

SCIENCE APPLICATIONS INTERNATIONAL CORPORATION  
Organic Data Review Checklist - Standard Validation

Project: Harley-Davidson

Page 1 of 11

SDG No:

180-47781-1

Analysis:

VOC

Laboratory:

TestAmerica Pittsburgh

Method:

8260 LL

Matrix:

Water

The above data package has been reviewed and the analytical quality control/quality assurance performance data have been summarized. The general criteria used to assess the analytical integrity of the data were based on an examination of the following:

Case Narrative	Analytical Surrogate Recoveries
Analytical Holding Times	Internal Standard Performance
Sample Preservation	MS/MSD Recoveries and Differences
Method Calibration	LCS Recoveries
Method and Project Blanks	Re-analysis and Secondary Dilution

Project Specific QA/QC or contract requirements may take priority over validation criteria in this procedure.

Overall Remarks:

See corresponding pages ms/msd, LCS,  
and calibration issues, as well as  
quality issues

Definition of Qualifiers:

"U", not detected at the associated level  
"UJ", not detected and associated value estimated  
"J", associated value estimated  
"R", associated value unusable or analyte identity unfounded  
"=", compound properly identified and value positive

Reviewed by:

[Signature] Alan G. Miller Sr.

Date:

11/13/15

QA Reviewed by:

[Signature]

Date:

1-25-16

FR AGM 12/2/15

Paul M. Gibney 11/20/15

**I. Case Narrative**

Verify direct statements made within the Laboratory Case Narrative (note discrepancies).

Remarks:           No major issues            
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**II. Re-analysis and Secondary Dilutions**

Verify that re-analysis and secondary dilutions were performed and reported as necessary. Determine appropriate results to report.

Remarks:           No issues            
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**III. Holding Times**

VOC - Waters - unpreserved: aromatic within 7 days, non-aromatic within 14 days of sample collection

VOC - Waters - preserved: aromatic and non-aromatic within 14 days of sample collection

VOC - Soils - preserve or analyze within 48 hours of sample collection; analyze within 14 days of preservation

SVOC, Pest., PCB - Waters - extract within 7 days of sample collection, analyze within 40 days of extraction

SVOC, Pest., PCB - Soils - extract within 14 days of sample collection, analyze within 40 days of extraction

**Deviations:**

Sample #	VOC		SVOC			Pest/PCB		
	Date Collected	Date Analyzed	Date Collected	Date Extracted	Date Analyzed	Date Collected	Date Extracted	Date Analyzed

**Actions:**

- 1. If holding times are exceeded, all results are qualified as estimated (J/UJ)
- 2. If holding times are exceeded by more than 2X, reviewer may qualify non-detected results as unusable (R)

**Remarks:**

*No issues*

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### IV. System Monitoring Compounds (SMC) Recoveries (VOC, SVOC, Pesticides, PCBs)

List SMC compounds with unacceptable recoveries:

**Deviations:**

Sample #	VOC			SVOC B/N Compounds			SVOC Acid Compounds			Pest	PCB
	TOL	BFB	DCE	NBZ	FBP	TPH	PHL	2FP	TBP	TCX	DCB
QC Limits											

**Actions:**

- 1. If any SMC recovery is <10%, qualify all positive results in associated fractions as estimated (J)
- 2. If any SMC recovery is <10%, qualify all nondetects in associated fractions as unusable (R)
- 3. If SMC recoveries fall between 10% and the lower recovery limit, qualify results as estimated (J/UJ)
- 4. If SMC recoveries fall above the upper recovery limit, qualify positive results as estimated (J)
- 5. Use professional judgement to qualify Pest/PCB results when SMC recoveries are >10%
- 6. Use professional judgement to qualify results when SMC recoveries have been diluted out of spec.
- 7. For SVOC, qualification of the data is required only when 2 or more SMC per fraction are not within control limits
- 8. Note: SMC formerly known as surrogates.

