

6C

SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: Alliance

Contract: CHEM02

Lab Code: ACE

SDG No.: Q2987

Instrument ID: BNA_P

Calibration Date(s): 09/08/2025 09/08/2025

Calibration Time(s): 14:41 18:07

LAB FILE ID:		RRFAL1 = BP025684.D			RRFAL2 = BP025685.D		RRFAL3 = BP025686.D	
		RRFAL4 = BP025687.D			RRFAL5 = BP025688.D		RRFAL6 = BP025689.D	
COMPOUND	RRFAL1	RRFAL2	RRFAL3	RRFAL4	RRFAL5	RRFAL6	RRF	% RSD
1,4-Dioxane	0.487	0.522	0.525	0.525	0.524		0.517	3.2
Benzaldehyde		0.981	0.979	0.814	0.578	0.369	0.744	35.8
Hexachloroethane	0.548	0.578	0.592	0.590	0.572		0.576	3.0
Nitrobenzene	0.321	0.368	0.382	0.400	0.391		0.372	8.4
Isophorone	0.704	0.760	0.792	0.792	0.764		0.762	4.7
2-Nitrophenol	0.155	0.177	0.187	0.197	0.193		0.182	9.2
2,4-Dimethylphenol	0.340	0.374	0.372	0.397	0.382		0.373	5.5
Bis(2-Chloroethoxy)methane	0.384	0.442	0.457	0.461	0.445		0.438	7.0
2,4-Dichlorophenol	0.263	0.302	0.317	0.329	0.321		0.306	8.6
Naphthalene	1.021	1.056	1.071	1.066	1.015		1.046	2.5
4-Chloroaniline		0.423	0.440	0.458	0.454	0.425	0.440	3.7
Hexachlorobutadiene	0.225	0.228	0.234	0.233	0.220		0.228	2.5
Phenol		1.498	1.675	1.722	1.773	1.629	1.659	6.3
Caprolactam		0.122	0.127	0.129	0.129	0.121	0.126	2.8
4-Chloro-3-methylphenol	0.321	0.345	0.381	0.378	0.377		0.360	7.3
2-Methylnaphthalene	0.712	0.751	0.776	0.760	0.735		0.747	3.2
Hexachlorocyclopentadiene		0.222	0.272	0.321	0.341	0.355	0.302	18.2
2,4,6-Trichlorophenol	0.338	0.378	0.403	0.420	0.406		0.389	8.3
2,4,5-Trichlorophenol	0.388	0.403	0.433	0.456	0.452		0.426	7.0
1,1-Biphenyl	1.332	1.412	1.450	1.459	1.357		1.402	4.0
2-Chloronaphthalene	1.011	1.114	1.137	1.157	1.099		1.104	5.1
2-Nitroaniline	0.292	0.333	0.356	0.386	0.384		0.350	11.1
Bis(2-Chloroethyl)ether		1.249	1.355	1.358	1.347	1.238	1.309	4.6
Dimethylphthalate	1.478	1.508	1.518	1.501	1.411		1.483	2.9
2,6-Dinitrotoluene	0.263	0.291	0.307	0.322	0.321		0.300	8.2
Acenaphthylene	1.781	1.863	1.902	1.931	1.817		1.859	3.3
3-Nitroaniline		0.253	0.293	0.292	0.264	0.234	0.267	9.5
Acenaphthene	1.157	1.220	1.209	1.224	1.162		1.194	2.7
2,4-Dinitrophenol		0.130	0.167	0.193	0.216	0.217	0.185	20.0
4-Nitrophenol		0.145	0.200	0.224	0.234	0.217	0.204	17.3
Dibenzofuran	1.672	1.756	1.723	1.740	1.648		1.708	2.7
2,4-Dinitrotoluene	0.385	0.443	0.461	0.485	0.477		0.450	8.9
Diethylphthalate	1.518	1.567	1.571	1.506	1.463		1.525	3.0
2-Chlorophenol	1.102	1.288	1.330	1.344	1.359		1.285	8.2

