

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name: CHEMTECH Contract: JAC005

Lab Code: CHEM Case No.: P3429 SAS No.: P3429 SDG No.: P3429

Instrument ID: BNA\_F Calibration Date/Time: 08/10/2024 10:41

Lab File ID: BF138901.D Init. Calib. Date(s): 07/30/2024 07/30/2024

EPA Sample No.: SSTDCCC040 Init. Calib. Time(s): 12:54 16:29

GC Column: DB-UI ID: 0.18 (mm)

| COMPOUND                     | RRF   | RRF040 | MIN<br>RRF | %D    | MAX%D |
|------------------------------|-------|--------|------------|-------|-------|
| Pyridine                     | 1.374 | 1.245  |            | -9.4  |       |
| 2-Fluorophenol               | 1.296 | 1.267  |            | -2.2  |       |
| Benzaldehyde                 | 1.043 | 0.916  |            | -12.2 |       |
| Phenol-d6                    | 1.740 | 1.736  |            | -0.2  |       |
| 2-Methylphenol               | 1.126 | 1.130  |            | 0.4   |       |
| Acetophenone                 | 0.490 | 0.511  |            | 4.3   |       |
| 3+4-Methylphenols            | 1.444 | 1.532  |            | 6.1   |       |
| Nitrobenzene-d5              | 0.409 | 0.425  |            | 3.9   |       |
| Nitrobenzene                 | 0.416 | 0.425  |            | 2.2   |       |
| 2,4-Dichlorophenol           | 0.275 | 0.287  |            | 4.4   | 20.0  |
| Naphthalene                  | 1.053 | 1.059  |            | 0.6   |       |
| Hexachlorobutadiene          | 0.192 | 0.208  |            | 8.3   | 20.0  |
| 2-Methylnaphthalene          | 0.665 | 0.691  |            | 3.9   |       |
| 2,4,6-Trichlorophenol        | 0.339 | 0.337  |            | -0.6  | 20.0  |
| 2-Fluorobiphenyl             | 1.331 | 1.376  |            | 3.4   |       |
| 2,4,5-Trichlorophenol        | 0.370 | 0.377  |            | 1.9   |       |
| Acenaphthylene               | 1.652 | 1.642  |            | -0.6  |       |
| Acenaphthene                 | 1.111 | 1.104  |            | -0.6  | 20.0  |
| Dibenzofuran                 | 1.568 | 1.590  |            | 1.4   |       |
| Fluorene                     | 1.249 | 1.309  |            | 4.8   |       |
| 2,4,6-Tribromophenol         | 0.164 | 0.177  |            | 7.9   |       |
| Hexachlorobenzene            | 0.224 | 0.229  |            | 2.2   |       |
| Pentachlorophenol            | 0.101 | 0.121  |            | 19.8  | 20.0  |
| Phenanthrene                 | 1.030 | 1.051  |            | 2.0   |       |
| Carbazole                    | 0.875 | 0.883  |            | 0.9   |       |
| Di-n-butylphthalate          | 0.984 | 1.131  |            | 14.9  |       |
| Fluoranthene                 | 0.961 | 0.969  |            | 0.8   | 20.0  |
| Pyrene                       | 1.883 | 2.088  |            | 10.9  |       |
| Terphenyl-d14                | 1.195 | 1.360  |            | 13.8  |       |
| Benzo (a) anthracene         | 1.377 | 1.402  |            | 1.8   |       |
| Chrysene                     | 1.243 | 1.236  |            | -0.6  |       |
| Bis (2-ethylhexyl) phthalate | 0.883 | 0.871  |            | -1.4  |       |
| Benzo (b) fluoranthene       | 1.240 | 1.227  |            | -1.0  |       |
| Benzo (k) fluoranthene       | 1.073 | 1.114  |            | 3.8   |       |
| Benzo (a) pyrene             | 1.043 | 1.061  |            | 1.7   | 20.0  |
| Indeno (1,2,3-cd) pyrene     | 1.433 | 1.363  |            | -4.9  |       |
| Dibenzo (a,h) anthracene     | 1.177 | 1.113  |            | -5.4  |       |
| Benzo (g,h,i) perylene       | 1.221 | 1.134  |            | -7.1  |       |
| 1,4-Dioxane                  | 0.567 | 0.483  |            | -14.8 | 20.0  |

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 GC Column: DB-UI ID: 0.18 (mm)

| COMPOUND            | RRF   | RRF040 | MIN<br>RRF | %D  | MAX%D |
|---------------------|-------|--------|------------|-----|-------|
| 1-Methylnaphthalene | 0.652 | 0.672  |            | 3.1 |       |

All other compounds must meet a minimum RRF of 0.010.