

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: TETR06
Lab Code: CHEM Case No.: Q1423 SAS No.: Q1423 SDG No.: Q1423
Instrument ID: MSVOA_X Calibration Date(s): 02/10/2025 02/10/2025
Heated Purge: (Y/N) N Calibration Time(s): 10:25 12:28
GC Column: DB-624UI ID: 0.18 (mm)

LAB FILE ID:		RRF001 = VX044868.D		RRF005 = VX044869.D		RRF020 = VX044870.D			
		RRF050 = VX044871.D		RRF100 = VX044872.D		RRF150 = VX044873.D			
COMPOUND	RRF001	RRF005	RRF020	RRF050	RRF100	RRF150	RRF	% RSD	
Chloromethane	0.958	0.860	0.862	0.843	0.805	0.793	0.854	6.9	
Vinyl Chloride	0.839	0.846	0.847	0.808	0.807	0.814	0.827	2.3	
Bromomethane		0.248	0.252	0.249	0.242	0.246	0.247	1.4	
Chloroethane	0.500	0.288	0.280	0.341	0.249	0.181	0.307	35.4	
Trichlorofluoromethane	1.062	1.066	1.096	1.029	1.013	1.008	1.046	3.3	
1,1,2-Trichlorotrifluoroethane	0.583	0.647	0.668	0.626	0.631	0.637	0.632	4.5	
1,1-Dichloroethene	0.647	0.639	0.661	0.630	0.632	0.657	0.644	2	
Acetone	0.305	0.292	0.298	0.293	0.285	0.292	0.294	2.3	
Carbon Disulfide	1.689	1.732	1.786	1.762	1.789	1.846	1.767	3	
Methyl tert-butyl Ether	1.941	2.065	2.130	2.046	2.011	2.110	2.050	3.4	
Methylene Chloride	0.747	0.717	0.741	0.704	0.695	0.720	0.721	2.8	
trans-1,2-Dichloroethene	0.608	0.622	0.657	0.640	0.633	0.644	0.634	2.7	
1,1-Dichloroethane	1.155	1.257	1.292	1.227	1.209	1.257	1.233	3.9	
2-Butanone	0.422	0.472	0.504	0.506	0.477	0.487	0.478	6.4	
Carbon Tetrachloride	0.457	0.466	0.478	0.453	0.445	0.459	0.460	2.4	
cis-1,2-Dichloroethene	0.680	0.783	0.812	0.758	0.758	0.779	0.762	5.9	
Chloroform	1.167	1.209	1.268	1.169	1.153	1.208	1.196	3.5	
1,1,1-Trichloroethane	1.014	1.003	1.051	1.005	0.984	1.028	1.014	2.3	
Methylcyclohexane	0.509	0.571	0.667	0.622	0.634	0.635	0.606	9.4	
Benzene	1.370	1.488	1.577	1.470	1.429	1.453	1.465	4.7	
1,2-Dichloroethane	0.417	0.465	0.502	0.472	0.462	0.482	0.467	6.1	
Trichloroethene	0.293	0.340	0.367	0.335	0.332	0.343	0.335	7.2	
1,2-Dichloropropane	0.343	0.354	0.389	0.367	0.360	0.372	0.364	4.3	
Bromodichloromethane	0.428	0.481	0.514	0.500	0.500	0.513	0.489	6.6	
4-Methyl-2-Pentanone	0.439	0.514	0.562	0.554	0.506	0.498	0.512	8.6	
Toluene	0.776	0.872	0.957	0.898	0.866	0.864	0.872	6.7	
t-1,3-Dichloropropene	0.417	0.451	0.518	0.514	0.528	0.543	0.495	10	
cis-1,3-Dichloropropene	0.452	0.511	0.587	0.577	0.587	0.599	0.552	10.5	
1,1,2-Trichloroethane	0.307	0.342	0.362	0.341	0.331	0.329	0.335	5.4	
2-Hexanone	0.313	0.360	0.406	0.404	0.369	0.362	0.369	9.3	

* 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	RRF001	RRF005	RRF020	RRF050	RRF100	RRF150	RRF	% RSD	
Dibromochloromethane	0.317	0.342	0.381	0.373	0.368	0.370	0.359	6.7	
Tetrachloroethene	0.306	0.311	0.343	0.310	0.307	0.314	0.315	4.4	
Chlorobenzene	0.969	1.093	1.140	1.096	1.071	1.076	1.074	5.3	
Ethyl Benzene	1.690	1.873	2.021	1.935	1.923	1.929	1.895	5.9	
m/p-Xylenes	0.616	0.700	0.754	0.724	0.706	0.694	0.699	6.6	
o-Xylene	0.661	0.721	0.747	0.707	0.691	0.681	0.701	4.4	
Styrene	0.909	1.124	1.249	1.199	1.161	1.139	1.130	10.4	
Bromoform	0.186	0.247	0.272	0.280	0.276	0.287	0.258	14.7	
Isopropylbenzene	3.735	4.012	4.347	4.045	3.940	4.076	4.026	4.9	
1,1,2,2-Tetrachloroethane	1.429	1.403	1.438	1.366	1.305	1.360	1.383	3.6	
1,3-Dichlorobenzene	1.616	1.669	1.741	1.679	1.663	1.703	1.678	2.5	
1,4-Dichlorobenzene	1.662	1.712	1.762	1.686	1.660	1.701	1.697	2.2	
1,2-Dichlorobenzene	1.512	1.713	1.763	1.666	1.604	1.639	1.650	5.3	
1,2-Dichloroethane-d4		0.764	0.718	0.723	0.707	0.747	0.732	3.2	
Dibromofluoromethane		0.335	0.322	0.320	0.320	0.328	0.325	2	
Toluene-d8		1.239	1.249	1.239	1.208	1.212	1.229	1.5	
4-Bromofluorobenzene		0.404	0.410	0.431	0.415	0.412	0.414	2.5	

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