Alliance TECHNICAL GROUP
284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
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SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: ALLI03

Lab Code: CHEM Case No.: Q1872 SAS No.: Q1872 SDG No.: Q1872

Instrument ID: BNA_P Calibration Date(s): 04/14/2025 04/14/2025

Calibration Time(s): 11:06 17:13

LAB FILE ID: RRF2.5 = BP024275.D RRF005 = BP024276.D RRF010 = BP024277.D
RRF020 = BP024278.D RRF040 = BP024279.D RRF050 = BP024280.D

COMPOUND	RRF2.5	RRF005	RRF010	RRF020	RRF040	RRF050	RRF	% RSD
n-Nitrosodimethylamine		0.424	0.443	0.469	0.457	0.442	0.448	3.3
Pyridine		1.265	1.280	1.422	1.396	1.333	1.351	4.4
2-Fluorophenol		1.123	1.145	1.262	1.254	1.210	1.210	4.9
Benzaldehyde		0.943	0.934	0.942	0.836	0.732	0.819	15.7
Aniline		1.439	1.487	1.636	1.587	1.403	1.480	6.5
Phenol-d6		1.463	1.552	1.718	1.748	1.697	1.656	6.7
Phenol		1.475	1.565	1.720	1.750	1.698	1.661	6.5
bis(2-Chloroethyl)ether		1.278	1.317	1.399	1.398	1.321	1.339	3.6
2-Chlorophenol		1.260	1.297	1.412	1.399	1.359	1.350	4.3
1,2-Dichlorobenzene		1.472	1.441	1.502	1.445	1.362	1.424	4.1
1,3-Dichlorobenzene		1.464	1.486	1.534	1.486	1.402	1.454	3.8
1,4-Dichlorobenzene		1.496	1.489	1.550	1.494	1.436	1.475	3.3
Benzyl Alcohol		0.857	0.935	1.106	1.158	1.142	1.071	11.6
2-Methylphenol		0.915	1.002	1.123	1.140	1.105	1.075	8.0
2,2-oxybis(1-Chloropropane)		1.452	1.496	1.517	1.514	1.405	1.447	4.5
Acetophenone		0.470	0.487	0.522	0.519	0.490	0.495	3.8
3+4-Methylphenols		1.224	1.365	1.560	1.592	1.567	1.491	9.6
n-Nitroso-di-n-propylamine	0.941	1.008	1.008	1.088	1.084	1.041	1.025	4.8
Nitrobenzene-d5		0.322	0.337	0.367	0.368	0.357	0.351	4.8
Hexachloroethane		0.532	0.537	0.558	0.542	0.521	0.533	2.9
Nitrobenzene		0.328	0.334	0.362	0.364	0.349	0.347	4.0
Isophorone		0.575	0.601	0.653	0.673	0.650	0.635	5.4
2-Nitrophenol		0.130	0.143	0.172	0.183	0.184	0.170	14.0
2,4-Dimethylphenol		0.180	0.194	0.219	0.226	0.225	0.213	8.8
bis(2-Chloroethoxy)methane		0.412	0.426	0.444	0.452	0.430	0.429	3.5
2,4-Dichlorophenol		0.242	0.268	0.297	0.308	0.307	0.291	8.9
1,2,4-Trichlorobenzene		0.308	0.305	0.318	0.321	0.309	0.311	2.0
Benzoic acid			0.185	0.226	0.242	0.270	0.248	15.6
Naphthalene		1.044	1.054	1.091	1.079	1.037	1.054	2.3
4-Chloroaniline		0.328	0.349	0.389	0.397	0.378	0.371	6.5
Hexachlorobutadiene		0.181	0.181	0.186	0.189	0.181	0.183	1.6
Caprolactam		0.085	0.092	0.111	0.112	0.115	0.107	12.5
4-Chloro-3-methylphenol		0.284	0.310	0.349	0.357	0.356	0.339	8.8
2-Methylnaphthalene		0.700	0.720	0.755	0.755	0.733	0.731	2.9

All other compounds must meet a minimum RRF of 0.010.

