284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: NOBI03

Lab Code: CHEM Case No.: P5306 SAS No.: P5306 SDG No.: P5306

Instrument ID: MSVOA_Y Calibration Date(s): 12/17/2024 12/17/2024

Heated Purge: (Y/N) Y Calibration Time(s): 10:01 14:51

GC Column: RXI-624 ID: 0.25 (mm)

LAB FILE ID: RRF010 = VY020613.D RRF020 = VY020614.D RRF050 = VY020615.D
RRF100 = VY020616.D RRF150 = VY020617.D RRF005 = VY020619.D

COMPOUND	RRF010	RRF020	RRF050	RRF100	RRF150	RRF005	$\overline{RRF}$	% RSD
Dichlorodifluoromethane	0.678	0.633	0.617	0.601	0.552	0.665	0.625	7.3
Chloromethane	0.490	0.506	0.428	0.426	0.372	0.555	0.463	14.3
Vinyl Chloride	0.443	0.469	0.371	0.440	0.374	0.487	0.431	11.3
Bromomethane	0.291	0.279	0.229	0.282	0.254	0.327	0.277	12.1
Chloroethane	0.273	0.287	0.212	0.272	0.251	0.302	0.266	11.8
Tetrahydrofuran	0.154	0.126	0.133	0.132	0.110	0.116	0.128	11.8
Trichlorofluoromethane	0.981	0.954	0.791	0.878	0.810	0.993	0.901	9.7
1,1,2-Trichlorotrifluoroethane	0.715	0.659	0.603	0.649	0.537	0.712	0.646	10.5
1,1-Dichloroethene	0.701	0.637	0.589	0.638	0.528	0.667	0.627	9.7
Acrylonitrile	0.179	0.144	0.155	0.153	0.132	0.144	0.151	10.4
Acetone	0.141	0.107	0.098	0.110		0.168	0.125	23.2
Carbon Disulfide	2.315	2.083	1.939	2.092	1.669	2.142	2.040	10.7
Methyl tert-butyl Ether	1.799	1.584	1.663	1.652	1.466	1.542	1.618	7.1
Methylene Chloride	0.758	0.683	0.617	0.665	0.596	0.766	0.681	10.3
trans-1,2-Dichloroethene	0.771	0.703	0.710	0.692	0.636	0.707	0.703	6.2
1,1-Dichloroethane	1.401	1.307	1.323	1.300	1.173	1.333	1.306	5.7
2-Butanone	0.245	0.191	0.197	0.195	0.166	0.228	0.204	13.9
Carbon Tetrachloride	0.768	0.715	0.701	0.697	0.656	0.730	0.711	5.2
2,2-Dichloropropane	1.366	1.219	1.184	1.146	1.041	1.451	1.235	12.1
cis-1,2-Dichloroethene	0.875	0.815	0.817	0.797	0.724	0.827	0.809	6.1
Chloroform	1.768	1.465	1.378	1.302	1.186	2.069	1.528	21.6
1,1,1-Trichloroethane	1.298	1.215	1.198	1.183	1.086	1.219	1.200	5.7
1,1-Dichloropropene	0.691	0.637	0.637	0.622	0.581	0.688	0.643	6.5
Benzene	2.055	1.900	1.859	1.847	1.713	1.927	1.883	5.9
1,2-Dichloroethane	0.546	0.499	0.500	0.498	0.452	0.502	0.500	6
Trichloroethene	0.501	0.462	0.458	0.453	0.422	0.481	0.463	5.8
1,2-Dichloropropane	0.473	0.450	0.434	0.440	0.401	0.442	0.440	5.3
Dibromomethane	0.262	0.230	0.231	0.233	0.209	0.232	0.233	7.2
Bromodichloromethane	0.672	0.632	0.632	0.627	0.581	0.626	0.628	4.6
4-Methyl-2-Pentanone	0.364	0.290	0.299	0.301	0.261	0.297	0.302	11.1

* Compounds with required minimum RRF and maximum %RSD values.
All other compounds must meet a minimum RRF of 0.010.
RRF of 1,4-Dioxane = Value should be divide by 1000.

