

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

Method File : SFAMXTR102920WMA.M

Title : TRACE VOA SFAM1.0

Last Update : Thu Oct 29 06:09:53 2020

Response Via : Initial Calibration

Calibration Files

0.5 =VX019164.D	1 =VX019165.D	5 =VX019170.D
10 =VX019171.D	20 =VX019168.D	

	Compound	0.5	1	5	10	20	Avg	%RSD
-----ISTD-----								
1) I	1,4-Difluorobenzene							
2) T	Dichlorodifluoromethane	0.374	0.394	0.425	0.440	0.433	0.413	6.78
3) T	Chloromethane	0.347	0.334	0.367	0.372	0.363	0.357	4.35
4) S	Vinyl Chloride-d3	0.303	0.313	0.323	0.333	0.345	0.323	5.14
5) T	Vinyl chloride	0.386	0.376	0.414	0.421	0.412	0.402	4.85
6) T	Bromomethane	0.253	0.243	0.262	0.271	0.266	0.259	4.33
7) S	Chloroethane-d5	0.261	0.262	0.272	0.275	0.279	0.270	2.93
8) T	Chloroethane	0.236	0.232	0.249	0.254	0.247	0.244	3.85
9) T	Trichlorofluoromethane	0.522	0.551	0.599	0.620	0.609	0.580	7.23
10) T	1,1,2-Trichloro-1,2-d	0.302	0.305	0.342	0.350	0.349	0.330	7.36
11) S	1,1-Dichloroethene	0.590	0.617	0.635	0.652	0.664	0.632	4.65
12) T	1,1-Dichloroethene	0.298	0.307	0.333	0.335	0.331	0.321	5.35
13) T	Acetone	0.036	0.038	0.043	0.043	0.042	0.040	7.59
14) T	Carbon disulfide	0.967	0.931	1.002	1.038	1.034	0.995	4.57
15) T	Methyl Acetate	0.091	0.099	0.116	0.118	0.113	0.108	11.11
16) T	Methylene chloride	0.444	0.444	0.344	0.346	0.334	0.382	14.74
17) T	Methyl tert-butyl E	0.699	0.725	0.799	0.810	0.804	0.768	6.70
18) T	trans-1,2-Dichloroethane	0.317	0.319	0.346	0.354	0.350	0.337	5.34
19) T	1,1-Dichloroethane	0.547	0.535	0.593	0.611	0.602	0.577	5.96
20) S	2-Butanone-d5	0.061	0.064	0.075	0.077	0.077	0.071	10.53
21) T	2-Butanone	0.069	0.063	0.075	0.074	0.073	0.071	6.72
22) T	cis-1,2-Dichloroethane	0.335	0.342	0.365	0.380	0.374	0.359	5.51
23) T	Bromochloromethane	0.136	0.137	0.156	0.159	0.158	0.149	7.79
24) S	Chloroform-d	0.561	0.570	0.622	0.644	0.652	0.610	6.90
25) T	Chloroform	0.556	0.575	0.611	0.634	0.626	0.600	5.56
26) S	1,2-Dichloroethane-d	0.396	0.342	0.328	0.325	0.325	0.343	8.81
27) T	1,2-Dichloroethane	0.358	0.351	0.389	0.396	0.385	0.376	5.26
28) I	Chlorobenzene-d5							
29) T	1,1,1-Trichloroethane	0.562	0.574	0.636	0.669	0.645	0.617	7.55
30) T	Cyclohexane	0.555	0.599	0.628	0.661	0.634	0.615	6.58
31) T	Carbon tetrachloride	0.469	0.498	0.535	0.570	0.558	0.526	8.05
32) S	Benzene-d6	1.264	1.308	1.373	1.439	1.430	1.363	5.58
33) T	Benzene	1.406	1.390	1.536	1.584	1.531	1.489	5.79
34) T	Trichloroethene	0.385	0.384	0.419	0.430	0.418	0.407	5.26
35) T	Methylcyclohexane	0.599	0.611	0.673	0.715	0.696	0.659	7.83
36) S	1,2-Dichloropropane	0.356	0.360	0.396	0.410	0.408	0.386	6.75
37) T	1,2-Dichloropropane	0.321	0.329	0.365	0.377	0.363	0.351	6.97
38) T	Bromodichloromethane	0.399	0.424	0.471	0.501	0.490	0.457	9.61
39) T	cis-1,3-Dichloropropane	0.475	0.481	0.559	0.596	0.585	0.539	10.69
40) T	4-Methyl-2-pentanone	0.166	0.168	0.200	0.206	0.198	0.187	10.31
41) S	Toluene-d8	1.392	1.327	1.312	1.337	1.339	1.341	2.26
42) T	Toluene	1.465	1.517	1.670	1.721	1.671	1.609	6.89
43) S	trans-1,3-Dichloropropene	0.143	0.150	0.171	0.182	0.185	0.166	11.53
44) T	trans-1,3-Dichloropropene	0.389	0.392	0.474	0.503	0.500	0.452	12.56
45) T	1,1,2-Trichloroethane	0.234	0.239	0.268	0.274	0.267	0.256	7.15
46) S	2-Hexanone-d5	0.060	0.065	0.077	0.082	0.080	0.073	13.30
47) T	Tetrachloroethene	0.283	0.296	0.320	0.331	0.326	0.311	6.65
48) T	2-Hexanone	0.113	0.117	0.141	0.146	0.139	0.131	11.59
49) T	Dibromochloromethane	0.254	0.260	0.308	0.330	0.332	0.297	12.62
50) T	1,2-Dibromoethane	0.212	0.216	0.249	0.259	0.252	0.238	9.23
51) T	Chlorobenzene	0.936	0.950	1.042	1.101	1.063	1.019	7.10
52) T	Ethylbenzene	1.608	1.668	1.863	1.957	1.903	1.800	8.48

