Site Characterization Investigation Work Plan Addendum, Essex-Hope Site in Jamestown, New York Site Number 907015

Prepared for

Essex Specialty Products, Inc.

June 2017



# **Professional Engineer Certification**

I, Key Rosebrook, certify that I am currently a New York State-registered professional engineer as defined in 6 New York Codes, Rules, and Regulations Part 375, and that this Site Characterization Investigation Work Plan Addendum was prepared in accordance with all applicable statutes and regulations, and in substantial conformance with NYSDEC DER-10: Technical Guidans for Site Investigation and Remediation.

Signature:

Date:

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# Acronyms and Abbreviations

bgs below ground surface

CBTEX cumene, benzene, toluene, ethylbenzene, and xylene

CH2M CH2M HILL Engineers, Inc.

CSM conceptual site model

DPT direct-push technology

EPA U.S. Environmental Protection Agency

ESP Essex Specialty Products, Inc.

MW monitoring well

LNAPL light nonaqueous phase liquid

NPLS North Parking Lot Sump

NYSDEC New York State Department of Environmental Conservation

NYSDOH New York State Department of Health

O&M operations and maintenance

PVC polyvinyl chloride

PZ piezometer

QAPP Quality Assurance Project Plan SOP standard operating procedure

TCE trichloroethene

UST underground storage tank
VOC volatile organic compound

WBZ water-bearing zone

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# Introduction

CH2M HILL Engineers, Inc. (CH2M) prepared this Work Plan Addendum on behalf of Essex Specialty Products, Inc. (ESP) to present proposed activities to supplement previous investigations for refining the conceptual site model (CSM) and planning for future remedial activities at the Essex-Hope State Superfund site located at 125 Blackstone Avenue in Jamestown, New York (site; Figure 1). The site, presently owned by Custom Production Manufacturing Inc., is listed under the New York Superfund Program (Site Number 907015) and managed by the New York State Department of Environmental Conservation (NYSDEC).

The work described herein is intended to supplement the investigation that was detailed in the *Site Characterization Investigation Work Plan, Essex-Hope Site in Jamestown, New York* (2016 Work Plan; CH2M, 2016). Results of this investigation were provided in the *2016 Data Gap Investigation Report for the Essex-Hope Site in Jamestown, New York* (2016 Data Gap Investigation Report; CH2M, 2017a), which is an appendix to the *2016 Annual Periodic Review Report, Essex/Hope Site, Jamestown, New York* (CH2M, 2017b). The 2016 Data Gap Investigation Report provides an updated understanding of hydrogeology as well as the nature and extent of the contaminants. As detailed in the report, further delineation of CBTEX (cumene, benzene, toluene, ethylbenzene, and xylene) in the Shallow Waterbearing Zone (WBZ) and trichloroethene (TCE) and its degradation products within both the Shallow and Deep WBZs is needed. Additionally, light nonaqueous phase liquid (LNAPL) was found during a follow-up sampling event, which requires additional assessment. This addendum to the 2016 Work Plan has been prepared to further refine this information. A full update to the CSM will be prepared upon collection of the results from this investigation. This Work Plan Addendum has been developed in general accordance with the NYSDEC's Division of Environmental Remediation (DER) *Technical Guidance for Site Investigation and Remediation* (DER-10), dated May 2010.

## 1.1 Purpose and Scope

The activities proposed in this Work Plan Addendum are being conducted to meet the following objectives:

- Objective 1 Define the lateral extent of CBTEX in the Shallow WBZ in the eastern Former Underground Storage Tank (UST) Area.
- Objective 2 Define the source and extent of the elevated TCE concentrations in the shallow WBZ near MW-104S on the Hope Windows property.
- Objective 3 Define the extent of the vadose zone soil TCE impacts in the North Parking Lot Sump (NPLS) Area near DPT-005 and MW-111D. Also, define the depth and extent of LNAPL found in PZ-4D.
- Objective 4 Define the extent of the TCE plume in the upper portion of the Deep WBZ to the north and east of MW-106D and MW-110D. Assess whether some of the existing Deep WBZ monitoring wells are screened too deeply to adequately assess and delineate the deep TCE plume.
- Objective 5 Collect and analyze samples to assist in assessment of remedial alternatives.
- Objective 6 Abandon and refurbish selected existing wells and piezometers.

To meet these objectives, the following activities are proposed:

Task 1: Utility Clearance.

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- Task 2: Soil and Groundwater Profiling. Continuous soil sampling at 22 soil borings (Figure 1). Seven shallow borings will be advanced to depths of 15 feet below ground surface (bgs) and an additional 15 deep borings will be advanced to depths of 45 feet bgs. Groundwater grab samples will be collected at 13 locations.
- Task 3: Installation, Development, and Sampling of Monitoring Wells. Installation and development of eight 2-inch monitoring wells (new borings) (Figure 2). Seven monitoring wells will are planned for installation to 45 feet bgs and 1 is planned for 15 feet bgs. Depths provided are estimates and actual depth will depend on field observations. The monitoring wells will be sampled after development. Soil samples will be collected at one location.
- Task 4: Abandonment and Redevelopment of Existing Wells. Abandonment of up to eight
  monitoring wells (Figure 3). Monitoring wells will be abandoned in accordance with NYSDEC CP-43:
  Groundwater Monitoring Well Decommissioning Policy (NYSDEC, 2009). Three existing monitoring
  wells and five recovery wells will be redeveloped, and eight wells will be repaired (Figure 4).
- Task 5: Survey the coordinates and elevations of the newly installed monitoring wells and new boring locations..

Proposed investigation and well abandonment/repair locations are shown on Figures 1 through 4.

## 1.2 Work Plan Organization

This work plan is organized as follows:

- Section 1 Introduction
- Section 2 Investigation Scope of Work
- Section 3 Other Planning Documents
- Section 4 Reporting and Schedule
- Section 5 References
- Appendix A Quality Assurance Project Plan

Supporting material in the form of tables, figures, and the appendix, are presented at the end of this Work Plan Addendum.

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# Investigation Scope of Work

This section presents a detailed scope of work for addressing the data needs detailed in Section 1.1. Locations of the proposed field investigation activities are shown on Figures 1 through 4.

## 2.1 Field Investigation Activities

### 2.1.1 Utility Clearance

At each drilling location, a third-party utility survey will be completed. An area approximately 5-foot-square around each boring as shown on Figure 1 will be evaluated using non-intrusive geophysical utility locator technologies (e.g., magnetic, ground-penetrating radar, or similar) as necessary to identify, mark out, and locate underground utilities and other subsurface obstructions within the survey area. Utilities will be flagged and marked for review prior to drill rig mobilization. CH2M personnel accompanying the utility locating contractor will take photographs of markings and draw rough sketches of the located subsurface utilities in the field logbook.

In addition to the private utility survey/mark out, Dig Safe System, Inc. will be contacted at least three working days and not more than 10 working days before initial mobilization of drilling equipment to the site.

#### 2.1.2 Soil and Groundwater Profiling

A total of 22 soil borings is planned for soil sample collection as part of this task. Depths and information of borings are provided in Table 1. Thirteen locations have been targeted for groundwater profiling. The proposed boring locations are shown on Figure 1. Specifically, the soil and groundwater profiling will include the following:

- Borings will be advanced using a 2-inch-diameter, 4-foot-long Macro-Core® sampler with a basket trap and a clean, disposable, acetate liner. Samples will be collected continuously during the boring and field screened by a CH2M representative using a photoionization detector at 6-inch intervals, in addition to any intervals showing staining or sheens.
- Soil will be logged and field-screened (including for NAPL using a Sudan IV dye test) in accordance
  with the 2016 Work Plan. To assess the potential of NAPL (found in PZ-4D), a Sudan IV dye test will
  be administered at locations DPT-24, -25, -26, and -27 at 6-inch intervals starting at 2-feet bgs, to 2feet below the water table, even if no signs of NAPL are detected.
- Soil samples will be collected for specific analysis dependent on each location as presented in Table 2. Laboratory analytical methods include the following:
  - To be sent to Alpha Analytical, Inc. of Westborough, Massachusetts:
    - Volatile organic compounds (VOCs) using U.S. Environmental Protection Agency (EPA)
       Method 8260C
    - Polychlorinated biphenyls using EPA Method 8082
  - To be sent to Microbial Insights, Inc. of Knoxville, Tennessee:
    - Magnetic susceptibility
  - To be sent to Test America Applied Sciences Laboratory of Corvallis, Oregon:
    - Total oxidant demand using permanganate

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- Total oxidant demand using alkaline-activated persulfate
- During soil sampling activities, groundwater samples will be collected at three wells (see Table
   2) for submittal to Applied Sciences Laboratory for use in the total oxidant demand analyses.
- Up to 25 groundwater grab samples will be taken at 13 boring locations from depths ranging from 15 feet bgs up to 45 feet bgs. Up to three groundwater grab samples may be collected from each boring at different depths. For each groundwater grab, a temporary 5-foot screened interval will be installed at the appropriate depth, as directed by the onsite CH2M representative, and a peristaltic pump and tubing for purging and sampling will be used. Groundwater samples will be sent to Alpha Analytical, Inc. for analysis for VOCs by EPA Method 8260C. Refer to Table 1 for a complete list of locations of groundwater profiling locations and depths of samples.
- Drill cuttings and liners will be containerized as detailed in Section 3.2 of the 2016 Work Plan (CH2M, 2016).
- Drilling equipment will be decontaminated as detailed in Section 3.2 of the 2016 Work Plan.

Depending on field observations, additional soil borings or soil samples to be collected for offsite analysis may be required. Any deviations to the plan noted above will be provided within the final report.

#### 2.1.3 Monitoring Well Installation, Development, and Sampling

#### 2.1.3.1 Monitoring Well Installations

Eight monitoring wells will be installed on the site and at adjacent properties. Six of the monitoring wells will be new wells and two existing monitoring wells will be replaced (MW-7D and PZ-7D). Potential monitoring well installation locations are shown on Figure 2; however, final locations will be selected based on the results of the DPT sampling. Boreholes at least 4 inches in diameter will be advanced to ensure sufficient annular space for well installation. To minimize potential downward migration of contaminants at MW-118D, a temporary outer casing or hollow stem auger flights may be used to seal off the Shallow WBZ. Monitoring well installation procedures will be consistent with the 2016 Work Plan and are briefly summarized below:

- Monitoring wells will be installed to depths ranging from 15 to 45 feet bgs as detailed in Table 3.
   Final depths will be based on field observations. Monitoring wells will be constructed with 2-inch-diameter schedule 40 polyvinyl chloride (PVC). Well materials will consist of 0.010-inch machine-slotted, flush-threaded, schedule 40 PVC screens and riser pipe. Screen lengths will generally be 10 feet unless site conditions indicate a 5-foot screen is required.
- At locations with planned soil sampling, as detailed in Section 2.1.2, monitoring wells will be installed to predetermined depths. Otherwise, continuous samples will be collected during monitoring well installations to determine final depth.
- After setting the well screen, riser, filter pack, and bentonite seal, the well will be grouted as the auger or temporary casing is withdrawn, preventing cross-contamination.
- Drill cuttings and liners will be containerized as detailed in Section 3.2 of the 2016 Work Plan.
- Drilling equipment will be decontaminated as detailed in the Section 3.2 of the 2016 Work Plan.

#### 2.1.3.2 Monitoring Well Development

The newly installed monitoring wells will be developed at least 48 hours after installation in accordance with NYSDEC protocols and the procedures outlined in the 2016 Work Plan.

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#### 2.1.3.3 Groundwater Sampling

Two rounds of groundwater sampling will be performed at all newly installed and replacement wells and will be submitted to Alpha Analytical, Inc. for analysis for VOCs (SW8260), to supplement the annual performance monitoring groundwater sampling. Groundwater samples will be collected in accordance with the *Low-Flow Groundwater Sampling Procedure* (Standard Operating Procedure [SOP]-07; EPA, 1996) and as detailed in Section 3.1.7 of the 2016 Work Plan.

During one groundwater sampling round, four wells will be sampled and submitted to Microbial Insights, Inc. for QuantArray-Chlor analysis to assess the microbial community composition.

