

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: <u>Alliance</u>	Contract: <u>TULL01</u>
Lab Code: <u>ACE</u>	SDG No.: <u>Q3281</u>
Instrument ID: <u>MSVOA_N</u>	Calibration Date(s): <u>10/08/2025</u> <u>10/08/2025</u>
Heated Purge: (Y/N) <u>N</u>	Calibration Time(s): <u>09:34</u> <u>11:12</u>
GC Column: <u>RXI-624</u> ID: <u>0.25</u> (mm)	

LAB FILE ID:		RRF005 = VN088030.D		RRF020 = VN088031.D		RRF050 = VN088032.D		
		RRF100 = VN088033.D		RRF150 = VN088034.D		RRF =		
COMPOUND	RRF005	RRF020	RRF050	RRF100	RRF150	RRF	RRF	% RSD
Benzene	1.491	1.402	1.366	1.489	1.384		1.426	4.2
Toluene	1.796	1.700	1.645	1.730	1.638		1.702	3.8
Ethyl Benzene	1.902	1.839	1.800	1.919	1.792		1.850	3.1
m/p-Xylenes	0.729	0.714	0.693	0.739	0.693		0.713	2.9
o-Xylene	0.746	0.710	0.685	0.732	0.677		0.710	4.2
1,2-Dichloroethane-d4	2.303	2.319	2.360	2.378	2.319		2.336	1.4
Toluene-d8	1.344	1.343	1.333	1.329	1.334		1.337	0.5
4-Bromofluorobenzene	0.488	0.497	0.493	0.507	0.489		0.495	1.5

* 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.