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VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: NOBI03

Lab Code: CHEM Case No.: P5306 SAS No.: P5306 SDG No.: P5306

Instrument ID: MSVOA_Y Calibration Date(s): 12/17/2024 12/17/2024

Heated Purge: (Y/N) Y Calibration Time(s): 10:01 14:51

GC Column: RXI-624 ID: 0.25 (mm)

LAB FILE ID: RRF010 = VY020613.D RRF020 = VY020614.D RRF050 = VY020615.D
RRF100 = VY020616.D RRF150 = VY020617.D RRF005 = VY020619.D

COMPOUND	RRF010	RRF020	RRF050	RRF100	RRF150	RRF005	$\overline{RRF}$	% RSD
Toluene	1.247	1.185	1.167	1.165	1.082	1.176	1.170	4.5
t-1,3-Dichloropropene	0.638	0.572	0.602	0.611	0.556	0.534	0.586	6.5
cis-1,3-Dichloropropene	0.742	0.694	0.706	0.714	0.646	0.661	0.694	5.1
1,1,2-Trichloroethane	0.344	0.306	0.303	0.309	0.276	0.308	0.308	7.1
1,3-Dichloropropane	0.609	0.544	0.555	0.554	0.495	0.534	0.548	6.7
2-Hexanone	0.225	0.183	0.199	0.202	0.175	0.180	0.194	9.5
Dibromochloromethane	0.447	0.402	0.420	0.421	0.381	0.396	0.411	5.7
1,2-Dibromoethane	0.319	0.281	0.288	0.290	0.261	0.277	0.286	6.7
Tetrachloroethene	0.505	0.464	0.452	0.446	0.422	0.533	0.470	8.7
Chlorobenzene	1.501	1.397	1.374	1.381	1.293	1.441	1.398	5
1,1,1,2-Tetrachloroethane	0.529	0.468	0.474	0.476	0.443	0.479	0.478	5.9
Ethyl Benzene	2.747	2.583	2.556	2.543	2.414	2.633	2.579	4.2
m/p-Xylenes	1.031	0.951	0.953	0.942	0.892	0.984	0.959	4.8
o-Xylene	0.956	0.892	0.898	0.887	0.842	0.894	0.895	4.1
Styrene	1.586	1.496	1.502	1.501	1.406	1.478	1.495	3.8
Bromoform	0.302	0.256	0.270	0.272	0.250	0.262	0.269	6.9
Isopropylbenzene	4.636	4.451	4.356	4.284	4.051	4.401	4.363	4.4
1,1,2,2-Tetrachloroethane	0.804	0.702	0.710	0.710	0.635	0.689	0.708	7.7
1,2,3-Trichloropropane	0.581	0.520	0.522	0.511	0.450	0.532	0.519	8.1
Bromobenzene	1.007	0.953	0.947	0.944	0.877	0.974	0.950	4.5
n-propylbenzene	5.566	5.397	5.283	5.151	4.837	5.307	5.257	4.7
2-Chlorotoluene	3.114	3.012	2.940	2.932	2.727	3.020	2.958	4.4
1,3,5-Trimethylbenzene	3.701	3.622	3.560	3.473	3.270	3.616	3.540	4.3
4-Chlorotoluene	3.185	3.123	3.059	2.993	2.786	3.264	3.068	5.5
tert-Butylbenzene	3.378	3.288	3.215	3.178	2.992	3.188	3.206	4
1,2,4-Trimethylbenzene	3.660	3.534	3.536	3.429	3.224	3.454	3.473	4.2
sec-Butylbenzene	4.894	4.777	4.690	4.543	4.263	4.806	4.662	4.9
p-Isopropyltoluene	3.975	3.951	3.916	3.809	3.574	3.838	3.844	3.8
1,3-Dichlorobenzene	1.982	1.912	1.864	1.832	1.702	1.966	1.876	5.5
1,4-Dichlorobenzene	1.980	1.852	1.832	1.797	1.680	2.008	1.858	6.5

* Compounds with required minimum RRF and maximum %RSD values.
All other compounds must meet a minimum RRF of 0.010.
RRF of 1,4-Dioxane = Value should be divide by 1000.

