

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: <u>Alliance</u>	Contract: <u>GENV01</u>
Lab Code: <u>ACE</u>	SDG No.: <u>Q2480</u>
Instrument ID: <u>MSVOA_X</u>	Calibration Date(s): <u>07/02/2025</u> <u>07/02/2025</u>
Heated Purge: (Y/N) <u>N</u>	Calibration Time(s): <u>12:11</u> <u>14:31</u>
GC Column: <u>DB-624UI</u> ID: <u>0.18</u> (mm)	

LAB FILE ID:		RRF001 = VX046860.D		RRF005 = VX046861.D		RRF020 = VX046862.D			
		RRF050 = VX046863.D		RRF100 = VX046864.D		RRF150 = VX046865.D			
COMPOUND	RRF001	RRF005	RRF020	RRF050	RRF100	RRF150	RRF	% RSD	
Dichlorodifluoromethane	0.445	0.433	0.512	0.487	0.503	0.490	0.478	6.7	
Chloromethane	0.505	0.507	0.541	0.508	0.543	0.530	0.522	3.4	
Vinyl Chloride	0.568	0.549	0.591	0.550	0.589	0.566	0.569	3.2	
Bromomethane		0.402	0.402	0.360	0.369	0.295	0.366	12	
Chloroethane	0.443	0.365	0.383	0.347	0.362	0.354	0.376	9.4	
Trichlorofluoromethane	0.847	0.874	0.921	0.857	0.893	0.885	0.880	3	
1,1,2-Trichlorotrifluoroethane	0.527	0.571	0.583	0.548	0.572	0.562	0.561	3.6	
1,1-Dichloroethene	0.543	0.555	0.546	0.524	0.546	0.540	0.542	1.9	
Acetone	0.236	0.210	0.195	0.193	0.197	0.200	0.205	7.9	
Carbon Disulfide	1.585	1.425	1.405	1.342	1.398	1.376	1.422	6	
Methyl tert-butyl Ether	1.447	1.470	1.540	1.528	1.587	1.615	1.531	4.2	
Methyl Acetate	0.488	0.474	0.560	0.553	0.570	0.599	0.541	9.1	
Methylene Chloride	0.623	0.625	0.624	0.585	0.596	0.588	0.607	3.1	
trans-1,2-Dichloroethene	0.552	0.564	0.575	0.541	0.551	0.545	0.555	2.3	
1,1-Dichloroethane	1.080	1.084	1.074	1.039	1.055	1.050	1.064	1.7	
Cyclohexane		0.973	0.977	0.918	0.926	0.927	0.944	3	
2-Butanone	0.260	0.270	0.275	0.275	0.276	0.286	0.274	3.1	
Carbon Tetrachloride	0.471	0.505	0.495	0.476	0.480	0.478	0.484	2.7	
cis-1,2-Dichloroethene	0.711	0.669	0.688	0.660	0.670	0.666	0.677	2.8	
Bromochloromethane	0.529	0.543	0.524	0.528	0.516	0.508	0.525	2.3	
Chloroform	1.150	1.106	1.112	1.048	1.052	1.050	1.087	3.9	
1,1,1-Trichloroethane	0.944	0.897	0.908	0.886	0.901	0.910	0.908	2.2	
Methylcyclohexane	0.539	0.588	0.589	0.567	0.578	0.576	0.573	3.2	
Benzene	1.391	1.445	1.452	1.356	1.342	1.324	1.385	3.9	
1,2-Dichloroethane	0.499	0.477	0.487	0.460	0.453	0.446	0.470	4.4	
Trichloroethene	0.386	0.377	0.365	0.344	0.347	0.344	0.361	5	
1,2-Dichloropropane	0.364	0.340	0.362	0.346	0.345	0.342	0.350	3	
Bromodichloromethane	0.538	0.508	0.535	0.510	0.508	0.507	0.518	2.9	
4-Methyl-2-Pentanone	0.330	0.340	0.368	0.366	0.353	0.357	0.353	4.3	
Toluene	0.855	0.894	0.906	0.841	0.831	0.823	0.858	3.9	

* 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:	<u>Alliance</u>	Contract:	<u>GENV01</u>
Lab Code:	<u>ACE</u>	SDG No.:	<u>Q2480</u>
Instrument ID:	<u>MSVOA_X</u>	Calibration Date(s):	<u>07/02/2025</u> <u>07/02/2025</u>
Heated Purge: (Y/N)	<u>N</u>	Calibration Time(s):	<u>12:11</u> <u>14:31</u>
GC Column:	<u>DB-624UI</u>	ID:	<u>0.18</u> (mm)

LAB FILE ID:		RRF001 = VX046860.D		RRF005 = VX046861.D		RRF020 = VX046862.D		
		RRF050 = VX046863.D		RRF100 = VX046864.D		RRF150 = VX046865.D		
COMPOUND	RRF001	RRF005	RRF020	RRF050	RRF100	RRF150	RRF	% RSD
t-1,3-Dichloropropene	0.408	0.431	0.467	0.472	0.490	0.504	0.462	7.8
cis-1,3-Dichloropropene	0.496	0.499	0.532	0.533	0.549	0.554	0.527	4.7
1,1,2-Trichloroethane	0.325	0.329	0.328	0.318	0.310	0.308	0.320	2.8
2-Hexanone	0.205	0.227	0.245	0.247	0.238	0.242	0.234	6.8
Dibromochloromethane	0.383	0.390	0.393	0.377	0.376	0.377	0.383	1.9
1,2-Dibromoethane	0.323	0.317	0.332	0.320	0.317	0.317	0.321	1.8
Tetrachloroethene	0.353	0.346	0.345	0.315	0.320	0.318	0.333	5.1
Chlorobenzene	1.111	1.117	1.113	1.049	1.051	1.045	1.081	3.3
Ethyl Benzene	1.810	1.868	1.929	1.830	1.846	1.824	1.851	2.3
m/p-Xylenes	0.702	0.710	0.731	0.688	0.683	0.672	0.698	3.1
o-Xylene	0.676	0.677	0.703	0.665	0.663	0.660	0.674	2.4
Styrene	1.070	1.152	1.228	1.175	1.157	1.134	1.153	4.5
Bromoform	0.269	0.261	0.273	0.269	0.275	0.275	0.270	1.9
Isopropylbenzene	3.267	3.428	3.611	3.601	3.592	3.590	3.515	4
1,1,2,2-Tetrachloroethane	1.000	0.948	0.981	0.963	0.939	0.968	0.966	2.3
1,3-Dichlorobenzene	1.615	1.664	1.673	1.627	1.625	1.619	1.637	1.5
1,4-Dichlorobenzene	1.852	1.776	1.711	1.632	1.622	1.632	1.704	5.5
1,2-Dichlorobenzene	1.626	1.592	1.625	1.568	1.560	1.550	1.587	2.1
1,2-Dibromo-3-Chloropropane	0.158	0.151	0.164	0.171	0.178	0.192	0.169	8.7
1,2,4-Trichlorobenzene	1.016	1.020	1.078	1.087	1.121	1.145	1.078	4.9
1,2,3-Trichlorobenzene	0.900	0.965	1.033	1.045	1.069	1.099	1.019	7.2
1,2-Dichloroethane-d4		0.722	0.652	0.647	0.641	0.640	0.661	5.3
Dibromofluoromethane		0.355	0.346	0.339	0.332	0.325	0.339	3.5
Toluene-d8		1.278	1.220	1.195	1.160	1.135	1.197	4.6
4-Bromofluorobenzene		0.493	0.469	0.455	0.433	0.426	0.455	6

* 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.