**Remarks:**         No issues        

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**V. Internal Standards Performance (VOC, SVOC)**

VOC internal standard area counts within -50% to +100% of standard (Y/N)

VOC internal standard retention times within  $\pm$  30 seconds of standard (Y/N)

SVOC internal standard area counts within -50% to +100% of standard (Y/N)

SVOC internal standard retention times within  $\pm$  30 seconds of standard (Y/N)

**Deviations:**

Sample #	IS Affected	Area Counts	Acceptable Range	RT	Std. RT Value

**Actions:**

1. If area counts are outside limits, qualify positive results associated with that IS as estimated (J)
2. Non-detected compounds quantitated using an IS area count >100% should not be qualified
3. Non-detected compounds quantitated using an IS area count <50%, qualify as estimated (UJ)
4. If extremely low area counts are reported (<50% of the lower limit), qualify non-detects as unusable (R)
5. If an IS retention time varies more than 30 seconds, review the chromatographic profile for shifts and irregularities. Use professional judgement to qualify the data estimated (J/UJ) or unusable (R)

**Remarks:**

*no issues*

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**VI. Blanks**

All blanks were reported per matrix per concentration level for each 12 hour period on each GC/ MS system used to analyze VOCs and SVOCs Yes  No

Review associated laboratory and project blank samples. List documented contamination below:

**Laboratory Method Blanks:**

<u>Date:</u>	<u>Lab ID #</u>	<u>Fraction</u>	<u>Compound</u>	<u>Conc. (ppb)</u>

**Associated Project Blanks (e.g., equipment rinsates, trip blanks, etc.)**

<u>Date</u>	<u>Lab ID #</u>	<u>Fraction</u>	<u>Compound</u>	<u>Conc. (ppb)</u>

Remarks: NO ISSUES

**VI. Blanks (continued)**

Calculate action levels based on 10X the highest blank concentration of "common laboratory solvents", VOCs (methylene chloride, acetone, toluene, 2-butanone, cyclohexane) or SVOCs (phthalates), and 5X the highest blank concentration for all other VOC, SVOC, Pesticides, and PCB compounds. Sample weights, volumes, and dilution factors must be taken into account when applying the 5X and 10X criteria. This allows the total amount of contaminant present to be considered.

**Deviations:**

Compound	Maximum Conc. Detected, (ppb)	Action Level (ppb)	Samples Affected

**Actions:**

1. If compound results exceed the action levels, the data are not qualified
2. If compound results are below the required reporting level, report results as non-detect (U) at the reporting level
3. If the compound is detected above the reporting level, but below the action level, qualify as not-detected (U)
4. If gross contamination exists in blanks (i.e., saturated peaks by GC/ MS), all affected compounds in the associated samles should be qualified as unusable (R) due to interference.
5. If blanks were not analyzed per matrix per concentration level for each 12 hour period on each GC/MS system used to analyze VOCs and SVOCs use professional judgement to qualify data. Data may be rejected (R).

**Remarks:** No Issues  
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**VII. Initial & Continuing Calibration (VOC, SVOC)**

GC/MS instrument performance checks (BFB / DFTPP) Acceptable Y or N  
 All compounds must have and RRF > 0.01, %RSD < 30, and %D < 25

VOC - Date of initial calibration: \_\_\_\_\_

*8/26/15 17:59*

VOC - Date(s) of continuing calibration: \_\_\_\_\_

*9/14/15 1226, 9/22/15 1322, 9/24/15 1105  
 9/24/15 1104*

Was the 12 hour criteria met?  Y or N

SVOC- Date of initial calibration: \_\_\_\_\_

SVOC - Date(s) of continuing calibration: \_\_\_\_\_

Was the 12 hour criteria met? Y or N

**Deviations:**

Compound	Date	RRF	%RSD	%D	Samples Affected

\* % Difference =  $((RF_{CCV} - RF_{ICAL\ AVG}) / RF_{ICAL\ AVG}) \times 100$ . In instances where the bias of the CCV impacts validation qualifiers, review the RF values or amount reported to confirm that the % Difference or % Drift are reported with the correct negative or positive value.