284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
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VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: NOBI03

Lab Code: CHEM Case No.: P5306 SAS No.: P5306 SDG No.: P5306

Instrument ID: MSVOA_Y Calibration Date(s): 12/17/2024 12/17/2024

Heated Purge: (Y/N) Y Calibration Time(s): 10:01 14:51

GC Column: RXI-624 ID: 0.25 (mm)

LAB FILE ID: RRF010 = VY020613.D RRF020 = VY020614.D RRF050 = VY020615.D
RRF100 = VY020616.D RRF150 = VY020617.D RRF005 = VY020619.D

COMPOUND	RRF010	RRF020	RRF050	RRF100	RRF150	RRF005	$\overline{RRF}$	% RSD
n-Butylbenzene	3.755	3.736	3.749	3.599	3.393	3.625	3.643	3.8
1,2-Dichlorobenzene	1.735	1.614	1.606	1.602	1.501	1.694	1.626	5
1,2-Dibromo-3-Chloropropane	0.133	0.107	0.113	0.114	0.107	0.116	0.115	8.4
1,2,4-Trichlorobenzene	0.975	0.927	1.014	1.030	0.990	0.964	0.983	3.8
Hexachlorobutadiene	0.650	0.636	0.657	0.634	0.613	0.667	0.643	3
1,2,3-Trichlorobenzene	0.798	0.791	0.864	0.874	0.846	0.813	0.831	4.2
1,2-Dichloroethane-d4	0.619	0.487	0.569	0.507	0.467	0.636	0.548	13
Dibromofluoromethane	0.385	0.328	0.360	0.325	0.313	0.404	0.352	10.4
Toluene-d8	1.314	1.127	1.271	1.143	1.100	1.445	1.233	10.9
4-Bromofluorobenzene	0.538	0.448	0.480	0.433	0.412	0.585	0.482	13.9
trans-1,4-Dichloro-2-butene	0.270	0.247	0.256	0.262	0.237	0.212	0.247	8.4

* Compounds with required minimum RRF and maximum %RSD values.
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VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: NOBI03

Lab Code: CHEM Case No.: P5306 SAS No.: P5306 SDG No.: P5306

Instrument ID: MSVOA_Y Calibration Date(s): 12/23/2024 12/23/2024

Heated Purge: (Y/N) Y Calibration Time(s): 11:30 13:24

GC Column: RXI-624 ID: 0.25 (mm)

LAB FILE ID: RRF005 = VY020684.D RRF010 = VY020685.D RRF020 = VY020686.D
RRF050 = VY020687.D RRF100 = VY020688.D RRF150 = VY020689.D

