

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: GENV01
 Lab Code: CHEM Case No.: Q1449 SAS No.: Q1449 SDG No.: Q1449
 Instrument ID: MSVOA_N Calibration Date(s): 02/18/2025 02/18/2025
 Heated Purge: (Y/N) N Calibration Time(s): 11:09 14:18
 GC Column: RXI-624 ID: 0.25 (mm)

LAB FILE ID:		RRF100 = VN085772.D		RRF050 = VN085773.D		RRF010 = VN085775.D			
		RRF005 = VN085776.D		RRF001 = VN085777.D		RRF020 = VN085779.D			
COMPOUND	RRF100	RRF050	RRF010	RRF005	RRF001	RRF020	RRF	% RSD	
Chloromethane	0.647	0.622	0.742	0.662	0.817	0.604	0.682	11.9	
Vinyl Chloride	0.698	0.678	0.796	0.729	0.761	0.674	0.723	6.7	
Bromomethane	0.434	0.424	0.543	0.496		0.413	0.462	12.1	
Chloroethane	0.447	0.417	0.516	0.470	0.584	0.438	0.479	12.9	
Tert butyl alcohol	0.072	0.078	0.092	0.083		0.073	0.080	9.9	
1,1-Dichloroethene	0.555	0.527	0.591	0.571	0.514	0.531	0.548	5.3	
Acetone	0.195	0.184	0.217	0.208	0.227	0.198	0.205	7.7	
Carbon Disulfide	1.565	1.473	1.804	1.604	1.848	1.455	1.625	10.2	
Methyl tert-butyl Ether	1.747	1.658	1.714	1.560	1.438	1.659	1.629	7	
Methylene Chloride	0.628	0.593	0.710	0.666	0.736	0.625	0.660	8.3	
trans-1,2-Dichloroethene	0.585	0.549	0.621	0.599	0.623	0.566	0.590	5.1	
1,1-Dichloroethane	1.097	1.035	1.216	1.125	1.073	1.089	1.106	5.6	
2-Butanone	0.308	0.294	0.331	0.299	0.279	0.299	0.302	5.6	
Carbon Tetrachloride	0.573	0.528	0.583	0.545	0.533	0.538	0.550	4.1	
cis-1,2-Dichloroethene	0.705	0.661	0.740	0.685	0.646	0.671	0.685	4.9	
Chloroform	1.120	1.070	1.269	1.164	1.171	1.141	1.156	5.7	
1,1,1-Trichloroethane	1.012	0.967	1.089	1.058	1.030	1.012	1.028	4.1	
Benzene	1.581	1.441	1.568	1.421	1.420	1.474	1.484	4.9	
1,2-Dichloroethane	0.499	0.456	0.527	0.464	0.472	0.483	0.483	5.4	
Trichloroethene	0.370	0.339	0.384	0.350	0.395	0.348	0.364	6.1	
1,2-Dichloropropane	0.377	0.347	0.385	0.344	0.339	0.357	0.358	5.2	
Bromodichloromethane	0.569	0.518	0.571	0.533	0.512	0.541	0.541	4.6	
4-Methyl-2-Pentanone	0.414	0.379	0.392	0.342	0.290	0.388	0.367	12.1	
Toluene	0.993	0.902	0.924	0.820	0.689	0.891	0.870	12	
t-1,3-Dichloropropene	0.576	0.521	0.518	0.471	0.408	0.505	0.500	11.3	
cis-1,3-Dichloropropene	0.626	0.566	0.582	0.513	0.442	0.559	0.548	11.6	
1,1,2-Trichloroethane	0.361	0.331	0.370	0.346	0.324	0.344	0.346	5	
2-Hexanone	0.296	0.273	0.266	0.227	0.184	0.270	0.253	16	
Dibromochloromethane	0.440	0.402	0.431	0.384	0.355	0.405	0.403	7.7	
Tetrachloroethene	0.384	0.361	0.420	0.391	0.393	0.375	0.387	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	RRF100	RRF050	RRF010	RRF005	RRF001	RRF020	RRF	% RSD	
Chlorobenzene	1.183	1.110	1.212	1.139	1.109	1.119	1.145	3.8	
Ethyl Benzene	2.105	1.904	1.830	1.665	1.424	1.838	1.794	12.8	
m/p-Xylenes	0.809	0.741	0.725	0.610	0.481	0.722	0.682	17.2	
o-Xylene	0.771	0.704	0.666	0.584	0.502	0.659	0.648	14.5	
Styrene	1.327	1.203	1.060	0.885	0.744	1.133	1.059	20.2	
Bromoform	0.327	0.303	0.322	0.301	0.274	0.309	0.306	6.2	
1,1,2,2-Tetrachloroethane	1.073	1.053	1.304	1.218	1.271	1.154	1.179	8.8	
1,2-Dichloroethane-d4	0.698	0.613	0.588	0.767		0.575	0.648	12.6	
Dibromofluoromethane	0.375	0.317	0.294	0.357		0.293	0.327	11.5	
Toluene-d8	1.444	1.199	1.017	1.212		1.080	1.190	13.8	
4-Bromofluorobenzene	0.503	0.419	0.323	0.370		0.346	0.392	18.2	

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