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

GC/MS SEMI-VOLATILE ANALYSIS CONFORMANCE/NON-CONFORMANCE SUMMARY

MATRIX: Solid

ORDER ID: Q3363

MET	HOD: 8270E/3541			
1.	Chromatograms Labeled/Compounds Identified. (Field samples and Method Blanks)	NA	NO	YES ✓
2.	GC/MS Tuning Specifications. DFTPP 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 within 30 days 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.		✓	
	The %RSD is greater than 20% in the Initial Calibration (Method 8270-BF100625.M) for 2,4-Dinitrophenol this Compounds is passing on Linear regression. The Continuous Calibration File ID BF143984.D met the requirements except for Di-n-			
	octyl phthalate, Associated sample does not have hit for this compound, Therefor no further corrective action was taken.			
	The Continuous Calibration File ID BG064531.D met the requirements except for Hexachlorocyclopentadiene, Associated sample does not have hit for this compound, Therefor no further corrective action was taken.			
6.	Blank Contamination - If yes, list compounds and concentrations in each blank:		✓	
7.	Surrogate Recoveries Meet Criteria			✓
	If not met, list those compounds and their recoveries which fall outside the acceptable ranges.			

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GC/MS SEMI-VOLATILE ANALYSIS 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 out range.	side the acceptable	
	The MSD {Q3363-05MSD} with File ID: BG064540.D recoveries requirements for all compounds except for Acenaphthylene[127%] and N-Nitrosodiphenylamine[120%], Recovery failed due to matrix	Carbazole[120%]	
	The Blank Spike met requirements for all compounds.		
	The RPD for {Q3363-05MSD} with File ID: BG064540.D met crit Dioxane[25%], 2,2-oxybis(1-Chloropropane)[22%], 2,4-Dichlorop Dimethylphenol[26%], 2-Chlorophenol[23%], 2-Methylphenol[23%] Methylphenols[26%], 4-Chloro-3-methylphenol[23%], Acetopheno Atrazine[21%], Benzaldehyde[28%], Benzo(a)anthracene[26%], Benzo(g,h,i)perylene[21%], Benzo(k)fluoranthene[23%], bis(2-Chloroethyl)ether[21%], bis(2-Ethylhexyl)phthalate[21%], Caprola Chrysene[22%], Dibenz(a,h)anthracene[21%], Hexachloroethane[2 Isophorone[21%], Nitrobenzene[23%], N-Nitroso-di-n-propylamin Nitrosodiphenylamine[21%] and Pyrene[24%], RPD failed due to c MS and MSD results.	henol[21%], 2,4- %], 3+4- one[21%], enzo(a)pyrene[21%], ctam[29%], 3%], e[25%], N-	
9.	Internal Standard Area/Retention Time Shift Meet Criteria		v
	Comments:		
10.	Extraction Holding Time Met		v
	If not met, list number of days exceeded for each sample:		
11.	Analysis Holding Time Met		•
	If not met, list number of days exceeded for each sample:		
ADDI	TIONAL COMMENTS:		
The F	orm 6 is not included in the data package because the Initial Calibration	n was performed using 7 points.	
The so	oil samples results are based on a dry weight basis.		
QA R	EVIEW	Date	