

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: RUTW01
Lab Code: CHEM Case No.: Q1215 SAS No.: Q1215 SDG No.: Q1215
Instrument ID: MSVOA_Y Calibration Date(s): 02/03/2025 02/03/2025
Heated Purge: (Y/N) Y Calibration Time(s): 10:35 12:44
GC Column: RXI-624 ID: 0.25 (mm)

LAB FILE ID:		RRF005 = VY021020.D		RRF010 = VY021021.D		RRF020 = VY021022.D		RRF050 = VY021023.D		RRF100 = VY021024.D		RRF150 = VY021025.D	
COMPOUND		RRF005	RRF010	RRF020	RRF050	RRF100	RRF150	RRF	% RSD				
Dichlorodifluoromethane		0.503	0.438	0.413	0.445	0.419	0.457	0.446	7.2				
Chloromethane		0.524	0.445	0.408	0.421	0.404	0.437	0.440	10.1				
Vinyl Chloride		0.535	0.441	0.404	0.424	0.408	0.443	0.443	10.9				
Bromomethane		0.356	0.275	0.249	0.258	0.246	0.257	0.274	15.2				
Chloroethane		0.316	0.270	0.240	0.255	0.239	0.261	0.263	10.7				
Trichlorofluoromethane		0.924	0.814	0.755	0.789	0.743	0.813	0.806	8				
1,1,2-Trichlorotrifluoroethane		0.587	0.507	0.472	0.494	0.466	0.510	0.506	8.6				
Tert butyl alcohol		0.042	0.039	0.035	0.032	0.028	0.032	0.035	14.8				
1,1-Dichloroethene		0.529	0.471	0.441	0.469	0.452	0.498	0.477	6.8				
Acetone		0.087	0.083	0.078	0.077	0.070	0.083	0.080	7.5				
Carbon Disulfide		1.673	1.456	1.389	1.475	1.421	1.556	1.495	7				
Methyl tert-butyl Ether		1.271	1.185	1.168	1.216	1.124	1.265	1.205	4.8				
Methyl Acetate		0.300	0.282	0.268	0.287	0.253	0.291	0.280	6.1				
Methylene Chloride		0.570	0.506	0.473	0.485	0.452	0.494	0.497	8.2				
trans-1,2-Dichloroethene		0.578	0.518	0.502	0.522	0.490	0.538	0.525	5.8				
1,1-Dichloroethane		1.104	0.948	0.923	0.947	0.911	0.989	0.970	7.3				
Cyclohexane		1.093	0.885	0.807	0.827	0.779	0.852	0.874	13				
2-Butanone		0.134	0.133	0.126	0.135	0.121	0.139	0.131	5.2				
Carbon Tetrachloride		0.604	0.551	0.522	0.547	0.522	0.581	0.555	5.9				
cis-1,2-Dichloroethene		0.662	0.584	0.577	0.599	0.575	0.626	0.604	5.7				
Bromochloromethane		0.373	0.368	0.381	0.413	0.374	0.394	0.384	4.4				
Chloroform		1.107	1.016	0.947	0.982	0.921	1.006	0.997	6.5				
1,1,1-Trichloroethane		1.067	0.901	0.867	0.912	0.863	0.949	0.927	8.2				
Methylcyclohexane		0.609	0.550	0.536	0.588	0.570	0.629	0.581	6.1				
Benzene		1.512	1.382	1.327	1.378	1.323	1.436	1.393	5.1				
1,2-Dichloroethane		0.435	0.390	0.367	0.379	0.352	0.395	0.386	7.4				
Trichloroethene		0.401	0.355	0.341	0.349	0.338	0.375	0.360	6.7				
1,2-Dichloropropane		0.357	0.329	0.319	0.327	0.311	0.341	0.331	5				
Bromodichloromethane		0.528	0.500	0.469	0.482	0.459	0.506	0.491	5.2				
4-Methyl-2-Pentanone		0.202	0.211	0.202	0.216	0.194	0.224	0.208	5.2				

* 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.

