

Method Path : Z:\VOASRV\HPCHEM1\MSVOA\_W\METHOD\

Method File : SOM2WLM020320S.M

Title : VOC Analysis

Last Update : Mon Feb 03 12:37:09 2020

Response Via : Initial Calibration

## Calibration Files

2.5 =VW014825.D 5 =VW014826.D 25 =VW014827.D  
 50 =VW014828.D 100 =VW014829.D

	Compound	2.5	5	25	50	100	Avg	%RSD
<hr/>								
1) I	1,4-Difluorobenzene			-----ISTD-----				
2) T	Dichlorodifluoromethane	0.354	0.333	0.289	0.288	0.310	0.315	9.10
3) T	Chloromethane	0.490	0.453	0.384	0.411	0.420	0.432	9.41
4) S	Vinyl Chloride-d3	0.282	0.353	0.452	0.439	0.432	0.392	18.56
5) T	Vinyl chloride	0.321	0.434	0.457	0.487	0.471	0.434	15.28
6) T	Bromomethane	0.300	0.273	0.226	0.238	0.233	0.254	12.43
7) S	Chloroethane-d5	0.432	0.386	0.375	0.345	0.333	0.374	10.38
8) T	Chloroethane	0.323	0.294	0.261	0.270	0.260	0.282	9.57
9) T	Trichlorofluoromethane	0.282	0.271	0.232	0.242	0.236	0.253	8.91
10) S	1,1-Dichloroethene	0.868	0.848	0.830	0.804	0.809	0.832	3.24
11) T	1,1,2-Trichloro-1,2	0.436	0.409	0.366	0.370	0.355	0.387	8.83
12) T	1,1-Dichloroethene	0.384	0.375	0.344	0.364	0.358	0.365	4.22
13) T	Acetone	0.233	0.178	0.141	0.126	0.117	0.159	29.95
14) T	Carbon disulfide	1.110	1.159	1.058	1.120	1.095	1.108	3.32
15) T	Methyl Acetate	0.219	0.236	0.252	0.242	0.242	0.238	5.15
16) T	Methylene chloride	0.603	0.577	0.388	0.379	0.362	0.462	25.53
17) T	Methyl tert-butyl E	0.460	0.444	0.461	0.481	0.471	0.463	2.96
18) T	trans-1,2-Dichloroethane	0.418	0.383	0.349	0.366	0.357	0.375	7.32
19) T	1,1-Dichloroethane	0.858	0.790	0.711	0.729	0.715	0.761	8.31
20) S	2-Butanone-d5	0.133	0.141	0.165	0.145	0.152	0.147	8.14
21)	2-Butanone	0.175	0.169	0.173	0.169	0.168	0.171	1.71
22) T	cis-1,2-Dichloroethane	0.404	0.382	0.372	0.390	0.388	0.387	2.99
23) T	Bromochloromethane	0.186	0.172	0.158	0.158	0.154	0.166	8.13
24) S	Chloroform-d	0.937	0.836	0.779	0.714	0.712	0.796	11.83
25) T	Chloroform	0.783	0.726	0.652	0.667	0.644	0.695	8.52
26) S	1,2-Dichloroethane	0.514	0.469	0.449	0.404	0.400	0.447	10.63
27) T	1,2-Dichloroethane	0.518	0.514	0.471	0.469	0.454	0.485	5.93
28) I	Chlorobenzene-d5			-----ISTD-----				
29) S	Benzene-d6	2.018	1.790	1.777	1.642	1.583	1.762	9.54
30) T	Cyclohexane	0.717	0.693	0.758	0.818	0.792	0.756	6.84
31) T	1,1,1-Trichloroethane	0.667	0.611	0.556	0.566	0.533	0.587	9.10
32) T	Carbon tetrachloride	0.599	0.547	0.501	0.519	0.498	0.533	7.84
33) S	1,2-Dichloroproppane	0.670	0.597	0.572	0.531	0.518	0.578	10.53
34) T	Benzene	1.938	1.809	1.682	1.718	1.605	1.751	7.32
35) T	Trichloroethene	0.483	0.440	0.406	0.424	0.409	0.432	7.19
36) T	Methylcyclohexane	0.745	0.736	0.742	0.794	0.768	0.757	3.17
37) S	Toluene-d8	1.652	1.533	1.589	1.483	1.435	1.538	5.57
38) S	trans-1,3-Dichloropropene	0.228	0.218	0.239	0.226	0.232	0.229	3.44
39) S	2-Hexanone-d5	0.081	0.093	0.135	0.125	0.130	0.113	21.51
40) T	1,2-Dichloroproppane	0.540	0.479	0.449	0.459	0.434	0.472	8.69
41) T	Bromodichloromethane	0.592	0.541	0.521	0.533	0.517	0.541	5.59
42) T	cis-1,3-Dichloropropane	0.635	0.616	0.654	0.694	0.686	0.657	5.07
43) T	4-Methyl-2-pentanone	0.286	0.321	0.378	0.376	0.373	0.347	12.02
44) T	Toluene	1.833	1.746	1.726	1.774	1.678	1.751	3.30
45) T	trans-1,3-Dichloropropene	0.513	0.518	0.544	0.565	0.564	0.541	4.58
46) T	1,1,2-Trichloroethane	0.315	0.325	0.309	0.312	0.301	0.312	2.84
47) T	Tetrachloroethene	0.362	0.330	0.306	0.313	0.301	0.323	7.70
48) S	1,1,2,2-Tetrachloroethane	0.453	0.467	0.482	0.427	0.422	0.450	5.75
49) T	2-Hexanone	0.175	0.206	0.276	0.272	0.267	0.239	19.21
50) T	Dibromochloromethane	0.334	0.336	0.328	0.342	0.339	0.336	1.52
51) T	1,2-Dibromoethane	0.284	0.303	0.292	0.300	0.290	0.294	2.54
52) T	Chlorobenzene	1.187	1.087	1.014	1.041	1.014	1.068	6.78

