

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: PSEG03
 Lab Code: CHEM Case No.: Q1214 SAS No.: Q1214 SDG No.: Q1214
 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								% RSD
Chloromethane		2.821	2.891	2.775	2.910	2.632		2.806		4					
Vinyl Chloride		2.718	2.750	2.650	2.953	2.656		2.745		4.5					
Bromomethane		0.720	0.648	0.619	0.683	0.647		0.663		5.9					
Chloroethane		0.337	0.431	0.443	0.506	0.459		0.435		14.2					
1,1-Dichloroethene		2.310	2.167	2.109	2.359	2.134		2.216		5					
Acrolein		0.481	0.295	0.325	0.353	0.335		0.358		20.1					
Acrylonitrile		1.225	1.195	1.163	1.234	1.078		1.179		5.3					
Methylene Chloride		2.516	2.531	2.426	2.551	2.375		2.480		3.1					
trans-1,2-Dichloroethene		2.273	2.202	2.143	2.306	2.094		2.203		4					
1,1-Dichloroethane		4.238	4.335	4.228	4.546	4.166		4.303		3.5					
Carbon Tetrachloride		0.616	0.587	0.558	0.605	0.551		0.583		4.8					
cis-1,2-Dichloroethene		2.682	2.813	2.656	2.863	2.622		2.727		3.8					
Chloroform		4.285	4.354	4.180	4.469	4.089		4.275		3.5					
1,1,1-Trichloroethane		0.729	0.702	0.665	0.708	0.654		0.692		4.5					
Benzene		1.720	1.726	1.635	1.681	1.562		1.665		4.1					
1,2-Dichloroethane		0.610	0.616	0.579	0.607	0.563		0.595		3.8					
Trichloroethene		0.394	0.405	0.394	0.424	0.393		0.402		3.3					
Bromodichloromethane		0.667	0.655	0.628	0.638	0.598		0.637		4.1					
Toluene		2.097	1.972	1.935	2.023	1.850		1.975		4.7					
cis-1,3-Dichloropropene		0.737	0.748	0.719	0.731	0.671		0.721		4.2					
1,1,2-Trichloroethane		0.425	0.419	0.398	0.391	0.355		0.397		7					
2-Chloroethyl vinyl ether		0.376	0.351	0.338	0.309	0.302		0.335		9.1					
Dibromochloromethane		0.507	0.503	0.469	0.468	0.428		0.475		6.8					
Tetrachloroethene		0.365	0.350	0.349	0.387	0.348		0.360		4.7					
Chlorobenzene		1.275	1.230	1.215	1.256	1.157		1.227		3.7					
Ethyl Benzene		2.213	2.181	2.126	2.228	2.031		2.156		3.7					
Bromoform		0.340	0.324	0.314	0.310	0.287		0.315		6.2					
1,1,2,2-Tetrachloroethane		0.749	0.684	0.650	0.671	0.607		0.672		7.7					
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					

* 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.

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COMPOUND	RRF005	RRF020	RRF050	RRF100	RRF150	RRF	% RSD
4-Bromofluorobenzene	0.579	0.578	0.588	0.586	0.582	0.583	0.7

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