

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: CAMP02
 Lab Code: CHEM Case No.: Q1914 SAS No.: Q1914 SDG No.: Q1914
 Instrument ID: MSVOA_Y Calibration Date(s): 04/22/2025 04/22/2025
 Heated Purge: (Y/N) Y Calibration Time(s): 13:39 16:15
 GC Column: RXI-624 ID: 0.25 (mm)

LAB FILE ID:		RRF005 = VY021953.D		RRF010 = VY021954.D		RRF020 = VY021955.D			
		RRF050 = VY021956.D		RRF100 = VY021957.D		RRF150 = VY021958.D			
COMPOUND	RRF005	RRF010	RRF020	RRF050	RRF100	RRF150	RRF	% RSD	
Benzene	1.423	1.439	1.351	1.337	1.415	1.358	1.387	3.1	
Toluene	0.838	0.915	0.875	0.894	0.952	0.913	0.898	4.4	
Ethyl Benzene	1.693	1.840	1.718	1.835	1.990	1.898	1.829	6.1	
m/p-Xylenes	0.664	0.720	0.699	0.734	0.797	0.754	0.728	6.3	
o-Xylene	0.599	0.639	0.635	0.689	0.747	0.716	0.671	8.3	
Isopropylbenzene	3.061	3.316	3.102	3.325	3.639	3.599	3.341	7.2	
n-propylbenzene	3.708	3.996	3.778	3.977	4.258	4.152	3.978	5.3	
1,3,5-Trimethylbenzene	2.467	2.757	2.614	2.800	2.978	2.898	2.752	6.8	
tert-Butylbenzene	2.246	2.382	2.263	2.494	2.725	2.701	2.469	8.5	
1,2,4-Trimethylbenzene	2.415	2.717	2.631	2.784	3.027	2.945	2.753	8	
sec-Butylbenzene	3.254	3.724	3.409	3.602	3.898	3.786	3.612	6.7	
p-Isopropyltoluene	2.717	3.016	2.869	3.094	3.386	3.321	3.067	8.4	
n-Butylbenzene	2.477	2.790	2.545	2.742	2.976	2.919	2.742	7.2	
1,2-Dichloroethane-d4	0.525	0.477	0.459	0.466	0.418	0.427	0.462	8.3	
Dibromofluoromethane	0.335	0.341	0.330	0.330	0.319	0.327	0.330	2.3	
Toluene-d8	1.220	1.260	1.211	1.276	1.244	1.261	1.245	2	
4-Bromofluorobenzene	0.408	0.429	0.401	0.426	0.423	0.428	0.419	2.8	

* 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.