

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name:	<u>CHEMTECH</u>	Contract:	<u>CHEM02</u>
Lab Code:	<u>CHEM</u>	Case No.:	<u>P5288</u>
Instrument ID:	<u>MSVOA_X</u>	SAS No.:	<u>P5288</u>
Heated Purge: (Y/N)	<u>N</u>	SDG No.:	<u>P5288</u>
GC Column:	<u>DB-624UI</u>	Calibration Date(s):	<u>12/11/2024</u>
ID:	<u>0.18 (mm)</u>	Calibration Time(s):	<u>10:41</u>
			<u>13:00</u>

LAB FILE ID:		RRF001 = VX044219.D		RRF005 = VX044220.D		RRF020 = VX044221.D		
		RRF050 = VX044222.D		RRF100 = VX044223.D		RRF150 = VX044224.D		
COMPOUND	RRF001	RRF005	RRF020	RRF050	RRF100	RRF150	RRF	% RSD
Dichlorodifluoromethane	0.877	0.749	0.721	0.788	0.745	0.759	0.773	7.2
Chloromethane	0.743	0.747	0.735	0.756	0.715	0.740	0.739	1.8
Vinyl Chloride	0.805	0.790	0.782	0.786	0.722	0.737	0.770	4.3
Ethyl Acetate	0.536	0.512	0.551	0.581	0.565	0.574	0.553	4.7
Bromomethane		0.590	0.513	0.551	0.529	0.548	0.546	5.3
Chloroethane	0.464	0.520	0.515	0.532	0.422	0.426	0.480	10.2
Trichlorofluoromethane	1.582	1.519	1.514	1.541	1.478	1.414	1.508	3.8
1,1,2-Trichlorotrifluoroethane	0.633	0.628	0.592	0.615	0.573	0.620	0.610	3.8
Tert butyl alcohol		0.103	0.111	0.120	0.133	0.151	0.124	15.4
1,1-Dichloroethene	0.512	0.560	0.544	0.596	0.554	0.596	0.560	5.7
Acrolein		0.179	0.163	0.185	0.174	0.193	0.179	6.3
Acrylonitrile	0.319	0.333	0.332	0.358	0.343	0.367	0.342	5.2
Acetone	0.359	0.353	0.331	0.381	0.348	0.372	0.357	5
Carbon Disulfide	0.841	0.872	0.937	1.150	1.230	1.362	1.065	20
Methyl tert-butyl Ether	2.054	2.188	2.127	2.305	2.170	2.279	2.187	4.3
Methyl Acetate	0.965	0.889	0.883	0.953	0.888	0.951	0.922	4.2
Methylene Chloride	0.679	0.678	0.666	0.689	0.648	0.668	0.671	2.1
trans-1,2-Dichloroethene	0.600	0.556	0.577	0.619	0.595	0.626	0.596	4.4
Vinyl Acetate	1.238	1.571	1.732	1.912	1.842	1.958	1.709	15.7
1,1-Dichloroethane	1.071	1.173	1.154	1.244	1.171	1.218	1.172	5.1
Cyclohexane		0.931	0.963	0.977	0.922	0.938	0.946	2.4
2-Butanone	0.449	0.491	0.510	0.548	0.509	0.541	0.508	7.1
Carbon Tetrachloride	0.484	0.443	0.432	0.501	0.502	0.529	0.482	7.8
2,2-Dichloropropane	0.782	0.896	0.956	1.045	0.992	1.039	0.952	10.5
cis-1,2-Dichloroethene	0.809	0.737	0.728	0.772	0.731	0.762	0.756	4.1
Bromochloromethane	0.578	0.577	0.576	0.592	0.556	0.568	0.575	2.1
Chloroform	1.389	1.312	1.272	1.343	1.272	1.310	1.316	3.4
1,1,1-Trichloroethane	0.993	1.019	1.077	1.176	1.119	1.173	1.093	7.1
Methylcyclohexane	0.514	0.503	0.527	0.555	0.550	0.562	0.535	4.5
1,1-Dichloropropene	0.489	0.442	0.422	0.444	0.445	0.444	0.448	4.9

* 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.

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: CHEM02
Lab Code: CHEM Case No.: P5288 SAS No.: P5288 SDG No.: P5288
Instrument ID: MSVOA_X Calibration Date(s): 12/11/2024 12/11/2024
Heated Purge: (Y/N) N Calibration Time(s): 10:41 13:00
GC Column: DB-624UI ID: 0.18 (mm)

