

Method Path : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\

Method File : SOMVLM030620WMA.M

Title : VOC Analysis

Last Update : Fri Mar 06 22:55:27 2020

Response Via : Initial Calibration

Calibration Files

5 =VV014760.D	10 =VV014761.D	50 =VV014762.D
100 =VV014763.D	200 =VV014764.D	

	Compound	5	10	50	100	200	Avg	%RSD
<hr/>								
1) I	1,4-Difluorobenzene			-----ISTD-----				
2) T	Dichlorodifluoromethane	0.528	0.476	0.492	0.487	0.488	0.494	4.00
3) T	Chloromethane	0.460	0.424	0.429	0.434	0.430	0.435	3.29
4) S	Vinyl Chloride-d3	0.316	0.318	0.345	0.358	0.354	0.338	5.80
5) T	Vinyl chloride	0.422	0.382	0.402	0.406	0.405	0.403	3.55
6) T	Bromomethane	0.259	0.240	0.248	0.250	0.254	0.250	2.75
7) S	Chloroethane-d5	0.265	0.258	0.267	0.274	0.273	0.268	2.39
8) T	Chloroethane	0.221	0.211	0.218	0.218	0.216	0.217	1.64
9) T	Trichlorofluoromethane	0.570	0.510	0.534	0.541	0.541	0.539	3.98
10) T	1,1,2-Trichloro-1,2	0.299	0.266	0.271	0.277	0.275	0.278	4.58
11) S	1,1-Dichloroethene	0.548	0.546	0.570	0.579	0.569	0.562	2.57
12) T	1,1-Dichloroethene	0.290	0.264	0.271	0.266	0.266	0.272	3.99
13) T	Acetone	0.243	0.205	0.217	0.201	0.188	0.211	9.90
14) T	Carbon disulfide	0.926	0.833	0.888	0.913	0.930	0.898	4.42
15) T	Methyl Acetate	0.390	0.384	0.404	0.401	0.399	0.396	2.10
16) T	Methylene chloride	0.397	0.355	0.365	0.364	0.364	0.369	4.37
17) T	trans-1,2-Dichloroethane	0.357	0.311	0.339	0.335	0.337	0.336	4.82
18) T	Methyl tert-butyl E	1.121	1.032	1.105	1.114	1.121	1.099	3.45
19) T	1,1-Dichloroethane	0.689	0.619	0.660	0.655	0.654	0.655	3.82
20) T	cis-1,2-Dichloroethane	0.380	0.354	0.378	0.380	0.385	0.375	3.24
21) S	2-Butanone-d5	0.191	0.238	0.270	0.281	0.281	0.252	15.23
22) T	2-Butanone	0.236	0.234	0.306	0.308	0.306	0.278	14.14
23) T	Bromochloromethane	0.197	0.186	0.189	0.192	0.191	0.191	2.00
24) S	Chloroform-d	0.619	0.582	0.644	0.676	0.682	0.641	6.48
25) T	Chloroform	0.683	0.685	0.700	0.670	0.658	0.679	2.35
26) S	1,2-Dichloroethane	0.410	0.429	0.443	0.451	0.444	0.435	3.75
27) T	1,2-Dichloroethane	0.555	0.499	0.525	0.525	0.526	0.526	3.76
28) I	Chlorobenzene-d5			-----ISTD-----				
29) T	Cyclohexane	0.613	0.558	0.621	0.624	0.630	0.609	4.76
30) T	1,1,1-Trichloroethane	0.611	0.560	0.573	0.579	0.575	0.580	3.24
31) T	Carbon tetrachloride	0.495	0.465	0.494	0.500	0.506	0.492	3.20
32) S	Benzene-d6	1.175	1.286	1.368	1.393	1.379	1.320	6.89
33) T	Benzene	1.497	1.379	1.463	1.458	1.445	1.449	2.99
34) T	Trichloroethene	0.391	0.365	0.376	0.378	0.383	0.379	2.56
35) T	Methylcyclohexane	0.636	0.582	0.625	0.639	0.642	0.625	3.99
36) S	1,2-Dichloropropane	0.398	0.437	0.444	0.458	0.451	0.438	5.39
37) T	1,2-Dichloropropane	0.401	0.374	0.397	0.393	0.393	0.391	2.66
38) T	Bromodichloromethane	0.502	0.470	0.500	0.508	0.510	0.498	3.21
39) T	cis-1,3-Dichloropropane	0.592	0.560	0.611	0.647	0.658	0.614	6.53
40) T	4-Methyl-2-pentanone	0.547	0.543	0.575	0.590	0.581	0.567	3.65
41) S	Toluene-d8	1.174	1.243	1.341	1.360	1.347	1.293	6.27
42) T	Toluene	1.602	1.544	1.613	1.623	1.609	1.598	1.96
43) S	trans-1,3-Dichloropropene	0.179	0.197	0.224	0.233	0.236	0.214	11.52
44) T	trans-1,3-Dichloropropene	0.540	0.520	0.569	0.580	0.588	0.559	5.10
45) T	1,1,2-Trichloroethane	0.364	0.360	0.356	0.363	0.360	0.361	0.91
46) T	Tetrachloroethene	0.317	0.301	0.306	0.311	0.314	0.310	2.04
47) S	2-Hexanone-d5	0.153	0.183	0.203	0.215	0.217	0.194	13.86
48) T	2-Hexanone	0.427	0.415	0.467	0.463	0.459	0.446	5.27
49) T	Dibromochloromethane	0.381	0.364	0.396	0.405	0.418	0.393	5.34
50) T	1,2-Dibromoethane	0.380	0.383	0.391	0.396	0.399	0.390	2.09
51) T	Chlorobenzene	1.032	0.989	1.031	1.042	1.047	1.028	2.25
52) T	Ethylbenzene	1.769	1.694	1.829	1.862	1.848	1.800	3.86

