

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

|                 |                   |                        |                              |
|-----------------|-------------------|------------------------|------------------------------|
| Lab Name:       | <u>Alliance</u>   | Contract:              | <u>CORE02</u>                |
| Lab Code:       | <u>ACE</u>        | SDG No.:               | <u>Q3615</u>                 |
| Instrument ID:  | <u>BNA_G</u>      | Calibration Date/Time: | <u>11/13/2025 14:35</u>      |
| Lab File ID:    | <u>BG064756.D</u> | Init. Calib. Date(s):  | <u>11/12/2025 11/12/2025</u> |
| EPA Sample No.: | <u>SSTDCCC040</u> | Init. Calib. Time(s):  | <u>11:17 16:04</u>           |
| GC Column:      | <u>ZB-GR</u>      | ID:                    | <u>0.25</u> (mm)             |

| COMPOUND                    | RRF   | RRF040 | MIN<br>RRF | %D   | MAX%D |
|-----------------------------|-------|--------|------------|------|-------|
| 2-Fluorophenol              | 1.145 | 1.131  |            | -1.2 |       |
| Benzaldehyde                | 0.997 | 0.963  |            | -3.4 |       |
| Phenol-d6                   | 1.523 | 1.539  |            | 1.1  |       |
| Phenol                      | 1.657 | 1.683  |            | 1.6  | 20.0  |
| bis(2-Chloroethyl)ether     | 1.179 | 1.218  |            | 3.3  |       |
| 2-Chlorophenol              | 1.258 | 1.295  |            | 2.9  |       |
| 2-Methylphenol              | 1.173 | 1.196  |            | 2.0  |       |
| 2,2-oxybis(1-Chloropropane) | 2.792 | 2.835  |            | 1.5  |       |
| Acetophenone                | 0.545 | 0.557  |            | 2.2  |       |
| 3+4-Methylphenols           | 1.607 | 1.614  |            | 0.4  |       |
| n-Nitroso-di-n-propylamine  | 1.030 | 1.080  | 0.050      | 4.9  |       |
| Nitrobenzene-d5             | 0.405 | 0.405  |            | 0.0  |       |
| Hexachloroethane            | 0.558 | 0.545  |            | -2.3 |       |
| Nitrobenzene                | 0.409 | 0.423  |            | 3.4  |       |
| Isophorone                  | 0.770 | 0.773  |            | 0.4  |       |
| 2-Nitrophenol               | 0.184 | 0.189  |            | 2.7  | 20.0  |
| 2,4-Dimethylphenol          | 0.327 | 0.339  |            | 3.7  |       |
| bis(2-Chloroethoxy)methane  | 0.423 | 0.434  |            | 2.6  |       |
| 2,4-Dichlorophenol          | 0.351 | 0.364  |            | 3.7  | 20.0  |
| Naphthalene                 | 1.134 | 1.123  |            | -1.0 |       |
| 4-Chloroaniline             | 0.454 | 0.461  |            | 1.5  |       |
| Hexachlorobutadiene         | 0.363 | 0.358  |            | -1.4 | 20.0  |
| Caprolactam                 | 0.105 | 0.103  |            | -1.9 |       |
| 4-Chloro-3-methylphenol     | 0.390 | 0.391  |            | 0.3  | 20.0  |
| 2-Methylnaphthalene         | 0.848 | 0.841  |            | -0.8 |       |
| Hexachlorocyclopentadiene   | 0.558 | 0.566  | 0.050      | 1.4  |       |
| 2,4,6-Trichlorophenol       | 0.457 | 0.468  |            | 2.4  | 20.0  |
| 2-Fluorobiphenyl            | 1.433 | 1.435  |            | 0.1  |       |
| 2,4,5-Trichlorophenol       | 0.512 | 0.526  |            | 2.7  |       |
| 1,1-Biphenyl                | 1.436 | 1.440  |            | 0.3  |       |
| 2-Chloronaphthalene         | 1.132 | 1.110  |            | -1.9 |       |
| 2-Nitroaniline              | 0.364 | 0.366  |            | 0.5  |       |
| Dimethylphthalate           | 1.580 | 1.576  |            | -0.3 |       |
| Acenaphthylene              | 1.933 | 1.930  |            | -0.2 |       |
| 2,6-Dinitrotoluene          | 0.304 | 0.316  |            | 3.9  |       |
| 3-Nitroaniline              | 0.300 | 0.315  |            | 5.0  |       |
| Acenaphthene                | 1.159 | 1.191  |            | 2.8  | 20.0  |
| 2,4-Dinitrophenol           | 0.158 | 0.168  | 0.050      | 6.3  |       |
| 4-Nitrophenol               | 0.238 | 0.242  | 0.050      | 1.7  |       |