COMPOUND	RRF005	RRF010	RRF020	RRF050	RRF100	RRF150	$\overline{RRF}$	% RSD
Dichlorodifluoromethane	0.338	0.391	0.389	0.424	0.391	0.378	0.385	7.2
Chloromethane	0.125	0.165	0.151	0.176	0.161	0.163	0.157	11.1
Vinyl Chloride	0.163	0.196	0.184	0.212	0.193	0.200	0.191	8.6
Bromomethane	0.111	0.133	0.123	0.150	0.144	0.151	0.135	11.9
Chloroethane	0.097	0.115	0.113	0.128	0.120	0.125	0.116	9.3
Tetrahydrofuran	0.035	0.039	0.040	0.053	0.039	0.044	0.042	15.4
Trichlorofluoromethane	0.597	0.687	0.671	0.765	0.698	0.711	0.688	8
1,1,2-Trichlorotrifluoroethane	0.316	0.386	0.379	0.413	0.379	0.381	0.376	8.5
1,1-Dichloroethene	0.279	0.341	0.315	0.371	0.349	0.346	0.333	9.6
Acrylonitrile	0.051	0.056	0.058	0.076	0.059	0.064	0.061	13.8
Acetone	0.059	0.049	0.047	0.058	0.043	0.048	0.051	12.4
Carbon Disulfide	0.654	0.768	0.769	0.978	0.907	0.913	0.831	14.6
Methyl tert-butyl Ether	0.769	0.869	0.904	1.042	0.908	0.949	0.907	9.9
Methylene Chloride	0.324	0.346	0.317	0.370	0.333	0.332	0.337	5.5
trans-1,2-Dichloroethene	0.310	0.350	0.350	0.404	0.370	0.374	0.360	8.7
1,1-Dichloroethane	0.500	0.583	0.597	0.668	0.616	0.611	0.596	9.3
2-Butanone	0.069	0.070	0.067	0.085	0.064	0.071	0.071	10.2
Carbon Tetrachloride	0.472	0.585	0.562	0.654	0.595	0.597	0.577	10.4
2,2-Dichloropropane	0.720	0.768	0.725	0.798	0.735	0.726	0.746	4.2
cis-1,2-Dichloroethene	0.389	0.415	0.418	0.472	0.433	0.442	0.428	6.6
Chloroform	0.787	0.799	0.753	0.841	0.760	0.760	0.783	4.2
1,1,1-Trichloroethane	0.706	0.835	0.809	0.903	0.829	0.828	0.818	7.8
1,1-Dichloropropene	0.302	0.358	0.352	0.402	0.364	0.364	0.357	8.9
Benzene	0.870	1.019	1.009	1.163	1.057	1.058	1.030	9.3
1,2-Dichloroethane	0.271	0.324	0.324	0.384	0.327	0.334	0.327	11
Trichloroethene	0.261	0.306	0.307	0.339	0.310	0.310	0.306	8.3
1,2-Dichloropropane	0.166	0.197	0.205	0.231	0.206	0.209	0.202	10.5
Dibromomethane	0.118	0.131	0.138	0.163	0.144	0.144	0.140	10.6
Bromodichloromethane	0.326	0.384	0.392	0.461	0.404	0.409	0.396	11
4-Methyl-2-Pentanone	0.110	0.111	0.116	0.148	0.114	0.127	0.121	12

* Compounds with required minimum RRF and maximum %RSD values.
All other compounds must meet a minimum RRF of 0.010.
RRF of 1,4-Dioxane = Value should be divide by 1000.

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VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: NOBI03

Lab Code: CHEM Case No.: P5306 SAS No.: P5306 SDG No.: P5306

Instrument ID: MSVOA_Y Calibration Date(s): 12/23/2024 12/23/2024

Heated Purge: (Y/N) Y Calibration Time(s): 11:30 13:24

GC Column: RXI-624 ID: 0.25 (mm)

LAB FILE ID: RRF005 = VY020684.D RRF010 = VY020685.D RRF020 = VY020686.D
RRF050 = VY020687.D RRF100 = VY020688.D RRF150 = VY020689.D

