

6C

SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: Alliance

Contract: ZUCC01

Lab Code: ACE

SDG No.: Q3032

Instrument ID: BNA_G

Calibration Date(s): 08/28/2025 08/28/2025

Calibration Time(s): 10:51 16:16

LAB FILE ID:		RRF2.5 = BG064218.D			RRF005 = BG064219.D		RRF010 = BG064220.D	
		RRF020 = BG064221.D			RRF040 = BG064222.D		RRF050 = BG064223.D	
COMPOUND	RRF2.5	RRF005	RRF010	RRF020	RRF040	RRF050	RRF	% RSD
2-Fluorophenol		1.126	1.094	1.155	1.104	1.075	1.115	2.4
Benzaldehyde			1.181	1.115	1.375	1.193	1.260	11.0
Phenol-d6		1.520	1.484	1.594	1.515	1.522	1.532	2.3
Phenol		1.590	1.627	1.750	1.669	1.739	1.689	3.7
bis(2-Chloroethyl)ether		1.232	1.263	1.323	1.281	1.255	1.269	2.2
2-Chlorophenol		1.318	1.287	1.381	1.418	1.360	1.362	3.4
2-Methylphenol		1.188	1.243	1.349	1.261	1.218	1.258	4.3
2,2-oxybis(1-Chloropropane)		2.923	2.920	2.979	2.897	2.826	2.897	2.1
Acetophenone		0.533	0.485	0.522	0.502	0.496	0.506	3.4
3+4-Methylphenols			1.734	1.765	1.702	1.750	1.749	1.6
n-Nitroso-di-n-propylamine	1.008	1.162	1.160	1.239	1.203	1.211	1.172	6.2
Nitrobenzene-d5		0.341	0.340	0.370	0.365	0.357	0.359	3.8
Hexachloroethane		0.555	0.578	0.574	0.583	0.572	0.571	1.6
Nitrobenzene		0.354	0.348	0.394	0.388	0.377	0.375	4.7
Isophorone		0.749	0.723	0.787	0.748	0.733	0.745	3.0
2-Nitrophenol		0.133	0.146	0.169	0.171	0.167	0.162	9.9
2,4-Dimethylphenol		0.275	0.271	0.300	0.297	0.276	0.284	4.0
bis(2-Chloroethoxy)methane		0.396	0.401	0.417	0.414	0.390	0.402	2.6
2,4-Dichlorophenol		0.289	0.278	0.319	0.310	0.299	0.300	4.7
Naphthalene		1.104	1.012	1.077	1.036	1.004	1.037	4.0
4-Chloroaniline		0.369	0.379	0.414	0.421	0.411	0.402	5.1
Hexachlorobutadiene		0.242	0.236	0.236	0.237	0.224	0.234	2.9
Caprolactam			0.108	0.129	0.123	0.112	0.116	7.1
4-Chloro-3-methylphenol		0.380	0.382	0.402	0.395	0.386	0.389	2.2
2-Methylnaphthalene		0.768	0.748	0.831	0.788	0.757	0.771	4.1
Hexachlorocyclopentadiene			0.221	0.256	0.275	0.290	0.274	11.6
2,4,6-Trichlorophenol		0.364	0.362	0.400	0.396	0.384	0.386	4.2
2-Fluorobiphenyl		1.364	1.286	1.358	1.328	1.288	1.310	3.3
2,4,5-Trichlorophenol		0.387	0.386	0.437	0.433	0.426	0.418	5.3
1,1-Biphenyl		1.459	1.412	1.488	1.494	1.449	1.457	2.1
2-Chloronaphthalene		1.090	1.059	1.113	1.113	1.096	1.095	1.9
2-Nitroaniline		0.303	0.329	0.372	0.410	0.394	0.373	11.1
Dimethylphthalate		1.672	1.476	1.569	1.555	1.462	1.514	5.9
Acenaphthylene		1.656	1.528	1.675	1.675	1.607	1.619	3.4

All other compounds must meet a minimum RRF of 0.010.