#### 2.1.4 Abandonment and Redevelopment of Existing Wells

#### 2.1.4.1 Monitoring Well Abandonment

Eight monitoring wells will be abandoned, as shown on Figure 3, in accordance with NYSDEC CP-43: *Groundwater Monitoring Well Decommissioning Policy* (NYSDEC, 2009). Construction details on the eight wells to be abandoned are provided in Table 4, as well as recent total depth measurements to provide an estimate of sediment thicknesses within the wells. Abandonment methods will comply with NYSDEC policies; wells will be abandoned by grouting the casing:

- Each well will be abandoned by tremie grouting (utilizing neat cement) to 5 feet bgs, removing well vault and casing materials to at least 5 feet bgs, and backfilling the remaining with materials similar to native materials for the upper 5 feet. The surface will be restored to match the condition of the area surrounding the borehole (i.e., an asphalt patch, concrete patch, or grass).
- The grout will conform with the requirements of NYSDEC CP-43, with a standard grout mix: of 1 bag Portland cement, 3.9 pounds of bentonite, and 7.8 gallons of potable water.

#### 2.1.4.2 Monitoring Well Redevelopment

Eight wells will be redeveloped, which includes all of the active recovery wells. Although these tasks are part of general operations and maintenance (O&M) at the site, these activities have been included within this work plan addendum to allow for a complete record. Monitoring/recovery wells to be redeveloped and repaired are shown on Figure 4 with details provided in Table 5. Redevelopment activities will be performed in accordance with NYSDEC protocols. Development water will be brought to the onsite treatment system for disposal.

#### 2.1.5 Survey

Following installation, a licensed surveyor will survey the completed soil borings and monitoring wells. The horizontal and vertical locations for the monitoring well ground surface, steel flush-mount lid, and top of PVC well riser will be surveyed.

## 2.2 Waste Management Plan

Waste will be managed in accordance with Section 3.2 of the 2016 Work Plan (CH2M, 2016).

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# Other Planning Documents

## 3.1 Quality Assurance Project Plan

An update to the previous Quality Assurance Project Plan (QAPP) is provided as Appendix A to this Work Plan Addendum.

## 3.2 Community Air Monitoring Plan

The Community Air Monitoring Plan detailed in the 2016 Work Plan remains relevant. No updates are needed.

## 3.3 Health and Safety Plan

The Health and Safety Plan detailed in the 2016 Work Plan remains relevant. No updates are needed.

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# Reporting and Schedule

## 4.1 Schedule

Schedule milestones have been established based on the assumption that an approved work plan is in place by July 2017. Field activities will not proceed until NYSDEC approves the work plan. In the event of changes in availability of regulatory staff, project approach, or site conditions, a revised schedule will be provided to NYSDEC. The anticipated schedule milestones are as follows:

Date	Milestone
June 2017	Submit work plan to NYSDEC and New York State Department of Health (NYSDOH) for review
July 2017	If necessary, meet with NYSDEC and NYSDOH to go over work plan
July 2017	Obtain NYSDEC approval on work plan
August/September 2017	Mobilize/Complete fieldwork
September 2017	Complete first groundwater sampling event in conjunction with performance monitoring sampling
March 2018	Complete second groundwater sampling event
March/April 2018	Submit report

## 4.2 Deliverables

CH2M will develop a report describing the investigation activities data and results; this report will be similar to the Data Investigation Report and will supplement that submittal. The report will include a summary of the field activities performed and present boring logs, summary tables and figures (including plume maps), provide updates to the CSM, and provide a preliminary identification of potential alternative remedial actions. The report will be submitted to NYSDEC and NYSDOH. Electronic data deliverables will be submitted to NSYDEC with the report.

A meeting with NYSDEC will be set up to discuss the results and proposed path forward.

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## References

CH2M HILL Engineers, Inc. (CH2M). 2016. *Site Characterization Investigation Work Plan, Essex-Hope Site in Jamestown, New York, Site Number 907015*. Prepared for Essex Specialty Products, Inc. July.

CH2M HILL Engineers, Inc. (CH2M). 2017a. 2016 Data Gap Investigation Report for the Essex-Hope Site in Jamestown, New York. Prepared for Essex Specialty Products, Inc. March.

CH2M HILL Engineers, Inc. (CH2M). 2017b. 2016 Annual Periodic Review Report, Essex/Hope Site, Jamestown, New York. Prepared for Essex Specialty Products, Inc. March.

New York State Department of Environmental Conservation (NYSDEC). 2009. *CP-43: Groundwater Monitoring Well Decommissioning Policy*.

New York State Department of Environmental Conservation (NYSDEC). 2010. *DER-10 Technical Guidance for Site Investigation and Remediation*. http://www.dec.ny.gov/docs/remediation\_hudson\_pdf/der10.pdf. May.

U.S. Environmental Protection Agency (EPA). 1996. Low-flow (minimal drawdown) ground-water sampling procedures. Office of Solid Waste and Emergency Response, EPA/540/S-95/504. April.

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Tables

#### Table 1. Soil Boring Information

Essex-Hope Site, Jamestown, New York

Sample Location	<b>Proposed Final Depth (ft</b>	<b>GW Profiling Depth</b>	
Name	bgs)	(ft bgs)	Rationale
DPT-05A	15	NA	Determine extent of TCE impacts observed at DPT-05. Quick turnaround to determine if
DPT-05B	15	NA	need to step out.
DPT-05C	15	NA	
DPT-05D	15	NA	
DDT 40	45	22-26	Assess if impacts in upper Deep WBZ, not intersected by well screen of MW-14D, exist.
DPT-10	45	32-36	
		10-15	Shallow: Assess level of impacts between MW-101S and 104S. Deep: Assess if impacts at
DPT-11	45	22-26	MW-15D are continuous with MW-110D (connected plume) or if location part of capture
DPT-II	45	32-36	zone/stagnation zone like MW-22D. Collect deep soil sample for NOD (permanganate).
		10-15	Assess northern extent of plume from MW-104S to west. Assess if any soil impacts from
DPT-12	45	22-26	offsite source near MW-104S.
		32-36	
		10-15	Assess northern extent of shallow and deep plume; quick turnaround on deep GW
DPT-13	45	22-26	analytical to determine if a deep well will be installed. Deep well may be installed if clean
		32-36	stepout to north (DPT-16, DPT-17, DPT-18) if impacted.
		22-26	Assess northern extent of deep plume – quick turnaround on analytical to determine if a
DPT-14	45	32-36	well will be installed. Deep well may be installed if clean; stepout to north (DPT-16, DPT-
		32 30	17, DPT-18) if impacted.
		22-26	Assess northern extent of deep plume – quick turnaround on analytical to determine if a
DPT-15	45	32-36	well will be installed. Well may be installed if clean; stepout to north (DPT-16, DPT-17,
			DPT-18) if impacted.
DPT-16	45	22-26	Install if DPT-13, 14, and 15 indicate impacts; quick turnaround on analytical results – we
DF 1-10	43	32-36	may be installed.
DPT-17	45	22-26	
DF 1-17	45	32-36	
DPT-18	45	22-26	
DI 1 10	45	32-36	
			Assess if impacts in upper Deep WBZ, not intersected by well screen MW-25D, exist.
DPT-19	45	22-26	Assess magnetic susceptibility of aquifer materials within shallow TCE plume.
DPT-20	15	10-15	Delineate CBTEX impacts observed around MW-26S
DPT-21	15	10-15	Defined to GBTEX impacts observed around NIW 203
DPT-22	15	10-15	
			Assess MNA amenability and oxidant demand (persulfate) in deep water bearing zone
DPT-23	45	NA	with high TCE and acetone concentrations
DPT-24	45	NA	Evalute extent of NAPL observed at PZ-4D.
DPT-25	45	NA	
DPT-26	45	NA	
DPT-27	45	NA	
MW-123D	45	NA	Assess vadose zone conditions in northern portion of North Parking Lot;
			Assess MNA amenability and oxidant demand (permanganate) in Deep WBZ with high
otes:			TCE concentrations.
			Tee concentrations.

ft bgs = feet below ground surface

All broundwater profiling samples to be analyzed for VOCs using EPA Method 8260.

DPT = direct-push technology

WBZ = water-bearing zone

TCE = trichloroethene

NOD = natural oxidant demand

MNA = monitored natural attenuation

CBTEX = cumene, benzene, toluene, ethylenzene, and xylene

PZ = piezometer

#### Table 2. Soil Sample Summary

Essex-Hope Site, Jamestown, New York

Proposed	Proposed Soil Sample Depth	VOCs	PCBs	Magnetic	Total Oxidant
Location Name	(ft bgs)	(SW8260)	(SW8082)	Susceptibillity	Demand
DPT-11 (MW- 110D)	30-35				Х
DPT-12	2-4 6-8	Х			
DPT-19	15-20			Х	
DPT-20	2-4 6-8	X			
DPT-21	2-4 6-8	X			
DPT-22	2-4 6-8	X			
DPT-05A	2-4 6-8	X	Х		
DPT-05B	2-4 6-8	Х	х		
DPT-05C	2-4 6-8	Х	х		
DPT-05D	2-4 6-8	Х	Х		
DPT-23 (MW- 021D)	25-35			Х	Х
DPT-24		Χ			
DPT-25	Continous based on NAPL	Χ			
DPT-26	observations	X			
DPT-27		Χ			
MW-123D	2-4 6-8	Х	Х		
	25-35			X	X

#### Note:

Additional VOC analysis may be performed at other locations not noted above based on field observations.

ft bgs = feet below ground surface

NAPL = nonaqueous phase liquid

PCB = polychlorinated biphenyl

VOC = volatile organic compound

#### Table 3. Monitoring Well Construction Details

Essex-Hope Site, Jamestown, New York

		Screen		Top of	<b>Bottom of</b>	Well
Well_ID	Material	Type/Slot	Filter Pack	Screen (ft	Screen (ft	Diameter
		Size (in.)		bgs)	bgs)	(in.)
MW-109D	Sched. 40 PVC	PVC/0.01	20/40 sieve size	35	45	2
IVIVV-109D	flush-threaded	PVC/0.01	(or equivalent)	33	43	2
MW-118S	Sched. 40 PVC	PVC/0.01	20/40 sieve size	5	15	2
IVIVV-1103	flush-threaded	PVC/0.01	(or equivalent)	3	15	2
MW-118D	Sched. 40 PVC	PVC/0.01	20/40 sieve size	35	45	2
INIM-TIRD	flush-threaded	PVC/0.01	(or equivalent)	35	43	۷
MW-119D	Sched. 40 PVC	PVC/0.01	20/40 sieve size	35	45	2
IVIVV-119D	flush-threaded	PVC/0.01	(or equivalent)	33	45	۷
MW-120D	Sched. 40 PVC	PVC/0.01	20/40 sieve size	35	45	2
IVIVV-120D	flush-threaded	PVC/0.01	(or equivalent)	33	43	2
MW-121D	Sched. 40 PVC	PVC/0.01	20/40 sieve size	35	45	2
IVIVV-121D	flush-threaded	PVC/0.01	(or equivalent)	33	43	2
MW-122D	Sched. 40 PVC	DVC /0.01	20/40 sieve size	25	45	2
IVIVV-122D	flush-threaded	PVC/0.01	(or equivalent)	35	45	2
MW-123D	Sched. 40 PVC	DVC/0.01	20/40 sieve size	35	45	2
IVIVV-123D	flush-threaded	PVC/0.01	(or equivalent)	33	45	2

ft bgs = feet below ground surface

in. = inch

PVC = polyvinyl chloride

#### Table 4. Wells to Abandon

Essex-Hope Site, Jamestown, New York

Well	Recommendation		Depth to Bottom of Screen (ft bgs)	Reported Depth to Bottom	Well Seal Depths (ft bgs)	Borehole Diam. (in.)	Casing Diameter (in.)	Screen/Riser Type	Outer Casing	Type of Completion	Well Vault	Identified Issue
HW-1	Abandon – Perforate and Grout, Pull or Overdrill	15	20	N/A	Unknown	Unknown	2	PVC	No	Stick up	None	Stick up destroyed, open well
MW-2	Abandon – Grout	9	14	< 9	6.5-8	4.25	2	Stainless Steel	No	Flush mount	Stainless Steel	Filled with sediment above water table
MW-4	Abandon – Grout	13	18	<13	7-10	4.25	2	Stainless Steel	No	Flush mount	Stainless Steel	Filled with sediment above water table
MW-7D	Abandon – Grout, Replace	35	45	19.5	32-33	8	2	PVC	No	Flush mount	Steel	> 20' of sediment, sediment above well screen but within deep WBZ
MW-8	Abandon – Grout	39.6	49.6	Obstruction at 30 and 40 ft bgs	34-35	8	2	PVC	No	Flush mount	Steel	Obstructions at 30 ft bgs and 40 ft bgs.
PZ-7D	Abandon – Grout, Replace	22.0	42.0	33.6	14.5-16.75	4	1	PVC	3 inch PVC, 0-18 ft	Flush mount	Steel	> 8' of sediment
PZ-11D	Abandon – Grout	21.3	41.3	15.8	19-21	2.25	1	PVC	3 inch PVC, 0-24 ft	Flush mount	Steel	> 25' of sediment; sediment above screened interval and confining clay but within sealed outer casing
VP-5D	Abandon – Grout	12.5	34.3	13.0	8-11.5	10 inch to 12.5 ft 4 inch 12.5 to 38 ft bgs	2	PVC	5 inch steel, 0-12.5 ft bgs	Flush mount	Steel	> 20' of sediment, but sediment within screened interval

ft bgs = feet below ground surface

in. = inch

PVC = polyvinyl chloride

#### Table 5. Wells to Repair and/or Redevelop

Essex-Hope Site, Jamestown, New York

Well	Recommendation				
MW-1	Cut down stick up well past "pinch point" and				
IVIVV-I	make flush mount well; redevelop				
MW-13	Ensure tight-fitting well plug; repair well vault,				
10100-13	redevelop				
VP-6D	Repair well vault; redevelop				
RW-1S	Repair well vault; Redevelop				
RW-2S	Recovery Well Redevelop				
RW-2D	Recovery Well Redevelop				
RW-3S	Recovery Well Redevelop				
RW-6D	Recovery Well Redevelop				
GP-5D	Replace well vault lid and well plug				
MW-6	Replace well vault lid and well plug				
MW-7S	Replace well vault lid and well plug				
MW-11	Replace well vault lid and well plug				
MW-19D	Replace well vault lid and well plug				