**Actions:**

1. If any compound has an initial or continuing RRF of < 0.01, qualify positive results as estimated (J)
2. If any compound has an initial or continuing RRF of < 0.01, qualify non-detects as unusable (R)
3. If any compound has a %RSD >30 or a %D >25, qualify positive results as estimated (J)
4. If any compound has a %RSD >40 or a %D >40, qualify non-detects as estimated (UJ)
5. If BFB or DFTPP mass assignment / ION abundance criteria are all associated data as unusable (R).
6. If samples were analyzed outside the 12 hour BFB or DFTPP performance check time period, qualify the affected sample data as estimated (J/UJ).
7. If separate calibration for water and soil were not performed, use professional judgement to evaluate the data. Data may be rejected (R).
8. If calibrations were not completed within the 12 hour criterion, qualify all associated data as estimated (J/UJ). If the 12 hour criterion was grossly exceeded, reject all associated data (R).

**Remarks:**

*See attached*



**VIII. Initial & Continuing Calibration (Pesticides, PCBs)**

Linearity evaluation, are %RSD <20? (Y/N)

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Is the RPD between calibration factors ≤25? (Y/N)

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Are multicomponent calibration data provided for each analysis date? (Y/N)

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Is the difference between columns check ≤ 25%D? (Y/N)

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Are 4, 4'- DDT and endrin breakdown (PEM) ≤ 20% and combined breakdown ≤ 30% (Y/N)

**Deviations:**

Compound	%RSD	RPD	Samples Affected

\* % Difference =  $((RF_{CCV} - RF_{ICAL\ AVG}) / RF_{ICAL\ AVG}) \times 100$ . In instances where the bias of the CCV impacts validation qualifiers, review the RF values or amount reported to confirm that the % Difference or % Drift are reported with the correct negative or positive value.

**Actions:**

- 1. If %RSD criteria are not met, qualify positive results as estimated (J) and non-detects as estimated (UJ)
- 2. If RPD criteria are not met, qualify positive results as estimated (J) and non-detects as estimated (UJ)
- 3. If %D criteria is not met, qualify positive results as estimated (J) and non-detects as estimated (UJ)
- 4. If breakdown criteria are not met, positive 4, 4'-DDT and endrin should be qualified as estimated (J).  
And non-detects should be rejected (R).

**Remarks:**

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**IX. Matrix Spike/Matrix Spike Duplicate Information**

General MS/MSD Criteria:

percent recovery (%R)

relative percent difference (RPD)

VOC	SVOC	Pest	PCB
70-130	45-135	40-140	40-140
<30	<50	<50	<50

Project Sample(s) Spiked: \_\_\_\_\_

**Deviations:**

Compound	%R	%R Limits	RPD	RPD Limits	Samples Affected

**Actions:**

1. If the spike recovery is above the upper control limit (UCL), qualify all positive values in the unspiked sample as estimated (J) and non-detects as estimated (UJ).
2. If the spike recovery is below the lower control limit (LCL), qualify positive values as estimated (J). And non-detects as estimated (UJ).
3. If the spike recovery is <10%, qualify non-detect values as unusable (R)
4. If the RPD does not meet criteria, qualify positive values in the unspiked sample as estimated (J)
5. Use professional judgement to qualify additional samples in the analytical group based on MS/MSD results
6. Use professional judgement for qualification of data for unspiked compounds

Remarks: \_\_\_\_\_

*See attached*

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X. Laboratory Control Sample Information

General LCS Criteria:  
percent recovery (%R)

VOC	SVOC	Pest	PCB
80-120	60-120	50-130	50-130

Laboratory LCS Identifications: \_\_\_\_\_

Deviations:

Compound	Date	%R	Samples Affected/Qualifiers Applied

Actions:

- Action should be based on both the number of compounds outside the criterion and the magnitude of the exceedance.
1. If the LCS recovery is below limits but > one-half the lower limit, qualify values as estimated (J/UJ).
  2. If the LCS recovery is < one-half the lower limit, qualify all data for that analyte as unusable (R).
  3. If the LCS recovery is greater than the upper limit, qualify positive values for that analyte as estimated (J).
  4. If more than half the compounds in this LCS are not within recovery criteria, then qualify associated detected compounds as estimated (J).
  5. Use professional judgement for qualification of data for compounds with no LCS information

