## ALLIANCE 284 Sheffield Street, Mountainside New Jersey 07092 NEW JERSEY LAB ID#: 20012: NEW YORK LAB ID#: 11376

## GC/MS VOA CONFORMANCE/NON-CONFORMANCE SUMMARY

ORDER ID: Q2230	
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MATRIX: Water

METHOD: 8260D

		NA	NO	YES
1.	Chromatograms Labeled/Compounds Identified. (Field samples and Method Blanks)			$\checkmark$
2.	GC/MS Tuning Specifications BFB Meet Criteria (NOTE THAT THERE ARE DIFFERENT CRITERIA FOR NY ASP CLP, CLP AND NJ)			✓
3.	GC/MS Tuning Frequency - Performed every 24 hours for 600 series and 12 hours for 8000 Series.			✓
4.	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.			✓
5.	GC/MS Calibration Requirements.			$\checkmark$
	The Initial Calibration met the requirements.			
6.	The Continuous Calibration met the requirements. Blank Contamination - If yes, list compounds and concentrations in each blank:		$\checkmark$	
7.	Surrogate Recoveries Meet Criteria			$\checkmark$
	If not met, list those compounds and their recoveries which fall outside the acceptable ranges.			
8.	Matrix Spike/Matrix Spike Duplicate Recoveries Meet Criteria		$\checkmark$	
	If not met, list those compounds and their recoveries which fall outside the acceptable range.			
	The MS {Q2230-03MS} with File ID: VX046641.D recoveries met the requirements for all compounds except for 1,2,4-Trimethylbenzene[148%], Isopropylbenzene[123%], N-propylbenzene[122%], tert-Butylbenzene[124%], Toluene[121%], m/p-Xylenes[134%] and o-Xylene[149%] due to matrix interference.			
	The MSD {Q2230-04MSD} with File ID: VX046642.D recoveries met the acceptable requirements except for 1,2,4-Trimethylbenzene[168%] and o-Xylenes [151%] due to matrix interference.			
9.	Internal Standard Area/Retention Time Shift Meet Criteria			$\checkmark$
	Comments:			
10.	Analysis Holding Time Met			$\checkmark$

If not met, list number of days exceeded for each sample:

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

NA NO YES

ADDITIONAL COMMENTS:

Please use %D calculated based on Avg RF and CCRF for all compounds using Average Response Factor when the %RSD value for a compound is <20% for the Initial Calibration curve and use %D calculated based on Amount added and Calculated amount for all compounds using Linear Regression when the %RSD value for a compound is >20% for the Initial Calibration curve for SW-846 analysis.

**QA REVIEW** 

Date