CHEMTECH 284 Sheffield Street, Mountainside New Jersey 07092

NEW JERSEY LAB ID#: 20012: NEW YORK LAB ID#: 11376

GC/MS VOA CONFORMANCE/NON-CONFORMANCE SUMMARY

TECH PROJECT NUMBER: O3631 MA	ATRIX: /Water			
DD: 8260-Low,8260D				
		NA	NO	YES
Chromatograms Labeled/Compounds Identified. (Field samples and	d Method Blanks)			\checkmark
GC/MS Tuning Specifications BFB Meet Criteria (NOTE THAT THERE ARE DIFFERENT CRIT ASP CLP, CLP AND NJ)	TERIA FOR NY			✓
GC/MS Tuning Frequency - Performed every 24 hours for 600 series.	es and 12 hours for			\checkmark
				✓
GC/MS Calibration Requirements.			\checkmark	
Bromomethane, Chloroethane, Methylene Chloride, 1,1,2,2-Tetrach	loroethane, 1,2-			
1	1			
Blank Contamination - If yes, list compounds and concentrations in	each blank:		\checkmark	
Surrogate Recoveries Meet Criteria				\checkmark
	 DD: 8260-Low,8260D Chromatograms Labeled/Compounds Identified. (Field samples and GC/MS Tuning Specifications BFB Meet Criteria (NOTE THAT THERE ARE DIFFERENT CRI ASP CLP, CLP AND NJ) GC/MS Tuning Frequency - Performed every 24 hours for 600 series 8000 Series. GC/MS Calibration - Initial Calibration performed before sample analysi 12 hours for 8000 series. GC/MS Calibration Requirements. The %RSD is greater than 15% in the Initial Calibration method (82 Bromomethane, Chloroethane, Methylene Chloride, 1,1,2,2-Tetrach Dibromo-3-Chloropropane, these compounds are passing on Linear The % RSD is greater than 15% in the Initial Calibration method (8 Methylene Chloride, this compound is passing on Linear Regression The Continuous Calibration File ID VY014663.D met the requirem Butanone,2-Hexanone and Methyl Acetate . But associated samples hit for these compounds therefore no corrective action was taken. Blank Contamination - If yes, list compounds and concentrations in 	 DD: 8260-Low,8260D Chromatograms Labeled/Compounds Identified. (Field samples and Method Blanks) GC/MS Tuning Specifications BFB Meet Criteria (NOTE THAT THERE ARE DIFFERENT CRITERIA FOR NY ASP CLP, CLP AND NJ) GC/MS Tuning Frequency - Performed every 24 hours for 600 series and 12 hours for 8000 Series. GC/MS Calibration - Initial Calibration performed before sample analysis and continuing calibration performed within 24 hours of sample analysis for 600 series and 12 hours for 8000 series. GC/MS Calibration Requirements. The %RSD is greater than 15% in the Initial Calibration method (82N071023S.M) for Bromomethane, Chloroethane, Methylene Chloride, 1,1,2,2-Tetrachloroethane, 1,2-Dibromo-3-Chloropropane, these compounds are passing on Linear Regression. The % RSD is greater than 15% in the Initial Calibration method (82Y071323S.M) for Methylene Chloride, this compound is passing on Linear Regression. The Continuous Calibration File ID VY014663.D met the requirements except for 2- Butanone, 2-Hexanone and Methyl Acetate . But associated samples have not positive hit for these compounds therefore no corrective action was taken. Blank Contamination - If yes, list compounds and concentrations in each blank:	DD: 8260-Low,8260D NA Chromatograms Labeled/Compounds Identified. (Field samples and Method Blanks) GC/MS Tuning Specifications BFB Meet Criteria (NOTE THAT THERE ARE DIFFERENT CRITERIA FOR NY ASP CLP, CLP AND NJ) GC/MS Tuning Frequency - Performed every 24 hours for 600 series and 12 hours for 8000 Series. GC/MS Calibration - Initial Calibration performed before sample analysis and continuing calibration performed within 24 hours of sample analysis for 600 series and 12 hours for 8000 series. GC/MS Calibration Requirements. The %RSD is greater than 15% in the Initial Calibration method (82N071023S.M) for Bromomethane, Chlorotehane, Methylene Chloride, 1,1,2,2-Tetrachloroethane, 1,2- Dibromo-3-Chloropropane, these compounds are passing on Linear Regression. The % RSD is greater than 15% in the Initial Calibration method (82Y071323S.M) for Methylene Chloride, this compound is passing on Linear Regression. The Continuous Calibration File ID VY014663.D met the requirements except for 2- Butanone,2-Hexanone and Methyl Acetate . But associated samples have not positive hit for these compounds therefore no corrective action was taken. Blank Contamination - If yes, list compounds and concentrations in each blank:	DD: 8260-Low, 8260D NA NO Chromatograms Labeled/Compounds Identified. (Field samples and Method Blanks) GC/MS Tuning Specifications BFB Meet Criteria (NOTE THAT THERE ARE DIFFERENT CRITERIA FOR NY ASP CLP, CLP AND NJ) GC/MS Tuning Frequency - Performed every 24 hours for 600 series and 12 hours for 8000 Series. GC/MS Calibration - Initial Calibration performed before sample analysis and continuing calibration performed within 24 hours of sample analysis for 600 series and 12 hours for 8000 series. GC/MS Calibration Requirements. The % RSD is greater than 15% in the Initial Calibration method (82N071023S.M) for Bromomethane, Chloroethane, Methylene Chloride, 1,1,2,2-Tetrachloroethane, 1,2- Dibromo-3-Chloropropane, these compounds are passing on Linear Regression. The % RSD is greater than 15% in the Initial Calibration method (82Y071323S.M) for Methylene Chloride, this compound is passing on Linear Regression. The Continuous Calibration File ID VY014663.D met the requirements except for 2- Butanone,2-Hexanone and Methyl Acctate . But associated samples have not positive hit for these compounds therefore no corrective action was taken. Blank Contamination - If yes, list compounds and concentrations in each blank:

