

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

|                 |                   |                        |                              |
|-----------------|-------------------|------------------------|------------------------------|
| Lab Name:       | <u>Alliance</u>   | Contract:              | <u>CHEM02</u>                |
| Lab Code:       | <u>ACE</u>        | SDG No.:               | <u>Q2987</u>                 |
| Instrument ID:  | <u>BNA_P</u>      | Calibration Date/Time: | <u>09/09/2025 10:31</u>      |
| Lab File ID:    | <u>BP025692.D</u> | Init. Calib. Date(s):  | <u>09/08/2025 09/08/2025</u> |
| EPA Sample No.: | <u>SSTD020678</u> | Init. Calib. Time(s):  | <u>14:41 18:07</u>           |
| GC Column:      | <u>ZB-GR</u>      | ID:                    | <u>0.25</u> (mm)             |

| COMPOUND                    | RRF   | RRFCAL | MIN RRF | %D    | MAX%D |
|-----------------------------|-------|--------|---------|-------|-------|
| 1,4-Dioxane                 | 0.517 | 0.560  | 0.010   | 8.5   | 40.0  |
| Benzaldehyde                | 0.744 | 0.975  | 0.010   | 31.1  | 40.0  |
| Phenol                      | 1.659 | 1.636  | 0.080   | -1.4  | 20.0  |
| Bis(2-Chloroethyl)ether     | 1.309 | 1.348  | 0.100   | 3.0   | 20.0  |
| 2-Chlorophenol              | 1.285 | 1.348  | 0.200   | 4.9   | 20.0  |
| 2-Methylphenol              | 1.272 | 1.266  | 0.010   | -0.5  | 20.0  |
| 2,2-oxybis(1-Chloropropane) | 2.076 | 2.140  | 0.010   | 3.1   | 40.0  |
| Acetophenone                | 2.069 | 2.156  | 0.060   | 4.2   | 20.0  |
| 4-Methylphenol              | 1.379 | 1.382  | 0.010   | 0.2   | 20.0  |
| N-Nitroso-di-n-propylamine  | 1.099 | 1.138  | 0.050   | 3.5   | 30.0  |
| Hexachloroethane            | 0.576 | 0.596  | 0.100   | 3.5   | 20.0  |
| Nitrobenzene                | 0.372 | 0.384  | 0.050   | 3.1   | 25.0  |
| Isophorone                  | 0.762 | 0.776  | 0.050   | 1.8   | 25.0  |
| 2-Nitrophenol               | 0.182 | 0.183  | 0.050   | 0.6   | 25.0  |
| 2,4-Dimethylphenol          | 0.373 | 0.368  | 0.050   | -1.4  | 25.0  |
| Bis(2-Chloroethoxy)methane  | 0.438 | 0.452  | 0.050   | 3.4   | 20.0  |
| 2,4-Dichlorophenol          | 0.306 | 0.315  | 0.060   | 2.8   | 20.0  |
| Naphthalene                 | 1.046 | 1.075  | 0.200   | 2.8   | 20.0  |
| 4-Chloroaniline             | 0.440 | 0.441  | 0.010   | 0.2   | 40.0  |
| Hexachlorobutadiene         | 0.228 | 0.233  | 0.040   | 2.0   | 30.0  |
| Caprolactam                 | 0.126 | 0.126  | 0.010   | 0.1   | 40.0  |
| 4-Chloro-3-methylphenol     | 0.360 | 0.366  | 0.040   | 1.4   | 25.0  |
| 2-Methylnaphthalene         | 0.747 | 0.754  | 0.100   | 0.9   | 20.0  |
| Hexachlorocyclopentadiene   | 0.302 | 0.271  | 0.010   | -10.5 | 40.0  |
| 2,4,6-Trichlorophenol       | 0.389 | 0.395  | 0.090   | 1.6   | 25.0  |
| 2,4,5-Trichlorophenol       | 0.426 | 0.424  | 0.100   | -0.5  | 25.0  |
| 1,1-Biphenyl                | 1.402 | 1.440  | 0.200   | 2.7   | 20.0  |
| 2-Chloronaphthalene         | 1.104 | 1.136  | 0.300   | 2.9   | 20.0  |
| 2-Nitroaniline              | 0.350 | 0.356  | 0.050   | 1.7   | 40.0  |
| Dimethylphthalate           | 1.483 | 1.505  | 0.300   | 1.5   | 20.0  |
| 2,6-Dinitrotoluene          | 0.300 | 0.302  | 0.080   | 0.3   | 30.0  |
| Acenaphthylene              | 1.859 | 1.889  | 0.400   | 1.7   | 20.0  |
| 3-Nitroaniline              | 0.267 | 0.296  | 0.010   | 10.9  | 40.0  |
| Acenaphthene                | 1.194 | 1.211  | 0.200   | 1.4   | 20.0  |
| 2,4-Dinitrophenol           | 0.185 | 0.155  | 0.010   | -16.1 | 40.0  |
| 4-Nitrophenol               | 0.204 | 0.192  | 0.010   | -5.8  | 40.0  |
| Dibenzofuran                | 1.708 | 1.749  | 0.300   | 2.4   | 20.0  |
| 2,4-Dinitrotoluene          | 0.450 | 0.464  | 0.070   | 3.0   | 30.0  |
| Diethylphthalate            | 1.525 | 1.531  | 0.300   | 0.4   | 20.0  |

