

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: POWE02
 Lab Code: CHEM Case No.: P4852 SAS No.: P4852 SDG No.: P4852
 Instrument ID: MSVOA_D Calibration Date(s): 11/15/2024 11/15/2024
 Heated Purge: (Y/N) Y Calibration Time(s): 10:04 12:18
 GC Column: RTX-VMS ID: 0.18 (mm)

LAB FILE ID:		RRF005 = VD080020.D		RRF010 = VD080021.D		RRF020 = VD080022.D			
		RRF050 = VD080023.D		RRF100 = VD080024.D		RRF150 = VD080025.D			
COMPOUND	RRF005	RRF010	RRF020	RRF050	RRF100	RRF150	RRF	% RSD	
cis-1,2-Dichloroethene	0.665	0.649	0.620	0.656	0.675	0.679	0.657	3.2	
1,1,1-Trichloroethane	0.970	0.930	0.851	0.869	0.864	0.867	0.892	5.3	
Benzene	1.391	1.430	1.351	1.436	1.439	1.463	1.418	2.8	
Trichloroethene	0.378	0.389	0.341	0.361	0.361	0.372	0.367	4.5	
Toluene	0.814	0.864	0.823	0.917	0.927	0.945	0.881	6.4	
Ethyl Benzene	1.740	1.714	1.693	1.866	1.965	2.002	1.830	7.3	
m/p-Xylenes	0.647	0.688	0.683	0.727	0.766	0.780	0.715	7.2	
o-Xylene	0.542	0.595	0.604	0.670	0.720	0.726	0.643	11.5	
Isopropylbenzene	3.045	3.093	2.974	3.271	3.463	3.501	3.224	6.9	
1,2-Dichloroethane-d4	0.554	0.532	0.451	0.526	0.504	0.523	0.515	6.8	
Dibromofluoromethane	0.335	0.357	0.285	0.338	0.329	0.354	0.333	7.8	
Toluene-d8	1.162	1.249	1.087	1.330	1.303	1.393	1.254	9	
4-Bromofluorobenzene	0.388	0.392	0.353	0.435	0.437	0.464	0.411	9.9	

* 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.