### 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

CHEMTECH PROJECT NUMBER: P5284 MATRIX: Water METHOD: 8270-Modified/3510 NO NA YES 1. Chromatograms Labeled/Compounds Identified. (Field samples and Method Blanks) 2. GC/MS Tuning Specifications. DFTPP 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 3. 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 Initial Calibration met the requirements. The Continuous Calibration File ID BN035731.D met the requirements except for 2,4,6-Tribromophenol, The failure compound not associated with the client parameters list, therefore no corrective action was taken. The Continuous Calibration File ID BN035747.D met the requirements except for 2,4,6-Tribromophenol,2-Fluorobiphenyl,2-Fluorophenol,Nitrobenzene-d5,Phenol-d6 and Terphenyl-d14, The failure compound not associated with the client parameters list, therefore no corrective action was taken. The Continuous Calibration File ID BN035749.D met the requirements except for Nitrobenzene-d5, The failure compound not associated with the client parameters list, therefore no corrective action was taken. The Continuous Calibration File ID BN035794.D met the requirements except for 2,4,6-Tribromophenol, The failure compound not associated with the client parameters list, therefore no corrective action was taken.. The Continuous Calibration File ID BN035845.D met the requirements except for Nitrobenzene-d5, The failure compound not associated with the client parameters list, therefore no corrective action was taken. .

Blank Contamination - If yes, list compounds and concentrations in each blank:

6.

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NA

NO

YES

7. Surrogate Recoveries Meet Criteria If not met, list those compounds and their recoveries which fall outside the acceptable ranges. The Surrogate recoveries met the acceptable criteria except for RW5-SP100-20241212 [Nitrobenzene-d5 - 123%, Terphenyl-d14 - 169%], RW5-SP100-20241212DL [2-Fluorobiphenyl - 139%, Fluoranthene-d10 - 179%, Nitrobenzene-d5 - 185%, Terphenyl-d14 - 231%], RW5-SP201-20241212 [Terphenyl-d14 - 154%], RW5-SP303-20241212 [Nitrobenzene-d5 - 114%, Terphenyl-d14 - 162%], PB165667BL [Terphenyl-d14 - 140%] and PB165667BS [Nitrobenzene-d5 - 116%], The failure surrogates not associated with the client parameters list, therefore no corrective action was taken. 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 {P5280-04MS} with File ID: BN035735.D recoveries met the requirements for all compounds except for 1,4-Dioxane[133%], marginally high due to matrix interference. therefore, no corrective action is required. The MSD {P5280-05MSD} with File ID: BN035736.D recoveries met the acceptable requirements except for 1,4-Dioxane[138%], marginally high due to matrix interference therefore, no corrective action is required. The Blank Spike met requirements for all samples. 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: ADDITIONAL COMMENTS: The laboratory certifies that the all-electronic diskette deliverable exactly match the data summary forms (i.e. Form Is)."

Sample RW5-SP100-20241212 was diluted due to high concentration.

The not QT review data is reported in the Miscellaneous.

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

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NA NO

YES

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 < 15\% \ for \ the \ Initial \ Calibration \ curve \ and \ use \ \% D \ calculated \ based \ on \ Amount \ add \ on \ Amount \ add \ on \ Amount \ add \ on \ and \ on \ $	ed
and Calculated amount for all compounds using Linear Regression when the %RSD value for a compound is $> 15\%$	for
the Initial Calibration curve for SW-846 analysis.	
QA REVIEW Date	