

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: Alliance

Contract: NOBI03

Lab Code: ACE

SDG No.: Q2879

Instrument ID: MSVOA_Y

Calibration Date(s): 08/12/2025 08/12/2025

Heated Purge: (Y/N) Y

Calibration Time(s): 14:54 16:47

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

LAB FILE ID:		RRF005 = VY023050.D		RRF010 = VY023051.D		RRF020 = VY023052.D			
		RRF050 = VY023053.D		RRF100 = VY023054.D		RRF150 = VY023055.D			
COMPOUND	RRF005	RRF010	RRF020	RRF050	RRF100	RRF150	RRF	% RSD	
Dichlorodifluoromethane	0.487	0.463	0.460	0.348	0.337	0.354	0.408	16.8	
Chloromethane	0.699	0.707	0.736	0.607	0.603	0.613	0.661	9	
Vinyl Chloride	0.868	0.870	0.875	0.769	0.754	0.766	0.817	7.3	
Bromomethane	0.673	0.638	0.653	0.540	0.576	0.581	0.610	8.5	
Chloroethane	0.585	0.567	0.569	0.515	0.508	0.516	0.543	6.2	
Tetrahydrofuran	0.084	0.091	0.089	0.087	0.081	0.091	0.087	4.6	
Trichlorofluoromethane	1.059	1.044	1.048	0.937	0.900	0.974	0.994	6.7	
1,1,2-Trichlorotrifluoroethane	0.522	0.519	0.521	0.462	0.445	0.472	0.490	7	
1,1-Dichloroethene	0.492	0.505	0.502	0.463	0.454	0.475	0.482	4.4	
Acrylonitrile	0.101	0.114	0.112	0.106	0.100	0.109	0.107	5.5	
Acetone	0.110	0.111	0.106	0.103	0.094	0.096	0.103	6.9	
Carbon Disulfide	1.656	1.663	1.662	1.491	1.453	1.501	1.571	6.3	
Methyl tert-butyl Ether	1.181	1.268	1.296	1.236	1.202	1.308	1.249	4.1	
Methylene Chloride	0.861	0.706	0.606	0.510	0.511	0.524	0.620	22.7	
trans-1,2-Dichloroethene	0.535	0.561	0.559	0.523	0.520	0.550	0.541	3.3	
1,1-Dichloroethane	0.929	0.969	0.984	0.918	0.907	0.945	0.942	3.1	
2-Butanone	0.135	0.152	0.143	0.142	0.129	0.142	0.140	5.5	
Carbon Tetrachloride	0.490	0.486	0.505	0.466	0.466	0.484	0.483	3.1	
2,2-Dichloropropane	0.837	0.830	0.844	0.790	0.776	0.800	0.813	3.4	
cis-1,2-Dichloroethene	0.621	0.624	0.645	0.607	0.610	0.648	0.626	2.8	
Chloroform	0.969	0.997	1.002	0.943	0.936	0.981	0.971	2.8	
1,1,1-Trichloroethane	0.862	0.877	0.899	0.837	0.833	0.864	0.862	2.8	
1,1-Dichloropropene	0.442	0.455	0.461	0.433	0.429	0.447	0.445	2.8	
Benzene	1.354	1.375	1.430	1.334	1.334	1.395	1.370	2.8	
1,2-Dichloroethane	0.353	0.370	0.371	0.347	0.339	0.359	0.356	3.6	
Trichloroethene	0.354	0.355	0.373	0.341	0.342	0.359	0.354	3.3	
1,2-Dichloropropane	0.303	0.321	0.326	0.309	0.309	0.325	0.315	3.1	
Dibromomethane	0.166	0.189	0.187	0.178	0.176	0.190	0.181	5.1	
Bromodichloromethane	0.446	0.464	0.489	0.459	0.456	0.481	0.466	3.5	
4-Methyl-2-Pentanone	0.168	0.201	0.200	0.199	0.190	0.211	0.195	7.7	

* Compounds with required minimum RRF and maximum %RSD values.
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RRF of 1,4-Dioxane = Value should be divide by 1000.

