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

CHEMTECH PROJECT NUMBER: P5164 MATRIX: Water

METHOD: 8260D

| | | 27.4 | 110 | MEG |
|----|---|------|-----|----------|
| 1. | Chromatograms Labeled/Compounds Identified. (Field samples and Method Blanks) | NA | NO | YES |
| 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. | | | √ |
| | The %RSD is greater than 20% in the Initial Calibration method (82X1121W.M) for Bromoform this compound is passing on Quadratic Regression. | | | · |
| | The Continuous Calibration File ID VX044174.D met the requirements except for Bromodichloromethane, Dibromochloromethane, t-1,3-Dichloropropene and Trichlorofluoromethane failing high but no positive hit in associated samples therefore no corrective action taken. | | | |
| | The Continuous Calibration File ID VX044197.D met the requirements except for 2-Hexanone and 4-Methyl-2-Pentanone failing high but no positive hit in associated samples therefore no corrective action 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. | | | |
| 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 Blank Spike met requirements for all samples . | | | |
| | The Blank Spike Duplicate met requirements for all samples . | | | |
| | The RPD for {VX1209WBSD01} with File ID: VX044194.D met criteria except for Chloroethane[24%] due to difference in results of BS-BSD. | | | |

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

YES

| | | NA | NO | Y |
|------------|--|----------|-----------|-------|
| 9. | Internal Standard Area/Retention Time Shift Meet Criteria | | | ✓ |
| | Comments: | | | |
| 10. | Analysis Holding Time Met | | | ✓ |
| | If not met, list number of days exceeded for each sample: | | | |
| ADDITIO | ONAL COMMENTS: | | | |
| Samples | RE108D2-20241206, RE105D2-20241206 was analyzed with straight 20X dilution due to | past his | tory of t | hese |
| samples o | containing high amounts of compound Trichloroethene. | | | |
| The Sam | ple #RE108D2-20241206, RE105D1-20241206 have the concentration of target compound | below | Method | |
| detection | n limits, therefore it is not reported as Hit in Form1. | | | |
| Samples | for MS/MSD for VOC analysis were not provided with this set of samples. The Blank Spik | e Dupli | cate is | |
| reported | with the data. | | | |
| The not (| QT review data is reported in the Miscellaneous. | | | |
| Please us | se %D calculated based on Avg RF and CCRF for all compounds using Average Response l | Factor v | when the | ; |
| %RSD v | value for a compound is <20% for the Initial Calibration curve and use %D calculated based | on Am | ount ado | led |
| and Calc | rulated amount for all compounds using Linear Regression when the %RSD value for a com- | pound | is > 20% | o for |
| the Initia | ll Calibration curve for SW-846 analysis. | | | |
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| | | | | |
| QA REV | VIEW Date | | | |