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

ORDER	ID:	01872
OILDDIC	· · ·	V 10/2

MATRIX: Solid

METHOD: 8270E/3541

		NA	NO	YES
1.	Chromatograms Labeled/Compounds Identified. (Field samples and Method Blanks)			\checkmark
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 (8270-BM042825.M) for 2,4- Dinitrophenol and 4-Nitrophenol these compound are passing on Linear Regression.			
	The % RSD is greater than 20% in the Initial Calibration (8270-BP041425.M) for Benzidine, only QC sample run under this Initial Calibration.			
	The % RSD is greater than 20% in the Initial Calibration (8270-BP051325.M) for Benzoic acid, 2,4-Dinitrophenol, 4-Nitrophenol these compound are passing on Linear Regression.			
	The Continuous Calibration File ID BP024568.D met the requirements except for 2,4,6- Tribromophenol . associated samples have no hit for this compound, Therefor no further corrective action was taken.			
6.	Blank Contamination - If yes, list compounds and concentrations in each blank:		\checkmark	

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		NA	NO	YES
7.	Surrogate Recoveries Meet Criteria		\checkmark	
	If not met, list those compounds and their recoveries which fall outside the acceptable ranges.			
	The Surrogate recoveries met the acceptable criteria except for PB167904BS [2,4 and6-Tribromophenol - 117%] is marginally biased high, Therefor no corrective action was taken.			
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 Blank Spike for {PB167904BS} with File ID: BP024570.D met requirements for all samples except for 2,4,5-Trichlorophenol[100%], 4-Bromophenyl-phenylether[106%], Benzidine[112%], Hexachlorobenzene[106%], Hexachlorobutadiene[100%], Hexachlorocyclopentadiene[185%] and Pentachlorophenol[106%]. Recovery of a very few compounds are slightly biased high but their hit is not present in any associated samples, Therefor no further corrective action was taken.			
9.	Internal Standard Area/Retention Time Shift Meet Criteria			\checkmark
	Comments:			
10.	Extraction Holding Time Met			\checkmark
	If not met, list number of days exceeded for each sample:			
11.	Analysis Holding Time Met			\checkmark
	If not met, list number of days exceeded for each sample:			

ADDITIONAL COMMENTS:

Sample HW0425-PT-BNA-SOIL was diluted due to high concentration.

The Form 6 is not included in the data package because the Initial Calibration was performed using 7 points.

The soil samples results are based on a dry weight basis.

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

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NA NO YES

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