LAB FILE ID:		RRF001 = VX044219.D		RRF005 = VX044220.D		RRF020 = VX044221.D			
		RRF050 = VX044222.D		RRF100 = VX044223.D		RRF150 = VX044224.D			
COMPOUND	RRF001	RRF005	RRF020	RRF050	RRF100	RRF150	RRF	% RSD	
Benzene	1.353	1.353	1.332	1.410	1.353	1.353	1.359	1.9	
1,2-Dichloroethane	0.554	0.557	0.545	0.584	0.556	0.564	0.560	2.4	
Trichloroethene	0.356	0.325	0.325	0.349	0.337	0.342	0.339	3.8	
1,2-Dichloropropane	0.368	0.345	0.330	0.357	0.341	0.342	0.347	3.8	
Dibromomethane	0.247	0.267	0.259	0.286	0.276	0.281	0.269	5.5	
Bromodichloromethane	0.380	0.411	0.436	0.508	0.518	0.538	0.465	13.9	
4-Methyl-2-Pentanone	0.506	0.551	0.547	0.590	0.561	0.582	0.556	5.4	
Toluene	0.908	0.830	0.816	0.877	0.841	0.849	0.853	4	
t-1,3-Dichloropropene	0.338	0.357	0.439	0.514	0.526	0.558	0.455	20.3	
cis-1,3-Dichloropropene	0.404	0.441	0.488	0.567	0.558	0.583	0.507	14.6	
1,1,2-Trichloroethane	0.337	0.332	0.335	0.358	0.334	0.340	0.339	2.8	
1,3-Dichloropropane	0.578	0.593	0.585	0.613	0.583	0.590	0.590	2.1	
2-Chloroethyl Vinyl ether	0.222	0.268	0.256	0.293	0.285	0.296	0.270	10.5	
2-Hexanone	0.323	0.386	0.408	0.439	0.420	0.438	0.403	10.8	
Dibromochloromethane	0.254	0.248	0.306	0.370	0.378	0.398	0.326	20.2	
1,2-Dibromoethane	0.303	0.322	0.344	0.369	0.364	0.370	0.345	8.1	
Tetrachloroethene	0.359	0.354	0.317	0.321	0.312	0.327	0.332	6	
Chlorobenzene	1.130	1.068	1.039	1.088	1.036	1.067	1.071	3.3	
1,1,1,2-Tetrachloroethane	0.301	0.305	0.334	0.375	0.368	0.383	0.344	10.5	
Hexachloroethane	0.358	0.396	0.406	0.507	0.542	0.567	0.463	18.9	
Ethyl Benzene	1.853	1.806	1.816	1.927	1.831	1.875	1.851	2.4	
m/p-Xylenes	0.653	0.660	0.665	0.707	0.683	0.707	0.679	3.5	
o-Xylene	0.686	0.683	0.685	0.703	0.671	0.697	0.687	1.6	
Styrene	1.007	1.019	1.099	1.162	1.148	1.177	1.102	6.7	
Bromoform	0.137	0.173	0.194	0.247	0.269	0.297	0.219	28	
Isopropylbenzene	3.665	3.976	3.920	3.912	3.747	3.747	3.828	3.2	
1,1,2,2-Tetrachloroethane	1.410	1.350	1.332	1.336	1.256	1.293	1.329	3.9	
1,2,3-Trichloropropane	1.272	1.075	1.113	1.124	1.059	1.055	1.116	7.3	
Bromobenzene	1.009	0.934	0.890	0.942	0.899	0.895	0.928	4.9	
n-propylbenzene	3.739	4.263	4.319	4.502	4.283	4.277	4.231	6.1	

* 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.

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name:	<u>CHEMTECH</u>	Contract:	<u>CHEM02</u>
Lab Code:	<u>CHEM</u>	Case No.:	<u>P5288</u>
Instrument ID:	<u>MSVOA_X</u>	SAS No.:	<u>P5288</u>
Heated Purge: (Y/N)	<u>N</u>	SDG No.:	<u>P5288</u>
GC Column:	<u>DB-624UI</u>	Calibration Date(s):	<u>12/11/2024</u>
ID:	<u>0.18 (mm)</u>	Calibration Time(s):	<u>10:41</u>
			<u>13:00</u>

LAB FILE ID:		RRF001 = VX044219.D		RRF005 = VX044220.D		RRF020 = VX044221.D		
		RRF050 = VX044222.D		RRF100 = VX044223.D		RRF150 = VX044224.D		
COMPOUND	RRF001	RRF005	RRF020	RRF050	RRF100	RRF150	RRF	% RSD
2-Chlorotoluene	2.940	2.847	2.707	2.771	2.582	2.569	2.736	5.4
1,3,5-Trimethylbenzene	3.066	3.222	3.224	3.308	3.097	3.128	3.174	2.9
4-Chlorotoluene	2.964	3.051	3.013	3.114	3.011	3.026	3.030	1.7
tert-Butylbenzene	3.170	3.077	3.105	3.259	3.106	3.106	3.137	2.1
1,2,4-Trimethylbenzene	2.699	3.254	3.207	3.290	3.148	3.105	3.117	6.9
sec-Butylbenzene	3.679	3.748	3.840	3.986	3.804	3.793	3.808	2.7
p-Isopropyltoluene	2.949	3.061	3.254	3.347	3.197	3.211	3.170	4.5
1,3-Dichlorobenzene	1.519	1.634	1.640	1.683	1.625	1.628	1.621	3.3
1,4-Dichlorobenzene	1.811	1.640	1.632	1.672	1.621	1.623	1.667	4.4
n-Butylbenzene	2.249	2.387	2.659	2.844	2.889	2.964	2.665	10.9
1,2-Dichlorobenzene	1.745	1.666	1.668	1.726	1.654	1.643	1.684	2.5
1,2-Dibromo-3-Chloropropane	0.222	0.226	0.241	0.273	0.279	0.306	0.258	13
1,2,4-Trichlorobenzene	0.770	0.827	0.891	0.978	1.013	1.058	0.923	12.2
Hexachlorobutadiene	0.390	0.349	0.367	0.369	0.371	0.381	0.371	3.8
Naphthalene	2.925	3.346	3.629	3.835	3.866	4.015	3.603	11.2
1,2,3-Trichlorobenzene	0.854	0.885	0.984	1.019	1.055	1.077	0.979	9.3
1,2-Dichloroethane-d4		0.934	0.778	0.921	0.862	0.888	0.877	7.1
Dibromofluoromethane		0.353	0.298	0.367	0.360	0.354	0.346	8
Toluene-d8		1.148	1.010	1.251	1.229	1.203	1.168	8.3
4-Bromofluorobenzene		0.390	0.351	0.442	0.432	0.426	0.408	9.2

* 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.