If not met, list those compounds and their recoveries which fall outside the acceptable ranges.

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GC/MS VOA CONFORMANCE/NON-CONFORMANCE SUMMARY (CONTINUED)

		NA	NO	YES
8.	Matrix Spike/Matrix Spike Duplicate Recoveries Meet Criteria If not met, list those compounds and their recoveries which fall outside the acceptable range.		✓	
	The MS {O3631-02MS} with File ID: VY014684.D recoveries met the requirements for all compounds except for 1,1,2-Trichloroethane[63%], 1,2-Dibromoethane[57%], 2-Butanone[34%], 2-Hexanone[38%], 4-Methyl-2-Pentanone[46%], Acetone[32%] and Bromoform[63%] due to matrix interference.			
	The MSD {O3631-03MSD} with File ID: VY014685.D recoveries met the acceptable requirements except for Methyl Acetate[178%] due to matrix interference .			
	The RPD for {O3631-03MSD} with File ID: VY014685.D met criteria except for 1,1,2,2-Tetrachloroethane[56%], 1,1,2-Trichloroethane[54%], 1,2,3- Trichlorobenzene[31%], 1,2-Dibromo-3-Chloropropane[79%], 1,2- Dibromoethane[60%], 1,2-Dichloroethane[43%], 1,2-Dichloropropane[24%], 2- Butanone[98%], 2-Hexanone[89%], 4-Methyl-2-Pentanone[81%], Acetone[95%], Bromochloromethane[41%], Bromodichloromethane[28%], Bromoform[57%], Chloroform[21%], cis-1,2-Dichloroethene[21%], cis-1,3-Dichloropropene[30%], Dibromochloromethane[46%], Methyl Acetate[84%], Methyl tert-butyl Ether[55%], Methylene Chloride[46%] and t-1,3-Dichloropropene[42%]. Due to difference in MS and MSD results.			
9.	Internal Standard Area/Retention Time Shift Meet Criteria Comments: The Internal Standards Areas met the acceptable requirements except for SB-02(2-4)MS. No internal standard failed in MSD therefore no corrective action was taken.		✓	
10.	Analysis Holding Time Met If not met, list number of days exceeded for each sample:			✓
ADDITIO	DNAL COMMENTS:			
This data	package has been revised to report TICs as per client request.			
This data	package has been revised due to data package type changed as per client request.			
Samples	for MS/MSD for VOC analysis were not provided with this set of samples. The Blank Spike	e Duplic	ate is	
reported	with the data.			
Trip Blan	ak was not provided with this set of samples.			
The soil s	samples results are based on a dry weight basis.			
Please us	e %D calculated based on Avg RF and CCRF for all compounds using Average Response F	Factor w	hen the	
%RSD va	alue for a compound is $<15\%$ for the Initial Calibration curve and use $\%D$ calculated based	on Amo	ount adde	d
and Calcu	ulated amount for all compounds using Linear Regression when the %RSD value for a com-	pound is	s > 15%	for
the Initial	l Calibration curve for SW-846 analysis.			