Form VI SV-1

6C

SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: Alliance

Contract: ZUCC01

Lab Code: ACE

SDG No.: Q3032

Instrument ID: BNA_G

Calibration Date(s): 08/28/2025 08/28/2025

Calibration Time(s): 10:51 16:16

LAB FILE ID:		RRF2.5 = BG064218.D			RRF005 = BG064219.D		RRF010 = BG064220.D	
		RRF020 = BG064221.D			RRF040 = BG064222.D		RRF050 = BG064223.D	
COMPOUND	RRF2.5	RRF005	RRF010	RRF020	RRF040	RRF050	RRF	% RSD
2,6-Dinitrotoluene		0.238	0.269	0.301	0.322	0.300	0.293	10.0
3-Nitroaniline		0.275	0.279	0.316	0.331	0.314	0.307	6.9
Acenaphthene		1.169	1.120	1.175	1.178	1.129	1.142	2.8
2,4-Dinitrophenol			0.096	0.137	0.178	0.184	0.163	23.3
4-Nitrophenol			0.199	0.250	0.286	0.258	0.252	11.5
Dibenzofuran		1.913	1.756	1.840	1.833	1.762	1.797	3.7
2,4-Dinitrotoluene		0.396	0.398	0.462	0.484	0.450	0.443	7.6
Diethylphthalate		1.830	1.577	1.665	1.688	1.536	1.620	7.1
4-Chlorophenyl-phenylether		0.781	0.717	0.759	0.755	0.720	0.734	4.2
Fluorene		1.567	1.478	1.549	1.529	1.416	1.473	5.3
4-Nitroaniline		0.267	0.275	0.311	0.351	0.332	0.313	9.9
4,6-Dinitro-2-methylphenol			0.085	0.107	0.118	0.121	0.115	14.4
n-Nitrosodiphenylamine		0.553	0.518	0.571	0.562	0.542	0.550	3.2
2,4,6-Tribromophenol		0.281	0.256	0.271	0.276	0.261	0.265	4.2
4-Bromophenyl-phenylether		0.195	0.213	0.218	0.215	0.212	0.212	3.8
Hexachlorobenzene		0.241	0.232	0.244	0.226	0.231	0.236	2.8
Atrazine		0.240	0.230	0.231	0.240	0.224	0.230	3.4
Pentachlorophenol			0.110	0.135	0.153	0.148	0.143	12.6
Phenanthrene		1.103	1.032	1.090	1.084	1.020	1.055	3.4
Anthracene		1.124	1.053	1.092	1.111	1.033	1.073	3.5
Carbazole		0.986	0.908	0.960	1.002	0.917	0.945	4.0
Di-n-butylphthalate		1.196	1.089	1.146	1.211	1.096	1.136	4.5
Fluoranthene		1.340	1.211	1.269	1.304	1.178	1.240	5.2
Pyrene		1.292	1.236	1.286	1.303	1.216	1.260	2.9
Terphenyl-d14		1.009	0.974	1.002	0.981	0.918	0.959	4.7
Butylbenzylphthalate		0.449	0.440	0.492	0.494	0.474	0.475	4.7
3,3-Dichlorobenzidine			0.401	0.440	0.429	0.410	0.417	3.7
Benzo (a) anthracene		1.323	1.238	1.312	1.296	1.248	1.280	2.7
Chrysene		1.277	1.197	1.254	1.248	1.204	1.228	2.7
Bis (2-ethylhexyl) phthalate		0.713	0.681	0.719	0.732	0.713	0.714	2.5
Di-n-octyl phthalate			1.160	1.266	1.263	1.228	1.243	3.9
Benzo (b) fluoranthene		1.079	1.081	1.145	1.176	1.104	1.129	3.6
Benzo (k) fluoranthene		1.161	1.127	1.178	1.141	1.131	1.150	1.9
Benzo (a) pyrene		1.019	0.978	1.076	1.084	1.047	1.050	3.8
Indeno (1,2,3-cd) pyrene		1.314	1.302	1.384	1.399	1.348	1.366	3.4

All other compounds must meet a minimum RRF of 0.010.

Form VI SV-1

6C

SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: Alliance

Contract: ZUCC01

Lab Code: ACE

SDG No.: Q3032

Instrument ID: BNA_G

Calibration Date(s): 08/28/2025 08/28/2025

Calibration Time(s): 10:51 16:16

LAB FILE ID:		RRF2.5 = BG064218.D			RRF005 = BG064219.D		RRF010 = BG064220.D	
		RRF020 = BG064221.D			RRF040 = BG064222.D		RRF050 = BG064223.D	
COMPOUND	RRF2.5	RRF005	RRF010	RRF020	RRF040	RRF050	RRF	% RSD
Dibenzo(a,h)anthracene		1.044	1.032	1.114	1.135	1.087	1.097	4.0
Benzo(g,h,i)perylene		1.067	1.029	1.070	1.083	1.042	1.067	2.3
1,2,4,5-Tetrachlorobenzene		0.554	0.556	0.597	0.582	0.582	0.577	2.9
1,4-Dioxane		0.574	0.524	0.479	0.387	0.430	0.465	14.0
2,3,4,6-Tetrachlorophenol		0.392	0.378	0.419	0.419	0.413	0.402	3.9