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SDG No.: Q2879

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Heated Purge: (Y/N) Y

Calibration Time(s): 14:54 16:47

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

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COMPOUND	RRF005	RRF010	RRF020	RRF050	RRF100	RRF150	RRF	% RSD	
Toluene	0.817	0.851	0.902	0.858	0.873	0.921	0.870	4.3	
t-1,3-Dichloropropene	0.377	0.410	0.435	0.420	0.428	0.459	0.422	6.5	
cis-1,3-Dichloropropene	0.457	0.486	0.513	0.496	0.500	0.540	0.499	5.5	
1,1,2-Trichloroethane	0.230	0.242	0.253	0.239	0.236	0.254	0.242	4	
1,3-Dichloropropane	0.385	0.421	0.429	0.405	0.395	0.429	0.411	4.5	
2-Hexanone	0.112	0.134	0.137	0.136	0.131	0.146	0.133	8.6	
Dibromochloromethane	0.300	0.317	0.326	0.318	0.316	0.340	0.320	4.1	
1,2-Dibromoethane	0.206	0.232	0.237	0.226	0.222	0.242	0.228	5.6	
Tetrachloroethene	0.477	0.463	0.475	0.435	0.393	0.437	0.446	7.2	
Chlorobenzene	1.062	1.070	1.116	1.055	1.057	1.111	1.079	2.6	
1,1,1,2-Tetrachloroethane	0.347	0.360	0.378	0.362	0.365	0.384	0.366	3.6	
Ethyl Benzene	1.775	1.795	1.894	1.859	1.858	1.956	1.856	3.6	
m/p-Xylenes	0.676	0.699	0.745	0.718	0.737	0.774	0.725	4.8	
o-Xylene	0.605	0.643	0.696	0.686	0.702	0.742	0.679	7.1	
Styrene	0.945	1.023	1.140	1.123	1.176	1.254	1.110	10	
Bromoform	0.197	0.213	0.216	0.214	0.213	0.231	0.214	5	
Isopropylbenzene	3.420	3.483	3.602	3.505	3.514	3.650	3.529	2.4	
1,1,2,2-Tetrachloroethane	0.584	0.647	0.605	0.585	0.581	0.613	0.603	4.2	
1,2,3-Trichloropropane	0.551	0.483	0.472	0.535	0.409	0.444	0.483	11.2	
Bromobenzene	0.824	0.830	0.866	0.828	0.831	0.873	0.842	2.6	
n-propylbenzene	4.052	4.149	4.317	4.217	4.158	4.315	4.201	2.5	
2-Chlorotoluene	2.349	2.398	2.464	2.382	2.390	2.461	2.407	1.9	
1,3,5-Trimethylbenzene	2.672	2.811	2.932	2.859	2.872	2.966	2.852	3.6	
4-Chlorotoluene	2.427	2.486	2.600	2.462	2.460	2.548	2.497	2.6	
tert-Butylbenzene	2.428	2.481	2.626	2.588	2.630	2.702	2.576	4	
1,2,4-Trimethylbenzene	2.596	2.760	2.955	2.837	2.894	2.993	2.839	5.1	
sec-Butylbenzene	3.637	3.749	3.881	3.773	3.781	3.882	3.784	2.4	
p-Isopropyltoluene	2.878	2.978	3.241	3.167	3.230	3.350	3.141	5.7	
1,3-Dichlorobenzene	1.600	1.642	1.701	1.610	1.656	1.752	1.660	3.5	
1,4-Dichlorobenzene	1.687	1.657	1.680	1.601	1.597	1.660	1.647	2.3	

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Lab Code: <u>ACE</u>	SDG No.: <u>Q2879</u>
Instrument ID: <u>MSVOA_Y</u>	Calibration Date(s): <u>08/12/2025</u> <u>08/12/2025</u>
Heated Purge: (Y/N) <u>Y</u>	Calibration Time(s): <u>14:54</u> <u>16:47</u>
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COMPOUND	RRF005	RRF010	RRF020	RRF050	RRF100	RRF150	RRF	% RSD
n-Butylbenzene	2.625	2.809	2.972	2.911	2.933	3.059	2.885	5.2
1,2-Dichlorobenzene	1.410	1.472	1.496	1.420	1.438	1.499	1.456	2.7
1,2-Dibromo-3-Chloropropane	0.093	0.096	0.094	0.095	0.086	0.098	0.094	4.3
1,2,4-Trichlorobenzene	0.788	0.839	0.881	0.847	0.856	0.955	0.861	6.4
Hexachlorobutadiene	0.533	0.506	0.522	0.491	0.480	0.504	0.506	3.8
1,2,3-Trichlorobenzene	0.685	0.733	0.752	0.720	0.727	0.834	0.742	6.7
1,2-Dichloroethane-d4	0.491	0.490	0.482	0.462	0.472	0.462	0.477	2.7
Dibromofluoromethane	0.290	0.299	0.301	0.293	0.307	0.305	0.299	2.1
Toluene-d8	1.151	1.157	1.173	1.143	1.202	1.186	1.169	1.9
4-Bromofluorobenzene	0.358	0.367	0.367	0.370	0.399	0.394	0.376	4.4
trans-1,4-Dichloro-2-butene	0.181	0.215	0.203	0.197	0.190	0.210	0.199	6.4

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