Figures

VP-5D

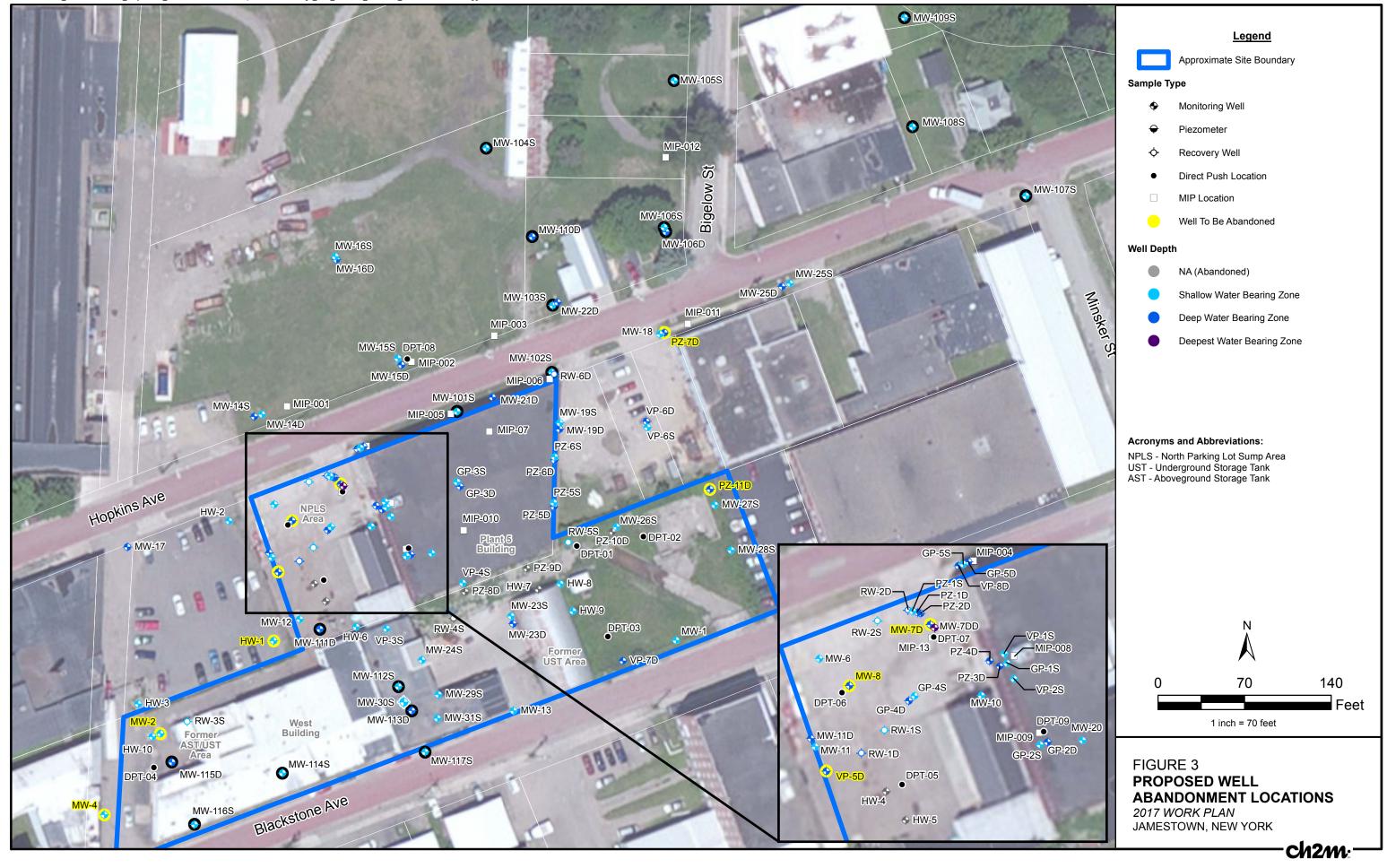
DPT-05

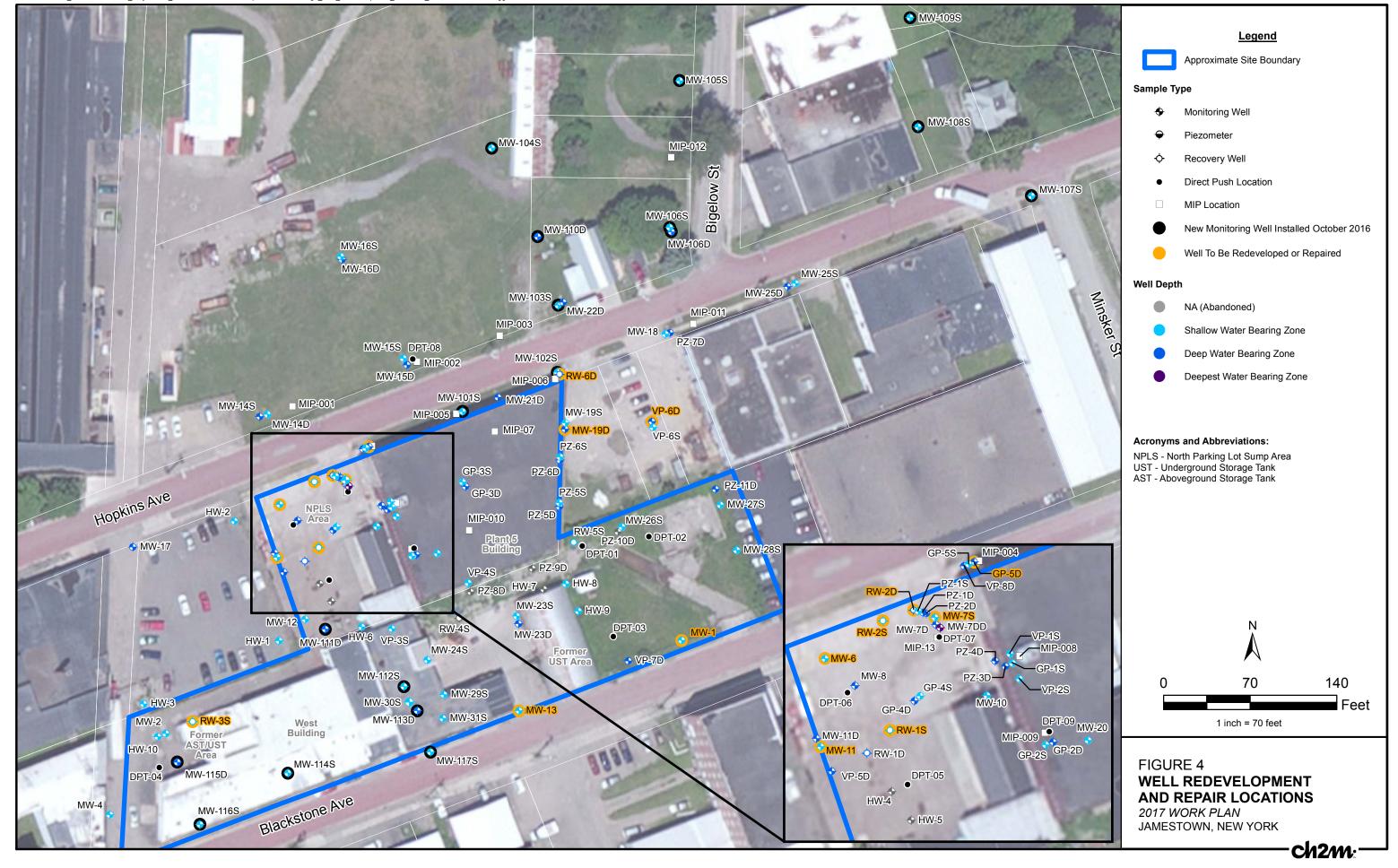
HW-4

FIGURE 1 PROPOSED DPT LOCATIONS 2017 WORK PLAN JAMESTOWN, NEW YORK

ch2m:

-ch2/m:-





Appendix A Quality Assurance Project Plan Addendum

## Quality Assurance Project Plan Addendum

# Site Characterization Investigation Activities

Essex-Hope Site, Jamestown, New York

Essex Specialty Products, Inc.

June 2017



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### **QUALITY ASSURANCE PROJECT PLAN ADDENDUM**

Essex Specialty Products, Inc., Essex-Hope Site

Jamestown, New York

# Site Characterization Investigation Activities June 2017

Rev. 0

Prepared by: CH2M HILL	Date: June 2017
Approved by:	
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## Introduction

This Quality Assurance Project Plan Addendum (QAPP Addendum) presents revisions to various sections of the Quality Assurance Project Plan, Site Characterization Investigation Activities, Essex-Hope Site, Jamestown, New York (CH2M HILL 2016) (QAPP) and presents additions and/or modifications to the quality assurance (QA)/quality control (QC) requirements specified in the QAPP (CH2M HILL 2016). The additions and/or modifications are necessary for current site investigation activities being conducted at The Essex Specialty Products, Inc. (ESP) Essex-Hope State Superfund site in Jamestown, New York.

The section numbers presented in this QAPP Addendum correspond to the pertinent sections in the QAPP (CH2M HILL 2016). All other sections of the QAPP (CH2M HILL 2016) will be followed by the project team.

## 1.1 Project Objectives and Background

The project objectives and background are provided in detail in Section 2 of the Site Characterization Investigation Work Plan (CH2M Hill 2016) (Work Plan).

## 1.3 Project Task/Organization

The investigation SOW and related project tasks are provided in detail in Sections 1.1 and 3 of the Work Plan.

## 3.3 Sampling Process Design and Rationale

Project-specific sampling plans will be developed following the logical process provided in USEPA's *Guidance* on Systemic Planning using the Data Quality Objective Process (USEPA 2006). The sampling design is a function of the matrix sampled, information about the sampling site, the type of data to be collected, and how the data are to be used. These sampling plans will be incorporated in the Work Plan.

## 3.4 Sample Handling and Custody Requirements

Laboratories will provide the required sample containers for all environmental and associated QC samples. Containers will be certified free of the analytes of concern for this project. No sample containers will be reused. The contract laboratory will add preservatives, if required, before shipping the sample containers to the field. Upon receipt of the samples, the laboratory will verify the adequacy of the preservation and add additional preservatives if necessary. Adjustments made by the laboratory will be documented on the appropriate sample receipt forms and noted in the case narrative. The analytical methods, sample containers, minimum quantities, required preservatives, and maximum holding times for select parameters are shown in Table 3-1. Extraction and preparatory methods are shown in Table 3-2. Tables are located at the end of this document.

Sample custody procedures will be followed as outlined in Section 3.5 of the QAPP (CH2M Hill 2016).

