROETHANE	107-06-2	0.15 U	0.41 U	2	2.1	1.2 U	0.29 U	0.41 U
1,2-DICHLOROPROPANE	78-87-5	0.15 U	0.41 U	4	3.6	1.2 U	0.29 U	3.9
ACETONE	67-64-1	17 U	17 U	180	250 U	130 U	56 U	17 U
ACRYLONITRILE	107-13-1	1.3	5.9	1.2 J	6.3 U	27	6.9	0.86 J
BENZENE	71-43-2	0.57	0.45	0.93	1.3 U	90	13	0.81
BROMODICHLOROMETHANE	75-27-4	0.15 U	0.41 U	0.34 U	1.3 U	1.2 U	0.29 U	0.41 U
BROMOFORM	75-25-2	0.76 U	2.1 U	1.7 U	6.3 U	6.2 U	1.5 U	2.1 U
BROMOMETHANE	74-83-9	0.15 U	0.41 U	0.34 U	1.3 U	1.2 U	0.29 U	0.41 U
CARBON DISULFIDE	75-15-0	0.71 U	1.2 U	10	8.6 U	34	9.5	2.2 U
CARBON TETRACHLORIDE	56-23-5	0.27	0.29 J	2.1	1.6	1.2 U	0.29 U	0.41 U
CHLOROBENZENE	108-90-7	0.15 U	0.41 U	0.34 U	1.3 U	1.2 U	0.29 U	0.41 U
CHLOROETHANE	75-00-3	0.15 U	0.41 U	0.34 U	1.3 U	1.2 U	0.72	0.41 U
CHLOROFORM	67-66-3	1.1	2	240	260	25	3.6	440
CHLOROMETHANE	74-87-3	0.29	1.3	0.34 U	1.3 U	1 J	0.39	0.62
cis-1,2-DICHLOROETHYLENE	156-59-2	0.15 U	0.41 U	0.34 U	1.3 U	1.2 U	1.4	0.41 U
cis-1,3-DICHLOROPROPENE	10061-01-5	0.76 U	2.1 U	1.7 U	6.3 U	6.2 U	1.5 U	2.1 U
DIBROMOCHLOROMETHANE	124-48-1	0.15 U	0.41 U	0.34 U	1.3 U	1.2 U	0.29 U	0.41 U
ETHYLBENZENE	100-41-4	0.24 J	2.1 U	0.54 J	6.3 U	0.91 J	0.43 J	0.61 J
METHYL ETHYL KETONE (2-BUTANONE)	78-93-3	3.1	3.9 U	4.7 U	2.5 U	2.8 U	4.3 U	3.7 U
METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	108-10-1	120	410	47 J	15 J	820	240	100
METHYLENE CHLORIDE	75-09-2	0.39 J	0.46 J	1.1 J	1.6 J	1.3 J	0.43 J	0.39 J
STYRENE	100-42-5	0.76 U	2.1 U	0.3 J	6.3 U	6.2 U	0.4 J	2.1 U
TETRACHLOROETHENE (PCE)	127-18-4	0.15 U	1.1	32	28	8.9	5.4	17
TOLUENE	108-88-3	1.2 U	1.5 U	5.7 J	2.1 U	16	26	5.8
TRANS-1,2-DICHLOROETHENE	156-60-5	0.15 U	0.41 U	0.34 U	1.3 U	1.2 U	0.29 U	0.41 U
TRANS-1,3-DICHLOROPROPENE	10061-02-6	0.76 U	2.1 U	1.7 U	6.3 U	6.2 U	1.5 U	2.1 U
TRICHLOROETHENE (TCE)	79-01-6	0.15 U	1.1	3.6	3.5	2.9	10	2.5
VINYL CHLORIDE	75-01-4	0.15 U	0.41 U	0.34 U	1.3 U	1.2 U	0.29 U	0.41 U
XYLENE, m,p-	108-38-3/1	0.91	0.77 J	3.3	6.3 U	4.4 J	2.3	4.3
XYLENE, o-	95-47-6	0.39 J	0.27 J	1.1 J	6.3 U	1.5 J	0.83 J	2.9
EPICHLOROHYDRIN	106-89-8	NF	NF	NF	NF	NF	NF	NF
Notes:								

All compounds and criteria are in µg/m<sup>3</sup>

NA = not applicable
NF = Not found by laboratory library search

J = estimated value

U = compound not detect at or above method detection limit

Shading in gray indicates that the constituent was detected above the site-specific ambient air concentration

TABLE 2 Subslab Soil Vapor Sample Results - Buildings 3, 4, and Tank Storage Area Soil Vapor Investigation - Buildings 1, 2, 3, 4, and Tank Storage Area Former Hampshire Chemical Corp. Facility, Waterloo, New York

-		Outside		Building 3			Building 4		Tank Storage
Field Sample ID		WAT-SG-B4-042408	WAT-SG-5-042408	WAT-SG-5a-042408	WAT-SG-6-042408	WAT-SG-7-042408	WAT-SG-7a-042408	WAT-SG-8-042408	WAT-SG-9-042408
Location ID		SG-B4	SG-5	SG-5a	SG-6	SG-7	SG-7A	SG-8	SG-9
Sample Date		4/24/2008	4/24/2008	4/24/2008	4/24/2008	4/24/2008	4/24/2008	4/24/2008	4/24/2008
Sample Type		Site-Specific Ambient Air	Normal	Normal	Normal	Normal	Normal	Normal	Normal
Lab Sample ID		P0801215-009	P0801215-010	P0801215-011	P0801215-012	P0801215-013	P0801215-014	P0801215-015	P0801215-016
Dilution		1.57	9.33	4.13	41.33	203.75	1220	16.9	4.88
Matrix		AIR	AIR	AIR	AIR	AIR	AIR	AIR	AIR
Parameter	CAS#								
Volatile Organics, by Method TO15 (UG/M3)									
1,1,1-TRICHLOROETHANE	71-55-6	0.16 U	0.93 U	0.41 U	4.1 U	20 U	120 U	1.7 U	0.49 U
1,1,2,2-TETRACHLOROETHANE	79-34-5	0.16 U	0.93 U	0.41 U	4.1 U	20 U	120 U	1.7 U	0.49 U
1,1,2-TRICHLOROETHANE	79-00-5	0.16 U	0.93 U	0.41 U	4.1 U	20 U	120 U	1.7 U	0.49 U
1,1-DICHLOROETHANE	75-34-3	0.16 U	1.4	0.41 U	4.1 U	84	120 U	1.7 U	0.49 U
1,1-DICHLOROETHENE	75-35-4	0.11 J	0.93 U	0.41 U	4.1 U	20 U	120 U	1.7 U	0.49 U
1,2-DICHLOROETHANE	107-06-2	0.16 U	0.93 U	0.41 U	4.1 U	20 U	120 U	1.7 U	0.49 U
1,2-DICHLOROPROPANE	78-87-5	0.16 U	0.93 U	0.41 U	4.1 U	52	120 U	2.1	15
ACETONE	67-64-1	35	10 U	12 U	26 U	1800 U	1600 U	230 U	13 U
ACRYLONITRILE	107-13-1	1.1	4.7 U	2.1 U	21 U	100 U	610 U	3.3 J	2.4 U
BENZENE	71-43-2	0.48	0.93 U	0.41 U	4.1 U	20 U	120 U	5	0.66
BROMODICHLOROMETHANE	75-27-4	0.16 U	0.93 U	0.43	4.1 U	20 U	120 U	1.7 U	0.49 U
BROMOFORM	75-25-2	0.79 U	4.7 U	2.1 U	21 U	100 U	610 U	8.5 U	2.4 U
BROMOMETHANE	74-83-9	0.16 U	0.93 U	0.41 U	4.1 U	20 U	120 U	1.7 U	0.49 U
CARBON DISULFIDE	75-15-0	0.92 U	2.4 U	20	2600	310 U	3800	43	3.8 U
CARBON TETRACHLORIDE	56-23-5	0.38	0.93 U	0.55	4.1 U	20 U	120 U	1.7 U	0.49 U
CHLOROBENZENE	108-90-7	0.16 U	0.93 U	0.41 U	4.1 U	20 U	120 U	2.3	0.49 U
CHLOROETHANE	75-00-3	0.086 J	0.93 U	0.41 U	4.1 U	20 U	120 U	1.7 U	0.49 U
CHLOROFORM	67-66-3	1.3	1100	450	76	24000	1300	170	270
CHLOROMETHANE	74-87-3	0.56	0.93 U	0.41 U	4.1 U	20 U	120 U	2.1	0.51
cis-1.2-DICHLOROETHYLENE	156-59-2	0.16 U	0.93 U	2.1	4.1 U	19 J	120 U	1.7 U	0.49 U
cis-1,3-DICHLOROPROPENE	10061-01-5	0.79 U	4.7 U	2.1 2.1 U	4.1 U 21 U	19 J	610 U	8.5 U	0.49 U 2.4 U
DIBROMOCHLOROMETHANE	124-48-1	0.79 U 0.16 U	0.93 U	0.41 U	4.1 U	20 U	120 U	6.5 U 1.7 U	2.4 U 0.49 U
ETHYLBENZENE	100-41-4	0.25 J	4.7 U	2.1 U	21 U	870	430 J	1.6 J	2.4 U
METHYL ETHYL KETONE (2-BUTANONE)	78-93-3	0.98 U	0.97 U	1.3 U	21 U	17 U	610 U	32	1.3 U
METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	108-10-1	94	3 U	2.3	110	12000	86000	380 J	11 J
METHYLENE CHLORIDE	75-09-2	0.36 J	4.7 U	0.29 J	2.6 J	35 J	240 J	2.3 J	0.31 J
STYRENE	100-42-5	0.12 J	4.7 U	2.1 U	21 U	100 U	610 U	8.5 U	2.4 U
TETRACHLOROETHENE (PCE)	127-18-4	0.16 U	3.8	4.1	3.4 J	26	120 U	3	6.4
TOLUENE	108-88-3	13	0.85 U	0.55 U	4.8 U	72 U	380 U	130	1.6 U
TRANS-1,2-DICHLOROETHENE	156-60-5	0.16 U	0.93 U	0.41 U	4.1 U	20 U	120 U	1.7 U	0.49 U
TRANS-1,3-DICHLOROPROPENE	10061-02-6	0.79 U	4.7 U	2.1 U	21 U	100 U	610 U	8.5 U	2.4 U
TRICHLOROETHENE (TCE)	79-01-6	0.45	2.4	20	4.1 U	54	120 U	1.7	520
VINYL CHLORIDE	75-01-4	0.16 U	0.93 U	0.41 U	4.1 U	20 U	120 U	1.7 U	0.49 U
XYLENE, m,p-	108-38-3/1	0.83	4.7 U	0.58 J	7.2 J	4600	8600	15	1.3 J
XYLENE, o-	95-47-6	0.21 J	4.7 U	0.27 J	2.9 J	950	1700	4.3 J	0.54 J
EPICHLOROHYDRIN	106-89-8	NF	NF	NF	NF	NF	NF	NF	NF

All compounds and criteria are in µg/m3
NA = not applicable
NF = Not found by laboratory library search

J = estimated value

U = compound not detect at or above method detection limit

Shading in gray indicates that the constituent was detected above the site specific ambient air concentration

TABLE 3 Building 2-B - COPIs Compared to NY State Background and USEPA Health Protective Levels Soil Vapor Investigation - Buildings 1, 2, 3, 4, and Tank Storage Area Former Hampshire Chemical Corp. Facility, Waterloo, New York

								NY:	SDOH 2003* (µg	g/m3)					
СОРІ	CAS#	Number of Soil Gas Analyses	Number of Soil Gas Detects	Minimum Detected Concentration (μg/m3)	Maximum Detected Concentration (MDC) (μg/m3)	Mean Detected Concentration (μg/m3)	Mean Detected Concentration (Attenuation Factor - 0.01) (μg/m3)	Minimum	90th Percentile	Maximum	NYSDOH Air Guideline Value** (µg/m3)	Target Shallow Gas Concentration (Attenuation Factor - 0.1) (µg/m3)	MDC Exceeds?	Target Shallow Gas Concentration (Attenuation Factor - 0.01) (µg/m3)	MDC Exceeds?
Summary Statistics Presented for COPIs Reported in Building 2	?-B (2007)														
ACRYLONITRILE	107-13-1	1	1	5.9	5.9	5.9	0.059	NA	NA	NA	NA	0.36	Yes	3.6	Yes
CARBON TETRACHLORIDE	56-23-5	1	1	0.29 J	0.29 J	0.29	0.0029	< 0.25	0.8	4.2	NA	1.6	No	16	No
CHLOROFORM	67-66-3	1	1	2	2	2	0.02	< 0.25	1.4	25	NA	1.1	Yes	11	No
CHLOROMETHANE	74-87-3	1	1	1.3	1.3	1.3	0.013	< 0.25	3.3	260	NA	24	No	240	No
METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	108-10-1	1	1	410	410	410	4.1	< 0.25	2.2	36	NA	800	No	8000	No
METHYLENE CHLORIDE	75-09-2	1	1	0.46 J	0.46 J	0.46	0.0046	< 0.25	22	2100	60	52	No	520	No
TETRACHLOROETHENE (PCE)	127-18-4	1	1	1.1	1.1	1.1	0.011	< 0.25	2.9	51	100	8.1	No	81	No
TRICHLOROETHENE (TCE)	79-01-6	1	1	1.1	1.1	1.1	0.011	< 0.25	0.5	25	5	0.22	Yes	2.2	No

**Background Indoor Air Value from** 

#### Notes:

\* = NYSDOH. 2003. Summary of Indoor and Outdoor Levels of Volatile Organic Compounds from Fuel Oil Heated Homes in NYS, 1997-2003. Revised November 14, 2005.

\*\* = NYSDOH Guidance for Evaluating Soil Vapor Intrusion in the State of New York. Table 3-1 - Air guideline values derived by the NYSDOH. October 2006.

Target shallow Gas Concentration = U.S. Environmental Protection Agency (USEPA). 2002. Draft Guidance for Evaluating the Vapor Intrusion to Indoor Air Pathway from Groundwater and Soils (Subsurface Vapor Intrusion Guidance). November 2002. MDC = Maximum detected concentration

COPI = chemical of potential interest

J = estimated value

TABLE 4 Building 2 - COPIs Compared to NY State Background and USEPA Health Protective Levels Soil Vapor Investigation - Buildings 1, 2, 3, 4, and Tank Storage Area Former Hampshire Chemical Corp. Facility, Waterloo, New York

							NYS	SDOH 2003* (µg	g/m3)					
COPI CAS	Number Soil Ga # Analyse	Soil Gas	Minimum Detected Concentration (µg/m3)	Maximum Detected Concentration (MDC) (µg/m3)	Mean Detected Concentration (µg/m3)	Mean Detected Concentration (Attenuation Factor - 0.01) (µg/m3)	Minimum	90th Percentile	Maximum	NYSDOH Air Guideline Value** (µg/m3)	Target Shallow Gas Concentration (Attenuation Factor - 0.1) (µg/m3)	MDC Exceeds?	Target Shallow Gas Concentration (Attenuation Factor - 0.01) (µg/m3)	MDC Exceeds?
Summary Statistics Presented for COPIs Reported in Building 2 (2007)	•		,,	,,		,,							, , , ,	
1,1-DICHLOROETHENE 75-38	-4 1	1	4.2	4.2	4.2	0.042	<0.25	< 0.25	430	NA	2000	No	20000	No
1,2-DICHLOROETHANE 107-0	5-2 1	1	2.1	2.1	2.1	0.021	< 0.25	< 0.25	4.9	NA	0.94	Yes	9.4	No
1,2-DICHLOROPROPANE 78-8	-5 1	1	4	4	4	0.04	< 0.25	< 0.25	34	NA	40	No	400	No
ACETONE 67-64	-1 1	1	180	180	180	1.8	< 0.25	110	690	NA	3500	No	35000	No
BENZENE 71-43	-2 1	1	0.93	0.93	0.93	0.0093	< 0.25	15	460	NA	3.1	No	31	No
CARBON DISULFIDE 75-19	-0 1	1	10	10	10	0.1	NA	NA	NA	NA	7000	No	70000	No
CARBON TETRACHLORIDE 56-23	-5 1	1	2.1	2.1	2.1	0.021	< 0.25	0.8	4.2	NA	1.6	Yes	16	No
CHLOROFORM 67-66	-3 1	1	260	260	260	2.6	< 0.25	1.4	25	NA	1.1	Yes	11	Yes
ETHYLBENZENE 100-4	-4 1	1	0.54 J	0.54 J	0.54	0.0054	< 0.25	7.3	340	NA	22	No	220	No
METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE) 108-1	)-1 1	1	47 J	47 J	47	0.47	< 0.25	2.2	36	NA	800	No	8000	No
METHYLENE CHLORIDE 75-09	-2 1	1	1.6 J	1.6 J	1.6	0.016	< 0.25	22	2100	60	52	No	520	No
TETRACHLOROETHENE (PCE) 127-1	3-4 1	1	32	32	32	0.32	< 0.25	2.9	51	100	8.1	Yes	81	No
TOLUENE 108-8	3-3 1	1	5.7 J	5.7 J	5.7	0.057	< 0.25	58	510	NA	4000	No	40000	No
TRICHLOROETHENE (TCE) 79-0	-6 1	1	3.6	3.6	3.6	0.036	<0.25	0.5	25	5	0.22	Yes	2.2	Yes
XYLENE, m,p- 108-38	3/1 1	1	3.3	3.3	3.3	0.033	< 0.25	12	550	NA	70000	No	700000	No
XYLENE, o- 95-4	-6 1	1	1.1 J	1.1 J	1.1	0.011	<0.25	7.6	310	NA	70000	No	700000	No

Background Indoor Air Value from

Notes:

\* = NYSDOH. 2003. Summary of Indoor and Outdoor Levels of Volatile Organic Compounds from Fuel Oil Heated Homes in NYS, 1997-2003. Revised November 14, 2005.

\*\* = NYSDOH Guidance for Evaluating Soil Vapor Intrusion in the State of New York. Table 3-1 - Air guideline values derived by the NYSDOH. October 2006.

Target shallow Gas Concentration = U.S. Environmental Protection Agency (USEPA). 2002. Draft Guidance for Evaluating the Vapor Intrusion to Indoor Air Pathway from Groundwater and Soils (Subsurface Vapor Intrusion Guidance). November 2002. MDC = Maximum detected concentration

COPI = chemical of potential interest

J = estimated value

TABLE 5 MPA Process Area (Building 2) - COPIs Compared to NY State Background and USEPA Health Protective Levels Soil Vapor Investigation - Buildings 1, 2, 3, 4, and Tank Storage Area Former Hampshire Chemical Corp. Facility, Waterloo, New York

								NYS	SDOH 2003* (µ	g/m3)					
СОРІ	CAS#	Number of Soil Gas Analyses	Number of Soil Gas Detects	Minimum Detected Concentration (μg/m3)	Maximum Detected Concentration (MDC) (μg/m3)	Mean Detected Concentration (μg/m3)	Mean Detected Concentration (Attenuation Factor - 0.01) (μg/m3)	Minimum	90th Percentile	Maximum	NYSDOH Air Guideline Value** (µg/m3)	Target Shallow Gas Concentration (Attenuation Factor - 0.1) (µg/m3)	MDC Exceeds?	Target Shallow Gas Concentration (Attenuation Factor - 0.01) (µg/m3)	MDC Exceeds?
Summary Statistics Presented for COPIs Reported in MPA Pr	ocess Area (Bu	ilding 2) (200	7)												
ACRYLONITRILE	107-13-1	2	2	6.9	27	16.95	0.1695	NA	NA	NA	NA	0.36	Yes	3.6	Yes
BENZENE	71-43-2	2	2	13	90	51.5	0.515	< 0.25	15	460	NA	3.1	Yes	31	Yes
CARBON DISULFIDE	75-15-0	2	2	9.5	34	21.75	0.2175	NA	NA	NA	NA	7000	No	70000	No
CHLOROETHANE	75-00-3	2	1	0.72	0.72	0.72	0.0072	< 0.25	< 0.25	4.5	NA	100000	No	1000000	No
CHLOROFORM	67-66-3	2	2	3.6	25	14.3	0.143	< 0.25	1.4	25	NA	1.1	Yes	11	Yes
CHLOROMETHANE	74-87-3	2	2	0.39	1 J	0.695	0.00695	< 0.25	3.3	260	NA	24	No	240	No
cis-1,2-DICHLOROETHYLENE	156-59-2	2	1	1.4	1.4	1.4	0.014	< 0.25	< 0.25	7.4	NA	350	No	3500	No
ETHYLBENZENE	100-41-4	2	2	0.43 J	0.91 J	0.67	0.0067	< 0.25	7.3	340	NA	22	No	220	No
METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	108-10-1	2	2	240	820	530	5.3	< 0.25	2.2	36	NA	800	No	8000	No
METHYLENE CHLORIDE	75-09-2	2	2	0.43 J	1.3 J	0.865	0.00865	< 0.25	22	2100	60	52	No	520	No
TETRACHLOROETHENE (PCE)	127-18-4	2	2	5.4	8.9	7.15	0.0715	< 0.25	2.9	51	100	8.1	No	81	No
TOLUENE	108-88-3	2	2	16	26	21	0.21	< 0.25	58	510	NA	4000	No	40000	No
TRICHLOROETHENE (TCE)	79-01-6	2	2	2.9	10	6.45	0.0645	< 0.25	0.5	25	5	0.22	Yes	2.2	Yes
XYLENE, m,p-	108-38-3/1	2	2	2.3	4.4 J	3.35	0.0335	< 0.25	12	550	NA	70000	No	700000	No
XYLENE, o-	95-47-6	2	2	0.83 J	1.5 J	1.165	0.01165	<0.25	7.6	310	NA	70000	No	700000	No

Background Indoor Air Value from

Notes:

\* = NYSDOH. 2003. Summary of Indoor and Outdoor Levels of Volatile Organic Compounds from Fuel Oil Heated Homes in NYS, 1997-2003. Revised November 14, 2005.

\*\* = NYSDOH Guidance for Evaluating Soil Vapor Intrusion in the State of New York. Table 3-1 - Air guideline values derived by the NYSDOH. October 2006.

Target shallow Gas Concentration = U.S. Environmental Protection Agency (USEPA). 2002. Draft Guidance for Evaluating the Vapor Intrusion to Indoor Air Pathway from Groundwater and Soils (Subsurface Vapor Intrusion Guidance). November 2002. MDC = Maximum detected concentration

COPI = chemical of potential interest

J = estimated value

TABLE 6 Building 1 - COPIs Compared to NY State Background and USEPA Health Protective Levels Soil Vapor Investigation - Buildings 1, 2, 3, 4, and Tank Storage Area Former Hampshire Chemical Corp. Facility, Waterloo, New York

								NY	SDOH 2003* (µ	g/m3)					
		Number of Soil Gas	Soil Gas	Minimum Detected Concentration	Maximum Detected Concentration (MDC)	Mean Detected Concentration	Mean Detected Concentration (Attenuation Factor - 0.01)		90th		NYSDOH Air Guideline Value**	Target Shallow Gas Concentration (Attenuation Factor - 0.1)	MDC	Target Shallow Gas Concentration (Attenuation	MDC
COPI	CAS#	Analyses	Detects	(µg/m3)	(µg/m3)	(µg/m3)	(µg/m3)	Minimum	Percentile	Maximum	(µg/m3)	(µg/m3)	Exceeds?	Factor - 0.01) (µg/m3)	Exceeds?
Summary Statistics Presented for COPIs Reported in Buildi	ing 1 (2007)														
1,1,1-TRICHLOROETHANE	71-55-6	1	1	12	12	12	0.12	< 0.25	3.1	110	NA	22000	No	220000	No
1,1-DICHLOROETHANE	75-34-3	1	1	2.1	2.1	2.1	0.021	< 0.25	< 0.25	4.4	NA	5000	No	50000	No
1,2-DICHLOROPROPANE	78-87-5	1	1	3.9	3.9	3.9	0.039	< 0.25	< 0.25	34	NA	40	No	400	No
BENZENE	71-43-2	1	1	0.81	0.81	0.81	0.0081	< 0.25	15	460	NA	3.1	No	31	No
CHLOROFORM	67-66-3	1	1	440	440	440	4.4	< 0.25	1.4	25	NA	1.1	Yes	11	Yes
CHLOROMETHANE	74-87-3	1	1	0.62	0.62	0.62	0.0062	< 0.25	3.3	260	NA	24	No	240	No
ETHYLBENZENE	100-41-4	1	1	0.61 J	0.61 J	0.61	0.0061	< 0.25	7.3	340	NA	22	No	220	No
METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	108-10-1	1	1	100	100	100	1	< 0.25	2.2	36	NA	800	No	8000	No
TETRACHLOROETHENE (PCE)	127-18-4	1	1	17	17	17	0.17	< 0.25	2.9	51	100	8.1	Yes	81	No
TOLUENE	108-88-3	1	1	5.8	5.8	5.8	0.058	< 0.25	58	510	NA	4000	No	40000	No
TRICHLOROETHENE (TCE)	79-01-6	1	1	2.5	2.5	2.5	0.025	< 0.25	0.5	25	5	0.22	Yes	2.2	Yes
XYLENE, m,p-	108-38-3/1	1	1	4.3	4.3	4.3	0.043	< 0.25	12	550	NA	70000	No	700000	No
XYLENE, o-	95-47-6	1	1	2.9	2.9	2.9	0.029	<0.25	7.6	310	NA	70000	No	700000	No

Background Indoor Air Value from

Notes:

\* = NYSDOH. 2003. Summary of Indoor and Outdoor Levels of Volatile Organic Compounds from Fuel Oil Heated Homes in NYS, 1997-2003. Revised November 14, 2005.

\*\* = NYSDOH Guidance for Evaluating Soil Vapor Intrusion in the State of New York. Table 3-1 - Air guideline values derived by the NYSDOH. October 2006.

Target shallow Gas Concentration = U.S. Environmental Protection Agency (USEPA). 2002. Draft Guidance for Evaluating the Vapor Intrusion to Indoor Air Pathway from Groundwater and Soils (Subsurface Vapor Intrusion Guidance). November 2002.

MDC = Maximum detected concentration

COPI = chemical of potential interest

J = estimated value

								NY	SDOH 2003* (µg	g/m3)	4				
СОРІ	CAS#	Number of Soil Gas Analyses	Number of Soil Gas Detects	Minimum Detected Concentration (µg/m3)	Maximum Detected Concentration (MDC) (µg/m3)	Mean Detected Concentration (μg/m3)	Mean Detected Concentration (Attenuation Factor - 0.01) (μg/m3)	Minimum	90th Percentile	Maximum	NYSDOH Air Guideline Value** (µg/m3)	Target Shallow Gas Concentration (Attenuation Factor0.1) (µg/m3)	MDC Exceeds?	Target Shallow Gas Concentration (Attenuation Factor - 0.01) (µg/m3)	MDC Exceeds?
Summary Statistics Presented for COPIs Reported in Building	g 3 (2007)														
1,1-DICHLOROETHANE	75-34-3	3	1	1.4	1.4	1.4	0.014	< 0.25	< 0.25	4.4	NA	5000	No	50000	No
BROMODICHLOROMETHANE	75-27-4	3	1	0.43	0.43	0.43	0.0043	NA	NA	NA	NA	1.4	No	14	No
CARBON DISULFIDE	75-15-0	3	2	20	2600	1310	13.1	NA	NA	NA	NA	7000	No	70000	No
CARBON TETRACHLORIDE	56-23-5	3	1	0.55	0.55	0.55	0.0055	< 0.25	8.0	4.2	NA	1.6	No	16	No
CHLOROFORM	67-66-3	3	3	76	1100	542	5.42	< 0.25	1.4	25	NA	1.1	Yes	11	Yes
cis-1,2-DICHLOROETHYLENE	156-59-2	3	1	2.1	2.1	2.1	0.021	< 0.25	< 0.25	7.4	NA	350	No	3500	No
METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	108-10-1	3	2	2.3	110	56.15	0.5615	< 0.25	2.2	36	NA	800	No	8000	No
METHYLENE CHLORIDE	75-09-2	3	1	26 J	2.6 J	2.6	0.026	< 0.25	22	2100	60	52	No	520	No
TETRACHLOROETHENE (PCE)	127-18-4	3	3	3.4 J	4.1	3.7667	0.0377	< 0.25	2.9	51	100	8.1	No	81	No
TRICHLOROETHENE (TCE)	79-01-6	3	2	2.4	20	11.2	0.112	< 0.25	0.5	25	5	0.22	Yes	2.2	Yes
XYLENE, m,p-	108-38-3/1	3	1	7.2 J	7.2	7.2	0.072	< 0.25	12	550	NA	70000	No	700000	No
XYLENE, o-	95-47-6	3	2	0.27 J	2.9 J	1.585	0.01585	<0.25	7.6	310	NA	70000	No	700000	No

Background Indoor Air Value from

COPI = chemical of potential interest

J = estimated value

Notes:
\* = NYSDOH. 2003. Summary of Indoor and Outdoor Levels of Volatile Organic Compounds from Fuel Oil Heated Homes in NYS, 1997-2003. Revised November 14, 2005.

<sup>\*\* =</sup> NYSDOH Guidance for Evaluating Soil Vapor Intrusion in the State of New York. Table 3-1 - Air guideline values derived by the NYSDOH. October 2006.

Target shallow Gas Concentration = U.S. Environmental Protection Agency (USEPA). 2002. Draft Guidance for Evaluating the Vapor Intrusion to Indoor Air Pathway from Groundwater and Soils (Subsurface Vapor Intrusion Guidance). November 2002. MDC = Maximum detected concentration

Building 4 - COPIs Compared to NY State Background and USEPA Health Protective Levels Soil Vapor Investigation - Buildings 1, 2, 3, 4, and Tank Storage Area Former Hampshire Chemical Corp. Facility, Waterloo, New York

								NYS	DOH 2003* (µg	/m3)	1				
COPI	CAS#	Number of Soil Gas Analyses	Number of Soil Gas Detects	Minimum Detected Concentration (μg/m3)	Maximum Detected Concentration (MDC) (μg/m3)	Mean Detected Concentration (μg/m3)	Mean Detected Concentration (Attenuation Factor - 0.01) (μg/m3)	Minimum	90th Percentile	Maximum	NYSDOH Air Guideline Value** (µg/m3)	Target Shallow Gas Concentration (Attenuation Factor - 0.1) (μg/m3)	MDC Exceeds?	Target Shallow Gas Concentration (Attenuation Factor - 0.01) (µg/m3)	MDC Exceeds?
Summary Statistics Presented for COPIs Reported in Building	g 4 (2007)	-			· · · · · ·									, <u>,</u>	
1,1-DICHLOROETHANE	75-34-3	3	1	84	84	84	0.84	<0.25	< 0.25	4.4	NA	5000	No	50000	No
1,2-DICHLOROPROPANE	78-87-5	3	2	2.1	52	27	0.27	<0.25	< 0.25	34	NA	40	No	400	No
ACRYLONITRILE	107-13-1	3	1	3.3 J	3.3 J	3.3	0.033	NA	NA	NA	NA	0.36	Yes	3.6	No
BENZENE	71-43-2	3	1	5	5	5	0.05	<0.25	15	460	NA	3.1	Yes	31	No
CARBON DISULFIDE	75-15-0	3	2	43	3800	1922	19.22	NA	NA	NA	NA	7000	No	70000	No
CHLOROBENZENE	108-90-7	3	1	2.3	2.3	2.3	0.023	<0.25	< 0.25	0.6	NA	600	No	6000	No
CHLOROFORM	67-66-3	3	3	170	24000	8490	84.9	<0.25	1.4	25	NA	1.1	Yes	11	Yes
CHLOROMETHANE	74-87-3	3	1	2.1	2.1	2.1	0.021	<0.25	3.3	260	NA	24	No	240	No
cis-1,2-DICHLOROETHYLENE	156-59-2	3	1	19 J	19 J	19	0.19	<0.25	< 0.25	7.4	NA	350	No	3500	No
ETHYLBENZENE	100-41-4	3	3	1.6 J	870	434	4.34	<0.25	7.3	340	NA	22	Yes	220	Yes
METHYL ETHYL KETONE (2-BUTANONE)	78-93-3	3	1	32	32	32	0.32	<0.25	16	180	NA	10000	No	100000	No
METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	108-10-1	3	3	380 J	86000	32793	327.93	<0.25	2.2	36	NA	800	Yes	8000	Yes
METHYLENE CHLORIDE	75-09-2	3	3	2.3 J	240 J	92	0.92	<0.25	22	2100	60	52	Yes	520	No
TETRACHLOROETHENE (PCE)	127-18-4	3	2	3	26	15	0.15	<0.25	2.9	51	100	8.1	Yes	81	No
TOLUENE	108-88-3	3	1	130	130	130	1.3	<0.25	58	510	NA	4000	No	40000	No
TRICHLOROETHENE (TCE)	79-01-6	3	2	1.7	54	28	0.28	<0.25	0.5	25	5	0.22	Yes	2.2	Yes
XYLENE, m,p-	108-38-3/1	3	3	15	8600	4405	44.05	<0.25	12	550	NA	70000	No	700000	No
XYLENE, o-	95-47-6	3	3	4.3 J	1700	885	8.85	<0.25	7.6	310	NA	70000	No	700000	No

Background Indoor Air Value from

Target shallow Gas Concentration = U.S. Environmental Protection Agency (USEPA). 2002. Draft Guidance for Evaluating the Vapor Intrusion to Indoor Air Pathway from Groundwater and Soils (Subsurface Vapor Intrusion Guidance). November 2002. COPI = chemical of potential interest

J = estimated value

Notes:

<sup>\* =</sup> NYSDOH. 2003. Summary of Indoor and Outdoor Levels of Volatile Organic Compounds from Fuel Oil Heated Homes in NYS, 1997-2003. Revised November 14, 2005.

\*\* = NYSDOH Guidance for Evaluating Soil Vapor Intrusion in the State of New York. Table 3-1 - Air guideline values derived by the NYSDOH. October 2006.

MDC = Maximum detected concentration

Tank Storage Area - COPIs Compared to NY State Background and USEPA Health Protective Levels Soil Vapor Investigation - Buildings 1, 2, 3, 4, and Tank Storage Area Former Hampshire Chemical Corp. Facility, Waterloo, New York

								NYS	SDOH 2003* (μο	g/m3)					
СОРІ	CAS#	Number of Soil Gas Analyses	Number of Soil Gas Detects	Minimum Detected Concentration (µg/m3)	Maximum Detected Concentration (MDC) (μg/m3)	Mean Detected Concentration (μg/m3)	Mean Detected Concentration (Attenuation Factor - 0.01) (μg/m3)	Minimum	90th Percentile	Maximum	NYSDOH Air Guideline Value** (µg/m3)	Target Shallow Gas Concentration (Attenuation Factor - 0.1) (µg/m3)	MDC Exceeds?	Target Shallow Gas Concentration (Attenuation Factor - 0.01) (µg/m3)	MDC Exceeds?
Summary Statistics Presented for COPIs Reported in Tank Sto				(1-9)	(µ3,	(1-97	(1-3)				(µg.me)	(1-9)		( -3,	
1,2-DICHLOROPROPANE	78-87-5	1	1	15	15	15	0.15	< 0.25	< 0.25	34	NA	40	No	400	No
BENZENE	71-43-2	1	1	0.66	0.66	0.66	0.0066	< 0.25	15	460	NA	3.1	No	31	No
CHLOROFORM	74-87-3	1	1	270	270	270	2.7	< 0.25	1.4	25	NA	1.1	Yes	11	Yes
METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	108-10-1	1	1	11 J	11 J	11	0.11	< 0.25	2.2	36	NA	800	No	8000	No
TETRACHLOROETHENE (PCE)	127-18-4	1	1	6.4	6.4	6.4	0.064	< 0.25	2.9	51	100	8.1	No	81	No
TRICHLOROETHENE (TCE)	79-01-6	1	1	520	520	520	5.2	< 0.25	0.5	25	5	0.22	Yes	2.2	Yes
XYLENE, m,p-	108-38-3/1	1	1	1.3 J	1.3 J	1.3	0.013	< 0.25	12	550	NA	70000	No	700000	No
XYLENE, o-	95-47-6	1	1	0.54 J	0.54 J	0.54	0.0054	< 0.25	7.6	310	NA	70000	No	700000	No

Background Indoor Air Value from

\* = NYSDOH. 2003. Summary of Indoor and Outdoor Levels of Volatile Organic Compounds from Fuel Oil Heated Homes in NYS, 1997-2003. Revised November 14, 2005.

\*\* = NYSDOH Guidance for Evaluating Soil Vapor Intrusion in the State of New York. Table 3-1 - Air guideline values derived by the NYSDOH. October 2006.

Target shallow Gas Concentration = U.S. Environmental Protection Agency (USEPA). 2002. Draft Guidance for Evaluating the Vapor Intrusion to Indoor Air Pathway from Groundwater and Soils (Subsurface Vapor Intrusion Guidance). November 2002. MDC = Maximum detected concentration
COPI = chemical of potential interest

J = estimated value

TABLE 10
MW-10 - COPIs Compared to Groundwater Screening Levels Protective of Human Health
Soil Vapor Investigation - Buildings 1, 2, 3, 4, and Tank Storage Area
Former Hampshire Chemical Corp. Facility, Waterloo, New York

		Number of	Number of	Minimum Detected	Maximum  Detected  Concentration		Target Groundwater	
		Groundwater	Groundwater	Concentration	(MDC)	Mean Detected	Concentration	MDC
COPI	CAS#	Detects	Analyses	(µg/L)	(µg/L)	Concentration	(µg/L)	Exceeds?
Summary Statistics Presented for COPIs Reported in MW-1	0 (2002 - 2005)							
ACRYLONITRILE	107-13-1	0	4	N/A	N/A	N/A	8.5	N/A
CARBON TETRACHLORIDE	56-23-5	0	4	N/A	N/A	N/A	5	N/A
CHLOROFORM	67-66-3	0	4	N/A	N/A	N/A	80	N/A
CHLOROMETHANE	74-87-3	0	4	N/A	N/A	N/A	6.7	N/A
METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	108-10-1	2	4	73	260	166.5	14000	No
METHYLENE CHLORIDE	75-09-2	0	4	N/A	N/A	N/A	58	N/A
TETRACHLOROETHENE (PCE)	127-18-4	0	4	N/A	N/A	N/A	5	N/A
TRICHLOROETHENE (TCE)	79-01-6	0	4	N/A	N/A	N/A	5	N/A

Target Groundwater Concentrations = U.S. Environmental Protection Agency (USEPA). 2002. Draft Guidance for Evaluating the Vapor Intrusion to Indoor Air Pathway from Groundwater and Soils (Subsurface Vapor Intrusion Guidance). November 2002. Table 2c Risk = 1x10^6

N/A = Not applicable

J = estimated value

Screen interval = 4 - 14 ft

MW-10 is located outside Building 2-B. SG-1 is located inside Building 2-B.

TABLE 11
MW-11S - COPIs Compared to Groundwater Screening Levels Protective of Human Health
Soil Vapor Investigation - Buildings 1, 2, 3, 4, and Tank Storage Area
Former Hampshire Chemical Corp. Facility, Waterloo, New York

lwater Grou		/A N//	C) Mean Detect Concentration A N/A		MDC Exceeds?
ects And	alyses (μg	<b>ι/L) (μg/</b> /A Ν/ <i>/</i>	/L) Concentrat	tion (µg/L)	Exceeds?
	4 N/	/A N//	A N/A	2200	N/A
				2200	N/A
1	4 N/	/A N//			. 4// (
		/A IN//	A N/A	2.1	N/A
•	4 2.	.4 10.	.8 5.5	560	No
)	4 N/	/A N//	A N/A	5	N/A
)	4 N/	/A N/A	A N/A	80	N/A
)	4 N/	/A N//	A N/A	210	N/A
1	4 5.9	9 J 10	8.5	14000	No
)	4 N/	/A N//	A N/A	58	N/A
)	4 N/	/A N//	A N/A	5	N/A
)	4 N/	/A N//	A N/A	5	N/A
Α	N/A N/	/A N//	A N/A	33000	N/A
Α	N/A N/	/A N//	A N/A	33000	N/A
<b>;</b>	4 1.	.7 20.	9 8.2	22000	No
/		0 4 N 0 4 N 0 4 N 7A N/A N 7A N/A N	0 4 N/A N/A 0 4 N/A N/A 0 4 N/A N/A 0/A N/A N/A N/A	0 4 N/A N/A N/A N/A 0/A N/A N/A N/A N/A N/A	0 4 N/A N/A N/A 58 0 4 N/A N/A N/A 5 0 4 N/A N/A N/A 5 0 4 N/A N/A N/A 5 0/A N/A N/A N/A N/A 33000 0/A N/A N/A N/A N/A N/A 33000

Target Groundwater Concentrations = U.S. Environmental Protection Agency (USEPA). 2002. Draft Guidance for Evaluating the Vapor Intrusion to Indoor Air Pathway from Groundwater and Soils (Subsurface Vapor Intrusion Guidance). November 2002. Table 2c Risk = 1x10^-6

N/A = Not applicable

J = estimated value

m,p-Xylenes was not reported. Total Xylenes were analyzed. The value shown for the m,p-xylenes and total xylenes is the most conservative value. Screen interval = 4 - 14 ft

MW-11S is located outside Building 3. SG-5a is located inside Building 3.

