

6C

SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: Alliance

Contract: GENV01

Lab Code: ACE

SDG No.: Q2664

Instrument ID: BNA_F

Calibration Date(s): 07/17/2025 07/17/2025

Calibration Time(s): 11:04 14:34

LAB FILE ID:		RRF2.5 = BF143140.D			RRF005 = BF143141.D		RRF010 = BF143142.D	
		RRF020 = BF143143.D			RRF040 = BF143144.D		RRF050 = BF143145.D	
COMPOUND	RRF2.5	RRF005	RRF010	RRF020	RRF040	RRF050	RRF	% RSD
2-Fluorophenol		1.385	1.313	1.319	1.231	1.284	1.264	7.0
Benzaldehyde			1.167	1.163	1.443	1.411	1.193	18.1
Phenol-d6		1.733	1.640	1.657	1.554	1.632	1.591	6.7
bis(2-Chloroethyl)ether		1.419	1.343	1.379	1.303	1.361	1.327	5.4
2,2-oxybis(1-Chloropropane)		2.670	2.553	2.545	2.421	2.517	2.461	6.5
Acetophenone		0.527	0.504	0.501	0.455	0.476	0.474	8.6
n-Nitroso-di-n-propylamine	1.101	1.111	1.051	1.028	0.986	1.023	1.021	6.8
Nitrobenzene-d5		0.407	0.394	0.417	0.399	0.417	0.401	3.5
Hexachloroethane		0.522	0.498	0.522	0.495	0.525	0.502	4.7
Nitrobenzene		0.380	0.370	0.388	0.375	0.388	0.374	3.9
Isophorone		0.748	0.707	0.711	0.692	0.734	0.708	3.8
bis(2-Chloroethoxy)methane		0.463	0.437	0.442	0.413	0.430	0.425	6.3
Naphthalene		1.118	1.038	1.032	0.954	0.986	0.985	8.9
4-Chloroaniline		0.437	0.417	0.419	0.390	0.408	0.402	6.3
Hexachlorobutadiene		0.193	0.190	0.191	0.181	0.188	0.184	4.6
Caprolactam			0.078	0.084	0.087	0.091	0.084	5.3
2-Methylnaphthalene		0.667	0.634	0.628	0.580	0.603	0.599	8.2
Hexachlorocyclopentadiene			0.337	0.373	0.373	0.404	0.376	6.1
2-Fluorobiphenyl		1.761	1.644	1.547	1.358	1.414	1.462	13.6
1,1-Biphenyl		1.748	1.667	1.640	1.489	1.553	1.560	8.6
2-Chloronaphthalene		1.268	1.223	1.192	1.103	1.166	1.155	7.1
2-Nitroaniline		0.305	0.333	0.372	0.377	0.396	0.362	8.9
Dimethylphthalate		1.422	1.340	1.340	1.279	1.325	1.304	6.2
Acenaphthylene		2.134	2.052	2.011	1.860	1.958	1.935	7.7
2,6-Dinitrotoluene		0.225	0.250	0.272	0.278	0.289	0.265	8.1
3-Nitroaniline		0.285	0.300	0.315	0.317	0.334	0.310	5.1
Acenaphthene		1.278	1.194	1.184	1.096	1.155	1.144	7.5
Dibenzofuran		1.929	1.807	1.766	1.633	1.698	1.698	9.0
2,4-Dinitrotoluene		0.269	0.296	0.342	0.354	0.372	0.333	11.0
Diethylphthalate		1.377	1.322	1.335	1.269	1.329	1.289	5.8
4-Chlorophenyl-phenylether		0.710	0.670	0.657	0.612	0.633	0.635	7.9
Fluorene		1.480	1.402	1.322	1.212	1.257	1.274	10.8
4-Nitroaniline		0.236	0.255	0.266	0.280	0.289	0.266	6.4
n-Nitrosodiphenylamine		0.761	0.728	0.721	0.674	0.726	0.704	6.0

All other compounds must meet a minimum RRF of 0.010.

Form VI SV-1

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2,4,6-Tribromophenol		0.200	0.202	0.212	0.210	0.218	0.207	3.5
4-Bromophenyl-phenylether		0.252	0.238	0.241	0.228	0.247	0.238	4.3
Hexachlorobenzene		0.269	0.248	0.252	0.239	0.258	0.249	4.8
Atrazine		0.184	0.187	0.198	0.196	0.208	0.194	4.3
Phenanthrene		1.191	1.100	1.117	1.020	1.069	1.062	8.0
Anthracene		1.184	1.136	1.134	1.038	1.095	1.083	7.1
Carbazole		1.042	0.969	0.982	0.929	0.963	0.946	6.9
Di-n-butylphthalate		1.076	1.054	1.129	1.074	1.119	1.070	4.4
Fluoranthene		1.085	0.987	1.018	0.953	0.979	0.975	6.9
Pyrene		1.909	1.783	1.783	1.715	1.740	1.707	8.7
Terphenyl-d14		1.567	1.456	1.419	1.311	1.342	1.344	11.4
Butylbenzylphthalate		0.438	0.472	0.536	0.544	0.582	0.523	9.6
3,3-Dichlorobenzidine			0.437	0.495	0.429	0.476	0.446	7.6
Benzo (a) anthracene		1.374	1.367	1.381	1.282	1.422	1.339	4.9
Chrysene		1.282	1.174	1.247	1.206	1.218	1.204	4.3
Bis (2-ethylhexyl) phthalate		0.691	0.755	0.824	0.789	0.862	0.791	7.0
Di-n-octyl phthalate			1.274	1.428	1.384	1.550	1.434	6.8
Benzo (b) fluoranthene		1.231	1.136	1.287	1.147	1.352	1.222	6.3
Benzo (k) fluoranthene		1.168	1.152	1.093	1.134	1.069	1.090	7.1
Benzo (a) pyrene		1.129	1.102	1.147	1.121	1.192	1.126	3.4
Indeno (1,2,3-cd) pyrene		1.424	1.336	1.430	1.413	1.505	1.410	3.9
Dibenzo (a,h) anthracene		1.159	1.103	1.179	1.147	1.221	1.148	3.9
Benzo (g,h,i) perylene		1.100	1.047	1.140	1.109	1.196	1.110	4.4
1,2,4,5-Tetrachlorobenzene		0.638	0.598	0.601	0.550	0.582	0.577	6.9
1,4-Dioxane		0.622	0.582	0.614	0.600	0.625	0.598	4.1

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