

Method Path : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\
 Method File : SFAM01-EPA-BG092820.M
 Title : SVOA CALIBRATION
 Last Update : Mon Sep 28 14:51:01 2020
 Response Via : Initial Calibration

Calibration Files

5 =BG046714.D 10 =BG046715.D 20 =BG046716.D
 40 =BG046717.D 80 =BG046718.D 160 =BG046719.D

	Compound	5	10	20	40	80	160	Avg	%RSD
1) I	1,4-Dichlorobenzene-d	-----ISTD-----							
2)	1,4-Dioxane	0.495	0.486	0.443	0.496	0.522		0.488	5.91
3) S	1,4-Dioxane-d8	0.540	0.406	0.343	0.414	0.431		0.427	16.74
4)	Benzaldehyde		1.249	1.166	1.303	1.224	1.027	1.194	8.82
5) S	Phenol-d5		1.610	1.535	1.716	1.912	1.636	1.682	8.57
6)	Phenol		1.758	1.613	1.850	1.961	1.700	1.776	7.57
7) S	Bis-(2-Chloroethy		1.123	0.982	1.064	1.149	0.990	1.062	7.10
8)	Bis(2-Chloroethyl		1.408	1.270	1.404	1.516	1.308	1.381	6.97
9) S	2-Chlorophenol-d4	1.148	1.186	1.117	1.283	1.381		1.223	8.84
10)	2-Chlorophenol	1.214	1.222	1.143	1.269	1.414		1.252	8.04
11)	2-Methylphenol		1.248	1.222	1.366	1.480	1.307	1.325	7.78
12)	2,2'-oxybis(1-Chl		2.261	2.083	2.211	2.370	2.017	2.188	6.43
13) S	4-Methylphenol-d8		1.259	1.197	1.366	1.479	1.307	1.321	8.17
14)	Acetophenone		2.216	2.010	2.194	2.363	2.040	2.165	6.63
15) P	N-Nitroso-di-n-pr	1.172	1.236	1.094	1.199	1.278		1.196	5.82
16)	4-Methylphenol		1.353	1.237	1.446	1.554	1.348	1.387	8.56
17)	Hexachloroethane	0.523	0.582	0.530	0.569	0.626		0.566	7.40
18) I	Naphthalene-d8	-----ISTD-----							
19) S	Nitrobenzene-d5	0.104	0.121	0.111	0.139	0.164		0.128	18.71
20)	Nitrobenzene	0.307	0.350	0.333	0.420	0.464		0.375	17.44
21)	Isophorone	0.829	0.891	0.729	0.855	0.881		0.837	7.76
22) S	2-Nitrophenol-d4	0.116	0.130	0.112	0.148	0.175		0.136	19.06
23) C	2-Nitrophenol	0.124	0.131	0.125	0.174	0.196		0.150	22.08#
24)	2,4-Dimethylpheno	0.389	0.407	0.362	0.420	0.443		0.404	7.61
25)	Bis(2-Chloroethox	0.484	0.511	0.424	0.484	0.507		0.482	7.20
26) S	2,4-Dichloropheno	0.279	0.348	0.305	0.362	0.395		0.338	13.60
27) C	2,4-Dichloropheno	0.292	0.341	0.294	0.353	0.383		0.333	11.79
28)	Naphthalene	1.112	1.173	0.954	1.108	1.159		1.101	7.93
29) S	4-Chloroaniline-d		0.467	0.381	0.445	0.453	0.348	0.419	12.33
30)	4-Chloroaniline		0.446	0.383	0.420	0.432	0.331	0.403	11.50
31) C	Hexachlorobutadie	0.306	0.312	0.256	0.299	0.314		0.297	8.02
32)	Caprolactam		0.121	0.101	0.124	0.131	0.117	0.119	9.42
33) C	4-Chloro-3-methyl	0.345	0.389	0.323	0.387	0.410		0.371	9.64
34)	2-Methylnaphthale	0.837	0.857	0.707	0.812	0.840		0.811	7.42
35)	1-Methylnaphthale	0.803	0.844	0.680	0.785	0.819		0.786	8.06
36) I	Acenaphthene-d10	-----ISTD-----							
37)	1,2,4,5-Tetrachlo	0.773	0.759	0.623	0.739	0.783		0.735	8.84
38)	Hexachlorocyclope		0.441	0.379	0.481	0.543	0.504	0.470	13.31
39) C	2,4,6-Trichloroph	0.363	0.419	0.355	0.442	0.501		0.416	14.51
40)	2,4,5-Trichloroph	0.379	0.432	0.377	0.470	0.526		0.437	14.54
41)	1,1'-Biphenyl	1.635	1.625	1.330	1.553	1.611		1.551	8.21
42)	2-Chloronaphthale	1.268	1.309	1.052	1.233	1.294		1.231	8.47
43)	2-Nitroaniline	0.200	0.248	0.239	0.330	0.400		0.283	28.49
44) S	Dimethylphthalate	1.508	1.607	1.282	1.542	1.624		1.513	9.06
45)	Dimethylphthalate	1.585	1.672	1.333	1.595	1.622		1.562	8.45
46)	2,6-Dinitrotoluen	0.209	0.230	0.203	0.277	0.320		0.248	20.04
47) S	Acenaphthylene-d8	1.860	1.938	1.551	1.825	1.880		1.811	8.34
48)	Acenaphthylene	1.889	1.928	1.555	1.823	1.858		1.811	8.19
49)	3-Nitroaniline		0.224	0.213	0.269	0.294	0.237	0.247	13.43
50) C	Acenaphthene	1.327	1.339	1.093	1.300	1.349		1.282	8.37
51)	2,4-Dinitrophenol		0.109	0.100	0.148	0.184	0.184	0.145	27.71