TABLE 12
PZ-1 - COPIs Compared to Groundwater Screening Levels Protective of Human Health Soil Vapor Investigation - Buildings 1, 2, 3, 4, and Tank Storage Area Former Hampshire Chemical Corp. Facility, Waterloo, New York

COPI	CAS#	Number of Groundwater Detects	Number of Groundwater Analyses	Minimum Detected Concentration (µg/L)	Maximum Detected Concentration (MDC) (μg/L)	Mean Detected Concentration	Target Groundwater Concentration (µg/L)	MDC Exceeds?
Summary Statistics Presented for COPIs Reported in PZ-1 (	2007)		•	<b>""</b>	,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,		(10)	
1,1-DICHLOROETHANE	75-34-3	1	1	1.49	1.49	1.49	2200	No
1,2-DICHLOROPROPANE	78-87-5	0	1	N/A	N/A	N/A	35	N/A
ACRYLONITRILE	107-13-1	0	1	N/A	N/A	N/A	8.5	N/A
BENZENE	71-43-2	0	1	N/A	N/A	N/A	5	N/A
CARBON DISULFIDE	75-15-0	1	1	1.56	1.56	1.56	560	No
CHLOROBENZENE	108-90-7	0	1	N/A	N/A	N/A	390	N/A
CHLOROFORM	67-66-3	1	1	101	101	101.0	80	Yes
CHLOROMETHANE	74-87-3	0	1	N/A	N/A	N/A	6.7	N/A
CIS-1,2-DICHLOROETHENE	156-59-2	1	1	18.6	18.6	18.6	210	No
ETHYLBENZENE	100-41-4	0	1	N/A	N/A	N/A	700	N/A
METHYL ETHYL KETONE (2-BUTANONE)	78-93-3	1	1	3.56 J	3.56 J	3.56	440000	No
METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	108-10-1	1	1	8070 J	8070 J	8070	14000	No
METHYLENE CHLORIDE	75-09-2	1	1	21.6	21.6	21.6	58	No
TETRACHLOROETHENE (PCE)	127-18-4	0	1	N/A	N/A	N/A	5	N/A
TOLUENE	108-88-3	0	1	N/A	N/A	N/A	1500	N/A
TRICHLOROETHENE (TCE)	79-01-6	0	1	N/A	N/A	N/A	5	N/A
XYLENES, m,p-	108-38-3/1	N/A	N/A	N/A	N/A	N/A	22000	N/A
XYLENES, o-	95-47-6	N/A	N/A	N/A	N/A	N/A	33000	N/A
XYLENES, total	1330-20-7	0	1	N/A	N/A	N/A	22000	N/A

Target Groundwater Concentrations = U.S. Environmental Protection Agency (USEPA). 2002. Draft Guidance for Evaluating the Vapor Intrusion to Indoor Air Pathway from Groundwater and Soils (Subsurface Vapor Intrusion Guidance). November 2002. Table 2c Risk = 1x10^6

N/A = Not applicable

J = estimated value

m,p-Xylenes was not reported. Total Xylenes were analyzed. The value shown for the m,p-xylenes and total xylenes is the most conservative value.

Screen interval = 9 - 9 ft

PZ-1 is located inside Building 4. SG-7 is located near this PZ-1.

TABLE 13

MW-03 - COPIs Compared to Groundwater Screening Levels Protective of Human Health

Soil Vapor Investigation - Buildings 1, 2, 3, 4, and Tank Storage Area

Former Hampshire Chemical Corp. Facility, Waterloo, New York

		Number of Groundwater	Number of Groundwater	Minimum Detected Concentration	Maximum Detected Concentration (MDC)	Mean Detected	Target Groundwater Concentration	MDC
COPI	CAS#	Detects	Analyses	(μg/L)	(µg/L)	Concentration	(µg/L)	Exceeds?
Summary Statistics Presented for COPIs Reported in MW-0	3 (2004 - 2007)							
1,1-DICHLOROETHANE	75-34-3	3	3	1.79 J	2.21	2.0	2200	No
1,2-DICHLOROPROPANE	78-87-5	1	3	1.77	1.77	1.77	35	No
ACRYLONITRILE	107-13-1	0	3	N/A	N/A	N/A	8.5	N/A
BENZENE	71-43-2	3	3	1.75	1.97	1.8	5	No
CARBON DISULFIDE	75-15-0	3	3	1.34	13.2	6.0	560	No
CHLOROBENZENE	108-90-7	3	3	5.49 J	5.99	5.7	390	No
CHLOROFORM	67-66-3	0	3	N/A	N/A	N/A	80	N/A
CHLOROMETHANE	74-87-3	0	3	N/A	N/A	N/A	6.7	N/A
CIS-1,2-DICHLOROETHENE	156-59-2	3	3	1.97 J	9.31	6.0	210	No
ETHYLBENZENE	100-41-4	3	3	1	1.13	1.0	700	No
METHYL ETHYL KETONE (2-BUTANONE)	78-93-3	3	3	4.9 J	28.8 J	15.4	440000	No
METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	108-10-1	3	3	4.58 J	29.6 J	14.7	14000	No
METHYLENE CHLORIDE	75-09-2	0	3	N/A	N/A	N/A	58	N/A
TETRACHLOROETHENE (PCE)	127-18-4	0	3	N/A	N/A	N/A	5	N/A
TOLUENE	108-88-3	3	3	10.2 J	35.7 J	26.4	1500	No
TRICHLOROETHENE (TCE)	79-01-6	1	3	0.292 J	0.292 J	0.3	5	No
XYLENES, m,p-	108-38-3/1	N/A	N/A	N/A	N/A	N/A	22000	N/A
XYLENES, o-	95-47-6	N/A	N/A	N/A	N/A	N/A	33000	N/A
XYLENES, total	1330-20-7	3	3	5.05	6.12	N/A	22000	N/A

Target Groundwater Concentrations = U.S. Environmental Protection Agency (USEPA). 2002. Draft Guidance for Evaluating the Vapor Intrusion to Indoor Air Pathway from Groundwater and Soils (Subsurface Vapor Intrusion Guidance). November 2002. Table 2c Risk = 1x10^6

N/A = Not applicable

J = estimated value

m,p-Xylenes was not reported. Total Xylenes were analyzed. The value shown for the m,p-xylenes and total xylenes is the most conservative value.

Screen interval = 3 - 16.5 ft (2004 & 2005); 14 - 14 ft (2007

MW-03 is located inside Building 4. SG-7a is located near MW-03.

TABLE 14
BLDG4-PIT-SSP - COPIs Compared to Groundwater Screening Levels Protective of Human Health Soil Vapor Investigation - Buildings 1, 2, 3, 4, and Tank Storage Area Former Hampshire Chemical Corp. Facility, Waterloo, New York

COPI	CAS#	Number of Groundwater Detects	Number of Groundwater Analyses	Minimum Detected Concentration (μg/L)	Maximum Detected Concentration (MDC) (μg/L)	Mean Detected Concentration	Target Groundwater Concentration (μg/L)	MDC Exceeds?
Summary Statistics Presented for COPIs Reported in BLDG			Analyses	(P9'-)	(MB) = )	Concontitution	(M9/ L)	Exceeds:
1,1-DICHLOROETHANE	75-34-3	2	4	0.556 J	2.47 J	1.5	2200	No
1,2-DICHLOROPROPANE	78-87-5	0	4	N/A	N/A	N/A	35	N/A
ACRYLONITRILE	107-13-1	0	4	N/A	N/A	N/A	8.5	N/A
BENZENE	71-43-2	1	4	0.208 J	0.208 J	0.208	5	No
CARBON DISULFIDE	75-15-0	4	4	14.1	123	46	560	No
CHLOROBENZENE	108-90-7	2	4	0.168 J	15.8 J	8.0	390	No
CHLOROFORM	67-66-3	4	4	3.05	611	184	80	Yes
CHLOROMETHANE	74-87-3	0	4	N/A	N/A	N/A	6.7	N/A
CIS-1,2-DICHLOROETHENE	156-59-2	1	4	1.25	1.25	1.25	210	No
ETHYLBENZENE	100-41-4	0	4	N/A	N/A	N/A	700	N/A
METHYL ETHYL KETONE (2-BUTANONE)	78-93-3	3	4	38.6	284	181	440000	No
METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	108-10-1	4	4	31.8	25500	7233	14000	No
METHYLENE CHLORIDE	75-09-2	2	4	16	98.5	57	58	No
TETRACHLOROETHENE (PCE)	127-18-4	0	4	N/A	N/A	N/A	5	N/A
TOLUENE	108-88-3	4	4	2.03	158 J	44	1500	No
TRICHLOROETHENE (TCE)	79-01-6	0	4	N/A	N/A	N/A	5	N/A
XYLENE, m p-	108-38-3/1	N/A	N/A	N/A	N/A	N/A	22000	N/A
XYLENES, o-	95-47-6	N/A	N/A	N/A	N/A	N/A	33000	N/A
XYLENES, total	1330-20-7	1	4	0.458 J	0.458 J	0.46	22000	No

Target Groundwater Concentrations = U.S. Environmental Protection Agency (USEPA). 2002. Draft Guidance for Evaluating the Vapor Intrusion to Indoor Air Pathway from Groundwater and Soils (Subsurface Vapor Intrusion Guidance). November 2002. Table 2c Risk = 1x10^-6

N/A = Not applicable

J = estimated value

m,p-Xylenes was not reported. Total Xylenes were analyzed. The value shown for the m,p-xylenes and total xylenes is the most conservative value.

Screen interval = 0 - 0.5 ft (2005); 2.67 - 2.67 ft (2007)

BLDG4-PIT SSP is a monitoring location inside Building 4. SG-7a was collected from a location near this BLDG4-PIT-SSP.

TABLE 15
PZ-3 - COPIs Compared to Groundwater Screening Levels Protective of Human Health Soil Vapor Investigation - Buildings 1, 2, 3, 4, and Tank Storage Area Former Hampshire Chemical Corp. Facility, Waterloo, New York

					Maximum			
					Detected		Target	
		Number of	Number of	Minimum Detected	Concentration		Groundwater	
		Groundwater	Groundwater	Concentration	(MDC)	Mean Detected	Concentration	MDC
COPI	CAS#	Detects	Analyses	(µg/L)	(µg/L)	Concentration	(µg/L)	Exceeds?
Summary Statistics Presented for COPIs Report	ed in PZ-3 (2007)							
1,2-DICHLOROPROPANE	78-87-5	1	1	0.323 J	0.323 J	0.323	35	No
BENZENE	71-43-2	0	1	N/A	N/A	N/A	5	N/A
CHLOROFORM	67-66-3	0	1	N/A	N/A	N/A	80	N/A
TETRACHLOROETHENE (PCE)	127-18-4	0	1	N/A	N/A	N/A	5	N/A
TRICHLOROETHENE (TCE)	79-01-6	0	1	N/A	N/A	N/A	5	N/A
XYLENES, m,p-	108-38-3/1	N/A	N/A	N/A	N/A	N/A	22000	N/A
XYLENES, o-	95-47-6	N/A	N/A	N/A	N/A	N/A	33000	N/A
XYLENES, total	1330-20-7	0	1	N/A	N/A	N/A	22000	N/A

Target Groundwater Concentrations = U.S. Environmental Protection Agency (USEPA). 2002. Draft Guidance for Evaluating the Vapor Intrusion to Indoor Air Pathway from Groundwater and Soils (Subsurface Vapor Intrusion Guidance). November 2002. Table 2c Risk = 1x10^-6

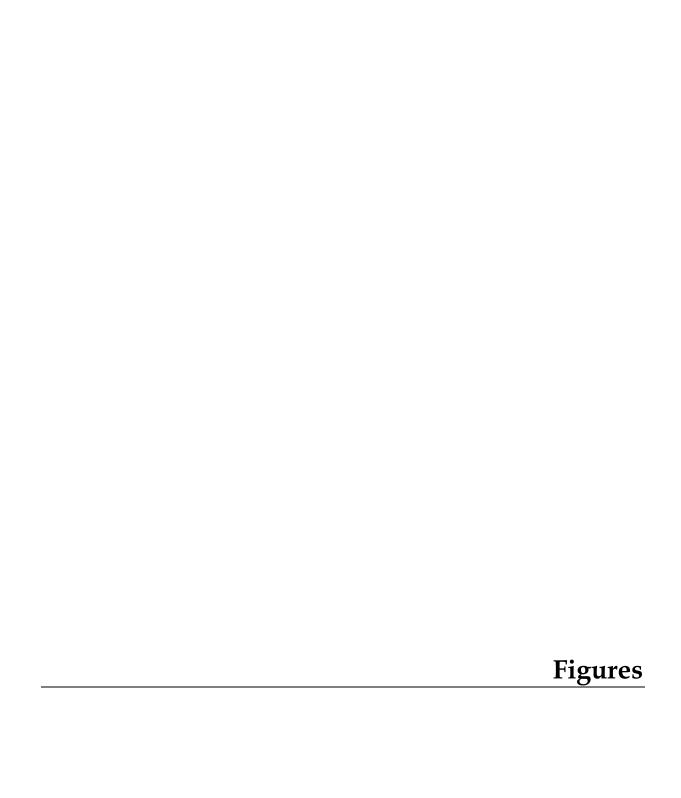
N/A = Not applicable

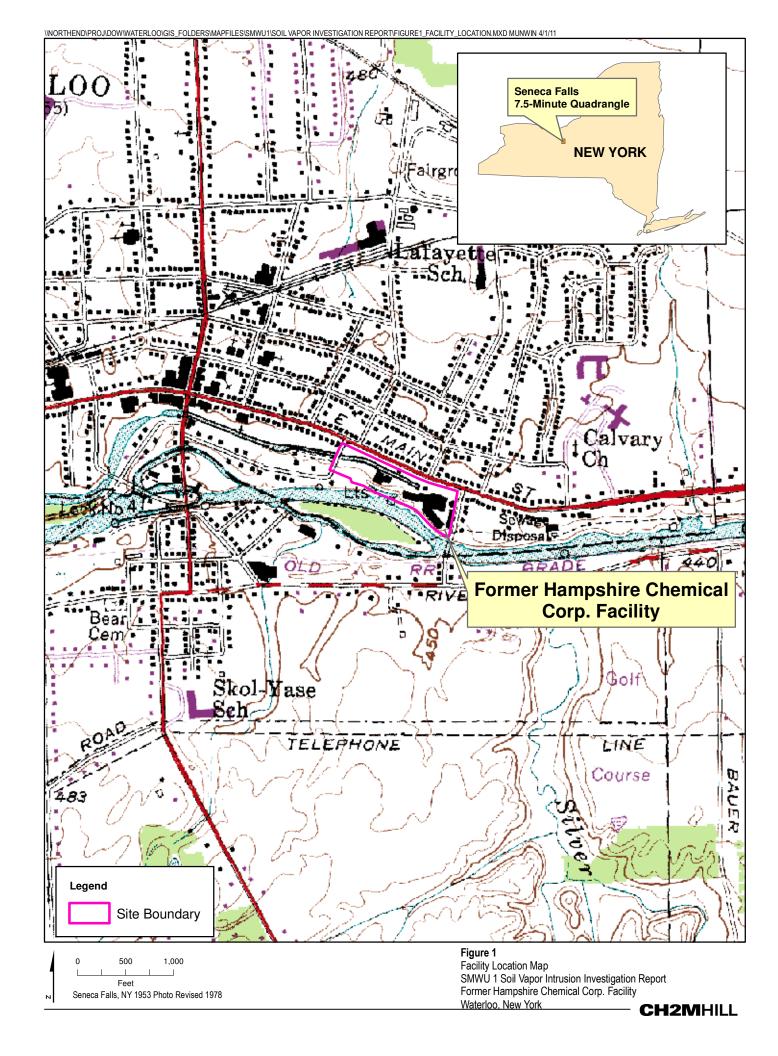
J = estimated value

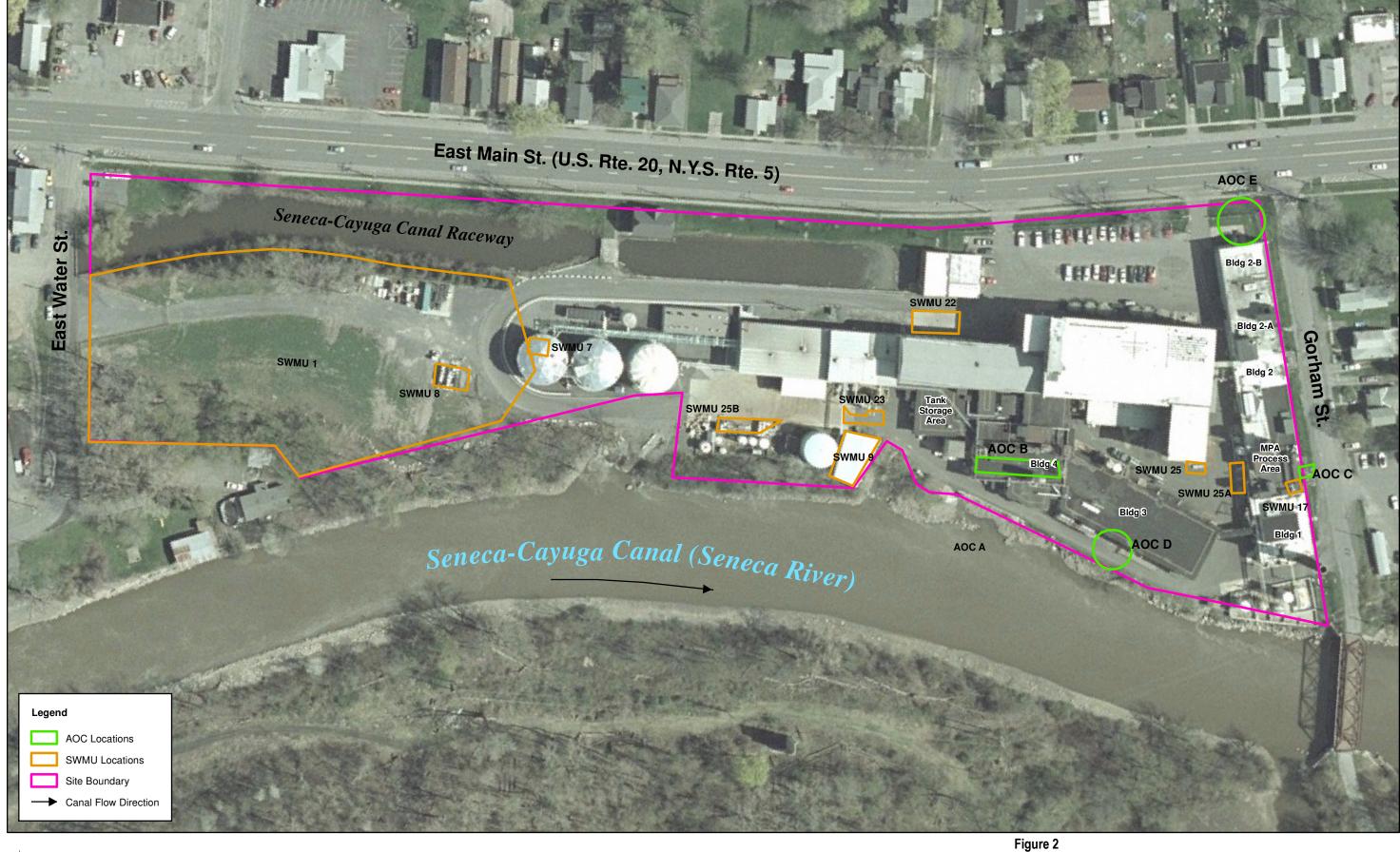
m,p-Xylenes was not reported. Total Xylenes were analyzed. The value shown for the m,p-xylenes and total xylenes is the most conservative value.

Screen interval = 11 - 11 ft

PZ-3 is located inside the Tank Storage area near SG-9.





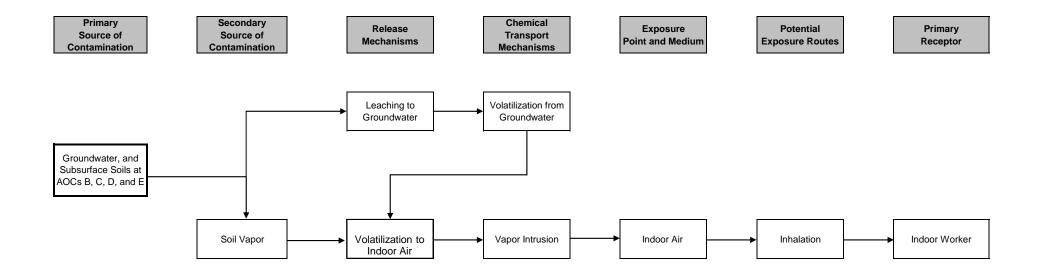


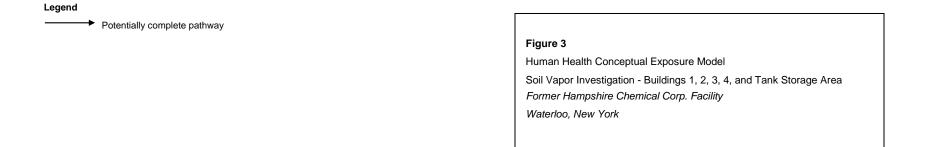


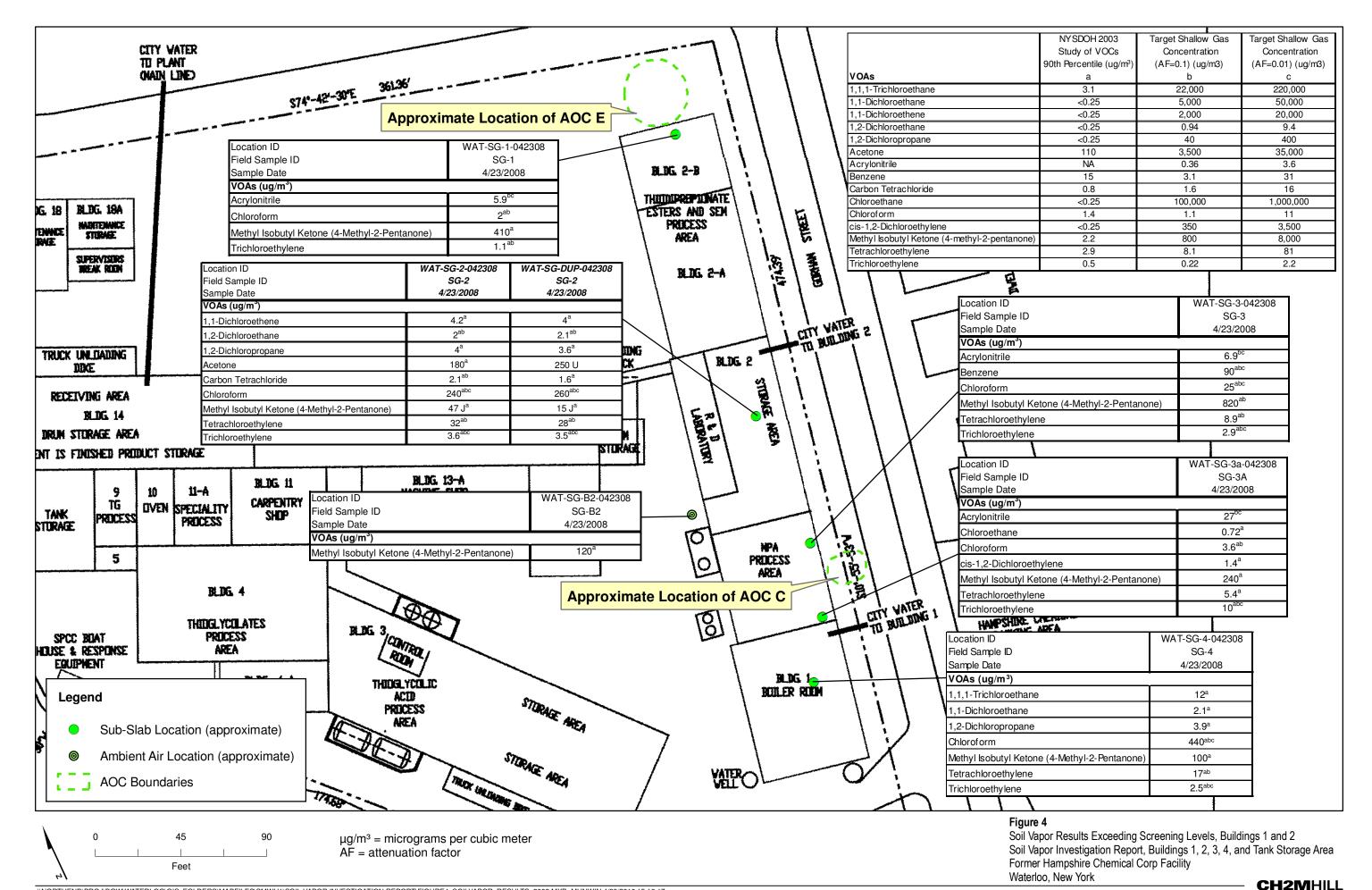
SWMU and AOC Locations

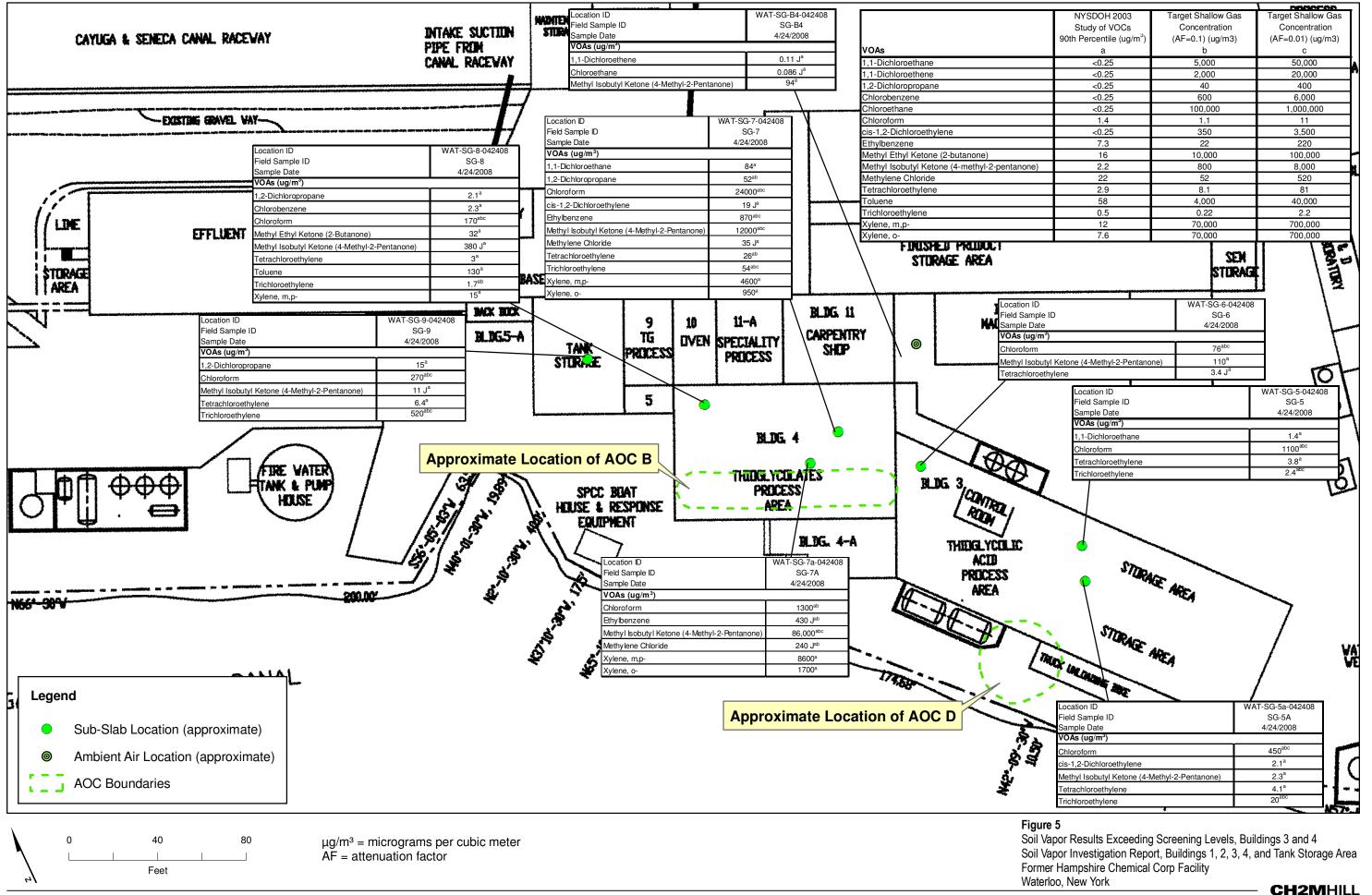
Soil Vapor Investigation Report, Buildings 1, 2, 3, 4, and Tank Storage Area Former Hampshire Chemical Corp Facility
Waterloo, New York

**CH2M**HILL

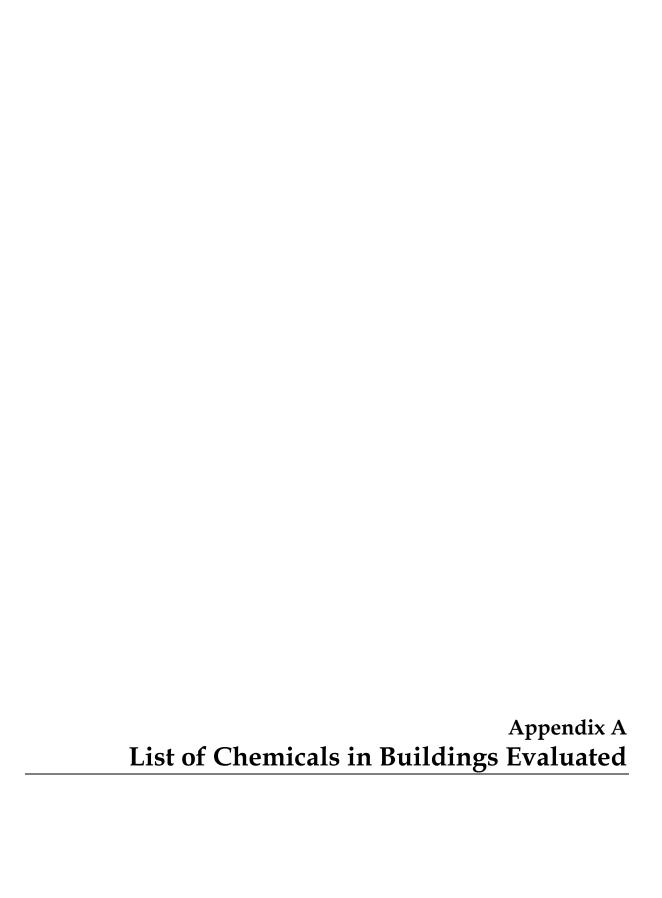












Chemical Inventory - Building 1 - April 2008

Soil Vapor Investigation - Buildings 1, 2, 3, 4, and Tank Storage Area

Former Hampshire Chemical Corp. Facility, Waterloo, New York

## Building 1 - Boiler Room

Quantity	ID	Units	Product Description
37	NA	50 lb bags	Solar salt for water softeners
1	NA	1 pint	Bostik - Never Seez Anti - seize & lubricating compound
1	NA	1 pint	Reactor seal pipe thread sealant
1	NA	16 oz	Top free excel cutting fluid by Winbro
1	NA	8 oz	WD - 40
2	NA	16 oz	Locite thread sealant with teflon
1	NA	18 oz can	ITW Dyman penetrating Oil & Moisture Remover
1	NA	NA	Nalco 1720 Oxygen scavenger
1	I-MT-102	NA	Brine tank
1	NA	NA	Nalco Tri - ACT ® 1820 Corrosion Inhibitor
1	NA	NA	Nex Guard 22310 Boiler water internal treatment

## Aboveground Storage Tanks South of Building 1

Quantity	ID	Units	Product Description
1	16 - HT - 104 - 01	NA	Epichlorohydrin
1	1 - HT - 97 - 01	NA	HCI
1	NA	NA	Acrylinitrile

## Aboveground Storage Tanks West of Buildings 1 and 2

Quantity	ID	Units	Product Description
1	2 - T - 534	NA	Ammonium bisulfate
1	2 - HT -11	Working Capacity - 1,800 gal	MIBK

Notes:

Chemical Inventory - MPA Process Area - April 2008

Soil Vapor Investigation - Buildings 1, 2, 3, 4, and Tank Storage Area

Former Hampshire Chemical Corp. Facility, Waterloo, New York

# Dept 69 - MPA Process Area

Quantity	ID	Units	Product Description
1	2 - AST - 6		Tank
1	AST - 30, 35		Line labeled Thiomalic acid *
		Gross 3900 lbs	
2	2A - PT-47	Tare 477 lbs	3-mercaptopropionic acid (99%)
5	NA	2,600 lbs	MPA Vac Dist
1	NA	NA	MIBK / MPA Water container
1	2 - HT - 78	NA	MPA crude
			Line near MIBK/MPA container labeled MIBK (from tote to
1	NA		acidification reactor)
1	NA	55 gal plastic drum	80% MPA
1	2-AT - 8	NA	MPA Blend Tank AST
1	2 - AST - 35	NA	MPA and Solvent MIBK
1	2-AST-30	NA	MPA and Solvent MIBK
1	2 - AST - 23	NA	MPA and Solvent MIBK
2	NA	2,600 lbs	MPA Cruel
1	NA	NA	Line labeled Sulfuric acid

#### Notes:

<sup>\*</sup> Thiomalic Acid (TMA) has not been produced since the late 1990s to early 2000s.

Chemical Inventory - Building 2 - April 2008

Soil Vapor Investigation - Buildings 1, 2, 3, 4, and Tank Storage Area

Former Hampshire Chemical Corp. Facility, Waterloo, New York

## **Building 2 - Storage Area**

Quantity	ID	Units	Product Description
1	NA	NA	Recycled HCl tank
1	2-T-44	NA	NaMPA Storage tank
1	2-AST-57	NA	Nitrile reactor
1	2-AST-55	NA	Hydro reactor
1	2-T-043	NA	Ammonia Scrubber
1	2-T-045	NA	Caustic Mix Tank
1	2-T-054	NA	Washed MIBK Tank
1	2-T-046	NA	Raffinate Tank
10	NA	20 kg bags	De - icing salt (Kwick melt safety salt, Sel Antiglance)
32	NA	2,600 lbs	VAC Dist MPA
1	2-HT-18	NA	Boiler Tank
1	NA	NA	Buffer solutions; 1.68 buffer; 4.0 buffer; 1.0 buffer

#### Lab Area

Quantity	ID	Units	Product Description
1	NA	5 gal container	NaSH + H2O container
1	NA	5 gal box container	0.1000 Normal NaOH by Anachemia
2	NA	6 gal box container	Sulfuric acid solution by GFS Chemicals Item # 1201
1	NA	4 Liters	Potassium iodate - iodide Item # 2269 by GFS Chemical
		1 gal glass	
1	NA	container	10% H2SO4
1	NA	NA	DI water
1	NA	500 ml	Phenolphthalem
1	NA	NA	Potassium iodide KIO3 crystals

Notes:

Chemical Inventory - Building 2-A - April 2008

Soil Vapor Investigation - Buildings 1, 2, 3, 4, and Tank Storage Area

Former Hampshire Chemical Corp. Facility, Waterloo, New York

Building 2-A - Dept. 69 Thiodipropionate & MPA Derivative Production Unit

Quantity	ID	Units	Product Description
1	2A-HT-74	15,000 gal	Ditridecyl Thiodipropionate AST
1	2A-AST-40	NA	These are reactors that can contain any of the raw materials in the building
1	2A-AST-41	NA	These are reactors that can contain any of the raw materials in the building
1	2A-AST-42	NA	These are reactors that can contain any of the raw materials in the building
1	NA	3,500 gal	Exxal ® 8, Product #769724
27	NA	50lb bags	Pike Creek all purpose granulated salt
1	NA		Ammonium bisulfate AST
1	2A-AST-45		AST
1	2A-AST-46		AST
1	NA	55 gal drum	E-Pure MSA
1	NA	55 gal drum	DTBSP Filter drain
1	NA	55 gal drum	DTBSP Pan and Filter
1	HT 133	NA	Tank
1	A-HT-111	NA	Dilite ITA AST
1	2A-R-38	NA	Conc ITA
1	NA	NA	Line labeled Metacrylic acid
1	2A-R-42	NA	Sulfoethyl methyacrylate (SEM) AST
1	NA	55 gal drum	DPPD Inhibitor
1	2A - H - 44	NA	Small tank
1	2A - H - 47	NA	Small tank
1	2A - HT - 109		
1	2A - HT - 110		
1	2A - HT - 22		Ammonium bisulfate
1	NA	NA	Raceway water line

Notes:

Chemical Inventory - Building 2-B - April 2008

Soil Vapor Investigation - Buildings 1, 2, 3, 4, and Tank Storage Area

Former Hampshire Chemical Corp. Facility, Waterloo, New York

**Building 2-B - Thiodipropionate Esters and SEM Process Area** 

Quantity	ID	Units	Product Description
1	2B HT 99	NA	Fatty Alcohol AST ALFOL - 18 Code 8098
1	2B WT - 1	NA	Fatty Alcohol AST ALFOL - 18
1	2B - T - 22	NA	S-19 Scrubber AST
1	2B - WT - 3	NA	TDPN Weight Tank AST
1	NA	10,500 lbs	Melt Tank AST
1	2B - T1	NA	Fatty Alcohol - Lauryl Alcohol - AST
		Total Capacity - 7600 gal	
1	2B - T2	Working Capacity - 7300 gal	Tridecyl Alcohol Storage Tank
1	NA	55 gal	E-Pure MSA plastic drum
1	2B - T4	NA	TDPN AST (medium)
			ALFOL 12 Code 2098 - Hook up connection outside building where
1	NA	NA	trucks unload

Notes:

Chemical Inventory - Building 3 - April 2008

Soil Vapor Investigation - Buildings 1, 2, 3, 4, and Tank Storage Area

Former Hampshire Chemical Corp. Facility, Waterloo, New York

## **Building 3 - OPWP Team Dept 68 Area**

Quantity	ID	Units	Product Description
1	3-HT-106	Total Capacity 1,500 gal	Sodium hydrosulfide
1	3-HT-106A	Total Capacity 1,000 gal	Sodium hydrosulfide
1	3-HT-66	Total Capacity 5,700 gal	Sodium hydrosulfide
1	3-HT 63	Total Capacity 5,700 gal	Sodium hydrosulfide
1	NA	NA	Line labeled alkali waste to ETP
			Area next to Former Tank Area called Hazardous Waste Scrubber Solution.
			Lines say caustic Feed to ETP (Flow out). Flow in Caustic Feed to Dept 68. Line
1	NA	NA	labeled NaSH filter 1B, NaSH Fill # 7A
4	NA	2600 lb totes	Ethanolamine solutions (UN2491), Monoethanolamine CI/Fe free
1	3-HT-122	NA	AST tank labeled Corrosive toxic
11	NA	2,600 lb totes	Ammonium thioglycolate 71%
1	NA	2,600 lb totes	Glycerol thioglycolate 7010
6	NA	2,600 lb totes	Glycerol thioglycolate 80%
2	NA	2,600 lb totes	Diammonium Dithioglycolate 48% and DAD totes 40%
4	NA	2,600 lb totes	Glyceryl thioglycolate - NK
8	NA	2,600 lb totes	MEAT totes
2	NA	2,600 lb totes	Glyceryl thioglycolate - 75%
1	NA	2,600 lb totes	60% ATG
1	NA	2,600 lb totes	90% TG - Blend
1	NA	2,600 lb totes	98% TG - Blend
1	NA	2000 lbs tote	Thioset - M
1	3-HT-113	Total Capacity 12,480 gal	Monochloroacidic acid
1	3-HT-59	Total Capacity 4,590 gal	Sodium hydrosulfide
1	3-AV-11	NA	NaSH dilution tank
1	3-AV-12	NA	NaSH holding tank
1	3-T-35	NA	Scrubber
1	3-AST-48	NA	H2S Generator

Chemical Inventory - Building 3 - April 2008

Soil Vapor Investigation - Buildings 1, 2, 3, 4, and Tank Storage Area

Former Hampshire Chemical Corp. Facility, Waterloo, New York

### Lab Area

Quantity	ID	Units	Product Description
1	NA	NA	Buffer solutions: 1.68, 4.0, 7.0, 1.0
1	NA	500 ml bottle	Sulfite Form Liquid by Nalco
1	NA	1 liter	Titrating solution by Nalco
1	NA	1 liter	Hardness buffer by Nalco

#### **Near Location #6**

Quantity	ID	Units	Product Description	
1	NA	1 gal	Clorox bleach	
1	NA	NA	Phenolphthalien	
1	NA	18 oz	State Fix "Terg-o-Cide in a can"	
1	NA	12 oz	Krylon Industrial Touch Coat Acylic Enamel Aerosol	
2	NA	21 oz	Trouble shooter by 3M	
1	NA	1 gal	Sherwin Williams Industrial Enamel HS Industrial Marine Coatings in yellow	

## **Outside Building 3 near Canal**

Quantity	ID	Units	Product Description
1	3 -HT- 50	Total Capacity 10,000 gal	Sulfuric acid
1	3-HT-13-1	Total Capacity 10,000 gal	Sulfuric acid

## Outside Building 3 between Bldg 3 and Machine Shop

Quantity	ID	Units	Product Description
1	3 -HT- 136	Total Capacity 9,290 gal	Sodium Hydroxide

Notes:

Chemical Inventory - Building 4-A - April 2008

Soil Vapor Investigation - Buildings 1, 2, 3, 4, and Tank Storage Area

Former Hampshire Chemical Corp. Facility, Waterloo, New York

# Buildings 4A and 4B

Quantity	ID	Units	Product Description		
1	4-HT-941	Total Capacity 5,200	Crude Thoglycolic Acid (TGA)		
1	4-HT-951	Total Capacity 5,200	Crude TGA		
1	NA	2,600 lb tote	1st Pass TGA Residue		
1	AV - 11	NA	T-Acid Residue Holding Tank		
1	AV - 8	NA	Could not see label; related to AV - 1		

## **Building 4 - Dept 68 T-Acid & Derivative Production Unit**

Quantity	ID	Units	Product Description	
1	4-V-6	NA	Recycled MIBK AST	
1	4-V-7	NA	Washed MIBK	
5	NA	2,000 lb tote	Thioglycolic acid (TGA) 96-99% totes	
4	NA	2,600 lb tote	DAD	
9	NA	55 gal drums	Vac Dist TG (blue plastic drums)	
9	NA	55 gal drums	TG concentrate	
1	4-HT-7	NA	Regular T-Acid Extract	
1	NA	5 gal container	Sulfuric acid	
1	NA	55 gal drum	TG -Forerun	
1	NA	5 gal container	TG NF	
2	NA	330 gal	Ammonium hydroxide 26 BE	
1	NA	2,600 lb tote	Glyceryl Thioglycolate	
1	4-AST-7	NA	TGA 16 - 99% AST	
1	HT-121	NA	T-Acid Extract Residue	

Chemical Inventory - Building 4-A - April 2008

Soil Vapor Investigation - Buildings 1, 2, 3, 4, and Tank Storage Area

Former Hampshire Chemical Corp. Facility, Waterloo, New York

### Lab Area

Quantity	ID	Units	Product Description		
1	NA	NA	Buffer solutions 1.68, 4.0, 1.0		
1	NA	5 gal container	Potassium iodate - Iodide by GFS Chemicals		
1	NA	NA	Potassium iodide (KI)		
1	NA	NA	10% sulfuric acid solution		

# Outside Building 4 near Canal

Quantity	ID	Units	Product Description	
2	NA	NA	Compressed nitrogen cylinders by Jackson Welding Supply Co.	

