

VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: CHEMTECH Contract: TETR16

Lab Code: CHEM Case No.: P5293 SAS No.: P5293 SDG No.: P5293

Instrument ID: MSVOA_L Calibration Date/Time: 12/18/2024 10:29

Lab File ID: VL041788.D Init. Calib. Date(s): 12/16/2024 12/16/2024

Heated Purge: (Y/N) N Init. Calib. Time(s): 13:30 16:41

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

COMPOUND	RRF	RRF010	MIN RRF	%D	MAX%D
Dichlorodifluoromethane	1.779	1.484		-16.58	30
Chloromethane	0.696	0.637		-8.48	30
Vinyl Chloride	0.701	0.613		-12.55	30
Bromomethane	0.331	0.292		-11.78	30
Chloroethane	0.272	0.253		-6.99	30
Tetrahydrofuran	0.451	0.377		-16.41	30
Trichlorofluoromethane	1.890	1.767		-6.51	30
1,1,2-Trichlorotrifluoroethane	1.438	1.269		-11.75	30
Dichlorotetrafluoroethane	1.593	1.468		-7.85	30
tert-Butyl alcohol	1.930	1.782		-7.67	30
Heptane	1.816	1.468		-19.16	30
1,1-Dichloroethene	0.603	0.542		-10.12	30
Acetone	1.946	1.464		-24.77	30
Carbon Disulfide	1.238	1.122		-9.37	30
Methyl tert-Butyl Ether	1.114	0.894		-19.75	30
Methylene Chloride	0.555	0.469		-15.49	30
trans-1,2-Dichloroethene	0.690	0.596		-13.62	30
1,1-Dichloroethane	1.396	1.260		-9.74	30
Cyclohexane	1.169	0.978		-16.34	30
2-Butanone	0.805	0.703		-12.67	30
Carbon Tetrachloride	0.710	0.745		4.93	30
cis-1,2-Dichloroethene	1.419	1.216		-14.31	30
Chloroform	2.093	1.914		-8.55	30
1,1,1-Trichloroethane	2.128	1.894		-11	30
2,2,4-Trimethylpentane	1.767	1.524		-13.75	30
Benzene	1.034	0.898		-13.15	30
1,2-Dichloroethane	0.644	0.612		-4.97	30
Trichloroethene	0.439	0.354		-19.36	30
1,2-Dichloropropane	0.374	0.323		-13.64	30
Bromodichloromethane	0.740	0.710		-4.05	30
4-Methyl-2-Pentanone	1.095	0.942		-13.97	30
Toluene	1.173	0.984		-16.11	30
t-1,3-Dichloropropene	0.408	0.351		-13.97	30
cis-1,3-Dichloropropene	0.518	0.450		-13.13	30
1,1,2-Trichloroethane	0.391	0.341		-12.79	30
Dibromochloromethane	0.534	0.514		-3.74	30
1,2-Dibromoethane	0.538	0.475		-11.71	30
Tetrachloroethene	0.385	0.301		-21.82	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: TETR16

Lab Code: CHEM Case No.: P5293 SAS No.: P5293 SDG No.: P5293

Instrument ID: MSVOA_L Calibration Date/Time: 12/18/2024 10:29

Lab File ID: VL041788.D Init. Calib. Date(s): 12/16/2024 12/16/2024

Heated Purge: (Y/N) N Init. Calib. Time(s): 13:30 16:41

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

COMPOUND	RRF	RRF010	MIN RRF	%D	MAX%D
Chlorobenzene	1.023	0.912		-10.85	30
Ethyl Benzene	1.800	1.634		-9.22	30
m/p-Xylene	1.448	1.355		-6.42	30
o-Xylene	1.452	1.348		-7.16	30
Styrene	0.628	0.599		-4.62	30
Bromoform	0.437	0.422		-3.43	30
1,1,2,2-Tetrachloroethane	0.909	0.837		-7.92	30
2-Chlorotoluene	1.665	1.538		-7.63	30
1,3,5-Trimethylbenzene	1.422	1.410		-0.84	30
1,2,4-Trimethylbenzene	1.529	1.529		0	30
1,3-Dichlorobenzene	0.975	0.899		-7.8	30
1,4-Dichlorobenzene	0.963	0.895		-7.06	30
1,2-Dichlorobenzene	0.938	0.844		-10.02	30
1,2,4-Trichlorobenzene	0.582	0.544		-6.53	30
Hexachloro-1,3-Butadiene	0.695	0.543		-21.87	30
1,3-Butadiene	0.766	0.701		-8.49	30
Naphthalene	1.065	1.163		9.2	30
4-Ethyltoluene	1.713	1.658		-3.21	30
1-Bromo-4-Fluorobenzene	0.833	0.875		5.04	30
Hexane	1.374	1.152		-16.16	30
Allyl Chloride	1.055	0.979		-7.2	30
1,4-Dioxane	182.757	152.553		-16.53	30
Methyl Methacrylate	0.394	0.351		-10.91	30

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