

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: <u>Alliance</u>	Contract: <u>WALS01</u>
Lab Code: <u>ACE</u>	SDG No.: <u>Q2487</u>
Instrument ID: <u>MSVOA_N</u>	Calibration Date(s): <u>06/30/2025</u> <u>06/30/2025</u>
Heated Purge: (Y/N) <u>N</u>	Calibration Time(s): <u>12:28</u> <u>14:51</u>
GC Column: <u>RXI-624</u> ID: <u>0.25</u> (mm)	

LAB FILE ID:		RRF001 = VN087238.D		RRF005 = VN087239.D		RRF020 = VN087240.D			
		RRF050 = VN087241.D		RRF100 = VN087242.D		RRF150 = VN087243.D			
COMPOUND	RRF001	RRF005	RRF020	RRF050	RRF100	RRF150	RRF	% RSD	
Dichlorodifluoromethane	0.442	0.449	0.479	0.507	0.508	0.534	0.487	7.5	
Chloromethane	0.362	0.374	0.412	0.393	0.402	0.428	0.395	6.2	
Vinyl Chloride	0.423	0.492	0.525	0.488	0.498	0.489	0.486	6.9	
Ethyl Acetate	0.421	0.423	0.499	0.465	0.490	0.510	0.468	8.2	
Bromomethane		0.444	0.451	0.402	0.439	0.480	0.443	6.3	
Chloroethane	0.362	0.383	0.424	0.382	0.383	0.374	0.385	5.4	
Trichlorofluoromethane	0.826	0.854	0.936	0.820	0.820	0.802	0.843	5.8	
1,1,2-Trichlorotrifluoroethane	0.453	0.501	0.534	0.479	0.486	0.483	0.489	5.5	
1,1-Dichloroethene	0.423	0.450	0.512	0.464	0.466	0.474	0.465	6.3	
Acrolein		0.071	0.076	0.076	0.083	0.090	0.079	9.2	
Acrylonitrile	0.267	0.278	0.309	0.276	0.286	0.292	0.285	5.2	
Acetone	0.260	0.237	0.263	0.226	0.227	0.226	0.240	7.2	
Carbon Disulfide	1.540	1.386	1.489	1.321	1.320	1.315	1.395	7	
Methyl tert-butyl Ether	1.448	1.548	1.759	1.606	1.677	1.739	1.629	7.3	
Methyl Acetate	0.501	0.542	0.616	0.540	0.544	0.554	0.549	6.8	
Methylene Chloride	0.526	0.552	0.594	0.506	0.508	0.503	0.531	6.7	
trans-1,2-Dichloroethene	0.579	0.552	0.600	0.531	0.550	0.560	0.562	4.3	
1,1-Dichloroethane	0.798	0.897	0.983	0.873	0.896	0.916	0.894	6.7	
Cyclohexane		0.982	0.873	0.783	0.808	0.827	0.855	9.2	
2-Butanone	0.330	0.354	0.406	0.370	0.390	0.400	0.375	7.8	
Carbon Tetrachloride	0.444	0.477	0.530	0.485	0.512	0.512	0.493	6.3	
2,2-Dichloropropane	0.796	0.854	0.927	0.826	0.859	0.871	0.855	5.2	
cis-1,2-Dichloroethene	0.617	0.613	0.702	0.633	0.656	0.672	0.649	5.3	
Bromochloromethane	0.362	0.397	0.384	0.401	0.404	0.420	0.395	5	
Chloroform	0.881	0.988	1.102	0.956	0.996	1.022	0.991	7.4	
1,1,1-Trichloroethane	0.915	0.919	0.994	0.887	0.898	0.924	0.923	4.1	
Methylcyclohexane	0.455	0.463	0.545	0.515	0.561	0.572	0.518	9.7	
1,1-Dichloropropene	0.396	0.399	0.467	0.425	0.459	0.466	0.435	7.6	
Benzene	1.284	1.270	1.434	1.310	1.381	1.414	1.349	5.2	
1,2-Dichloroethane	0.400	0.437	0.470	0.443	0.470	0.471	0.449	6.3	