## 3.6 Analytical Method Requirements

Analytical methods will be performed in accordance with the QAPP (CH2M Hill 2016) and this QAPP Addendum, and will be reported as definitive data results. The following analytical methods will be used:

SW8260C; Volatile organic compounds (VOCs) in groundwater/soil

In addition, the following methods, which are not included in the QAPP (CH2M Hill 2016), will be performed:

- SW8082; Polychlorinated Biphenyls (PCB) in soil
- Quantarray-Chlor; qPCR<sup>1</sup> in groundwater
- Magnetic Suspectibility in soil
- Total Oxidant Demand in soil

Target analyte lists and RL objectives for each method are specified in Attachment A. Requested RLs will be based on meeting New York State Department of Environmental Conservation (NYSDEC) unrestricted use soil cleanup objectives or water quality standards.

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<sup>&</sup>lt;sup>1</sup> qPCR = quantitative polymerase chain reaction

## 3.8 Quality Control Requirements

The QC samples to be included in the site investigation activities are outlined in Section 3.8 of the QAPP (CH2M Hill 2016).

Required QC checks, minimum frequencies, acceptance criteria, and corrective actions are included in Attachment A, when applicable. No specific QC checks are required for Quantarray-Chlor, magnetic suspectibility, or total oxidant demand.

#### 5.2 Data Verification and Validation Methods

All analytical results of the data collection effort associated with the site investigation activities will be validated by CH2M. All data for this program will undergo **Level 3** validation unless otherwise noted as specified in Section 5.2.3 of the QAPP (CH2M Hill 2016).

# References

CH2M HILL (CH2M). 2016. Site Characterization Investigation Work Plan, Essex-Hope Site in Jamestown, New York. June.

CH2M HILL (CH2M). 2016. *Quality Assurance Project Plan, Site Characterization Investigation Activities, Essex-Hope Site in Jamestown, New York*. June.

# **Tables**

Table 3-1. Required Analytical Method, Sample Containers, Preservation, and Holding Times Essex-Hope Site, Jamestown, New York

Analyses	Analytical Method	Sample Matrix <sup>a</sup>	Container <sup>b</sup>	Qty	Preservative <sup>c</sup>	Holding Time <sup>d</sup>
Volatile Organic Compounds	SW-846 8260C	W	40-mL, glass	3	HCl, pH<2; cool to 4°C	14 days
		S	5 g–Encore or equivalent sampling technique	3	Cool 4°C	48 hours from collection to preservation, 14 days to analysis
Polychlorinated Biphenyls	SW-846 8082	S	4-oz glass	1	Cool 4°C	14 days to extraction; 40 days from extraction to analysis
Quantarray-Chlor	Lab SOP	W	1-L, poly	1	Cool 4°C	24 – 48hours
Magnetic Suspectibility	Lab SOP	S	TBD	1	None	None
Total Oxidant Demand	Lab SOP	S	Large Ziplock Bag containing 2 kg soil	1	Cool 4°C	None

#### Notes:

Sample container, and volume requirements will be specified by the analytical laboratory performing the tests.

Three times the required volume should be collected for samples designated as MS/MSD samples.

g = Grams

mL = Milliliter

kg = Kilograms

HCl = Hydrochloric acid

oz = Ounce

TBD = to be determined

Table 3-2. Extraction and Digestion Methods

Essex-Hope Site, Jamestown, New York

ESSEX Trope Site, Sumestown, New	TOTA	
Analytical Method	Parameter	Preparatory Methods
SW8082	Polychlorinated Biphenyls (water and soil)	SW3010C, SW3050C
SW8260C	Volatile Organic Compounds (water and soil)	SW5030C, SW5035
Quantarray-Chlor	qPCR (water)	See analytical method
Lab SOP	Magnetic Suspectibility	See analytical method
Lab SOP	Total Oxidant Demand	See analytical method

<sup>&</sup>lt;sup>a</sup> Sample matrix: S = surface soil, subsurface soil; W = surface water

<sup>&</sup>lt;sup>b</sup> All containers will be sealed with Teflon-lined screw caps.

<sup>&</sup>lt;sup>c</sup> All samples will be stored promptly at 4°C in an insulated chest.

<sup>&</sup>lt;sup>d</sup> Holding times are from the time of sample collection unless otherwise specified.

<sup>°</sup>C = Degrees Celsius



Table A-1. Reporting Limit Objectives for Volatile Organic Compounds in Groundwater by Method SW8260C *Essex-Hope Site, Jamestown, New York* 

Parameter	CAS Number	Site-Specific Remedial Action Objective	NYSDEC Water Quality Standards <sup>a</sup> (μg/L)	Method Detection Limit (μg/L)	Reporting Limit (µg/L)
1,1,1,2-Tetrachloroethane	630-20-6			0.7	2.5
1,1,1-Trichloroethane	71-55-6		5.0	0.7	2.5
1,1,2,2-Tetrachloroethane	79-34-5		5.0	0.14	0.5
1,1,2-Trichloroethane	79-00-5		1.0	0.5	1.5
1,1-Dichloroethane	75-34-3		5.0	0.7	2.5
1,1-Dichloroethene	75-35-4		5.0	0.14	0.5
1,1-Dichloropropene	563-58-6			0.7	2.5
1,2,3-Trichlorobenzene	87-61-6		5.0	0.7	2.5
1,2,3-Trichloropropane	96-18-4			0.7	2.5
1,2,4-Trichlorobenzene	120-82-1		5.0	0.7	2.5
1,2,4-Trimethylbenzene	95-63-6			0.7	2.5
1,2-Dibromo-3-chloropropane	96-12-8		0.04	0.7	2.5
1,2-Dibromoethane	106-93-4		0.0006	0.65	2.0
1,2-Dichlorobenzene	95-50-1		3.0	0.7	2.5
1,2-Dichloroethane	107-06-2		0.6	0.13	0.5
1,2-Dichloropropane	78-87-5		1.0	0.13	1.0
1,3,5-Trimethylbenzene	108-67-8			0.7	2.5
1,3-Dichlorobenzene	541-73-1		3.0	0.7	2.5
1,3-Dichloropropane	142-28-9			0.7	2.5
1,4-Dichlorobenzene	106-46-7		3.0	0.7	2.5
1,4-Dioxane	123-91-1			41	250
2,2-Dichloropropane	594-20-7			0.7	2.5
2-Butanone	78-93-3		50	1.9	5.0
2-Chlorotoluene	95-49-8			0.7	2.5
2-Hexanone	591-78-6		50	1.0	5.0
4-Chlorotoluene	106-43-4			0.7	2.5
4-Methyl-2-pentanone	108-10-1			1.0	5.0
Acetone	67-64-1		50	1.5	5.0
Benzene	71-43-2		1.0	0.16	0.5
Bromobenzene	108-86-1			0.7	2.5
Bromochloromethane	74-97-5		5.0	0.7	2.5
Bromodichloromethane	75-27-4		50	0.19	0.5
Bromoform	75-25-2		50	0.65	2.0
Bromomethane	74-83-9		5.0	0.7	2.5
Carbon disulfide	75-15-0		60	1.0	5.0
Carbon tetrachloride	56-23-5		5.0	0.13	0.5
Chlorobenzene	108-90-7		5.0	0.7	2.5
Chloroethane	75-00-3		5.0	0.7	2.5
Chloroform	67-66-3		7.0	0.7	2.5
Chloromethane	74-87-3			0.7	2.5

Table A-1. Reporting Limit Objectives for Volatile Organic Compounds in Groundwater by Method SW8260C *Essex-Hope Site, Jamestown, New York* 

Lassex-Hope Site, Juliestowii, New		Site-Specific Remedial Action	NYSDEC Water Quality Standards <sup>a</sup>	Method Detection Limit	Reporting Limit
Parameter	CAS Number	Objective	(μg/L)	(μg/L)	(μg/L)
cis-1,2-Dichloroethene	156-59-2		5.0	0.7	2.5
cis-1,3-Dichloropropene	10061-01-5		0.4	0.14	0.5
Chlorodibromomethane	124-48-1		50	0.15	0.5
Dibromomethane	74-95-3			1.0	5.0
Dichlorodifluoromethane	75-71-8		5.0	1.0	5.0
Ethylbenzene	100-41-4	5.0	5.0	0.7	2.5
Hexachlorobutadiene	87-68-3			0.7	2.5
Isopropylbenzene	98-82-8		5.0	0.7	2.5
m-,p-Xylene	179601-23-1	5.0	5.0	0.7	2.5
Methylene chloride	75-09-2		5.0	0.7	2.5
Methyl t-Butyl Ether (MTBE)	1634-04-4			0.7	2.5
n-Butylbenzene	104-51-8			0.7	2.5
n-Propylbenzene	103-65-1			0.7	2.5
Naphthalene	91-20-3			0.7	2.5
o-Xylene	95-47-6	5.0	5.0	0.7	2.5
p-Isopropyltoluene	99-87-6			0.7	2.5
sec-Butylbenzene	135-98-8			0.7	2.5
Styrene	100-42-5		5.0	0.7	2.5
tert-Butylbenzene	98-06-6			0.7	2.5
Tetrachloroethene	127-18-4		5.0	0.18	0.5
Toluene	108-88-3	5.0	5.0	0.7	2.5
trans-1,2-Dichloroethene	156-60-5	5.0	5.0	0.7	2.5
trans-1,3-Dichloropropene	10061-02-6		0.4	0.16	0.5
Trichloroethene	79-01-6	5.9	5.0	0.18	0.5
Trichlorofluoromethane	75-69-4		5.0	0.7	2.5
Vinyl acetate	108-05-4			1.0	5.0
Vinyl chloride	75-01-4	5.0	2.0	0.07	1

 $<sup>^{\</sup>rm a}$  NYSDEC Ambient Water Quality Standards and Guidance Values (1998) and 2000 Addendum  $\mu g/L$  – micrograms per liter

Table A-2. Reporting Limit Objectives for Volatile Organic Compounds in Soils by Method SW8260C *Essex-Hope Site, Jamestown, New York* 

Parameter	CAS Number	Site Specific Remedial Action Objective	NYSDEC Unrestricted Use SCO <sup>a</sup> Screening Criteria (µg/kg)	Method Detection Limit (μg/kg)	Reporting Limit (μg/kg)
1,1,1,2-Tetrachloroethane	630-20-6	1,000		0.32	1.0
1,1,1-Trichloroethane	71-55-6	1,000	680	0.11	1.0
1,1,2,2-Tetrachloroethane	79-34-5	1,000		0.10	1.0
1,1,2-Trichloroethane	79-00-5	1,000		0.30	1.5
1,1-Dichloroethane	75-34-3	1,000	270	0.086	1.5
1,1-Dichloroethene	75-35-4	1,000	330	0.26	1.0
1,1-Dichloropropene	563-58-6	1,000		0.14	5.0
1,2,3-Trichlorobenzene	87-61-6	1,000		0.15	5.0
1,2,3-Trichloropropane	96-18-4	1,000		0.16	10
1,2,4-Trichlorobenzene	120-82-1	1,000		0.18	2.0
1,2,4-Trimethylbenzene	95-63-6	1,000	3,600	0.14	5.0
1,2-Dibromo-3-chloropropane	96-12-8	1,000		0.40	5.0
1,2-Dibromoethane	106-93-4	1,000		0.17	4.0
1,2-Dichlorobenzene	95-50-1	1,000	1,100	0.15	5.0
1,2-Dichloroethane	107-06-2	1,000	20	0.11	1.0
1,2-Dichloropropane	78-87-5	1,000		0.23	3.5
1,3,5-Trimethylbenzene	108-67-8	1,000	8,400	0.14	5.0
1,3-Dichlorobenzene	541-73-1	1,000	2,400	0.14	5.0
1,3-Dichloropropane	142-28-9	1,000		0.15	5.0
1,4-Dichlorobenzene	106-46-7	1,000	1,800	0.14	5.0
1,4-Dioxane	123-91-1	1,000	100	14	100
2,2-Dichloropropane	594-20-7	1,000		0.23	5.0
2-Butanone	78-93-3	1,000	120	0.27	10
2-Chlorotoluene	95-49-8	1,000		0.16	5.0
2-Hexanone	591-78-6	1,000		0.67	10
4-Chlorotoluene	106-43-4	1,000		0.13	5.0
4-Methyl-2-pentanone	108-10-1	1,000		0.24	10
Acetone	67-64-1	1,000	50	1.0	10
Benzene	71-43-2	1,000	60	0.12	1.0
Bromobenzene	108-86-1	1,000		0.21	5.0
Bromochloromethane	74-97-5	1,000		0.28	5.0
Bromodichloromethane	75-27-4	1,000		0.17	1.0
Bromoform	75-25-2	1,000		0.24	4.0
Bromomethane	74-83-9	1,000		0.34	2.0
Carbon disulfide	75-15-0	1,000		1.1	10
Carbon tetrachloride	56-23-5	1,000	760	0.21	1.0
Chlorobenzene	108-90-7	1,000	1,100	0.35	1.0
Chloroethane	75-00-3	1,000		0.32	2.0
Chloroform	67-66-3	1,000	370	0.37	1.5

