

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: <u>CHEMTECH</u>	Contract: <u>TETR06</u>
Lab Code: <u>CHEM</u> Case No.: <u>Q1401</u>	SAS No.: <u>Q1401</u> SDG No.: <u>Q1401</u>
Instrument ID: <u>MSVOA_Y</u>	Calibration Date(s): <u>02/25/2025</u> <u>02/25/2025</u>
Heated Purge: (Y/N) <u>Y</u>	Calibration Time(s): <u>12:40</u> <u>16:49</u>
GC Column: <u>RXI-624</u> ID: <u>0.25</u> (mm)	

LAB FILE ID:	RRF005 = VY021309.D	RRF010 = VY021310.D	RRF020 = VY021311.D	RRF050 = VY021312.D	RRF150 = VY021314.D	RRF100 = VY021316.D		
COMPOUND	RRF005	RRF010	RRF020	RRF050	RRF150	RRF100	RRF	% RSD
Chloromethane	0.709	0.628	0.566	0.602	0.597	0.570	0.612	8.6
Vinyl Chloride	0.648	0.591	0.554	0.627	0.606	0.580	0.601	5.6
Bromomethane	0.505	0.414	0.361	0.387	0.371	0.359	0.399	13.9
Chloroethane	0.414	0.365	0.347	0.390	0.365	0.355	0.373	6.7
Trichlorofluoromethane	0.976	0.915	0.838	0.936	0.909	0.870	0.908	5.3
1,1,2-Trichlorotrifluoroethane	0.546	0.529	0.491	0.537	0.541	0.508	0.525	4.1
1,1-Dichloroethene	0.545	0.489	0.462	0.515	0.518	0.490	0.503	5.7
Acetone	0.122	0.102	0.094	0.128	0.104	0.103	0.109	11.9
Carbon Disulfide	1.800	1.652	1.540	1.734	1.678	1.620	1.671	5.4
Methyl tert-butyl Ether	1.312	1.268	1.267	1.410	1.410	1.307	1.329	4.9
Methylene Chloride	0.733	0.598	0.539	0.585	0.547	0.529	0.588	12.8
trans-1,2-Dichloroethene	0.589	0.536	0.507	0.573	0.570	0.547	0.554	5.4
1,1-Dichloroethane	1.101	1.032	0.961	1.083	1.058	1.007	1.040	4.9
2-Butanone	0.162	0.155	0.156	0.183	0.171	0.159	0.164	6.5
Carbon Tetrachloride	0.567	0.562	0.509	0.581	0.567	0.535	0.554	4.8
cis-1,2-Dichloroethene	0.649	0.610	0.588	0.665	0.656	0.626	0.632	4.7
Chloroform	1.128	1.046	0.977	1.095	1.062	1.017	1.054	5.1
1,1,1-Trichloroethane	1.020	0.956	0.889	0.986	0.979	0.927	0.960	4.8
Methylcyclohexane	0.597	0.613	0.584	0.662	0.663	0.618	0.623	5.3
Benzene	1.515	1.445	1.359	1.533	1.472	1.409	1.456	4.5
1,2-Dichloroethane	0.422	0.417	0.393	0.434	0.419	0.394	0.413	3.9
Trichloroethene	0.375	0.360	0.341	0.382	0.375	0.354	0.364	4.3
1,2-Dichloropropane	0.360	0.350	0.329	0.371	0.357	0.336	0.350	4.5
Bromodichloromethane	0.519	0.503	0.477	0.541	0.528	0.495	0.511	4.5
4-Methyl-2-Pentanone	0.220	0.227	0.239	0.267	0.264	0.238	0.243	7.9
Toluene	0.887	0.888	0.852	0.965	0.936	0.897	0.904	4.4
t-1,3-Dichloropropene	0.438	0.436	0.434	0.515	0.509	0.474	0.468	8
cis-1,3-Dichloropropene	0.541	0.526	0.528	0.591	0.581	0.546	0.552	5
1,1,2-Trichloroethane	0.250	0.251	0.238	0.267	0.254	0.237	0.249	4.4
2-Hexanone	0.133	0.146	0.155	0.183	0.178	0.162	0.159	11.9

* 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 Name: CHEMTECH Contract: TETR06
 Lab Code: CHEM Case No.: Q1401 SAS No.: Q1401 SDG No.: Q1401
 Instrument ID: MSVOA_Y Calibration Date(s): 02/25/2025 02/25/2025
 Heated Purge: (Y/N) Y Calibration Time(s): 12:40 16:49
 GC Column: RXI-624 ID: 0.25 (mm)

LAB FILE ID:		RRF005 = VY021309.D		RRF010 = VY021310.D		RRF020 = VY021311.D			
		RRF050 = VY021312.D		RRF150 = VY021314.D		RRF100 = VY021316.D			
COMPOUND	RRF005	RRF010	RRF020	RRF050	RRF150	RRF100	RRF	% RSD	
Dibromochloromethane	0.340	0.325	0.322	0.368	0.352	0.332	0.340	5.2	
Tetrachloroethene	0.405	0.380	0.351	0.389	0.386	0.355	0.378	5.5	
Chlorobenzene	1.171	1.111	1.039	1.176	1.159	1.078	1.122	5	
Ethyl Benzene	1.938	1.941	1.826	2.130	2.115	1.979	1.988	5.8	
m/p-Xylenes	0.735	0.738	0.696	0.802	0.783	0.733	0.748	5.1	
o-Xylene	0.689	0.675	0.648	0.758	0.739	0.694	0.701	5.9	
Styrene	1.081	1.128	1.091	1.277	1.257	1.159	1.166	7.2	
Bromoform	0.221	0.221	0.211	0.246	0.236	0.214	0.225	5.9	
Isopropylbenzene	3.744	3.605	3.437	3.976	4.009	3.783	3.759	5.8	
1,1,2,2-Tetrachloroethane	0.700	0.649	0.621	0.697	0.689	0.624	0.663	5.5	
1,3-Dichlorobenzene	1.812	1.729	1.598	1.792	1.777	1.668	1.730	4.8	
1,4-Dichlorobenzene	1.825	1.710	1.593	1.751	1.744	1.629	1.709	5	
1,2-Dichlorobenzene	1.599	1.499	1.408	1.575	1.564	1.446	1.515	5.1	
1,2-Dichloroethane-d4	0.681	0.510	0.548	0.551	0.566	0.547	0.567	10.3	
Dibromofluoromethane	0.370	0.299	0.317	0.325	0.334	0.322	0.328	7.2	
Toluene-d8	1.420	1.158	1.230	1.251	1.290	1.253	1.267	6.8	
4-Bromofluorobenzene	0.490	0.387	0.404	0.423	0.435	0.420	0.426	8.2	

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