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COMPOUND	RRF2.5	RRF005	RRF010	RRF020	RRF040	RRF050	RRF	% RSD
Hexachlorocyclopentadiene		0.162	0.185	0.197	0.218	0.199	0.195	8.9
2,4,6-Trichlorophenol		0.301	0.335	0.369	0.389	0.386	0.366	9.5
2-Fluorobiphenyl		1.332	1.334	1.359	1.366	1.263	1.306	4.1
2,4,5-Trichlorophenol		0.332	0.371	0.409	0.427	0.429	0.405	9.6
1,1-Biphenyl		1.454	1.489	1.528	1.529	1.434	1.470	3.1
2-Chloronaphthalene		1.063	1.100	1.135	1.139	1.085	1.099	2.6
2-Nitroaniline		0.250	0.270	0.317	0.330	0.324	0.308	11.0
Dimethylphthalate		1.429	1.416	1.495	1.495	1.460	1.454	2.2
Acenaphthylene		1.604	1.680	1.793	1.814	1.726	1.730	4.1
2,6-Dinitrotoluene		0.264	0.290	0.319	0.325	0.319	0.309	7.6
3-Nitroaniline		0.250	0.289	0.338	0.354	0.342	0.325	12.2
Acenaphthene		1.087	1.087	1.141	1.126	1.081	1.097	2.4
2,4-Dinitrophenol			0.125	0.163	0.186	0.196	0.182	18.2
4-Nitrophenol		0.190	0.231	0.289	0.305	0.306	0.280	17.9
Dibenzofuran		1.801	1.807	1.871	1.833	1.746	1.793	2.8
2,4-Dinitrotoluene		0.334	0.368	0.439	0.448	0.442	0.421	11.8
Diethylphthalate		1.451	1.484	1.528	1.550	1.499	1.499	2.2
4-Chlorophenyl-phenylether		0.667	0.675	0.699	0.680	0.667	0.674	2.0
Fluorene		1.376	1.397	1.454	1.398	1.368	1.388	2.4
4-Nitroaniline		0.263	0.300	0.353	0.369	0.364	0.344	13.0
4,6-Dinitro-2-methylphenol			0.095	0.118	0.131	0.134	0.126	13.5
n-Nitrosodiphenylamine		0.567	0.585	0.616	0.630	0.591	0.597	3.7
Azobenzene		1.269	1.383	1.446	1.431	1.385	1.378	4.3
2,4,6-Tribromophenol		0.235	0.256	0.278	0.289	0.289	0.277	8.3
4-Bromophenyl-phenylether		0.203	0.212	0.220	0.230	0.220	0.219	4.1
Hexachlorobenzene		0.247	0.247	0.260	0.271	0.261	0.260	3.7
Atrazine		0.176	0.167	0.134	0.121	0.162	0.152	15.6
Pentachlorophenol		0.143	0.155	0.175	0.196	0.197	0.181	13.3
Phenanthrene		1.098	1.086	1.117	1.127	1.070	1.091	2.3
Anthracene		0.995	1.017	1.084	1.111	1.053	1.052	3.9
Carbazole		0.968	0.998	1.060	1.084	1.014	1.027	3.8
Di-n-butylphthalate		1.147	1.230	1.253	1.398	1.310	1.280	6.3
Fluoranthene		1.289	1.274	1.316	1.347	1.280	1.298	2.1
Benzidine		0.144	0.141	0.109	0.436	0.333	0.254	48.4
Pyrene		1.192	1.236	1.345	1.305	1.294	1.271	4.4

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COMPOUND	RRF2.5	RRF005	RRF010	RRF020	RRF040	RRF050	RRF	% RSD
Terphenyl-d14		0.967	0.998	1.055	1.032	1.005	0.992	4.4
Butylbenzylphthalate		0.434	0.488	0.546	0.596	0.580	0.544	11.2
3,3-Dichlorobenzidine		0.345	0.383	0.431	0.486	0.463	0.436	12.3
Benzo(a)anthracene		1.195	1.214	1.292	1.293	1.234	1.243	3.1
Chrysene		1.180	1.181	1.229	1.213	1.173	1.191	1.9
Bis(2-ethylhexyl)phthalate		0.639	0.735	0.779	0.893	0.853	0.798	11.1
Di-n-octyl phthalate		0.945	1.093	1.222	1.461	1.430	1.297	16.5
Benzo(b)fluoranthene		1.130	1.174	1.217	1.258	1.196	1.199	3.3
Benzo(k)fluoranthene		1.088	1.131	1.217	1.212	1.154	1.158	3.9
Benzo(a)pyrene		0.939	0.971	1.057	1.091	1.053	1.034	5.4
Indeno(1,2,3-cd)pyrene		1.252	1.304	1.407	1.444	1.401	1.388	5.8
Dibenzo(a,h)anthracene		1.046	1.089	1.172	1.203	1.169	1.152	5.3
Benzo(g,h,i)perylene		1.073	1.110	1.198	1.218	1.180	1.173	5.0
1,2,4,5-Tetrachlorobenzene		0.522	0.538	0.568	0.576	0.544	0.550	3.3
1,4-Dioxane		0.551	0.517	0.552	0.512	0.481	0.512	6.1
2,3,4,6-Tetrachlorophenol		0.327	0.351	0.380	0.389	0.387	0.373	6.6
1-Methylnaphthalene		0.700	0.714	0.737	0.731	0.712	0.715	2.3

All other compounds must meet a minimum RRF of 0.010.