Remarks: See attached,  
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# Hold Time Summary

SDG 180-47781-1

Sample Number	Sample Name	Method	Date Collected	Analysis Date	Date Extracted	Days to Analysis
180-47781-1	HD-COD-SW-6-0/1-0	SW846 8260C	9/15/2015	9/22/2015		7
180-47781-10	HD-COD-SW-26-0/1-0	SW846 8260C	9/15/2015	9/23/2015		8
180-47781-11	HD-COD-SW-27-0/1-0	SW846 8260C	9/15/2015	9/23/2015		8
180-47781-12	HD-COD-SW-28-0/1-0	SW846 8260C	9/15/2015	9/23/2015		8
180-47781-13	HD-COD-SW-29-0/1-0	SW846 8260C	9/15/2015	9/23/2015		8
180-47781-14	HD-QC1-0/1-1	SW846 8260C	9/15/2015	9/23/2015		8
180-47781-15	HD-QC1-0/1-2	SW846 8260C	9/15/2015	9/18/2015		3
180-47781-16	HD-MW-136A-459.5/460-0	SW846 8260C	9/15/2015	9/23/2015		8
180-47781-16	HD-MW-136A-459.5/460-0	SW846 8260C	9/15/2015	9/24/2015		9
180-47781-2	HD-COD-SW-7-0/1-0	SW846 8260C	9/15/2015	9/22/2015		7
180-47781-3	HD-COD-SW-8-0/1-0	SW846 8260C	9/15/2015	9/22/2015		7
180-47781-4	HD-COD-SW-9-0/1-0	SW846 8260C	9/15/2015	9/22/2015		7
180-47781-5	HD-COD-SW-12-0/1-0	SW846 8260C	9/15/2015	9/23/2015		8
180-47781-6	HD-COD-SW-13-0/1-0	SW846 8260C	9/15/2015	9/23/2015		8
180-47781-7	HD-COD-SW-15-0/1-0	SW846 8260C	9/15/2015	9/23/2015		8
180-47781-8	HD-COD-SW-16-0/1-0	SW846 8260C	9/15/2015	9/23/2015		8
180-47781-9	HD-COD-SW-17-0/1-0	SW846 8260C	9/15/2015	9/18/2015		3

# Blank Detections

SDG

Sample ID	Sample	Analyte	Result	Method	Units	Qual
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# Qualifier Check