Table A-2. Reporting Limit Objectives for Volatile Organic Compounds in Soils by Method SW8260C Essex-Hope Site, Jamestown, New York

Parameter	CAS Number	Site Specific Remedial Action Objective	NYSDEC Unrestricted Use SCO® Screening Criteria (µg/kg)	Method Detection Limit (μg/kg)	Reporting Limit (μg/kg)
Chloromethane	74-87-3	1,000		0.29	5.0
cis-1,2-Dichloroethene	156-59-2	1,000	250	0.14	1.0
cis-1,3-Dichloropropene	10061-01-5	1,000		0.12	1.0
Chlorodibromomethane	124-48-1	1,000		0.15	1.0
Dibromomethane	74-95-3	1,000		0.16	10
Dichlorodifluoromethane	75-71-8	1,000		0.19	10
Ethylbenzene	100-41-4	1,000	1,000	0.13	1.0
Hexachlorobutadiene	87-68-3	1,000		0.23	5.0
Isopropylbenzene	98-82-8	1,000	2,300	0.10	1.0
m-,p-Xylene	179601-23-1	1,000	260	0.20	2.0
Methylene chloride	75-09-2	1,000	50	1.1	10
Methyl t-Butyl Ether (MTBE)	1634-04-4	1,000	930	0.084	2.0
n-Butylbenzene	104-51-8	1,000	12,000	0.11	1.0
n-Propylbenzene	103-65-1	1,000	3,900	0.11	1.0
Naphthalene	91-20-3	1,000	12,000	0.14	5.0
o-Xylene	95-47-6	1,000	260	0.17	2.0
p-Isopropyltoluene	99-87-6	1,000	10,000	0.13	1.0
sec-Butylbenzene	135-98-8	1,000	11,000	0.12	1.0
Styrene	100-42-5	1,000		0.40	2.0
tert-Butylbenzene	98-06-6	1,000	5,900	0.14	5.0
Tetrachloroethene	127-18-4	1,000	1,300	0.14	1.0
Toluene	108-88-3	1,000	700	0.19	1.5
trans-1,2-Dichloroethene	156-60-5	1,000	190	0.21	1.5
trans-1,3-Dichloropropene	10061-02-6	1,000		0.12	1.0
Trichloroethene	79-01-6	1,000	470	0.13	1.0
Trichlorofluoromethane	75-69-4	1,000		0.39	5.0
Vinyl acetate	108-05-4	1,000		0.13	10
Vinyl chloride	75-01-4	1,000	20	0.12	2.0

 $<sup>^{</sup>a}$  NYSDEC Soil Cleanup Objectives (SCO) for Unrestricted Use (NYSDEC 2006, 2010) are presented for selecting appropriate Method Detection Limits and Reporting Limits. Site-specific RAOs were presented in the 1994 ROD and are the site cleanup goals  $\mu g/kg - micrograms$  per kilogram

Table A-3. Reporting Limit Objectives for Polychlorinated Biphenyls in Soil by Method SW8082 Essex-Hope Site, Jamestown, New York

Analyte	NYSDEC Unrestricted Use SCO <sup>a</sup> Screening Criteria (μg/kg)	Method Detection Limits (µg/kg)	Reporting Limits (μg/kg)
Aroclor 1016	100	2.6	33.5
Aroclor 1221	100	3.1	33.5
Aroclor 1232	100	3.9	33.5
Aroclor 1242	100	4.1	33.5
Aroclor 1248	100	2.8	33.5
Aroclor 1254	100	2.8	33.5
Aroclor 1260	100	2.6	33.5
Aroclor 1262	100	1.7	33.5
Aroclor 1268	100	4.9	33.5

 $<sup>^{</sup>a}$  NYSDEC Soil Cleanup Objectives (SCO) for Unrestricted Use (NYSDEC 2006 and 2010) are presented for selecting appropriate Method Detection Limits and Reporting Limits. Site-specific RAOs were presented in the 1994 ROD and are the site cleanup goals  $\mu g/kg - micrograms$  per kilogram

Table A-4. Reporting Limit Objectives for Other Parameters in Groundwater/Soil by Various Methods Essex-Hope Site, Jamestown, New York

Analyte	Method	NYSDEC Water Quality Standards <sup>a</sup>	NYSDEC Unrestricted Use SCO <sup>b</sup> Screening Criteria	Reporting Limits
Dehalococcoides spp. (DHC)	QuantArray-Chlor			4.2 cells/mL
tceA Reductase (TCE)	QuantArray-Chlor			4.2 cells/mL
BAV1 Vinyl Chloride Reductase (BVC)	QuantArray-Chlor			4.2 cells/mL
Vinyl Chloride Reductase (VCR)	QuantArray-Chlor			4.2 cells/mL
Dehalobacter sp. (DHBt)	QuantArray-Chlor			4.2 cells/mL
Dehalobacter DCM (DCM)	QuantArray-Chlor			4.2 cells/mL
Dehalogenimonas spp. (DHG)	QuantArray-Chlor			4.2 cells/mL
Desulfitobacterium spp (DSB)	QuantArray-Chlor			4.2 cells/mL
Dehalobium chlorocoercia (DECO)	QuantArray-Chlor			4.2 cells/mL
Desulfuromonas spp. (DSM)	QuantArray-Chlor			4.2 cells/mL
Chloroform reductase (CFR)	QuantArray-Chlor			4.2 cells/mL
1,1-DCA Reductase (DCA)	QuantArray-Chlor			4.2 cells/mL
1,2-DCA Reductase (DCAR)	QuantArray-Chlor			4.2 cells/mL
Soluble Methane Monooxygenase (SMMO)	QuantArray-Chlor			4.2 cells/mL
Particulate Methane Monooxygenase (PMMO)	QuantArray-Chlor			4.2 cells/mL
Toluene Dioxygenase (TOD)	QuantArray-Chlor			4.2 cells/mL
Phenol Hydroxylase (PHE)	QuantArray-Chlor			4.2 cells/mL
Trichlorobenzene Dioxygenase (TCBO)	QuantArray-Chlor			4.2 cells/mL
Toluene Monooxygenase 2 (RDEG)	QuantArray-Chlor			4.2 cells/mL
Toluene Monooxygenase (RMO)	QuantArray-Chlor			4.2 cells/mL
Ethene Monooxygenase (EtnC)	QuantArray-Chlor			4.2 cells/mL
Epoxyalkane transferase (EtnE)	QuantArray-Chlor			4.2 cells/mL
Dichlormethane dehalogenase (DCMA)	QuantArray-Chlor			4.2 cells/mL
Total Eubacteria (EBAC)	QuantArray-Chlor			4.2 cells/mL
Sulfate Reducing Bacteria (APS)	QuantArray-Chlor			4.2 cells/mL
Methanogens (MGN)	QuantArray-Chlor			4.2 cells/mL
Total Oxidant Demand	Lab SOP			0.5 g/Kg

<sup>&</sup>lt;sup>a</sup> NYSDEC Water Quality Standards and Guidance Values (1998)

<sup>&</sup>lt;sup>b</sup> NYSDEC Soil Cleanup Objectives (SCO) for Unrestricted Use (NYSDEC 2006 and 2010) are presented for selecting appropriate Method Detection Limits and Reporting Limits.

g – grams

Kg – kilograms mL – milliliter

Table A-5. Accuracy and Precision for Volatile Organic Compounds by Method SW8260C *Essex-Hope Site, Jamestown, New York* 

