

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name:	<u>Alliance</u>	Contract:	<u>REMI02</u>
Lab Code:	<u>ACE</u>	SDG No.:	<u>Q3495</u>
Instrument ID:	<u>BNA_P</u>	Calibration Date/Time:	<u>11/03/2025 11:06</u>
Lab File ID:	<u>BP026053.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	1.217		0.4	
Benzaldehyde	0.802	0.703		-12.3	
Phenol-d6	1.543	1.579		2.3	
Phenol	1.713	1.731		1.1	20.0
bis(2-Chloroethyl)ether	1.352	1.405		3.9	
2-Chlorophenol	1.342	1.364		1.6	
2-Methylphenol	1.121	1.153		2.9	
2,2-oxybis(1-Chloropropane)	2.043	2.143		4.9	
Acetophenone	0.506	0.516		2.0	
3+4-Methylphenols	1.560	1.560		0.0	
n-Nitroso-di-n-propylamine	0.958	1.020	0.050	6.5	
Nitrobenzene-d5	0.321	0.336		4.7	
Hexachloroethane	0.533	0.534		0.2	
Nitrobenzene	0.339	0.351		3.5	
Isophorone	0.687	0.709		3.2	
2-Nitrophenol	0.104	0.124		19.2	20.0
2,4-Dimethylphenol	0.289	0.287		-0.7	
bis(2-Chloroethoxy)methane	0.433	0.442		2.1	
2,4-Dichlorophenol	0.305	0.308		1.0	20.0
Naphthalene	1.091	1.101		0.9	
4-Chloroaniline	0.419	0.430		2.6	
Hexachlorobutadiene	0.211	0.207		-1.9	20.0
Caprolactam	0.107	0.108		0.9	
4-Chloro-3-methylphenol	0.320	0.333		4.1	20.0
2-Methylnaphthalene	0.763	0.765		0.3	
Hexachlorocyclopentadiene	0.227	0.232	0.050	2.2	
2,4,6-Trichlorophenol	0.367	0.381		3.8	20.0
2-Fluorobiphenyl	1.392	1.389		-0.2	
2,4,5-Trichlorophenol	0.419	0.436		4.1	
1,1-Biphenyl	1.531	1.559		1.8	
2-Chloronaphthalene	1.205	1.213		0.7	
2-Nitroaniline	0.272	0.298		9.6	
Dimethylphthalate	1.455	1.469		1.0	
Acenaphthylene	1.756	1.802		2.6	
2,6-Dinitrotoluene	0.237	0.261		10.1	
3-Nitroaniline	0.291	0.328		12.7	
Acenaphthene	1.206	1.235		2.4	20.0
2,4-Dinitrophenol	0.100	0.115	0.050	15.0	
4-Nitrophenol	0.290	0.277	0.050	-4.5	

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COMPOUND	RRF	RRF040	MIN RRF	%D	MAX%D
Dibenzofuran	1.822	1.828		0.3	
2,4-Dinitrotoluene	0.347	0.392		13.0	
Diethylphthalate	1.463	1.489		1.8	
4-Chlorophenyl-phenylether	0.713	0.723		1.4	
Fluorene	1.427	1.458		2.2	
4-Nitroaniline	0.311	0.349		12.2	
4,6-Dinitro-2-methylphenol	0.077	0.083		7.8	
n-Nitrosodiphenylamine	0.625	0.635		1.6	20.0
2,4,6-Tribromophenol	0.216	0.215		-0.5	
4-Bromophenyl-phenylether	0.222	0.225		1.4	
Hexachlorobenzene	0.253	0.250		-1.2	
Atrazine	0.211	0.209		-0.9	
Pentachlorophenol	0.154	0.151		-1.9	20.0
Phenanthrene	1.164	1.162		-0.2	
Anthracene	1.154	1.174		1.7	
Carbazole	1.092	1.126		3.1	
Di-n-butylphthalate	1.222	1.248		2.1	
Fluoranthene	1.358	1.380		1.6	20.0
Pyrene	1.353	1.388		2.6	
Terphenyl-d14	0.984	1.002		1.8	
Butylbenzylphthalate	0.432	0.480		11.1	
3,3-Dichlorobenzidine	0.443	0.455		2.7	
Benzo (a) anthracene	1.374	1.402		2.0	
Chrysene	1.302	1.324		1.7	
Bis(2-ethylhexyl)phthalate	0.718	0.765		6.5	
Di-n-octyl phthalate	1.191	1.230		3.3	20.0
Benzo (b) fluoranthene	1.242	1.277		2.8	
Benzo (k) fluoranthene	1.268	1.281		1.0	
Benzo (a) pyrene	1.141	1.181		3.5	20.0
Indeno (1,2,3-cd) pyrene	1.476	1.531		3.7	
Dibenzo (a,h) anthracene	1.202	1.240		3.2	
Benzo (g,h,i) perylene	1.156	1.186		2.6	
1,2,4,5-Tetrachlorobenzene	0.621	0.614		-1.1	
1,4-Dioxane	0.511	0.528		3.3	20.0
2,3,4,6-Tetrachlorophenol	0.330	0.342		3.6	

All other compounds must meet a minimum RRF of 0.010.