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name:	<u>CHEMTECH</u>	Contract:	<u>RUTW01</u>
Lab Code:	<u>CHEM</u>	Case No.:	<u>Q1215</u>
Instrument ID:	<u>MSVOA_Y</u>	SAS No.:	<u>Q1215</u>
Heated Purge: (Y/N)	<u>Y</u>	SDG No.:	<u>Q1215</u>
GC Column:	<u>RXI-624</u>	Calibration Date(s):	<u>02/03/2025</u>
ID:	<u>0.25 (mm)</u>	Calibration Time(s):	<u>10:35</u>
			<u>12:44</u>

LAB FILE ID:	RRF005 = VY021020.D		RRF010 = VY021021.D		RRF020 = VY021022.D			
	RRF050 = VY021023.D		RRF100 = VY021024.D		RRF150 = VY021025.D			
COMPOUND	RRF005	RRF010	RRF020	RRF050	RRF100	RRF150	RRF	% RSD
Toluene	0.913	0.856	0.840	0.876	0.844	0.918	0.875	3.9
t-1,3-Dichloropropene	0.441	0.426	0.419	0.449	0.427	0.483	0.441	5.3
cis-1,3-Dichloropropene	0.528	0.516	0.502	0.525	0.501	0.560	0.522	4.2
1,1,2-Trichloroethane	0.261	0.235	0.231	0.235	0.217	0.242	0.237	6.1
2-Hexanone	0.124	0.131	0.128	0.144	0.129	0.149	0.134	7.4
Dibromochloromethane	0.359	0.330	0.324	0.329	0.310	0.346	0.333	5.2
1,2-Dibromoethane	0.232	0.223	0.215	0.225	0.203	0.233	0.222	5
Tetrachloroethene	0.409	0.367	0.352	0.361	0.351	0.387	0.371	6.2
Chlorobenzene	1.237	1.115	1.054	1.080	1.048	1.146	1.113	6.4
Ethyl Benzene	2.101	1.899	1.865	1.943	1.914	2.093	1.969	5.2
m/p-Xylenes	0.772	0.723	0.709	0.733	0.711	0.773	0.737	3.9
o-Xylene	0.723	0.658	0.660	0.681	0.670	0.729	0.687	4.6
Styrene	1.154	1.118	1.105	1.156	1.128	1.208	1.145	3.2
Bromoform	0.231	0.222	0.217	0.216	0.205	0.229	0.220	4.4
Isopropylbenzene	4.122	3.695	3.673	3.809	3.871	4.263	3.906	6.1
1,1,2,2-Tetrachloroethane	0.707	0.646	0.635	0.632	0.597	0.680	0.650	6
1,3-Dichlorobenzene	1.937	1.686	1.660	1.683	1.652	1.824	1.740	6.6
1,4-Dichlorobenzene	1.906	1.702	1.632	1.665	1.599	1.778	1.714	6.6
1,2-Dichlorobenzene	1.668	1.524	1.440	1.460	1.405	1.567	1.510	6.4
1,2-Dibromo-3-Chloropropane	0.099	0.099	0.089	0.097	0.089	0.109	0.097	7.6
1,2,4-Trichlorobenzene	0.796	0.754	0.743	0.875	0.873	1.004	0.841	11.6
1,2,3-Trichlorobenzene	0.641	0.640	0.626	0.733	0.726	0.832	0.700	11.4
1,2-Dichloroethane-d4	0.566	0.557	0.478	0.503	0.454	0.516	0.512	8.5
Dibromofluoromethane	0.351	0.326	0.301	0.313	0.297	0.333	0.320	6.4
Toluene-d8	1.295	1.237	1.139	1.202	1.158	1.288	1.220	5.3
4-Bromofluorobenzene	0.417	0.408	0.371	0.403	0.375	0.418	0.399	5.2

* 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.