

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name:	<u>Alliance</u>	Contract:	<u>TULL01</u>
Lab Code:	<u>ACE</u>	SDG No.:	<u>Q3760</u>
Instrument ID:	<u>MSVOA_N</u>	Calibration Date(s):	<u>11/13/2025</u> <u>11/13/2025</u>
Heated Purge: (Y/N)	<u>N</u>	Calibration Time(s):	<u>10:14</u> <u>11:50</u>
GC Column:	<u>RXI-624</u>	ID:	<u>0.25</u> (mm)

LAB FILE ID:		RRF005 = VN088274.D		RRF020 = VN088275.D		RRF050 = VN088276.D	
		RRF100 = VN088277.D		RRF150 = VN088278.D		RRF =	
COMPOUND	RRF005	RRF020	RRF050	RRF100	RRF150	RRF	% RSD
Benzene	1.560	1.495	1.392	1.499	1.535	1.496	4.3
Toluene	1.716	1.777	1.619	1.714	1.782	1.722	3.8
Ethyl Benzene	1.763	1.927	1.813	1.933	2.052	1.897	6
m/p-Xylenes	0.660	0.723	0.687	0.718	0.760	0.710	5.4
o-Xylene	0.619	0.665	0.643	0.698	0.721	0.669	6.2
1,2-Dichloroethane-d4	2.725	2.740	2.769	2.745	2.723	2.740	0.7
Toluene-d8	1.418	1.455	1.418	1.385	1.426	1.420	1.8
4-Bromofluorobenzene	0.479	0.494	0.513	0.515	0.534	0.507	4.1

* 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.