

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: PSEG03
 Lab Code: CHEM Case No.: P2687 SAS No.: P2687 SDG No.: P2687
 Instrument ID: MSVOA_X Calibration Date(s): 05/29/2024 05/29/2024
 Heated Purge: (Y/N) N Calibration Time(s): 09:34 11:29
 GC Column: DB-624UI ID: 0.18 (mm)

LAB FILE ID:		RRF001 = VX041662.D		RRF005 = VX041663.D		RRF020 = VX041664.D			
		RRF050 = VX041665.D		RRF100 = VX041666.D		RRF150 = VX041667.D			
COMPOUND	RRF001	RRF005	RRF020	RRF050	RRF100	RRF150	RRF	% RSD	
Vinyl Chloride	0.490	0.547	0.484	0.481	0.487	0.473	0.493	5.4	
1,1-Dichloroethene	0.437	0.473	0.430	0.439	0.451	0.445	0.446	3.4	
2-Butanone	0.339	0.404	0.391	0.388	0.390	0.376	0.381	5.9	
Carbon Tetrachloride	0.434	0.469	0.470	0.482	0.500	0.498	0.476	5.1	
Chloroform	0.818	0.953	0.896	0.890	0.898	0.883	0.890	4.9	
Benzene	1.218	1.388	1.320	1.311	1.317	1.296	1.308	4.2	
1,2-Dichloroethane	0.443	0.534	0.493	0.492	0.488	0.480	0.488	6	
Trichloroethene	0.324	0.370	0.352	0.358	0.366	0.365	0.356	4.7	
Tetrachloroethene	0.359	0.396	0.381	0.366	0.379	0.369	0.375	3.5	
Chlorobenzene	1.040	1.133	1.068	1.058	1.084	1.056	1.073	3.1	
1,2-Dichloroethane-d4		0.743	0.555	0.601	0.586	0.580	0.613	12.1	
Dibromofluoromethane		0.397	0.316	0.339	0.340	0.337	0.346	8.7	
Toluene-d8		1.310	1.121	1.215	1.202	1.191	1.208	5.6	
4-Bromofluorobenzene		0.443	0.408	0.454	0.466	0.458	0.446	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.