

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name:	<u>Alliance</u>	Contract:	<u>TETR06</u>
Lab Code:	<u>ACE</u>	SDG No.:	<u>Q3776</u>
Instrument ID:	<u>BNA_P</u>	Calibration Date/Time:	<u>12/08/2025 17:49</u>
Lab File ID:	<u>BP026256.D</u>	Init. Calib. Date(s):	<u>11/18/2025 11/18/2025</u>
EPA Sample No.:	<u>SSTDCCC040EC</u>	Init. Calib. Time(s):	<u>14:31 21:22</u>
GC Column:	<u>ZB-GR</u>	ID:	<u>0.25</u> (mm)

COMPOUND	RRF	RRF040	MIN RRF	%D	MAX%D
2-Fluorophenol	1.246	1.259		1.0	50.0
Benzaldehyde	0.831	1.058		27.3	50.0
Phenol-d6	1.586	1.518		-4.3	50.0
Phenol	1.709	1.615		-5.5	50.0
bis(2-Chloroethyl)ether	1.338	1.268		-5.2	50.0
2-Chlorophenol	1.418	1.414		-0.3	50.0
2-Methylphenol	1.163	1.109		-4.6	50.0
2,2-oxybis(1-Chloropropane)	1.879	1.787		-4.9	50.0
Acetophenone	0.508	0.522		2.8	50.0
3+4-Methylphenols	1.612	1.501		-6.9	50.0
n-Nitroso-di-n-propylamine	0.985	0.929	0.050	-5.7	50.0
Nitrobenzene-d5	0.352	0.371		5.4	50.0
Hexachloroethane	0.560	0.599		7.0	50.0
Nitrobenzene	0.356	0.371		4.2	50.0
Isophorone	0.698	0.719		3.0	50.0
2-Nitrophenol	0.180	0.196		8.9	50.0
2,4-Dimethylphenol	0.325	0.314		-3.4	50.0
bis(2-Chloroethoxy)methane	0.437	0.439		0.5	50.0
2,4-Dichlorophenol	0.325	0.344		5.8	50.0
Naphthalene	1.081	1.124		4.0	50.0
4-Chloroaniline	0.460	0.459		-0.2	50.0
Hexachlorobutadiene	0.215	0.243		13.0	50.0
Caprolactam	0.116	0.116		0.0	50.0
4-Chloro-3-methylphenol	0.345	0.356		3.2	50.0
2-Methylnaphthalene	0.766	0.792		3.4	50.0
Hexachlorocyclopentadiene	0.396	0.411	0.050	3.8	50.0
2,4,6-Trichlorophenol	0.415	0.445		7.2	50.0
2-Fluorobiphenyl	1.365	1.480		8.4	50.0
2,4,5-Trichlorophenol	0.463	0.494		6.7	50.0
1,1-Biphenyl	1.506	1.556		3.3	50.0
2-Chloronaphthalene	1.184	1.236		4.4	50.0
2-Nitroaniline	0.329	0.342		4.0	50.0
Dimethylphthalate	1.494	1.573		5.3	50.0
Acenaphthylene	1.953	2.045		4.7	50.0
2,6-Dinitrotoluene	0.296	0.324		9.5	50.0
3-Nitroaniline	0.349	0.345		-1.1	50.0
Acenaphthene	1.145	1.168		2.0	50.0
2,4-Dinitrophenol	0.179	0.207	0.050	15.6	50.0
4-Nitrophenol	0.315	0.289	0.050	-8.3	50.0

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COMPOUND	RRF	RRF040	MIN RRF	%D	MAX%D
Dibenzofuran	1.773	1.867		5.3	50.0
2,4-Dinitrotoluene	0.442	0.469		6.1	50.0
Diethylphthalate	1.505	1.570		4.3	50.0
4-Chlorophenyl-phenylether	0.697	0.748		7.3	50.0
Fluorene	1.401	1.494		6.6	50.0
4-Nitroaniline	0.362	0.339		-6.4	50.0
4,6-Dinitro-2-methylphenol	0.127	0.142		11.8	50.0
n-Nitrosodiphenylamine	0.619	0.654		5.7	50.0
2,4,6-Tribromophenol	0.259	0.287		10.8	50.0
4-Bromophenyl-phenylether	0.221	0.244		10.4	50.0
Hexachlorobenzene	0.257	0.287		11.7	50.0
Atrazine	0.233	0.248		6.4	50.0
Pentachlorophenol	0.180	0.193		7.2	50.0
Phenanthrene	1.127	1.175		4.3	50.0
Anthracene	1.149	1.208		5.1	50.0
Carbazole	1.091	1.101		0.9	50.0
Di-n-butylphthalate	1.294	1.410		9.0	50.0
Fluoranthene	1.369	1.450		5.9	50.0
Pyrene	1.311	1.381		5.3	50.0
Terphenyl-d14	0.981	1.086		10.7	50.0
Butylbenzylphthalate	0.578	0.616		6.6	50.0
3,3-Dichlorobenzidine	0.500	0.514		2.8	50.0
Benzo (a) anthracene	1.374	1.404		2.2	50.0
Chrysene	1.295	1.363		5.3	50.0
Bis(2-ethylhexyl)phthalate	0.874	0.926		5.9	50.0
Di-n-octyl phthalate	1.529	1.660		8.6	50.0
Benzo (b) fluoranthene	1.218	1.291		6.0	50.0
Benzo (k) fluoranthene	1.217	1.268		4.2	50.0
Benzo (a) pyrene	1.154	1.209		4.8	50.0
Indeno (1,2,3-cd) pyrene	1.500	1.593		6.2	50.0
Dibenzo (a,h) anthracene	1.228	1.296		5.5	50.0
Benzo (g,h,i) perylene	1.220	1.308		7.2	50.0
1,2,4,5-Tetrachlorobenzene	0.605	0.653		7.9	50.0
1,4-Dioxane	0.505	0.547		8.3	50.0
2,3,4,6-Tetrachlorophenol	0.393	0.417		6.1	50.0

All other compounds must meet a minimum RRF of 0.010.