

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: <u>Alliance</u>	Contract: <u>POWE02</u>
Lab Code: <u>ACE</u>	SDG No.: <u>Q3155</u>
Instrument ID: <u>MSVOA_W</u>	Calibration Date(s): <u>08/26/2025</u> <u>08/26/2025</u>
Heated Purge: (Y/N) <u>Y</u>	Calibration Time(s): <u>09:39</u> <u>12:34</u>
GC Column: <u>RXI-624</u> ID: <u>0.25</u> (mm)	

LAB FILE ID:		RRF005 = VW032137.D		RRF010 = VW032138.D		RRF020 = VW032139.D		
		RRF100 = VW032141.D		RRF150 = VW032142.D		RRF050 = VW032144.D		
COMPOUND	RRF005	RRF010	RRF020	RRF100	RRF150	RRF050	RRF	% RSD
cis-1,2-Dichloroethene	0.734	0.721	0.722	0.705	0.689	0.739	0.718	2.6
1,1,1-Trichloroethane	0.938	0.907	0.901	0.847	0.839	0.917	0.891	4.5
Benzene	1.381	1.421	1.411	1.363	1.334	1.525	1.406	4.7
Trichloroethene	0.350	0.359	0.352	0.343	0.343	0.374	0.354	3.3
Toluene	0.890	0.920	0.912	0.902	0.879	1.010	0.919	5.1
Ethyl Benzene	1.676	1.823	1.920	1.880	1.862	2.027	1.865	6.2
m/p-Xylenes	0.627	0.700	0.751	0.730	0.699	0.798	0.718	8
o-Xylene	0.581	0.659	0.695	0.688	0.699	0.752	0.679	8.4
Isopropylbenzene	2.833	3.266	3.258	3.570	3.666	3.788	3.397	10.3
1,2-Dichloroethane-d4	0.752	0.715	0.723	0.690	0.713	0.697	0.715	3.1
Dibromofluoromethane	0.367	0.350	0.347	0.335	0.346	0.360	0.351	3.2
Toluene-d8	1.173	1.215	1.235	1.191	1.253	1.322	1.231	4.3
4-Bromofluorobenzene	0.448	0.431	0.450	0.447	0.468	0.489	0.455	4.5

* 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.