

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: TULL01
 Lab Code: CHEM Case No.: Q1083 SAS No.: Q1083 SDG No.: Q1083
 Instrument ID: MSVOA_X Calibration Date(s): 01/16/2025 01/16/2025
 Heated Purge: (Y/N) N Calibration Time(s): 08:39 10:12
 GC Column: DB-624UI ID: 0.18 (mm)

LAB FILE ID:		RRF005 = VX044659.D		RRF020 = VX044660.D		RRF050 = VX044661.D			
		RRF100 = VX044662.D		RRF150 = VX044663.D		RRF =			
COMPOUND	RRF005	RRF020	RRF050	RRF100	RRF150	RRF	RRF	% RSD	
Benzene	1.720	1.726	1.635	1.681	1.562		1.665	4.1	
Toluene	2.097	1.972	1.935	2.023	1.850		1.975	4.7	
Ethyl Benzene	2.213	2.181	2.126	2.228	2.031		2.156	3.7	
m/p-Xylenes	0.838	0.817	0.780	0.808	0.719		0.792	5.8	
o-Xylene	0.841	0.823	0.787	0.800	0.726		0.795	5.5	
1,2-Dichloroethane-d4	2.935	2.867	2.905	2.916	2.857		2.896	1.1	
Toluene-d8	1.674	1.620	1.622	1.634	1.609		1.632	1.5	
4-Bromofluorobenzene	0.579	0.578	0.588	0.586	0.582		0.583	0.7	

* 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.