Notes:

Chemical Inventory - Tank Storage Area - April 2008

Soil Vapor Investigation - Buildings 1, 2, 3, 4, and Tank Storage Area

Former Hampshire Chemical Corp. Facility, Waterloo, New York

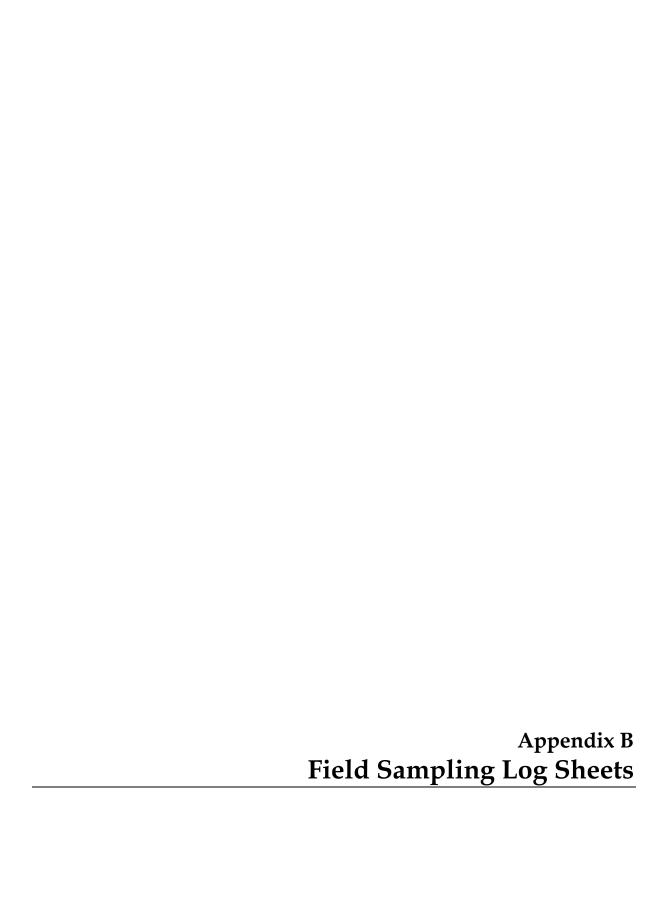
# **Tank Storage Area**

Quantity	ID	Units	Product Description
1	7-HT-105	Total Capacity 15,000 gal	Sodium hydrosulfide
28	NA	55 gal drums	Hydrogen peroxide
1	7-HT-128	Total Capacity 6,000 gal	Glycerin
1	7-HT-127	Total Capacity 6,000 gal	Glycerol monothioglycolate
1	7-HT-4	Total Capacity 5,000 gal	Ammonium thioglycolate
1	7-HT-5	Total Capacity 2,600 gal	I-MCH AST
4	NA	55 gal drums	Thioglycerol (TGNF)
4	NA	NA	Compressed Nitrogen cylinders by Jackson Welding Supply Co.

# Outside Tank Storage Area

Quantity	ID	Units	Product Description
NA	NA	NA	MPA Fill Hopper Storage Area

Notes:



# Indoor Vapor Intrusion Assessment Sub Slab Vapor Sampling Field Log

Project Info							
Project Name: Former Hamp  By: Rachel Kopec	oshire Chemical Corp Fa c/NJO; Lisa La Fortune/N			369548.B2.RP 4/23/2008			
Structure							
	amples and field blank						
Address:							
Sample Location type:							
concrete slab on grade		Yard or Driveway	•				
☐ concrete footing w/crawl space		other (describe)	See identification				
☐ basement							
Soil Vapor Sampling System							
Probe type (describe):	n/a						
Probe to sample interface syste	em (describe):						
Sample collection type:	Syringe	☐ Tedla	ır bag	mma canister			
Other info (describe other aspects)							
Soil Vapor Probe Purging & S	ampling I og						
Sample location (show in diagram)	AA 1	FB 1	AA 2				
		<b>.</b>					
Sample Identification (field ID)	WAT-SG-B2-042308	WAT-SG-FB-042308	WAT-SG-B4-042408				
Date Installed*	n/a	n/a	n/a				
Depth of installed probe (feet bgs)	Ambient Air	Field Blank	Ambient Air				
Leak check (probe/sampling interface)	n/a	n/a	n/a				
Calculated dead volume (1 purge volume), ml							
Purge rate, ml/min.							
Purge duration, min. (3 volumes)							
Purge vacuum, " Hg							
Max PID Reading, ppmv (optional)							
Purge completed (time of day)							
Sampling period started (time of day)	8:27	9:55	7:55				
Sampling rate, ml/min	11.25	n/a	11.25				
Sampling vacuum, " Hg (initial/end)	30/6	0/0	30/6.5				
Sampling period ended (time of day)	16:27	17:55	15:55				
Observations and Comments:							
n/a - Not applicable							
AA 1 corresponds to grass surf							
AA 2 correspond to asphalt paved area between Building 4 and the machine shop.							

# Indoor Vapor Intrusion Assessment Sub Slab Vapor Sampling Field Log

Project Info								
Project Name: Former Hamp	Project Name: Former Hampshire Chemical Corp Facility, Waterloo, NY Project #: 369548.B2.RP							
By: Rachel Koped	/NJO; Lisa La Fortune/N	NO	Date:	4/23/2008				
Structure								
Identification: Buildings 2-B,	2, and MPA Process Ar	ea						
Address:								
Sample Location type:  Concrete slab on grade  Concrete slab on grade  Yard or Driveway								
concrete footing w/crawl space	· 🗆	other (describe)						
basement								
Soil Vapor Sampling System								
	Stainless steel tubing, fi	nished as a flushmount						
Probe to sample interface syste	em (describe):	Swagelock and stainles	s steel tubing					
Sample collection type:	☐ Syringe	☐ Tedla	ar bag 🔟 Su	mma canister				
Other info (describe other aspects)								
Soil Vapor Probe Purging & Sa	ampling Log							
Sample location (show in diagram)	1	2	2	3				
Sample Identification (field ID)	WAT-SG-1-042308	WAT-SG-2-042308	WAT-SG-DUP-042308	WAT-SG-3-042308				
Date Installed*	4/23/2008	4/23/2008	4/23/2008	4/23/2008				
Depth of installed probe (inches bgs)	8.75	10.0	10.0	9.25				
Leak check (probe/sampling interface)	\ !! /	ok (0 ppm)	ok (0 ppm)	ok (0 ppm)				
Calculated dead volume (1 purge volume), ml	17	17	17	17				
Purge rate, ml/min.	100	100	100	100				
Purge duration, min. (3 volumes)	30 sec	30 sec	30 sec	30 sec				
Purge vacuum, " Hg								
Max PID Reading, ppmv (optional)								
Purge completed (time of day)	8:33	8:59 	-	9:17				
Sampling period started (time of day)	8:38	9:04	-	9:21				
Sampling rate, ml/min	11.25	11.25	11.25	11.25				
Sampling vacuum, " Hg (initial/end)	30/0	30/8	30/0	30/0				
Sampling period ended (time of day)	16:38	17:04	-	17:21				
Observations and Comments:	WAT-SG-DUP-042308	is a duplicate sample of	WAT-SG-2-042308					

### Indoor Vapor Intrusion Assessment Sub Slab Vapor Sampling Field Log

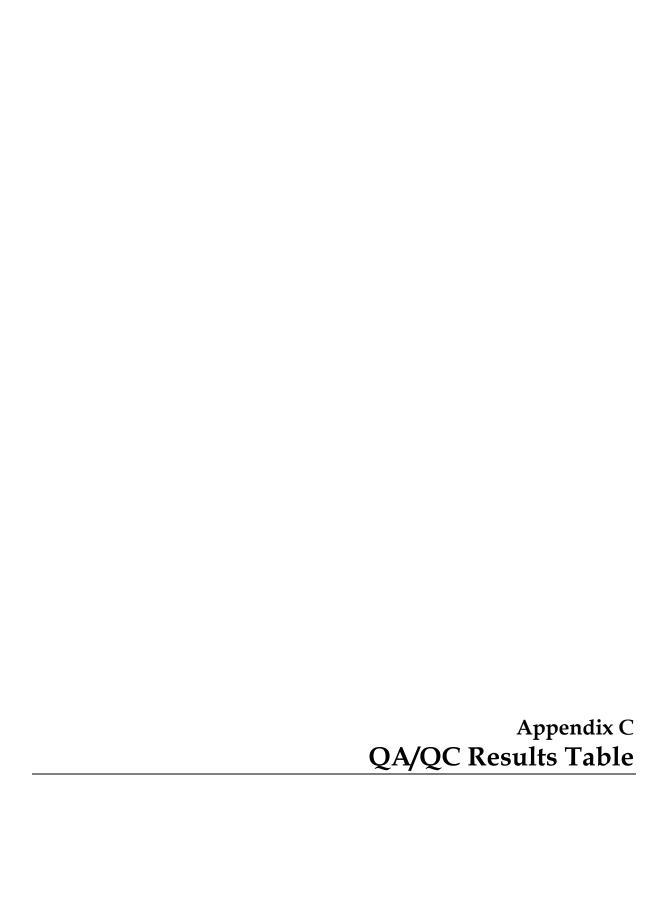
Project Info						
	Project Name: Former Hampshire Chemical Corp Facility, Waterloo, NY       Project #: 369548.B2.RP         By: Rachel Kopec/NJO; Lisa La Fortune/NJO       Date: 4/23/2008					
Structure						
Identification: MPA Process Area and Building 1 Address: Sample Location type:						
<ul><li>concrete slab on grade</li><li>concrete footing w/crawl space</li><li>basement</li></ul>	□ • □	Yard or Driveway other (describe)				
Soil Vapor Sampling System						
Probe type (describe):	Stainless steel tubing, fi	nished as a flushmount				
Probe to sample interface syste	em (describe):	Swagelock and stainles	s steel tubing			
Sample collection type:	Syringe	☐ Tedla	ar bag 🔣 S	umma canister		
Other info (describe other aspects)						
Soil Vapor Probe Purging & Sa	ampling I og					
Sample location (show in diagram)	4	5				
Sample Identification (field ID)	WAT-SG-3a-042308			<b></b>		
Date Installed*	4/23/2008	4/23/2008				
Depth of installed probe (feet bgs)	10.0	9.5				
Leak check (probe/sampling interface)	ok (0 ppm)	ok (0 ppm)				
Calculated dead volume (1 purge volume), ml	17	17				
Purge rate, ml/min.	100	100				
Purge duration, min. (3 volumes)	30 sec	30 sec				
Purge vacuum, " Hg						
Max PID Reading, ppmv (optional)						
Purge completed (time of day)	9:29	9:49		. <b></b>		
Sampling period started (time of day)	9:33	9:54				
Sampling rate, ml/min	11.25	11.25				
Sampling vacuum, " Hg (initial/end)	26/5	30/0				
Sampling period ended (time of day)	17:33	17:54				
Observations and Comments:						

### Indoor Vapor Intrusion Assessment Sub Slab Vapor Sampling Field Log

Project Info							
Project Name: Former Hamp	Project Name: Former Hampshire Chemical Corp Facility, Waterloo, NY Project #: 369548.B2.RP						
By: Rachel Koped	:/NJO; Lisa La Fortune/N	NO	Date:	4/24/2008			
Structure							
Identification: Buildings 3 ar	nd 4						
Address:							
Sample Location type:   concrete slab on grade		Yard or Driveway					
concrete footing w/crawl space	· 🗆	other (describe)					
basement							
Soil Vapor Sampling System							
Probe type (describe):	Stainless steel tubing, fi	nished as a flushmount					
Probe to sample interface syste	em (describe):	Swagelock and stainless	steel tubing				
Sample collection type:	☐ Syringe	☐ Tedla	r bag	ımma canister			
Other info (describe other aspects)							
Soil Vapor Probe Purging & Sa	ampling Log						
Sample location (show in diagram)	5	5a	6	7			
Sample Identification (field ID)	WAT-SG-5-042408	WAT-SG-5a-042408	WAT-SG-6-042408	WAT-SG-7-042408			
Date Installed*	4/24/2008	4/24/2008	4/24/2008	4/24/2008			
Depth of installed probe (inches bgs)	7.0	8.5	10.0	13.5			
Leak check (probe/sampling interface)	ok (0 ppm)	ok (0 ppm)	ok (0 ppm)	ok (0 ppm)			
Calculated dead volume (1 purge volume), ml	16	17	17	18			
Purge rate, ml/min.	100	100	100	100			
Purge duration, min. (3 volumes)	27 sec	30 sec	29 sec	30 sec			
Purge vacuum, " Hg							
Max PID Reading, ppmv (optional)							
Purge completed (time of day)	7:55	8:03	8:17	8:27			
Sampling period started (time of day)	7:59	8:08	8:21	8:32			
Sampling rate, ml/min	11.25	11.25	11.25	11.25			
Sampling vacuum, " Hg (initial/end)	27/4	30/0	29/0	30/11			
Sampling period ended (time of day)	15:59	16:08	16:21	16:32			
Observations and Comments:							

### Indoor Vapor Intrusion Assessment Sub Slab Vapor Sampling Field Log

Project Info						
Project Name:       Former Hampshire Chemical Corp Facility, Waterloo, NY       Project #: 369548.B2.RP         By:       Rachel Kopec/NJO; Lisa La Fortune/NJO       Date: 4/24/2008						
Structure						
Identification: Building 4 and Tank Storage Area  Address:  Sample Location type:  Concrete slab on grade  Concrete slab on grade  Concrete slab on grade  Concrete slab on grade						
concrete footing w/crawl space		other (describe)				
Soil Vapor Sampling System						
Probe type (describe):	Stainless steel tubing, fi	nished as a flushmount				
Probe to sample interface syste	em (describe):	Swagelock and stainless	s steel tubing			
Sample collection type:	☐ Syringe	☐ Tedla	r bag	ımma canister		
Other info (describe other aspects)						
Soil Vapor Probe Purging & Sa	ampling Log					
Sample location (show in diagram)	7a	8*	9			
Sample Identification (field ID)	WAT-SG-7a-042408	WAT-SG-8-042408	WAT-SG-9-042408			
Date Installed*	4/24/2008	4/24/2008	4/24/2008			
Depth of installed probe (inches bgs)	15.00	16.75	8.00			
Leak check (probe/sampling interface)	ok (0 ppm)	ok (0 ppm)	ok (0 ppm)			
Calculated dead volume (1 purge volume), ml	19	20	16			
Purge rate, ml/min.	100	100	100			
Purge duration, min. (3 volumes)	30 sec	30 sec	28 sec			
Purge vacuum, " Hg						
Max PID Reading, ppmv (optional)						
Purge completed (time of day)	8:50	8:59	9:20			
Sampling period started (time of day)	8:54	9:06	9:24			
Sampling rate, ml/min	11.25	11.25	11.25			
Sampling vacuum, " Hg (initial/end)	30/0	30/10	28/0			
Sampling period ended (time of day)	16:54	17:30	17:24			
Observations and Comments:  * Water was noted in tubing flow controller were switched		ble collection. Gauge ind	icated sample was not co	ollecting. Gauge and		



Appendix C Quality Assurance/Quality Control Sample Results - April 2008 Sampling Event Soil Vapor Investigation - Buildings 1, 2, 3, 4, and Tank Storage Area Former Hampshire Chemical Corp. Facility, Waterloo, New York

Field Sample ID				WAT-SG-2-042308	WAT-SG-DUP-042308	WAT-SG-FB-042308
Location ID				SG-2	SG-2	QC
Sample Date				4/23/2008	4/23/2008	4/23/2008
Sample Type Lab Sample ID		NYSDOH	NYSDOH	Normal P0801215-003	Duplicate P0801215-007	Equip Blank P0801215-008
Dilution		2003	2003	3.4	12.5	1
Matrix		Guidance	Guidance	AIR	AIR	AIR
		а	b			
Volatile Organics, by Method TO15 (UG/M3)	CAS#	90th Percentile	Maximum			
1,1,1-TRICHLOROETHANE	71-55-6	3.1	110	0.34 U	1.3 U	0.1 U
1,1,2,2-TETRACHLOROETHANE 1,1,2-TRICHLOROETHANE	79-34-5 79-00-5	<0.25 <0.25	2.7 6.2	0.34 U 0.34 U	1.3 U 1.3 U	0.1 U 0.1 U
1,1-DICHLOROETHANE	75-34-3	<0.25	4.4	0.34 U	1.3 U	0.1 U
1,1-DICHLOROETHENE	75-35-4	<0.25	430	4.2ª	<b>4</b> <sup>a</sup>	0.1 U
1,2-DICHLOROETHANE	107-06-2	<0.25	4.9	2ª	2.1ª	0.1 U
1,2-DICHLOROPROPANE	78-87-5	<0.25	34	4 <sup>a</sup>	3.6ª	0.1 U
ACETONE	67-64-1	110	690	180ª	250 U	2.2 JB
ACRYLONITRILE BENZENE	107-13-1 71-43-2	NA 15	NA 460	1.2 J 0.93	6.3 U 1.3 U	0.5 U 0.1 U
BROMODICHLOROMETHANE	75-27-4	NA	NA	0.93 0.34 U	1.3 U	0.1 U
BROMOFORM	75-25-2	NA NA	NA NA	1.7 U	6.3 U	0.5 U
BROMOMETHANE	74-83-9	0.6	23	0.34 U	1.3 U	0.1 U
CARBON DISULFIDE	75-15-0	NA	NA	10	8.6 U	0.44 J
CARBON TETRACHLORIDE	56-23-5	0.8	4.2	2.1ª	1.6ª	0.1 U
CHLOROBENZENE CHLOROETHANE	108-90-7 75-00-3	<0.25 <0.25	0.6 4.5	0.34 U 0.34 U	1.3 U 1.3 U	0.1 U 0.1 U
CHLOROFORM	67-66-3	<0.25 1.4	4.5 25	240 <sup>ab</sup>	260 <sup>ab</sup>	0.1 U
CHLOROMETHANE	74-87-3	3.3	260	0.34 U	1.3 U	0.1 U
cis-1,2-DICHLOROETHYLENE	156-59-2	<0.25	7.4	0.34 U	1.3 U	0.1 U
cis-1,3-DICHLOROPROPENE	10061-01-5	<0.25	3.5	1.7 U	6.3 U	0.5 U
DIBROMOCHLOROMETHANE	124-48-1	NA	NA	0.34 U	1.3 U	0.1 U
ETHYLBENZENE	100-41-4	7.3	340	0.54 J	6.3 U	0.5 U
METHYL ETHYL KETONE (2-BUTANONE) METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE	78-93-3 108-10-1	16 2.2	180 36	4.7 U <b>47 J</b> <sup>ab</sup>	2.5 U <b>15 J</b> <sup>a</sup>	0.36 J 0.096 J
METHYLENE CHLORIDE	75-09-2	22	2100	1.1 J	1.6 J	0.090 J
STYRENE	100-42-5	1.3	50	0.3 J	6.3 U	0.5 U
TETRACHLOROETHYLENE (PCE)	127-18-4	2.9	51	32 a	28ª	0.1 U
TOLUENE	108-88-3	58	510	5.7 J	2.1 U	0.24 J
TRANS-1,2-DICHLOROETHENE	156-60-5	NA	NA	0.34 U	1.3 U	0.1 U
TRANS-1,3-DICHLOROPROPENE TRICHLOROETHYLENE (TCE)	10061-02-6 79-01-6	<0.25 0.5	<0.25 25	1.7 U 3.6 <sup>a</sup>	6.3 U <b>3.5</b> <sup>a</sup>	0.5 U 0.1 U
VINYL CHLORIDE	75-01-6	<0.25	1	0.34 U	1.3 U	0.1 U
XYLENE, m,p-	108-38-3/1	12	550	3.3	6.3 U	0.5 U
XYLENE, o-	95-47-6	7.6	310	1.1 J	6.3 U	0.5 U
EPICHLOROHYDRIN	106-89-8	NA	NA	0 UJ	0 UJ	0 UJ
Tentatively Identified Compounds (UG/M3)						
1,1'-Oxybispentane;Amyl Ether (C10H22O)	TIC	NA	NA			
1-METHYLNAPHTHALENE 2,2'-Oxybispentane (C10H22O)	90-12-0 TIC	NA NA	NA NA			
2-METHYLNAPHTHALENE	91-57-6	NA NA	NA NA			
3-METHYLPENTANE	96-14-0	NA NA	NA NA			
Acetaldehyde + Isobutane	TIC	NA	NA		-	
BENZALDEHYDE	100-52-7	NA	NA			
BUTANE	106-97-8	NA	NA	-	-	
C11H24 Branched Alkane	TIC	NA NA	NA NA		-	
C12H26 Branched Alkane C14H28 Compound	TIC	NA NA	NA NA	 40 J		
C6HF13 (Tridecafluorohexane)	TIC	NA NA	NA NA	100 J	50 J	
Chlorodifluoromethane + Propene + Propane	TIC	NA NA	NA NA			
Dimethyl Sulfide	75-18-3	NA	NA			
HEXAMETHYLCYCLOTRISILOXANE	541-05-9	NA	NA	30 J		
Isooctanol	26952-21-6	NA NA	NA			
NAPHTHALENE n-Butyl Ether	91-20-3 TIC	NA NA	NA NA			
n-Butyl Ether n-DODECANE	112-40-3	11	420	 10 J		
n-Nonanal	124-19-6	NA	NA			
n-Nonanal + C10H14 Aromatic	TIC	NA	NA	20 J		
n-UNDECANE	1120-21-4	12	290			
PROPANE	74-98-6	NA	NA		-	
Propane + Propene	TIC	NA NA	NA			
Propane +Carbonyl Sulfide Unidentified Oxygenated Compound	TIC	NA NA	NA NA			
отпоститей Охуденатей Соттроини	110	INA	INA	-	-	-

<sup>\* =</sup> New York State Department of Health (NYSDOH) 2003: Study of volatile organic chemicals in air of fuel oil heated homes.

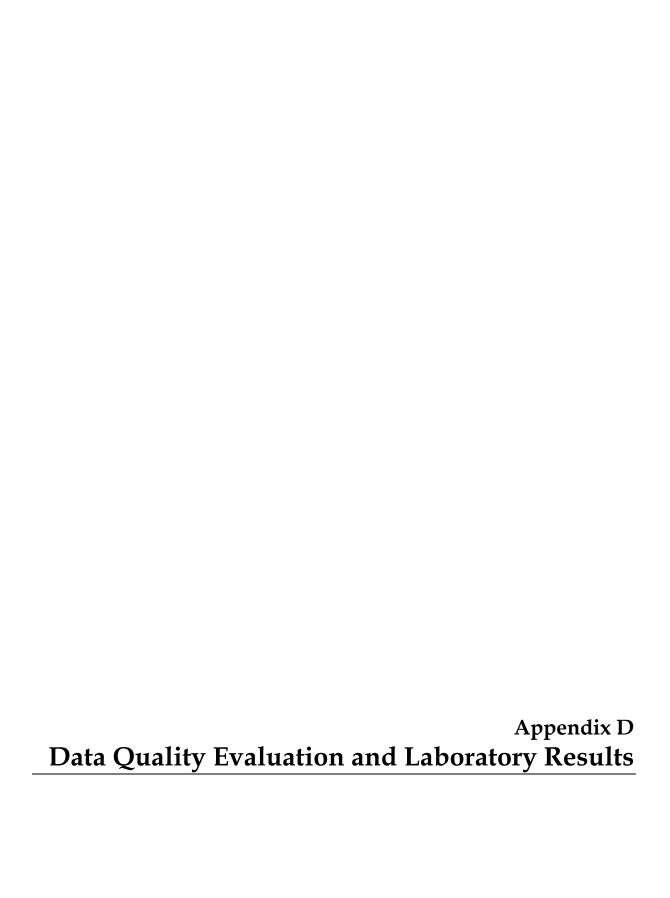
All compounds and criteria are in ug/m3

Shading in gray indicates that constituent was detected above the guidance value NA = not applicable ND = compound not detected at or above method detection limit

NF = Compound was searched for, but not found

J = estimated value

U = compound not detected at or above method detection limit



### Data Quality Evaluation for the April 2008 Soil Vapor Investigation, Waterloo, New York

PREPARED FOR: Paul Taylor/CH2M HILL

PREPARED BY: Berney Kidd/CH2M HILL

DATE: May 30, 2008

### Introduction

The objective of this Data Quality Evaluation (DQE) report is to assess the data quality of analytical results for soil vapor samples collected at the former Hampshire Chemical Corp. facility located in Waterloo, New York. Guidance for this DQE report came from the *U.S. Environmental Protection Agency (USEPA) Contract Laboratory National Functional Guidelines (NFG) for Organic Data Review, October 1999*, individual method requirements and historical laboratory quality control limits.

This report is intended as a general data quality assessment designed to summarize data issues.

### **Analytical Data**

This DQE report covers 14 soil vapor samples, one field duplicate (FD), and one ambient blank (FB). Samples were collected on April 23 and April 24, 2008. The samples were reported as one sample delivery group, P0801215. Samples were collected and shipped by overnight carrier to Columbia Analytical Services in Simi Valley, California. The samples were analyzed for volatile organic compounds by method TO-15.

The sample delivery groups were assessed by reviewing the following: (1) the chain-of-custody documentation; (2) holding-time compliance; (3) initial and continuing calibration criteria; (4) method blanks and field blanks; (5) laboratory control samples (LCS); (6) surrogate spike recoveries; (7) internal standard recoveries; and (8) the required quality control (QC) samples at the specified frequencies.

Data flags were assigned according to the NFG. Multiple flags are routinely applied to specific sample method/matrix/analyte combinations, but there will only be one final flag. A final flag is applied to the data and is the most conservative of the applied validation flags. The final flag also includes matrix and blank sample impacts.

The data flags are those listed in the NFG and are defined below:

• J = The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.

1

- R = The sample result was rejected due to serious deficiencies in the ability to analyze the sample and meet the QC criteria. The presence or absence of the analyte could not be verified.
- U = The analyte was analyzed for but was not detected above the reported sample quantitation limit.
- UJ = The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.

### **Findings**

The overall summaries of the data validation are contained in the following sections and table below.

### **Holding Time**

All holding time criteria were met.

### Calibration

Initial and continuing calibration analyses were performed as required by the methods and all acceptance criteria were met.

The 4-methyl-2-pentanone results for samples WAT-SG-8-042408 and WAT-SG-9-042408 were reported above the linear calibration range. The results were qualified as estimated and flagged "J".

### Method Blanks

Method blanks were analyzed at the required frequency and were free of contamination with the following exception:

Acetone was detected at a concentration less than the reporting limit (RL) in the method blanks. The associated data were qualified as non-detected results and flagged "U" when the sample concentrations were less than 10 times the blank concentrations.

### Field Blanks

An ambient blank was collected and analyzed at the required frequency and was free of contamination with the following exceptions:

Acetone; 2-butanone; 4-methyl-2-pentanone; carbon disulfide; and, toluene were detected at concentrations less than the RL in the ambient blank. The associated data were qualified as non-detected results and flagged "U" when the sample concentrations were less than five times (10 times for acetone and 2-butanone) the blank concentrations.

### **Laboratory Control Samples**

LCSs were analyzed as required and all accuracy criteria were met.

### **Surrogates**

All surrogate acceptance criteria were met.

### Internal Standards

All internal standard acceptance criteria were met.

### Field Duplicates

A FD set was collected and analyzed and precision criteria were met with the following exceptions:

The relative percent difference (RPD) between the native sample and FD exceeded criteria for several 2-butanone, toluene and 4-methyl-2-pentanone. The detected results in the normal and FD were qualified as estimated and flagged "J".

### **Tentatively Identified Compounds**

Tentatively identified compounds (TIC) were reported by the laboratory, either as detected TICs or as non-detected compounds for which a library search was performed. All TIC data were qualified as estimated; non-detected results were flagged "UJ" and detected results were flagged "J".

### **Chain of Custody**

Required procedures were followed and were free of errors.

### **Overall Assessment**

The goal of this assessment is to demonstrate that a sufficient number of representative samples were collected and the resulting analytical data can be used to support the decision making process. The procedures for assessing the precision, accuracy, representativeness, completeness and comparability were modeled after the *USEPA Contract Laboratory NFG for Organic Data Review, October 1999.* The following summary highlights the PARCC findings for the above-defined events:

Precision of the data was verified through the review of the field and laboratory data quality indicators that include FD RPD. Precision was generally acceptable with three analytes being qualified as estimated due to FD RPD exceedances.

Accuracy of the data was verified through the review of the calibration data, LCSs, and surrogate standards. Accuracy was generally acceptable. Two results were reported as estimated due to calibration issues. TIC data were reported are considered estimated.

Representativeness of the data was verified through the sample's collection, storage and preservation procedures, verification of holding-time compliance and evaluation of method/field blank data. All data were reported from analyses within the EPA recommended holding time. The method/field blank samples were generally free of

contamination with a few compounds being qualified as non-detected results due to low level detections in the method and/or ambient blanks.

Comparability of the data was ensured through the use of standard EPA analytical procedures and standard units for reporting. Results obtained are comparable to industry standards in that the collection and analytical techniques followed approved, documented procedures.

Completeness is a measure of the number of valid measurements obtained in relation to the total number of measurements planned. Completeness is expressed as the percentage of valid or usable measurements compared to planned measurements. Valid data are defined as all data that are not rejected for project use. All data were considered valid.

**Table 1 - Validation Flags** 

Field ID	Method	Analyte	Final Result	Units	Final Flag	Reason
WAT-SG-DUP-042308	TO15	Acetone	250	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-DUP-042308	TO15	4-Methyl-2-pentanone	15	UG/M3	J	FD>RPD
WAT-SG-DUP-042308	TO15	C6HF13 (Tridecafluorohexane)	50	UG/M3	J	TIC
WAT-SG-DUP-042308	TO15	Epichlorohydrin	0	UG/M3	UJ	TIC
WAT-SG-DUP-042308	TO15	Carbon Disulfide	8.6	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-DUP-042308	TO15	2-Butanone	2.5	UG/M3	U	FD>RPD
WAT-SG-DUP-042308	TO15	2-Butanone	2.5	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-DUP-042308	TO15	Toluene	2.1	UG/M3	U	FD>RPD
WAT-SG-DUP-042308	TO15	Toluene	2.1	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-7a-042408	TO15	Acetone	1600	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-7a-042408	TO15	Acetone	1600	UG/M3	U	LB <rl< td=""></rl<>
WAT-SG-7a-042408	TO15	Toluene	380	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-7a-042408	TO15	n-Butyl Ether	30000	UG/M3	J	TIC
WAT-SG-7a-042408	TO15	2,2'-Oxybispentane	30000	UG/M3	J	TIC
WAT-SG-7a-042408	TO15	Unidentified Oxygenated Compound	10000	UG/M3	J	TIC
WAT-SG-7a-042408	TO15	1,1'-Oxybispentane;Amyl Ether	10000	UG/M3	J	TIC
WAT-SG-7a-042408	TO15	Epichlorohydrin	0	UG/M3	UJ	TIC
WAT-SG-B2-042308	TO15	Acetone	17	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-B2-042308	TO15	Benzaldehyde	20	UG/M3	J	TIC
WAT-SG-B2-042308	TO15	Epichlorohydrin	0	UG/M3	UJ	TIC
WAT-SG-B2-042308	TO15	Toluene	1.2	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-B2-042308	TO15	Carbon Disulfide	0.71	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-7-042408	TO15	Acetone	1800	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-7-042408	TO15	Carbon Disulfide	310	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-7-042408	TO15	2-Butanone	17	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-7-042408	TO15	Toluene	72	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-7-042408	TO15	Epichlorohydrin	0	UG/M3	UJ	TIC
WAT-SG-5a-042408	TO15	Acetone	12	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-5a-042408	TO15	Acetone	12	UG/M3	U	LB <rl< td=""></rl<>
WAT-SG-5a-042408	TO15	C6HF13 (Tridecafluorohexane)	50	UG/M3	J	TIC
WAT-SG-5a-042408	TO15	n-Nonanal	20	UG/M3	J	TIC
WAT-SG-5a-042408	TO15	Naphthalene	50	UG/M3	J	TIC
WAT-SG-5a-042408	TO15	2-Methylnaphthalene	70	UG/M3	J	TIC
WAT-SG-5a-042408	TO15	1- Methylnaphthalene	70	UG/M3	J	TIC
WAT-SG-5a-042408	TO15	Epichlorohydrin	0	UG/M3	UJ	TIC
WAT-SG-5a-042408	TO15	2-Butanone	1.3	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-5a-042408	TO15	Toluene	0.55	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-9-042408	TO15	Acetone	13	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-9-042408	TO15	Acetone	13	UG/M3	U	LB <rl< td=""></rl<>
WAT-SG-9-042408	TO15	4-Methyl-2-pentanone	11	UG/M3	J	>ICLinearRange
WAT-SG-9-042408	TO15	Propane	10	UG/M3	J	TIC
WAT-SG-9-042408	TO15	Acetaldehyde + Isobutane	30	UG/M3	J	TIC
WAT-SG-9-042408	TO15	C6HF13 (Tridecafluorohexane)	70	UG/M3	J	TIC

Field ID	Method	Analyte	Final Result	Units	Final Flag	Reason
WAT-SG-9-042408	TO15	Unidentified Oxygenated Compound	30	UG/M3	J	TIC
WAT-SG-9-042408	TO15	Epichlorohydrin	0	UG/M3	UJ	TIC
WAT-SG-9-042408	TO15	Carbon Disulfide	3.8	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-9-042408	TO15	2-Butanone	1.3	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-9-042408	TO15	Toluene	1.6	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-2-042308	TO15	4-Methyl-2-pentanone	47	UG/M3	J	FD>RPD
WAT-SG-2-042308	TO15	C6HF13 (Tridecafluorohexane)	100	UG/M3	J	TIC
WAT-SG-2-042308	TO15	n-Nonanal + C10H14 Aromatic	20	UG/M3	J	TIC
WAT-SG-2-042308	TO15	Hexamethylcyclotrisiloxane	30	UG/M3	J	TIC
WAT-SG-2-042308	TO15	n-Dodecane	10	UG/M3	J	TIC
WAT-SG-2-042308	TO15	C14H28 Compound	40	UG/M3	J	TIC
WAT-SG-2-042308	TO15	Epichlorohydrin	0	UG/M3	UJ	TIC
WAT-SG-2-042308	TO15	2-Butanone	4.7	UG/M3	U	FD>RPD
WAT-SG-2-042308	TO15	2-Butanone	4.7	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-2-042308	TO15	Toluene	5.7	UG/M3	J	FD>RPD
WAT-SG-4-042308	TO15	Acetone	17	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-4-042308	TO15	Acetone	17	UG/M3	U	LB <rl< td=""></rl<>
WAT-SG-4-042308	TO15	C6HF13 (Tridecafluorohexane)	80	UG/M3	J	TIC
WAT-SG-4-042308	TO15	C12H26 Branched Alkane	40	UG/M3	J	TIC
WAT-SG-4-042308	TO15	C11H24 Branched Alkane	50	UG/M3	J	TIC
WAT-SG-4-042308	TO15	C11H24 Branched Alkane	60	UG/M3	J	TIC
WAT-SG-4-042308	TO15	n-Undecane	70	UG/M3	J	TIC
WAT-SG-4-042308	TO15	Epichlorohydrin	0	UG/M3	UJ	TIC
WAT-SG-4-042308	TO15	Carbon Disulfide	2.2	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-4-042308	TO15	2-Butanone	3.7	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-6-042408	TO15	Acetone	26	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-6-042408	TO15	Acetone	26	UG/M3	U	LB <rl< td=""></rl<>
WAT-SG-6-042408	TO15	Propane +Carbonyl Sulfide	100	UG/M3	J	TIC
WAT-SG-6-042408	TO15	Dimethyl Sulfide	1000	UG/M3	J	TIC
WAT-SG-6-042408	TO15	Epichlorohydrin	0	UG/M3	UJ	TIC
WAT-SG-6-042408	TO15	Toluene	4.8	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-3a-042308	TO15	Acetone	56	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-3a-042308	TO15	Propane + Propene	30	UG/M3	J	TIC
WAT-SG-3a-042308	TO15	C6HF13 (Tridecafluorohexane)	200	UG/M3	J	TIC
WAT-SG-3a-042308	TO15	n-Nonanal	10	UG/M3	J	TIC
WAT-SG-3a-042308	TO15	Hexamethylcyclotrisiloxane	20	UG/M3	J	TIC
WAT-SG-3a-042308	TO15	C14H28 Compound	40	UG/M3	J	TIC
WAT-SG-3a-042308	TO15	Epichlorohydrin	0	UG/M3	UJ	TIC
WAT-SG-3a-042308	TO15	2-Butanone	4.3	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-1-042308	TO15	Acetone	17	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-1-042308	TO15	Acetone	17	UG/M3	U	LB <rl< td=""></rl<>
WAT-SG-1-042308	TO15	Chlorodifluoromethane + Propene + Propane	20	UG/M3	J	TIC
WAT-SG-1-042308	TO15	Hexamethylcyclotrisiloxane	10	UG/M3	J	TIC
WAT-SG-1-042308	TO15	C6HF13 (Tridecafluorohexane)	10	UG/M3	J	TIC
WAT-SG-1-042308	TO15	Hexamethylcyclotrisiloxane	40	UG/M3	J	TIC

Field ID	Method	Analyte	Final Result	Units	Final Flag	Reason
WAT-SG-1-042308	TO15	Hexamethylcyclotrisiloxane	50	UG/M3	J	TIC
WAT-SG-1-042308	TO15	Epichlorohydrin	0	UG/M3	UJ	TIC
WAT-SG-1-042308	TO15	Carbon Disulfide	1.2	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-1-042308	TO15	2-Butanone	3.9	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-1-042308	TO15	Toluene	1.5	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-5-042408	TO15	Acetone	10	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-5-042408	TO15	Acetone	10	UG/M3	U	LB <rl< td=""></rl<>
WAT-SG-5-042408	TO15	4-Methyl-2-pentanone	3	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-5-042408	TO15	C6HF13 (Tridecafluorohexane)	80	UG/M3	J	TIC
WAT-SG-5-042408	TO15	Hexamethylcyclotrisiloxane	100	UG/M3	J	TIC
WAT-SG-5-042408	TO15	Hexamethylcyclotrisiloxane	30	UG/M3	J	TIC
WAT-SG-5-042408	TO15	Epichlorohydrin	0	UG/M3	UJ	TIC
WAT-SG-5-042408	TO15	Carbon Disulfide	2.4	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-5-042408	TO15	2-Butanone	0.97	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-5-042408	TO15	Toluene	0.85	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-B4-042408	TO15	Propane + Propene	20	UG/M3	J	TIC
WAT-SG-B4-042408	TO15	Butane	10	UG/M3	J	TIC
WAT-SG-B4-042408	TO15	Hexamethylcyclotrisiloxane	10	UG/M3	J	TIC
WAT-SG-B4-042408	TO15	Benzaldehyde	9	UG/M3	J	TIC
WAT-SG-B4-042408	TO15	n-Nonanal	40	UG/M3	J	TIC
WAT-SG-B4-042408	TO15	Epichlorohydrin	0	UG/M3	UJ	TIC
WAT-SG-B4-042408	TO15	Carbon Disulfide	0.92	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-B4-042408	TO15	2-Butanone	0.98	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-3-042308	TO15	Acetone	130	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-3-042308	TO15	C6HF13 (Tridecafluorohexane)	70	UG/M3	J	TIC
WAT-SG-3-042308	TO15	Epichlorohydrin	0	UG/M3	UJ	TIC
WAT-SG-3-042308	TO15	2-Butanone	2.8	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-FB-042308	TO15	Epichlorohydrin	0	UG/M3	UJ	TIC
WAT-SG-8-042408	TO15	Acetone	230	UG/M3	U	FB <rl< td=""></rl<>
WAT-SG-8-042408	TO15	4-Methyl-2-pentanone	380	UG/M3	J	>ICLinearRange
WAT-SG-8-042408	TO15	Isooctanol	400	UG/M3	J	TIC
WAT-SG-8-042408	TO15	3-Methylpentane	300	UG/M3	J	TIC
WAT-SG-8-042408	TO15	Unidentified Oxygenated Compound	600	UG/M3	J	TIC
WAT-SG-8-042408	TO15	C12H26 Branched Alkane	200	UG/M3	J	TIC
WAT-SG-8-042408	TO15	C11H24 Branched Alkane	100	UG/M3	J	TIC
WAT-SG-8-042408	TO15	Epichlorohydrin	0	UG/M3	UJ	TIC

LB<RL = Laboratory blank concentration below reporting limit
FB<RL = Field blank concentration below reporting limit
FD>RPD = Field duplicate relative percent difference criteria exceeded
>ICLinearRange = Result above instrument linear calibration range
TIC = Tentatively identified compound



### LABORATORY REPORT

Simi Valley, California 93065

May 14, 2008

Lisa La Fortune CH2M Hill 119 Cherry Hill Road, Suite 300 Parsippany, NJ 07054

RE: DOW - Waterloo, NY / 369548.B2.FI

Dear Lisa:

Enclosed are the results of the samples submitted to our laboratory on April 28, 2008. For your reference, these analyses have been assigned our service request number P0801215.

All analyses were performed in accordance with our laboratory's quality assurance program. Results are intended to be considered in their entirety and apply only to the samples analyzed and reported herein. Your report contains **%1** pages.

Columbia Analytical Services, Inc. is certified by the California Department of Health Services, NELAP Laboratory Certificate No. 02115CA; Arizona Department of Health Services, Certificate No. AZ0694; Florida Department of Health, NELAP Certification E871020; New Jersey Department of Environmental Protection, NELAP Laboratory Certification ID #CA009; New York State Department of Health, NELAP NY Lab ID No: 11221; Oregon Environmental Laboratory Accreditation Program, NELAP ID: CA20007; The American Industrial Hygiene Association, Laboratory #101661; Department of the Navy (NFESC); Pennsylvania Registration No. 68-03307. Each of the certifications listed above have an explicit Scope of Accreditation that applies to specific matrices/methods/analytes; therefore, please contact me for information corresponding to a particular certification.

