

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

|                 |                   |                        |                              |
|-----------------|-------------------|------------------------|------------------------------|
| Lab Name:       | <u>Alliance</u>   | Contract:              | <u>ENTA05</u>                |
| Lab Code:       | <u>ACE</u>        | SDG No.:               | <u>Q3618</u>                 |
| Instrument ID:  | <u>BNA_P</u>      | Calibration Date/Time: | <u>11/13/2025 13:16</u>      |
| Lab File ID:    | <u>BP026126.D</u> | Init. Calib. Date(s):  | <u>10/29/2025 10/29/2025</u> |
| EPA Sample No.: | <u>SSTDCCC040</u> | Init. Calib. Time(s):  | <u>09:26 14:13</u>           |
| GC Column:      | <u>ZB-GR</u>      | ID:                    | <u>0.25</u> (mm)             |

| COMPOUND                    | RRF   | RRF040 | MIN RRF | %D    | MAX%D |
|-----------------------------|-------|--------|---------|-------|-------|
| 2-Fluorophenol              | 1.212 | 1.262  |         | 4.1   |       |
| Benzaldehyde                | 0.802 | 0.738  |         | -8.0  |       |
| Phenol-d6                   | 1.543 | 1.607  |         | 4.1   |       |
| Phenol                      | 1.713 | 1.751  |         | 2.2   | 20.0  |
| bis(2-Chloroethyl)ether     | 1.352 | 1.349  |         | -0.2  |       |
| 2-Chlorophenol              | 1.342 | 1.454  |         | 8.3   |       |
| 2-Methylphenol              | 1.121 | 1.157  |         | 3.2   |       |
| 2,2-oxybis(1-Chloropropane) | 2.043 | 1.829  |         | -10.5 |       |
| Acetophenone                | 0.506 | 0.502  |         | -0.8  |       |
| 3+4-Methylphenols           | 1.560 | 1.591  |         | 2.0   |       |
| n-Nitroso-di-n-propylamine  | 0.958 | 0.967  | 0.050   | 0.9   |       |
| Nitrobenzene-d5             | 0.321 | 0.351  |         | 9.3   |       |
| Hexachloroethane            | 0.533 | 0.568  |         | 6.6   |       |
| Nitrobenzene                | 0.339 | 0.359  |         | 5.9   |       |
| Isophorone                  | 0.687 | 0.684  |         | -0.4  |       |
| 2-Nitrophenol               | 0.104 | 0.183  |         | 76.0  | 20.0  |
| 2,4-Dimethylphenol          | 0.289 | 0.290  |         | 0.3   |       |
| bis(2-Chloroethoxy)methane  | 0.433 | 0.432  |         | -0.2  |       |
| 2,4-Dichlorophenol          | 0.305 | 0.336  |         | 10.2  | 20.0  |
| Naphthalene                 | 1.091 | 1.092  |         | 0.1   |       |
| 4-Chloroaniline             | 0.419 | 0.421  |         | 0.5   |       |
| Hexachlorobutadiene         | 0.211 | 0.220  |         | 4.3   | 20.0  |
| Caprolactam                 | 0.107 | 0.113  |         | 5.6   |       |
| 4-Chloro-3-methylphenol     | 0.320 | 0.343  |         | 7.2   | 20.0  |
| 2-Methylnaphthalene         | 0.763 | 0.760  |         | -0.4  |       |
| Hexachlorocyclopentadiene   | 0.227 | 0.250  | 0.050   | 10.1  |       |
| 2,4,6-Trichlorophenol       | 0.367 | 0.429  |         | 16.9  | 20.0  |
| 2-Fluorobiphenyl            | 1.392 | 1.381  |         | -0.8  |       |
| 2,4,5-Trichlorophenol       | 0.419 | 0.476  |         | 13.6  |       |
| 1,1-Biphenyl                | 1.531 | 1.521  |         | -0.7  |       |
| 2-Chloronaphthalene         | 1.205 | 1.207  |         | 0.2   |       |
| 2-Nitroaniline              | 0.272 | 0.333  |         | 22.4  |       |
| Dimethylphthalate           | 1.455 | 1.498  |         | 3.0   |       |
| Acenaphthylene              | 1.756 | 1.777  |         | 1.2   |       |
| 2,6-Dinitrotoluene          | 0.237 | 0.307  |         | 29.5  |       |
| 3-Nitroaniline              | 0.291 | 0.350  |         | 20.3  |       |
| Acenaphthene                | 1.206 | 1.181  |         | -2.1  | 20.0  |
| 2,4-Dinitrophenol           | 0.100 | 0.156  | 0.050   | 56.0  |       |
| 4-Nitrophenol               | 0.290 | 0.315  | 0.050   | 8.6   |       |

