

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name:	<u>Alliance</u>	Contract:	<u>NOBI03</u>
Lab Code:	<u>ACE</u>	SDG No.:	<u>Q2558</u>
Instrument ID:	<u>MSVOA_W</u>	Calibration Date(s):	<u>06/30/2025</u> <u>06/30/2025</u>
Heated Purge: (Y/N)	<u>Y</u>	Calibration Time(s):	<u>09:54</u> <u>12:55</u>
GC Column:	<u>RXI-624</u> ID: <u>0.25</u> (mm)		

LAB FILE ID:		RRF005 = VW031729.D		RRF010 = VW031730.D		RRF020 = VW031731.D		
		RRF050 = VW031732.D		RRF100 = VW031733.D		RRF150 = VW031734.D		
COMPOUND	RRF005	RRF010	RRF020	RRF050	RRF100	RRF150	RRF	% RSD
Dichlorodifluoromethane	0.330	0.371	0.352	0.310	0.329	0.351	0.341	6.3
Chloromethane	0.409	0.410	0.428	0.375	0.405	0.436	0.411	5.1
Vinyl Chloride	0.523	0.534	0.576	0.514	0.545	0.546	0.540	4
Bromomethane	0.429	0.431	0.443	0.396	0.412	0.418	0.421	3.9
Chloroethane	0.370	0.372	0.379	0.347	0.360	0.372	0.367	3.1
Tetrahydrofuran	0.151	0.160	0.156	0.168	0.156	0.158	0.158	3.6
Trichlorofluoromethane	0.509	0.508	0.442	0.469	0.466	0.546	0.490	7.7
1,1,2-Trichlorotrifluoroethane	0.577	0.563	0.569	0.508	0.517	0.530	0.544	5.4
1,1-Dichloroethene	0.602	0.625	0.622	0.556	0.581	0.588	0.596	4.4
Acrylonitrile	0.177	0.186	0.182	0.191	0.181	0.181	0.183	2.7
Acetone	0.238	0.191	0.171	0.163	0.157	0.144	0.177	18.8
Carbon Disulfide	1.571	1.588	1.647	1.536	1.590	1.633	1.594	2.5
Methyl tert-butyl Ether	1.007	1.018	1.038	1.033	0.982	0.988	1.011	2.3
Methylene Chloride	1.029	0.980	0.905	0.692	0.655	0.626	0.814	21.7
trans-1,2-Dichloroethene	0.635	0.623	0.667	0.613	0.631	0.629	0.633	2.9
1,1-Dichloroethane	1.154	1.163	1.219	1.116	1.131	1.153	1.156	3.1
2-Butanone	0.224	0.224	0.221	0.249	0.231	0.237	0.231	4.6
Carbon Tetrachloride	0.477	0.498	0.498	0.461	0.468	0.485	0.481	3.2
2,2-Dichloropropane	0.693	0.683	0.698	0.657	0.643	0.672	0.674	3.2
cis-1,2-Dichloroethene	0.711	0.709	0.754	0.716	0.736	0.740	0.728	2.5
Chloroform	1.230	1.239	1.299	1.204	1.189	1.208	1.228	3.2
1,1,1-Trichloroethane	0.933	0.983	0.971	0.925	0.907	0.959	0.946	3.1
1,1-Dichloropropene	0.477	0.473	0.497	0.456	0.463	0.467	0.472	3
Benzene	1.410	1.394	1.483	1.375	1.349	1.371	1.397	3.4
1,2-Dichloroethane	0.490	0.490	0.493	0.464	0.449	0.445	0.472	4.7
Trichloroethene	0.344	0.346	0.367	0.344	0.337	0.354	0.349	3.1
1,2-Dichloropropane	0.348	0.344	0.353	0.331	0.324	0.325	0.338	3.7
Dibromomethane	0.228	0.217	0.227	0.219	0.206	0.210	0.218	4.1
Bromodichloromethane	0.509	0.512	0.524	0.511	0.502	0.510	0.511	1.4
4-Methyl-2-Pentanone	0.285	0.301	0.295	0.313	0.289	0.287	0.295	3.5