If you have any questions, please call me at (805) 526-7161.

Respectfully submitted,

Late Squel

Columbia Analytical Services, Inc.

Kate Aguilera Project Manager



Client: Project: CH2M Hill

DOW - Waterloo, NY / 369548.B2.FI

CAS Project No: NJ Certification ID: P0801215

CA009

### **CASE NARRATIVE**

The samples were received intact under chain of custody on April 28, 2008 and were stored in accordance with the analytical method requirements. Please refer to the sample acceptance check form for additional information. The results reported herein are applicable only to the condition of the samples at the time of sample receipt.

### Volatile Organic Compound Analysis

The samples were analyzed for selected volatile organic compounds and tentatively identified compounds in accordance with EPA Method TO-15 from the Compendium of Methods for the Determination of Toxic Organic Compounds in Ambient Air, Second Edition (EPA/625/R-96/010b), January, 1999. The analytical system was comprised of a gas chromatograph / mass spectrometer (GC/MS) interfaced to a whole-air preconcentrator.

The Method Blank associated with the samples analyzed on 5/2/08 and 5/5/08 contained low levels of Acetone and Carbon Disulfide below the Method Reporting Limit (MRL). Concentrations reported below the MRL are considered estimated and below what would generally be reported; therefore, the sample results are not significantly affected. However, the data is flagged to indicate the presence of the target analytes.

The results of analyses are given in the attached laboratory report. All results are intended to be considered in their entirety, and Columbia Analytical Services, Inc. (CAS) is not responsible for utilization of less than the complete report.

Client:

CH2M Hill DOW - Waterloo, NY 369548.B2.FI Project:

### Detailed Sample Information

Folder: P0801215

Order#	8489	8489				antiba.	8489		8489	8489		8489	8489	
FC ID	FC00693	oa00585 OA00611	oa00594	oa00590 oa00624	oa00621	oa00617	FC00659	oa00572	OA00591	OA00589	oa00616	OA00587	OA00601	oa00584
Order#	8489	8489 8489	8489	8489 <b>8489</b>	8489	8489	8489	8489	8489	8489	8489	8489	8489	8489
Cont ID	AC00812	AC01311 AC00950	AC01189	AC01387 AC00648	AC01168	AC01209	AC00097	AC00802	AC00615	AC00814	AC01276	AC01211	AC01235	AC01020
Pi2 Pi2 (Hg) (psig) Pf2														
Pf1 (F	3.5	<b>3.5</b> 3.5	3.5	3.5 3.6	3.6	3.5	3.5	3.5	3.5	3.5	3.5	3.5	3.6	3.6
Pi1 (psig)	-2.7	0.0	0.1	-2.2 0.0	-0.1	-14.3	-3.1	-1.7	0.0	0.0	-3.5	0.2	-3.9	0.3
Pi1 (Hg)	-5.4	0.0 -5.6		4.4 -0.1	-0.3	-29.2	-6.4	-3.5	0.0	0:0	-7.2		-8.0	
CAS Sample ID Client Sample ID Container Type	1	308 6.0 L-Summa Canister Ambient 308 6.0 L-Summa Canister Ambient	308 6.0 L-Summa Canister Ambient		WAT-SG-DUP-04230 6.0 L-Summa Canister Ambient	WAT-SG-FB-042308 6.0 L-Summa Canister Ambient	2408 6.0 L-Summa Canister Ambient		2408 6.0 L-Summa Canister Ambient	1650023	į.	2408 6.0 L-Summa Canister Ambient	408 6.0 L-Summa Canister Ambient	408 6.0 L-Summa Canister Ambient
Client Sample	WAT-SG-B2-042308	WAT-SG-1-042308 WAT-SG-2-042308	WAT-SG-3-042308	WA1-SG-3a-042308 WAT-SG-4-042308	WAT-SG-DUP-C	WAT-SG-FB-04.	WAT-SG-B4-042408	WAT-SG-5-042408	WAT-SG-5a-042408	WAT-SG-6-042408	WAT-SG-7-042408	WAT-SG-7a-042408	WAT-SG-8-042408	WAT-SG-9-042408
CAS Sample ID	P0801215-001.01	P0801215-002.01 P0801215-003.01	P0801215-004.01	P0801215-005.01 P0801215-006.01	P0801215-007.01	P0801215-008.01	P0801215-009.01	P0801215-010.01	P0801215-011.01	P0801215-012.01	P0801215-013.01	P0801215-014.01	P0801215-015.01	P0801215-016.01

CH2M Hill Client:

DOW - Waterloo, NY 369548.B2.FI Project: Detailed Sample Information

CAS Sample ID Client Sample ID Container Type

Miscellaneous Items - received

AVG00528 AVG00807 AVG00794

AVG00582 AVG00633

AVG00463

OA00607

AVG00466 AVG00484 AVG00793 AVG00683

AVG00758 AVG00754 AVG00469 AVG00462 AVG00333 AVG00696

AVG00686

AC00940

Pi2 (Hg) Pf1 Pi1 (psig)

띮

Pi2 (psig)

Order#

Order #

FC ID

Folder: P0801215

Cont ID

Pf2

\$5/12/2008 8:58:49AM



# Air - Chain of Custody Record & Analytical Service Request

Page of

2655 Park Center Drive, Suite A Simi Valley, California 93065 Phone (805) 526-7161 Fax (805) 526-7270

80801215 CAS Project No. Requested Turnaround Time in Business Days (Surcharges) please circle 1 Day (100%) 2 Day (75%) 3 Day (50%) 4 Day (35%) 5 Day (25%) C10 Day - Standard

to of cogo 7 e.g. Actual Preservative or specific instructions Analysis Method and/or Analytes Late Agailera CAS Contact Perobose croler) Sample Volume 9 0 14601335 101400601 ACCOUNT OF 000 659 400076 0A00616 OAGOSTIA 13000814 000589 Acoust 10Acoss7 400000 Flow Controller 0400550 AC01209 0480617 PCOOF S OASST ACOURT 1019 00594 14.50000 ACO 1168 0400631 AC00950 0400611 Down Woterdoo, n.Y 世 (Bar Code # -AC, SC, etc.) A01387 AC00684 Accord Acooli Canister ID 6180021 ACO1311 369548.82. 927815 P.O. # / Billing Information Sample Type (Air/Tube/ Solid) Sampler (Print & Sign) Rochelhoger 子で T.J. Air 4 Project Number Project Name Date Time Collected Collected 0911 22 (0) solhely 0838 40%0 0933 12/1/03 10808 元式 6520 post57 19/106 OS32 7290 1660 1080 124/00 Jossi 4/03/06 4/23/05 413/68 4/23/63 30/86/2 100 VOV 123/08 20/80/12 4/23/08 Company Name & Address (Reporting Information) 119 Ohemy HITZA, Sulk 300 973-334-11847 Rechel. Kopec QCHDM.com Laboratory ID Number WATER -SG-DVO-HOOR 9-03 6.19.L WAT-SG-R-04406(5)-80 WAT-SG-Ta-043464 19402 O 54 0.0-77 (3).7110-0.0 (D-C) A. to. ( 84 H 10-10 H 0.35 (0-0i through No cooperat Stecay Felth-Back Email Address for Result Reporting WAT- SG-22-04338 WAT-S0-6-04248 WAT-SG-4-043808 1AT-56-52-042406 Report Tier Levels - please select WAT-SC+FB-043308 WAT-56-5-042408 WAT-SG-7-0470F 30/ACHO-1/8-475-14M 808060-1-98-14M WAT-SU-3-04208 MAT-S6-2-042308 WAT-S6-82-042309 CHOW THE 972-316-3300 Project Manager Client Sample ID

Ter 1 - (Results/Default if not specified) ier II - (Results + QC)

Received by: (Signature) Received by: (Signature) Time: 800 Time: Date: Relinquished by: (Signature) Relinquished by: (Signature)

Defrance 编 FED EX

Cooler / Blank

Date: Time:

EDD Units: & S

EDD required (Yes// No Type: Lob Auc L

Tier III - (Data Validation Package) 10% Surcharge \_\_\_\_\_ Tier V - (client specified) \_\_\_\_\_\_

Date:#

Project Requirements (MRLs, QAPP)

CAS Project No. POSO (US

# Air - Chain of Custody Record & Analytical Service Request

2655 Park Center Drive, Suite A Simi Valley, California 93065 Phone (805) 526-7161

Columbia Analytical Services <sup>MC</sup>

n Employee - Owned Company

Requested Turnaround Time in Business Days (Surcharges) please eircle 1 Day (100%) 2 Day (75%) 3 Day (50%) 4 Day (35%) 5 Day (25% 10 Day - Standard Fax (805) 526-7270

e.g. Actual Preservative or specific instructions Project Requirements (MRLs, QAPP) Comments Cooler / Blank 朝に上て Time: Time: Analysis Method and/or Analytes Time: EDD Units: Last / M. 3 Sole Aguillepec CAS Contact EDD required Kes I No Type: Cop Spec 7 CO Hinnouse Sample Volume 70 Received by: (Signature) (Signature) Canister ID <sup>4</sup> Flow Controller Bar Code # - (Bar Code -CASOSTY Received by: (Signature) Down Worker 100, NY 369548, BJ. FT (Bar Code # -AC, SC, etc.) SCOLODA OCOLODA Tier III - (Data Validation Package) 10% Surcharge \_ Tier V - (client specified) 927815 P.O. # / Billing Information Sample Type (Air/Tube/ Solid) Sampler (Print & Sign) Date 1/1/01 Time; for Kahal Karac ARC Project Number Project Name Date Time Collected Collected 4000 40/4/4 504/11) Company Name & Address (Reporting Information)
CHOM HM
CHOM HM
CHOM HM Ref. Sutt 30c Kichel Kope Octom. com 973-316-9360 978-334-5847 Laboratory ID Number Procious, No oral Staces Felts-Book Email Address for Result Reporting 90/cho-6-95-15M Report Tier Levels - please select Fier 1 - (Results/Default if not specified) Relinquished by: (Signature) Relinquisheday: (Signature) Fier II - (Results + QC) Project Manager Client Sample ID

Temperature

Received by: (Signature)

Time:

Relinquished by: (Signature)

### Columbia Analytical Services, Inc. Sample Acceptance Check Form

Cliente	CHOMICH		Samp	le Acceptance	Check Fori		D0001016			
	CH2M Hill	erloo, NY / 369548.B2	EI		-	Work order:	P0801215			
-	(s) received on		.I'I		Date opened	• 4/28/2008	by:	MZAN	AOD A	<del></del>
-		all samples received by CA	S. The use of this	– form for custody	-		letted .			n of
		y. Thermal preservation an				_				11 01
		_					•	<u>Yes</u>	<u>No</u>	<u>N/A</u>
1	Were sample	e containers properly	marked with c	lient sample II	<b>)</b> ?			$\times$		
2	Container(s)	supplied by CAS?						X		
3	Did sample o	containers arrive in go	ood condition?					$\times$		
4	Were chain-	of-custody papers use	d and filled ou	t?				X		
5	Did sample o	container labels and/o	r tags agree w	ith custody pa	pers?			$\times$		
6	Was sample	volume received adeq	uate for analys	sis?				X		
7	Are samples	within specified holding	ng times?					X		
8	Was proper t	emperature (thermal	preservation)	of cooler at rec	eipt adhered	to?				X
		Cooler Temperature			Temperature		°C			
9.	Was a trip bl	ank received?		-					X	
	Trip blank	supplied by CAS: Seri	al#		-TB					
10	Were custody	y seals on outside of c	ooler/Box?				-		X	
	Location of	seal(s)?					Sealing Lid?			X
	Were signar	ture and date included	?							X
	Were seals	intact?							X	
	Were custody	seals on outside of sa	mple containe	r?						X
	Location of		•				Sealing Lid?			X
		ture and date included	?			**************************************				X
	Were seals			· ·						X
11		s have appropriate pre	eservation, acc	cording to met	hod/SOP or C	Client specified in	formation?			$\boxtimes$
		ent indication that the								$\boxtimes$
		vials checked for prese								$\boxtimes$
		nt/method/SOP requir			amplo nU an	d ifnooggamiali	ton it?			$\boxtimes$
12	Tubes:	Are the tubes cap			ampie pri am	u <u>ii necessary</u> an	er it?			
12	Tubes.	•	•							X
10		Do they contain r								$\overline{\times}$
13	Badges:	Are the badges p								$\boxtimes$
		Are dual bed bad	ges separated a	and individuall	y capped and	l intact?	7 1000000000000000000000000000000000000			X
Lab S	Sample ID	Container	Required	Received	Adjusted	VOA Headspace	Receip	t / Pres	ervation	
		Description	pH *	pН	р <b>Н</b>	(Presence/Absence)	C	ommer	its	
P0801215		6.0 L Ambient Can								
P0801215		6.0 L Ambient Can								
P0801215		6.0 L Ambient Can							····	
P0801215 P0801215		6.0 L Ambient Can		<del></del>	*****					
P0801215		6.0 L Ambient Can 6.0 L Ambient Can								
P0801215		6.0 L Ambient Can								
Explain a	ny discrepancies	s: (include lab sample ID	numbers):							الحصيح
***		•	, , , , , , , , , , , , , , , , , , ,			***				

Sample -006 canister serial number is listed as AC00684 on the COC, we received canister AC00648.

<sup>\*</sup>Required pH: Phenois/COD/NH3/TOC/TOX/NO3+NO2/TKN/T.PHOS, H2SO4 (pH<2); Metals, HNO3 (pH<2); CN (NaOH or NaOH/Asc Acid) (pH>12); Phenois/COD/NH3/TOC/TOX/NO3+NO2/TKN/T.PHOS, H2SO4 (pH<2); Metals, HNO3 (pH<2); CN (NaOH or NaOH/Asc Acid) (pH>12); Phenois/COD/NH3/TOC/TOX/NO3+NO2/TKN/T.PHOS, H2SO4 (pH<2); Metals, HNO3 (pH<2); CN (NaOH or NaOH/Asc Acid) (pH>12); Phenois/COD/NH3/TOC/TOX/NO3+NO2/TKN/T.PHOS, H2SO4 (pH<2); Metals, HNO3 (pH<2); CN (NaOH or NaOH/Asc Acid) (pH>12); Phenois/COD/NH3/TOC/TOX/NO3+NO2/TKN/T.PHOS, H2SO4 (pH<2); Metals, HNO3 (pH<2); CN (NaOH or NaOH/Asc Acid) (pH>12); Phenois/COD/NH3/TOC/TOX/NO3+NO2/TKN/T.PHOS, H2SO4 (pH<2); Metals, HNO3 (pH<2); CN (NaOH or NaOH/Asc Acid) (pH>12); Phenois/COD/NH3/TOC/TOX/NO3+NO2/TKN/T.PHOS, H2SO4 (pH<2); Metals, HNO3 (pH<2); CN (NaOH or NaOH/Asc Acid) (pH>12); Phenois/COD/NH3/TOC/TOX/NO3+NO2/TKN/T.PHOS, H2SO4 (pH<2); Metals, HNO3 (pH<2); CN (NaOH or NaOH/Asc Acid) (pH<2); Phenois/COD/NH3/TOC/TOX/NO3+NO2/TKN/T.PHOS, H2SO4 (pH<2); Phenois/COD/NH3/TOC/TOX/NO3+NO2/TKN/T.PHOS (pH<2); Phenois/COD/NH3/TCD/TOX/NO3+NO2/TKN/T.PHOS (pH<2); Phenois/COD/NH3/TCD/TOX/NO3+NO2/TKN/T.PHOS (phenois/COD/NH3/TCD/TOX/NO3+NO2/TKN/T.PHOS (phenois/COD/NH3/TCD/TOX/ND3/TCD/TOX/ND3/TCD/TOX/ND3/TCD/TOX/ND3/TCD/TOX/ND3/TCD/TOX/ND3/TCD/TOX/ND3/TCD/TOX/ND3/TCD/TOX/ND3/TCD/TOX/ND3/TCD/TOX/ND3/TCD/TOX/ND3/TCD/TOX/ND3/

### Columbia Analytical Services, Inc. Sample Acceptance Check Form

Client: CH2M Hill	Work order:	P0801215	
Project: DOW - Waterloo, NY / 369548.B2.FI	,	***************************************	
Sample(s) received on: 4/28/2008	Date opened: 4/28/2008	hv.	MZAMORA

Lab Sample ID	Container	Required	Received	Adjusted	VOA Headspace	Receipt / Preservation
	Description	pH *	pН	pН	(Presence/Absence)	Comments
P0801215-008.01	6.0 L Ambient Can					
P0801215-009.01	6.0 L Ambient Can					
P0801215-010.01	6.0 L Ambient Can					
P0801215-011.01	6.0 L Ambient Can					the state of the s
P0801215-012.01	6.0 L Ambient Can					
P0801215-013.01	6.0 L Ambient Can					
P0801215-014.01	6.0 L Ambient Can					
P0801215-015.01	6.0 L Ambient Can					
P0801215-016.01	6.0 L Ambient Can					
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Explain any discrepancies: (include lab sample ID numbers	):

### **RESULTS OF ANALYSIS**

Page 1 of 3

**Client:** 

**CH2M Hill** 

Client Sample ID: WAT-SG-B2-042308

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-001

Test Code:

EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/23/08

Date Received: 4/28/08

Instrument ID: Analyst:

Wida Ang

Date Analyzed: 5/2/08

Sampling Media:

6.0 L Summa Canister

Volume(s) Analyzed:

1.00 Liter(s)

Test Notes:

Container ID:

AC00812

Initial Pressure (psig):

-2.7

Final Pressure (psig):

3.5

Canister Dilution Factor: 1.52

Compound	Result	MRL	MDL	Result	MRL	MDL	Data
	$\mu g/m^3$	$\mu g/m^3$	$\mu g/m^3$	ppbV	ppbV	ppbV	Qualifier
Chloromethane	0.29	0.15	0.076	0.14	0.074	0.037	,
Vinyl Chloride	0.15	0.15	0.076	0.059	0.059	0.030	$\mathbf{U}$
Bromomethane	0.15	0.15	0.076	0.039	0.039	0.020	$\mathbf{U}$
Chloroethane	0.15	0.15	0.076	0.058	0.058	0.029	$\mathbf{U}$
Acetone	17	7.6	0.11	7.2	3.2	0.047	В .
Acrylonitrile	1.3	0.76	0.11	0.59	0.35	0.049	
1,1-Dichloroethene	0.15	0.15	0.076	0.038	0.038	0.019	$\mathbf{U}$
Methylene Chloride	0.39	0.76	0.076	0.11	0.22	0.022	J
Carbon Disulfide	0.71	0.76	0.18	0.23	0.24	0.059	J
trans-1,2-Dichloroethene	0.15	0.15	0.076	0.038	0.038	0.019	U
1,1-Dichloroethane	0.15	0.15	0.076	0.038	0.038	0.019	U
2-Butanone (MEK)	3.1	0.76	0.076	1.0	0.26	0.026	
cis-1,2-Dichloroethene	0.15	0.15	0.076	0.038	0.038	0.019	$\mathbf{U}$
Chloroform	1.1	0.15	0.090	0.23	0.031	0.018	
1,2-Dichloroethane	0.15	0.15	0.076	0.038	0.038	0.019	U
1,1,1-Trichloroethane	0.15	0.15	0.076	0.028	0.028	0.014	U
Benzene	0.57	0.15	0.076	0.18	0.048	0.024	
Carbon Tetrachloride	0.27	0.15	0.076	0.043	0.024	0.012	
	Chloromethane Vinyl Chloride Bromomethane Chloroethane Acetone Acrylonitrile 1,1-Dichloroethene Methylene Chloride Carbon Disulfide trans-1,2-Dichloroethene 1,1-Dichloroethane 2-Butanone (MEK) cis-1,2-Dichloroethene Chloroform 1,2-Dichloroethane 1,1,1-Trichloroethane 1,1,1-Trichloroethane Benzene	Chloromethane         μg/m³           Vinyl Chloride         0.15           Bromomethane         0.15           Chloroethane         0.15           Acetone         17           Acrylonitrile         1.3           1,1-Dichloroethene         0.15           Methylene Chloride         0.39           Carbon Disulfide         0.71           trans-1,2-Dichloroethene         0.15           1,1-Dichloroethane         0.15           2-Butanone (MEK)         3.1           cis-1,2-Dichloroethene         0.15           Chloroform         1.1           1,2-Dichloroethane         0.15           1,1,1-Trichloroethane         0.15           Benzene         0.57	Chloromethane         μg/m³         μg/m³           Vinyl Chloride         0.15         0.15           Bromomethane         0.15         0.15           Chloroethane         0.15         0.15           Chloroethane         0.15         0.15           Acetone         17         7.6           Acrylonitrile         1.3         0.76           1,1-Dichloroethene         0.15         0.15           Methylene Chloride         0.39         0.76           Carbon Disulfide         0.71         0.76           trans-1,2-Dichloroethene         0.15         0.15           1,1-Dichloroethane         0.15         0.15           2-Butanone (MEK)         3.1         0.76           cis-1,2-Dichloroethene         0.15         0.15           Chloroform         1.1         0.15           1,2-Dichloroethane         0.15         0.15           1,1,1-Trichloroethane         0.15         0.15           Benzene         0.57         0.15	Chloromethane         μg/m³         μg/m³         μg/m³           Vinyl Chloride         0.15         0.15         0.076           Bromomethane         0.15         0.15         0.076           Chloroethane         0.15         0.15         0.076           Chloroethane         0.15         0.15         0.076           Acetone         17         7.6         0.11           Acrylonitrile         1.3         0.76         0.11           1,1-Dichloroethene         0.15         0.15         0.076           Methylene Chloride         0.39         0.76         0.076           Carbon Disulfide         0.71         0.76         0.18           trans-1,2-Dichloroethene         0.15         0.15         0.076           1,1-Dichloroethane         0.15         0.15         0.076           2-Butanone (MEK)         3.1         0.76         0.076           Chloroform         1.1         0.15         0.076           Chloroform         1.1         0.15         0.076           1,1,1-Trichloroethane         0.15         0.15         0.076           Benzene         0.57         0.15         0.076	μg/m³         μg/m³         μg/m³         ppbV           Chloromethane         0.29         0.15         0.076         0.14           Vinyl Chloride         0.15         0.15         0.076         0.059           Bromomethane         0.15         0.15         0.076         0.039           Chloroethane         0.15         0.15         0.076         0.039           Chloroethane         17         7.6         0.11         0.59           1,1-Dichloroethene         0.15         0.15         0.076         0.038           Methylene Chloride         0.39         0.76         0.076         0.011           Carbon Disulfide         0.71         0.76         0.18         0.23           trans-1,2-Dichloroethene         0.15         0.15         0.076         0.038           1,1-Dichloroethane         0.15         0.15         0.076         0.038           2-Butanone (MEK)         3.1         0.76         0.076         0.038           Chloroform         1.1         0.15         0.076         0.038           Chloroform         1.1         0.15         0.076         0.038           1,1,1-Trichloroethane         0.15         0.15	μg/m³         μg/m³         μg/m³         ppbV         ppbV           Chloromethane         0.29         0.15         0.076         0.14         0.074           Vinyl Chloride         0.15         0.15         0.076         0.059         0.059           Bromomethane         0.15         0.15         0.076         0.039         0.039           Chloroethane         0.15         0.15         0.076         0.058         0.058           Acetone         17         7.6         0.11         7.2         3.2           Acrylonitrile         1.3         0.76         0.11         0.59         0.35           1,1-Dichloroethene         0.15         0.15         0.076         0.038         0.038           Methylene Chloride         0.39         0.76         0.076         0.011         0.22           Carbon Disulfide         0.71         0.76         0.18         0.23         0.24           trans-1,2-Dichloroethene         0.15         0.15         0.076         0.038         0.038           1,1-Dichloroethane         0.15         0.15         0.076         0.038         0.038           2-Butanone (MEK)         3.1         0.76         0.076	μg/m³         μg/m³         μg/m³         ppbV         ppbV         ppbV           Chloromethane         0.29         0.15         0.076         0.14         0.074         0.037           Vinyl Chloride         0.15         0.15         0.076         0.059         0.059         0.030           Bromomethane         0.15         0.15         0.076         0.039         0.039         0.020           Chloroethane         0.15         0.15         0.076         0.038         0.058         0.029           Acetone         17         7.6         0.11         7.2         3.2         0.047           Acrylonitrile         1.3         0.76         0.11         0.59         0.35         0.049           1,1-Dichloroethene         0.15         0.15         0.076         0.11         0.22         0.024           Methylene Chloride         0.39         0.76         0.076         0.11         0.22         0.022           Carbon Disulfide         0.71         0.76         0.18         0.23         0.24         0.059           trans-1,2-Dichloroethene         0.15         0.15         0.076         0.038         0.038         0.019           2-Butanone

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

B = Analyte was found in the method blank.

### **RESULTS OF ANALYSIS**

Page 2 of 3

Client:

CH2M Hill

Client Sample ID: WAT-SG-B2-042308

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-001

Test Code:

EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/23/08 Date Received: 4/28/08

Instrument ID: Analyst:

Date Analyzed: 5/2/08

6.0 L Summa Canister Sampling Media:

Wida Ang

Volume(s) Analyzed:

1.00 Liter(s)

Test Notes:

Container ID:

AC00812

Initial Pressure (psig):

-2.7

Final Pressure (psig):

3.5

Canister Dilution Factor: 1.52

CAS#	Compound	Result	MRL	MDL	Result	MRL	MDL	Data
		μg/m³	μg/m³	μg/m³	ppbV	ppbV	ppbV	Qualifier
78-87-5	1,2-Dichloropropane	0.15	0.15	0.076	0.033	0.033	0.016	$\mathbf{U}$
75-27-4	Bromodichloromethane	0.15	0.15	0.076	0.023	0.023	0.011	$\mathbf{U}$
79-01-6	Trichloroethene	0.15	0.15	0.076	0.028	0.028	0.014	$\mathbf{U}$
10061-01-5	cis-1,3-Dichloropropene	0.76	0.76	0.079	0.17	0.17	0.017	$\mathbf{U}$
108-10-1	4-Methyl-2-pentanone	120	0.76	0.085	29	0.19	0.021	
10061-02-6	trans-1,3-Dichloropropene	0.76	0.76	0.096	0.17	0.17	0.021	U
79-00-5	1,1,2-Trichloroethane	0.15	0.15	0.076	0.028	0.028	0.014	$\mathbf{U}$
108-88-3	Toluene	1.2	0.76	0.076	0.32	0.20	0.020	
124-48-1	Dibromochloromethane	0.15	0.15	0.10	0.018	0.018	0.012	$\mathbf{U}$
127-18-4	Tetrachloroethene	0.15	0.15	0.076	0.022	0.022	0.011	$\mathbf{U}$
108-90-7	Chlorobenzene	0.15	0.15	0.078	0.033	0.033	0.017	U
100-41-4	Ethylbenzene	0.24	0.76	0.094	0.055	0.18	0.022	J
179601-23-1	m,p-Xylenes	0.91	0.76	0.20	0.21	0.18	0.046	
75-25-2	Bromoform	0.76	0.76	0.12	0.074	0.074	0.011	U
100-42-5	Styrene	0.76	0.76	0.12	0.18	0.18	0.027	U
95-47-6	o-Xylene	0.39	0.76	0.096	0.089	0.18	0.022	J
79-34-5	1,1,2,2-Tetrachloroethane	0.15	0.15	0.097	0.022	0.022	0.014	<u>U</u>

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

### RESULTS OF ANALYSIS

Page 3 of 3

Client:

CH2M Hill

Client Sample ID: WAT-SG-B2-042308

CAS Project ID: P0801215

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Sample ID: P0801215-001

**Tentatively Identified Compounds** 

Test Code:

EPA TO-15

Date Collected: 4/23/08

Instrument ID:

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Received: 4/28/08

Analyst:

Wida Ang

Date Analyzed: 5/2/08

Sampling Media:

6.0 L Summa Canister

Volume(s) Analyzed:

1.00 Liter(s)

Test Notes:

T

Container ID:

AC00812

Initial Pressure (psig):

-2.7

Final Pressure (psig):

3.5

Canister Dilution Factor: 1.52

GC/MS	Compound Identification	Concentration	Data
Retention Time		$\mu \mathrm{g}/\mathrm{m}^3$	Qualifier
18.87	Benzaldehyde	20	
	Epichlorohydrin	NF	

T = Analyte is a tentatively identified compound, result is estimated.

NF = Compound was searched for, but not found.

### **RESULTS OF ANALYSIS**

Page 1 of 3

Client: **CH2M Hill** 

Client Sample ID: WAT-SG-1-042308 CAS Project ID: P0801215

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI CAS Sample ID: P0801215-002

Test Code: EPA TO-15

Date Collected: 4/23/08 Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16 Instrument ID: Date Received: 4/28/08 Analyst: Wida Ang Date Analyzed: 5/2/08

6.0 L Summa Canister Volume(s) Analyzed: Sampling Media: 0.30 Liter(s)

Test Notes:

Container ID: AC01311

> Initial Pressure (psig): 0.0 Final Pressure (psig): 3.5

> > Canister Dilution Factor: 1.24

CAS#	Compound	Result μg/m³	MRL μg/m³	$MDL$ $\mu g/m^3$	Result ppbV	MRL ppbV	MDL ppbV	Data Qualifier
74-87-3	Chloromethane	1.3	0.41	0.21	0.61	0.20	0.10	
75-01-4	Vinyl Chloride	0.41	0.41	0.21	0.16	0.16	0.081	U
74-83-9	Bromomethane	0.41	0.41	0.21	0.11	0.11	0.053	$\mathbf{U}$
75-00-3	Chloroethane	0.41	0.41	0.21	0.16	0.16	0.078	$\mathbf{U}$
67-64-1	Acetone	17	21	0.30	7.0	8.7	0.13	J, B
107-13-1	Acrylonitrile	5.9	2.1	0.29	2.7	0.95	0.13	The section of the section is
75-35-4	1,1-Dichloroethene	0.41	0.41	0.21	0.10	0.10	0.052	$\mathbf{U}$
75-09-2	Methylene Chloride	0.46	2.1	0.21	0.13	0.60	0.060	J
75-15-0	Carbon Disulfide	1.2	2.1	0.50	0.40	0.66	0.16	J
156-60-5	trans-1,2-Dichloroethene	0.41	0.41	0.21	0.10	0.10	0.052	$\mathbf{U}$
75-34-3	1,1-Dichloroethane	0.41	0.41	0.21	0.10	0.10	0.051	U
78-93-3	2-Butanone (MEK)	3.9	2.1	0.21	1.3	0.70	0.070	
156-59-2	cis-1,2-Dichloroethene	0.41	0.41	0.21	0.10	0.10	0.052	U
67-66-3	Chloroform	2.0	0.41	0.24	0.40	0.085	0.050	
107-06-2	1,2-Dichloroethane	0.41	0.41	0.21	0.10	0.10	0.051	$\mathbf{U}$
71-55-6	1,1,1-Trichloroethane	0.41	0.41	0.21	0.076	0.076	0.038	U
71-43-2	Benzene	0.45	0.41	0.21	0.14	0.13	0.065	
56-23-5	Carbon Tetrachloride	0.29	0.41	0.21	0.045	0.066	0.033	J

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

B = Analyte was found in the method blank.

### **RESULTS OF ANALYSIS**

Page 2 of 3

**Client:** 

**CH2M Hill** 

Client Sample ID: WAT-SG-1-042308

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-002

Test Code:

EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/23/08

Instrument ID: Analyst:

Wida Ang

Date Received: 4/28/08

Date Analyzed: 5/2/08

Sampling Media:

6.0 L Summa Canister

Volume(s) Analyzed:

0.30 Liter(s)

Test Notes:

Container ID:

AC01311

Initial Pressure (psig):

0.0

Final Pressure (psig):

3.5

Canister Dilution Factor: 1.24

CAS#	Compound	Result μg/m³	MRL μg/m³	MDL μg/m³	Result ppbV	MRL ppbV	MDL ppbV	Data Qualifier
78-87-5	1,2-Dichloropropane	0.41	0.41	0.21	0.089	0.089	0.045	U
75-27-4	Bromodichloromethane	0.41	0.41	0.21	0.062	0.062	0.031	$\mathbf{U}$
79-01-6	Trichloroethene	1.1	0.41	0.21	0.21	0.077	0.038	
10061-01-5	cis-1,3-Dichloropropene	2.1	2.1	0.21	0.46	0.46	0.047	$\mathbf{U}$
108-10-1	4-Methyl-2-pentanone	410	2.1	0.23	100	0.50	0.057	
10061-02-6	trans-1,3-Dichloropropene	2.1	2.1	0.26	0.46	0.46	0.057	U
79-00-5	1,1,2-Trichloroethane	0.41	0.41	0.21	0.076	0.076	0.038	U
108-88-3	Toluene	1.5	2.1	0.21	0.39	0.55	0.055	J
124-48-1	Dibromochloromethane	0.41	0.41	0.28	0.049	0.049	0.033	$\mathbf{U}$
127-18-4	Tetrachloroethene	1.1	0.41	0.21	0.16	0.061	0.030	
108-90-7	Chlorobenzene	0.41	0.41	0.21	0.090	0.090	0.046	U
100-41-4	Ethylbenzene	2.1	2.1	0.26	0.48	0.48	0.059	$\mathbf{U}$
179601-23-1	m,p-Xylenes	0.77	2.1	0.54	0.18	0.48	0.12	J
75-25-2	Bromoform	2.1	2.1	0.31	0.20	0.20	0.030	$\mathbf{U}$
100-42-5	Styrene	2.1	2.1	0.31	0.49	0.49	0.074	$\mathbf{U}$
95-47-6	o-Xylene	0.27	2.1	0.26	0.063	0.48	0.060	J
79-34-5	1,1,2,2-Tetrachloroethane	0.41	0.41	0.26	0.060	0.060	0.039	U

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

### RESULTS OF ANALYSIS

Page 3 of 3

**Client:** 

**CH2M Hill** 

Client Sample ID: WAT-SG-1-042308

6.0 L Summa Canister

CAS Project ID: P0801215

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Sample ID: P0801215-002

**Tentatively Identified Compounds** 

EPA TO-15

Date Collected: 4/23/08

Test Code: Instrument ID:

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Received: 4/28/08

Analyst:

Wida Ang

Date Analyzed: 5/2/08

Volume(s) Analyzed:

0.30 Liter(s)

Sampling Media: Test Notes:

T

Container ID:

AC01311

Initial Pressure (psig):

0.0

Final Pressure (psig):

3.5

Canister Dilution Factor: 1.24

GC/MS Retention Time	Compound Identification	Concentration µg/m³	Data Qualifier
4.46	Chlorodifluoromethane + Propene + Propane	20	
15.92	Hexamethylcyclotrisiloxane	10	*
16.27	C <sub>6</sub> HF <sub>13</sub> (Tridecafluorohexane)	10	
19.72	Hexamethylcyclotrisiloxane	40	*
22.06	Hexamethylcyclotrisiloxane	50	*
	Epichlorohydrin	NF	

T = Analyte is a tentatively identified compound, result is estimated.

NF = Compound was searched for, but not found.

<sup>\* =</sup> Possible artifact

### RESULTS OF ANALYSIS

Page 1 of 3

Client: CH2M Hill

Client Sample ID: WAT-SG-2-042308 CAS Project ID: P0801215

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI CAS Sample ID: P0801215-003

Test Code: EPA TO-15 Date Collected: 4/23/08
Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16 Date Received: 4/28/08

Analyst: Wida Ang Date Analyzed: 5/2/08

Sampling Media: 6.0 L Summa Canister Volume(s) Analyzed: 0.45 Liter(s)

Test Notes:
Container ID: AC00950

Initial Pressure (psig): -2.8 Final Pressure (psig): 3.5

Canister Dilution Factor: 1.53

CAS#	Compound	Result μg/m³	MRL μg/m³	MDL μg/m³	Result ppbV	MRL ppbV	MDL ppbV	Data Qualifier
74-87-3	Chloromethane	0.34	0.34	0.17	0.16	0.16	0.082	U
75-01-4	Vinyl Chloride	0.34	0.34	0.17	0.13	0.13	0.067	$\mathbf{U}$
74-83-9	Bromomethane	0.34	0.34	0.17	0.088	0.088	0.044	U
75-00-3	Chloroethane	0.34	0.34	0.17	0.13	0.13	0.064	U
67-64-1	Acetone	180	17	0.25	77	7.2	0.10	В
107-13-1	Acrylonitrile	1.2	1.7	0.24	0.53	0.78	0.11	J
75-35-4	1,1-Dichloroethene	4.2	0.34	0.17	1.1	0.086	0.043	
75-09-2	Methylene Chloride	1.1	1.7	0.17	0.31	0.49	0.049	J
75-15-0	Carbon Disulfide	10	1.7	0.41	3.3	0.55	0.13	
156-60-5	trans-1,2-Dichloroethene	0.34	0.34	0.17	0.086	0.086	0.043	$\mathbf{U}$
75-34-3	1,1-Dichloroethane	0.34	0.34	0.17	0.084	0.084	0.042	U
78-93-3	2-Butanone (MEK)	4.7	1.7	0.17	1.6	0.58	0.058	
156-59-2	cis-1,2-Dichloroethene	0.34	0.34	0.17	0.086	0.086	0.043	U
67-66-3	Chloroform	240	0.34	0.20	50	0.070	0.041	
107-06-2	1,2-Dichloroethane	2.0	0.34	0.17	0.49	0.084	0.042	
71-55-6	1,1,1-Trichloroethane	0.34	0.34	0.17	0.062	0.062	0.031	U
71-43-2	Benzene	0.93	0.34	0.17	0.29	0.11	0.053	
56-23-5	Carbon Tetrachloride	2.1	0.34	0.17	0.33	0.054	0.027	

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

B = Analyte was found in the method blank.

### RESULTS OF ANALYSIS

Page 2 of 3

Client: CH2M Hill

Client Sample ID: WAT-SG-2-042308 CAS Project ID: P0801215

CAS Sample ID: P0801215-003 Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

Test Code: EPA TO-15

Date Received: 4/28/08 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Analyst: Wida Ang Date Analyzed: 5/2/08 Volume(s) Analyzed:

6.0 L Summa Canister Sampling Media: Test Notes:

AC00950 Container ID:

Initial Pressure (psig): Final Pressure (psig): -2.83.5

Canister Dilution Factor: 1.53

0.45 Liter(s)

Date Collected: 4/23/08

CAS#	Compound	Result μg/m³	MRL μg/m³	MDL μg/m³	Result ppbV	MRL ppbV	MDL ppbV	Data Qualifier
78-87-5	1,2-Dichloropropane	$\frac{\mu g/m}{4.0}$	$\frac{\mu g/m}{0.34}$	$\frac{\mu g/m}{0.17}$	0.87	0.074	$\frac{pp0.7}{0.037}$	Quanner
75-27-4	Bromodichloromethane	0.34	0.34	0.17	0.051	0.051	0.025	$\mathbf{U}$
79-01-6	Trichloroethene	3.6	0.34	0.17	0.66	0.063	0.032	
10061-01-5	cis-1,3-Dichloropropene	1.7	1.7	0.18	0.37	0.37	0.039	$\mathbf{U}$
108-10-1	4-Methyl-2-pentanone	47	1.7	0.19	11	0.41	0.046	
10061-02-6	trans-1,3-Dichloropropene	1.7	1.7	0.21	0.37	0.37	0.047	U
79-00-5	1,1,2-Trichloroethane	0.34	0.34	0.17	0.062	0.062	0.031	$\mathbf{U}$
108-88-3	Toluene	5.7	1.7	0.17	1.5	0.45	0.045	
124-48-1	Dibromochloromethane	0.34	0.34	0.23	0.040	0.040	0.027	$\mathbf{U}$
127-18-4	Tetrachloroethene	32	0.34	0.17	4.7	0.050	0.025	
108-90-7	Chlorobenzene	0.34	0.34	0.17	0.074	0.074	0.038	U
100-41-4	Ethylbenzene	0.54	1.7	0.21	0.13	0.39	0.049	J
179601-23-1	m,p-Xylenes	3.3	1.7	0.44	0.75	0.39	0.10	
75-25-2	Bromoform	1.7	1.7	0.26	0.16	0.16	0.025	U
100-42-5	Styrene	0.30	1.7	0.26	0.070	0.40	0.061	J
95-47-6	o-Xylene	1.1	1.7	0.21	0.26	0.39	0.049	J
79-34-5	1,1,2,2-Tetrachloroethane	0.34	0.34	0.22	0.050	0.050	0.032	$\mathbf{U}$

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

### RESULTS OF ANALYSIS

Page 3 of 3

Client:

**CH2M Hill** 

Client Sample ID: WAT-SG-2-042308

6.0 L Summa Canister

CAS Project ID: P0801215

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Sample ID: P0801215-003

**Tentatively Identified Compounds** 

Test Code:

EPA TO-15

Date Collected: 4/23/08

Instrument ID:

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Received: 4/28/08

Date Analyzed: 5/2/08

Analyst: Sampling Media: Wida Ang

Volume(s) Analyzed:

0.45 Liter(s)

Test Notes:

 $\mathbf{T}$ 

Container ID:

AC00950

Initial Pressure (psig):

-2.8

Final Pressure (psig):

3.5

Canister Dilution Factor: 1.53

GC/MS Retention Time	Compound Identification	Concentration μg/m³	Data Qualifier
16.27	C <sub>6</sub> HF <sub>13</sub> (Tridecafluorohexane)	100	
21.27	n-Nonanal + C <sub>10</sub> H <sub>14</sub> Aromatic	20	
22.06	Hexamethylcyclotrisiloxane	30	*
22.66	n-Dodecane	10	
24.81	C <sub>14</sub> H <sub>28</sub> Compound	40	
	Epichlorohydrin	NF	

T = Analyte is a tentatively identified compound, result is estimated.

NF = Compound was searched for, but not found.