Method Path : Z:\VOASRV\HPCHEM1\MSVOA\_W\METHOD\

Method File : SOM2WLM020320S.M

Title : VOC Analysis

Last Update : Mon Feb 03 12:37:09 2020

Response Via : Initial Calibration

## Calibration Files

2.5 =VW014825.D	5 =VW014826.D	25 =VW014827.D
50 =VW014828.D	100 =VW014829.D	

	Compound	2.5	5	25	50	100	Avg	%RSD
53) T	Ethylbenzene	1.925	1.844	1.880	1.955	1.884	1.898	2.28
54) T	m,p-Xylene	0.695	0.668	0.685	0.727	0.697	0.694	3.09
55) T	o-xylene	0.631	0.602	0.649	0.683	0.669	0.647	4.97
56) T	Styrene	1.051	1.021	1.125	1.174	1.138	1.102	5.76
57) T	Isopropylbenzene	1.689	1.642	1.773	1.852	1.791	1.750	4.79
58) T	1,1,2,2-Tetrachloro	0.390	0.414	0.424	0.418	0.404	0.410	3.26
59)	1,2,3-Trichloroprop	0.298	0.320	0.329	0.316	0.307	0.314	3.81
60) I	1,4-Dichlorobenzene-d	-----ISTD-----						
61) S	1,2-Dichlorobenzene	1.177	1.117	1.035	0.954	0.953	1.047	9.47
62) T	Bromoform	0.398	0.425	0.404	0.417	0.423	0.413	2.80
63) T	1,3-Dichlorobenzene	1.825	1.763	1.582	1.626	1.587	1.677	6.62
64) T	1,4-Dichlorobenzene	1.889	1.762	1.587	1.616	1.598	1.690	7.79
65) T	1,2-Dichlorobenzene	1.659	1.573	1.456	1.490	1.455	1.527	5.78
66) T	1,2-Dibromo-3-chlor	0.140	0.152	0.156	0.148	0.150	0.149	4.02
67)	1,3,5-Trichlorobenz	1.264	1.218	1.148	1.168	1.143	1.188	4.35
68) T	1,2,4-trichlorobenz	0.923	0.884	0.927	0.950	0.957	0.928	3.08
69)	Naphthalene	1.466	1.562	2.070	2.144	2.174	1.883	18.11
70) T	1,2,3-Trichlorobenz	0.857	0.830	0.874	0.870	0.858	0.858	2.00

(#) = Out of Range