

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: RUTW01
 Lab Code: CHEM Case No.: Q1232 SAS No.: Q1232 SDG No.: Q1232
 Instrument ID: MSVOA_X Calibration Date(s): 01/15/2025 01/15/2025
 Heated Purge: (Y/N) N Calibration Time(s): 10:31 13:37
 GC Column: DB-624UI ID: 0.18 (mm)

LAB FILE ID:		RRF001 = VX044648.D		RRF005 = VX044649.D		RRF020 = VX044650.D			
		RRF050 = VX044651.D		RRF100 = VX044652.D		RRF150 = VX044653.D			
COMPOUND	RRF001	RRF005	RRF020	RRF050	RRF100	RRF150	RRF	% RSD	
Vinyl Chloride	0.801	0.681	0.681	0.702	0.699	0.749	0.719	6.6	
1,1-Dichloroethene	0.638	0.546	0.572	0.566	0.571	0.618	0.585	6	
2-Butanone	0.451	0.444	0.444	0.426	0.449	0.450	0.444	2.1	
Carbon Tetrachloride	0.585	0.503	0.468	0.460	0.465	0.502	0.497	9.4	
Chloroform	1.261	1.226	1.185	1.142	1.095	1.206	1.186	5	
Benzene	1.517	1.376	1.364	1.306	1.274	1.384	1.370	6.1	
1,2-Dichloroethane	0.552	0.494	0.491	0.477	0.473	0.513	0.500	5.8	
Trichloroethene	0.338	0.319	0.324	0.316	0.324	0.355	0.329	4.5	
Tetrachloroethene	0.346	0.309	0.299	0.296	0.308	0.326	0.314	5.9	
Chlorobenzene	1.178	1.085	1.036	1.008	0.998	1.067	1.062	6.2	
1,2-Dichloroethane-d4		0.863	0.691	0.745	0.689	0.785	0.755	9.6	
Dibromofluoromethane		0.355	0.289	0.318	0.293	0.333	0.318	8.8	
Toluene-d8		1.270	1.051	1.111	1.001	1.104	1.108	9.1	
4-Bromofluorobenzene		0.445	0.394	0.421	0.383	0.421	0.413	6	

* 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.