

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: ARDM01

Lab Code: CHEM Case No.: Q2264 SAS No.: Q2264 SDG No.: Q2264

Instrument ID: MSVOA_X Calibration Date(s): 06/10/2025 06/10/2025

Heated Purge: (Y/N) N Calibration Time(s): 20:19 21:44

GC Column: DB-624UI ID: 0.18 (mm)

LAB FILE ID:		RRF005 = VX046614.D		RRF020 = VX046615.D		RRF050 = VX046616.D		RRF100 = VX046617.D		RRF150 = VX046618.D		RRF =	
COMPOUND		RRF005	RRF020	RRF050	RRF100	RRF150	RRF						% RSD
Chloromethane		2.074	2.119	1.934	2.075	1.846		2.010					5.7
Vinyl Chloride		1.964	2.039	1.866	2.045	1.772		1.937					6
Bromomethane		1.292	1.283	1.129	1.091	0.990		1.157					11.2
Chloroethane		1.143	1.108	1.106	1.162	1.046		1.113					4
Trichlorofluoromethane		2.825	2.970	2.572	2.960	2.547		2.775					7.4
1,1-Dichloroethene		1.851	1.842	1.661	1.911	1.690		1.791					6.1
Acrolein		0.228	0.207	0.189	0.261	0.221		0.221					12.2
Acrylonitrile		0.972	0.976	0.832	0.985	0.873		0.928					7.5
Methylene Chloride		2.161	2.078	1.819	2.024	1.799		1.976					8.1
trans-1,2-Dichloroethene		1.935	1.920	1.756	1.939	1.731		1.856					5.6
1,1-Dichloroethane		3.710	3.704	3.211	3.694	3.291		3.522					7.1
Carbon Tetrachloride		0.524	0.556	0.490	0.542	0.477		0.518					6.5
Chloroform		3.918	3.837	3.244	3.783	3.345		3.625					8.5
1,1,1-Trichloroethane		0.599	0.628	0.537	0.613	0.539		0.583					7.3
Benzene		1.464	1.500	1.310	1.415	1.244		1.387					7.7
1,2-Dichloroethane		0.570	0.582	0.468	0.536	0.467		0.525					10.4
Trichloroethene		0.375	0.366	0.339	0.352	0.308		0.348					7.5
1,2-Dichloropropane		0.361	0.369	0.334	0.352	0.314		0.346					6.4
Bromodichloromethane		0.526	0.576	0.504	0.548	0.480		0.527					7.1
Toluene		1.663	1.749	1.544	1.682	1.492		1.626					6.5
t-1,3-Dichloropropene		0.444	0.475	0.463	0.517	0.463		0.472					5.8
cis-1,3-Dichloropropene		0.503	0.547	0.515	0.566	0.505		0.527					5.3
1,1,2-Trichloroethane		0.341	0.352	0.322	0.324	0.291		0.326					7.1
2-Chloroethyl vinyl ether		0.252	0.286	0.268	0.296	0.193		0.259					15.6
Dibromochloromethane		0.375	0.406	0.388	0.398	0.354		0.384					5.3
Tetrachloroethene		0.346	0.332	0.299	0.312	0.283		0.314					8
Chlorobenzene		1.106	1.099	1.014	1.067	0.957		1.049					6
Ethyl Benzene		1.878	1.964	1.821	1.919	1.700		1.856					5.5
m/p-Xylenes		0.688	0.734	0.681	0.704	0.626		0.687					5.7
o-Xylene		0.665	0.703	0.652	0.676	0.610		0.661					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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COMPOUND	RRF005	RRF020	RRF050	RRF100	RRF150	RRF	RRF	% RSD	
Bromoform	0.226	0.249	0.241	0.249	0.222		0.237	5.3	
1,1,2,2-Tetrachloroethane	0.571	0.565	0.502	0.513	0.456		0.521	9.1	
1,3-Dichlorobenzene	0.833	0.825	0.753	0.768	0.691		0.774	7.5	
1,4-Dichlorobenzene	0.850	0.837	0.745	0.766	0.695		0.779	8.3	
1,2-Dichlorobenzene	0.808	0.799	0.721	0.740	0.656		0.745	8.3	
1,2-Dichloroethane-d4	2.549	2.442	2.160	2.504	2.486		2.428	6.4	
Toluene-d8	1.359	1.354	1.307	1.342	1.337		1.340	1.5	
4-Bromofluorobenzene	0.515	0.506	0.505	0.497	0.505		0.506	1.2	

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