

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: <u>Alliance</u>	Contract: <u>FIRS02</u>
Lab Code: <u>ACE</u>	SDG No.: <u>Q2814</u>
Instrument ID: <u>MSVOA_X</u>	Calibration Date(s): <u>07/21/2025</u> <u>07/21/2025</u>
Heated Purge: (Y/N) <u>N</u>	Calibration Time(s): <u>09:47</u> <u>11:32</u>
GC Column: <u>DB-624UI</u> ID: <u>0.18</u> (mm)	

LAB FILE ID:		RRF001 = VX047055.D		RRF005 = VX047056.D		RRF020 = VX047057.D			
		RRF050 = VX047058.D		RRF100 = VX047059.D		RRF150 = VX047060.D			
COMPOUND	RRF001	RRF005	RRF020	RRF050	RRF100	RRF150	RRF	% RSD	
Dichlorodifluoromethane	0.516	0.640	0.610	0.768	0.721	0.729	0.664	14.1	
Chloromethane	0.582	0.668	0.626	0.699	0.667	0.701	0.657	7	
Vinyl Chloride	0.578	0.739	0.700	0.780	0.716	0.723	0.706	9.7	
Bromomethane		0.408	0.400	0.448	0.353	0.309	0.384	14	
Chloroethane	0.547	0.482	0.436	0.470	0.423	0.424	0.464	10.3	
Trichlorofluoromethane	0.853	1.123	1.058	1.158	1.054	1.071	1.053	10.1	
1,1,2-Trichlorotrifluoroethane	0.558	0.671	0.646	0.675	0.627	0.631	0.634	6.7	
1,1-Dichloroethene	0.553	0.659	0.635	0.674	0.624	0.625	0.628	6.7	
Acetone	0.268	0.267	0.246	0.252	0.232	0.231	0.250	6.5	
Carbon Disulfide	1.633	1.926	1.823	1.936	1.783	1.805	1.818	6.1	
Methyl tert-butyl Ether	1.640	2.011	1.960	2.055	1.891	1.916	1.912	7.6	
Methyl Acetate	0.657	0.686	0.636	0.688	0.651	0.662	0.663	3.1	
Methylene Chloride	0.635	0.751	0.708	0.722	0.653	0.651	0.686	6.8	
trans-1,2-Dichloroethene	0.531	0.701	0.667	0.695	0.632	0.632	0.643	9.7	
1,1-Dichloroethane	1.044	1.327	1.245	1.300	1.177	1.193	1.214	8.4	
Cyclohexane		1.221	1.187	1.215	1.099	1.101	1.164	5.2	
2-Butanone	0.317	0.376	0.371	0.391	0.363	0.362	0.363	6.8	
Carbon Tetrachloride	0.526	0.612	0.566	0.595	0.531	0.540	0.562	6.4	
cis-1,2-Dichloroethene	0.670	0.832	0.777	0.807	0.739	0.744	0.761	7.6	
Bromochloromethane	0.457	0.608	0.568	0.609	0.559	0.550	0.559	9.9	
Chloroform	1.180	1.357	1.252	1.289	1.175	1.164	1.236	6.2	
1,1,1-Trichloroethane	0.961	1.156	1.101	1.127	1.041	1.057	1.074	6.5	
Methylcyclohexane	0.583	0.693	0.674	0.699	0.640	0.653	0.657	6.5	
Benzene	1.327	1.656	1.573	1.650	1.451	1.480	1.523	8.4	
1,2-Dichloroethane	0.468	0.587	0.558	0.576	0.512	0.517	0.536	8.5	
Trichloroethene	0.350	0.427	0.392	0.420	0.371	0.383	0.391	7.5	
1,2-Dichloropropane	0.334	0.419	0.393	0.408	0.363	0.371	0.381	8.2	
Bromodichloromethane	0.494	0.626	0.579	0.617	0.543	0.557	0.569	8.7	
4-Methyl-2-Pentanone	0.371	0.477	0.470	0.497	0.440	0.440	0.449	9.8	
Toluene	0.762	1.047	0.987	1.017	0.896	0.910	0.936	11.1	

* 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.

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Contract: FIRS02

Lab Code: ACE

SDG No.: Q2814

Instrument ID: MSVOA_X

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Heated Purge: (Y/N) N

Calibration Time(s): 09:47 11:32

GC Column: DB-624UI ID: 0.18 (mm)

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		RRF050 = VX047058.D		RRF100 = VX047059.D		RRF150 = VX047060.D			
COMPOUND	RRF001	RRF005	RRF020	RRF050	RRF100	RRF150	RRF	% RSD	
t-1,3-Dichloropropene	0.429	0.549	0.546	0.592	0.548	0.562	0.538	10.4	
cis-1,3-Dichloropropene	0.442	0.611	0.618	0.653	0.593	0.614	0.588	12.6	
1,1,2-Trichloroethane	0.300	0.391	0.373	0.376	0.332	0.335	0.351	9.9	
2-Hexanone	0.192	0.312	0.324	0.343	0.305	0.304	0.297	17.9	
Dibromochloromethane	0.343	0.449	0.428	0.444	0.398	0.409	0.412	9.4	
1,2-Dibromoethane	0.334	0.398	0.377	0.395	0.349	0.351	0.367	7.3	
Tetrachloroethene	0.322	0.403	0.381	0.376	0.337	0.344	0.360	8.6	
Chlorobenzene	1.029	1.295	1.225	1.231	1.114	1.135	1.171	8.2	
Ethyl Benzene	1.814	2.189	2.119	2.204	1.975	1.987	2.048	7.4	
m/p-Xylenes	0.693	0.827	0.795	0.818	0.734	0.732	0.766	7.1	
o-Xylene	0.586	0.820	0.769	0.785	0.704	0.711	0.729	11.4	
Styrene	1.024	1.356	1.332	1.361	1.209	1.219	1.250	10.4	
Bromoform	0.257	0.323	0.302	0.325	0.292	0.299	0.300	8.2	
Isopropylbenzene	3.180	4.148	4.146	4.309	3.907	3.979	3.945	10.2	
1,1,2,2-Tetrachloroethane	0.979	1.203	1.182	1.210	1.085	1.103	1.127	7.9	
1,3-Dichlorobenzene	1.562	1.968	1.869	1.908	1.704	1.776	1.798	8.3	
1,4-Dichlorobenzene	1.659	1.982	1.903	1.893	1.723	1.767	1.821	6.8	
1,2-Dichlorobenzene	1.533	1.873	1.764	1.824	1.625	1.691	1.718	7.4	
1,2-Dibromo-3-Chloropropane	0.140	0.223	0.223	0.239	0.231	0.242	0.216	17.6	
1,2,4-Trichlorobenzene	0.919	1.189	1.159	1.224	1.144	1.176	1.135	9.6	
1,2,3-Trichlorobenzene	0.925	1.120	1.117	1.172	1.102	1.126	1.094	7.9	
1,2-Dichloroethane-d4		0.768	0.511	0.575	0.611	0.649	0.623	15.4	
Dibromofluoromethane		0.355	0.262	0.296	0.302	0.328	0.309	11.3	
Toluene-d8		1.293	0.821	0.924	0.946	1.014	1.000	17.8	
4-Bromofluorobenzene		0.514	0.369	0.409	0.417	0.445	0.431	12.5	

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