Method Path : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\

Method File : SOMVLM030620WMA.M

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Response Via : Initial Calibration

Calibration Files

5 =VV014760.D	10 =VV014761.D	50 =VV014762.D
100 =VV014763.D	200 =VV014764.D	

	Compound	5	10	50	100	200	Avg	%RSD
53) T	m,p-Xylene	0.663	0.629	0.683	0.701	0.701	0.676	4.46
54) T	o-xylene	0.646	0.616	0.677	0.687	0.692	0.664	4.84
55) T	Styrene	1.055	1.046	1.176	1.201	1.196	1.135	6.84
56) T	Isopropylbenzene	1.769	1.646	1.796	1.842	1.835	1.778	4.47
57) S	1,1,2,2-Tetrachloro	0.543	0.579	0.615	0.635	0.637	0.602	6.69
58) T	1,1,2,2-Tetrachloro	0.627	0.598	0.618	0.629	0.632	0.621	2.21
59)	1,2,3-Trichloroprop	0.506	0.487	0.502	0.504	0.501	0.500	1.49
60) I	1,4-Dichlorobenzene-d	-----ISTD-----						
61) T	Bromoform	0.531	0.516	0.537	0.554	0.575	0.542	4.17
62) T	1,3-Dichlorobenzene	1.674	1.567	1.619	1.613	1.612	1.617	2.34
63) T	1,4-Dichlorobenzene	1.656	1.585	1.645	1.634	1.626	1.629	1.68
64) S	1,2-Dichlorobenzene	0.932	0.992	0.995	1.020	1.017	0.991	3.56
65) T	1,2-Dichlorobenzene	1.710	1.566	1.586	1.598	1.588	1.610	3.56
66) T	1,2-Dibromo-3-chlor	0.271	0.253	0.272	0.282	0.288	0.273	4.86
67)	1,3,5-Trichlorobenz	1.273	1.190	1.232	1.262	1.296	1.250	3.29
68) T	1,2,4-trichlorobenz	1.129	1.081	1.142	1.173	1.222	1.149	4.55
69)	Naphthalene	1.590	1.607	1.745	1.844	1.865	1.730	7.43
70) T	1,2,3-Trichlorobenz	1.117	1.070	1.153	1.187	1.202	1.146	4.69

(#) = Out of Range