SDG 180-47781-1

Sample ID	Sample	Analyte	Result	5x	10x	Method	Units	Qual
MB 180-142252/1-A	MB 180-142252/1-A	Sodium				SW846 6020A	ug/L	U
180-47781-7	HD-COD-SW-15-0/1-0	1,1,1-Trichloroethane	0.57	2.85	5.7	SW846 8260C	ug/L	J
180-47781-14	HD-QC1-0/1-1	1,1,1-Trichloroethane	0.59	2.95	5.9	SW846 8260C	ug/L	J
180-47781-7	HD-COD-SW-15-0/1-0	1,1-Dichloroethane	0.26	1.3	2.6	SW846 8260C	ug/L	J
180-47781-9	HD-COD-SW-17-0/1-0	1,1-Dichloroethane	0.67	3.35	6.7	SW846 8260C	ug/L	J
180-47781-16	HD-MW-136A-459.5/460-0	1,1-Dichloroethane	3.2	16	32	SW846 8260C	ug/L	J
180-47781-14	HD-QC1-0/1-1	1,1-Dichloroethane	0.25	1.25	2.5	SW846 8260C	ug/L	J
180-47781-7	HD-COD-SW-15-0/1-0	1,1-Dichloroethene	0.77	3.85	7.7	SW846 8260C	ug/L	J
180-47781-9	HD-COD-SW-17-0/1-0	1,1-Dichloroethene	0.84	4.2	8.4	SW846 8260C	ug/L	J
180-47781-16	HD-MW-136A-459.5/460-0	1,1-Dichloroethene	4.0	20	40	SW846 8260C	ug/L	J
180-47781-14	HD-QC1-0/1-1	1,1-Dichloroethene	0.65	3.25	6.5	SW846 8260C	ug/L	J
180-47781-1	HD-COD-SW-6-0/1-0	1,4-Dioxane				SW846 8260C	ug/L	^c
180-47781-2	HD-COD-SW-7-0/1-0	1,4-Dioxane				SW846 8260C	ug/L	^c
180-47781-3	HD-COD-SW-8-0/1-0	1,4-Dioxane				SW846 8260C	ug/L	^c
180-47781-4	HD-COD-SW-9-0/1-0	1,4-Dioxane				SW846 8260C	ug/L	^c
180-47781-1	HD-COD-SW-6-0/1-0	2-Butanone (MEK)				SW846 8260C	ug/L	^c
180-47781-2	HD-COD-SW-7-0/1-0	2-Butanone (MEK)				SW846 8260C	ug/L	^c
180-47781-3	HD-COD-SW-8-0/1-0	2-Butanone (MEK)				SW846 8260C	ug/L	^c
180-47781-4	HD-COD-SW-9-0/1-0	2-Butanone (MEK)				SW846 8260C	ug/L	^c
180-47781-16	HD-MW-136A-459.5/460-0	2-Butanone (MEK)	180	900	1800	SW846 8260C	ug/L	^c
180-47781-1	HD-COD-SW-6-0/1-0	2-Hexanone				SW846 8260C	ug/L	^c
180-47781-2	HD-COD-SW-7-0/1-0	2-Hexanone				SW846 8260C	ug/L	^c
180-47781-3	HD-COD-SW-8-0/1-0	2-Hexanone				SW846 8260C	ug/L	^c
180-47781-4	HD-COD-SW-9-0/1-0	2-Hexanone				SW846 8260C	ug/L	^c
180-47781-1	HD-COD-SW-6-0/1-0	Acetone				SW846 8260C	ug/L	^c
180-47781-3	HD-COD-SW-8-0/1-0	Acetone				SW846 8260C	ug/L	^c
180-47781-11	HD-COD-SW-27-0/1-0	Acetone	4.3	21.5	43	SW846 8260C	ug/L	J
180-47781-12	HD-COD-SW-28-0/1-0	Acetone	3.3	16.5	33	SW846 8260C	ug/L	J
180-47781-2	HD-COD-SW-7-0/1-0	Acetone	2.8	14	28	SW846 8260C	ug/L	J ^c
180-47781-4	HD-COD-SW-9-0/1-0	Acetone	2.7	13.5	27	SW846 8260C	ug/L	J ^c
180-47781-16	HD-MW-136A-459.5/460-0	Acetone	42	210	420	SW846 8260C	ug/L	J ^c
180-47781-16	HD-MW-136A-459.5/460-0	Acrylonitrile	740	3700	7400	SW846 8260C	ug/L	J
180-47781-9	HD-COD-SW-17-0/1-0	Bromomethane				SW846 8260C	ug/L	^c

