

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: <u>Alliance</u>	Contract: <u>TULL01</u>
Lab Code: <u>ACE</u>	SDG No.: <u>Q2730</u>
Instrument ID: <u>MSVOA_N</u>	Calibration Date(s): <u>07/28/2025</u> <u>07/28/2025</u>
Heated Purge: (Y/N) <u>N</u>	Calibration Time(s): <u>09:06</u> <u>11:09</u>
GC Column: <u>RXI-624</u> ID: <u>0.25</u> (mm)	

LAB FILE ID:		RRF005 = VN087419.D		RRF020 = VN087420.D		RRF050 = VN087421.D	
		RRF100 = VN087422.D		RRF150 = VN087423.D		RRF =	
COMPOUND	RRF005	RRF020	RRF050	RRF100	RRF150	RRF	% RSD
Benzene	1.354	1.412	1.148	1.483	1.569	1.393	11.4
Toluene	1.674	1.685	1.419	1.814	1.873	1.693	10.3
Ethyl Benzene	1.787	1.887	1.561	1.979	2.036	1.850	10.1
m/p-Xylenes	0.687	0.704	0.591	0.736	0.747	0.693	9
o-Xylene	0.651	0.704	0.570	0.730	0.736	0.678	10.2
1,2-Dichloroethane-d4	2.246	2.272	2.251	2.339	2.362	2.294	2.3
Toluene-d8	1.368	1.381	1.376	1.393	1.371	1.378	0.7
4-Bromofluorobenzene	0.484	0.493	0.486	0.505	0.495	0.493	1.7

* 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.