

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name:	<u>Alliance</u>	Contract:	<u>YANN01</u>
Lab Code:	<u>ACE</u>	SDG No.:	<u>Q3238</u>
Instrument ID:	<u>BNA_F</u>	Calibration Date/Time:	<u>09/30/2025 13:36</u>
Lab File ID:	<u>BF143830.D</u>	Init. Calib. Date(s):	<u>09/23/2025 09/23/2025</u>
EPA Sample No.:	<u>SSTDCCC040</u>	Init. Calib. Time(s):	<u>10:28 15:27</u>
GC Column:	<u>DB-UI</u>	ID:	<u>0.18</u> (mm)

COMPOUND	RRF	RRF040	MIN RRF	%D	MAX%D
2-Fluorophenol	1.231	1.238		0.6	
Benzaldehyde	1.246	1.159		-7.0	
Phenol-d6	1.422	1.418		-0.3	
Phenol	1.653	1.739		5.2	20.0
bis(2-Chloroethyl)ether	1.296	1.278		-1.4	
2-Chlorophenol	1.260	1.285		2.0	
2-Methylphenol	0.976	0.992		1.6	
2,2-oxybis(1-Chloropropane)	3.073	2.569		-16.4	
Acetophenone	0.439	0.430		-2.0	
3+4-Methylphenols	1.185	1.233		4.1	
n-Nitroso-di-n-propylamine	0.915	0.880	0.050	-3.8	
Nitrobenzene-d5	0.327	0.340		4.0	
Hexachloroethane	0.514	0.502		-2.3	
Nitrobenzene	0.362	0.371		2.5	
Isophorone	0.637	0.640		0.5	
2-Nitrophenol	0.132	0.165		25.0	20.0
2,4-Dimethylphenol	0.285	0.289		1.4	
bis(2-Chloroethoxy)methane	0.414	0.409		-1.2	
2,4-Dichlorophenol	0.280	0.290		3.6	20.0
Naphthalene	0.957	0.950		-0.7	
4-Chloroaniline	0.336	0.344		2.4	
Hexachlorobutadiene	0.217	0.231		6.5	20.0
Caprolactam	0.072	0.076		5.6	
4-Chloro-3-methylphenol	0.263	0.273		3.8	20.0
2-Methylnaphthalene	0.611	0.620		1.5	
Hexachlorocyclopentadiene	0.383	0.363	0.050	-5.2	
2,4,6-Trichlorophenol	0.378	0.398		5.3	20.0
2-Fluorobiphenyl	1.282	1.279		-0.2	
2,4,5-Trichlorophenol	0.395	0.418		5.8	
1,1-Biphenyl	1.448	1.406		-2.9	
2-Chloronaphthalene	1.165	1.134		-2.7	
2-Nitroaniline	0.352	0.398		13.1	
Dimethylphthalate	1.219	1.232		1.1	
Acenaphthylene	1.604	1.578		-1.6	
2,6-Dinitrotoluene	0.200	0.235		17.5	
3-Nitroaniline	0.243	0.275		13.2	
Acenaphthene	1.056	1.037		-1.8	20.0
2,4-Dinitrophenol	0.064	0.076	0.050	18.8	
4-Nitrophenol	0.196	0.189	0.050	-3.6	

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COMPOUND	RRF	RRF040	MIN RRF	%D	MAX%D
Dibenzofuran	1.493	1.484		-0.6	
2,4-Dinitrotoluene	0.261	0.311		19.2	
Diethylphthalate	1.141	1.133		-0.7	
4-Chlorophenyl-phenylether	0.566	0.582		2.8	
Fluorene	1.110	1.116		0.5	
4-Nitroaniline	0.228	0.247		8.3	
4,6-Dinitro-2-methylphenol	0.067	0.082		22.4	
n-Nitrosodiphenylamine	0.658	0.657		-0.2	20.0
2,4,6-Tribromophenol	0.296	0.339		14.5	
4-Bromophenyl-phenylether	0.341	0.380		11.4	
Hexachlorobenzene	0.470	0.541		15.1	
Atrazine	0.190	0.192		1.1	
Pentachlorophenol	0.234	0.266		13.7	20.0
Phenanthrene	1.027	1.014		-1.3	
Anthracene	1.030	1.027		-0.3	
Carbazole	0.901	0.875		-2.9	
Di-n-butylphthalate	1.036	1.018		-1.7	
Fluoranthene	1.033	0.998		-3.4	20.0
Pyrene	0.983	1.061		7.9	
Terphenyl-d14	0.983	1.123		14.2	
Butylbenzylphthalate	0.340	0.359		5.6	
3,3-Dichlorobenzidine	0.465	0.471		1.3	
Benzo (a) anthracene	1.074	1.098		2.2	
Chrysene	0.972	0.954		-1.9	
Bis (2-ethylhexyl) phthalate	0.499	0.481		-3.6	
Di-n-octyl phthalate	0.905	0.816		-9.8	20.0
Benzo (b) fluoranthene	1.019	1.000		-1.8	
Benzo (k) fluoranthene	1.023	1.018		-0.5	
Benzo (a) pyrene	0.957	0.960		0.3	20.0
Indeno (1,2,3-cd) pyrene	1.517	1.699		12.0	
Dibenzo (a,h) anthracene	1.241	1.389		11.9	
Benzo (g,h,i) perylene	1.196	1.349		12.8	
1,2,4,5-Tetrachlorobenzene	0.604	0.612		1.3	
1,4-Dioxane	0.589	0.581		-1.4	20.0
2,3,4,6-Tetrachlorophenol	0.339	0.391		15.3	

All other compounds must meet a minimum RRF of 0.010.