Sample ID	Sample	Analyte	Result	5x	10x	Method	Units	Qual
180-47781-15	HD-QC1-0/1-2	Bromomethane				SW846 8260C	ug/L	^c
180-47781-9	HD-COD-SW-17-0/1-0	Chloroethane				SW846 8260C	ug/L	^c
180-47781-15	HD-QC1-0/1-2	Chloroethane				SW846 8260C	ug/L	^c
180-47781-5	HD-COD-SW-12-0/1-0	Chloroform	0.20	1	2	SW846 8260C	ug/L	J
180-47781-7	HD-COD-SW-15-0/1-0	Chloroform	0.21	1.05	2.1	SW846 8260C	ug/L	J
180-47781-9	HD-COD-SW-17-0/1-0	Chloroform	0.20	1	2	SW846 8260C	ug/L	J
180-47781-12	HD-COD-SW-28-0/1-0	Chloroform	0.19	0.95	1.9	SW846 8260C	ug/L	J
180-47781-4	HD-COD-SW-9-0/1-0	Chloroform	0.17	0.85	1.7	SW846 8260C	ug/L	J
180-47781-14	HD-QC1-0/1-1	Chloroform	0.23	1.15	2.3	SW846 8260C	ug/L	J
180-47781-1	HD-COD-SW-6-0/1-0	Chloromethane				SW846 8260C	ug/L	^c
180-47781-2	HD-COD-SW-7-0/1-0	Chloromethane				SW846 8260C	ug/L	^c
180-47781-3	HD-COD-SW-8-0/1-0	Chloromethane				SW846 8260C	ug/L	^c
180-47781-4	HD-COD-SW-9-0/1-0	Chloromethane				SW846 8260C	ug/L	^c
180-47781-16	HD-MW-136A-459.5/460-0	Chloromethane				SW846 8260C	ug/L	^c
180-47781-16	HD-MW-136A-459.5/460-0	cis-1,2-Dichloroethene	2500	12500	25000	SW846 8260C	ug/L	F
180-47781-9	HD-COD-SW-17-0/1-0	cis-1,2-Dichloroethene	15	75	150	SW846 8260C	ug/L	F1
180-47781-6	HD-COD-SW-13-0/1-0	cis-1,2-Dichloroethene	0.35	1.75	3.5	SW846 8260C	ug/L	J
180-47781-10	HD-COD-SW-26-0/1-0	cis-1,2-Dichloroethene	0.28	1.4	2.8	SW846 8260C	ug/L	J
180-47781-11	HD-COD-SW-27-0/1-0	cis-1,2-Dichloroethene	0.85	4.25	8.5	SW846 8260C	ug/L	J
180-47781-12	HD-COD-SW-28-0/1-0	cis-1,2-Dichloroethene	0.61	3.05	6.1	SW846 8260C	ug/L	J
180-47781-4	HD-COD-SW-9-0/1-0	cis-1,2-Dichloroethene	0.53	2.65	5.3	SW846 8260C	ug/L	J
180-47781-6	HD-COD-SW-13-0/1-0	Tetrachloroethene	0.37	1.85	3.7	SW846 8260C	ug/L	J
180-47781-8	HD-COD-SW-16-0/1-0	Tetrachloroethene	0.40	2	4	SW846 8260C	ug/L	J
180-47781-11	HD-COD-SW-27-0/1-0	Tetrachloroethene	0.48	2.4	4.8	SW846 8260C	ug/L	J
180-47781-12	HD-COD-SW-28-0/1-0	Tetrachloroethene	0.33	1.65	3.3	SW846 8260C	ug/L	J
180-47781-3	HD-COD-SW-8-0/1-0	Tetrachloroethene	0.23	1.15	2.3	SW846 8260C	ug/L	J
180-47781-4	HD-COD-SW-9-0/1-0	Tetrachloroethene	0.29	1.45	2.9	SW846 8260C	ug/L	J
180-47781-16	HD-MW-136A-459.5/460-0	Tetrachloroethene	2.1	10.5	21	SW846 8260C	ug/L	J
180-47781-16	HD-MW-136A-459.5/460-0	Toluene	37	185	370	SW846 8260C	ug/L	J
180-47781-16	HD-MW-136A-459.5/460-0	trans-1,2-Dichloroethene	4.4	22	44	SW846 8260C	ug/L	J
180-47781-9	HD-COD-SW-17-0/1-0	Trichloroethene	16	80	160	SW846 8260C	ug/L	F1
180-47781-6	HD-COD-SW-13-0/1-0	Trichloroethene	0.57	2.85	5.7	SW846 8260C	ug/L	J
180-47781-8	HD-COD-SW-16-0/1-0	Trichloroethene	0.37	1.85	3.7	SW846 8260C	ug/L	J
180-47781-10	HD-COD-SW-26-0/1-0	Trichloroethene	0.44	2.2	4.4	SW846 8260C	ug/L	J
180-47781-12	HD-COD-SW-28-0/1-0	Trichloroethene	0.83	4.15	8.3	SW846 8260C	ug/L	J
180-47781-13	HD-COD-SW-29-0/1-0	Trichloroethene	0.22	1.1	2.2	SW846 8260C	ug/L	J

