

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: <u>Alliance</u>	Contract: <u>TULL01</u>
Lab Code: <u>ACE</u>	SDG No.: <u>Q2845</u>
Instrument ID: <u>MSVOA_N</u>	Calibration Date(s): <u>08/05/2025</u> <u>08/05/2025</u>
Heated Purge: (Y/N) <u>N</u>	Calibration Time(s): <u>13:37</u> <u>15:01</u>
GC Column: <u>RXI-624</u> ID: <u>0.25</u> (mm)	

LAB FILE ID:		RRF005 = VN087452.D		RRF020 = VN087453.D		RRF050 = VN087454.D		
		RRF100 = VN087455.D		RRF150 = VN087456.D		RRF =		
COMPOUND	RRF005	RRF020	RRF050	RRF100	RRF150	RRF	RRF	% RSD
Benzene	1.608	1.540	1.410	1.580	1.584		1.545	5.1
Toluene	1.975	1.905	1.689	1.854	1.835		1.852	5.7
Ethyl Benzene	2.154	2.115	1.892	2.094	2.064		2.064	4.9
m/p-Xylenes	0.796	0.768	0.702	0.768	0.753		0.758	4.6
o-Xylene	0.791	0.754	0.681	0.756	0.745		0.745	5.4
1,2-Dichloroethane-d4	2.772	2.838	2.776	2.704	2.746		2.767	1.8
Toluene-d8	1.413	1.382	1.386	1.386	1.336		1.380	2
4-Bromofluorobenzene	0.512	0.520	0.522	0.520	0.508		0.516	1.1

* 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.