* 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: Alliance

Contract: NOBI03

Lab Code: ACE

SDG No.: Q2558

Instrument ID: MSVOA_W

Calibration Date(s): 06/30/2025 06/30/2025

Heated Purge: (Y/N) Y

Calibration Time(s): 09:54 12:55

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

LAB FILE ID:		RRF005 = VW031729.D		RRF010 = VW031730.D		RRF020 = VW031731.D			
		RRF050 = VW031732.D		RRF100 = VW031733.D		RRF150 = VW031734.D			
COMPOUND	RRF005	RRF010	RRF020	RRF050	RRF100	RRF150	RRF	% RSD	
Toluene	0.882	0.898	0.938	0.876	0.874	0.899	0.894	2.7	
t-1,3-Dichloropropene	0.432	0.460	0.491	0.496	0.497	0.500	0.480	5.7	
cis-1,3-Dichloropropene	0.514	0.521	0.571	0.549	0.547	0.560	0.544	4.1	
1,1,2-Trichloroethane	0.299	0.282	0.299	0.285	0.275	0.273	0.286	4	
1,3-Dichloropropane	0.522	0.524	0.518	0.491	0.477	0.476	0.501	4.5	
2-Hexanone	0.185	0.201	0.195	0.219	0.201	0.199	0.200	5.6	
Dibromochloromethane	0.329	0.332	0.353	0.352	0.336	0.345	0.341	3	
1,2-Dibromoethane	0.288	0.285	0.296	0.290	0.274	0.281	0.286	2.7	
Tetrachloroethene	0.331	0.321	0.324	0.314	0.329	0.344	0.327	3.2	
Chlorobenzene	1.128	1.116	1.116	1.062	1.078	1.120	1.103	2.4	
1,1,1,2-Tetrachloroethane	0.347	0.351	0.345	0.350	0.352	0.365	0.352	2	
Ethyl Benzene	1.860	1.899	1.911	1.885	1.934	1.989	1.913	2.3	
m/p-Xylenes	0.690	0.730	0.751	0.727	0.750	0.764	0.735	3.6	
o-Xylene	0.636	0.654	0.685	0.682	0.705	0.723	0.681	4.7	
Styrene	1.066	1.154	1.195	1.209	1.206	1.232	1.177	5.1	
Bromoform	0.197	0.202	0.203	0.219	0.215	0.223	0.210	5.1	
Isopropylbenzene	3.643	3.549	3.683	3.732	4.080	4.132	3.803	6.4	
1,1,2,2-Tetrachloroethane	0.891	0.839	0.812	0.843	0.829	0.852	0.844	3.2	
1,2,3-Trichloropropane	0.693	0.651	0.611	0.652	0.617	0.652	0.646	4.6	
Bromobenzene	0.883	0.821	0.816	0.781	0.886	0.902	0.848	5.7	
n-propylbenzene	4.370	4.433	4.418	4.463	4.733	4.909	4.554	4.7	
2-Chlorotoluene	2.717	2.648	2.721	2.684	2.800	2.885	2.743	3.1	
1,3,5-Trimethylbenzene	2.947	2.987	3.121	3.155	3.314	3.365	3.148	5.4	
4-Chlorotoluene	2.935	2.802	2.837	2.822	2.901	2.978	2.879	2.4	
tert-Butylbenzene	2.532	2.560	2.642	2.613	2.838	2.995	2.697	6.7	
1,2,4-Trimethylbenzene	3.054	3.078	3.192	3.122	3.292	3.399	3.190	4.2	
sec-Butylbenzene	3.914	3.877	3.937	3.987	4.185	4.386	4.048	4.9	
p-Isopropyltoluene	3.123	3.166	3.292	3.275	3.517	3.643	3.336	6.1	
1,3-Dichlorobenzene	1.771	1.705	1.683	1.644	1.681	1.723	1.701	2.5	
1,4-Dichlorobenzene	1.815	1.750	1.700	1.702	1.735	1.745	1.741	2.4	

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SDG No.: Q2558

Instrument ID: MSVOA_W

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Heated Purge: (Y/N) Y

Calibration Time(s): 09:54 12:55

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

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COMPOUND	RRF005	RRF010	RRF020	RRF050	RRF100	RRF150	RRF	% RSD	
n-Butylbenzene	3.127	3.110	3.134	3.176	3.389	3.511	3.241	5.2	
1,2-Dichlorobenzene	1.599	1.518	1.511	1.575	1.563	1.563	1.555	2.2	
1,2-Dibromo-3-Chloropropane	0.169	0.159	0.153	0.166	0.167	0.174	0.165	4.6	
1,2,4-Trichlorobenzene	0.965	0.899	0.929	0.933	1.000	1.028	0.959	5	
Hexachlorobutadiene	0.474	0.469	0.438	0.414	0.493	0.489	0.463	6.7	
1,2,3-Trichlorobenzene	0.863	0.865	0.827	0.862	0.944	0.926	0.881	5	
1,2-Dichloroethane-d4	0.731	0.732	0.739	0.715	0.702	0.687	0.718	2.8	
Dibromofluoromethane	0.322	0.324	0.348	0.324	0.327	0.312	0.326	3.6	
Toluene-d8	1.121	1.245	1.290	1.208	1.229	1.193	1.214	4.7	
4-Bromofluorobenzene	0.436	0.452	0.463	0.447	0.450	0.434	0.447	2.4	
trans-1,4-Dichloro-2-butene	0.245	0.272	0.249	0.291	0.299	0.309	0.277	9.6	

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