Sample ID	Sample	Analyte	Result	5x	10x	Method	Units	Qual
180-47781-2	HD-COD-SW-7-0/1-0	Trichloroethene	0.16	0.8	1.6	SW846 8260C	ug/L	J
180-47781-3	HD-COD-SW-8-0/1-0	Trichloroethene	0.24	1.2	2.4	SW846 8260C	ug/L	J
180-47781-4	HD-COD-SW-9-0/1-0	Trichloroethene	0.53	2.65	5.3	SW846 8260C	ug/L	J
180-47781-9	HD-COD-SW-17-0/1-0	Vinyl chloride				SW846 8260C	ug/L	^c
180-47781-15	HD-QC1-0/1-2	Vinyl chloride				SW846 8260C	ug/L	^c



Initial & Continuing Calibration

Compound	Date	RRF	%RSD	%D	Samples Affected
1,4-Dioxane	8/26/2015	0.0022			R- 1,2,3,4,5,6,7,8,9,10,11,12,13,14,15, - code 3
Isobutyl alcohol	9/18/2015 12:26	0.0095			NA
1,4-Dioxane	9/18/2015 12:26	0.0019			See above
1,4-Dioxane	9/22/2015 13:22	0.0027			See above
1,4-Dioxane	9/24/2015 11:05	0.0026			See above
2,3,6-Trichlorotoluene	9/18/2015 12:26			36.9	NA
2,4,5-Trichlorotoluene	9/18/2015 12:26			25.8	NA
1,2,3-Trichlorobenzene	9/18/2015 12:26			25.5	NA
Chloromethane	9/22/2015 13:22			26.8	None
Acetone	9/22/2015 13:22			47.2	1,3-UJ 2,4-J - code 4
Vinyl acetate	9/22/2015 13:22			54.1	NA
2-Hexanone	9/22/2015 13:22			28.2	None
Dichlorodifluoromethane	9/24/2015 11:05			26	NA
Chloromethane	9/24/2015 11:05			26.1	None
1,3-Butadiene	9/24/2015 11:05			37.6	NA
Acetone	9/24/2015 11:05			44.1	16-UJ - code 4
Vinyl acetate	9/24/2015 11:05			39.1	NA
2-Butanone (MEK)	9/24/2015 11:05			27.6	None
Isobutyl alcohol	9/24/2015 11:05			26.3	NA

Laboratory Control Sample			
Compound	Date	%R	Samples Affected
Chloromethane	9/24/2015	129	16-UJ
1,1,2,2-Tetrachloroethane	9/24/2015	125	16-UJ
1,4-Dioxane	9/24/2015	125	16-UJ
1,4-Dioxane	9/22/2015	124	1,2,3,4- <del>16</del> UJ
Bromomethane	9/18/2015	78	15,9-UJ
Bromomethane	9/23/2015	78	5,6,7,8,10,11,12,13,14,16-UJ
Carbon disulfide	9/23/2015	74	5,6,7,8,10,11,12,13,14,16-UJ
Carbon disulfide	9/18/2015	71	15,9-UJ
Chloroethane	9/18/2015	64	15,9-UJ

code 11

## MS MSD

Compound	%R	%R Limits	RPD	RPD Limits	Samples Affected
Carbon disulfide	69				None
Bromomethane	69				
Chloroethane	67				
1,1-Dichloroethene	66				
Trichloroethene	62				
Carbon disulfide	59				
Chloroethane	59				
Trichloroethene	56				
cis-1,2-Dichloroethene	56				
cis-1,2-Dichloroethene	51				
Tetrachloroethene	10				9-J
Tetrachloroethene	6				9-J
2-Hexanone	132				9,15-UJ

} code 10