

6C

SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: PSEG03  
Lab Code: CHEM Case No.: P2687 SAS No.: P2687 SDG No.: P2687  
Instrument ID: BNA\_P Calibration Date(s): 05/29/2024 05/29/2024  
Calibration Time(s): 09:59 15:04

| LAB FILE ID:          |        | RRF2.5 = BP020469.D |        |        | RRF005 = BP020470.D |        | RRF010 = BP020471.D |       |
|-----------------------|--------|---------------------|--------|--------|---------------------|--------|---------------------|-------|
|                       |        | RRF020 = BP020472.D |        |        | RRF040 = BP020473.D |        | RRF050 = BP020474.D |       |
| COMPOUND              | RRF2.5 | RRF005              | RRF010 | RRF020 | RRF040              | RRF050 | RRF                 | % RSD |
| Pyridine              |        | 1.188               | 1.265  | 1.330  | 1.291               | 1.266  | 1.269               | 3.6   |
| 2-Fluorophenol        |        | 1.090               | 1.121  | 1.167  | 1.131               | 1.104  | 1.117               | 2.7   |
| Phenol-d6             |        | 1.561               | 1.575  | 1.656  | 1.600               | 1.551  | 1.575               | 3.2   |
| 1,4-Dichlorobenzene   |        | 1.568               | 1.534  | 1.578  | 1.504               | 1.433  | 1.499               | 4.4   |
| 2-Methylphenol        |        | 1.061               | 1.091  | 1.171  | 1.122               | 1.073  | 1.095               | 4.0   |
| 3+4-Methylphenols     |        | 1.398               | 1.442  | 1.589  | 1.535               | 1.476  | 1.489               | 4.5   |
| Nitrobenzene-d5       |        | 0.355               | 0.363  | 0.379  | 0.371               | 0.362  | 0.365               | 2.6   |
| Hexachloroethane      |        | 0.563               | 0.553  | 0.575  | 0.552               | 0.534  | 0.548               | 3.5   |
| Nitrobenzene          |        | 0.375               | 0.368  | 0.390  | 0.372               | 0.363  | 0.371               | 2.9   |
| Hexachlorobutadiene   |        | 0.218               | 0.220  | 0.218  | 0.213               | 0.209  | 0.215               | 1.9   |
| 2,4,6-Trichlorophenol |        | 0.283               | 0.311  | 0.365  | 0.364               | 0.366  | 0.350               | 10.9  |
| 2-Fluorobiphenyl      |        | 1.455               | 1.393  | 1.455  | 1.312               | 1.270  | 1.342               | 7.3   |
| 2,4,5-Trichlorophenol |        | 0.351               | 0.375  | 0.418  | 0.420               | 0.411  | 0.404               | 7.4   |
| 2,4-Dinitrotoluene    |        | 0.296               | 0.351  | 0.416  | 0.416               | 0.425  | 0.396               | 13.3  |
| 2,4,6-Tribromophenol  |        | 0.177               | 0.195  | 0.214  | 0.220               | 0.227  | 0.208               | 8.8   |
| Hexachlorobenzene     |        | 0.241               | 0.238  | 0.243  | 0.236               | 0.234  | 0.241               | 2.4   |
| Pentachlorophenol     |        |                     | 0.120  | 0.136  | 0.146               | 0.150  | 0.146               | 11.3  |
| Terphenyl-d14         |        | 1.238               | 1.205  | 1.218  | 1.277               | 1.170  | 1.201               | 5.0   |

All other compounds must meet a minimum RRF of 0.010.

Form VI SV-1