

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: TETR06

Lab Code: CHEM Case No.: P5316 SAS No.: P5316 SDG No.: P5316

Instrument ID: MSVOA_Y Calibration Date(s): 12/17/2024 12/17/2024

Heated Purge: (Y/N) Y Calibration Time(s): 10:01 14:51

GC Column: RXI-624 ID: 0.25 (mm)

LAB FILE ID:	RRF010 = VY020613.D			RRF020 = VY020614.D		RRF050 = VY020615.D		RRF100 = VY020616.D		RRF150 = VY020617.D		RRF005 = VY020619.D	
COMPOUND	RRF010	RRF020	RRF050	RRF100	RRF150	RRF005	RRF	% RSD					
Vinyl Chloride	0.443	0.469	0.371	0.440	0.374	0.487	0.431	11.3					
1,1-Dichloroethene	0.701	0.637	0.589	0.638	0.528	0.667	0.627	9.7					
Acetone	0.141	0.107	0.098	0.110		0.168	0.125	23.2					
Methyl tert-butyl Ether	1.799	1.584	1.663	1.652	1.466	1.542	1.618	7.1					
Methylene Chloride	0.758	0.683	0.617	0.665	0.596	0.766	0.681	10.3					
trans-1,2-Dichloroethene	0.771	0.703	0.710	0.692	0.636	0.707	0.703	6.2					
1,1-Dichloroethane	1.401	1.307	1.323	1.300	1.173	1.333	1.306	5.7					
2-Butanone	0.245	0.191	0.197	0.195	0.166	0.228	0.204	13.9					
Carbon Tetrachloride	0.768	0.715	0.701	0.697	0.656	0.730	0.711	5.2					
cis-1,2-Dichloroethene	0.875	0.815	0.817	0.797	0.724	0.827	0.809	6.1					
Chloroform	1.768	1.465	1.378	1.302	1.186	2.069	1.528	21.6					
1,1,1-Trichloroethane	1.298	1.215	1.198	1.183	1.086	1.219	1.200	5.7					
Benzene	2.055	1.900	1.859	1.847	1.713	1.927	1.883	5.9					
1,2-Dichloroethane	0.546	0.499	0.500	0.498	0.452	0.502	0.500	6					
Trichloroethene	0.501	0.462	0.458	0.453	0.422	0.481	0.463	5.8					
Toluene	1.247	1.185	1.167	1.165	1.082	1.176	1.170	4.5					
Tetrachloroethene	0.505	0.464	0.452	0.446	0.422	0.533	0.470	8.7					
Chlorobenzene	1.501	1.397	1.374	1.381	1.293	1.441	1.398	5					
Ethyl Benzene	2.747	2.583	2.556	2.543	2.414	2.633	2.579	4.2					
m/p-Xylenes	1.031	0.951	0.953	0.942	0.892	0.984	0.959	4.8					
o-Xylene	0.956	0.892	0.898	0.887	0.842	0.894	0.895	4.1					
n-propylbenzene	5.566	5.397	5.283	5.151	4.837	5.307	5.257	4.7					
tert-Butylbenzene	3.378	3.288	3.215	3.178	2.992	3.188	3.206	4					
1,2,4-Trimethylbenzene	3.660	3.534	3.536	3.429	3.224	3.454	3.473	4.2					
sec-Butylbenzene	4.894	4.777	4.690	4.543	4.263	4.806	4.662	4.9					
1,3-Dichlorobenzene	1.982	1.912	1.864	1.832	1.702	1.966	1.876	5.5					
1,4-Dichlorobenzene	1.980	1.852	1.832	1.797	1.680	2.008	1.858	6.5					
n-Butylbenzene	3.755	3.736	3.749	3.599	3.393	3.625	3.643	3.8					
1,2-Dichlorobenzene	1.735	1.614	1.606	1.602	1.501	1.694	1.626	5					
1,2-Dichloroethane-d4	0.619	0.487	0.569	0.507	0.467	0.636	0.548	13					

* 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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Dibromofluoromethane	0.385	0.328	0.360	0.325	0.313	0.404	0.352	10.4
Toluene-d8	1.314	1.127	1.271	1.143	1.100	1.445	1.233	10.9
4-Bromofluorobenzene	0.538	0.448	0.480	0.433	0.412	0.585	0.482	13.9
1,4-Dioxane	3.173	2.476	2.673	2.757	2.467	2.462	2.668	10.4

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