Parameter   CL   UC   (KRPD)   CL   UC   CL   CL   CL   CL   CL   CL		Acc W	D/MS/MSD curacy /ater %R)	Precision	LCS/LCSD/MS/MSD Accuracy Soil (%R)		Precision
1,1,1-Trichloroethane         67         130         \$ 20         70         130         \$ 30           1,1,2,2-Tertachloroethane         67         130         \$ 20         70         130         \$ 30           1,1,2-Trichloroethane         70         130         \$ 20         65         135         \$ 30           1,1-Dichloroethane         61         145         \$ 20         70         130         \$ 30           1,1-Dichloropropene         70         130         \$ 20         70         130         \$ 30           1,2,3-Trichlorobenzene         70         130         \$ 20         70         130         \$ 30           1,2,3-Trichloropopane         64         130         \$ 20         70         130         \$ 30           1,2,4-Trichlorobenzene         70         130         \$ 20         70         130         \$ 30           1,2,4-Trichlorobenzene         70         130         \$ 20         70         130         \$ 30           1,2-Dichromo-3-chloropropane         41         144         \$ 20         68         130         \$ 30           1,2-Dichromo-3-chloropropane         40         130         \$ 20         70         130         \$ 30	Parameter		<u> </u>	_ Water (%RPD)		•	Soil (%RPD)
1,1,2,2-Tetrachloroethane         67         130         ≤ 20         70         130         ≤ 30           1,1,2-Trichloroethane         70         130         ≤ 20         70         130         ≤ 30           1,1-Dichloroethane         70         130         ≤ 20         70         130         ≤ 30           1,1-Dichloroethene         61         145         ≤ 20         70         130         ≤ 30           1,2,3-Trichloroperopene         70         130         ≤ 20         70         130         ≤ 30           1,2,3-Trichloroperopane         64         130         ≤ 20         70         130         ≤ 30           1,2,4-Trinchloroperopane         64         130         ≤ 20         70         130         ≤ 30           1,2,4-Trinchloroperopane         70         130         ≤ 20         70         130         ≤ 30           1,2-Dichloroperopane         41         144         ≤ 20         68         130         ≤ 30           1,2-Dichloroperopane         70         130         ≤ 20         70         130         ≤ 30           1,2-Dichloroperopane         70         130         ≤ 20         70         130         ≤ 30	1,1,1,2-Tetrachloroethane	64	130	≤ 20	70	130	≤ 30
1,1,2-Trichloroethane         70         130         ≤ 20         70         130         ≤ 30           1,1-Dichloroethane         70         130         ≤ 20         65         135         ≤ 30           1,1-Dichloroethene         61         145         ≤ 20         70         130         ≤ 30           1,2,3-Trichloropene         70         130         ≤ 20         70         130         ≤ 30           1,2,3-Trichloropene         64         130         ≤ 20         70         130         ≤ 30           1,2,4-Trichlorobenzene         70         130         ≤ 20         70         130         ≤ 30           1,2,4-Trichlorobenzene         70         130         ≤ 20         70         130         ≤ 30           1,2-Trichlorobenzene         70         130         ≤ 20         70         130         ≤ 30           1,2-Dichrome-Sachloropropane         41         144         ≤ 20         68         130         ≤ 30           1,2-Dichlorobenzene         70         130         ≤ 20         70         130         ≤ 30           1,2-Dichlorobenzene         70         130         ≤ 20         70         130         ≤ 30           1,3-	1,1,1-Trichloroethane	67	130	≤ 20	70	130	≤ 30
1,1-Dichloroethane         70         130         ≤ 20         65         135         ≤ 30           1,1-Dichloroethene         61         145         ≤ 20         70         130         ≤ 30           1,1-Dichloropropene         70         130         ≤ 20         70         130         ≤ 30           1,2,3-Trichloropropane         64         130         ≤ 20         70         130         ≤ 30           1,2,4-Trichloropenzene         70         130         ≤ 20         70         130         ≤ 30           1,2,4-Trichlorobenzene         70         130         ≤ 20         70         130         ≤ 30           1,2-Dichromo-3-chloropropane         41         144         ≤ 20         70         130         ≤ 30           1,2-Dichlorobenzene         70         130         ≤ 20         70         130         ≤ 30           1,2-Dichloroperane         70         130         ≤ 20         70         130         ≤ 30           1,2-Dichloroperane         70         130         ≤ 20         70         130         ≤ 30           1,2-Dichloroperopane         70         130         ≤ 20         70         130         ≤ 30           1,3-5	1,1,2,2-Tetrachloroethane	67	130	≤ 20	70	130	≤ 30
1,1-Dichloroethene       61       145       \$20       70       130       \$30         1,1-Dichloropropene       70       130       \$20       70       130       \$30         1,2,3-Trichlorobenzene       70       130       \$20       70       130       \$30         1,2,3-Trichlorobenzene       70       130       \$20       70       130       \$30         1,2,4-Trichlorobenzene       70       130       \$20       70       130       \$30         1,2,4-Trimethylbenzene       70       130       \$20       70       130       \$30         1,2-Dibromo-3-chloropropane       41       144       \$20       68       130       \$30         1,2-Dichlorobenzene       70       130       \$20       70       130       \$30         1,2-Dichloroethane       70       130       \$20       70       130       \$30         1,2-Dichloroptopane       70       130       \$20       70       130       \$30         1,3-Dichlorobenzene       70       130       \$20       70       130       \$30         1,3-Dichlorobenzene       70       130       \$20       70       130       \$30         1,4-D	1,1,2-Trichloroethane	70	130	≤ 20	70	130	≤ 30
1,1-Dichloropropene         70         130         ≤ 20         70         130         ≤ 30           1,2,3-Trichlorobenzene         70         130         ≤ 20         70         130         ≤ 30           1,2,3-Trichloropropane         64         130         ≤ 20         70         130         ≤ 30           1,2,4-Trinchlorobenzene         70         130         ≤ 20         70         130         ≤ 30           1,2,4-Trinchlorobenzene         70         130         ≤ 20         70         130         ≤ 30           1,2-Dibromo-3-chloropropane         41         144         ≤ 20         70         130         ≤ 30           1,2-Dichlorobenzene         70         130         ≤ 20         70         130         ≤ 30           1,2-Dichloropropane         70         130         ≤ 20         70         130         ≤ 30           1,2-Dichloropropane         70         130         ≤ 20         70         130         ≤ 30           1,3-Dichloropropane         70         130         ≤ 20         70         130         ≤ 30           1,3-Dichloropropane         70         130         ≤ 20         65         130         ≤ 30 <td< td=""><td>1,1-Dichloroethane</td><td>70</td><td>130</td><td>≤ 20</td><td>65</td><td>135</td><td>≤ 30</td></td<>	1,1-Dichloroethane	70	130	≤ 20	65	135	≤ 30
1,2,3-Trichlorobenzene         70         130         \$ 20         70         130         \$ 30           1,2,3-Trichloropropane         64         130         \$ 20         68         130         \$ 30           1,2,4-Trichlorobenzene         70         130         \$ 20         70         130         \$ 30           1,2-Hrimethylbenzene         70         130         \$ 20         70         130         \$ 30           1,2-Dibromo-3-chloropropane         41         144         \$ 20         68         130         \$ 30           1,2-Dichlorobenzene         70         130         \$ 20         70         130         \$ 30           1,2-Dichlorobenzene         70         130         \$ 20         70         130         \$ 30           1,2-Dichloropropane         70         130         \$ 20         70         130         \$ 30           1,3-5-Trimethylbenzene         64         130         \$ 20         70         130         \$ 30           1,3-Dichloropenzene         70         130         \$ 20         70         130         \$ 30           1,4-Dichlorobenzene         70         130         \$ 20         70         130         \$ 30	1,1-Dichloroethene	61	145	≤ 20	70	130	≤ 30
1,2,3-Trichloropropane       64       130       \$ 20       68       130       \$ 30         1,2,4-Trichlorobenzene       70       130       \$ 20       70       130       \$ 30         1,2,4-Trimethylbenzene       70       130       \$ 20       70       130       \$ 30         1,2-Dichloromoethane       70       130       \$ 20       70       130       \$ 30         1,2-Dichlorobenzene       70       130       \$ 20       70       130       \$ 30         1,2-Dichloroethane       70       130       \$ 20       70       130       \$ 30         1,2-Dichloroptopane       70       130       \$ 20       70       130       \$ 30         1,2-Dichloroptopane       70       130       \$ 20       70       130       \$ 30         1,3-Dichlorobenzene       70       130       \$ 20       70       130       \$ 30         1,3-Dichlorobenzene       70       130       \$ 20       70       130       \$ 30         1,4-Dichlorobenzene       70       130       \$ 20       70       130       \$ 30         1,4-Dichlorobenzene       70       130       \$ 20       70       130       \$ 30	1,1-Dichloropropene	70	130	≤ 20	70	130	≤ 30
1,2,4-Trichlorobenzene         70         130         ≤ 20         70         130         ≤ 30           1,2,4-Trimethylbenzene         70         130         ≤ 20         70         130         ≤ 30           1,2-Dibromo-3-chloropropane         41         144         ≤ 20         68         130         ≤ 30           1,2-Dibromoethane         70         130         ≤ 20         70         130         ≤ 30           1,2-Dibromoethane         70         130         ≤ 20         70         130         ≤ 30           1,2-Dibroropropane         70         130         ≤ 20         70         130         ≤ 30           1,3-Dichloropropane         70         130         ≤ 20         70         130         ≤ 30           1,3-Dichloropropane         70         130         ≤ 20         70         130         ≤ 30           1,3-Dichloropropane         70         130         ≤ 20         70         130         ≤ 30           1,3-Dichloropropane         70         130         ≤ 20         70         130         ≤ 30           1,3-Dichloropropane         70         130         ≤ 20         70         130         ≤ 30           1,4-Dichloro	1,2,3-Trichlorobenzene	70	130	≤ 20	70	130	≤ 30
1,2,4-Trimethylbenzene         70         130         ≤ 20         70         130         ≤ 30           1,2-Dibromo-3-chloropropane         41         144         ≤ 20         68         130         ≤ 30           1,2-Dibromoethane         70         130         ≤ 20         70         130         ≤ 30           1,2-Dichlorobenzene         70         130         ≤ 20         70         130         ≤ 30           1,2-Dichloropenane         70         130         ≤ 20         70         130         ≤ 30           1,3-Dichlorobenzene         70         130         ≤ 20         70         130         ≤ 30           1,3-Dichloropenzene         70         130         ≤ 20         70         130         ≤ 30           1,3-Dichloropropane         70         130         ≤ 20         70         130         ≤ 30           1,4-Dichlorobenzene         70         130         ≤ 20         69         130         ≤ 30           1,4-Dichloropropane         63         133         ≤ 20         70         130         ≤ 30           2,2-Dichloropropane         63         133         ≤ 20         70         130         ≤ 30           2-Butanone (	1,2,3-Trichloropropane	64	130	≤ 20	68	130	≤ 30
1,2-Dibromo-3-chloropropane         41         144         ≤ 20         70         130         ≤ 30           1,2-Dibromoethane         70         130         ≤ 20         70         130         ≤ 30           1,2-Dichlorobenzene         70         130         ≤ 20         70         130         ≤ 30           1,2-Dichloropenae         70         130         ≤ 20         70         130         ≤ 30           1,3-5-Trimethylbenzene         64         130         ≤ 20         70         130         ≤ 30           1,3-Dichlorobenzene         70         130         ≤ 20         70         130         ≤ 30           1,3-Dichlorobenzene         70         130         ≤ 20         70         130         ≤ 30           1,4-Dichlorobenzene         70         130         ≤ 20         69         130         ≤ 30           1,4-Dichlorobenzene         70         130         ≤ 20         70         130         ≤ 30           1,4-Dichlorobenzene         70         130         ≤ 20         70         130         ≤ 30           2,2-Dichloropropane         63         133         ≤ 20         70         130         ≤ 30           2-Dichloroben	1,2,4-Trichlorobenzene	70	130	≤ 20	70	130	≤ 30
1,2-Dibromoethane       70       130       ≤ 20       70       130       ≤ 30         1,2-Dichlorobenzene       70       130       ≤ 20       70       130       ≤ 30         1,2-Dichloroethane       70       130       ≤ 20       70       130       ≤ 30         1,2-Dichloropropane       70       130       ≤ 20       70       130       ≤ 30         1,3,5-Trimethylbenzene       64       130       ≤ 20       70       130       ≤ 30         1,3-Dichlorobenzene       70       130       ≤ 20       70       130       ≤ 30         1,3-Dichloropropane       70       130       ≤ 20       69       130       ≤ 30         1,4-Dichlorobenzene       70       130       ≤ 20       70       130       ≤ 30         1,4-Dichloropropane       56       162       ≤ 20       65       136       ≤ 30         1,4-Dioxane       56       162       ≤ 20       65       130       ≤ 30         2,2-Dichloropropane       63       133       ≤ 20       70       130       ≤ 30         2-Butanone (MEK)       63       138       ≤ 20       70       130       ≤ 30         2-Hexanon	1,2,4-Trimethylbenzene	70	130	≤ 20	70	130	≤ 30
1,2-Dichlorobenzene       70       130       ≤ 20       70       130       ≤ 30         1,2-Dichloroethane       70       130       ≤ 20       70       130       ≤ 30         1,2-Dichloropropane       70       130       ≤ 20       70       130       ≤ 30         1,3-Dichlorobenzene       64       130       ≤ 20       70       130       ≤ 30         1,3-Dichloropropane       70       130       ≤ 20       69       130       ≤ 30         1,4-Dichlorobenzene       70       130       ≤ 20       69       130       ≤ 30         1,4-Dichlorobenzene       70       130       ≤ 20       69       130       ≤ 30         1,4-Dichloropropane       56       162       ≤ 20       65       136       ≤ 30         1,4-Dioxane       663       133       ≤ 20       70       130       ≤ 30         2,2-Dichloropropane       63       133       ≤ 20       70       130       ≤ 30         2,2-Dichloropropane       63       133       ≤ 20       70       130       ≤ 30         2-Butanone (MEK)       63       138       ≤ 20       70       130       ≤ 30         2-Hexanon	1,2-Dibromo-3-chloropropane	41	144	≤ 20	68	130	≤ 30
1,2-Dichloroethane         70         130         ≤ 20         70         130         ≤ 30           1,2-Dichloropropane         70         130         ≤ 20         70         130         ≤ 30           1,3,5-Trimethylbenzene         64         130         ≤ 20         70         130         ≤ 30           1,3-Dichlorobenzene         70         130         ≤ 20         69         130         ≤ 30           1,4-Dichlorobenzene         70         130         ≤ 20         69         130         ≤ 30           1,4-Dichlorobenzene         70         130         ≤ 20         69         130         ≤ 30           1,4-Dichlorobenzene         56         162         ≤ 20         65         136         ≤ 30           1,4-Dioxane         56         162         ≤ 20         65         136         ≤ 30           2,2-Dichloropropane         63         133         ≤ 20         70         130         ≤ 30           2,2-Dichloropropane         63         138         ≤ 20         70         130         ≤ 30           2-Chlorobluene         70         130         ≤ 20         70         130         ≤ 30           2-Hexanone         70<	1,2-Dibromoethane	70	130	≤ 20	70	130	≤ 30
