

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: LIR001
Lab Code: CHEM Case No.: P5021 SAS No.: P5021 SDG No.: P5021
Instrument ID: MSVOA_X Calibration Date(s): 11/21/2024 11/21/2024
Heated Purge: (Y/N) N Calibration Time(s): 09:47 12:03
GC Column: DB-624UI ID: 0.18 (mm)

LAB FILE ID:		RRF001 = VX043925.D		RRF005 = VX043926.D		RRF020 = VX043927.D			
		RRF050 = VX043928.D		RRF100 = VX043929.D		RRF150 = VX043930.D			
COMPOUND	RRF001	RRF005	RRF020	RRF050	RRF100	RRF150	RRF	% RSD	
Methyl tert-butyl Ether	1.732	1.991	2.212	2.264	2.160	2.192	2.092	9.5	
Benzene	1.211	1.369	1.457	1.494	1.388	1.402	1.387	7	
Toluene	0.636	0.839	0.883	0.921	0.850	0.852	0.830	12	
Ethyl Benzene	1.579	1.831	1.910	2.024	1.884	1.938	1.861	8.2	
m/p-Xylenes	0.610	0.663	0.704	0.762	0.700	0.725	0.694	7.6	
o-Xylene	0.543	0.665	0.691	0.757	0.693	0.716	0.678	10.7	
Isopropylbenzene	3.435	3.927	3.996	4.097	3.679	3.927	3.843	6.3	
n-propylbenzene	3.667	4.115	4.469	4.733	4.286	4.486	4.293	8.6	
1,3,5-Trimethylbenzene	2.704	3.115	3.352	3.444	3.118	3.238	3.162	8.2	
tert-Butylbenzene	2.581	3.097	3.266	3.440	3.152	3.269	3.134	9.4	
1,2,4-Trimethylbenzene	2.557	3.173	3.366	3.440	3.124	3.228	3.148	9.9	
sec-Butylbenzene	3.116	3.921	3.968	4.249	3.812	3.974	3.840	10	
p-Isopropyltoluene	2.255	3.043	3.316	3.521	3.219	3.362	3.119	14.5	
n-Butylbenzene	1.897	2.580	2.805	3.184	2.909	3.126	2.750	17.2	
Naphthalene	2.605	3.361	3.698	4.005	3.733	4.058	3.576	15	
1,2-Dichloroethane-d4		0.926	0.876	0.751	0.855	0.880	0.858	7.6	
Dibromofluoromethane		0.371	0.353	0.327	0.359	0.376	0.357	5.4	
Toluene-d8		1.327	1.237	1.104	1.232	1.257	1.232	6.6	
4-Bromofluorobenzene		0.401	0.414	0.387	0.438	0.455	0.419	6.6	

* 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.