<sup>\* =</sup> Possible artifact

### RESULTS OF ANALYSIS

Page 1 of 3

Client:

**CH2M Hill** 

Client Sample ID: WAT-SG-3-042308

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-004

Test Code:

EPA TO-15

Instrument ID:

Sampling Media:

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/23/08 Date Received: 4/28/08

Analyst:

Wida Ang

6.0 L Summa Canister

Date Analyzed: 5/2/08 Volume(s) Analyzed:

0.10 Liter(s)

Test Notes:

Container ID:

AC01189

Initial Pressure (psig):

0.1

Final Pressure (psig):

3.5

Canister Dilution Factor: 1.23

CAC#	Commound	Dogult	MDI	MDI	Dogulf	MDI	MDI	Data
CAS#	Compound	Result	MRL	MDL	Result	MRL	MDL	Data
		μg/m³	<u>μg/m³</u>	$\mu g/m^3$	ppbV	ppbV	ppbV	
74-87-3	Chloromethane	1.0	1.2	0.62	0.51	0.60	0.30	J
75-01-4	Vinyl Chloride	1.2	1.2	0.62	0.48	0.48	0.24	$\mathbf{U}$
74-83-9	Bromomethane	1.2	1.2	0.62	0.32	0.32	0.16	$\mathbf{U}$
75-00-3	Chloroethane	1.2	1.2	0.62	0.47	0.47	0.23	$\mathbf{U}$
67-64-1	Acetone	130	62	0.90	53	26	0.38	В
107-13-1	Acrylonitrile	27	6.2	0.86	12	2.8	0.40	
75-35-4	1,1-Dichloroethene	1.2	1.2	0.62	0.31	0.31	0.16	$\mathbf{U}$
75-09-2	Methylene Chloride	1.3	6.2	0.62	0.39	1.8	0.18	J
75-15-0	Carbon Disulfide	34	6.2	1.5	11	2.0	0.47	
156-60-5	trans-1,2-Dichloroethene	1.2	1.2	0.62	0.31	0.31	0.16	$\mathbf{U}$
75-34-3	1,1-Dichloroethane	1.2	1.2	0.62	0.30	0.30	0.15	U
78-93-3	2-Butanone (MEK)	2.8	6.2	0.62	0.96	2.1	0.21	J
156-59-2	cis-1,2-Dichloroethene	1.2	1.2	0.62	0.31	0.31	0.16	$\mathbf{U}^{-1}$
67-66-3	Chloroform	25	1.2	0.73	5.1	0.25	0.15	
107-06-2	1,2-Dichloroethane	1.2	1.2	0.62	0.30	0.30	0.15	$\mathbf{U}$
71-55-6	1,1,1-Trichloroethane	1.2	1.2	0.62	0.23	0.23	0.11	U
71-43-2	Benzene	90	1.2	0.62	28	0.39	0.19	
56-23-5	Carbon Tetrachloride	1.2	1.2	0.62	0.20	0.20	0.098	U

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

B = Analyte was found in the method blank.

### RESULTS OF ANALYSIS

Page 2 of 3

**Client: CH2M Hill** 

CAS Project ID: P0801215 Client Sample ID: WAT-SG-3-042308

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI CAS Sample ID: P0801215-004

Test Code: EPA TO-15

Date Collected: 4/23/08 Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16 Date Received: 4/28/08 Instrument ID:

Date Analyzed: 5/2/08 Analyst: Wida Ang

6.0 L Summa Canister Volume(s) Analyzed: 0.10 Liter(s) Sampling Media:

Test Notes: Container ID: AC01189

Final Pressure (psig): Initial Pressure (psig): 0.1 3.5

Canister Dilution Factor: 1.23

CAS#	Compound	Result μg/m³	MRL μg/m³	$\begin{array}{c} MDL \\ \mu g/m^3 \end{array}$	Result ppbV	MRL ppbV	MDL ppbV	Data Qualifier
78-87-5	1,2-Dichloropropane	1.2	1.2	0.62	0.27	0.27	0.13	U
75-27-4	Bromodichloromethane	1.2	1.2	0.62	0.18	0.18	0.092	$\mathbf{U}$
79-01-6	Trichloroethene	2.9	1.2	0.62	0.54	0.23	0.11	
10061-01-5	cis-1,3-Dichloropropene	6.2	6.2	0.64	1.4	1.4	0.14	$\mathbf{U}$
108-10-1	4-Methyl-2-pentanone	820	6.2	0.69	200	1.5	0.17	
10061-02-6	trans-1,3-Dichloropropene	6.2	6.2	0.77	1.4	1.4	0.17	U
79-00-5	1,1,2-Trichloroethane	1.2	1.2	0.62	0.23	0.23	0.11	$\mathbf{U}$
108-88-3	Toluene	16	6.2	0.62	4.1	1.6	0.16	
124-48-1	Dibromochloromethane	1.2	1.2	0.84	0.14	0.14	0.098	U
127-18-4	Tetrachloroethene	8.9	1.2	0.62	1.3	0.18	0.091	
108-90-7	Chlorobenzene	1.2	1.2	0.63	0.27	0.27	0.14	U
100-41-4	Ethylbenzene	0.91	6.2	0.76	0.21	1.4	0.18	${f J}$
179601-23-1	m,p-Xylenes	4.4	6.2	1.6	1.0	1.4	0.37	J
75-25-2	Bromoform	6.2	6.2	0.93	0.60	0.60	0.090	$\mathbf{U}$
100-42-5	Styrene	6.2	6.2	0.93	1.4	1.4	0.22	U
95-47-6	o-Xylene	1.5	6.2	0.77	0.34	1.4	0.18	J
79-34-5	1,1,2,2-Tetrachloroethane	1.2	1.2	0.79	0.18	0.18	0.11	U

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

### RESULTS OF ANALYSIS

Page 3 of 3

**Tentatively Identified Compounds** 

Client:

**CH2M Hill** 

Client Sample ID: WAT-SG-3-042308

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-004

Test Code:

EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/23/08

Date Received: 4/28/08

Instrument ID: Analyst:

Wida Ang

Date Analyzed: 5/2/08 Volume(s) Analyzed:

0.10 Liter(s)

Sampling Media: Test Notes:

6.0 L Summa Canister T

Container ID:

AC01189

Initial Pressure (psig):

0.1

Final Pressure (psig):

3.5

Canister Dilution Factor: 1.23

GC/MS	Compound Identification	Concentration	Data
Retention Time	•	$\mu \mathrm{g}/\mathrm{m}^3$	Qualifier
16.27	C <sub>6</sub> HF <sub>13</sub> (Tridecafluorohexane)	70	
	Epichlorohydrin	NF	

T = Analyte is a tentatively identified compound, result is estimated.

NF = Compound was searched for, but not found.

### RESULTS OF ANALYSIS

Page 1 of 3

Client: **CH2M Hill** 

CAS Project ID: P0801215 Client Sample ID: WAT-SG-3a-042308

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI CAS Sample ID: P0801215-005

Test Code: EPA TO-15

Date Collected: 4/23/08 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16 Date Received: 4/28/08

Analyst: Wida Ang Date Analyzed: 5/5/08

Sampling Media: Test Notes:

6.0 L Summa Canister Volume(s) Analyzed: 0.50 Liter(s)

Container ID: AC01387

> Initial Pressure (psig): Final Pressure (psig): -2.23.5

> > Canister Dilution Factor: 1.46

CAS#	Compound	Result μg/m³	MRL μg/m³	MDL μg/m³	Result ppbV	MRL ppbV	MDL ppbV	Data Qualifier
74-87-3	Chloromethane	0.39	0.29	0.15	0.19	0.14	0.071	.,
75-01-4	Vinyl Chloride	0.29	0.29	0.15	0.11	0.11	0.057	$\mathbf{U}$
74-83-9	Bromomethane	0.29	0.29	0.15	0.075	0.075	0.038	$\mathbf{U}$
75-00-3	Chloroethane	0.72	0.29	0.15	0.27	0.11	0.055	
67-64-1	Acetone	<b>56</b>	15	0.21	23	6.1	0.090	В
107-13-1	Acrylonitrile	6.9	1.5	0.20	3.2	0.67	0.094	
75-35-4	1,1-Dichloroethene	0.29	0.29	0.15	0.074	0.074	0.037	$\mathbf{U}$
75-09-2	Methylene Chloride	0.43	1.5	0.15	0.12	0.42	0.042	J
75-15-0	Carbon Disulfide	9.5	1.5	0.35	3.0	0.47	0.11	
156-60-5	trans-1,2-Dichloroethene	0.29	0.29	0.15	0.074	0.074	0.037	$\mathbf{U}$
75-34-3	1,1-Dichloroethane	0.29	0.29	0.15	0.072	0.072	0.036	U
78-93-3	2-Butanone (MEK)	4.3	1.5	0.15	1.4	0.50	0.050	
156-59-2	cis-1,2-Dichloroethene	1.4	0.29	0.15	0.35	0.074	0.037	
67-66-3	Chloroform	3.6	0.29	0.17	0.74	0.060	0.035	
107-06-2	1,2-Dichloroethane	0.29	0.29	0.15	0.072	0.072	0.036	$\mathbf{U}$
71-55-6	1,1,1-Trichloroethane	0.29	0.29	0.15	0.054	0.054	0.027	U
71-43-2	Benzene	13	0.29	0.15	4.0	0.091	0.046	
56-23-5	Carbon Tetrachloride	0.29	0.29	0.15	0.046	0.046	0.023	$\mathbf{U}$

U = Compound was analyzed for, but not detected above the laboratory detection limit.

Verified By: Re

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

B = Analyte was found in the method blank.

### **RESULTS OF ANALYSIS**

Page 2 of 3

**Client:** 

**CH2M Hill** 

Client Sample ID: WAT-SG-3a-042308

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-005

Test Code:

EPA TO-15

Instrument ID:

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Received: 4/28/08

Date Collected: 4/23/08

Analyst: Sampling Media: Wida Ang 6.0 L Summa Canister

Date Analyzed: 5/5/08 Volume(s) Analyzed:

0.50 Liter(s)

Test Notes:

Container ID:

AC01387

Initial Pressure (psig):

-2.2

Final Pressure (psig):

3.5

Canister Dilution Factor: 1.46

CAS#	Compound	Result μg/m³	MRL μg/m³	MDL μg/m³	Result ppbV	MRL ppbV	MDL ppbV	Data Qualifier
78-87-5	1,2-Dichloropropane	0.29	0.29	0.15	0.063	0.063	0.032	U
75-27-4	Bromodichloromethane	0.29	0.29	0.15	0.044	0.044	0.022	$\mathbf{U}$
79-01-6	Trichloroethene	10	0.29	0.15	1.9	0.054	0.027	
10061-01-5	cis-1,3-Dichloropropene	1.5	1.5	0.15	0.32	0.32	0.033	$\mathbf{U}$
108-10-1	4-Methyl-2-pentanone	240	1.5	0.16	58	0.36	0.040	
10061-02-6	trans-1,3-Dichloropropene	1.5	1.5	0.18	0.32	0.32	0.041	U
79-00-5	1,1,2-Trichloroethane	0.29	0.29	0.15	0.054	0.054	0.027	$\mathbf{U}$
108-88-3	Toluene	26	1.5	0.15	6.8	0.39	0.039	
124-48-1	Dibromochloromethane	0.29	0.29	0.20	0.034	0.034	0.023	$\mathbf{U}$
127-18-4	Tetrachloroethene	5.4	0.29	0.15	0.79	0.043	0.022	
108-90-7	Chlorobenzene	0.29	0.29	0.15	0.063	0.063	0.032	U
100-41-4	Ethylbenzene	0.43	1.5	0.18	0.10	0.34	0.042	J
179601-23-1	m,p-Xylenes	2.3	1.5	0.38	0.53	0.34	0.087	
75-25-2	Bromoform	1.5	1.5	0.22	0.14	0.14	0.021	$\mathbf{U}$
100-42-5	Styrene	0.40	1.5	0.22	0.095	0.34	0.052	J
95-47-6	o-Xylene	0.83	1.5	0.18	0.19	0.34	0.042	J
79-34-5	1,1,2,2-Tetrachloroethane	0.29	0.29	0.19	0.043	0.043	0.027	<u>U</u>

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

### **RESULTS OF ANALYSIS**

Page 3 of 3

**Client:** 

**CH2M Hill** 

Client Sample ID: WAT-SG-3a-042308

6.0 L Summa Canister

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-005

**Tentatively Identified Compounds** 

Test Code: Instrument ID: EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/23/08

Date Received: 4/28/08

Analyst:

Wida Ang

Date Analyzed: 5/5/08

Volume(s) Analyzed:

0.50 Liter(s)

Sampling Media: Test Notes:

Container ID:

AC01387

Initial Pressure (psig):

-2.2

Final Pressure (psig):

3.5

Canister Dilution Factor: 1.46

GC/MS	Compound Identification	Concentration	Data	
Retention Time		$\mu \mathrm{g}/\mathrm{m}^3$	Qualifier	
4.46	Propane + Propene	30		
16.27	C <sub>6</sub> HF <sub>13</sub> (Tridecafluorohexane)	200		
21.26	n-Nonanal	10		
22.06	Hexamethylcyclotrisiloxane	20	*	
24.81	$C_{14}H_{28}$ Compound	40		
	Epichlorohydrin	NF		

T = Analyte is a tentatively identified compound, result is estimated.

NF = Compound was searched for, but not found.

<sup>\* =</sup> Possible artifact

#### RESULTS OF ANALYSIS

Page 1 of 3

Client: CH2M Hill

AC00648

Client Sample ID: WAT-SG-4-042308 CAS Project ID: P0801215

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI CAS Sample ID: P0801215-006

Test Code: EPA TO-15 Date Collected: 4/23/08

Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16 Date Received: 4/28/08

Analyst: Wida Ang Date Analyzed: 5/2/08 Sampling Media: 6.0 L Summa Canister Volume(s) Analyzed: 0.30 Liter(s)

Test Notes:

Container ID:

Initial Pressure (psig): 0.0 Final Pressure (psig): 3.6

CAS#	Compound	Result μg/m³	MRL μg/m³	MDL μg/m³	Result ppbV	MRL ppbV	MDL ppbV	Data Qualifier
74-87-3	Chloromethane	0.62	0.41	0.21	0.30	0.20	0.10	
75-01-4	Vinyl Chloride	0.41	0.41	0.21	0.16	0.16	0.081	$\mathbf{U}$
74-83-9	Bromomethane	0.41	0.41	0.21	0.11	0.11	0.053	$\mathbf{U}$
75-00-3	Chloroethane	0.41	0.41	0.21	0.16	0.16	0.078	$\mathbf{U}$
67-64-1	Acetone	. 17	21	0.30	7.0	8.7	0.13	J, B
107-13-1	Acrylonitrile	0.86	2.1	0.29	0.39	0.95	0.13	J
75-35-4	1,1-Dichloroethene	0.41	0.41	0.21	0.10	0.10	0.052	$\mathbf{U}$
75-09-2	Methylene Chloride	0.39	2.1	0.21	0.11	0.60	0.060	J
75-15-0	Carbon Disulfide	2.2	2.1	0.50	0.71	0.66	0.16	
156-60-5	trans-1,2-Dichloroethene	0.41	0.41	0.21	0.10	0.10	0.052	U
75-34-3	1,1-Dichloroethane	2.1	0.41	0.21	0.53	0.10	0.051	
78-93-3	2-Butanone (MEK)	3.7	2.1	0.21	1.2	0.70	0.070	
156-59-2	cis-1,2-Dichloroethene	0.41	0.41	0.21	0.10	0.10	0.052	$\mathbf{U}$
67-66-3	Chloroform	440	0.41	0.24	90	0.085	0.050	
107-06-2	1,2-Dichloroethane	0.41	0.41	0.21	0.10	0.10	0.051	$\mathbf{U}$
71-55-6	1,1,1-Trichloroethane	12	0.41	0.21	2.2	0.076	0.038	halderden floren er sembler er e
71-43-2	Benzene	0.81	0.41	0.21	0.25	0.13	0.065	
56-23-5	Carbon Tetrachloride	0.41	0.41	0.21	0.066	0.066	0.033	$\mathbf{U}$

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

B = Analyte was found in the method blank.

# RESULTS OF ANALYSIS

Page 2 of 3

Client: CH2M Hill

Client Sample ID: WAT-SG-4-042308 CAS Project ID: P0801215

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI CAS Sample ID: P0801215-006

Test Code: EPA TO-15

Date Collected: 4/23/08 Date Received: 4/28/08 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Analyst: Wida Ang Date Analyzed: 5/2/08

Volume(s) Analyzed: Sampling Media: 6.0 L Summa Canister 0.30 Liter(s)

Test Notes:

Container ID: AC00648

Initial Pressure (psig): 0.0 Final Pressure (psig): 3.6

Canister Dilution Factor: 1.24

CAS#	Compound	Result	MRL	MDL	Result	MRL	MDL	Data Qualifier
78-87-5	1.2 Diablevenyenene	$\frac{\mu g/m^3}{3.9}$	$\frac{\mu g/m^3}{0.41}$	$\frac{\mu g/m^3}{0.21}$	0.84	ppbV 0.089	ppbV 0.045	Quanner
	1,2-Dichloropropane							**
75-27-4	Bromodichloromethane	0.41	0.41	0.21	0.062	0.062	0.031	U
79-01-6	Trichloroethene	2.5	0.41	0.21	0.46	0.077	0.038	
10061-01-5	cis-1,3-Dichloropropene	2.1	2.1	0.21	0.46	0.46	0.047	U
108-10-1	4-Methyl-2-pentanone	100	2.1	0.23	25	0.50	0.057	
10061-02-6	trans-1,3-Dichloropropene	2.1	2.1	0.26	0.46	0.46	0.057	$\mathbf{U}$
79-00-5	1,1,2-Trichloroethane	0.41	0.41	0.21	0.076	0.076	0.038	$\mathbf{U}$
108-88-3	Toluene	5.8	2.1	0.21	1.5	0.55	0.055	
124-48-1	Dibromochloromethane	0.41	0.41	0.28	0.049	0.049	0.033	$\mathbf{U}$
127-18-4	Tetrachloroethene	17	0.41	0.21	2.6	0.061	0.030	
108-90-7	Chlorobenzene	0.41	0.41	0.21	0.090	0.090	0.046	U
100-41-4	Ethylbenzene	0.61	2.1	0.26	0.14	0.48	0.059	J
179601-23-1	m,p-Xylenes	4.3	2.1	0.54	0.99	0.48	0.12	
75-25-2	Bromoform	2.1	2.1	0.31	0.20	0.20	0.030	$\mathbf{U}$
100-42-5	Styrene	2.1	2.1	0.31	0.49	0.49	0.074	$\mathbf{U}$
95-47-6	o-Xylene	2.9	2.1	0.26	0.66	0.48	0.060	
79-34-5	1,1,2,2-Tetrachloroethane	0.41	0.41	0.26	0.060	0.060	0.039	$\mathbf{U}$

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

#### **RESULTS OF ANALYSIS**

Page 3 of 3

Client:

CH2M Hill

Client Sample ID: WAT-SG-4-042308

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-006

**Tentatively Identified Compounds** 

Test Code: Instrument ID: EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/23/08

Date Received: 4/28/08

Analyst:

Wida Ang

Date Analyzed: 5/2/08 Volume(s) Analyzed:

0.30 Liter(s)

Sampling Media: Test Notes:

6.0 L Summa Canister T

Container ID:

AC00648

Initial Pressure (psig):

0.0

Final Pressure (psig):

3.6

GC/MS	Compound Identification	Concentration	Data
Retention Time		$\mu \mathrm{g}/\mathrm{m}^3$	Qualifier
16.28	C <sub>6</sub> HF <sub>13</sub> (Tridecafluorohexane)	80	
20.73	C <sub>12</sub> H <sub>26</sub> Branched Alkane	40	
21.02	C <sub>11</sub> H <sub>24</sub> Branched Alkane	50	
21.05	C <sub>11</sub> H <sub>24</sub> Branched Alkane	60	
21.44	n-Undecane	70	
	Epichlorohydrin	NF	

T = Analyte is a tentatively identified compound, result is estimated.

NF = Compound was searched for, but not found.

# **RESULTS OF ANALYSIS**

Page 1 of 3

Client:

**CH2M Hill** 

Client Sample ID: WAT-SG-DUP-042308

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-007

Test Code: Instrument ID: EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/23/08

Date Received: 4/28/08

Analyst:

Wida Ang

Date Analyzed: 5/5/08

Sampling Media:

6.0 L Summa Canister

Volume(s) Analyzed:

0.10 Liter(s)

Test Notes:

Container ID:

AC01168

Initial Pressure (psig):

-0.1

Final Pressure (psig):

3.6

CAS#	Compound	Result µg/m³	MRL μg/m³	$MDL$ $\mu g/m^3$	Result ppbV	MRL ppbV	MDL ppbV	Data Qualifier
74-87-3	Chloromethane	1.3	1.3	0.63	0.61	0.61	0.30	U
75-01-4	Vinyl Chloride	1.3	1.3	0.63	0.49	0.49	0.24	U
74-83-9	Bromomethane	1.3	1.3	0.63	0.32	0.32	0.16	U
75-00-3	Chloroethane	1.3	1.3	0.63	0.47	0.47	0.24	Ü
67-64-1	Acetone	250	63	0.91	100	26	0.38	В
107-13-1	Acrylonitrile	6.3	6.3	0.88	2.9	2.9	0.40	U
75-35-4	1,1-Dichloroethene	4.0	1.3	0.63	1.0	0.32	0.16	
75-09-2	Methylene Chloride	1.6	6.3	0.63	0.45	1.8	0.18	J
75-15-0	Carbon Disulfide	8.6	6.3	1.5	2.8	2.0	0.48	
156-60-5	trans-1,2-Dichloroethene	1.3	1.3	0.63	0.32	0.32	0.16	$\mathbf{U}$
75-34-3	1,1-Dichloroethane	1.3	1.3	0.63	0.31	0.31	0.15	U
78-93-3	2-Butanone (MEK)	2.5	6.3	0.63	0.83	2.1	0.21	J
156-59-2	cis-1,2-Dichloroethene	1.3	1.3	0.63	0.32	0.32	0.16	U
67-66-3	Chloroform	260	1.3	0.74	53	0.26	0.15	
107-06-2	1,2-Dichloroethane	2.1	1.3	0.63	0.51	0.31	0.15	
71-55-6	1,1,1-Trichloroethane	1.3	1.3	0.63	0.23	0.23	0.11	U
71-43-2	Benzene	1.3	1.3	0.63	0.39	0.39	0.20	$\mathbf{U}$
56-23-5	Carbon Tetrachloride	1.6	1.3	0.63	0.26	0.20	0.099	

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

B = Analyte was found in the method blank.

### **RESULTS OF ANALYSIS** Page 2 of 3

**CH2M Hill** 

Client Sample ID: WAT-SG-DUP-042308 CAS Project ID: P0801215

CAS Sample ID: P0801215-007 Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

Test Code: EPA TO-15

Date Collected: 4/23/08 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16 Date Received: 4/28/08

Date Analyzed: 5/5/08 Analyst: Wida Ang

6.0 L Summa Canister Volume(s) Analyzed: 0.10 Liter(s) Sampling Media:

Test Notes:

Client:

AC01168 Container ID:

Initial Pressure (psig): -0.1Final Pressure (psig): 3.6

Canister Dilution Factor: 1.25

CAS#	Compound	Result μg/m³	MRL μg/m³	MDL μg/m³	Result ppbV	MRL ppbV	MDL ppbV	Data Qualifier
78-87-5	1,2-Dichloropropane	3.6	1.3	0.63	0.78	0.27	0.14	
75-27-4	Bromodichloromethane	1.3	1.3	0.63	0.19	0.19	0.093	$\mathbf{U}$
79-01-6	Trichloroethene	3.5	1.3	0.63	0.65	0.23	0.12	
10061-01-5	cis-1,3-Dichloropropene	6.3	6.3	0.65	1.4	1.4	0.14	$\mathbf{U}$
108-10-1	4-Methyl-2-pentanone	15	6.3	0.70	3.6	1.5	0.17	
10061-02-6	trans-1,3-Dichloropropene	6.3	6.3	0.79	1.4	1.4	0.17	U
79-00-5	1,1,2-Trichloroethane	1.3	1.3	0.63	0.23	0.23	0.11	$\mathbf{U}$
108-88-3	Toluene	2.1	6.3	0.63	0.56	1.7	0.17	J
124-48-1	Dibromochloromethane	1.3	1.3	0.85	0.15	0.15	0.10	$\mathbf{U}$
127-18-4	Tetrachloroethene	28	1.3	0.63	4.1	0.18	0.092	
108-90-7	Chlorobenzene	1.3	1.3	0.64	0.27	0.27	0.14	U
100-41-4	Ethylbenzene	6.3	6.3	0.78	1.4	1.4	0.18	$\mathbf{U}$
179601-23-1	m,p-Xylenes	6.3	6.3	1.6	1.4	1.4	0.37	U
75-25-2	Bromoform	6.3	6.3	0.95	0.60	0.60	0.092	$\mathbf{U}$
100-42-5	Styrene	6.3	6.3	0.95	1.5	1.5	0.22	$\mathbf{U}$
95-47-6	o-Xylene	6.3	6.3	0.79	1.4	1.4	0.18	U
79-34-5	1,1,2,2-Tetrachloroethane	1.3	1.3	0.80	0.18	0.18	0.12	U

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

Verified By:

#### RESULTS OF ANALYSIS

Page 3 of 3

**Tentatively Identified Compounds** 

**Client:** 

**CH2M Hill** 

Client Sample ID: WAT-SG-DUP-042308

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-007

Test Code:

EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/23/08

Date Received: 4/28/08

Instrument ID: Analyst:

Wida Ang

Date Analyzed: 5/5/08

Sampling Media:

6.0 L Summa Canister

Volume(s) Analyzed:

0.10 Liter(s)

Test Notes:

T

Container ID:

AC01168

Initial Pressure (psig):

-0.1

Final Pressure (psig):

3.6

Canister Dilution Factor: 1.25

GC/MS	Compound Identification	Concentration	Data
Retention Time		$\mu \mathrm{g}/\mathrm{m}^3$	Qualifier
16.26	C <sub>6</sub> HF <sub>13</sub> (Tridecafluorohexane)	50	
	Epichlorohydrin	NF	

T = Analyte is a tentatively identified compound, result is estimated.

NF = Compound was searched for, but not found.

# **RESULTS OF ANALYSIS**

Page 1 of 3

Client:

CH2M Hill

Client Sample ID: WAT-SG-FB-042308

6.0 L Summa Canister

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-008

Test Code:

EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/23/08 Date Received: 4/28/08

Instrument ID: Analyst:

Wida Ang

Date Analyzed: 5/2/08

Volume(s) Analyzed:

1.00 Liter(s)

Sampling Media: Test Notes:

Container ID:

AC01209

CAS#	Compound	Result	MRL	MDL	Result	MRL	MDL	Data
	-	$\mu g/m^3$	$\mu g/m^3$	$\mu g/m^3$	$\mathbf{p}\mathbf{p}\mathbf{b}\mathbf{V}$	ppbV	ppbV	Qualifier
74-87-3	Chloromethane	0.10	0.10	0.050	0.048	0.048	0.024	U
75-01-4	Vinyl Chloride	0.10	0.10	0.050	0.039	0.039	0.020	$\mathbf{U}$
74-83-9	Bromomethane	0.10	0.10	0.050	0.026	0.026	0.013	$\mathbf{U}$
75-00-3	Chloroethane	0.10	0.10	0.050	0.038	0.038	0.019	$\mathbf{U}_{\perp}$
67-64-1	Acetone	2.2	5.0	0.073	0.92	2.1	0.031	J, B
107-13-1	Acrylonitrile	0.50	0.50	0.070	0.23	0.23	0.032	U
75-35-4	1,1-Dichloroethene	0.10	0.10	0.050	0.025	0.025	0.013	$\mathbf{U}$
75-09-2	Methylene Chloride	0.50	0.50	0.050	0.14	0.14	0.014	$\mathbf{U}$
75-15-0	Carbon Disulfide	0.44	0.50	0.12	0.14	0.16	0.039	J
156-60-5	trans-1,2-Dichloroethene	0.10	0.10	0.050	0.025	0.025	0.013	U
75-34-3	1,1-Dichloroethane	0.10	0.10	0.050	0.025	0.025	0.012	U
78-93-3	2-Butanone (MEK)	0.36	0.50	0.050	0.12	0.17	0.017	J
156-59-2	cis-1,2-Dichloroethene	0.10	0.10	0.050	0.025	0.025	0.013	$\mathbf{U}$
67-66-3	Chloroform	0.10	0.10	0.059	0.020	0.020	0.012	$\mathbf{U}$
107-06-2	1,2-Dichloroethane	0.10	0.10	0.050	0.025	0.025	0.012	U
71-55-6	1,1,1-Trichloroethane	0.10	0.10	0.050	0.018	0.018	0.0092	U
71-43-2	Benzene	0.10	0.10	0.050	0.031	0.031	0.016	$\mathbf{U}$
56-23-5	Carbon Tetrachloride	0.10	0.10	0.050	0.016	0.016	0.0080	U

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

B = Analyte was found in the method blank.

### RESULTS OF ANALYSIS

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**Client:** 

**CH2M Hill** 

Client Sample ID: WAT-SG-FB-042308

6.0 L Summa Canister

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-008

Test Code:

EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/23/08

Date Received: 4/28/08

Instrument ID: Analyst:

Wida Ang

Date Analyzed: 5/2/08 Volume(s) Analyzed:

1.00 Liter(s)

Sampling Media: Test Notes:

Container ID:

AC01209

Canister Dilution Factor: 1.00

CAS#	Compound	Result μg/m³	MRL μg/m³	MDL μg/m³	Result ppbV	MRL ppbV	MDL ppbV	Data Oualifier
78-87-5	1,2-Dichloropropane	0.10	$\frac{\mu g/m}{0.10}$	$\frac{\mu g m}{0.050}$	0.022	0.022	0.011	U
75-27-4	Bromodichloromethane	0.10	0.10	0.050	0.015	0.015	0.0075	U
79-01-6	Trichloroethene	0.10	0.10	0.050	0.019	0.019	0.0093	$\mathbf{U}$
10061-01-5	cis-1,3-Dichloropropene	0.50	0.50	0.052	0.11	0.11	0.011	U
108-10-1	4-Methyl-2-pentanone	0.096	0.50	0.056	0.023	0.12	0.014	J
10061-02-6	trans-1,3-Dichloropropene	0.50	0.50	0.063	0.11	0.11	0.014	U
79-00-5	1,1,2-Trichloroethane	0.10	0.10	0.050	0.018	0.018	0.0092	$\mathbf{U}$
108-88-3	Toluene	0.24	0.50	0.050	0.065	0.13	0.013	J
124-48-1	Dibromochloromethane	0.10	0.10	0.068	0.012	0.012	0.0080	$\mathbf{U}$
127-18-4	Tetrachloroethene	0.10	0.10	0.050	0.015	0.015	0.0074	U
108-90-7	Chlorobenzene	0.10	0.10	0.051	0.022	0.022	0.011	U
100-41-4	Ethylbenzene	0.50	0.50	0.062	0.12	0.12	0.014	$\mathbf{U}$
179601-23-1	m,p-Xylenes	0.50	0.50	0.13	0.12	0.12	0.030	U
75-25-2	Bromoform	0.50	0.50	0.076	0.048	0.048	0.0074	$\mathbf{U}$
100-42-5	Styrene	0.50	0.50	0.076	0.12	0.12	0.018	U
95-47-6	o-Xylene	0.50	0.50	0.063	0.12	0.12	0.015	U
79-34-5	1,1,2,2-Tetrachloroethane	0.10	0.10	0.064	0.015	0.015	0.0093	U

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

#### RESULTS OF ANALYSIS

Page 3 of 3

**Tentatively Identified Compounds** 

Client:

**CH2M Hill** 

Client Sample ID: WAT-SG-FB-042308

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-008

Test Code:

EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/23/08

Date Received: 4/28/08

Instrument ID: Analyst:

Wida Ang

Date Analyzed: 5/2/08

Sampling Media:

Test Notes:

6.0 L Summa Canister

Volume(s) Analyzed:

1.00 Liter(s)

Container ID:

AC01209

Canister Dilution Factor: 1.00

GC/MS

**Compound Identification** 

Concentration

Data

Retention Time

Epichlorohydrin

 $\mu g/m^3$ 

Qualifier

NF

NF = Compound was searched for, but not found.

# RESULTS OF ANALYSIS

Page 1 of 3

Client: **CH2M Hill** 

Client Sample ID: WAT-SG-B4-042408 CAS Project ID: P0801215

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI CAS Sample ID: P0801215-009

Test Code: EPA TO-15

Date Collected: 4/24/08 Date Received: 4/28/08 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Analyst: Wida Ang Date Analyzed: 5/2/08

6.0 L Summa Canister Volume(s) Analyzed: Sampling Media: 1.00 Liter(s)

Test Notes: Container ID: AC00097

Initial Pressure (psig): -3.1 Final Pressure (psig): 3.5

C 1 C #	C	D a guyl4	MDI	MDI	D a guil4	MDI	MDI	Data
CAS#	Compound	Result	MRL	MDL	Result	MRL	MDL	Data
		μg/m³	μg/m³	μg/m³	ppbV	ppbV	ppbV	Qualifier
74-87-3	Chloromethane	0.56	0.16	0.079	0.27	0.076	0.038	
75-01-4	Vinyl Chloride	0.16	0.16	0.079	0.061	0.061	0.031	$\mathbf{U}_{-}$
74-83-9	Bromomethane	0.16	0.16	0.079	0.040	0.040	0.020	$\mathbf{U}$
75-00-3	Chloroethane	0.086	0.16	0.079	0.033	0.060	0.030	J
67-64-1	Acetone	35	7.9	0.11	15	3.3	0.048	В
107-13-1	Acrylonitrile	1.1	0.79	0.11	0.48	0.36	0.051	
75-35-4	1,1-Dichloroethene	0.11	0.16	0.079	0.029	0.040	0.020	J
75-09-2	Methylene Chloride	0.36	0.79	0.079	0.10	0.23	0.023	$\mathbf{J}$
75-15-0	Carbon Disulfide	0.92	0.79	0.19	0.29	0.25	0.061	
156-60-5	trans-1,2-Dichloroethene	0.16	0.16	0.079	0.040	0.040	0.020	$\mathbf{U}$
75-34-3	1,1-Dichloroethane	0.16	0.16	0.079	0.039	0.039	0.019	U
78-93-3	2-Butanone (MEK)	0.98	0.79	0.079	0.33	0.27	0.027	
156-59-2	cis-1,2-Dichloroethene	0.16	0.16	0.079	0.040	0.040	0.020	$\mathbf{U}$
67-66-3	Chloroform	1.3	0.16	0.093	0.26	0.032	0.019	
107-06-2	1,2-Dichloroethane	0.16	0.16	0.079	0.039	0.039	0.019	$\mathbf{U}$
71-55-6	1,1,1-Trichloroethane	0.16	0.16	0.079	0.029	0.029	0.014	U
71-43-2	Benzene	0.48	0.16	0.079	0.15	0.049	0.025	
56-23-5	Carbon Tetrachloride	0.38	0.16	0.079	0.060	0.025	0.012	

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

B = Analyte was found in the method blank.

### **RESULTS OF ANALYSIS**

Page 2 of 3

**Client:** 

**CH2M Hill** 

Client Sample ID: WAT-SG-B4-042408

6.0 L Summa Canister

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-009

Test Code:

EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/24/08 Date Received: 4/28/08

Instrument ID: Analyst:

Sampling Media:

Wida Ang

Date Analyzed: 5/2/08

Volume(s) Analyzed:

1.00 Liter(s)

Test Notes:

Container ID:

AC00097

Initial Pressure (psig):

-3.1

Final Pressure (psig):

3.5

CAS#	Compound	Result μg/m³	MRL μg/m³	MDL μg/m³	Result ppbV	MRL ppbV	MDL ppbV	Data Qualifier
78-87-5	1,2-Dichloropropane	<u>μg/III</u> 0.16	0.16	$\frac{\mu g/m}{0.079}$	0.034	0.034	$\frac{pp0  V}{0.017}$	U
75-27-4	Bromodichloromethane	0.16	0.16	0.079	0.023	0.023	0.012	Ü
79-01-6	Trichloroethene	0.45	0.16	0.079	0.083	0.029	0.015	
10061-01-5	cis-1,3-Dichloropropene	0.79	0.79	0.082	0.17	0.17	0.018	U
108-10-1	4-Methyl-2-pentanone	94	0.79	0.088	23	0.19	0.021	
10061-02-6	trans-1,3-Dichloropropene	0.79	0.79	0.099	0.17	0.17	0.022	U
79-00-5	1,1,2-Trichloroethane	0.16	0.16	0.079	0.029	0.029	0.014	$\mathbf{U}$
108-88-3	Toluene	13	0.79	0.079	3.5	0.21	0.021	
124-48-1	Dibromochloromethane	0.16	0.16	0.11	0.018	0.018	0.013	$\mathbf{U}$
127-18-4	Tetrachloroethene	0.16	0.16	0.079	0.023	0.023	0.012	$\mathbf{U}$
108-90-7	Chlorobenzene	0.16	0.16	0.080	0.034	0.034	0.017	U
100-41-4	Ethylbenzene	0.25	0.79	0.097	0.057	0.18	0.022	J
179601-23-1	m,p-Xylenes	0.83	0.79	0.20	0.19	0.18	0.047	
75-25-2	Bromoform	0.79	0.79	0.12	0.076	0.076	0.012	U
100-42-5	Styrene	0.12	0.79	0.12	0.028	0.18	0.028	J
95-47-6	o-Xylene	0.21	0.79	0.099	0.049	0.18	0.023	J
79-34-5	1,1,2,2-Tetrachloroethane	0.16	0.16	0.10	0.023	0.023	0.015	U

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

#### **RESULTS OF ANALYSIS**

Page 3 of 3

**Client:** 

**CH2M Hill** 

Client Sample ID: WAT-SG-B4-042408

6.0 L Summa Canister

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-009

**Tentatively Identified Compounds** 

Date Collected: 4/24/08

Date Received: 4/28/08

Instrument ID:

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

EPA TO-15

Date Analyzed: 5/2/08

Analyst:

Test Code:

Wida Ang

Volume(s) Analyzed:

1.00 Liter(s)

Sampling Media: Test Notes:

 $\mathbf{T}$ 

Container ID:

AC00097

Initial Pressure (psig):

-3.1

Final Pressure (psig):

3.5

GC/MS Compound Identification		Concentration	Data
Retention Time		$\mu \mathrm{g}/\mathrm{m}^3$	Qualifier
4.46	Propane + Propene	20	
5.18	n-Butane	10	
16.45	Hexamethylcyclotrisiloxane	10	*
18.87	Benzaldehyde	9	
21.27	n-Nonanal	40	
	Epichlorohydrin	NF	

T = Analyte is a tentatively identified compound, result is estimated.

NF = Compound was searched for, but not found.

<sup>\* =</sup> Possible artifact

#### **RESULTS OF ANALYSIS**

Page 1 of 3

Client: CH2M Hill

Client Sample ID: WAT-SG-5-042408 CAS Project ID: P0801215

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI CAS Sample ID: P0801215-010

Test Code: EPA TO-15 Date Collected: 4/24/08

Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16 Date Received: 4/28/08
Analyst: Wida Ang Date Analyzed: 5/2/08

Sampling Media: 6.0 L Summa Canister Volume(s) Analyzed: 0.15 Liter(s)

Test Notes:

Container ID: AC00802

Initial Pressure (psig): -1.7 Final Pressure (psig): 3.5

CAS#	Compound	Result μg/m³	MRL μg/m³	MDL μg/m³	Result ppbV	MRL ppbV	MDL ppbV	Data Qualifier
74-87-3	Chloromethane	0.93	0.93	0.47	0.45	0.45	0.23	U
75-01-4	Vinyl Chloride	0.93	0.93	0.47	0.37	0.37	0.18	${f U}$
74-83-9	Bromomethane	0.93	0.93	0.47	0.24	0.24	0.12	$\mathbf{U}$
75-00-3	Chloroethane	0.93	0.93	0.47	0.35	0.35	0.18	U
67-64-1	Acetone	10	47	0.68	4.3	20	0.29	J, B
107-13-1	Acrylonitrile	4.7	4.7	0.65	2.2	2.2	0.30	U
75-35-4	1,1-Dichloroethene	0.93	0.93	0.47	0.24	0.24	0.12	U
75-09-2	Methylene Chloride	4.7	4.7	0.47	1.3	1.3	0.13	U
75-15-0	Carbon Disulfide	2.4	4.7	1.1	0.78	1.5	0.36	J
156-60-5	trans-1,2-Dichloroethene	0.93	0.93	0.47	0.24	0.24	0.12	$\mathbf{U}$
75-34-3	1,1-Dichloroethane	1.4	0.93	0.47	0.33	0.23	0.12	
78-93-3	2-Butanone (MEK)	0.97	4.7	0.47	0.33	1.6	0.16	J
156-59-2	cis-1,2-Dichloroethene	0.93	0.93	0.47	0.24	0.24	0.12	$\mathbf{U}$
67-66-3	Chloroform	1,100	0.93	0.55	230	0.19	0.11	
107-06-2	1,2-Dichloroethane	0.93	0.93	0.47	0.23	0.23	0.12	U
71-55-6	1,1,1-Trichloroethane	0.93	0.93	0.47	0.17	0.17	0.086	U
71-43-2	Benzene	0.93	0.93	0.47	0.29	0.29	0.15	$\mathbf{U}$
56-23-5	Carbon Tetrachloride	0.93	0.93	0.47	0.15	0.15	0.074	U

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

B = Analyte was found in the method blank.