1,2-Dichloropropane       70       130       ≤ 20       70       130       ≤ 30         1,3,5-Trimethylbenzene       64       130       ≤ 20       70       130       ≤ 30         1,3-Dichlorobenzene       70       130       ≤ 20       69       130       ≤ 30         1,4-Dichloropropane       70       130       ≤ 20       69       130       ≤ 30         1,4-Dichlorobenzene       70       130       ≤ 20       65       136       ≤ 30         1,4-Dichloropropane       56       162       ≤ 20       65       136       ≤ 30         2,2-Dichloropropane       63       133       ≤ 20       70       130       ≤ 30         2-Butanone (MEK)       63       138       ≤ 20       70       130       ≤ 30         2-Butanone (MEK)       63       138       ≤ 20       70       130       ≤ 30         2-Hexanone       70       130       ≤ 20       70       130       ≤ 30         4-Chlorotoluene       70       130       ≤ 20       70       130       ≤ 30         4-Methyl-2-pentanone (MIBK)       59       130       ≤ 20       70       130       ≤ 30         Bromob	1,2-Dichlorobenzene	70	130	≤ 20	70	130	≤ 30
1,3,5-Trimethylbenzene       64       130       ≤ 20       70       130       ≤ 30         1,3-Dichlorobenzene       70       130       ≤ 20       70       130       ≤ 30         1,3-Dichloropropane       70       130       ≤ 20       69       130       ≤ 30         1,4-Dichlorobenzene       70       130       ≤ 20       70       130       ≤ 30         1,4-Dioxane       56       162       ≤ 20       65       136       ≤ 30         2,2-Dichloropropane       63       133       ≤ 20       70       130       ≤ 30         2-Butanone (MEK)       63       138       ≤ 20       70       130       ≤ 30         2-Hexanone       70       130       ≤ 20       70       130       ≤ 30         4-Chlorotoluene       70       130       ≤ 20       70       130       ≤ 30         4-Methyl-2-pentanone (MIBK)       59       130       ≤ 20       70       130       ≤ 30         Benzene       70       130       ≤ 20       70       130       ≤ 30         Bromobenzene       70       130       ≤ 20       70       130       ≤ 30         Bromochloromethane	1,2-Dichloroethane	70	130	≤ 20	70	130	≤ 30
1,3-Dichlorobenzene       70       130       ≤ 20       70       130       ≤ 30         1,3-Dichloropropane       70       130       ≤ 20       69       130       ≤ 30         1,4-Dichlorobenzene       70       130       ≤ 20       70       130       ≤ 30         1.4-Dioxane       56       162       ≤ 20       65       136       ≤ 30         2,2-Dichloropropane       63       133       ≤ 20       70       130       ≤ 30         2-Butanone (MEK)       63       138       ≤ 20       70       130       ≤ 30         2-Chlorotoluene       70       130       ≤ 20       70       130       ≤ 30         2-Hexanone       57       130       ≤ 20       70       130       ≤ 30         4-Chlorotoluene       70       130       ≤ 20       70       130       ≤ 30         4-Methyl-2-pentanone (MIBK)       59       130       ≤ 20       70       130       ≤ 30         Acetone       58       148       ≤ 20       54       140       ≤ 30         Benzene       70       130       ≤ 20       70       130       ≤ 30         Bromobenzene       70	1,2-Dichloropropane	70	130	≤ 20	70	130	≤ 30
1,3-Dichloropropane         70         130         ≤ 20         69         130         ≤ 30           1,4-Dichlorobenzene         70         130         ≤ 20         70         130         ≤ 30           1.4-Dioxane         56         162         ≤ 20         65         136         ≤ 30           2,2-Dichloropropane         63         133         ≤ 20         70         130         ≤ 30           2-Butanone (MEK)         63         138         ≤ 20         70         130         ≤ 30           2-Chlorotoluene         70         130         ≤ 20         70         130         ≤ 30           2-Hexanone         57         130         ≤ 20         70         130         ≤ 30           4-Chlorotoluene         70         130         ≤ 20         70         130         ≤ 30           4-Methyl-2-pentanone (MIBK)         59         130         ≤ 20         70         130         ≤ 30           Acetone         58         148         ≤ 20         70         130         ≤ 30           Benzene         70         130         ≤ 20         70         130         ≤ 30           Bromochloromethane         67         130	1,3,5-Trimethylbenzene	64	130	≤ 20	70	130	≤ 30
1,4-Dicklorobenzene       70       130       ≤ 20       70       130       ≤ 30         1.4-Dioxane       56       162       ≤ 20       65       136       ≤ 30         2,2-Dickloropropane       63       133       ≤ 20       70       130       ≤ 30         2-Butanone (MEK)       63       138       ≤ 20       70       130       ≤ 30         2-Chlorotoluene       70       130       ≤ 20       70       130       ≤ 30         2-Hexanone       57       130       ≤ 20       70       130       ≤ 30         4-Chlorotoluene       70       130       ≤ 20       70       130       ≤ 30         4-Methyl-2-pentanone (MIBK)       59       130       ≤ 20       70       130       ≤ 30         Acetone       58       148       ≤ 20       54       140       ≤ 30         Benzene       70       130       ≤ 20       70       130       ≤ 30         Bromobenzene       70       130       ≤ 20       70       130       ≤ 30         Bromochloromethane       67       130       ≤ 20       70       130       ≤ 30         Bromoform       54       136 <td>1,3-Dichlorobenzene</td> <td>70</td> <td>130</td> <td>≤ 20</td> <td>70</td> <td>130</td> <td>≤ 30</td>	1,3-Dichlorobenzene	70	130	≤ 20	70	130	≤ 30
1.4-Dioxane       56       162       ≤ 20       65       136       ≤ 30         2,2-Dichloropropane       63       133       ≤ 20       70       130       ≤ 30         2-Butanone (MEK)       63       138       ≤ 20       70       130       ≤ 30         2-Chlorotoluene       70       130       ≤ 20       70       130       ≤ 30         2-Hexanone       57       130       ≤ 20       70       130       ≤ 30         4-Chlorotoluene       70       130       ≤ 20       70       130       ≤ 30         4-Methyl-2-pentanone (MIBK)       59       130       ≤ 20       70       130       ≤ 30         Acetone       58       148       ≤ 20       54       140       ≤ 30         Benzene       70       130       ≤ 20       70       130       ≤ 30         Bromobenzene       70       130       ≤ 20       70       130       ≤ 30         Bromodichloromethane       67       130       ≤ 20       70       130       ≤ 30         Bromoform       54       136       ≤ 20       70       130       ≤ 30         Bromodichloromethane       51       130<	1,3-Dichloropropane	70	130	≤ 20	69	130	≤ 30
2,2-Dichloropropane       63       133       ≤ 20       70       130       ≤ 30         2-Butanone (MEK)       63       138       ≤ 20       70       130       ≤ 30         2-Chlorotoluene       70       130       ≤ 20       70       130       ≤ 30         2-Hexanone       57       130       ≤ 20       70       130       ≤ 30         4-Chlorotoluene       70       130       ≤ 20       70       130       ≤ 30         4-Methyl-2-pentanone (MIBK)       59       130       ≤ 20       70       130       ≤ 30         Acetone       58       148       ≤ 20       54       140       ≤ 30         Benzene       70       130       ≤ 20       70       130       ≤ 30         Bromobenzene       70       130       ≤ 20       70       130       ≤ 30         Bromochloromethane       67       130       ≤ 20       70       130       ≤ 30         Bromoform       54       136       ≤ 20       70       130       ≤ 30         Bromomethane       39       139       ≤ 20       57       147       ≤ 30         Carbon tetrachloride       63       132 </td <td>1,4-Dichlorobenzene</td> <td>70</td> <td>130</td> <td>≤ 20</td> <td>70</td> <td>130</td> <td>≤ 30</td>	1,4-Dichlorobenzene	70	130	≤ 20	70	130	≤ 30
2-Butanone (MEK) 63 138 ≤20 70 130 ≤30 2-Chlorotoluene 70 130 ≤20 70 130 ≤30 2-Hexanone 57 130 ≤20 70 130 ≤30 4-Chlorotoluene 70 130 ≤20 70 130 ≤30 4-Chlorotoluene 70 130 ≤20 70 130 ≤30 4-Methyl-2-pentanone (MIBK) 59 130 ≤20 70 130 ≤30 Acetone 58 148 ≤20 54 140 ≤30 Benzene 70 130 ≤20 70 130 ≤30 Bromobenzene 70 130 ≤20 70 130 ≤30 Bromochloromethane 70 130 ≤20 70 130 ≤30 Bromochloromethane 67 130 ≤20 70 130 ≤30 Bromodichloromethane 67 130 ≤20 70 130 ≤30 Carbon disulfide 51 130 ≤20 57 147 ≤30 Carbon tetrachloride 63 132 ≤20 70 130 ≤30 Chlorobenzene 75 130 ≤20 70 130 ≤30 Chlorodibromomethane	1.4-Dioxane	56	162	≤ 20	65	136	≤ 30
2-Chlorotoluene       70       130       ≤ 20       70       130       ≤ 30         2-Hexanone       57       130       ≤ 20       70       130       ≤ 30         4-Chlorotoluene       70       130       ≤ 20       70       130       ≤ 30         4-Methyl-2-pentanone (MIBK)       59       130       ≤ 20       70       130       ≤ 30         Acetone       58       148       ≤ 20       54       140       ≤ 30         Benzene       70       130       ≤ 20       70       130       ≤ 30         Bromobenzene       70       130       ≤ 20       70       130       ≤ 30         Bromochloromethane       70       130       ≤ 20       70       130       ≤ 30         Bromoform       67       130       ≤ 20       70       130       ≤ 30         Bromoform       54       136       ≤ 20       70       130       ≤ 30         Bromomethane       39       139       ≤ 20       57       147       ≤ 30         Carbon disulfide       51       130       ≤ 20       59       130       ≤ 30         Chlorobenzene       75       130       ≤ 2	2,2-Dichloropropane	63	133	≤ 20	70	130	≤ 30
2-Hexanone 57 130 ≤20 70 130 ≤30 4-Chlorotoluene 70 130 ≤20 70 130 ≤30 4-Methyl-2-pentanone (MIBK) 59 130 ≤20 70 130 ≤30 Acetone 58 148 ≤20 54 140 ≤30 Benzene 70 130 ≤20 70 130 ≤30 Bromobenzene 70 130 ≤20 70 130 ≤30 Bromochloromethane 70 130 ≤20 70 130 ≤30 Bromochloromethane 67 130 ≤20 70 130 ≤30 Bromoform 54 136 ≤20 70 130 ≤30 Bromomethane 39 139 ≤20 70 130 ≤30 Carbon disulfide 51 130 ≤20 57 147 ≤30 Carbon tetrachloride 63 132 ≤20 70 130 ≤30 Chlorobenzene 75 130 ≤20 70 130 ≤30 Chlorodibromomethane 63 130 ≤20 70 130 ≤30	2-Butanone (MEK)	63	138	≤ 20	70	130	≤ 30
4-Chlorotoluene       70       130       ≤ 20       70       130       ≤ 30         4-Methyl-2-pentanone (MIBK)       59       130       ≤ 20       70       130       ≤ 30         Acetone       58       148       ≤ 20       54       140       ≤ 30         Benzene       70       130       ≤ 20       70       130       ≤ 30         Bromobenzene       70       130       ≤ 20       70       130       ≤ 30         Bromochloromethane       70       130       ≤ 20       70       130       ≤ 30         Bromodichloromethane       67       130       ≤ 20       70       130       ≤ 30         Bromoform       54       136       ≤ 20       70       130       ≤ 30         Bromomethane       39       139       ≤ 20       57       147       ≤ 30         Carbon disulfide       51       130       ≤ 20       59       130       ≤ 30         Chlorobenzene       75       130       ≤ 20       70       130       ≤ 30         Chlorodibromomethane       63       130       ≤ 20       70       130       ≤ 30         Chlorodibromomethane       63	2-Chlorotoluene	70	130	≤ 20	70	130	≤ 30
4-Methyl-2-pentanone (MIBK)59130 $\leq 20$ 70130 $\leq 30$ Acetone58148 $\leq 20$ 54140 $\leq 30$ Benzene70130 $\leq 20$ 70130 $\leq 30$ Bromobenzene70130 $\leq 20$ 70130 $\leq 30$ Bromochloromethane70130 $\leq 20$ 70130 $\leq 30$ Bromoform67130 $\leq 20$ 70130 $\leq 30$ Bromomethane39139 $\leq 20$ 70130 $\leq 30$ Carbon disulfide51130 $\leq 20$ 59130 $\leq 30$ Carbon tetrachloride63132 $\leq 20$ 70130 $\leq 30$ Chlorobenzene75130 $\leq 20$ 70130 $\leq 30$ Chlorodibromomethane63130 $\leq 20$ 70130 $\leq 30$	2-Hexanone	57	130	≤ 20	70	130	≤ 30
Acetone       58       148 $\leq 20$ 54       140 $\leq 30$ Benzene       70       130 $\leq 20$ 70       130 $\leq 30$ Bromobenzene       70       130 $\leq 20$ 70       130 $\leq 30$ Bromochloromethane       67       130 $\leq 20$ 70       130 $\leq 30$ Bromoform       54       136 $\leq 20$ 70       130 $\leq 30$ Bromomethane       39       139 $\leq 20$ 57       147 $\leq 30$ Carbon disulfide       51       130 $\leq 20$ 59       130 $\leq 30$ Carbon tetrachloride       63       132 $\leq 20$ 70       130 $\leq 30$ Chlorobenzene       75       130 $\leq 20$ 70       130 $\leq 30$ Chlorodibromomethane       63       130 $\leq 20$ 70       130 $\leq 30$	4-Chlorotoluene	70	130	≤ 20	70	130	≤ 30
Benzene       70       130       ≤ 20       70       130       ≤ 30         Bromobenzene       70       130       ≤ 20       70       130       ≤ 30         Bromochloromethane       70       130       ≤ 20       70       130       ≤ 30         Bromodichloromethane       67       130       ≤ 20       70       130       ≤ 30         Bromoform       54       136       ≤ 20       70       130       ≤ 30         Bromomethane       39       139       ≤ 20       57       147       ≤ 30         Carbon disulfide       51       130       ≤ 20       59       130       ≤ 30         Carbon tetrachloride       63       132       ≤ 20       70       130       ≤ 30         Chlorobenzene       75       130       ≤ 20       70       130       ≤ 30         Chlorodibromomethane       63       130       ≤ 20       70       130       ≤ 30	4-Methyl-2-pentanone (MIBK)	59	130	≤ 20	70	130	≤ 30
Bromobenzene       70       130       ≤ 20       70       130       ≤ 30         Bromochloromethane       70       130       ≤ 20       70       130       ≤ 30         Bromodichloromethane       67       130       ≤ 20       70       130       ≤ 30         Bromoform       54       136       ≤ 20       70       130       ≤ 30         Bromomethane       39       139       ≤ 20       57       147       ≤ 30         Carbon disulfide       51       130       ≤ 20       59       130       ≤ 30         Carbon tetrachloride       63       132       ≤ 20       70       130       ≤ 30         Chlorobenzene       75       130       ≤ 20       70       130       ≤ 30         Chlorodibromomethane       63       130       ≤ 20       70       130       ≤ 30	Acetone	58	148	≤ 20	54	140	≤ 30
Bromochloromethane       70       130       ≤ 20       70       130       ≤ 30         Bromodichloromethane       67       130       ≤ 20       70       130       ≤ 30         Bromoform       54       136       ≤ 20       70       130       ≤ 30         Bromomethane       39       139       ≤ 20       57       147       ≤ 30         Carbon disulfide       51       130       ≤ 20       59       130       ≤ 30         Carbon tetrachloride       63       132       ≤ 20       70       130       ≤ 30         Chlorobenzene       75       130       ≤ 20       70       130       ≤ 30         Chlorodibromomethane       63       130       ≤ 20       70       130       ≤ 30	Benzene	70	130	≤ 20	70	130	≤ 30
Bromodichloromethane $67$ $130$ $\leq 20$ $70$ $130$ $\leq 30$ Bromoform $54$ $136$ $\leq 20$ $70$ $130$ $\leq 30$ Bromomethane $39$ $139$ $\leq 20$ $57$ $147$ $\leq 30$ Carbon disulfide $51$ $130$ $\leq 20$ $59$ $130$ $\leq 30$ Carbon tetrachloride $63$ $132$ $\leq 20$ $70$ $130$ $\leq 30$ Chlorobenzene $75$ $130$ $\leq 20$ $70$ $130$ $\leq 30$ Chlorodibromomethane $63$ $130$ $\leq 20$ $70$ $130$ $\leq 30$	Bromobenzene	70	130	≤ 20	70	130	≤ 30
Bromoform       54       136       ≤ 20       70       130       ≤ 30         Bromomethane       39       139       ≤ 20       57       147       ≤ 30         Carbon disulfide       51       130       ≤ 20       59       130       ≤ 30         Carbon tetrachloride       63       132       ≤ 20       70       130       ≤ 30         Chlorobenzene       75       130       ≤ 20       70       130       ≤ 30         Chlorodibromomethane       63       130       ≤ 20       70       130       ≤ 30	Bromochloromethane	70	130	≤ 20	70	130	≤ 30
Bromomethane       39       139       ≤ 20       57       147       ≤ 30         Carbon disulfide       51       130       ≤ 20       59       130       ≤ 30         Carbon tetrachloride       63       132       ≤ 20       70       130       ≤ 30         Chlorobenzene       75       130       ≤ 20       70       130       ≤ 30         Chlorodibromomethane       63       130       ≤ 20       70       130       ≤ 30	Bromodichloromethane	67	130	≤ 20	70	130	≤ 30
Carbon disulfide       51       130       ≤ 20       59       130       ≤ 30         Carbon tetrachloride       63       132       ≤ 20       70       130       ≤ 30         Chlorobenzene       75       130       ≤ 20       70       130       ≤ 30         Chlorodibromomethane       63       130       ≤ 20       70       130       ≤ 30	Bromoform	54	136	≤ 20	70	130	≤ 30
Carbon tetrachloride       63       132       ≤ 20       70       130       ≤ 30         Chlorobenzene       75       130       ≤ 20       70       130       ≤ 30         Chlorodibromomethane       63       130       ≤ 20       70       130       ≤ 30	Bromomethane	39	139	≤ 20	57	147	≤ 30
Chlorobenzene       75       130       ≤ 20       70       130       ≤ 30         Chlorodibromomethane       63       130       ≤ 20       70       130       ≤ 30	Carbon disulfide	51	130	≤ 20	59	130	≤ 30
Chlorodibromomethane         63         130         ≤ 20         70         130         ≤ 30	Carbon tetrachloride	63	132	≤ 20	70	130	≤ 30
	Chlorobenzene	75	130	≤ 20	70	130	≤ 30
Chloroethane 55 138 ≤ 20 50 151 ≤ 30	Chlorodibromomethane	63	130	≤ 20	70	130	≤ 30
	Chloroethane	55	138	≤ 20	50	151	≤ 30