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| Lab Code:       | <u>ACE</u>        | SDG No.:               | <u>Q2987</u>                 |
| Instrument ID:  | <u>BNA_P</u>      | Calibration Date/Time: | <u>09/09/2025 10:31</u>      |
| Lab File ID:    | <u>BP025692.D</u> | Init. Calib. Date(s):  | <u>09/08/2025 09/08/2025</u> |
| EPA Sample No.: | <u>SSTD020678</u> | Init. Calib. Time(s):  | <u>14:41 18:07</u>           |
| GC Column:      | <u>ZB-GR</u>      | ID:                    | <u>0.25</u> (mm)             |

| COMPOUND                    | RRF   | RRFCAL | MIN RRF | %D    | MAX%D |
|-----------------------------|-------|--------|---------|-------|-------|
| Fluorene                    | 1.395 | 1.429  | 0.200   | 2.5   | 20.0  |
| 4-Chlorophenyl-phenylether  | 0.714 | 0.725  | 0.100   | 1.5   | 20.0  |
| 4-Nitroaniline              | 0.263 | 0.293  | 0.010   | 11.2  | 40.0  |
| 4,6-Dinitro-2-methylphenol  | 0.138 | 0.124  | 0.010   | -10.2 | 40.0  |
| N-Nitrosodiphenylamine      | 0.578 | 0.558  | 0.050   | -3.5  | 20.0  |
| 1,2,4,5-Tetrachlorobenzene  | 0.595 | 0.606  | 0.100   | 1.9   | 20.0  |
| 4-Bromophenyl-phenylether   | 0.213 | 0.216  | 0.070   | 1.1   | 20.0  |
| Hexachlorobenzene           | 0.244 | 0.246  | 0.050   | 0.9   | 25.0  |
| Atrazine                    | 0.237 | 0.250  | 0.010   | 5.4   | 25.0  |
| Pentachlorophenol           | 0.156 | 0.150  | 0.010   | -4.0  | 40.0  |
| Phenanthrene                | 1.097 | 1.116  | 0.200   | 1.7   | 20.0  |
| Anthracene                  | 1.113 | 1.133  | 0.200   | 1.9   | 20.0  |
| Carbazole                   | 1.000 | 1.046  | 0.050   | 4.6   | 40.0  |
| Di-n-butylphthalate         | 1.285 | 1.314  | 0.500   | 2.2   | 25.0  |
| Fluoranthene                | 1.246 | 1.262  | 0.400   | 1.3   | 25.0  |
| Pyrene                      | 1.303 | 1.328  | 0.400   | 2.0   | 25.0  |
| Butylbenzylphthalate        | 0.580 | 0.597  | 0.100   | 2.8   | 40.0  |
| 3,3-Dichlorobenzidine       | 0.429 | 0.446  | 0.010   | 4.1   | 40.0  |
| Benzo(a)anthracene          | 1.332 | 1.350  | 0.300   | 1.4   | 30.0  |
| Chrysene                    | 1.254 | 1.287  | 0.200   | 2.7   | 30.0  |
| Bis(2-ethylhexyl)phthalate  | 0.853 | 0.893  | 0.200   | 4.7   | 40.0  |
| Di-n-octyl phthalate        | 1.350 | 1.420  | 0.010   | 5.2   | 40.0  |