* 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: WALS01

Lab Code: ACE

SDG No.: Q2487

Instrument ID: MSVOA_N

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

Heated Purge: (Y/N) N

Calibration Time(s): 12:28 14:51

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

LAB FILE ID:		RRF001 = VN087238.D		RRF005 = VN087239.D		RRF020 = VN087240.D			
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COMPOUND	RRF001	RRF005	RRF020	RRF050	RRF100	RRF150	RRF	% RSD	
Trichloroethene	0.322	0.363	0.385	0.346	0.364	0.370	0.358	6.1	
1,2-Dichloropropane	0.247	0.275	0.321	0.290	0.311	0.320	0.294	9.9	
Dibromomethane	0.238	0.234	0.271	0.241	0.255	0.261	0.250	5.9	
Bromodichloromethane	0.408	0.489	0.521	0.465	0.486	0.495	0.477	8.1	
4-Methyl-2-Pentanone	0.313	0.377	0.471	0.454	0.516	0.534	0.444	19.1	
Toluene	0.741	0.812	0.928	0.889	0.956	0.980	0.885	10.4	
t-1,3-Dichloropropene	0.403	0.466	0.536	0.514	0.569	0.583	0.512	13.2	
cis-1,3-Dichloropropene	0.414	0.494	0.557	0.519	0.566	0.579	0.521	11.8	
1,1,2-Trichloroethane	0.298	0.331	0.354	0.325	0.350	0.348	0.334	6.3	
1,3-Dichloropropane	0.492	0.524	0.578	0.527	0.573	0.578	0.545	6.6	
2-Chloroethyl Vinyl ether	0.200	0.234	0.256	0.266	0.290	0.306	0.259	14.8	
2-Hexanone		0.268	0.268	0.307	0.366	0.381	0.318	16.7	
Dibromochloromethane	0.328	0.376	0.420	0.393	0.425	0.420	0.394	9.5	
1,2-Dibromoethane	0.294	0.340	0.388	0.357	0.390	0.388	0.359	10.6	
Tetrachloroethene	0.304	0.326	0.362	0.323	0.342	0.344	0.333	6	
Chlorobenzene	1.107	1.082	1.194	1.071	1.131	1.145	1.122	4	
1,1,1,2-Tetrachloroethane	0.368	0.370	0.415	0.378	0.394	0.399	0.387	4.9	
Hexachloroethane	0.476	0.552	0.548	0.516	0.543	0.556	0.532	5.8	
Ethyl Benzene	1.482	1.659	1.945	1.801	1.968	2.038	1.815	11.7	
m/p-Xylenes	0.585	0.655	0.782	0.734	0.787	0.809	0.725	12.1	
o-Xylene	0.499	0.625	0.722	0.695	0.751	0.778	0.679	15.1	
Styrene	0.803	0.978	1.258	1.203	1.313	1.365	1.153	18.9	
Bromoform	0.217	0.289	0.320	0.308	0.326	0.333	0.299	14.3	
Isopropylbenzene	2.651	2.894	3.493	3.144	3.508	3.640	3.222	12.2	
1,1,2,2-Tetrachloroethane	1.002	1.049	1.137	0.994	1.073	1.104	1.060	5.3	
1,2,3-Trichloropropane	0.859	1.016	1.118	0.851	0.925	0.938	0.951	10.6	
Bromobenzene	0.716	0.823	0.909	0.827	0.878	0.893	0.841	8.4	
2-Chlorotoluene	2.129	2.149	2.449	2.239	2.426	2.504	2.316	7.1	
1,3,5-Trimethylbenzene	2.169	2.469	2.895	2.718	2.931	2.976	2.693	11.8	
4-Chlorotoluene	2.197	2.230	2.573	2.332	2.587	2.628	2.424	8	

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COMPOUND	RRF001	RRF005	RRF020	RRF050	RRF100	RRF150	RRF	% RSD
1,2,4-Trimethylbenzene	2.004	2.381	2.970	2.779	2.989	3.063	2.698	15.5
1,3-Dichlorobenzene	1.572	1.612	1.789	1.611	1.716	1.749	1.675	5.3
1,4-Dichlorobenzene	1.883	1.715	1.869	1.625	1.680	1.721	1.749	6
1,2-Dichlorobenzene	1.444	1.531	1.720	1.508	1.573	1.622	1.566	6.1
1,2-Dibromo-3-Chloropropane	0.270	0.236	0.260	0.237	0.257	0.268	0.255	5.8
1,2,4-Trichlorobenzene	0.853	0.890	0.992	0.894	0.936	0.992	0.926	6.2
Hexachlorobutadiene	0.355	0.281	0.310	0.273	0.263	0.284	0.294	11.5
1,2,3-Trichlorobenzene	0.834	0.838	0.967	0.876	0.917	0.982	0.902	7
1,2-Dichloroethane-d4		0.655	0.598	0.592	0.636	0.656	0.628	4.9
Dibromofluoromethane		0.313	0.303	0.307	0.334	0.345	0.320	5.7
Toluene-d8		1.137	1.150	1.205	1.363	1.384	1.248	9.4
4-Bromofluorobenzene		0.390	0.403	0.432	0.490	0.505	0.444	11.6
Methyl Iodide		0.332	0.466	0.526	0.554	0.550	0.486	19.1
1,4-Dioxane	2.618	5.085	6.447	6.464	6.208	6.526	5.558	27.7

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