

Method Path : Z:\VOASRV\HPCHEM1\MSVOA\_V\METHOD\  
 Method File : SOMVLM102318WMA.M  
 Title : VOC Analysis  
 Last Update : Wed Oct 24 02:41:57 2018  
 Response Via : Initial Calibration

Instrument :  
 MSVOA\_V  
 ClientSampleId :  
 BFB48

## Calibration Files

5 =VV008381.D 10 =VV008382.D 50 =VV008383.D  
 100 =VV008384.D 200 =VV008385.D

	Compound	5	10	50	100	200	Avg	%RSD
-----ISTD-----								
1) I	1,4-Difluorobenzene							
2) T	Dichlorodifluoromet	0.372	0.361	0.354	0.347	0.351	0.357	2.75
3) T	Chloromethane	0.411	0.395	0.377	0.366	0.363	0.382	5.37
4) S	Vinyl Chloride-d3	0.297	0.290	0.269	0.262	0.266	0.277	5.55
5) T	Vinyl chloride	0.368	0.358	0.350	0.346	0.350	0.354	2.45
6) T	Bromomethane	0.109	0.094	0.093	0.093	0.096	0.097	7.13
7) S	Chloroethane-d5	0.219	0.211	0.192	0.191	0.187	0.200	7.14
8) T	Chloroethane	0.178	0.186	0.178	0.202	0.187	0.186	5.17
9) T	Trichlorofluorometh	0.472	0.462	0.454	0.443	0.449	0.456	2.50
10) T	1,1,2-Trichloro-1,2	0.244	0.237	0.239	0.235	0.235	0.238	1.53
11) S	1,1-Dichloroethene-	0.530	0.507	0.489	0.486	0.488	0.500	3.74
12) T	1,1-Dichloroethene	0.235	0.227	0.220	0.219	0.221	0.224	3.03
13) T	Acetone	0.317	0.211	0.217	0.223	0.200	0.234	20.44
14) T	Carbon disulfide	1.062	0.994	0.994	0.994	1.008	1.010	2.89
15) T	Methyl Acetate	0.498	0.476	0.457	0.451	0.447	0.466	4.51
16) T	Methylene chloride	0.439	0.393	0.373	0.368	0.369	0.389	7.72
17) T	trans-1,2-Dichloroe	0.362	0.344	0.335	0.333	0.335	0.342	3.47
18) T	Methyl tert-butyl E	1.304	1.213	1.213	1.192	1.198	1.224	3.73
19) T	1,1-Dichloroethane	0.746	0.709	0.697	0.683	0.691	0.705	3.47
20) T	cis-1,2-Dichloroeth	0.394	0.389	0.388	0.384	0.393	0.390	1.02
21) S	2-Butanone-d5	0.299	0.306	0.302	0.300	0.298	0.301	1.04
22) T	2-Butanone	0.360	0.340	0.333	0.328	0.325	0.337	4.13
23) T	Bromochloromethane	0.197	0.183	0.180	0.179	0.181	0.184	3.99
24) S	Chloroform-d	0.705	0.706	0.669	0.667	0.667	0.683	3.01
25) T	Chloroform	0.723	0.686	0.679	0.675	0.680	0.689	2.84
26) S	1,2-Dichloroethane-	0.480	0.473	0.442	0.443	0.445	0.457	4.05
27) T	1,2-Dichloroethane	0.622	0.576	0.566	0.550	0.562	0.575	4.84
-----ISTD-----								
28) I	Chlorobenzene-d5							
29) T	Cyclohexane	0.800	0.736	0.705	0.698	0.700	0.728	5.93
30) T	1,1,1-Trichloroetha	0.694	0.653	0.637	0.625	0.633	0.648	4.24
31) T	Carbon tetrachlorid	0.584	0.546	0.541	0.536	0.545	0.550	3.52
32) S	Benzene-d6	1.525	1.487	1.385	1.369	1.358	1.425	5.33
33) T	Benzene	1.776	1.649	1.607	1.577	1.563	1.634	5.26
34) T	Trichloroethene	0.457	0.422	0.406	0.400	0.407	0.418	5.48
35) T	Methylcyclohexane	0.736	0.718	0.709	0.698	0.696	0.711	2.29
36) S	1,2-Dichloropropane	0.531	0.529	0.475	0.472	0.467	0.495	6.55
37) T	1,2-Dichloropropane	0.484	0.457	0.441	0.438	0.437	0.451	4.40
38) T	Bromodichloromethan	0.617	0.571	0.558	0.562	0.569	0.575	4.11
39) T	cis-1,3-Dichloropro	0.703	0.722	0.704	0.706	0.714	0.710	1.12
40) T	4-Methyl-2-pentanon	0.714	0.675	0.649	0.651	0.652	0.668	4.15
41) S	Toluene-d8	1.355	1.348	1.268	1.253	1.245	1.294	4.15
42) T	Toluene	1.810	1.725	1.680	1.642	1.631	1.698	4.27
43) S	trans-1,3-Dichlorop	0.247	0.233	0.238	0.240	0.243	0.240	2.23
44) T	trans-1,3-Dichlorop	0.682	0.640	0.660	0.665	0.673	0.664	2.37
45) T	1,1,2-Trichloroetha	0.427	0.405	0.396	0.383	0.384	0.399	4.54
46) T	Tetrachloroethene	0.283	0.268	0.272	0.269	0.273	0.273	2.19
47) S	2-Hexanone-d5	0.263	0.259	0.254	0.258	0.257	0.258	1.22
48) T	2-Hexanone	0.510	0.505	0.503	0.531	0.530	0.516	2.66
49) T	Dibromochloromethan	0.448	0.402	0.427	0.430	0.438	0.429	3.94
50) T	1,2-Dibromoethane	0.449	0.414	0.414	0.407	0.409	0.419	4.05
51) T	Chlorobenzene	1.096	1.045	1.033	1.024	1.028	1.045	2.82
52) T	Ethylbenzene	1.975	1.884	1.898	1.875	1.876	1.902	2.19