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| Lab Code:       | <u>ACE</u>        | SDG No.:               | <u>Q3615</u>                 |
| Instrument ID:  | <u>BNA_G</u>      | Calibration Date/Time: | <u>11/13/2025 14:35</u>      |
| Lab File ID:    | <u>BG064756.D</u> | Init. Calib. Date(s):  | <u>11/12/2025 11/12/2025</u> |
| EPA Sample No.: | <u>SSTDCCC040</u> | Init. Calib. Time(s):  | <u>11:17 16:04</u>           |
| GC Column:      | <u>ZB-GR</u>      | ID:                    | <u>0.25</u> (mm)             |

| COMPOUND                   | RRF   | RRF040 | MIN<br>RRF | %D   | MAX%D |
|----------------------------|-------|--------|------------|------|-------|
| Dibenzofuran               | 1.916 | 1.896  |            | -1.0 |       |
| 2,4-Dinitrotoluene         | 0.456 | 0.482  |            | 5.7  |       |
| Diethylphthalate           | 1.513 | 1.532  |            | 1.3  |       |
| 4-Chlorophenyl-phenylether | 0.902 | 0.902  |            | 0.0  |       |
| Fluorene                   | 1.503 | 1.475  |            | -1.9 |       |
| 4-Nitroaniline             | 0.311 | 0.324  |            | 4.2  |       |
| 4,6-Dinitro-2-methylphenol | 0.111 | 0.123  |            | 10.8 |       |
| n-Nitrosodiphenylamine     | 0.524 | 0.550  |            | 5.0  | 20.0  |
| 2,4,6-Tribromophenol       | 0.390 | 0.387  |            | -0.8 |       |
| 4-Bromophenyl-phenylether  | 0.263 | 0.273  |            | 3.8  |       |
| Hexachlorobenzene          | 0.318 | 0.328  |            | 3.1  |       |
| Atrazine                   | 0.253 | 0.255  |            | 0.8  |       |
| Pentachlorophenol          | 0.157 | 0.168  |            | 7.0  | 20.0  |
| Phenanthrene               | 1.094 | 1.107  |            | 1.2  |       |
| Anthracene                 | 1.116 | 1.128  |            | 1.1  |       |
| Carbazole                  | 0.959 | 0.957  |            | -0.2 |       |
| Di-n-butylphthalate        | 1.070 | 1.087  |            | 1.6  |       |
| Fluoranthene               | 1.496 | 1.486  |            | -0.7 | 20.0  |
| Pyrene                     | 1.121 | 1.181  |            | 5.4  |       |
| Terphenyl-d14              | 1.017 | 1.067  |            | 4.9  |       |
| Butylbenzylphthalate       | 0.394 | 0.404  |            | 2.5  |       |
| 3,3-Dichlorobenzidine      | 0.501 | 0.488  |            | -2.6 |       |
| Benzo (a) anthracene       | 1.360 | 1.363  |            | 0.2  |       |
| Chrysene                   | 1.282 | 1.292  |            | 0.8  |       |
| Bis(2-ethylhexyl)phthalate | 0.576 | 0.582  |            | 1.0  |       |
| Di-n-octyl phthalate       | 0.996 | 0.995  |            | -0.1 | 20.0  |
| Benzo (b) fluoranthene     | 1.273 | 1.277  |            | 0.3  |       |
| Benzo (k) fluoranthene     | 1.279 | 1.262  |            | -1.3 |       |
| Benzo (a) pyrene           | 1.181 | 1.176  |            | -0.4 | 20.0  |
| Indeno (1,2,3-cd) pyrene   | 1.374 | 1.421  |            | 3.4  |       |
| Dibenzo (a,h) anthracene   | 1.126 | 1.168  |            | 3.7  |       |
| Benzo (g,h,i) perylene     | 1.083 | 1.119  |            | 3.3  |       |
| 1,2,4,5-Tetrachlorobenzene | 0.778 | 0.773  |            | -0.6 |       |
| 1,4-Dioxane                | 0.443 | 0.450  |            | 1.6  | 20.0  |
| 2,3,4,6-Tetrachlorophenol  | 0.506 | 0.525  |            | 3.8  |       |

All other compounds must meet a minimum RRF of 0.010.