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: TCLP

ORDER ID: Q2641

NA amples and Method Blanks) FOR NY ASP CLP,	NO	YES ✓
or 600 series and 12 hours for		✓
a 30 days before sample hours of sample analysis for		✓
		✓
ntrations in each blank:	✓	
ch fall outside the acceptable		✓
	amples and Method Blanks) OR NY ASP CLP, or 600 series and 12 hours for a 30 days before sample hours of sample analysis for attrations in each blank:	amples and Method Blanks) FOR NY ASP CLP, or 600 series and 12 hours for a 30 days before sample hours of sample analysis for attrations in each blank:

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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 outside the acceptable range.
	The MS {Q2646-03MS} with File ID: BF143199.D recoveries met the requirements for all compounds except for 2,4,6-Trichlorophenol[72%], 2-Methylphenol[158%] and Pyridine[0%]. Recovery failed due to matrix interference, Therefor no further corrective action was taken.
	The MSD {Q2646-03MSD} with File ID: BF143200.D recoveries met the requirements for all compounds except for 2,4,6-Trichlorophenol[76%], 2-Methylphenol[152%] and Pyridine[0%]. Recovery failed due to matrix interference, Therefor no further corrective action was taken.
	The Blank Spike met requirements for all compounds.
9.	Internal Standard Area/Retention Time Shift Meet Criteria ✓
	Comments:
10.	Extraction Holding Time Met
	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:
ADDIT	IONAL COMMENTS:
	m 6 is not included in the data package because the Initial Calibration was performed using 7 points.
	se %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 Cal	culated amount for all compounds using Linear Regression when the %RSD value for a compound is > 20% for
the Initi	al Calibration curve for SW-846 analysis.

Date

QA REVIEW