

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
Vinyl Chloride	0.423	0.492	0.525	0.488	0.498	0.489	0.486	6.9
1,1-Dichloroethene	0.423	0.450	0.512	0.464	0.466	0.474	0.465	6.3
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
Chloroform	0.881	0.988	1.102	0.956	0.996	1.022	0.991	7.4
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
Trichloroethene	0.322	0.363	0.385	0.346	0.364	0.370	0.358	6.1
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
1,2-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

* 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.