#### **RESULTS OF ANALYSIS**

Page 2 of 3

Client:

**CH2M Hill** 

Client Sample ID: WAT-SG-5-042408

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-010

Test Code:

Analyst:

EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/24/08 Date Received: 4/28/08

Date Analyzed: 5/2/08

Sampling Media:

Instrument ID:

Wida Ang 6.0 L Summa Canister

Volume(s) Analyzed:

0.15 Liter(s)

Test Notes:

Container ID:

AC00802

Initial Pressure (psig):

-1.7

Final Pressure (psig):

3.5

CAS#	Compound	Result μg/m³	MRL μg/m³	MDL μg/m³	Result ppbV	MRL ppbV	MDL ppbV	Data Qualifier
78-87-5	1,2-Dichloropropane	0.93	0.93	0.47	0.20	0.20	0.10	U
75-27-4	Bromodichloromethane	0.93	0.93	0.47	0.14	0.14	0.070	${f U}$
79-01-6	Trichloroethene	2.4	0.93	0.47	0.45	0.17	0.087	
10061-01-5	cis-1,3-Dichloropropene	4.7	4.7	0.49	1.0	1.0	0.11	$\mathbf{U}$
108-10-1	4-Methyl-2-pentanone	3.0	4.7	0.52	0.74	1.1	0.13	$\mathbf{J}$
10061-02-6	trans-1,3-Dichloropropene	4.7	4.7	0.59	1.0	1.0	0.13	U
79-00-5	1,1,2-Trichloroethane	0.93	0.93	0.47	0.17	0.17	0.086	U
108-88-3	Toluene	0.85	4.7	0.47	0.23	1.2	0.12	J
124-48-1	Dibromochloromethane	0.93	0.93	0.63	0.11	0.11	0.075	$\mathbf{U}$
127-18-4	Tetrachloroethene	3.8	0.93	0.47	0.56	0.14	0.069	
108-90-7	Chlorobenzene	0.93	0.93	0.48	0.20	0.20	0.10	U
100-41-4	Ethylbenzene	4.7	4.7	0.58	1.1	1.1	0.13	$\mathbf{U}$
179601-23-1	m,p-Xylenes	4.7	4.7	1.2	1.1	1.1	0.28	$\mathbf{U}$
75-25-2	Bromoform	4.7	4.7	0.71	0.45	0.45	0.069	$\mathbf{U}$
100-42-5	Styrene	4.7	4.7	0.71	1.1	1.1	0.17	U
95-47-6	o-Xylene	4.7	4.7	0.59	1.1	1.1	0.14	U
79-34-5	1,1,2,2-Tetrachloroethane	0.93	0.93	0.60	0.14	0.14	0.087	U

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

#### **RESULTS OF ANALYSIS**

Page 3 of 3

Client:

**CH2M Hill** 

Client Sample ID: WAT-SG-5-042408

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-010

**Tentatively Identified Compounds** 

Test Code:

EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/24/08

Instrument ID: Analyst:

Wida Ang

Date Received: 4/28/08

Date Analyzed: 5/2/08

0.15 Liter(s)

Sampling Media:

Test Notes:

6.0 L Summa Canister T

Container ID:

AC00802

Initial Pressure (psig):

-1.7

Final Pressure (psig):

3.5

Volume(s) Analyzed:

GC/MS	Compound Identification	Concentration	Data
Retention Time		$\mu \mathrm{g}/\mathrm{m}^3$	Qualifier
16.27	C <sub>6</sub> HF <sub>13</sub> (Tridecafluorohexane)	80	
22.07	Hexamethylcyclotrisiloxane	100	*
24.09	Hexamethylcyclotrisiloxane	30	*
	Epichlorohydrin	NF	

T = Analyte is a tentatively identified compound, result is estimated.

NF = Compound was searched for, but not found.

<sup>\* =</sup> Possible artifact

#### RESULTS OF ANALYSIS

Page 1 of 3

Client:

**CH2M Hill** 

Client Sample ID: WAT-SG-5a-042408

6.0 L Summa Canister

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-011

Test Code:

Analyst:

Instrument ID:

Sampling Media:

EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/24/08 Date Received: 4/28/08

Wida Ang

Date Analyzed: 5/5/08

Volume(s) Analyzed:

0.30 Liter(s)

Test Notes:

Container ID:

AC00615

Initial Pressure (psig):

0.0

Final Pressure (psig):

3.5

CAS#	Compound	Result	MRL	MDL	Result	MRL	MDL	Data
		$\mu g/m^3$	$\mu g/m^3$	$\mu g/m^3$	${f ppbV}$	ppbV	ppbV	Qualifier
74-87-3	Chloromethane	0.41	0.41	0.21	0.20	0.20	0.10	U
75-01-4	Vinyl Chloride	0.41	0.41	0.21	0.16	0.16	0.081	$\mathbf{U}$
74-83-9	Bromomethane	0.41	0.41	0.21	0.11	0.11	0.053	$\mathbf{U}$
75-00-3	Chloroethane	0.41	0.41	0.21	0.16	0.16	0.078	$\mathbf{U}$
67-64-1	Acetone	12	21	0.30	5.1	8.7	0.13	J, B
107-13-1	Acrylonitrile	2.1	2.1	0.29	0.95	0.95	0.13	U
75-35-4	1,1-Dichloroethene	0.41	0.41	0.21	0.10	0.10	0.052	$\mathbf{U}$
75-09-2	Methylene Chloride	0.29	2.1	0.21	0.083	0.60	0.060	J
75-15-0	Carbon Disulfide	20	2.1	0.50	6.5	0.66	0.16	
156-60-5	trans-1,2-Dichloroethene	0.41	0.41	0.21	0.10	0.10	0.052	$\mathbf{U}$
75-34-3	1,1-Dichloroethane	0.41	0.41	0.21	0.10	0.10	0.051	U
78-93-3	2-Butanone (MEK)	1.3	2.1	0.21	0.45	0.70	0.070	J
156-59-2	cis-1,2-Dichloroethene	2.1	0.41	0.21	0.54	0.10	0.052	
67-66-3	Chloroform	450	0.41	0.24	93	0.085	0.050	
107-06-2	1,2-Dichloroethane	0.41	0.41	0.21	0.10	0.10	0.051	$\mathbf{U}$
71-55-6	1,1,1-Trichloroethane	0.41	0.41	0.21	0.076	0.076	0.038	U
71-43-2	Benzene	0.41	0.41	0.21	0.13	0.13	0.065	$\mathbf{U}$
56-23-5	Carbon Tetrachloride	0.55	0.41	0.21	0.087	0.066	0.033	

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

B = Analyte was found in the method blank.

#### **RESULTS OF ANALYSIS**

Page 2 of 3

**Client:** 

**CH2M Hill** 

Client Sample ID: WAT-SG-5a-042408

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-011

Test Code:

EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/24/08 Date Received: 4/28/08

Instrument ID: Analyst:

Wida Ang

Date Analyzed: 5/5/08

Sampling Media:

6.0 L Summa Canister

Volume(s) Analyzed:

0.30 Liter(s)

Test Notes:

Container ID:

AC00615

Initial Pressure (psig):

0.0

Final Pressure (psig):

3.5

CAS#	Compound	Result μg/m³	MRL μg/m³	$\begin{array}{c} MDL \\ \mu g/m^3 \end{array}$	Result ppbV	MRL ppbV	MDL ppbV	Data Qualifier
78-87-5	1,2-Dichloropropane	0.41	$\frac{\mu g/m}{0.41}$	0.21	0.089	0.089	$\frac{pp0.7}{0.045}$	U
75-27-4	Bromodichloromethane	0.43	0.41	0.21	0.064	0.062	0.031	
79-01-6	Trichloroethene	20	0.41	0.21	3.6	0.077	0.038	
10061-01-5	cis-1,3-Dichloropropene	2.1	2.1	0.21	0.46	0.46	0.047	U
108-10-1	4-Methyl-2-pentanone	2.3	2.1	0.23	0.55	0.50	0.057	
10061-02-6	trans-1,3-Dichloropropene	2.1	2.1	0.26	0.46	0.46	0.057	U
79-00-5	1,1,2-Trichloroethane	0.41	0.41	0.21	0.076	0.076	0.038	$\mathbf{U}$
108-88-3	Toluene	0.55	2.1	0.21	0.15	0.55	0.055	J
124-48-1	Dibromochloromethane	0.41	0.41	0.28	0.049	0.049	0.033	$\mathbf{U}$
127-18-4	Tetrachloroethene	4.1	0.41	0.21	0.61	0.061	0.030	
108-90-7	Chlorobenzene	0.41	0.41	0.21	0.090	0.090	0.046	U
100-41-4	Ethylbenzene	2.1	2.1	0.26	0.48	0.48	0.059	$\mathbf{U}$
179601-23-1	m,p-Xylenes	0.58	2.1	0.54	0.13	0.48	0.12	J
75-25-2	Bromoform	2.1	2.1	0.31	0.20	0.20	0.030	U
100-42-5	Styrene	2.1	2.1	0.31	0.49	0.49	0.074	$\mathbf{U}$
95-47-6	o-Xylene	0.27	2.1	0.26	0.062	0.48	0.060	J
79-34-5	1,1,2,2-Tetrachloroethane	0.41	0.41	0.26	0.060	0.060	0.039	U

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

#### **RESULTS OF ANALYSIS**

Page 3 of 3

**Client:** 

**CH2M Hill** 

Client Sample ID: WAT-SG-5a-042408

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-011

**Tentatively Identified Compounds** 

Test Code: Instrument ID: EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/24/08

Date Received: 4/28/08

Analyst:

Wida Ang

Date Analyzed: 5/5/08 Volume(s) Analyzed:

0.30 Liter(s)

Sampling Media: Test Notes:

6.0 L Summa Canister T

Container ID:

AC00615

Initial Pressure (psig):

0.0

Final Pressure (psig):

3.5

GC/MS	Compound Identification	Concentration	Data
Retention Time		$\mu \mathrm{g}/\mathrm{m}^3$	Qualifier
16.27	C <sub>6</sub> HF <sub>13</sub> (Tridecafluorohexane)	50	
21.27	n-Nonanal	20	
22.69	Naphthalene	50	
24.03	2-Methylnaphthalene	70	
24.25	1-Methylnaphthalene	70	
	Epichlorohydrin	NF	

T = Analyte is a tentatively identified compound, result is estimated.

NF = Compound was searched for, but not found.

#### RESULTS OF ANALYSIS

Page 1 of 3

**Client:** 

CH2M Hill

Client Sample ID: WAT-SG-6-042408

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-012

Test Code:

EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/24/08

Date Received: 4/28/08

Instrument ID: Analyst:

Wida Ang

Date Analyzed: 5/5/08

Sampling Media:

6.0 L Summa Canister

Volume(s) Analyzed:

0.030 Liter(s)

Test Notes:

Container ID:

AC00814

Initial Pressure (psig):

0.0

Final Pressure (psig):

3.5

CAS#         Compound         Result µg/m³         MRL µg/m³         MDL µg/m³         Result ppbV         MRL ppbV ppbV ppbV Qualifier         Data ppbV Qualifier           74-87-3         Chloromethane         4.1         4.1         2.1         2.0         2.0         1.0         U           75-01-4         Vinyl Chloride         4.1         4.1         2.1         1.6         1.6         0.81         U           75-01-3         Bromomethane         4.1         4.1         2.1         1.1         1.1         0.53         U           75-00-3         Chloroethane         4.1         4.1         2.1         1.1         1.1         0.53         U           67-64-1         Acetone         26         210         3.0         11         87         1.3         J,8           107-13-1         Acrylonitrile         21         2.1         2.9         9.5         9.5         1.3         U           75-35-4         1,1-Dichloroethene         4.1         4.1         2.1         1.0         1.0         0.52         U           75-15-0         Carbon Disulfide         2,600         21         5.0         840         6.6         1.6         1.5									
74-87-3         Chloromethane         4.1         4.1         2.1         2.0         2.0         1.0         U           75-01-4         Vinyl Chloride         4.1         4.1         2.1         1.6         1.6         0.81         U           74-83-9         Bromomethane         4.1         4.1         2.1         1.1         1.1         0.53         U           75-00-3         Chloroethane         4.1         4.1         2.1         1.6         1.6         0.78         U           67-64-1         Acetone         26         210         3.0         11         87         1.3         J,8           107-13-1         Acrylonitrile         21         21         2.9         9.5         9.5         1.3         U           75-35-4         1,1-Dichloroethene         4.1         4.1         2.1         1.0         1.0         0.52         U           75-15-0         Methylene Chloride         2.6         21         2.1         0.74         6.0         0.60         J           75-15-0         Carbon Disulfide         2,600         21         5.0         840         6.6         1.6           156-60-5         trans-1,2-Dichloroethane </th <th>CAS#</th> <th>Compound</th> <th></th> <th></th> <th></th> <th></th> <th></th> <th></th> <th></th>	CAS#	Compound							
74-83-9         Bromomethane         4.1         4.1         2.1         1.1         1.1         0.53         U           75-00-3         Chloroethane         4.1         4.1         2.1         1.6         1.6         0.78         U           67-64-1         Acetone         26         210         3.0         11         87         1.3         J, B           107-13-1         Acrylonitrile         21         21         2.9         9.5         9.5         1.3         U           75-35-4         1,1-Dichloroethene         4.1         4.1         2.1         1.0         1.0         0.52         U           75-09-2         Methylene Chloride         2.6         21         2.1         0.74         6.0         0.60         J           75-15-0         Carbon Disulfide         2,600         21         5.0         840         6.6         1.6           156-60-5         trans-1,2-Dichloroethene         4.1         4.1         2.1         1.0         1.0         0.52         U           78-93-3         2-Butanone (MEK)         21         21         2.1         7.0         7.0         0.70         U           156-59-2         cis-1,2-Dich	74-87-3	Chloromethane	4.1	4.1	2.1	2.0	2.0	1.0	U
75-00-3         Chloroethane         4.1         4.1         2.1         1.6         1.6         0.78         U           67-64-1         Acetone         26         210         3.0         11         87         1.3         J, B           107-13-1         Acrylonitrile         21         21         2.9         9.5         9.5         1.3         U           75-35-4         1,1-Dichloroethene         4.1         4.1         2.1         1.0         1.0         0.52         U           75-09-2         Methylene Chloride         2.6         21         2.1         0.74         6.0         0.60         J           75-15-0         Carbon Disulfide         2.600         21         5.0         840         6.6         1.6           156-60-5         trans-1,2-Dichloroethene         4.1         4.1         2.1         1.0         1.0         0.52         U           75-34-3         1,1-Dichloroethane         4.1         4.1         2.1         1.0         1.0         0.51         U           78-93-3         2-Butanone (MEK)         21         21         2.1         7.0         7.0         0.70         0.70         U           67-66-3 <td>75-01-4</td> <td>Vinyl Chloride</td> <td>4.1</td> <td>4.1</td> <td>2.1</td> <td>1.6</td> <td>1.6</td> <td>0.81</td> <td><math>\mathbf{U}</math></td>	75-01-4	Vinyl Chloride	4.1	4.1	2.1	1.6	1.6	0.81	$\mathbf{U}$
67-64-1         Acetone         26         210         3.0         11         87         1.3         J, B           107-13-1         Acrylonitrile         21         21         2.9         9.5         9.5         1.3         U           75-35-4         1,1-Dichloroethene         4.1         4.1         2.1         1.0         1.0         0.52         U           75-09-2         Methylene Chloride         2.6         21         2.1         0.74         6.0         0.60         J           75-15-0         Carbon Disulfide         2,600         21         5.0         840         6.6         1.6           156-60-5         trans-1,2-Dichloroethene         4.1         4.1         2.1         1.0         1.0         0.52         U           75-34-3         1,1-Dichloroethane         4.1         4.1         2.1         1.0         1.0         0.51         U           78-93-3         2-Butanone (MEK)         21         21         2.1         7.0         7.0         0.70         U           67-66-3         Chloroform         76         4.1         2.4         16         0.85         0.50           107-06-2         1,2-Dichloroethane	74-83-9	Bromomethane	4.1	4.1	2.1	1.1	1.1	0.53	$\mathbf{U}$
107-13-1         Acrylonitrile         21         21         2.9         9.5         9.5         1.3         U           75-35-4         1,1-Dichloroethene         4.1         4.1         2.1         1.0         1.0         0.52         U           75-09-2         Methylene Chloride         2.6         21         2.1         0.74         6.0         0.60         J           75-15-0         Carbon Disulfide         2,600         21         5.0         840         6.6         1.6           156-60-5         trans-1,2-Dichloroethene         4.1         4.1         2.1         1.0         1.0         0.52         U           75-34-3         1,1-Dichloroethane         4.1         4.1         2.1         1.0         1.0         0.51         U           78-93-3         2-Butanone (MEK)         21         21         2.1         7.0         7.0         0.70         U           156-59-2         cis-1,2-Dichloroethene         4.1         4.1         2.1         1.0         1.0         0.52         U           67-66-3         Chloroform         76         4.1         2.4         16         0.85         0.50           107-06-2         1,2-Dichlor	75-00-3	Chloroethane	4.1	4.1	2.1	1.6	1.6	0.78	$\mathbf{U}$
75-35-4         1,1-Dichloroethene         4.1         4.1         2.1         1.0         1.0         0.52         U           75-09-2         Methylene Chloride         2.6         21         2.1         0.74         6.0         0.60         J           75-15-0         Carbon Disulfide         2,600         21         5.0         840         6.6         1.6           156-60-5         trans-1,2-Dichloroethene         4.1         4.1         2.1         1.0         1.0         0.52         U           75-34-3         1,1-Dichloroethane         4.1         4.1         2.1         1.0         1.0         0.51         U           78-93-3         2-Butanone (MEK)         21         21         2.1         7.0         7.0         0.70         U           156-59-2         cis-1,2-Dichloroethene         4.1         4.1         2.1         1.0         1.0         0.52         U           67-66-3         Chloroform         76         4.1         2.4         16         0.85         0.50           107-06-2         1,2-Dichloroethane         4.1         4.1         2.1         1.0         1.0         0.51         U           71-55-6         1,1,	67-64-1	Acetone	26	210	3.0	11	87	1.3	J, B
75-09-2         Methylene Chloride         2.6         21         2.1         0.74         6.0         0.60         J           75-15-0         Carbon Disulfide         2,600         21         5.0         840         6.6         1.6           156-60-5         trans-1,2-Dichloroethene         4.1         4.1         2.1         1.0         1.0         0.52         U           75-34-3         1,1-Dichloroethane         4.1         4.1         2.1         1.0         1.0         0.51         U           78-93-3         2-Butanone (MEK)         21         21         2.1         7.0         7.0         0.70         U           156-59-2         cis-1,2-Dichloroethene         4.1         4.1         2.1         1.0         1.0         0.52         U           67-66-3         Chloroform         76         4.1         2.4         16         0.85         0.50           107-06-2         1,2-Dichloroethane         4.1         4.1         2.1         1.0         1.0         0.51         U           71-43-2         Benzene         4.1         4.1         2.1         1.3         1.3         0.65         U	107-13-1	Acrylonitrile	21	21	2.9	9.5	9.5	1.3	U
75-15-0         Carbon Disulfide         2,600         21         5.0         840         6.6         1.6           156-60-5         trans-1,2-Dichloroethene         4.1         4.1         2.1         1.0         1.0         0.52         U           75-34-3         1,1-Dichloroethane         4.1         4.1         2.1         1.0         1.0         0.51         U           78-93-3         2-Butanone (MEK)         21         21         2.1         7.0         7.0         0.70         U           156-59-2         cis-1,2-Dichloroethene         4.1         4.1         2.1         1.0         1.0         0.52         U           67-66-3         Chloroform         76         4.1         2.4         16         0.85         0.50           107-06-2         1,2-Dichloroethane         4.1         4.1         2.1         1.0         1.0         0.51         U           71-55-6         1,1,1-Trichloroethane         4.1         4.1         2.1         0.76         0.76         0.38         U           71-43-2         Benzene         4.1         4.1         2.1         1.3         1.3         0.65         U	75-35-4	1,1-Dichloroethene	4.1	4.1	2.1	1.0	1.0	0.52	$\mathbf{U}$
156-60-5         trans-1,2-Dichloroethene         4.1         4.1         2.1         1.0         1.0         0.52         U           75-34-3         1,1-Dichloroethane         4.1         4.1         2.1         1.0         1.0         0.51         U           78-93-3         2-Butanone (MEK)         21         21         2.1         7.0         7.0         0.70         U           156-59-2         cis-1,2-Dichloroethene         4.1         4.1         2.1         1.0         1.0         0.52         U           67-66-3         Chloroform         76         4.1         2.4         16         0.85         0.50           107-06-2         1,2-Dichloroethane         4.1         4.1         2.1         1.0         1.0         0.51         U           71-55-6         1,1,1-Trichloroethane         4.1         4.1         2.1         0.76         0.76         0.38         U           71-43-2         Benzene         4.1         4.1         2.1         1.3         1.3         0.65         U	75-09-2	Methylene Chloride	2.6	21	2.1	0.74	6.0	0.60	J
75-34-3       1,1-Dichloroethane       4.1       4.1       2.1       1.0       1.0       0.51       U         78-93-3       2-Butanone (MEK)       21       21       2.1       7.0       7.0       0.70       U         156-59-2       cis-1,2-Dichloroethene       4.1       4.1       2.1       1.0       1.0       0.52       U         67-66-3       Chloroform       76       4.1       2.4       16       0.85       0.50         107-06-2       1,2-Dichloroethane       4.1       4.1       2.1       1.0       1.0       0.51       U         71-55-6       1,1,1-Trichloroethane       4.1       4.1       2.1       0.76       0.76       0.38       U         71-43-2       Benzene       4.1       4.1       2.1       1.3       1.3       0.65       U	75-15-0	Carbon Disulfide	2,600	21	5.0	840	6.6	1.6	
78-93-3       2-Butanone (MEK)       21       21       2.1       7.0       7.0       0.70       U         156-59-2       cis-1,2-Dichloroethene       4.1       4.1       2.1       1.0       1.0       0.52       U         67-66-3       Chloroform       76       4.1       2.4       16       0.85       0.50         107-06-2       1,2-Dichloroethane       4.1       4.1       2.1       1.0       1.0       0.51       U         71-55-6       1,1,1-Trichloroethane       4.1       4.1       2.1       0.76       0.76       0.38       U         71-43-2       Benzene       4.1       4.1       2.1       1.3       1.3       0.65       U	156-60-5	trans-1,2-Dichloroethene	4.1	4.1	2.1	1.0	1.0	0.52	$\mathbf{U}$
156-59-2         cis-1,2-Dichloroethene         4.1         4.1         2.1         1.0         1.0         0.52         U           67-66-3         Chloroform         76         4.1         2.4         16         0.85         0.50           107-06-2         1,2-Dichloroethane         4.1         4.1         2.1         1.0         1.0         0.51         U           71-55-6         1,1,1-Trichloroethane         4.1         4.1         2.1         0.76         0.76         0.38         U           71-43-2         Benzene         4.1         4.1         2.1         1.3         1.3         0.65         U	75-34-3	1,1-Dichloroethane	4.1	4.1	2.1	1.0	1.0	0.51	U
67-66-3         Chloroform         76         4.1         2.4         16         0.85         0.50           107-06-2         1,2-Dichloroethane         4.1         4.1         2.1         1.0         1.0         0.51         U           71-55-6         1,1,1-Trichloroethane         4.1         4.1         2.1         0.76         0.76         0.38         U           71-43-2         Benzene         4.1         4.1         2.1         1.3         1.3         0.65         U	78-93-3	2-Butanone (MEK)	21	21	2.1	7.0	7.0	0.70	$\mathbf{U}$
107-06-2       1,2-Dichloroethane       4.1       4.1       2.1       1.0       1.0       0.51       U         71-55-6       1,1,1-Trichloroethane       4.1       4.1       2.1       0.76       0.76       0.38       U         71-43-2       Benzene       4.1       4.1       2.1       1.3       1.3       0.65       U	156-59-2	cis-1,2-Dichloroethene	4.1	4.1	2.1	1.0	1.0	0.52	$\mathbf{U}$
71-55-6 1,1,1-Trichloroethane 4.1 4.1 2.1 0.76 0.76 0.38 U 71-43-2 Benzene 4.1 4.1 2.1 1.3 1.3 0.65 U	67-66-3	Chloroform	76	4.1	2.4	16	0.85	0.50	
71-43-2 Benzene 4.1 4.1 2.1 1.3 1.3 0.65 U	107-06-2	1,2-Dichloroethane	4.1	4.1	2.1	1.0	1.0	0.51	U
	71-55-6	1,1,1-Trichloroethane	4.1	4.1	2.1	0.76	0.76	0.38	U
56-23-5 Carbon Tetrachloride 4.1 4.1 2.1 0.66 0.66 0.33 U	71-43-2	Benzene	4.1	4.1	2.1	1.3	1.3	0.65	U
	56-23-5	Carbon Tetrachloride	4.1	4.1	2.1	0.66	0.66	0.33	U

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

B = Analyte was found in the method blank.

# RESULTS OF ANALYSIS

Page 2 of 3

Client:

CH2M Hill

Client Sample ID: WAT-SG-6-042408

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-012

Test Code:

EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/24/08

Date Received: 4/28/08

Instrument ID: Analyst:

Wida Ang

Date Analyzed: 5/5/08

Sampling Media:

6.0 L Summa Canister

Volume(s) Analyzed:

0.030 Liter(s)

Test Notes:

Container ID:

AC00814

Initial Pressure (psig):

0.0

Final Pressure (psig):

3.5

Canister Dilution Factor: 1.24

CAS#	Compound	Result μg/m³	MRL μg/m³	MDL μg/m³	Result ppbV	MRL ppbV	MDL ppbV	Data Qualifier
78-87-5	1,2-Dichloropropane	4.1	4.1	2.1	0.89	0.89	0.45	U
75-27-4	Bromodichloromethane	4.1	4.1	2.1	0.62	0.62	0.31	$\mathbf{U}$
79-01-6	Trichloroethene	4.1	4.1	2.1	0.77	0.77	0.38	$\mathbf{U}$
10061-01-5	cis-1,3-Dichloropropene	21	21	2.1	4.6	4.6	0.47	$\mathbf{U}$
108-10-1	4-Methyl-2-pentanone	110	21	2.3	26	5.0	0.57	
10061-02-6	trans-1,3-Dichloropropene	21	21	2.6	4.6	4.6	0.57	U
79-00-5	1,1,2-Trichloroethane	4.1	4.1	2.1	0.76	0.76	0.38	$\mathbf{U}$
108-88-3	Toluene	4.8	21	2.1	1.3	5.5	0.55	J
124-48-1	Dibromochloromethane	4.1	4.1	2.8	0.49	0.49	0.33	U
127-18-4	Tetrachloroethene	3.4	4.1	2.1	0.51	0.61	0.30	J
108-90-7	Chlorobenzene	4.1	4.1	2.1	0.90	0.90	0.46	U
100-41-4	Ethylbenzene	21	21	2.6	4.8	4.8	0.59	$\mathbf{U}$
179601-23-1	m,p-Xylenes	7.2	21	5.4	1.6	4.8	1.2	J
75-25-2	Bromoform	21	21	3.1	2.0	2.0	0.30	$\mathbf{U}$
100-42-5	Styrene	21	21	3.1	4.9	4.9	0.74	$\mathbf{U}$
95-47-6	o-Xylene	2.9	21	2.6	0.66	4.8	0.60	J
79-34-5	1,1,2,2-Tetrachloroethane	4.1	4.1	2.6	0.60	0.60	0.39	U

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

#### RESULTS OF ANALYSIS

Page 3 of 3

Client:

CH2M Hill

Client Sample ID: WAT-SG-6-042408

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-012

**Tentatively Identified Compounds** 

Test Code:

EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/24/08

Date Received: 4/28/08

Instrument ID: Analyst:

Date Analyzed: 5/5/08

Sampling Media:

Wida Ang 6.0 L Summa Canister

Volume(s) Analyzed:

0.030 Liter(s)

Test Notes:

T

Container ID:

AC00814

Initial Pressure (psig):

0.0

Final Pressure (psig):

3.5

Canister Dilution Factor: 1.24

GC/MS	Compound Identification	Concentration	Data
Retention Time		$\mu \mathrm{g}/\mathrm{m}^3$	Qualifier
4.45	Propane +Carbonyl Sulfide	100	
13.51	Dimethyl Sulfide	1,000	
	Epichlorohydrin	NF	

T = Analyte is a tentatively identified compound, result is estimated.

NF = Compound was searched for, but not found.

Verified By:

TO15SCAN.XLT - NL - PageNo .:

# **RESULTS OF ANALYSIS**

Page 1 of 3

**Client:** 

**CH2M Hill** 

Client Sample ID: WAT-SG-7-042408

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-013

Test Code: Instrument ID: EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/24/08 Date Received: 4/28/08

Wida Ang

Date Analyzed: 5/2/08

Sampling Media:

6.0 L Summa Canister

Volume(s) Analyzed: 0.0080 Liter(s)

Test Notes:

Analyst:

Container ID:

AC01276

Initial Pressure (psig):

-3.5

Final Pressure (psig):

3.5

Canister Dilution Factor: 1.63

Compound	Result	MRL	MDL	Result	MRL	MDL	Data
	$\mu g/m^3$	$\mu g/m^3$	$\mu g/m^3$	${f ppbV}$	ppbV	ppbV	Qualifier
Chloromethane	20	20	10	9.9	9.9	4.9	U
Vinyl Chloride	20	20	10	8.0	8.0	4.0	$\mathbf{U}$
Bromomethane	20	20	10	5.2	5.2	2.6	$\mathbf{U}$
Chloroethane	20	20	10	7.7	7.7	3.9	$\mathbf{U}$
Acetone	1,800	1,000	15	760	430	6.3	В
Acrylonitrile	100	100	14	47	47	6.6	U
1,1-Dichloroethene	20	20	10	5.1	5.1	2.6	U
Methylene Chloride	35	100	10	10	29	2.9	J
Carbon Disulfide	310	100	24	100	33	7.9	
trans-1,2-Dichloroethene	20	20	10	5.1	5.1	2.6	$\mathbf{U}$
1,1-Dichloroethane	84	20	10	21	5.0	2.5	
2-Butanone (MEK)	17	100	10	5.8	35	3.5	J
cis-1,2-Dichloroethene	19	20	10	4.9	5.1	2.6	J
Chloroform	24,000	20	12	4,800	4.2	2.5	
1,2-Dichloroethane	20	20	10	5.0	5.0	2.5	U
1,1,1-Trichloroethane	20	20	10	3.7	3.7	1.9	U
Benzene	20	20	10	6.4	6.4	3.2	$\mathbf{U}$
Carbon Tetrachloride	20	20	10	3.2	3.2	1.6	U
	Chloromethane Vinyl Chloride Bromomethane Chloroethane Acetone Acrylonitrile 1,1-Dichloroethene Methylene Chloride Carbon Disulfide trans-1,2-Dichloroethene 1,1-Dichloroethane 2-Butanone (MEK) cis-1,2-Dichloroethene Chloroform 1,2-Dichloroethane 1,1,1-Trichloroethane Benzene	μg/m³           Chloromethane         20           Vinyl Chloride         20           Bromomethane         20           Chloroethane         20           Acetone         1,800           Acrylonitrile         100           1,1-Dichloroethene         20           Methylene Chloride         35           Carbon Disulfide         310           trans-1,2-Dichloroethene         20           1,1-Dichloroethane         84           2-Butanone (MEK)         17           cis-1,2-Dichloroethene         19           Chloroform         24,000           1,2-Dichloroethane         20           1,1,1-Trichloroethane         20           Benzene         20	μg/m³         μg/m³           Chloromethane         20         20           Vinyl Chloride         20         20           Bromomethane         20         20           Chloroethane         20         20           Acetone         1,800         1,000           Acrylonitrile         100         100           1,1-Dichloroethene         20         20           Methylene Chloride         35         100           Carbon Disulfide         310         100           trans-1,2-Dichloroethene         20         20           1,1-Dichloroethane         84         20           2-Butanone (MEK)         17         100           cis-1,2-Dichloroethene         19         20           Chloroform         24,000         20           1,2-Dichloroethane         20         20           1,1,1-Trichloroethane         20         20           Benzene         20         20	μg/m³         μg/m³         μg/m³         μg/m³           Chloromethane         20         20         10           Vinyl Chloride         20         20         10           Bromomethane         20         20         10           Chloroethane         20         20         10           Acetone         1,800         1,000         15           Acrylonitrile         100         100         14           1,1-Dichloroethene         20         20         10           Methylene Chloride         35         100         10           Carbon Disulfide         310         100         24           trans-1,2-Dichloroethene         20         20         10           1,1-Dichloroethane         84         20         10           2-Butanone (MEK)         17         100         10           cis-1,2-Dichloroethene         19         20         10           Chloroform         24,000         20         12           1,2-Dichloroethane         20         20         10           1,1,1-Trichloroethane         20         20         10           Benzene         20         20         10	μg/m³         μg/m³         μg/m³         ppbV           Chloromethane         20         20         10         9.9           Vinyl Chloride         20         20         10         8.0           Bromomethane         20         20         10         5.2           Chloroethane         20         20         10         7.7           Acetone         1,800         1,000         15         760           Acrylonitrile         100         100         14         47           1,1-Dichloroethene         20         20         10         5.1           Methylene Chloride         35         100         10         10           Carbon Disulfide         310         100         24         100           trans-1,2-Dichloroethene         20         20         10         5.1           1,1-Dichloroethane         84         20         10         5.8           cis-1,2-Dichloroethene         17         100         10         5.8           cis-1,2-Dichloroethene         19         20         10         4.9           Chloroform         24,000         20         12         4,800           1,1,1-Trichloroethane <td>μg/m³         μg/m³         μg/m³         ppbV         ppbV           Chloromethane         20         20         10         9.9         9.9           Vinyl Chloride         20         20         10         8.0         8.0           Bromomethane         20         20         10         5.2         5.2           Chloroethane         20         20         10         7.7         7.7           Acetone         1,800         1,000         15         760         430           Acrylonitrile         100         10         14         47         47           1,1-Dichloroethene         20         20         10         5.1         5.1           Methylene Chloride         35         100         10         10         29           Carbon Disulfide         310         100         24         100         33           trans-1,2-Dichloroethene         20         20         10         5.1         5.1           1,1-Dichloroethane         84         20         10         5.8         35           cis-1,2-Dichloroethene         19         20         10         4.9         5.1           Chloroform         24,</td> <td>μg/m³         μg/m³         μg/m³         ppbV         ppbV         ppbV           Chloromethane         20         20         10         9.9         9.9         4.9           Vinyl Chloride         20         20         10         8.0         8.0         4.0           Bromomethane         20         20         10         5.2         5.2         2.6           Chloroethane         20         20         10         7.7         7.7         3.9           Acetone         1,800         1,000         15         760         430         6.3           Acrylonitrile         100         10         14         47         47         6.6           1,1-Dichloroethene         20         20         10         5.1         5.1         2.6           Methylene Chloride         35         100         10         10         29         2.9           Carbon Disulfide         310         100         24         100         33         7.9           trans-1,2-Dichloroethene         84         20         10         5.1         5.1         2.6           1,1-Dichloroethane         84         20         10         5.8         <td< td=""></td<></td>	μg/m³         μg/m³         μg/m³         ppbV         ppbV           Chloromethane         20         20         10         9.9         9.9           Vinyl Chloride         20         20         10         8.0         8.0           Bromomethane         20         20         10         5.2         5.2           Chloroethane         20         20         10         7.7         7.7           Acetone         1,800         1,000         15         760         430           Acrylonitrile         100         10         14         47         47           1,1-Dichloroethene         20         20         10         5.1         5.1           Methylene Chloride         35         100         10         10         29           Carbon Disulfide         310         100         24         100         33           trans-1,2-Dichloroethene         20         20         10         5.1         5.1           1,1-Dichloroethane         84         20         10         5.8         35           cis-1,2-Dichloroethene         19         20         10         4.9         5.1           Chloroform         24,	μg/m³         μg/m³         μg/m³         ppbV         ppbV         ppbV           Chloromethane         20         20         10         9.9         9.9         4.9           Vinyl Chloride         20         20         10         8.0         8.0         4.0           Bromomethane         20         20         10         5.2         5.2         2.6           Chloroethane         20         20         10         7.7         7.7         3.9           Acetone         1,800         1,000         15         760         430         6.3           Acrylonitrile         100         10         14         47         47         6.6           1,1-Dichloroethene         20         20         10         5.1         5.1         2.6           Methylene Chloride         35         100         10         10         29         2.9           Carbon Disulfide         310         100         24         100         33         7.9           trans-1,2-Dichloroethene         84         20         10         5.1         5.1         2.6           1,1-Dichloroethane         84         20         10         5.8 <td< td=""></td<>

U = Compound was analyzed for, but not detected above the laboratory detection limit.

Verified By:

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

B = Analyte was found in the method blank.

# **RESULTS OF ANALYSIS**

Page 2 of 3

**Client:** 

CH2M Hill

Client Sample ID: WAT-SG-7-042408

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-013

Test Code:

EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/24/08 Date Received: 4/28/08

Instrument ID: Analyst:

Wida Ang

Date Analyzed: 5/2/08

Sampling Media:

6.0 L Summa Canister

Volume(s) Analyzed: 0.0080 Liter(s)

Test Notes:

Container ID:

AC01276

Initial Pressure (psig):

-3.5

Final Pressure (psig):

3.5

CAS#	Compound	Result μg/m³	MRL μg/m³	MDL μg/m³	Result ppbV	MRL ppbV	MDL ppbV	Data Qualifier
78-87-5	1,2-Dichloropropane	52	20	10	11	4.4	2.2	- Quantiter
75-27-4	Bromodichloromethane	20	20	10	3.0	3.0	1.5	$\mathbf{U}$
79-01-6	Trichloroethene	54	20	10	10	3.8	1.9	
10061-01-5	cis-1,3-Dichloropropene	100	100	11	22	22	2.3	$\mathbf{U}$
108-10-1	4-Methyl-2-pentanone	12,000	100	11	2,900	25	2.8	
10061-02-6	trans-1,3-Dichloropropene	100	100	13	22	22	2.8	U
79-00-5	1,1,2-Trichloroethane	20	20	10	3.7	3.7	1.9	$\mathbf{U}$
108-88-3	Toluene	72	100	10	19	27	2.7	J
124-48-1	Dibromochloromethane	20	20	14	2.4	2.4	1.6	$\mathbf{U}$
127-18-4	Tetrachloroethene	26	20	10	3.9	3.0	1.5	
108-90-7	Chlorobenzene	20	20	10	4.4	4.4	2.3	U
100-41-4	Ethylbenzene	870	100	13	200	23	2.9	
179601-23-1	m,p-Xylenes	4,600	100	26	1,100	23	6.1	
75-25-2	Bromoform	100	100	15	9.9	9.9	1.5	$\mathbf{U}$
100-42-5	Styrene	100	100	15	24	24	3.6	$\mathbf{U}$
95-47-6	o-Xylene	950	100	13	220	23	3.0	NATION OF STATE AND A SHARE A
79-34-5	1,1,2,2-Tetrachloroethane	20	20	13	3.0	3.0	1.9	U

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

# RESULTS OF ANALYSIS

Page 3 of 3

**Client:** 

**CH2M Hill** 

Client Sample ID: WAT-SG-7-042408

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-013

**Tentatively Identified Compounds** 

Test Code: Instrument ID: EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/24/08 Date Received: 4/28/08

Date Analyzed: 5/2/08

Analyst: Sampling Media: Wida Ang 6.0 L Summa Canister

Volume(s) Analyzed: 0.0080 Liter(s)

Test Notes:

Container ID:

AC01276

Initial Pressure (psig):

-3.5

Final Pressure (psig):

3.5

Canister Dilution Factor: 1.63

GC/MS Retention Time **Compound Identification** 

Concentration

Data

Epichlorohydrin

 $\mu g/m^3$ NF Qualifier

NF = Compound was searched for, but not found.

#### **RESULTS OF ANALYSIS**

Page 1 of 3

**Client:** 

**CH2M Hill** 

Client Sample ID: WAT-SG-7a-042408

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-014

Test Code:

EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/24/08 Date Received: 4/28/08

Instrument ID: Analyst:

Wida Ang

Date Analyzed: 5/5/08

6.0 L Summa Canister

Volume(s) Analyzed: 0.0010 Liter(s)

Test Notes:

Container ID:

Sampling Media:

AC01211

Initial Pressure (psig):

0.2

Final Pressure (psig):

3.5

Canister Dilution Factor: 1.22

CAS#	Compound	Result	MRL	MDL	Result	MRL	MDL	Data
		$\mu g/m^3$	$\mu g/m^3$	$\mu g/m^3$	ppbV	ppbV	ppbV	Qualifier
74-87-3	Chloromethane	120	120	61	59	59	30	U
75-01-4	Vinyl Chloride	120	120	61	48	48	24	$\mathbf{U}$
74-83-9	Bromomethane	120	120	61	31	31	16	$\mathbf{U}$
75-00-3	Chloroethane	120	120	61	46	46	23	$\mathbf{U}$
67-64-1	Acetone	1,600	6,100	89	690	2,600	38	J, B
107-13-1	Acrylonitrile	610	610	85	280	280	39	U
75-35-4	1,1-Dichloroethene	120	120	61	31	31	15	$\mathbf{U}$
75-09-2	Methylene Chloride	240	610	61	69	180	18	J
75-15-0	Carbon Disulfide	3,800	610	150	1,200	200	47	
156-60-5	trans-1,2-Dichloroethene	120	120	61	31	31	15	$\mathbf{U}$
75-34-3	1,1-Dichloroethane	120	120	61	30	30	15	U
78-93-3	2-Butanone (MEK)	610	610	61	210	210	21	U
156-59-2	cis-1,2-Dichloroethene	120	120	61	31	31	15	U
67-66-3	Chloroform	1,300	120	72	270	25	15	
107-06-2	1,2-Dichloroethane	120	120	61	30	30	15	$\mathbf{U}$
71-55-6	1,1,1-Trichloroethane	120	120	61	22	22	11	U
71-43-2	Benzene	120	120	61	38	38	19	U
56-23-5	Carbon Tetrachloride	120	. 120	61	19	19	9.7	$\mathbf{U}$

U = Compound was analyzed for, but not detected above the laboratory detection limit.

Verified By: Res

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

B = Analyte was found in the method blank.

