

VOLATILE CONTINUING CALIBRATION CHECK

Lab Name:	<u>Alliance</u>	Contract:	<u>GENV01</u>
Lab Code:	<u>ACE</u>	SDG No.:	<u>Q3172</u>
Instrument ID:	<u>MSVOA_V</u>	Calibration Date/Time:	<u>09/25/2025 09:59</u>
Lab File ID:	<u>VV039168.D</u>	Init. Calib. Date(s):	<u>09/22/2025 09/22/2025</u>
Heated Purge: (Y/N)	<u>N</u>	Init. Calib. Time(s):	<u>12:26 16:35</u>
GC Column:	<u>DB-624UI</u>	ID:	<u>0.18</u> (mm)

COMPOUND	RRF	RRF050	MIN RRF	%D	MAX%D
Dichlorodifluoromethane	0.792	0.682		-13.89	20
Chloromethane	0.850	0.729	0.1	-14.23	20
Vinyl Chloride	0.910	0.819		-10	20
Bromomethane	0.573	0.458		-20.07	20
Chloroethane	0.631	0.507		-19.65	20
Trichlorofluoromethane	1.317	1.199		-8.96	20
1,1,2-Trichlorotrifluoroethane	0.784	0.710		-9.44	20
Tert butyl alcohol	0.153	0.153		0	20
1,1-Dichloroethene	0.754	0.698		-7.43	20
Acetone	0.523	0.406		-22.37	20
Carbon Disulfide	2.306	2.004		-13.1	20
Methyl tert-butyl Ether	2.402	2.247		-6.45	20
Methyl Acetate	1.072	0.995		-7.18	20
Methylene Chloride	1.066	0.788		-26.08	20
trans-1,2-Dichloroethene	0.803	0.724		-9.84	20
1,1-Dichloroethane	1.576	1.367	0.1	-13.26	20
Cyclohexane	1.291	1.147		-11.15	20
2-Butanone	0.579	0.544		-6.05	20
Carbon Tetrachloride	0.743	0.689		-7.27	20
cis-1,2-Dichloroethene	0.885	0.825		-6.78	20
Bromochloromethane	0.696	0.612		-12.07	20
Chloroform	1.645	1.426		-13.31	20
1,1,1-Trichloroethane	1.424	1.255		-11.87	20
Methylcyclohexane	0.734	0.797		8.58	20
Benzene	1.979	1.926		-2.68	20
1,2-Dichloroethane	0.693	0.639		-7.79	20
Trichloroethene	0.457	0.463		1.31	20
1,2-Dichloropropane	0.492	0.493		0.2	20
Bromodichloromethane	0.778	0.729		-6.3	20
4-Methyl-2-Pentanone	0.661	0.702		6.2	20
Toluene	1.147	1.204		4.97	20
t-1,3-Dichloropropene	0.678	0.709		4.57	20
cis-1,3-Dichloropropene	0.749	0.783		4.54	20
1,1,2-Trichloroethane	0.514	0.498		-3.11	20
2-Hexanone	0.463	0.528		14.04	20
Dibromochloromethane	0.586	0.574		-2.05	20
1,2-Dibromoethane	0.526	0.534		1.52	20
Tetrachloroethene	0.419	0.418		-0.24	20

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:	<u>Alliance</u>	Contract:	<u>GENV01</u>
Lab Code:	<u>ACE</u>	SDG No.:	<u>Q3172</u>
Instrument ID:	<u>MSVOA_V</u>	Calibration Date/Time:	<u>09/25/2025 09:59</u>
Lab File ID:	<u>VV039168.D</u>	Init. Calib. Date(s):	<u>09/22/2025 09/22/2025</u>
Heated Purge: (Y/N)	<u>N</u>	Init. Calib. Time(s):	<u>12:26 16:35</u>
GC Column:	<u>DB-624UI</u>	ID:	<u>0.18</u> (mm)

COMPOUND	RRF	RRF050	MIN RRF	%D	MAX%D
Chlorobenzene	1.393	1.377	0.3	-1.15	20
Ethyl Benzene	2.108	2.320		10.06	20
m/p-Xylenes	0.814	0.936		14.99	20
o-Xylene	0.720	0.841		16.81	20
Styrene	1.263	1.530		21.14	20
Bromoform	0.393	0.402	0.1	2.29	20
Isopropylbenzene	3.627	3.946		8.8	20
1,1,2,2-Tetrachloroethane	1.589	1.396	0.3	-12.15	20
1,3-Dichlorobenzene	1.904	1.857		-2.47	20
1,4-Dichlorobenzene	2.089	1.903		-8.9	20
1,2-Dichlorobenzene	1.779	1.736		-2.42	20
1,2-Dibromo-3-Chloropropane	0.330	0.292		-11.52	20
1,2,4-Trichlorobenzene	1.020	1.062		4.12	20
1,2,3-Trichlorobenzene	1.043	1.082		3.74	20
1,2-Dichloroethane-d4	0.724	0.609		-15.88	20
Dibromofluoromethane	0.402	0.384		-4.48	20
Toluene-d8	1.181	1.121		-5.08	20
4-Bromofluorobenzene	0.486	0.508		4.53	20

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