

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name:	CHEMTECH	Contract:	PORT06
Lab Code:	CHEM	SAS No.:	<u>Q1194</u>
Instrument ID:	MSVOA_Y	SDG No.:	<u>Q1194</u>
Heated Purge:	(Y/N) Y	Calibration Date(s):	<u>01/24/2025</u>
GC Column:	RXI-624	Calibration Time(s):	<u>14:03</u> <u>16:21</u>
	ID: 0.25 (mm)		

LAB FILE ID:	RRF050 = VY020927.D	RRF100 = VY020928.D	RRF150 = VY020929.D					
COMPOUND	RRF050	RRF100	RRF150	RRF020	RRF010	RRF005	RRF	% RSD
Dichlorodifluoromethane	0.384	0.335	0.373	0.388	0.395	0.394	0.378	6
Chloromethane	0.281	0.237	0.277	0.283	0.309	0.301	0.281	8.9
Vinyl Chloride	0.313	0.271	0.310	0.310	0.327	0.309	0.307	6
Bromomethane	0.199	0.167	0.191	0.203	0.241	0.244	0.208	14.4
Chloroethane	0.185	0.158	0.180	0.183	0.190	0.182	0.180	6.2
Trichlorofluoromethane	0.665	0.582	0.669	0.669	0.685	0.638	0.651	5.7
1,1,2-Trichlorotrifluoroethane	0.436	0.382	0.438	0.432	0.454	0.437	0.430	5.7
1,1-Dichloroethene	0.408	0.365	0.421	0.412	0.428	0.401	0.406	5.4
Acetone	0.058	0.044	0.052	0.060	0.058	0.052	0.054	10.7
Carbon Disulfide	1.276	1.131	1.297	1.274	1.304	1.237	1.253	5.1
Methyl tert-butyl Ether	0.977	0.843	0.981	1.017	0.982	0.841	0.940	8.2
Methyl Acetate	0.208	0.168	0.194	0.210	0.198	0.192	0.195	7.7
Methylene Chloride	0.404	0.349	0.403	0.413	0.427	0.415	0.402	6.8
trans-1,2-Dichloroethene	0.446	0.388	0.448	0.443	0.460	0.440	0.438	5.8
1,1-Dichloroethane	0.740	0.647	0.745	0.742	0.752	0.708	0.722	5.5
Cyclohexane	0.680	0.605	0.681	0.714	0.762	0.817	0.710	10.3
2-Butanone	0.097	0.078	0.091	0.104	0.100	0.079	0.091	12
Carbon Tetrachloride	0.493	0.438	0.500	0.487	0.502	0.479	0.483	4.9
cis-1,2-Dichloroethene	0.501	0.449	0.510	0.500	0.516	0.486	0.494	4.9
Bromochloromethane	0.298	0.296	0.297	0.308	0.313	0.232	0.291	10.1
Chloroform	0.777	0.682	0.790	0.786	0.792	0.759	0.764	5.5
1,1,1-Trichloroethane	0.755	0.663	0.766	0.753	0.774	0.738	0.742	5.4
Methylcyclohexane	0.559	0.502	0.572	0.545	0.553	0.524	0.543	4.7
Benzene	1.202	1.061	1.205	1.197	1.234	1.156	1.176	5.2
1,2-Dichloroethane	0.305	0.262	0.300	0.307	0.302	0.281	0.293	6.1
Trichloroethene	0.322	0.290	0.332	0.326	0.336	0.309	0.319	5.3
1,2-Dichloropropane	0.268	0.236	0.270	0.267	0.270	0.246	0.260	5.8
Bromodichloromethane	0.406	0.358	0.411	0.398	0.412	0.371	0.393	5.8
4-Methyl-2-Pentanone	0.162	0.132	0.155	0.170	0.152	0.120	0.149	12.7
Toluene	0.779	0.692	0.785	0.774	0.786	0.722	0.756	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.

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Lab Code:	CHEM	Case No.:	<u>Q1194</u>
Instrument ID:	MSVOA_Y	Calibration Date(s):	<u>01/24/2025</u>
Heated Purge:	(Y/N) <u>Y</u>	Calibration Time(s):	<u>14:03</u> <u>16:21</u>
GC Column:	RXI-624	ID:	0.25 (mm)

LAB FILE ID:	RRF050 = VY020927.D	RRF100 = VY020928.D	RRF150 = VY020929.D					
COMPOUND	RRF050	RRF100	RRF150	RRF020	RRF010	RRF005	RRF	% RSD
t-1,3-Dichloropropene	0.379	0.336	0.387	0.381	0.365	0.312	0.360	8.3
cis-1,3-Dichloropropene	0.446	0.396	0.455	0.446	0.436	0.399	0.430	6
1,1,2-Trichloroethane	0.207	0.179	0.205	0.217	0.201	0.189	0.199	6.9
2-Hexanone	0.106	0.089	0.103	0.110	0.096	0.070	0.096	15.4
Dibromochloromethane	0.290	0.257	0.296	0.296	0.289	0.257	0.281	6.7
1,2-Dibromoethane	0.199	0.169	0.198	0.203	0.200	0.174	0.191	7.8
Tetrachloroethene	0.352	0.312	0.358	0.348	0.361	0.358	0.348	5.3
Chlorobenzene	0.992	0.881	1.012	0.991	1.019	0.972	0.978	5.1
Ethyl Benzene	1.761	1.565	1.777	1.722	1.762	1.650	1.706	4.9
m/p-Xylenes	0.674	0.596	0.679	0.667	0.677	0.627	0.653	5.2
o-Xylene	0.632	0.559	0.635	0.619	0.634	0.583	0.610	5.3
Styrene	1.054	0.926	1.051	1.027	1.043	0.899	1.000	6.9
Bromoform	0.208	0.177	0.202	0.213	0.203	0.178	0.197	7.9
Isopropylbenzene	3.562	3.149	3.639	3.423	3.496	3.394	3.444	4.9
1,1,2,2-Tetrachloroethane	0.570	0.469	0.548	0.577	0.566	0.485	0.536	8.8
1,3-Dichlorobenzene	1.588	1.371	1.583	1.554	1.638	1.540	1.546	6
1,4-Dichlorobenzene	1.555	1.334	1.553	1.523	1.602	1.544	1.518	6.2
1,2-Dichlorobenzene	1.386	1.182	1.378	1.346	1.408	1.335	1.339	6.1
1,2-Dibromo-3-Chloropropane	0.085	0.070	0.086	0.090	0.083	0.069	0.080	10.9
1,2,4-Trichlorobenzene	0.840	0.782	0.938	0.792	0.786	0.697	0.806	9.9
1,2,3-Trichlorobenzene	0.704	0.651	0.797	0.672	0.648	0.589	0.677	10.3
1,2-Dichloroethane-d4	0.446	0.369	0.386	0.409	0.404	0.378	0.399	7
Dibromofluoromethane	0.320	0.276	0.284	0.287	0.291	0.264	0.287	6.6
Toluene-d8	1.254	1.081	1.109	1.121	1.135	0.981	1.113	7.9
4-Bromofluorobenzene	0.416	0.359	0.364	0.390	0.375	0.319	0.371	8.8

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