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VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: REMI02
 Lab Code: CHEM Case No.: O2505 SAS No.: O2505 SDG No.: O2505
 Instrument ID: MSVOA_Y Calibration Date(s): 04/24/2023 04/24/2023
 Heated Purge: (Y/N) Y Calibration Time(s): 11:26 13:19
 GC Column: RXI-624 ID: 0.25 (mm)

LAB FILE ID:		RRF005 = VY013393.D		RRF010 = VY013394.D		RRF020 = VY013395.D			
		RRF050 = VY013396.D		RRF100 = VY013397.D		RRF150 = VY013398.D			
COMPOUND	RRF005	RRF010	RRF020	RRF050	RRF100	RRF150	RRF	% RSD	
Dichlorodifluoromethane	0.389	0.497	0.490	0.489	0.500	0.487	0.475	9	
Chloromethane	0.634	0.672	0.622	0.623	0.615	0.656	0.637	3.5	
Vinyl Chloride	0.628	0.631	0.616	0.620	0.643	0.668	0.634	3	
Bromomethane	0.542	0.524	0.484	0.472	0.483	0.528	0.506	5.8	
Chloroethane	0.432	0.422	0.414	0.415	0.447	0.500	0.438	7.5	
Trichlorofluoromethane	0.938	0.978	0.979	0.953	0.969	0.949	0.961	1.8	
1,1,2-Trichlorotrifluoroethane	0.536	0.542	0.549	0.523	0.546	0.539	0.539	1.7	
1,1-Dichloroethene	0.506	0.512	0.522	0.525	0.523	0.529	0.520	1.7	
Acetone	0.225	0.198	0.189	0.237	0.211	0.235	0.216	9.2	
Carbon Disulfide	1.555	1.728	1.695	1.727	1.701	1.722	1.688	3.9	
Methyl tert-butyl Ether	1.455	1.450	1.431	1.556	1.489	1.621	1.500	4.9	
Methyl Acetate	0.749	0.705	0.663	0.724	0.667	0.770	0.713	6	
Methylene Chloride	1.590	1.194	0.879	0.665	0.600	0.613	0.924	42.9	
trans-1,2-Dichloroethene	0.536	0.588	0.564	0.581	0.558	0.581	0.568	3.4	
1,1-Dichloroethane	1.034	1.071	1.034	1.057	1.019	1.089	1.050	2.5	
Cyclohexane	1.062	0.970	0.959	0.943	0.973	0.982	0.981	4.2	
2-Butanone	0.272	0.264	0.247	0.288	0.265	0.306	0.274	7.5	
Carbon Tetrachloride	0.521	0.557	0.546	0.553	0.560	0.545	0.547	2.6	
cis-1,2-Dichloroethene	0.615	0.634	0.618	0.633	0.613	0.653	0.628	2.5	
Bromochloromethane	0.366	0.399	0.391	0.369	0.339	0.373	0.373	5.7	
Chloroform	1.033	1.085	1.032	1.040	1.009	1.066	1.044	2.6	
1,1,1-Trichloroethane	0.926	0.942	0.922	0.938	0.923	0.940	0.932	1	
Methylcyclohexane	0.566	0.573	0.596	0.628	0.667	0.662	0.615	7.1	
Benzene	1.438	1.501	1.447	1.487	1.456	1.492	1.470	1.8	
1,2-Dichloroethane	0.476	0.501	0.468	0.491	0.464	0.485	0.481	2.9	
Trichloroethene	0.397	0.392	0.389	0.395	0.386	0.387	0.391	1.1	
1,2-Dichloropropane	0.373	0.391	0.380	0.387	0.377	0.398	0.384	2.4	
Bromodichloromethane	0.503	0.524	0.504	0.522	0.505	0.529	0.515	2.3	
4-Methyl-2-Pentanone	0.348	0.358	0.344	0.390	0.371	0.405	0.369	6.5	
Toluene	0.830	0.903	0.877	0.904	0.881	0.913	0.885	3.4	
t-1,3-Dichloropropene	0.499	0.547	0.537	0.564	0.546	0.576	0.545	4.9	
cis-1,3-Dichloropropene	0.595	0.612	0.600	0.624	0.609	0.637	0.613	2.5	
1,1,2-Trichloroethane	0.311	0.310	0.297	0.308	0.293	0.308	0.305	2.4	

* 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 Code:	CHEM	Case No.:	O2505
Instrument ID:	MSVOA_Y	SAS No.:	O2505
Heated Purge: (Y/N)	Y	SDG No.:	O2505
GC Column:	RXI-624	Calibration Date(s):	04/24/2023
ID:	0.25 (mm)	Calibration Time(s):	11:26
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COMPOUND	RRF005	RRF010	RRF020	RRF050	RRF100	RRF150	RRF	% RSD
2-Hexanone	0.245	0.259	0.249	0.291	0.273	0.294	0.269	7.8
Dibromochloromethane	0.356	0.377	0.367	0.383	0.360	0.376	0.370	2.8
1,2-Dibromoethane	0.303	0.311	0.279	0.308	0.289	0.306	0.299	4.2
Tetrachloroethene	0.384	0.384	0.368	0.383	0.386	0.375	0.380	1.8
Chlorobenzene	1.077	1.116	1.074	1.094	1.087	1.087	1.089	1.4
Ethyl Benzene	1.759	1.878	1.878	1.955	1.996	2.009	1.913	4.9
m/p-Xylenes	0.654	0.725	0.726	0.744	0.761	0.754	0.727	5.3
o-Xylene	0.634	0.672	0.680	0.714	0.713	0.713	0.688	4.7
Styrene	1.029	1.147	1.139	1.187	1.189	1.208	1.150	5.6
Bromoform	0.294	0.287	0.277	0.291	0.279	0.285	0.285	2.3
Isopropylbenzene	3.365	3.588	3.680	3.898	4.055	4.105	3.782	7.6
1,1,2,2-Tetrachloroethane	0.939	0.908	0.867	0.953	0.918	0.961	0.924	3.7
1,3-Dichlorobenzene	1.777	1.763	1.716	1.774	1.721	1.724	1.746	1.6
1,4-Dichlorobenzene	1.832	1.830	1.710	1.772	1.721	1.716	1.764	3.2
1,2-Dichlorobenzene	1.580	1.595	1.521	1.622	1.563	1.579	1.577	2.1
1,2-Dibromo-3-Chloropropane	0.166	0.165	0.145	0.164	0.162	0.177	0.163	6.2
1,2,4-Trichlorobenzene	0.974	0.935	0.873	0.951	0.950	0.957	0.940	3.8
1,2,3-Trichlorobenzene	0.837	0.830	0.788	0.861	0.859	0.853	0.838	3.3
1,2-Dichloroethane-d4	0.619	0.590	0.575	0.520	0.476	0.508	0.548	10
Dibromofluoromethane	0.340	0.336	0.327	0.296	0.291	0.297	0.314	7
Toluene-d8	1.174	1.219	1.193	1.049	1.034	1.043	1.119	7.6
4-Bromofluorobenzene	0.413	0.412	0.394	0.352	0.345	0.347	0.377	8.7

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