### **RESULTS OF ANALYSIS**

Page 2 of 3

**Client:** CH2M Hill

Client Sample ID: WAT-SG-7a-042408 CAS Project ID: P0801215

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI CAS Sample ID: P0801215-014

Test Code: EPA TO-15

Date Collected: 4/24/08 Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16 Date Received: 4/28/08 Analyst: Wida Ang Date Analyzed: 5/5/08

Sampling Media: 6.0 L Summa Canister Volume(s) Analyzed: 0.0010 Liter(s)

Test Notes:

Container ID: AC01211

> Initial Pressure (psig): 0.2 Final Pressure (psig): 3.5

> > Canister Dilution Factor: 1.22

CAS#	Compound	Result	MRL	MDL	Result	MRL	MDL	Data
		μg/m³	μg/m³	μg/m³	ppbV	ppbV	ppbV	Qualifier
78-87-5	1,2-Dichloropropane	120	120	61	26	26	13	$\mathbf{U}$
75-27-4	Bromodichloromethane	120	120	61	18	18	9.1	U
79-01-6	Trichloroethene	120	120	61	23	23	11	$\mathbf{U}$
10061-01-5	cis-1,3-Dichloropropene	610	610	63	130	130	14	$\mathbf{U}$
108-10-1	4-Methyl-2-pentanone	86,000	610	68	21,000	150	17	
10061-02-6	trans-1,3-Dichloropropene	610	610	77	130	130	17	U
79-00-5	1,1,2-Trichloroethane	120	120	61	22	22	11	$\mathbf{U}$
108-88-3	Toluene	380	610	61	100	160	16	J
124-48-1	Dibromochloromethane	120	120	83	14	14	9.7	$\mathbf{U}$
127-18-4	Tetrachloroethene	120	120	61	18	18	9.0	$\mathbf{U}$
108-90-7	Chlorobenzene	120	120	62	27	27	14	U
100-41-4	Ethylbenzene	430	610	76	. 99	140	17	J
179601-23-1	m,p-Xylenes	8,600	610	160	2,000	140	37	
75-25-2	Bromoform	610	610	93	59	59	9.0	$\mathbf{U}$
100-42-5	Styrene	610	610	93	140	140	22	U
95-47-6	o-Xylene	1,700	610	77	390	140	18	
79-34-5	1,1,2,2-Tetrachloroethane	120	120	78	18	18	11	$\mathbf{U}$

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

#### **RESULTS OF ANALYSIS**

Page 3 of 3

Client:

**CH2M Hill** 

Client Sample ID: WAT-SG-7a-042408

CAS Project ID: P0801215

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Sample ID: P0801215-014

**Tentatively Identified Compounds** 

EPA TO-15

Date Collected: 4/24/08

Test Code: Instrument ID:

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Received: 4/28/08

Analyst:

Wida Ang

Date Analyzed: 5/5/08

Sampling Media:

6.0 L Summa Canister

Volume(s) Analyzed: 0.0010 Liter(s)

Test Notes:

T

Container ID:

AC01211

Initial Pressure (psig):

0.2

Final Pressure (psig):

3.5

GC/MS	Compound Identification	Concentration	Data
Retention Time		$\mu \mathrm{g}/\mathrm{m}^3$	Qualifier
17.38	n-Butyl Ether	30,000	
20.17	2,2'-Oxybispentane ( $C_{10}H_{22}O$ )	30,000	
20.24	Unidentified Oxygenated Compound	10,000	
20.47	Unidentified Oxygenated Compound	10,000	
21.03	1,1'-Oxybispentane; Amyl Ether (C <sub>10</sub> H <sub>22</sub> O)	10,000	
	Epichlorohydrin	NF	

T = Analyte is a tentatively identified compound, result is estimated.

NF = Compound was searched for, but not found.

#### RESULTS OF ANALYSIS

Page 1 of 3

Client:

CH2M Hill

Client Sample ID: WAT-SG-8-042408

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-015

Test Code:

EPA TO-15

Instrument ID:

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

6.0 L Summa Canister

Date Received: 4/28/08

Date Collected: 4/24/08 Date Analyzed: 5/5/08

Analyst: Sampling Media: Wida Ang

Volume(s) Analyzed:

0.10 Liter(s)

Test Notes:

Container ID:

AC01235

Initial Pressure (psig):

-3.9

Final Pressure (psig):

3.6

CAC#	Comment	Dogule	MDI	MDI	Result	MRL	MDL	Data
CAS#	Compound	Result	MRL	MDL				Data
		μg/m³	μg/m³	μg/m³	ppbV	ppbV	ppbV	Qualifier
74-87-3	Chloromethane	2.1	1.7	0.85	1.0	0.82	0.41	
75-01-4	Vinyl Chloride	1.7	1.7	0.85	0.66	0.66	0.33	U
74-83-9	Bromomethane	1.7	1.7	0.85	0.44	0.44	0.22	$\mathbf{U}$
75-00-3	Chloroethane	1.7	1.7	0.85	0.64	0.64	0.32	$\mathbf{U}$
67-64-1	Acetone	230	85	1.2	98	36	0.52	В
107-13-1	Acrylonitrile	3.3	8.5	1.2	1.5	3.9	0.55	J
75-35-4	1,1-Dichloroethene	1.7	1.7	0.85	0.43	0.43	0.21	U
75-09-2	Methylene Chloride	2.3	8.5	0.85	0.66	2.4	0.24	J
75-15-0	Carbon Disulfide	43	8.5	2.0	14	2.7	0.65	
156-60-5	trans-1,2-Dichloroethene	1.7	1.7	0.85	0.43	0.43	0.21	$\mathbf{U}$
75-34-3	1,1-Dichloroethane	1.7	1.7	0.85	0.42	0.42	0.21	U
78-93-3	2-Butanone (MEK)	32	8.5	0.85	11	2.9	0.29	
156-59-2	cis-1,2-Dichloroethene	1.7	1.7	0.85	0.43	0.43	0.21	$\mathbf{U}$
67-66-3	Chloroform	170	1.7	1.0	34	0.35	0.20	
107-06-2	1,2-Dichloroethane	1.7	1.7	0.85	0.42	0.42	0.21	U
71-55-6	1,1,1-Trichloroethane	1.7	1.7	0.85	0.31	0.31	0.15	U
71-43-2	Benzene	5.0	1.7	0.85	1.6	0.53	0.26	
56-23-5	Carbon Tetrachloride	1.7	1.7	0.85	0.27	0.27	0.13	U

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

B = Analyte was found in the method blank.

# **RESULTS OF ANALYSIS**

Page 2 of 3

Client:

**CH2M Hill** 

Client Sample ID: WAT-SG-8-042408

6.0 L Summa Canister

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-015

Test Code:

EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/24/08 Date Received: 4/28/08

Date Analyzed: 5/5/08

Instrument ID: Analyst:

Wida Ang

Volume(s) Analyzed:

0.10 Liter(s)

Sampling Media: Test Notes:

Container ID:

AC01235

Initial Pressure (psig):

-3.9

Final Pressure (psig):

3.6

Canister Dilution Factor: 1.69

CAS#	Compound	Result μg/m³	MRL μg/m³	MDL μg/m³	Result ppbV	MRL ppbV	MDL ppbV	Data Qualifier
78-87-5	1,2-Dichloropropane	2.1	1.7	0.85	0.46	0.37	0.18	
75-27-4	Bromodichloromethane	1.7	1.7	0.85	0.25	0.25	0.13	$\mathbf{U}$
79-01-6	Trichloroethene	1.7	1.7	0.85	0.32	0.31	0.16	
10061-01-5	cis-1,3-Dichloropropene	8.5	8.5	0.88	1.9	1.9	0.19	$\mathbf{U}$
108-10-1	4-Methyl-2-pentanone	380	8.5	0.95	92	2.1	0.23	E
10061-02-6	trans-1,3-Dichloropropene	8.5	8.5	1.1	1.9	1.9	0.23	U
79-00-5	1,1,2-Trichloroethane	1.7	1.7	0.85	0.31	0.31	0.15	$\mathbf{U}$
108-88-3	Toluene	130	8.5	0.85	33	2.2	0.22	
124-48-1	Dibromochloromethane	1.7	1.7	1.1	0.20	0.20	0.13	$\mathbf{U}$
127-18-4	Tetrachloroethene	3.0	1.7	0.85	0.44	0.25	0.12	
108-90-7	Chlorobenzene	2.3	1.7	0.86	0.51	0.37	0.19	
100-41-4	Ethylbenzene	1.6	8.5	1.0	0.38	1.9	0.24	J
179601-23-1	m,p-Xylenes	15	8.5	2.2	3.4	1.9	0.51	
75-25-2	Bromoform	8.5	8.5	1.3	0.82	0.82	0.12	${f U}$
100-42-5	Styrene	8.5	8.5	1.3	2.0	2.0	0.30	$\mathbf{U}$
95-47-6	o-Xylene	4.3	8.5	1.1	0.98	1.9	0.25	J
79-34-5	1,1,2,2-Tetrachloroethane	1.7	1.7	1.1	0.25	0.25	0.16	U

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

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#### RESULTS OF ANALYSIS

Page 3 of 3

**Client:** 

**CH2M Hill** 

Client Sample ID: WAT-SG-8-042408

CAS Project ID: P0801215

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Sample ID: P0801215-015

**Tentatively Identified Compounds** 

Test Code:

EPA TO-15

Date Collected: 4/24/08

Instrument ID:

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Received: 4/28/08

Analyst:

Date Analyzed: 5/5/08

Sampling Media:

Wida Ang 6.0 L Summa Canister

Volume(s) Analyzed:

0.10 Liter(s)

Test Notes:

 $\mathbf{T}$ 

Container ID:

AC01235

Initial Pressure (psig):

-3.9

Final Pressure (psig):

3.6

GC/MS	Compound Identification	Concentration	Data
Retention Time		$\mu \mathrm{g}/\mathrm{m}^3$	Qualifier
19.52	Isooctanol	400	
20.18	3-Methylpentane	300	
20.24	Unidentified Oxygenated Compound	600	
20.72	C <sub>12</sub> H <sub>26</sub> Branched Alkane	200	
21.01	C <sub>11</sub> H <sub>24</sub> Branched Alkane	100	
	Epichlorohydrin	NF	

T = Analyte is a tentatively identified compound, result is estimated.

NF = Compound was searched for, but not found.

### **RESULTS OF ANALYSIS**

Page 1 of 3

**Client:** 

CH2M Hill

Client Sample ID: WAT-SG-9-042408

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-016

Test Code:

EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/24/08

Date Received: 4/28/08

Instrument ID: Analyst:

Wida Ang

Date Analyzed: 5/5/08

Sampling Media:

6.0 L Summa Canister

Volume(s) Analyzed:

0.25 Liter(s)

Test Notes:

Container ID:

AC01020

Initial Pressure (psig):

0.3

Final Pressure (psig):

3.6

CAS#	Compound	Result	MRL	MDL	Result	MRL	MDL	Data
САЗ П	Compound	μg/m³	μg/m³	μg/m <sup>3</sup>	ppbV	ppbV	ppbV	Qualifier
74-87-3	Chloromethane	0.51	0.49	0.24	0.25	0.24	0.12	
75-01-4	Vinyl Chloride	0.49	0.49	0.24	0.19	0.19	0.095	U
74-83-9	Bromomethane	0.49	0.49	0.24	0.13	0.13	0.063	$\mathbf{U}$
75-00-3	Chloroethane	0.49	0.49	0.24	0.19	0.19	0.093	$\mathbf{U}$
67-64-1	Acetone	13	24	0.36	5.6	10	0.15	J, B
107-13-1	Acrylonitrile	2.4	2.4	0.34	1.1	1.1	0.16	U
75-35-4	1,1-Dichloroethene	0.49	0.49	0.24	0.12	0.12	0.062	$\mathbf{U}$
75-09-2	Methylene Chloride	0.31	2.4	0.24	0.089	0.70	0.070	J
75-15-0	Carbon Disulfide	3.8	2.4	0.59	1.2	0.78	0.19	
156-60-5	trans-1,2-Dichloroethene	0.49	0.49	0.24	0.12	0.12	0.062	$\mathbf{U}$
75-34-3	1,1-Dichloroethane	0.49	0.49	0.24	0.12	0.12	0.060	U
78-93-3	2-Butanone (MEK)	1.3	2.4	0.24	0.44	0.83	0.083	J
156-59-2	cis-1,2-Dichloroethene	0.49	0.49	0.24	0.12	0.12	0.062	$\mathbf{U}$
67-66-3	Chloroform	270	0.49	0.29	56	0.10	0.059	
107-06-2	1,2-Dichloroethane	0.49	0.49	0.24	0.12	0.12	0.060	$\mathbf{U}$
71-55-6	1,1,1-Trichloroethane	0.49	0.49	0.24	0.089	0.089	0.045	U
71-43-2	Benzene	0.66	0.49	0.24	0.21	0.15	0.076	
56-23-5	Carbon Tetrachloride	0.49	0.49	0.24	0.078	0.078	0.039	U

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

B = Analyte was found in the method blank.

# RESULTS OF ANALYSIS

Page 2 of 3

Client: CH2M Hill

Client Sample ID: WAT-SG-9-042408 CAS Project ID: P0801215

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI CAS Sample ID: P0801215-016

Test Code: EPA TO-15 Date Collected: 4/24/08

Instrument ID: Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16 Date Received: 4/28/08
Analyst: Wida Ang Date Analyzed: 5/5/08

Sampling Media: 6.0 L Summa Canister Volume(s) Analyzed: 0.25 Liter(s)

Test Notes:
Container ID: AC01020

Initial Pressure (psig): 0.3 Final Pressure (psig): 3.6

Canister Dilution Factor: 1.22

CAS#	Compound	Result μg/m³	MRL μg/m³	MDL μg/m³	Result ppbV	MRL ppbV	MDL ppbV	Data Qualifier
78-87-5	1,2-Dichloropropane	15	0.49	0.24	3.3	0.11	$\frac{pp0.7}{0.053}$	Quanner
75-27-4	Bromodichloromethane	0.49	0.49	0.24	0.073	0.073	0.036	$\mathbf{U}$
79-01-6	Trichloroethene	520	0.49	0.24	96	0.091	0.045	
10061-01-5	cis-1,3-Dichloropropene	2.4	2.4	0.25	0.54	0.54	0.056	U
108-10-1	4-Methyl-2-pentanone	<b>11</b>	2.4	0.27	2.6	0.60	0.067	$\mathbf{E}$
10061-02-6	trans-1,3-Dichloropropene	2.4	2.4	0.31	0.54	0.54	0.068	U
79-00-5	1,1,2-Trichloroethane	0.49	0.49	0.24	0.089	0.089	0.045	U
108-88-3	Toluene	1.6	2.4	0.24	0.42	0.65	0.065	J
124-48-1	Dibromochloromethane	0.49	0.49	0.33	0.057	0.057	0.039	U
127-18-4	Tetrachloroethene	6.4	0.49	0.24	0.94	0.072	0.036	
108-90-7	Chlorobenzene	0.49	0.49	0.25	0.11	0.11	0.054	U
100-41-4	Ethylbenzene	2.4	2.4	0.30	0.56	0.56	0.070	U
179601-23-1	m,p-Xylenes	1.3	2.4	0.63	0.29	0.56	0.15	J
75-25-2	Bromoform	2.4	2.4	0.37	0.24	0.24	0.036	U
100-42-5	Styrene	2.4	2.4	0.37	0.57	0.57	0.087	$\mathbf{U}$
95-47-6	o-Xylene	0.54	2.4	0.31	0.12	0.56	0.071	J
79-34-5	1,1,2,2-Tetrachloroethane	0.49	0.49	0.31	0.071	0.071	0.045	U

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

# **RESULTS OF ANALYSIS**

Page 3 of 3

**Tentatively Identified Compounds** 

Client:

CH2M Hill

Client Sample ID: WAT-SG-9-042408

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-016

Instrument ID:

Test Code:

EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: 4/24/08

Date Received: 4/28/08

Analyst:

Wida Ang

Date Analyzed: 5/5/08 Volume(s) Analyzed:

0.25 Liter(s)

Sampling Media:

6.0 L Summa Canister

Test Notes: Container ID: Т

AC01020

Initial Pressure (psig):

0.3

Final Pressure (psig):

3.6

GC/MS	Compound Identification	Concentration	Data
Retention Time		$\mu \mathrm{g}/\mathrm{m}^3$	Qualifier
4.46	Propane	10	
4.87	Acetaldehyde + Isobutane	30	
16.27	C6HF13 (Tridecafluorohexane)	70	
20.23	Unidentified Oxygenated Compound	30	
	Epichlorohydrin	NF	

T = Analyte is a tentatively identified compound, result is estimated.

NF = Compound was searched for, but not found.

### RESULTS OF ANALYSIS Page 1 of 3

**Client:** 

**CH2M Hill** 

Client Sample ID: Method Blank

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215 CAS Sample ID: P080502-MB

Test Code:

EPA TO-15

Instrument ID:

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Wida Ang

Sampling Media:

6.0 L Summa Canister

Date Received: NA

Volume(s) Analyzed:

Date Analyzed: 5/2/08

Date Collected: NA

1.00 Liter(s)

Test Notes:

Analyst:

Canister Dilution Factor: 1.00

CAS#	Compound	Result μg/m³	MRL μg/m³	$MDL$ $\mu g/m^3$	Result ppbV	MRL ppbV	MDL ppbV	Data Qualifier
74-87-3	Chloromethane	0.10	0.10	0.050	0.048	0.048	0.024	U
75-01-4	Vinyl Chloride	0.10	0.10	0.050	0.039	0.039	0.020	$\mathbf{U}$
74-83-9	Bromomethane	0.10	0.10	0.050	0.026	0.026	0.013	$\mathbf{U}$
75-00-3	Chloroethane	0.10	0.10	0.050	0.038	0.038	0.019	$\mathbf{U}$
67-64-1	Acetone	0.78	5.0	0.073	0.33	2.1	0.031	J
107-13-1	Acrylonitrile	0.50	0.50	0.070	0.23	0.23	0.032	U
75-35-4	1,1-Dichloroethene	0.10	0.10	0.050	0.025	0.025	0.013	$\mathbf{U}^{r}$
75-09-2	Methylene Chloride	0.50	0.50	0.050	0.14	0.14	0.014	$\mathbf{U}$
75-15-0	Carbon Disulfide	0.50	0.50	0.12	0.16	0.16	0.039	$\mathbf{U}$
156-60-5	trans-1,2-Dichloroethene	0.10	0.10	0.050	0.025	0.025	0.013	${f U}$
75-34-3	1,1-Dichloroethane	0.10	0.10	0.050	0.025	0.025	0.012	U
78-93-3	2-Butanone (MEK)	0.50	0.50	0.050	0.17	0.17	0.017	$\mathbf{U}$
156-59-2	cis-1,2-Dichloroethene	0.10	0.10	0.050	0.025	0.025	0.013	$\mathbf{U}$
67-66-3	Chloroform	0.10	0.10	0.059	0.020	0.020	0.012	$\mathbf{U}$
107-06-2	1,2-Dichloroethane	0.10	0.10	0.050	0.025	0.025	0.012	$\mathbf{U}$
71-55-6	1,1,1-Trichloroethane	0.10	0.10	0.050	0.018	0.018	0.0092	U
71-43-2	Benzene	0.10	0.10	0.050	0.031	0.031	0.016	$\mathbf{U}$
56-23-5	Carbon Tetrachloride	0.10	0.10	0.050	0.016	0.016	0.0080	U

ND = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method. J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

Verified By:\_\_\_

# **RESULTS OF ANALYSIS** Page 2 of 3

Client:

**CH2M Hill** 

Client Sample ID: Method Blank

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

Date Collected: NA

CAS Sample ID: P080502-MB

Test Code:

EPA TO-15

Instrument ID:

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Wida Ang

Analyst: Sampling Media:

6.0 L Summa Canister

Date Received: NA Date Analyzed: 5/2/08

Volume(s) Analyzed:

1.00 Liter(s)

Test Notes:

Canister Dilution Factor: 1.00

CAS#	Compound	Result μg/m³	MRL μg/m³	$MDL \ \mu g/m^3$	Result ppbV	MRL ppbV	MDL ppbV	Data Qualifier
78-87-5	1,2-Dichloropropane	0.10	0.10	0.050	0.022	0.022	0.011	U
75-27-4	Bromodichloromethane	0.10	0.10	0.050	0.015	0.015	0.0075	U
79-01-6	Trichloroethene	0.10	0.10	0.050	0.019	0.019	0.0093	$\mathbf{U}$
10061-01-5	cis-1,3-Dichloropropene	0.50	0.50	0.052	0.11	0.11	0.011	$\mathbf{U}$
108-10-1	4-Methyl-2-pentanone	0.50	0.50	0.056	0.12	0.12	0.014	$\mathbf{U}$
10061-02-6	trans-1,3-Dichloropropene	0.50	0.50	0.063	0.11	0.11	0.014	U
79-00-5	1,1,2-Trichloroethane	0.10	0.10	0.050	0.018	0.018	0.0092	U
108-88-3	Toluene	0.50	0.50	0.050	0.13	0.13	0.013	U
124-48-1	Dibromochloromethane	0.10	0.10	0.068	0.012	0.012	0.0080	U
127-18-4	Tetrachloroethene	0.10	0.10	0.050	0.015	0.015	0.0074	$\mathbf{U}$
108-90-7	Chlorobenzene	0.10	0.10	0.051	0.022	0.022	0.011	U
100-41-4	Ethylbenzene	0.50	0.50	0.062	0.12	0.12	0.014	U
179601-23-1	m,p-Xylenes	0.50	0.50	0.13	0.12	0.12	0.030	$\mathbf{U}$
75-25-2	Bromoform	0.50	0.50	0.076	0.048	0.048	0.0074	U
100-42-5	Styrene	0.50	0.50	0.076	0.12	0.12	0.018	$\mathbf{U}$
95-47-6	o-Xylene	0.50	0.50	0.063	0.12	0.12	0.015	U
79-34-5	1,1,2,2-Tetrachloroethane	0.10	0.10	0.064	0.015	0.015	0.0093	U

ND = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

Verified By:\_

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# **RESULTS OF ANALYSIS**

Page 3 of 3

Client:

**CH2M Hill** 

Client Sample ID: Method Blank

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P080502-MB

**Tentatively Identified Compounds** 

Test Code:

EPA TO-15

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Collected: NA

Date Received: NA

Instrument ID: Analyst:

Wida Ang

Date Analyzed: 5/2/08

1.00 Liter(s)

Sampling Media:

6.0 L Summa Canister

Volume(s) Analyzed:

Test Notes:

Canister Dilution Factor: 1.00

GC/MS

**Compound Identification** 

Concentration

Data

59

Retention Time

Epichlorohydrin

 $\mu g/m^3$ NF Qualifier

NF = Compound was searched for, but not found.

Verified By: TO15SCAN.XLT - NL - PageNo.:

### **RESULTS OF ANALYSIS** Page 1 of 3

**Client:** 

CH2M Hill

Client Sample ID: Method Blank

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

Date Collected: NA

Date Received: NA

CAS Sample ID: P080505-MB

Test Code:

EPA TO-15

Instrument ID:

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Analyst:

Wida Ang

Sampling Media:

6.0 L Summa Canister

Date Analyzed: 5/5/08 Volume(s) Analyzed:

1.00 Liter(s)

Test Notes:

Canister Dilution Factor: 1.00

CAS#	Compound	Result μg/m³	MRL μg/m³	MDL μg/m³	Result ppbV	MRL ppbV	MDL ppbV	Data Qualifier
74-87-3	Chloromethane	0.10	0.10	0.050	0.048	0.048	0.024	U
75-01-4	Vinyl Chloride	0.10	0.10	0.050	0.039	0.039	0.020	$\mathbf{U}$
74-83-9	Bromomethane	0.10	0.10	0.050	0.026	0.026	0.013	$\mathbf{U}$
75-00-3	Chloroethane	0.10	0.10	0.050	0.038	0.038	0.019	$\mathbf{U}$
67-64-1	Acetone	0.45	5.0	0.073	0.19	2.1	0.031	J
107-13-1	Acrylonitrile	0.50	0.50	0.070	0.23	0.23	0.032	U
75-35-4	1,1-Dichloroethene	0.10	0.10	0.050	0.025	0.025	0.013	U
75-09-2	Methylene Chloride	0.50	0.50	0.050	0.14	0.14	0.014	$\mathbf{U}$
75-15-0	Carbon Disulfide	0.50	0.50	0.12	0.16	0.16	0.039	$\mathbf{U}$
156-60-5	trans-1,2-Dichloroethene	0.10	0.10	0.050	0.025	0.025	0.013	$\mathbf{U}$
75-34-3	1,1-Dichloroethane	0.10	0.10	0.050	0.025	0.025	0.012	U
78-93-3	2-Butanone (MEK)	0.50	0.50	0.050	0.17	0.17	0.017	$\mathbf{U}$
156-59-2	cis-1,2-Dichloroethene	0.10	0.10	0.050	0.025	0.025	0.013	$\mathbf{U}$
67-66-3	Chloroform	0.10	0.10	0.059	0.020	0.020	0.012	$\mathbf{U}$
107-06-2	1,2-Dichloroethane	0.10	0.10	0.050	0.025	0.025	0.012	$\mathbf{U}$
71-55-6	1,1,1-Trichloroethane	0.10	0.10	0.050	0.018	0.018	0.0092	U
71-43-2	Benzene	0.10	0.10	0.050	0.031	0.031	0.016	$\mathbf{U}$
56-23-5	Carbon Tetrachloride	0.10	0.10	0.050	0.016	0.016	0.0080	U

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method. J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

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### RESULTS OF ANALYSIS Page 2 of 3

**Client:** 

**CH2M Hill** 

Client Sample ID: Method Blank

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215 CAS Sample ID: P080505-MB

Date Collected: NA

Date Received: NA

Test Code:

EPA TO-15

Instrument ID:

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Analyst:

Wida Ang

Sampling Media:

6.0 L Summa Canister

Date Analyzed: 5/5/08 Volume(s) Analyzed:

1.00 Liter(s)

Test Notes:

Canister Dilution Factor: 1.00

CAS#	Compound	Result μg/m³	MRL μg/m³	MDL μg/m³	Result ppbV	MRL ppbV	MDL ppbV	Data Qualifier
78-87-5	1,2-Dichloropropane	0.10	0.10	0.050	0.022	0.022	0.011	U
75-27-4	Bromodichloromethane	0.10	0.10	0.050	0.015	0.015	0.0075	$\mathbf{U}$
79-01-6	Trichloroethene	0.10	0.10	0.050	0.019	0.019	0.0093	$\mathbf{U}$
10061-01-5	cis-1,3-Dichloropropene	0.50	0.50	0.052	0.11	0.11	0.011	$\mathbf{U}$
108-10-1	4-Methyl-2-pentanone	0.50	0.50	0.056	0.12	0.12	0.014	$\mathbf{U}$
10061-02-6	trans-1,3-Dichloropropene	0.50	0.50	0.063	0.11	0.11	0.014	U
79-00-5	1,1,2-Trichloroethane	0.10	0.10	0.050	0.018	0.018	0.0092	$\mathbf{U}$
108-88-3	Toluene	0.50	0.50	0.050	0.13	0.13	0.013	U
124-48-1	Dibromochloromethane	0.10	0.10	0.068	0.012	0.012	0.0080	$\mathbf{U}$
127-18-4	Tetrachloroethene	0.10	0.10	0.050	0.015	0.015	0.0074	U
108-90-7	Chlorobenzene	0.10	0.10	0.051	0.022	0.022	0.011	U
100-41-4	Ethylbenzene	0.50	0.50	0.062	0.12	0.12	0.014	U
179601-23-1	m,p-Xylenes	0.50	0.50	0.13	0.12	0.12	0.030	$\mathbf{U}$
75-25-2	Bromoform	0.50	0.50	0.076	0.048	0.048	0.0074	$\mathbf{U}$
100-42-5	Styrene	0.50	0.50	0.076	0.12	0.12	0.018	U
95-47-6	o-Xylene	0.50	0.50	0.063	0.12	0.12	0.015	U
79-34-5	1,1,2,2-Tetrachloroethane	0.10	0.10	0.064	0.015	0.015	0.0093	U

U = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

Verified By:

### **RESULTS OF ANALYSIS**

Page 3 of 3

**Client:** 

**CH2M Hill** 

Client Sample ID: Method Blank

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P080505-MB

**Tentatively Identified Compounds** 

Test Code:

EPA TO-15

Date Collected: NA

Instrument ID:

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Received: NA

Wida Ang

Date Analyzed: 5/5/08

Sampling Media:

6.0 L Summa Canister

Volume(s) Analyzed:

1.00 Liter(s)

Test Notes:

Analyst:

Canister Dilution Factor: 1.00

GC/MS

**Compound Identification** 

Concentration

Data

Retention Time

Epichlorohydrin

 $\mu g/m^3$ 

Qualifier

NF = Compound was searched for, but not found.

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### SURROGATE SPIKE RECOVERY RESULTS

Page 1 of 1

**Client:** 

**CH2M Hill** 

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

Test Code:

EPA TO-15

Instrument ID:

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Analyst:

Wida Ang

Sampling Media:

6.0 L Summa Canister(s)

Test Notes:

Date(s) Collected: 4/23 - 4/24/08

Date(s) Received: 4/28/08

Date(s) Analyzed: 5/2 - 5/5/08

		1,2-Dichlor	oethane-d4	Tolue	ene-d8	Bromofluo	robenzene	
Client Sample ID	<b>CAS Sample ID</b>	%	Acceptance	%	Acceptance	%	Acceptance	Data
		Recovered	Limits	Recovered	Limits	Recovered	Limits	Qualifier
Method Blank	P080502-MB	93	70-130	97	70-130	105	70-130	
Method Blank	P080505-MB	98	70-130	98	70-130	102	70-130	
Lab Control Sample	P080502-LCS	93	70-130	98	70-130	106	70-130	
Lab Control Sample	P080505-LCS	99	70-130	96	70-130	106	70-130	
WAT-SG-B2-042308	P0801215-001	92	70-130	97	70-130	104	70-130	
WAT-SG-1-042308	P0801215-002	94	70-130	97	70-130	103	70-130	
WAT-SG-1-042308	P0801215-002DUP	95	70-130	97	70-130	103	70-130	
WAT-SG-2-042308	P0801215-003	98	70-130	97	70-130	103	70-130	
WAT-SG-3-042308	P0801215-004	93	70-130	98	70-130	105	70-130	
WAT-SG-3a-042308	P0801215-005	101	70-130	97	70-130	103	70-130	
WAT-SG-4-042308	P0801215-006	95	70-130	97	70-130	104	70-130	
WAT-SG-DUP-042308	P0801215-007	99	70-130	98	70-130	105	70-130	
WAT-SG-FB-042308	P0801215-008	94	70-130	98	70-130	103	70-130	
WAT-SG-B4-042408	P0801215-009	96	70-130	96	70-130	102	70-130	
WAT-SG-5-042408	P0801215-010	100	70-130	98	70-130	102	70-130	
WAT-SG-5a-042408	P0801215-011	105	70-130	96	70-130	101	70-130	
WAT-SG-6-042408	P0801215-012	97	70-130	98	70-130	104	70-130	
WAT-SG-7-042408	P0801215-013	96	70-130	97	70-130	104	70-130	
WAT-SG-7a-042408	P0801215-014	100	70-130	97	70-130	103	70-130	
WAT-SG-8-042408	P0801215-015	98	70-130	97	70-130	104	70-130	
WAT-SG-9-042408	P0801215-016	101	70-130	98	70-130	102	70-130	

### LABORATORY CONTROL SAMPLE SUMMARY Page 1 of 2

Client: **CH2M Hill** 

Client Sample ID: Lab Control Sample

CAS Project ID: P0801215 Client Project ID: DOW - Waterloo, NY / 369548.B2.FI CAS Sample ID: P080502-LCS

Test Code:

EPA TO-15

Instrument ID:

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Analyst:

Test Notes:

Wida Ang

Sampling Media:

6.0 L Summa Canister

Date Analyzed: 5/02/08 Volume(s) Analyzed: NA Liter(s)

Date Collected: NA

Date Received: NA

					Project	
CAS#	Compound	Spike Amount	Result	% Recovery	Acceptance	Data
		ng	ng		Limits	Qualifier
74-87-3	Chloromethane	24.5	24.8	101	70-130	
75-01-4	Vinyl Chloride	24.8	29.5	119	70-130	
74-83-9	Bromomethane	25.0	28.1	112	70-130	
75-00-3	Chloroethane	25.0	29.9	120	70-130	
67-64-1	Acetone	26.8	29.5	110	70-130	
107-13-1	Acrylonitrile	25.5	30.3	119	70-130	
75-35-4	1,1-Dichloroethene	27.8	32.2	116	70-130	
75-09-2	Methylene Chloride	27.8	29.6	106	70-130	
75-15-0	Carbon Disulfide	25.0	27.7	111	70-130	
156-60-5	trans-1,2-Dichloroethene	26.5	29.8	112	70-130	
75-34-3	1,1-Dichloroethane	26.8	28.3	106	70-130	
78-93-3	2-Butanone (MEK)	27.0	28.2	104	70-130	
156-59-2	cis-1,2-Dichloroethene	27.0	29.8	110	70-130	
67-66-3	Chloroform	29.8	33.0	111	70-130	
107-06-2	1,2-Dichloroethane	26.3	27.8	106	70-130	
71-55-6	1,1,1-Trichloroethane	26.8	28.8	107	70-130	
71-43-2	Benzene	27.0	29.6	110	70-130	
56-23-5	Carbon Tetrachloride	26.0	30.1	116	70-130	

### LABORATORY CONTROL SAMPLE SUMMARY Page 2 of 2

Client:

**CH2M Hill** 

Client Sample ID: Lab Control Sample

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P080502-LCS

Test Code:

EPA TO-15

Instrument ID:

Test Notes:

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Analyst:

Sampling Media:

6.0 L Summa Canister

Date Collected: NA

Date Received: NA Date Analyzed: 5/02/08

Volume(s) Analyzed:

NA Liter(s)

CAS#	Compound	Spike Amount	Result	% Recovery	Acceptance	Data
		ng	ng		Limits	Qualifier
78-87-5	1,2-Dichloropropane	26.5	29.4	111	70-130	
75-27-4	Bromodichloromethane	27.8	31.5	113	70-130	
79-01-6	Trichloroethene	27.3	30.3	111	70-130	
10061-01-5	cis-1,3-Dichloropropene	25.0	27.9	112	70-130	
108-10-1	4-Methyl-2-pentanone	27.5	29.8	108	70-130	
10061-02-6	trans-1,3-Dichloropropene	28.0	31.2	111	70-130	
79-00-5	1,1,2-Trichloroethane	26.3	29.3	111	70-130	
108-88-3	Toluene	26.5	29.2	110	70-130	
124-48-1	Dibromochloromethane	27.0	29.6	110	70-130	
127-18-4	Tetrachloroethene	26.0	28.7	110	70-130	
108-90-7	Chlorobenzene	26.5	29.2	110	70-130	
100-41-4	Ethylbenzene	26.3	29.0	110	70-130	
179601-23-1	m,p-Xylenes	62.5	<b>68.7</b>	110	70-130	
75-25-2	Bromoform	31.3	36.2	116	70-130	
100-42-5	Styrene	26.3	29.8	113	70-130	
95-47-6	o-Xylene	29.8	32.7	110	70-130	
79-34-5	1,1,2,2-Tetrachloroethane	29.8	33.9	114	70-130	

### LABORATORY CONTROL SAMPLE SUMMARY Page 1 of 2

Client:

**CH2M Hill** 

Client Sample ID: Lab Control Sample

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P080505-LCS

Test Code:

EPA TO-15

Instrument ID:

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Analyst:

Sampling Media:

6.0 L Summa Canister

Date Analyzed: 5/05/08

Volume(s) Analyzed:

Date Collected: NA

Date Received: NA

NA Liter(s)

Test Notes:

					Project	
CAS#	Compound	Spike Amount	Result	% Recovery	Acceptance	Data
		ng	ng		Limits	Qualifier
74-87-3	Chloromethane	24.5	26.6	109	70-130	
75-01-4	Vinyl Chloride	24.8	28.4	115	70-130	
74-83-9	Bromomethane	25.0	28.2	113	70-130	
75-00-3	Chloroethane	25.0	29.1	116	70-130	
67-64-1	Acetone	26.8	30.0	112	70-130	
107-13-1	Acrylonitrile	25.5	30.2	118	70-130	handa milar risposition della seriesta a seriesta del seriesta della seriesta della seriesta della seriesta del
75-35-4	1,1-Dichloroethene	27.8	29.8	107	70-130	
75-09-2	Methylene Chloride	27.8	28.4	102	70-130	
75-15-0	Carbon Disulfide	25.0	27.1	108	70-130	
156-60-5	trans-1,2-Dichloroethene	26.5	28.8	109	70-130	
75-34-3	1,1-Dichloroethane	26.8	27.6	103	70-130	and the state of t
78-93-3	2-Butanone (MEK)	27.0	28.8	107	70-130	
156-59-2	cis-1,2-Dichloroethene	27.0	29.0	107	70-130	
67-66-3	Chloroform	29.8	32.1	108	70-130	
107-06-2	1,2-Dichloroethane	26.3	27.5	105	70-130	
71-55-6	1,1,1-Trichloroethane	26.8	28.5	106	70-130	
71-43-2	Benzene	27.0	28.6	106	70-130	
56-23-5	Carbon Tetrachloride	26.0	29.3	113	70-130	

### LABORATORY CONTROL SAMPLE SUMMARY Page 2 of 2

**Client:** 

**CH2M Hill** 

Client Sample ID: Lab Control Sample

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P080505-LCS

Test Code:

EPA TO-15

Instrument ID:

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Analyst:

Wida Ang

Sampling Media:

6.0 L Summa Canister

Date Received: NA Date Analyzed: 5/05/08

Volume(s) Analyzed:

Date Collected: NA

NA Liter(s)

Test Notes:

CAS#	Compound	Spike Amount ng	Result ng	% Recovery	Project Acceptance Limits	Data Qualifier
78-87-5	1,2-Dichloropropane	26.5	28.8	109	70-130	
75-27-4	Bromodichloromethane	27.8	31.2	112	70-130	
79-01-6	Trichloroethene	27.3	28.4	104	70-130	
10061-01-5	cis-1,3-Dichloropropene	25.0	28.1	112	70-130	
108-10-1	4-Methyl-2-pentanone	27.5	28.4	103	70-130	
10061-02-6	trans-1,3-Dichloropropene	28.0	31.5	113	70-130	
79-00-5	1,1,2-Trichloroethane	26.3	28.3	108	70-130	
108-88-3	Toluene	26.5	26.6	100	70-130	
124-48-1	Dibromochloromethane	27.0	28.1	104	70-130	•
127-18-4	Tetrachloroethene	26.0	25.7	99	70-130	
108-90-7	Chlorobenzene	26.5	26.5	100	70-130	
100-41-4	Ethylbenzene	26.3	26.4	100	70-130	
179601-23-1	m,p-Xylenes	62.5	62.8	100	70-130	
75-25-2	Bromoform	31.3	34.9	112	70-130	
100-42-5	Styrene	26.3	27.0	103	70-130	
95-47-6	o-Xylene	29.8	29.8	100	70-130	
79-34-5	1,1,2,2-Tetrachloroethane	29.8	31.9	107	70-130	

### LABORATORY DUPLICATE SUMMARY RESULTS

Page 1 of 2

**Client:** 

CH2M Hill

Client Sample ID: WAT-SG-1-042308

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-002DUP

Test Code:

EPA TO-15

Date Collected: 4/23/08

Instrument ID:

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Received: 4/28/08

Wida Ang

Date Analyzed: 5/2/08

Sampling Media:

6.0 L Summa Canister

Volume(s) Analyzed:

0.30 Liter(s)

Test Notes:

Analyst:

Container ID:

AC01311

Initial Pressure (psig):

0.0

Final Pressure (psig): 3.5

Canister Dilution Factor: 1.24

	Duplicate							
Compound	Sample	Result	Sample	Result	Average	% RPD	RPD	Data
	$\mu g/m^3$	ppbV	$\mu g/m^3$	ppbV	$\mu g/m^3$		Limit	Qualifier
Chloromethane	1.26	0.609	1.17	0.565	1.215	7	25	
Vinyl Chloride	ND	ND	ND	ND	-	-	25	$\mathbf{U}$
Bromomethane	ND	ND	ND	ND	-	-	25	$\mathbf{U}$
Chloroethane	ND	ND	ND	ND	-	-	25	$\mathbf{U}$
Acetone	16.5	6.97	16.7	7.05	16.6	1	25	J, B
Acrylonitrile	5.90	2.72	5.91	2.72	5.905	0.2	25	
1,1-Dichloroethene	ND	ND	ND	ND	-	-	25	${f U}$
Methylene Chloride	0.463	0.133	0.413	0.119	0.438	11	25	J
Carbon Disulfide	1.23	0.396	1.32	0.425	1.275	7	25	J
trans-1,2-Dichloroethene	ND	ND	ND	ND	-	-	25	U
1,1-Dichloroethane	ND	ND	ND	ND	-	-	25	$\mathbf{U}$
2-Butanone (MEK)	3.90	1.32	3.91	1.32	3.905	0.3	25	
cis-1,2-Dichloroethene	ND	ND	ND	ND	-	-	25	$\mathbf{U}$
Chloroform	1.96	0.401	1.98	0.406	1.97	1	25	
1,2-Dichloroethane	ND	ND	ND	ND	-	-	25	U
1,1,1-Trichloroethane	ND	ND	ND	ND		<u>-</u> ·	25	$\mathbf{U}$
Benzene	0.446	0.140	0.417	0.131	0.4315	7	25	
Carbon Tetrachloride	0.285	0.0454	0.277	0.0440	0.281	3	25	J

U = Compound was analyzed for, but not detected above the laboratory detection limit.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

### LABORATORY DUPLICATE SUMMARY RESULTS

Page 2 of 2

**Client:** 

CH2M Hill

Client Sample ID: WAT-SG-1-042308

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

CAS Project ID: P0801215

CAS Sample ID: P0801215-002DUP

Test Code:

EPA TO-15

Date Collected: 4/23/08

Instrument ID:

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Date Received: 4/28/08

Wida Ang

Date Analyzed: 5/2/08

Sampling Media:

6.0 L Summa Canister

Volume(s) Analyzed:

0.30 Liter(s)

Test Notes:

Analyst:

Container ID:

AC01311

Initial Pressure (psig):

0.0

Final Pressure (psig): 3.5

Canister Dilution Factor: 1.24

$ \begin{array}{cccccccccccccccccccccccccccccccccccc$	Data Qualifier
μg/m³ ppbV <b>μg/m³ ppbV</b> μg/m³ Limit	
1,2-Dichloropropane ND ND ND 25	$\mathbf{U}$
Bromodichloromethane ND ND ND 25	$\mathbf{U}$
<b>Trichloroethene</b> 1.10 0.205 1.13 0.210 1.115 <b>3</b> 25	
cis-1,3-Dichloropropene ND ND ND 25	$\mathbf{U}$
<b>4-Methyl-2-pentanone</b> 408 99.6 405 99.0 406.5 <b>0.7</b> 25	
trans-1,3-Dichloropropene ND ND ND 25	U
1,1,2-Trichloroethane ND ND ND 25	$\mathbf{U}$
<b>Toluene</b> 1.45 0.386 1.40 0.371 1.425 4 25	J
Dibromochloromethane ND ND ND 25	U
<b>Tetrachloroethene</b> 1.10 0.163 1.11 0.164 1.105 <b>0.9</b> 25	
Chlorobenzene ND ND ND 25	U
Ethylbenzene ND ND ND 25	$\mathbf{U}$
m,p-Xylenes 0.773 0.178 0.789 0.182 0.781 <b>2</b> 25	J
Bromoform ND ND ND 25	$\mathbf{U}$
Styrene ND ND ND 25	U
o-Xylene 0.273 0.0628 0.277 0.0638 0.275 <b>1</b> 25	J
1,1,2,2-Tetrachloroethane ND ND ND 25	U