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| Instrument ID:  | <u>BNA_P</u>      | Calibration Date/Time: | <u>11/13/2025 13:16</u>      |
| Lab File ID:    | <u>BP026126.D</u> | Init. Calib. Date(s):  | <u>10/29/2025 10/29/2025</u> |
| EPA Sample No.: | <u>SSTDCCC040</u> | Init. Calib. Time(s):  | <u>09:26 14:13</u>           |
| GC Column:      | <u>ZB-GR</u>      | ID:                    | <u>0.25</u> (mm)             |

| COMPOUND                   | RRF   | RRF040 | MIN<br>RRF | %D   | MAX%D |
|----------------------------|-------|--------|------------|------|-------|
| Dibenzofuran               | 1.822 | 1.821  |            | -0.1 |       |
| 2,4-Dinitrotoluene         | 0.347 | 0.455  |            | 31.1 |       |
| Diethylphthalate           | 1.463 | 1.510  |            | 3.2  |       |
| 4-Chlorophenyl-phenylether | 0.713 | 0.716  |            | 0.4  |       |
| Fluorene                   | 1.427 | 1.425  |            | -0.1 |       |
| 4-Nitroaniline             | 0.311 | 0.363  |            | 16.7 |       |
| 4,6-Dinitro-2-methylphenol | 0.077 | 0.118  |            | 53.2 |       |
| n-Nitrosodiphenylamine     | 0.625 | 0.616  |            | -1.4 | 20.0  |
| 2,4,6-Tribromophenol       | 0.216 | 0.264  |            | 22.2 |       |
| 4-Bromophenyl-phenylether  | 0.222 | 0.228  |            | 2.7  |       |
| Hexachlorobenzene          | 0.253 | 0.264  |            | 4.3  |       |
| Atrazine                   | 0.211 | 0.218  |            | 3.3  |       |
| Pentachlorophenol          | 0.154 | 0.179  |            | 16.2 | 20.0  |
| Phenanthrene               | 1.164 | 1.146  |            | -1.5 |       |
| Anthracene                 | 1.154 | 1.163  |            | 0.8  |       |
| Carbazole                  | 1.092 | 1.095  |            | 0.3  |       |
| Di-n-butylphthalate        | 1.222 | 1.322  |            | 8.2  |       |
| Fluoranthene               | 1.358 | 1.373  |            | 1.1  | 20.0  |
| Pyrene                     | 1.353 | 1.357  |            | 0.3  |       |
| Terphenyl-d14              | 0.984 | 1.001  |            | 1.7  |       |
| Butylbenzylphthalate       | 0.432 | 0.579  |            | 34.0 |       |
| 3,3-Dichlorobenzidine      | 0.443 | 0.473  |            | 6.8  |       |
| Benzo (a) anthracene       | 1.374 | 1.382  |            | 0.6  |       |
| Chrysene                   | 1.302 | 1.288  |            | -1.1 |       |
| Bis(2-ethylhexyl)phthalate | 0.718 | 0.863  |            | 20.2 |       |
| Di-n-octyl phthalate       | 1.191 | 1.511  |            | 26.9 | 20.0  |
| Benzo (b) fluoranthene     | 1.242 | 1.268  |            | 2.1  |       |
| Benzo (k) fluoranthene     | 1.268 | 1.258  |            | -0.8 |       |
| Benzo (a) pyrene           | 1.141 | 1.182  |            | 3.6  | 20.0  |
| Indeno (1,2,3-cd) pyrene   | 1.476 | 1.539  |            | 4.3  |       |
| Dibenzo (a,h) anthracene   | 1.202 | 1.253  |            | 4.2  |       |
| Benzo (g,h,i) perylene     | 1.156 | 1.203  |            | 4.1  |       |
| 1,2,4,5-Tetrachlorobenzene | 0.621 | 0.643  |            | 3.5  |       |
| 1,4-Dioxane                | 0.511 | 0.503  |            | -1.6 | 20.0  |
| 2,3,4,6-Tetrachlorophenol  | 0.330 | 0.400  |            | 21.2 |       |

All other compounds must meet a minimum RRF of 0.010.