

VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: CHEMTECH Contract: WOOD06

Lab Code: CHEM Case No.: Q1665 SAS No.: Q1665 SDG No.: Q1665

Instrument ID: MSVOA_L Calibration Date/Time: 04/28/2025 08:44

Lab File ID: VL042416.D Init. Calib. Date(s): 04/17/2025 04/17/2025

Heated Purge: (Y/N) N Init. Calib. Time(s): 09:44 12:54

GC Column: RTX-1 ID: 0.32 (mm)

COMPOUND	RRF	RRF010	MIN RRF	%D	MAX%D
Dichlorodifluoromethane	1.767	1.796		1.64	30
Ethanol	0.122	0.038		-68.85	30
Chloromethane	0.699	0.698		-0.14	30
Vinyl Chloride	0.761	0.679		-10.77	30
Bromomethane	0.327	0.307		-6.12	30
Chloroethane	0.275	0.264		-4	30
Tetrahydrofuran	0.461	0.386		-16.27	30
Trichlorofluoromethane	1.680	1.728		2.86	30
1,1,2-Trichlorotrifluoroethane	1.374	1.337		-2.69	30
Dichlorotetrafluoroethane	1.517	1.551		2.24	30
Bromoethene	0.518	0.518		0	30
tert-Butyl alcohol	2.029	1.859		-8.38	30
Heptane	2.153	1.687		-21.64	30
Isopropyl Alcohol	0.909	0.828		-8.91	30
1,1-Dichloroethene	0.642	0.605		-5.76	30
Acetone	1.637	1.412		-13.74	30
Carbon Disulfide	1.749	1.685		-3.66	30
Methyl tert-Butyl Ether	1.070	0.955		-10.75	30
Methylene Chloride	0.641	0.512		-20.13	30
trans-1,2-Dichloroethene	0.731	0.669		-8.48	30
1,1-Dichloroethane	1.388	1.337		-3.67	30
Cyclohexane	1.320	1.168		-11.52	30
2-Butanone	0.791	0.699		-11.63	30
Carbon Tetrachloride	0.905	0.885		-2.21	30
cis-1,2-Dichloroethene	1.553	1.488		-4.18	30
Chloroform	2.309	2.303		-0.26	30
1,1,1-Trichloroethane	2.599	2.470		-4.96	30
2,2,4-Trimethylpentane	1.863	1.630		-12.51	30
Benzene	1.079	0.982		-8.99	30
1,2-Dichloroethane	0.647	0.684		5.72	30
Trichloroethene	0.471	0.396		-15.92	30
1,2-Dichloropropane	0.402	0.354		-11.94	30
Bromodichloromethane	0.901	0.901		0	30
4-Methyl-2-Pentanone	1.314	1.032		-21.46	30
Toluene	1.294	1.154		-10.82	30
t-1,3-Dichloropropene	0.497	0.456		-8.25	30
cis-1,3-Dichloropropene	0.608	0.554		-8.88	30
1,1,2-Trichloroethane	0.410	0.388		-5.37	30

All other compounds must meet a minimum RRF of 0.010.
RRF of 1,4-Dioxane = Value should be divide by 1000.

VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: CHEMTECH Contract: WOOD06

Lab Code: CHEM Case No.: Q1665 SAS No.: Q1665 SDG No.: Q1665

Instrument ID: MSVOA_L Calibration Date/Time: 04/28/2025 08:44

Lab File ID: VL042416.D Init. Calib. Date(s): 04/17/2025 04/17/2025

Heated Purge: (Y/N) N Init. Calib. Time(s): 09:44 12:54

GC Column: RTX-1 ID: 0.32 (mm)

COMPOUND	RRF	RRF010	MIN RRF	%D	MAX%D
Dibromochloromethane	0.723	0.711		-1.66	30
1,2-Dibromoethane	0.586	0.550		-6.14	30
Tetrachloroethene	0.417	0.331		-20.62	30
Chlorobenzene	1.040	0.985		-5.29	30
Ethyl Benzene	1.981	1.861		-6.06	30
m/p-Xylene	1.596	1.522		-4.64	30
o-Xylene	1.607	1.528		-4.92	30
Styrene	0.767	0.698		-9	30
Bromoform	0.633	0.577		-8.85	30
1,1,2,2-Tetrachloroethane	1.042	0.900		-13.63	30
2-Chlorotoluene	1.851	1.752		-5.35	30
1,3,5-Trimethylbenzene	1.735	1.596		-8.01	30
1,2,4-Trimethylbenzene	2.026	1.775		-12.39	30
1,3-Dichlorobenzene	1.053	0.941		-10.64	30
1,4-Dichlorobenzene	1.059	0.952		-10.1	30
1,2-Dichlorobenzene	1.024	0.917		-10.45	30
1,2,4-Trichlorobenzene	1.040	0.884		-15	30
Hexachloro-1,3-Butadiene	1.034	0.871		-15.76	30
1,3-Butadiene	0.783	0.695		-11.24	30
Naphthalene	2.423	2.040		-15.81	30
4-Ethyltoluene	1.965	1.823		-7.23	30
1-Bromo-4-Fluorobenzene	0.822	0.866		5.35	30
Hexane	1.633	1.367		-16.29	30
Allyl Chloride	1.130	1.026		-9.2	30
1,4-Dioxane	220.893	170.527		-22.8	30
Methyl Methacrylate	0.486	0.400		-17.69	30

All other compounds must meet a minimum RRF of 0.010.
RRF of 1,4-Dioxane = Value should be divide by 1000.