U = Compound was analyzed for, but not detected above the laboratory detection limit.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

## RESULTS OF ANALYSIS

Page 1 of 1

Client:

CH2M Hill

CAS Project ID: P0801215

Client Project ID:

DOW - Waterloo, NY / 369548.B2.FI

**Method Blank Summary** 

Test Code:

EPA TO-15

Instrument ID:

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Analyst:

Wida Ang

Sampling Media:

6.0 L Summa Canister(s)

Lab File ID: 05020803.D Date Analyzed: 5/02/08

Time Analyzed: 08:11

Test Notes:

Client Sample ID	CAS Sample ID	Lab File ID	Time Analyzed
T.1.C. + 101	D000700 I CC	05020005 D	00.27
Lab Control Sample	P080502-LCS	05020805.D	09:27
WAT-SG-3-042308	P0801215-004	05020806.D	10:25
WAT-SG-B2-042308	P0801215-001	05020810.D	13:00
WAT-SG-7-042408	P0801215-013	05020811.D	13:38
WAT-SG-FB-042308	P0801215-008	05020814.D	15:46
WAT-SG-1-042308	P0801215-002	05020815.D	16:23
WAT-SG-1-042308 (Lab Duplicate)	P0801215-002DUP	05020816.D	17:01
WAT-SG-2-042308	P0801215-003	05020817.D	17:39
WAT-SG-4-042308	P0801215-006	05020819.D	18:54
WAT-SG-B4-042408	P0801215-009	05020821.D	20:15
WAT-SG-5-042408	P0801215-010	05020822.D	20:53

# RESULTS OF ANALYSIS Page 1 of 1

Client:

CH2M Hill

CAS Project ID: P0801215

**Client Project ID:** 

DOW - Waterloo, NY / 369548.B2.FI

**Method Blank Summary** 

Test Code:

EPA TO-15

Instrument ID:

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Lab File ID:

Lab File ID: 05050806.D

Analyst:

Wida Ang

Date Analyzed: 5/05/08

Sampling Media:

6.0 L Summa Canister(s)

Time Analyzed: 08:54

Test Notes:

Client Sample ID	CAS Sample ID	Lab File ID	Time Analyzed	
WAT-SG-7a-042408	P0801215-014	05050807.D	09:32	
WAT-SG-9-042408	P0801215-016	05050809.D	10:55	
Lab Control Sample	P080505-LCS	05050815.D	14:54	
WAT-SG-6-042408	P0801215-012	05050816.D	15:32	
WAT-SG-3a-042308	P0801215-005	05050819.D	17:50	
WAT-SG-5a-042408	P0801215-011	05050821.D	19:06	
WAT-SG-DUP-042308	P0801215-007	05050827.D	23:14	
WAT-SG-8-042408	P0801215-015	05050828.D	23:52	

Data Path : J:\MS16\DATA\2008 04\14\

Data File : 04140805.D

Acq On : 14 Apr 2008 10:53 am

Operator : CB Sample : BFB

Misc : S20-04070801

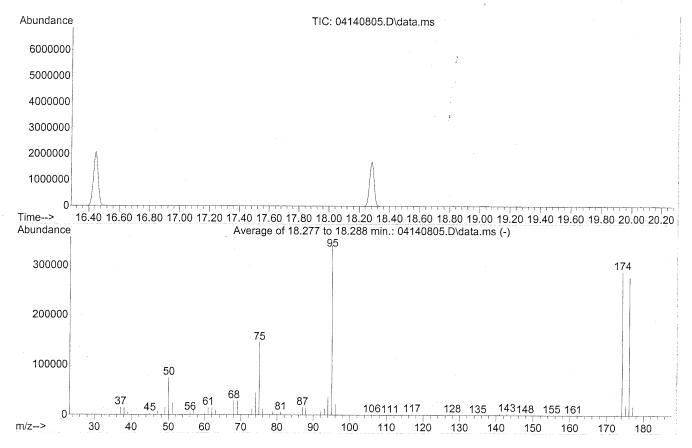
ALS Vial : 1 Sample Multiplier: 1

Integration File: RTEINT.P

Method: J:\MS16\METHODS\R16041408.M

Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)

Last Update : Thu Apr 10 07:29:28 2008



AutoFind: Scans 2479, 2480, 2481; Background Corrected with Scan 2469

Target	Rel. to	Lower	Upper	Rel.	Raw	Result
Mass		Limit%	Limit%	Abn%	Abn	Pass/Fail
50 75	95	8 · 30	40 66	21.9	74461 145195	PASS PASS
95 96	95 95	100	100 9	100.0	340565 22379	PASS PASS
173	174	0.00	2	0.7	2058	PASS
174	95	50	120	84.9	289109	PASS
175	174	93	9	7.0	20245	PASS
176	174		101	96.1	277717	PASS
177	176   	5	9	6.5	18179	PASS

CB04/14/08

Data Path : J:\MS16\DATA\2008 05\02\

Data File : 05020801.D

Acq On : 2 May 2008 6:39 am

Operator : WA

Sample : 25ng TO-15 CCV STD

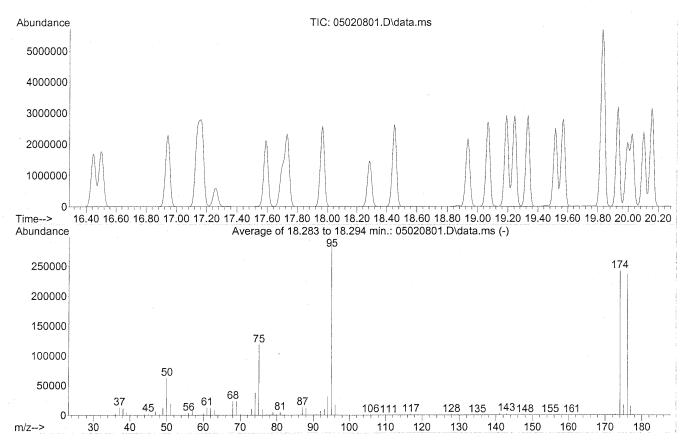
Misc : S20-04290801/S20-04110802 ALS Vial : 1 Sample Multiplier: 1

Integration File: RTEINT.P

Method : J:\MS16\METHODS\R16041408.M

Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)

Last Update : Tue Apr 15 07:25:17 2008



AutoFind: Scans 2480, 2481, 2482; Background Corrected with Scan 2469

Target	Rel. to	Lower	Upper	Rel.	Raw	Result	
Mass	Mass	Limit%	Limit%	Abn%	Abn	Pass/Fail	
50 75 95 96 173 174 175 176	95 95 95 95 174 95 174 174	8 30 100 5 0.00 50 4 93	40 66 100 9 2 120 9	22.0 41.6 100.0 6.3 0.7 85.1 7.3 97.8 6.4	62621 118403 284779 17874 1782 242325 17623 236907 15197	PASS PASS PASS PASS PASS PASS PASS	

NOA 5/2/08

Data Path : J:\MS16\DATA\2008 05\05\

Data File : 05050802.D

Acq On : 5 May 2008 5:41 am

Operator : WA

Sample : 25ng TO-15 CCV STD

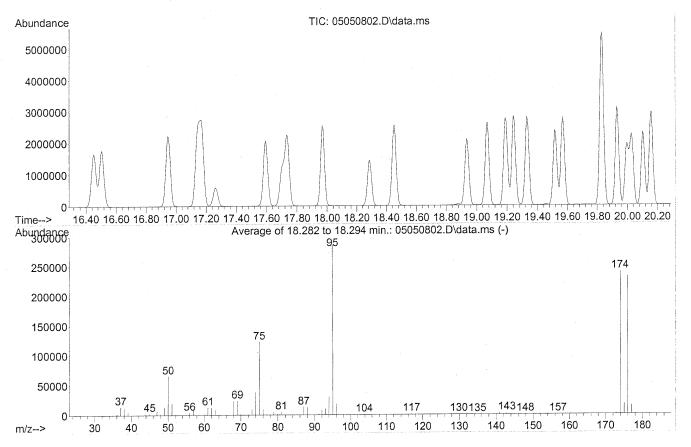
Misc : S20-04290801/S20-04110802 ALS Vial : 1 Sample Multiplier: 1

Integration File: rteint.p

Method : J:\MS03\METHODS\R3041708.M

Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)

Last Update : Thu Apr 17 14:25:29 2008



AutoFind: Scans 2480, 2481, 2482; Background Corrected with Scan 2470

Target	Rel. to	Lower	Upper	Rel.	Raw	Result
Mass	Mass	Limit%	Limit%	Abn%	Abn	Pass/Fail
50 75 95 96 173 174 175 176	95 95 95 174 95 174 174 176	8 30 100 5 0.00 50 4 93 5	40 66 100 9 2 120 9 101	22.9 42.9 100.0 6.4 0.8 84.0 7.1 96.8 6.5	65459 122848 286080 18408 1898 240384 17112 232597 15091	PASS PASS PASS PASS PASS PASS PASS PASS

: J:\MS16\METHODS\ : R16041408.M Method Path

Method File

Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)
Last Update : Tue Apr 15 07:25:17 2008
Response Via : Initial Calibration

Calibration Files

=04140811.D																																
D 20	%RSD	1 ! !																														
=04140810.	Avg	! ! ! ! ! ! !	4.2	13.83	6.5	3.5	1.5	1.9	6.0	6.6	9.	3.2	5.5	0.2	9.5	9	9.	1.	6.6	-	1.8	1.3	3.1	4.	2.5	1.1	0.7	1.7	0.2	0.	3.6	2.0
25	100	 	.86	2.385	.91	.38	.80	.30	.89	.93	.05	.02	.95	.26	.16	.34	.37	.09	.18	.09	.02	.02	.10	.10	.43	.30	.20	. 74	.95	.93	.48	. 82
0809.D	20		. 79	2.100	. 59	. 22	.71	.44	.72	.77	. 23	. 12	80	.04	.04	.49	.32	.01	. 74	96.	.99	.97	.83	96.	.08	.02	. 23	. 68	. 79	.87	.47	. 55
=0414	25		.67	2.055	.59	. 24	.61	. 24	. 73	89	.21	.21	.87	.06	.04	.75	.35	.01	. 79	.98	.84	.94	. 82	. 98	.20	.03	.24	.68	.84	.90	.48	. 53
0 2.0	5.0		1.65	2.138	.91	. 25	.67	.25	.75	.91	.23	.33	.95	. 21	.06	.61	.42	.02	.90	.02	.97	.94	. 92	00.	.30	.07	.24	.70	. 88	.92	.49	. 62
40808.L	1.0	S	.671	$\alpha$	. 55	.31	.70	.03	. 68	.79	.85	.50	.82	.33	.01	.80	.02	.99	.70	.03	.71	.95	.77	.02	.33	.15	. 22	.76	. 88	. 92	4.9	.64
0 =0414	0.5	1 1	.77	2.431	.14	.33	.77	. 28	. 93	.96	.88	. 52	.94	.67	.05	.43	.14	.12	.80	.12	. 92	.97	.12	.99	.43	.19	.17	. 72	. 91	5	4.9	.76
D 1.0	0.1	1	.25	2.893	. 92	. 68	.21	.40	. 25	. 24	.76	.4	.24		.54	.72	.63	.33	. 88	.46	.34	. 25	.18	.50	.97	.98	.13	.91	.38	1.150	. 58	.44
140807.	   	•	2.24	2.754	. 65	.61	. 94	.48	. 18	.96	. 18				2.355	. 59	99.	.15	.44		2.346	.10		. 23	2.688	,64			99	0,823	.36	. I 4
04140806.D 0.5 =041	Compound	ochlorome	opene	chlorodifl	hloro	reon 114	iny	1,3-Butadiene	LC.	Chloroethane	thanol	ceton	Acrolein	cetone	orof	sopropanol	cry]	,1-Dichlor	rt-Butanol	thylene C	lyl Chlorid	chlorotrifl	rbon Disulf	ans-1,	,1-Dich	ethy	nY	-Butanon	is-1,2-Dichlo	isopropyl	thyl	n-Hexane
00 = 0	 	) IR	T (			_ '	_	<u> </u>			<u> </u>	<u> </u>	_		<u> </u>	<u> </u>	<u> </u>	_	<u> </u>	$\overline{}$	<u> </u>	<u> </u>	<u> </u>	_	$\overline{}$	_	<u> </u>	<u> </u>	· .			_
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SOP VOA-TO15 (CASS TO-15/GC-MS) .723 2.026 1.615 1.563 1.535 1.492 1.465 1.631 .474 1.488 1.470 1.521 1.505 1.507 1.499 1.495 .436 0.895 0.686 0.680 0.706 0.675 0.655 0.676 .305 1.695 1.333 1.314 1.359 1.331 1.313 1.379 .070 2.184 1.761 1.726 1.720 1.656 1.619 1.819	ISTD	559 2.529 2.516 2.526 2.490 2.526 2.519 2.524 197 3.468 2.823 2.644 2.614 2.553 2.506 2.829 284 2.927 2.378 2.351 2.414 2.295 2.163 2.545 844 0.967 0.759 0.766 0.798 0.790 0.778 0.815 818 0.954 0.761 0.737 0.735 0.726 0.707 0.777 616 2.904 2.281 2.218 2.570 2.467 2.335 2.484 805 1.003 0.863 0.796 0.827 0.798 0.764 0.837 943 1.037 0.863 0.777 0.779 0.761 0.742 0.837 257 2.495 2.048 1.908 1.895 1.848 1.800 2.036 740 3.979 3.214 3.043 3.079 3.011 2.938 3.286 411 2.575 2.097 1.991 2.005 1.967 1.930 2.139 345 2.445 2.005 1.995 2.053 2.013 1.963 2.117
Method Path: J:\MS16\METHO Method File: R16041408.M Title: EPA TO-15 per S' 2) T Chloroform 3) S 1,2-Dichloroet 1. 4) T Tetrahydrofuran 0. 5) T Ethyl tert-But 1. 6) T 1,2-Dichloroet 2.	7) IR 1,4-Difluorobenzen  8) T 1,1,1-Trichlor 0  9) T Isopropyl Acetate 0  1) T Benzene 1  2) T Carbon Tetrach 0  3) T Cyclohexane 0  4) T tert-Amyl Meth 0  5) T 1,2-Dichloropr 0  7) T Trichloroethene 0  8) T 1,4-Dioxane 0  9) T Esooctane 1  1) T n-Heptane 0  2) T cis-1,3-Dichlo 0  3) T 4-Methyl 2-pen 0  4) T trans-1,3-Dichlo 0  4) T trans-1,3-Dichlo 0	6) I Chlorobenzene-d5 ( 7) S Toluene-d8 (SS2) 2. 8) T Toluene 3. 9) T 2-Hexanone 3. 0) T Dibromochlorom 0. 1) T 1,2-Dibromoethane 0. 2) T Butyl Acetate 2. 3) T n-Octane 0. 4) T Tetrachloroethene 0. 5) T Chlorobenzene 2. 6) T Ethylbenzene 3. 7) T m- & p-Xylene 2. 8) T Styrene 2. 9) T Styrene 2.
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Page: 3

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	2.20	6 0.946	0.87	3.31	1.59	3.85	3.59	3.34	2.91	1.61	3.53	2.95	2.13	2.32	1.88	1.82	3.81	3.32	2.78	1.71	0.90	0.51	2.10	0.30	3.96	1.72	0.50
	.84	0.84	. 88	.07	.48	. 59	.39	.01	.67	.61	.30	.65	.91	.25	.67	.63	.58	1	.64	.49	.87	.55	.97	.28	.05	.61	.46
	00	0.877	.88	.12	.53	.66	.44	.06	.70	.64	.37	.71	90	.30	.71	.68	.67	.19	.68	.56	.90	.56	.07	.29	.12	.72	.48
) GC	2.118	.90	.87	.21	. 58	. 74	47	.15	. 74	.68	.43	.76	.06	.33	.76	.72	. 73	.27	. 74	. 63	.94	.57	.13	.30	.10	. 73	.49
T0-1		.876	.878	.022	.508	.503	.195	860.	.707	.493	.165	.693	.972	.095	.714	.654	.506	.974	.473	.569	.850	.423	.946	.256	.923	.514	.429
(CAS	.178	.899	.870	.138	.529	689.	.305	.256	.796	.476	.338	.863	.013	.116	.785	.723	.583	.109	.555	.639	.855	.426	976.	.270	.041	. 568	.459
A-T01		.121	.876	. 983	.889	,604	.218	.045	.591	.814	.215	.557	.480	.593	.293	.193	.583	.891	.289	.097	.072	.561	.478	.357	.054	. 944	.610
_ >	.50	.102	.870	.629	.644	.203	.141	.754	.172	.578	.890	.433	.476	.575	.251	.125	.030	.729	.116	.989	.830	.510	.179	.387	.441	.986	.602
J:\MS16\METHODS R16041408.M PA TO-15 per SOP		Tetr	luoroben (		nene	ylbenzene	ltoluene	ne	$hy \dots$	Methylst	ltoluene	$\mathtt{methy}$		loride	orobe	orobe	ne	:	$hy \cdots$	orobe	Φ	no-3	4)	richlor	lene	cane 1	loro-1,3 (
Path : File :	n-Nonar	,1,2	Bromof]	umene	alpha-F	1	-Ethy	-Eth}	,3,5-	lpha-	2-Ethy]	-		Benzyl Ch		,4-D	sec-But	-Isog	, 2, 3-	-D	-Limc	,2-Di	1		Naphtha	n-Dode	Hexach]
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<sup>(#) =</sup> Out of Range

Data Path : J:\MS16\DATA\2008 05\02\

Data File : 05020801.D

Acq On : 2 May 2008 6:39 am

Operator : WA

Sample : 25ng TO-15 CCV STD

Misc : S20-04290801/S20-04110802 ALS Vial : 1 Sample Multiplier: 1

Quant Time: May 02 12:20:28 2008

Quant Method : J:\MS16\METHODS\R16041408.M

Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)

QLast Update : Tue Apr 15 07:25:17 2008

Response via: Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min

Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev Area%	Dev(min)
1 T T T T T T T T T T T T T T T T T T T	Bromochloromethane (IS1) Propene Dichlorodifluoromethane Chloromethane Freon 114 Vinyl Chloride 1,3-Butadiene Bromomethane Chloroethane Ethanol Acetonitrile Acrolein Acetone Trichlorofluoromethane Isopropanol Acrylonitrile 1,1-Dichloroethene tert-Butanol Methylene Chloride Allyl Chloride Trichlorotrifluoroethane Carbon Disulfide trans-1,2-Dichloroethene 1,1-Dichloroethane Methyl tert-Butyl Ether Vinyl Acetate 2-Butanone cis-1,2-Dichloroethene Diisopropyl Ether Ethyl Acetate n-Hexane Chloroform 1,2-Dichloroethane-d4 (SS1) Tetrahydrofuran Ethyl tert-Butyl Ether 1,2-Dichloroethane	AvgRF 1.000 1.867 2.385 2.913 1.383 1.805 1.308 0.895 1.020 0.956 1.266 2.160 4.346 2.370 1.098 2.160 4.181 1.098 2.103 2.103 2.109 2.103 2.109 2.103 2.109 2.103 2.109 2.109 2.103 2.109 2.109 2.109 2.109 2.109 2.109 2.103 2.109 2.	1.000 1.830 2.463 3.106 1.996 1.838 0.951 1.351 3.477 1.103 1.2516 2.619 1.154 4.016 1.154 4.016 1.154 4.016 1.154 4.016 1.153 2.2516 2.187 2.450 3.388 0.772 2.450 3.388 0.772 2.450 3.388 0.772 1.017 0.547 3.077 1.468 1.468 1.787	*Dev Area*  0.0 87  2.0 90  -3.3 100  -3.6 4 92  -3.4 10.6 # 128  -9.9 113  -12.4 -28.5 -15.4 100  -4.5 9 91  -15.4 106  -4.5 9 95  -15.5 9 94  -11.0 1 95  -37.8 # 10.6  -3.8 995  -11.0 1 95  -37.8 # 10.6  -3.8 995  -1.0 8	0.01 -0.03 -0.02 -0.01 -0.02 -0.01 0.00 0.00 0.00 0.03 0.02 0.00 0.03 0.02 0.00 0.03 0.02 0.00 0.03 0.02 0.00 0.03 0.02 0.00 0.03
37 IR 38 T	1,4-Difluorobenzene (IS2) 1,1,1-Trichloroethane	1.000	1.000	0.0 83 -5.5 92	

Page: 1

Data Path : J:\MS16\DATA\2008 05\02\

Data File : 05020801.D

Acq On : 2 May 2008 6:39 am

Operator : WA

Sample : 25ng TO-15 CCV STD

Misc : S20-04290801/S20-04110802 ALS Vial : 1 Sample Multiplier: 1

Quant Time: May 02 12:20:28 2008

Quant Method : J:\MS16\METHODS\R16041408.M

Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)

QLast Update : Tue Apr 15 07:25:17 2008

Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min

Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev Area	% Dev(min)
39 T 40 T 41 T 42 T 43 T 44 T 45 T 46 T 47 T 48 T 50 T 51 T 51 T 52 T 55 T	Isopropyl Acetate 1-Butanol Benzene Carbon Tetrachloride Cyclohexane tert-Amyl Methyl Ether 1,2-Dichloropropane Bromodichloromethane Trichloroethene 1,4-Dioxane Isooctane Methyl Methacrylate n-Heptane cis-1,3-Dichloropropene 4-Methyl-2-pentanone trans-1,3-Dichloropropene 1,1,2-Trichloroethane	0.212 0.284 1.026 0.378 0.412 0.719 0.321 0.311 0.340 0.207 1.693 0.111 0.267 0.407 0.370 0.365 0.271	0.239 0.310 1.059 0.438 0.419 0.774 0.336 0.334 0.213 1.822 0.126 0.282 0.443 0.382 0.400 0.283	-9.2 6 -3.2 9 -15.9 9 -1.7 9 -7.6 9 -4.7 9 -4.1 9 -2.9 8 -7.6 9 -13.5 8 -5.6 9 -8.8 9 -3.2 8	1 0.00 3 0.00 9 0.00 3 0.00 7 0.00 2 0.00 2 0.00 7 0.00 1 0.00
1 S T T T T T T T T T T T T T T T T T T	Chlorobenzene-d5 (IS3) Toluene-d8 (SS2) Toluene 2-Hexanone Dibromochloromethane 1,2-Dibromoethane Butyl Acetate n-Octane Tetrachloroethene Chlorobenzene Ethylbenzene m- & p-Xylene Bromoform Styrene o-Xylene n-Nonane 1,1,2,2-Tetrachloroethane Bromofluorobenzene (SS3) Cumene alpha-Pinene n-Propylbenzene	1.000 2.524 2.829 2.545 0.815 0.777 2.484 0.837 0.839 2.036 3.286 2.139 0.481 2.117 2.304 2.202 0.946 0.877 3.313 1.596 3.858	1.000 2.475 2.862 2.272 0.867 0.806 2.455 0.891 0.843 2.049 3.314 2.155 0.522 2.164 2.313 2.211 0.986 0.921 3.434 1.646 4.014	0.0 1.9 8 1.2 9 10.7 -6.4 9 -3.7 9 1.2 -6.5 9 -0.5 9 -0.6 9 -0.7 -8.5 9 -0.7 -8.5 -0.4 9 -0.4 -4.2 -5.0 8 -3.7 8 -3.7 -	4 0.00 3 0.00 2 0.00 9 0.00 1 0.00 0 0.00 0 0.00 1 0.00 1 0.00 0 0.00 0 0.00 0 0.00 0 0.00 8 0.00 0 0.00 8 0.00 8 0.00 9 0.00 7 0.00

Data Path : J:\MS16\DATA\2008 05\02\

Data File : 05020801.D

Acq On : 2 May 2008 6:39 am

Operator : WA

Sample : 25ng TO-15 CCV STD Misc : S20-04290801/S20-04 : S20-04290801/S20-04110802 ALS Vial : 1 Sample Multiplier: 1

Quant Time: May 02 12:20:28 2008
Quant Method : J:\MS16\METHODS\R16041408.M

Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)

QLast Update: Tue Apr 15 07:25:17 2008 Response via: Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min

Max. RRF Dev: 30% Max. Rel. Area: 200%

	Compound	AvgRF	CCRF	%Dev A	rea%	Dev(min)
77 T	3-Ethyltoluene	3.596	3.623	-0.8	87	0.00
78 T	4-Ethyltoluene	3.341	3.427	-2.6	91	0.00
79 T	1,3,5-Trimethylbenzene	2.912	2.920	-0.3	89	0.00
80 T	alpha-Methylstyrene	1.615	1.617	-0.1	80	0.00
81 T	2-Ethyltoluene	3.532	3.645	-3.2	89	0.00
82 T	1,2,4-Trimethylbenzene	2.954	2.941	0.4	89	0.00
83 T	n-Decane	2.131	2.207	-3.6	90	0.00
84 T	Benzyl Chloride	2.323	2.421	-4.2	87	0.00
85 T	1,3-Dichlorobenzene	1.885	1.906	-1.1	90	0.00
86 T	1,4-Dichlorobenzene	1.820	1.850	-1.6	90	0.00
87 T	sec-Butylbenzene	3.813	3.982	-4.4	89	0.00
T 88	p-Isopropyltoluene	3.327	3.442	-3.5	88	0.00
89 T	1,2,3-Trimethylbenzene	2.787	2.899	-4.0	89	0.00
90 T	1,2-Dichlorobenzene	1.712	1.761	-2.9	91	0.00
91 T	d-Limonene	0.904	0.870	3.8	77	0.00
92 T	1,2-Dibromo-3-Chloropropane	0.518	0.617	-19.1	89	0.00
93 T	n-Undecane	2.109	2.319	-10.0	91	0.00
94 T	1,2,4-Trichlorobenzene	0.308	0.328	-6.5	90	0.00
95 T	Naphthalene	3.962	4.244	-7.1	87	0.00
96 T	n-Dodecane	1.728	2.173	-25.8	105	0.00
97 T	Hexachloro-1,3-butadiene	0.508	0.520	-2.4	88	0.00

<sup>(#) =</sup> Out of Range

SPCC's out = 0 CCC's out = 0

Data Path : J:\MS16\DATA\2008 05\05\

Data File : 05050802.D

Acq On : 5 May 2008 5:41 am

Operator : WA

Sample : 25ng TO-15 CCV STD

Misc : S20-04290801/S20-04110802 ALS Vial : 1 Sample Multiplier: 1

Quant Time: May 05 09:24:03 2008

Quant Method: J:\MS16\METHODS\R16041408.M

Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)

QLast Update: Tue Apr 15 07:25:17 2008

Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min

Max. RRF Dev: 30% Max. Rel. Area: 200%

	Compound	AvgRF	CCRF	%Dev Are	ea%	Dev(min)
1 IR 2 T T T T T T T T T T T T T T T T T T T	Bromochloromethane (IS1) Propene Dichlorodifluoromethane Chloromethane Freon 114 Vinyl Chloride 1,3-Butadiene Bromomethane Chloroethane Ethanol Acetonitrile Acrolein Acetone Trichlorofluoromethane Isopropanol Acrylonitrile 1,1-Dichloroethene tert-Butanol Methylene Chloride Allyl Chloride Trichlorotrifluoroethane Carbon Disulfide trans-1,2-Dichloroethene 1,1-Dichloroethane Methyl tert-Butyl Ether Vinyl Acetate 2-Butanone cis-1,2-Dichloroethene Diisopropyl Ether Ethyl Acetate	1.000 1.867 2.385 2.913 1.383 1.805 1.308 0.895 0.956 1.051 3.020 0.956 1.266 2.160 4.346 2.370 1.094 4.181 1.098 2.021 1.020 4.109 2.103 2.431 3.304 0.209 0.747 1.955 0.936 0.485	1.000 1.930 2.513 3.240 1.446 2.035 1.871 0.926 0.972 1.398 3.648 1.107 1.518 2.292 4.628 2.668 1.156 4.076 1.124 2.203 1.142 4.471 2.220 2.505 3.412 0.283 0.779 2.038 1.027 0.560	0.0 -3.4 -5.4 -11.2 -4.6 -12.7 -43.0 -3.5 -4.0 -3.5 -12.6 -5.7 -2.4 -9.0 -12.0 -3.3 -3.5 -4.3 -4.3 -4.3 -4.3 -4.3 -4.2 -9.7 -15.5	-899990208999098998999999999999999999999	0.02 -0.02 0.00 0.00 0.00 0.00 0.00 0.00 0.01 0.04 0.03 0.01 0.04 0.03 0.01 0.04 0.03 0.01 0.04 0.03 0.01 0.04 0.03 0.01 0.04 0.03
	Ethyl Acetate n-Hexane Chloroform			-15.5 -8.8		
33 S 34 T 35 T 36 T	1,2-Dichloroethane-d4(SS1) Tetrahydrofuran Ethyl tert-Butyl Ether 1,2-Dichloroethane	1.495 0.676 1.379 1.819	1.438 0.742 1.470 1.820	3.8 -9.8 -6.6	79 87 90 88	0.02 0.01 0.01 0.02
37 IR 38 T	1,4-Difluorobenzene (IS2) 1,1,1-Trichloroethane	1.000	1.000		81 89	0.01

Data Path : J:\MS16\DATA\2008 05\05\

Data File : 05050802.D

Acq On : 5 May 2008 5:41 am

Operator : WA

Sample : 25ng TO-15 CCV STD

: S20-04290801/S20-04110802 Misc ALS Vial : 1 Sample Multiplier: 1

Quant Time: May 05 09:24:03 2008
Quant Method : J:\MS16\METHODS\R16041408.M

Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)

QLast Update: Tue Apr 15 07:25:17 2008 Response via: Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min

Max. RRF Dev: 30% Max. Rel. Area: 200%

	Compound	AvgRF	CCRF	%Dev Area%	Dev(min)
39 T 40 T 41 T 42 T 43 T 45 T 46 T 47 T 48 T 50 T 51 T 52 T 53 T 54 T	Isopropyl Acetate 1-Butanol Benzene Carbon Tetrachloride Cyclohexane tert-Amyl Methyl Ether 1,2-Dichloropropane Bromodichloromethane Trichloroethene 1,4-Dioxane Isooctane Methyl Methacrylate n-Heptane cis-1,3-Dichloropropene 4-Methyl-2-pentanone trans-1,3-Dichloropropene 1,1,2-Trichloroethane	0.212 0.284 1.026 0.378 0.412 0.719 0.321 0.311 0.340 0.207 1.693 0.111 0.267 0.407 0.370 0.365 0.271	0.240 0.295 1.055 0.430 0.418 0.766 0.335 0.338 0.351 0.209 1.828 0.124 0.281 0.439 0.380 0.397 0.285	-13.2 89 -3.9 61 -2.8 90 -13.8 87 -1.5 92 -6.5 88 -4.4 90 -8.7 89 -1.0 84 -8.0 90 -11.7 83 -5.2 89 -2.7 84 -8.8 87 -5.2 90	0.02 0.02 0.02 0.02 0.01 0.01 0.01 0.01
56 I S T T T T T T T T T T T T T T T T T T	Chlorobenzene-d5 (IS3) Toluene-d8 (SS2) Toluene 2-Hexanone Dibromochloromethane 1,2-Dibromoethane Butyl Acetate n-Octane Tetrachloroethene Chlorobenzene Ethylbenzene m- & p-Xylene Bromoform Styrene o-Xylene n-Nonane 1,1,2,2-Tetrachloroethane Bromofluorobenzene (SS3) Cumene alpha-Pinene n-Propylbenzene	1.000 2.524 2.829 2.545 0.815 0.777 2.484 0.837 0.839 2.036 3.286 2.139 0.481 2.117 2.304 2.202 0.946 0.877 3.313 1.596 3.858	1.000 2.454 2.839 2.271 0.859 0.795 2.488 0.903 0.825 2.041 3.286 2.129 0.518 2.142 2.284 2.279 0.987 0.911 3.382 1.621 4.001	0.0 82 2.8 81 -0.4 89 10.8 77 -5.4 88 -2.3 89 -0.2 79 -7.9 89 1.7 87 -0.2 88 0.0 87 0.5 87 -7.7 87 -1.2 85 0.9 87 -3.5 88 -4.3 90 -3.5 88 -4.3 90 -3.6 84 -3.7 87	0.00 0.01 0.00 0.00 0.01 0.00 0.00 0.00 0.00 0.00 0.00 0.00 0.00 0.00 0.00 0.00

### matuace conclinating carratacton reporc

Data Path : J:\MS16\DATA\2008 05\05\

Data File : 05050802.D

Acq On : 5 May 2008 5:41 am

Operator : WA

Sample : 25ng TO-15 CCV STD

: S20-04290801/S20-04110802 Misc ALS Vial : 1 Sample Multiplier: 1

Quant Time: May 05 09:24:03 2008

Quant Method: J:\MS16\METHODS\R16041408.M

Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)

QLast Update: Tue Apr 15 07:25:17 2008 Response via: Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min

Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev Ai	cea%	Dev(min)
77 T	3-Ethyltoluene	3.596	3.667	-2.0	87	0.00
78 T	4-Ethyltoluene	3.341	3.309	1.0	86	0.00
79 T	1,3,5-Trimethylbenzene	2.912	2.870	1.4	86	0.00
80 T	alpha-Methylstyrene	1.615	1.573	2.6	76	0.00
81 T	2-Ethyltoluene	3.532	3.634	-2.9	87	0.00
82 T	1,2,4-Trimethylbenzene	2.954	2.908	1.6	86	0.00
83 T	n-Decane	2.131	2.221	-4.2	88	0.00
84 T	Benzyl Chloride	2.323	2.413	-3.9	85	0.00
85 T	1,3-Dichlorobenzene	1.885	1.886	-0.1	88	0.00
86 T	1,4-Dichlorobenzene	1.820	1.830	-0.5	87	0.00
87 T	sec-Butylbenzene	3.813	3.966	-4.0	87	0.00
88 T	p-Isopropyltoluene	3.327	3.408	-2.4	85	0.00
89 T	1,2,3-Trimethylbenzene	2.787	2.890	-3.7	86	0.00
90 T	1,2-Dichlorobenzene	1.712	1.752	-2.3	88	0.00
91 T	d-Limonene	0.904	0.851	5.9	74	0.00
92 T	1,2-Dibromo-3-Chloropropane	0.518	0.620	-19.7	88	0.00
93 T	n-Undecane	2.109	2.331	-10.5	90	0.00
94 T	1,2,4-Trichlorobenzene	0.308	0.330	-7.1	89	0.00
95 T	Naphthalene	3.962	4.151	-4.8	83	0.00
96 T	n-Dodecane	1.728	2.162	-25.1	102	0.00
97 T	Hexachloro-1,3-butadiene	0.508	0.523	-3.0	86	0.00

<sup>(#) =</sup> Out of Range

SPCC's out = 0 CCC's out = 0

**RESULTS OF ANALYSIS** Page 1 of 1

Client:

**CH2M Hill** 

CAS Project ID: P0801215

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

### **Internal Standard Area and RT Summary**

Test Code:

EPA TO-15

Instrument ID:

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Lab File ID: 05020801.D 5/2/08

Analyst:

Wida Ang

Date Analyzed:

Sampling Media:

6.0 L Summa Canister(s)

Time Analyzed: 06:39

Test Notes:

		IS1 (BCM) IS2 (DFB				IS3 (CBZ)	
		AREA #	RT #	AREA #	RT #	AREA #	RT #
	24 Hour Standard	328364	9.22	1386260	11.36	583340	16.45
	Upper Limit	459710	9.55	1940764	11.69	816676	16.78
	Lower Limit	197018	8.89	831756	11.03	350004	16.12
			-				
	Client Sample ID						
01	Method Blank	319561	9.18	1348291	11.34	568631	16.44
02	Lab Control Sample	325686	9.23	1361442	11.37	-571978	16.45
03	WAT-SG-3-042308	322553	9.19	1364976	11.34	575596	16.45
04	WAT-SG-B2-042308	320779	9.19	1340869	11.35	565181	16.45
05	WAT-SG-7-042408	313548	9.20	1337401	11.35	564597	16.45
06	WAT-SG-FB-042308	308726	9.19	1310306	11.34	550257	16.44
07	WAT-SG-1-042308	309028	9.19	1291100	11.35	548762	16.45
08	WAT-SG-1-042308 (Lab Duplicate)	306175	9.20	1300972	11.35	547690	16.45
09	WAT-SG-2-042308	297029	9.21	1276048	11.36	540559	16.45
10	WAT-SG-4-042308	312570	9.20	1310803	11.35	556700	16.45
11	WAT-SG-B4-042408	275684	9.20	1153300	11.34	501911	16.45
12	WAT-SG-5-042408	276441	9.20	1190544	11.34	508037	16.44
13							
14							
15							
16							
17							
18							
19							
20							

IS1 (BCM) = Bromochloromethane

IS2 (DFB) = 1,4-Difluorobenzene

IS3 (CBZ) = Chlorobenzene-d5

AREA UPPER LIMIT = 140% of internal standard area

AREA LOWER LIMIT = 60% of internal standard area

RT UPPER LIMIT = 0.33 minutes of internal standard RT

RT LOWER LIMIT = 0.33 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.

Verified By:\_\_\_\_

**RESULTS OF ANALYSIS** Page 1 of 1

Client:

**CH2M Hill** 

CAS Project ID: P0801215

Client Project ID: DOW - Waterloo, NY / 369548.B2.FI

### **Internal Standard Area and RT Summary**

Test Code:

EPA TO-15

Instrument ID:

Tekmar AUTOCAN/Agilent 5975Cinert/6890N/MS16

Lab File ID: 05050802.D 5/5/08

Analyst:

Wida Ang

Date Analyzed:

Sampling Media:

6.0 L Summa Canister(s)

Time Analyzed: 05:41

Test Notes:

		IS1 (BCM)		IS2 (DFB)		IS3 (CBZ)	
		AREA #	RT #	AREA #	RT #	AREA #	RT #
	24 Hour Standard	314157	9.23	1343679	11.37	570129	16.45
	Upper Limit	439820	9.56	1881151	11.70	798181	16.78
	Lower Limit	188494	8.90	806207	11.04	342077	16.12
	Client Sample ID						
01	Method Blank	283657	9.19	1218732	11.34	515790	16.44
02	WAT-SG-7a-042408	278358	9.19	1195074	11.34	512911	16.45
03	WAT-SG-9-042408	276807	9.19	1198759	11.35	509213	16.44
04	Lab Control Sample	282838	9.23	1207055	11.37	525387	16.45
05	WAT-SG-6-042408	298964	9.19	1278940	11.34	541935	16.45
06	WAT-SG-3a-042308	270641	9.20	1164207	11.35	503590	16.45
07	WAT-SG-5a-042408	232996	9.20	1023989	11.35	451363	16.45
08	WAT-SG-DUP-042308	309518	9.19	1359291	11.34	570320	16.44
09	WAT-SG-8-042408	311326	9.19	1357811	11.34	575259	16.45
10							
11							
12							
13							
14							
15							
16							
17							
18							
19							
20							

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RT LOWER LIMIT = 0.33 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.

Verified By:

# QC Certification



Columbia Analytical Services, Inc. 2655 Park Center Drive, Suite A Simi Valley, CA 93065 Ph. 805-526-7161 Fax 805-526-7270

Comments	EPA TO-15 (Client Specified)	SIM	EPA TO-15 (Client Specified)	SIM	EPA TO-15 (Client Specified)	SIM	EPA TO-15 (Client Specified)	SIM	SIM	SIM	SIM	SIM												
<b>QC</b> Results	Pass w/ Conditions	Pass w/ Conditions	Pass w/ Conditions	Pass w/ Conditions	Pass w/ Conditions	Pass w/ Conditions	Pass w/ Conditions	Pass w/ Conditions	Pass w/ Conditions	Pass w/ Conditions	Pass w/ Conditions	Pass w/ Conditions	Pass w/ Conditions	Pass w/ Conditions	Pass w/ Conditions	Pass w/ Conditions	Pass w/ Conditions							
Date Analyzed	04/21/2008	04/21/2008	04/21/2008	04/21/2008	04/21/2008	04/21/2008	04/21/2008	04/21/2008	04/21/2008	04/21/2008	04/21/2008	04/21/2008	04/21/2008	04/21/2008	04/21/2008	04/21/2008	04/21/2008	04/09/2008	04/10/2008	04/21/2008	04/21/2008	04/21/2008	04/21/2008	04/21/2008
Cleaned Date	04/11/2008	04/17/2008	04/02/2008	04/17/2008	04/11/2008	04/02/2008	04/02/2008	04/02/2008	04/11/2008	04/02/2008	04/17/2008	04/02/2008	04/17/2008	04/17/2008	04/17/2008	04/17/2008	04/17/2008	04/09/2008	04/10/2008	04/18/2008	04/18/2008	04/18/2008	04/18/2008	04/18/2008
Container IDs	AC00097*	AC00615	AC00648	AC00802	AC00812*	AC00814	AC00940	AC00950	AC01020*	AC01168	AC01189	AC01209	AC01211	AC01235	AC01276	AC01311	AC01387	FC00659	FC00693	OA00572	OA00584	OA00585	OA00587	OA00589

\* QC Canister

Container IDs	Cleaned Date	Date Analyzed	OC Results	Comments
OA00590	04/18/2008	04/21/2008		
OA00591	04/18/2008	04/21/2008		
OA00594	04/18/2008	04/21/2008		
OA00601	04/18/2008	04/21/2008		
OA00607	04/18/2008	04/21/2008		
OA00611	04/18/2008	04/21/2008		
OA00616	04/18/2008	04/21/2008		
OA00617	04/18/2008	04/21/2008		
OA00621	04/18/2008	04/21/2008		
OA00624	04/18/2008	04/21/2008		