Method Path : Z:\VOASRV\HPCHEM1\MSVOA\_V\METHOD\  
 Method File : SOMVLM102318WMA.M  
 Title : VOC Analysis  
 Last Update : Wed Oct 24 02:41:57 2018  
 Response Via : Initial Calibration

Instrument :  
 MSVOA\_V  
 ClientSampleId :  
 BFB48

## Calibration Files

5 =VV008381.D 10 =VV008382.D 50 =VV008383.D  
 100 =VV008384.D 200 =VV008385.D

	Compound	5	10	50	100	200	Avg	%RSD
53) T	m,p-Xylene	0.720	0.671	0.694	0.694	0.695	0.695	2.48
54) T	o-xylene	0.741	0.698	0.681	0.677	0.672	0.694	4.02
55) T	Styrene	1.181	1.134	1.165	1.171	1.194	1.169	1.91
56) T	Isopropylbenzene	1.846	1.779	1.808	1.823	1.820	1.815	1.33
57) S	1,1,2,2-Tetrachloro	0.678	0.687	0.649	0.662	0.673	0.670	2.22
58) T	1,1,2,2-Tetrachloro	0.737	0.682	0.687	0.689	0.699	0.699	3.17
59) T	1,2,3-Trichloroprop	0.579	0.571	0.556	0.559	0.563	0.566	1.67
60) I	1,4-Dichlorobenzene-d	-----ISTD-----						
61) T	Bromoform	0.671	0.628	0.623	0.614	0.619	0.631	3.65
62) T	1,3-Dichlorobenzene	1.792	1.628	1.636	1.615	1.595	1.653	4.78
63) T	1,4-Dichlorobenzene	1.774	1.660	1.642	1.608	1.600	1.657	4.23
64) S	1,2-Dichlorobenzene	1.134	1.081	0.991	0.971	0.937	1.023	8.02
65) T	1,2-Dichlorobenzene	1.847	1.696	1.680	1.627	1.578	1.686	6.01
66) T	1,2-Dibromo-3-chlor	0.363	0.353	0.344	0.339	0.335	0.347	3.21
67) T	1,3,5-Trichlorobenz	1.091	1.032	1.085	1.093	1.097	1.079	2.49
68) T	1,2,4-trichlorobenz	0.796	0.774	0.894	0.939	0.964	0.873	9.74
69) T	Naphthalene	2.602	2.477	2.852	3.042	3.148	2.824	10.06
70) T	1,2,3-Trichlorobenz	0.812	0.809	0.889	0.914	0.940	0.873	6.88

(#) = Out of Range