Table A-5. Accuracy and Precision for Volatile Organic Compounds by Method SW8260C *Essex-Hope Site, Jamestown, New York* 

	Acc W	D/MS/MSD uracy /ater %R)	Precision Water	LCS/LCSD/MS/MSD Accuracy Soil (%R)		Precision _ Soil
Parameter	LCL	UCL	(%RPD)	LCL	UCL	(%RPD)
Chloroform	70	130	≤ 20	70	130	≤ 30
Chloromethane	64	130	≤ 20	52	130	≤ 30
cis-1,2-Dichloroethene	70	130	≤ 20	70	130	≤ 30
cis-1,3-Dichloropropene	70	130	≤ 20	70	130	≤ 30
Dibromomethane	70	130	≤ 20	70	130	≤ 30
Dichlorodifluoromethane	36	147	≤ 20	30	146	≤ 30
Ethylbenzene	70	130	≤ 20	70	130	≤ 30
Hexachlorobutadiene	63	130	≤ 20	67	130	≤ 30
Isopropylbenzene	70	130	≤ 20	70	130	≤ 30
Methylene Chloride	70	130	≤ 20	70	130	≤ 30
Methyl-t butyl ether	63	130	≤ 20	66	130	≤ 30
m,p-xylene	70	130	≤ 20	70	130	≤ 30
n-Butylbenzene	53	136	≤ 20	70	130	≤ 30
N-Propylbenzene	69	130	≤ 20	70	130	≤ 30
Naphthalene	70	130	≤ 20	70	130	≤ 30
o-xylene	70	130	≤ 20	70	130	≤ 30
p-Isopropyltoluene	70	130	≤ 20	70	130	≤ 30
sec-Butylbenzene	70	130	≤ 20	70	130	≤ 30
Styrene	70	130	≤ 20	70	130	≤ 30
tert-Butylbenzene	70	130	≤ 20	70	130	≤ 30
Tetrachloroethene	70	130	≤ 20	70	130	≤ 30
Toluene	70	130	≤ 20	70	130	≤ 30
trans-1,2-Dichloroethene	70	130	≤ 20	70	130	≤ 30
trans-1,3-Dichloropropene	70	130	≤ 20	70	130	≤ 30
Trichloroethene	70	130	≤ 20	70	130	≤ 30
Trichlorofluoromethane	62	150	≤ 20	70	139	≤ 30
Vinyl Acetate	70	130	≤ 20	70	130	≤ 30
Vinyl chloride	55	140	≤ 20	67	130	≤ 30
Surrogates						
1,2-Dichloroethane-d4	70	130		70	130	
4-Bromofluorobenzene	70	130		70	130	
Dibromofluoromethane	70	130		70	130	
Toluene-d8	70	130		70	130	

%R = percent recovery

RPD = relative percent difference

Table A-6. Accuracy and Precision Limits for Polychlorinated Biphenyls by Method SW8082 (Soil) Essex-Hope Site, Jamestown, New York

	S	SD Accuracy oil 6R)	Precision Soil
Analyte	LCL	UCL	(% RPD)
Aroclor 1016	40	140	≤ 50
Aroclor 1260	40	140	≤ 50
Surrogates			
2,4,5,6-Tetrachloro-m-xylene	30	150	
Decachlorobiphenyl	30	150	

<sup>%</sup>R = percent recovery

RPD = relative percent difference

Table A-7. Calibration and QC Requirements for Volatile Organic Compounds by SW8260C *Essex-Hope Site, Jamestown, New York* 

QC Check	Frequency	Criteria	<b>Corrective Action</b>
BFB Tuning	Prior to initial calibration and calibration (every 12 hours)	Refer to criteria listed in the method	Retune instrument and verify
Multi-point initial calibration (minimum five points)	Prior to sample analysis, or when calibration fails	Average RF for all analytes ≥ 0.10 <sup>a</sup> and one option below:  Option 1:  %RSD for all analytes ≤ 20%  Option 2:  Least squares regression r ≥ 0.990	Correct the problem and repeat the initial calibration.
Second-source calibration verification	Once for each multi-point initial calibration	All analytes within ±30% of expected value	Correct the problem and repeat initial calibration.
Continuing calibration verification	At the start of each analytical sequence and every 12 hours thereafter	Average RF for all analytes $\geq 0.10^a$ All analytes within + 20% of expected value.	Correct the problem, then recalibrate and reanalyze all samples since the last acceptable continuing calibration verification.
Retention time window calculated for each analyte	Each analyte	Relative retention time of each analyte within + 0.06 relative retention time units of the continuing calibration verification	Not applicable (used for identification of analyte)
Internal Standards	Each sample and QC sample, method blank, MS/MSD and LCS	Retention time within ±30 seconds from retention time of the daily continuing calibration verification standard.  EICP area within –50% to +100% of the daily continuing calibration	Inspect mass spectrometer and gas chromatography for malfunctions; reanalyze all affected samples
Method Blank	At least one per analytical batch	No analytes detected at or above the reporting limit	Correct the problem, then re-prep and reanalyze all associated samples
Surrogate spike	Every standard, sample, method blank, MS/MSD and LCS	All surrogates in samples, method blank and LCS within limits specified in Accuracy and Precision table	Correct the problem and reanalyze (re-prep if necessary).
MS/MSD	One set per 20 project-specific samples	Within limits specified in Accuracy and Precision table	None
LCS	At least one per analytical batch	Within limits specified in Accuracy and Precision table	Correct the problem, then re-prep and reanalyze the LCS and all samples in the analytical batch.

<sup>&</sup>lt;sup>a</sup>Average relative response factor (RRF) specific for each comment. See Table 4 in SW8260C method.

Table A-8. Calibration and QC Requirements for Polychlorinated Biphenyls by SW8082 Essex-Hope Site, Jamestown, New York

QC Check	Frequency	Criteria	<b>Corrective Action</b>
Multi-point initial calibration (minimum five points) for Aroclors	Prior to sample analysis, or when calibration verification fails	Option 1: %RSD for all analytes ≤ 20%	Correct the problem and repeat the initial calibration.
1016/1260 only		Option 2: Least squares regression $r \ge 0.990$	
Second-source calibration verification	Once for each multi-point initial calibration	All analytes within ±20% of expected value	Correct the problem and repeat initial calibration.
Continuing calibration verification	At the start of each analytical sequence and every 12 hours thereafter	All analytes within ±15% of expected value.	Correct the problem, then recalibrate and reanalyze all samples since the last acceptable continuing calibration verification.
Method Blank	At least one per analytical batch	No analytes detected at or above the reporting limit	Correct the problem, then re-prep and reanalyze all associated samples
Surrogate spike	Every standard, sample, method blank, MS/MSD and LCS	All surrogates in samples, method blank and LCS within limits specified in Accuracy and Precision table	Correct the problem and reanalyze (re-prep if necessary).
MS/MSD	One set per 20 project-specific samples	Within limits specified in Accuracy and Precision table	None
LCS	At least one per analytical batch	Within limits specified in Accuracy and Precision table	Correct the problem, then re-prep and reanalyze the LCS and all samples in the analytical batch.