COMPOUND	RRF005	RRF010	RRF020	RRF050	RRF100	RRF150	$\overline{RRF}$	% RSD
Toluene	0.582	0.687	0.693	0.792	0.723	0.729	0.701	9.8
t-1,3-Dichloropropene	0.277	0.323	0.349	0.411	0.366	0.376	0.351	13.3
cis-1,3-Dichloropropene	0.304	0.369	0.386	0.452	0.402	0.405	0.386	12.7
1,1,2-Trichloroethane	0.154	0.175	0.176	0.214	0.180	0.186	0.181	10.8
1,3-Dichloropropane	0.242	0.286	0.291	0.353	0.296	0.305	0.296	12.1
2-Hexanone	0.057	0.067	0.073	0.100	0.078	0.087	0.077	19.4
Dibromochloromethane	0.244	0.271	0.287	0.335	0.292	0.303	0.289	10.6
1,2-Dibromoethane	0.132	0.172	0.168	0.203	0.166	0.176	0.170	13.5
Tetrachloroethene	0.274	0.339	0.331	0.362	0.341	0.346	0.332	9.1
Chlorobenzene	0.811	0.932	0.924	1.022	0.956	0.961	0.934	7.5
1,1,1,2-Tetrachloroethane	0.321	0.368	0.360	0.398	0.369	0.373	0.365	6.9
Ethyl Benzene	1.397	1.698	1.668	1.841	1.722	1.731	1.676	8.9
m/p-Xylenes	0.531	0.645	0.644	0.703	0.656	0.656	0.639	8.9
o-Xylene	0.493	0.610	0.599	0.670	0.621	0.611	0.600	9.7
Styrene	0.823	1.026	1.020	1.119	1.039	1.033	1.010	9.8
Bromoform	0.180	0.195	0.208	0.247	0.206	0.219	0.209	11
Isopropylbenzene	2.983	3.415	3.372	3.667	3.515	3.453	3.401	6.7
1,1,2,2-Tetrachloroethane	0.386	0.444	0.430	0.498	0.424	0.445	0.438	8.3
1,2,3-Trichloropropane	0.305	0.372	0.341	0.349	0.346	0.357	0.345	6.5
Bromobenzene	0.667	0.779	0.752	0.853	0.794	0.802	0.774	8
n-propylbenzene	3.382	3.920	3.786	4.131	3.928	3.859	3.834	6.5
2-Chlorotoluene	1.923	2.228	2.169	2.368	2.221	2.188	2.183	6.7
1,3,5-Trimethylbenzene	2.458	2.840	2.814	3.058	2.911	2.845	2.821	7
4-Chlorotoluene	1.938	2.214	2.232	2.421	2.262	2.290	2.226	7.1
tert-Butylbenzene	2.340	2.609	2.608	2.842	2.703	2.647	2.625	6.3
1,2,4-Trimethylbenzene	2.330	2.749	2.772	3.039	2.858	2.786	2.756	8.5
sec-Butylbenzene	3.092	3.493	3.553	3.813	3.671	3.584	3.534	6.9
p-Isopropyltoluene	2.598	3.016	3.111	3.376	3.260	3.178	3.090	8.8
1,3-Dichlorobenzene	1.340	1.499	1.475	1.675	1.528	1.510	1.504	7.1
1,4-Dichlorobenzene	1.290	1.485	1.473	1.605	1.502	1.489	1.474	6.9

* Compounds with required minimum RRF and maximum %RSD values.
All other compounds must meet a minimum RRF of 0.010.
RRF of 1,4-Dioxane = Value should be divide by 1000.

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Lab Code: CHEM Case No.: P5306 SAS No.: P5306 SDG No.: P5306

Instrument ID: MSVOA_Y Calibration Date(s): 12/23/2024 12/23/2024

Heated Purge: (Y/N) Y Calibration Time(s): 11:30 13:24

GC Column: RXI-624 ID: 0.25 (mm)

LAB FILE ID: RRF005 = VY020684.D RRF010 = VY020685.D RRF020 = VY020686.D
RRF050 = VY020687.D RRF100 = VY020688.D RRF150 = VY020689.D

COMPOUND	RRF005	RRF010	RRF020	RRF050	RRF100	RRF150	$\overline{RRF}$	% RSD
n-Butylbenzene	2.232	2.504	2.602	2.825	2.750	2.705	2.603	8.2
1,2-Dichlorobenzene	1.121	1.291	1.286	1.431	1.311	1.308	1.291	7.7
1,2-Dibromo-3-Chloropropane	0.075	0.080	0.079	0.096	0.080	0.090	0.083	9.3
1,2,4-Trichlorobenzene	0.563	0.667	0.713	0.882	0.866	0.908	0.767	18.2
Hexachlorobutadiene	0.466	0.530	0.561	0.628	0.622	0.618	0.571	11.3
1,2,3-Trichlorobenzene	0.465	0.545	0.607	0.747	0.721	0.760	0.641	18.8
1,2-Dichloroethane-d4		0.407	0.359	0.511	0.469	0.468	0.443	13.5
Dibromofluoromethane		0.286	0.250	0.330	0.303	0.299	0.294	10
Toluene-d8		0.891	0.815	1.244	1.209	1.174	1.066	18.6
4-Bromofluorobenzene		0.373	0.339	0.424	0.408	0.396	0.388	8.6
trans-1,4-Dichloro-2-butene	0.128	0.153	0.155	0.195	0.154	0.167	0.159	13.8

* Compounds with required minimum RRF and maximum %RSD values.
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