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Title : TRACE VOA SFAM1.0

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Response Via : Initial Calibration

Calibration Files

0.5 =VX019164.D	1 =VX019165.D	5 =VX019170.D
10 =VX019171.D	20 =VX019168.D	

	Compound	0.5	1	5	10	20	Avg	%RSD
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53) T	m,p-xylene	0.608	0.638	0.717	0.761	0.729	0.691	9.35
54) T	o-xylene	0.570	0.618	0.688	0.721	0.710	0.662	9.80
55) T	Styrene	0.913	0.947	1.134	1.222	1.188	1.081	13.11
56) S	1,1,2,2-Tetrachloro	0.260	0.269	0.305	0.325	0.323	0.296	10.23
57) T	1,1,2,2-Tetrachloro	0.258	0.263	0.307	0.321	0.311	0.292	10.04
58) I	1,4-Dichlorobenzene-d	-----ISTD-----						
59) T	Bromoform	0.288	0.290	0.341	0.356	0.366	0.328	11.16
60) T	Isopropylbenzene	3.360	3.454	3.805	3.923	3.821	3.673	6.77
61) T	1,2,3-Trichloroprop	0.432	0.421	0.472	0.465	0.453	0.449	4.81
62) T	1,3,5-Trimethylbenz	2.706	2.821	3.225	3.360	3.265	3.075	9.48
63) T	1,2,4-Trimethylbenz	2.712	2.893	3.267	3.390	3.326	3.118	9.55
64) T	1,3-Dichlorobenzene	1.508	1.541	1.705	1.724	1.680	1.632	6.13
65) T	1,4-Dichlorobenzene	1.567	1.542	1.661	1.710	1.665	1.629	4.36
66) S	1,2-Dichlorobenzene	0.850	0.880	0.951	0.976	0.983	0.928	6.43
67) T	1,2-Dichlorobenzene	1.393	1.418	1.558	1.590	1.540	1.500	5.88
68) T	1,2-Dibromo-3-chlor	0.091	0.094	0.109	0.108	0.110	0.103	8.93
69)	1,3,5-Trichlorobenz	1.026	1.092	1.273	1.305	1.292	1.198	10.78
70) T	1,2,4-trichlorobenz	0.941	0.950	1.083	1.137	1.115	1.045	8.93
71) T	Naphthalene	1.585	1.639	2.017	2.078	2.081	1.880	13.12
72) T	1,2,3-Trichlorobenz	0.823	0.851	0.963	0.994	0.981	0.922	8.63

(#) = Out of Range