| Benzo(b)fluoranthene        | 1.171 | 1.204  | 0.200   | 2.8   | 25.0  |
| Benzo(k)fluoranthene        | 1.188 | 1.237  | 0.200   | 4.2   | 25.0  |
| Benzo(a)pyrene              | 1.125 | 1.151  | 0.200   | 2.3   | 20.0  |
| Indeno(1,2,3-cd)pyrene      | 1.407 | 1.423  | 0.200   | 1.1   | 25.0  |
| Dibenzo(a,h)anthracene      | 1.140 | 1.151  | 0.200   | 1.0   | 30.0  |
| Benzo(g,h,i)perylene        | 1.141 | 1.144  | 0.200   | 0.2   | 30.0  |
| 2,3,4,6-Tetrachlorophenol   | 0.374 | 0.380  | 0.040   | 1.7   | 20.0  |
| 1,4-Dioxane-d8              | 0.489 | 0.538  | 0.010   | 10.1  | 25.0  |
| Pyridine-d5                 | 1.417 | 1.508  | 0.010   | 6.4   | 40.0  |
| Phenol-d5                   | 1.616 | 1.597  | 0.010   | -1.2  | 25.0  |
| Bis-(2-Chloroethyl)ether-d8 | 0.988 | 1.009  | 0.050   | 2.1   | 25.0  |
| 2-Chlorophenol-d4           | 1.258 | 1.304  | 0.200   | 3.7   | 20.0  |
| 4-Methylphenol-d8           | 1.338 | 1.334  | 0.010   | -0.2  | 20.0  |
| Nitrobenzene-d5             | 0.154 | 0.158  | 0.050   | 2.5   | 20.0  |
| 2-Nitrophenol-d4            | 0.175 | 0.177  | 0.050   | 1.4   | 30.0  |
| 2,4-Dichlorophenol-d3       | 0.315 | 0.323  | 0.060   | 2.4   | 20.0  |
| 4-Chloroaniline-d4          | 0.442 | 0.443  | 0.010   | 0.4   | 40.0  |

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 Lab File ID: BP025692.D Init. Calib. Date(s): 09/08/2025 09/08/2025  
 EPA Sample No.: SSTD020678 Init. Calib. Time(s): 14:41 18:07  
 GC Column: ZB-GR ID: 0.25 (mm)

| COMPOUND                      | RRF   | RRFCAL | MIN<br>RRF | %D    | MAX%D |
|-------------------------------|-------|--------|------------|-------|-------|
| Dimethylphthalate-d6          | 1.512 | 1.537  | 0.300      | 1.6   | 20.0  |
| Acenaphthylene-d8             | 1.660 | 1.697  | 0.400      | 2.2   | 20.0  |
| 4-Nitrophenol-d4              | 0.251 | 0.243  | 0.010      | -3.3  | 40.0  |
| Fluorene-d10                  | 1.259 | 1.284  | 0.100      | 2.0   | 20.0  |
| 4,6-Dinitro-2-methylphenol-d2 | 0.136 | 0.120  | 0.010      | -11.5 | 40.0  |
| Anthracene-d10                | 0.963 | 0.982  | 0.300      | 2.1   | 20.0  |
| Pyrene-d10                    | 1.095 | 1.108  | 0.300      | 1.2   | 25.0  |
| Benzo(a)pyrene-d12            | 1.027 | 1.056  | 0.010      | 2.8   | 20.0  |

All other compounds must meet a minimum RRF of 0.010.