

VOLATILE CONTINUING CALIBRATION CHECK

Lab Name:	<u>Alliance</u>	Contract:	<u>TETR06</u>
Lab Code:	<u>ACE</u>	SDG No.:	<u>Q3246</u>
Instrument ID:	<u>MSVOA_X</u>	Calibration Date/Time:	<u>10/02/2025 18:58</u>
Lab File ID:	<u>VX047979.D</u>	Init. Calib. Date(s):	<u>09/16/2025 09/16/2025</u>
Heated Purge: (Y/N)	<u>N</u>	Init. Calib. Time(s):	<u>09:23 12:01</u>
GC Column:	<u>DB-624UI</u>	ID:	<u>0.18</u> (mm)

COMPOUND	RRF	RRF050	MIN RRF	%D	MAX%D
Dichlorodifluoromethane	0.518	0.504		-2.7	50
Chloromethane	0.680	0.565	0.1	-16.91	50
Vinyl Chloride	0.666	0.606		-9.01	50
Bromomethane	0.422	0.397		-5.92	50
Chloroethane	0.445	0.403		-9.44	50
Trichlorofluoromethane	0.989	0.943		-4.65	50
1,1,2-Trichlorotrifluoroethane	0.578	0.576		-0.35	50
1,1-Dichloroethene	0.594	0.555		-6.57	50
Acetone	0.346	0.233		-32.66	50
Carbon Disulfide	1.626	1.369		-15.81	50
Methyl tert-butyl Ether	2.169	1.936		-10.74	50
Methyl Acetate	0.770	0.739		-4.03	50
Methylene Chloride	0.732	0.653		-10.79	50
trans-1,2-Dichloroethene	0.640	0.580		-9.38	50
1,1-Dichloroethane	1.263	1.188	0.1	-5.94	50
Cyclohexane	1.052	0.923		-12.26	50
2-Butanone	0.432	0.356		-17.59	50
Carbon Tetrachloride	0.532	0.512		-3.76	50
cis-1,2-Dichloroethene	0.783	0.717		-8.43	50
Bromochloromethane	0.551	0.552		0.18	50
Chloroform	1.276	1.229		-3.68	50
1,1,1-Trichloroethane	1.082	1.053		-2.68	50
Methylcyclohexane	0.556	0.527		-5.22	50
Benzene	1.488	1.435		-3.56	50
1,2-Dichloroethane	0.566	0.530		-6.36	50
Trichloroethene	0.360	0.351		-2.5	50
1,2-Dichloropropane	0.381	0.380		-0.26	50
Bromodichloromethane	0.572	0.568		-0.7	50
4-Methyl-2-Pentanone	0.477	0.455		-4.61	50
Toluene	0.917	0.896		-2.29	50
t-1,3-Dichloropropene	0.584	0.560		-4.11	50
cis-1,3-Dichloropropene	0.624	0.603		-3.37	50
1,1,2-Trichloroethane	0.354	0.355		0.28	50
2-Hexanone	0.343	0.312		-9.04	50
Dibromochloromethane	0.407	0.417		2.46	50
1,2-Dibromoethane	0.360	0.365		1.39	50
Tetrachloroethene	0.326	0.310		-4.91	50
Chlorobenzene	1.136	1.099	0.3	-3.26	50

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>TETR06</u>
Lab Code:	<u>ACE</u>	SDG No.:	<u>Q3246</u>
Instrument ID:	<u>MSVOA_X</u>	Calibration Date/Time:	<u>10/02/2025</u> <u>18:58</u>
Lab File ID:	<u>VX047979.D</u>	Init. Calib. Date(s):	<u>09/16/2025</u> <u>09/16/2025</u>
Heated Purge: (Y/N)	<u>N</u>	Init. Calib. Time(s):	<u>09:23</u> <u>12:01</u>
GC Column:	<u>DB-624UI</u>	ID:	<u>0.18</u> (mm)

COMPOUND	RRF	RRF050	MIN RRF	%D	MAX%D
Ethyl Benzene	1.960	1.899		-3.11	50
m/p-Xylenes	0.729	0.712		-2.33	50
o-Xylene	0.706	0.685		-2.97	50
Styrene	1.236	1.215		-1.7	50
Bromoform	0.293	0.285	0.1	-2.73	50
Isopropylbenzene	3.801	3.548		-6.66	50
1,1,2,2-Tetrachloroethane	1.150	1.078	0.3	-6.26	50
1,3-Dichlorobenzene	1.739	1.605		-7.71	50
1,4-Dichlorobenzene	1.785	1.628		-8.8	50
1,2-Dichlorobenzene	1.657	1.520		-8.27	50
1,2-Dibromo-3-Chloropropane	0.233	0.198		-15.02	50
1,2,4-Trichlorobenzene	1.077	0.969		-10.03	50
1,2,3-Trichlorobenzene	1.002	0.923		-7.88	50
1,2-Dichloroethane-d4	0.796	0.736		-7.54	50
Dibromofluoromethane	0.335	0.338		0.9	50
Toluene-d8	1.160	1.196		3.1	50
4-Bromofluorobenzene	0.440	0.452		2.73	50

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