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: CHEM02

Lab Code: ACE

SDG No.: Q2987

Instrument ID: BNA_P

Calibration Date(s): 09/08/2025 09/08/2025

Calibration Time(s): 14:41 18:07

LAB FILE ID:		RRFAL1 = BP025684.D			RRFAL2 = BP025685.D		RRFAL3 = BP025686.D	
		RRFAL4 = BP025687.D			RRFAL5 = BP025688.D		RRFAL6 = BP025689.D	
COMPOUND	RRFAL1	RRFAL2	RRFAL3	RRFAL4	RRFAL5	RRFAL6	RRF	% RSD
Fluorene	1.389	1.447	1.428	1.401	1.310		1.395	3.8
4-Chlorophenyl-phenylether	0.723	0.741	0.723	0.709	0.674		0.714	3.5
4-Nitroaniline		0.281	0.296	0.290	0.248	0.202	0.263	14.9
4,6-Dinitro-2-methylphenol		0.125	0.136	0.147	0.145	0.137	0.138	6.3
N-Nitrosodiphenylamine	0.534	0.595	0.588	0.612	0.563		0.578	5.3
1,2,4,5-Tetrachlorobenzene	0.547	0.586	0.607	0.635	0.600		0.595	5.4
4-Bromophenyl-phenylether	0.197	0.219	0.215	0.223	0.212		0.213	4.7
Hexachlorobenzene	0.231	0.251	0.251	0.247	0.241		0.244	3.4
Atrazine		0.239	0.251	0.250	0.230	0.216	0.237	6.2
Pentachlorophenol		0.140	0.154	0.164	0.167	0.157	0.156	6.8
2-Methylphenol		1.185	1.295	1.313	1.343	1.225	1.272	5.1
Phenanthrene	1.064	1.150	1.114	1.119	1.040		1.097	4.1
Anthracene	1.064	1.149	1.140	1.151	1.060		1.113	4.2
Carbazole		1.026	1.055	1.055	0.996	0.867	1.000	7.8
Di-n-butylphthalate	1.237	1.322	1.351	1.296	1.221		1.285	4.3
Fluoranthene	1.192	1.269	1.271	1.265	1.234		1.246	2.7
Pyrene	1.281	1.323	1.327	1.327	1.254		1.303	2.6
Butylbenzylphthalate	0.567	0.582	0.596	0.594	0.563		0.580	2.6
3,3-Dichlorobenzidine		0.452	0.454	0.461	0.417	0.360	0.429	9.8
2,2-oxybis(1-Chloropropane)		2.063	2.173	2.119	2.101	1.925	2.076	4.5
Benzo(a)anthracene	1.258	1.347	1.358	1.370	1.328		1.332	3.3
Chrysene	1.195	1.284	1.271	1.298	1.221		1.254	3.5
Bis(2-ethylhexyl)phthalate	0.839	0.891	0.895	0.838	0.801		0.853	4.7
Di-n-octyl phthalate		1.409	1.419	1.340	1.320	1.261	1.350	4.8
Benzo(b)fluoranthene	1.082	1.187	1.191	1.225	1.169		1.171	4.6
Benzo(k)fluoranthene	1.162	1.193	1.208	1.195	1.179		1.188	1.5
Benzo(a)pyrene	1.077	1.132	1.119	1.163	1.134		1.125	2.8
Indeno(1,2,3-cd)pyrene	1.303	1.415	1.409	1.464	1.446		1.407	4.4
Dibenzo(a,h)anthracene	1.010	1.153	1.159	1.207	1.172		1.140	6.6
Benzo(g,h,i)perylene	1.068	1.133	1.130	1.200	1.175		1.141	4.4
Acetophenone		2.017	2.153	2.128	2.175	1.874	2.069	6.0
2,3,4,6-Tetrachlorophenol	0.343	0.370	0.388	0.387	0.382		0.374	5.0
1,4-Dioxane-d8	0.476	0.496	0.488	0.485	0.498		0.489	1.8
Pyridine-d5		1.338	1.381	1.532	1.483	1.350	1.417	6.1
Phenol-d5		1.465	1.591	1.678	1.732	1.612	1.616	6.3

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: CHEM02

Lab Code: ACE

SDG No.: Q2987

Instrument ID: BNA_P

Calibration Date(s): 09/08/2025 09/08/2025

Calibration Time(s): 14:41 18:07

LAB FILE ID:		RRFAL1 = BP025684.D			RRFAL2 = BP025685.D		RRFAL3 = BP025686.D	
		RRFAL4 = BP025687.D			RRFAL5 = BP025688.D		RRFAL6 = BP025689.D	
COMPOUND	RRFAL1	RRFAL2	RRFAL3	RRFAL4	RRFAL5	RRFAL6	RRF	% RSD
Bis-(2-Chloroethyl) ether-d8		0.949	1.015	1.021	1.018	0.936	0.988	4.2
2-Chlorophenol-d4	1.121	1.207	1.313	1.319	1.332		1.258	7.3
4-Methylphenol-d8		1.225	1.363	1.391	1.423	1.287	1.338	6.0
Nitrobenzene-d5	0.133	0.150	0.157	0.168	0.161		0.154	8.8
2-Nitrophenol-d4	0.151	0.164	0.180	0.190	0.189		0.175	9.7
2,4-Dichlorophenol-d3	0.267	0.311	0.329	0.339	0.329		0.315	9.2
4-Chloroaniline-d4		0.409	0.450	0.457	0.459	0.432	0.442	4.7
4-Methylphenol		1.264	1.416	1.429	1.471	1.314	1.379	6.3
Dimethylphthalate-d6	1.477	1.552	1.542	1.537	1.451		1.512	3.0
Acenaphthylene-d8	1.559	1.657	1.701	1.712	1.672		1.660	3.7
4-Nitrophenol-d4		0.186	0.248	0.274	0.283	0.263	0.251	15.3
Fluorene-d10	1.218	1.285	1.297	1.278	1.218		1.259	3.0
4,6-Dinitro-2-methylphenol		0.122	0.134	0.143	0.143	0.139	0.136	6.6
Anthracene-d10	0.927	0.998	0.978	0.982	0.928		0.963	3.4
Pyrene-d10	1.053	1.124	1.119	1.105	1.071		1.095	2.8
Benzo(a)pyrene-d12	0.975	1.017	1.040	1.061	1.040		1.027	3.2
N-Nitroso-di-n-propylamine	0.980	1.082	1.165	1.130	1.138		1.099	6.7

All other compounds must meet a minimum RRF of 0.010.

Form VI SV-1