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	Compound	5	10	20	40	80	160	Avg	%RSD
52) S	4-Nitrophenol-d4		0.209	0.187	0.255	0.288	0.266	0.241	17.38
53)	4-Nitrophenol		0.200	0.186	0.248	0.289	0.265	0.238	18.36
54)	Dibenzofuran	2.010	1.993	1.619	1.871	1.909		1.880	8.34
55)	2,4-Dinitrotoluen	0.280	0.288	0.285	0.382	0.454		0.338	23.05
56)	2,3,4,6-Tetrachlo	0.345	0.396	0.356	0.450	0.505		0.411	16.26
57)	Diethylphthalate	1.541	1.636	1.324	1.587	1.635		1.544	8.37
58) S	Fluorene-d10	1.426	1.476	1.221	1.418	1.493		1.407	7.73
59)	Fluorene	1.646	1.691	1.289	1.532	1.583		1.548	10.15
60)	4-Chlorophenyl-ph	0.908	0.942	0.735	0.856	0.932		0.874	9.71
61)	4-Nitroaniline		0.278	0.243	0.303	0.320	0.272	0.283	10.44
62) I	Phenanthrene-d10		-----ISTD-----						
63) S	4,6-Dinitro-2-met		0.073	0.065	0.090	0.110	0.114	0.090	24.11
64)	4,6-Dinitro-2-met		0.077	0.070	0.100	0.126	0.124	0.099	26.03
65)	N-Nitrosodiphenyl	0.588	0.613	0.489	0.584	0.600		0.575	8.57
66)	4-Bromophenyl-phe	0.253	0.264	0.208	0.255	0.272		0.251	9.94
67)	Hexachlorobenzene	0.127	0.251	0.121	0.257	0.268		0.205	36.19
68)	Atrazine		0.250	0.199	0.245	0.260	0.226	0.236	10.26
69) C	Pentachlorophenol		0.130	0.119	0.159	0.181	0.166	0.151	17.01
70)	Phenanthrene	1.162	1.188	0.933	1.095	1.106		1.097	9.04
71) S	Anthracene-d10	0.968	0.991	0.789	0.912	0.925		0.917	8.54
72)	Anthracene	1.195	1.200	0.943	1.093	1.093		1.105	9.46
73)	1,2,3,4-Tetrachlo	0.316	0.319	0.266	0.317	0.336		0.311	8.55
74)	Pentachlorobenzen	0.336	0.333	0.267	0.317	0.339		0.319	9.42
75)	Carbazole		1.002	0.795	0.939	0.937	0.774	0.889	11.18
76)	Di-n-butylphthala	1.060	1.171	0.955	1.127	1.115		1.085	7.66
77) C	Fluoranthene	1.476	1.522	1.197	1.381	1.331		1.381	9.26
78) I	Chrysene-d12		-----ISTD-----						
79) S	Pyrene-d10	1.048	1.063	0.918	1.084	1.092		1.041	6.83
80)	Pyrene	1.337	1.375	1.132	1.304	1.287		1.287	7.23
81)	Butylbenzylphthal	0.405	0.449	0.398	0.489	0.520		0.452	11.61
82)	3,3'-Dichlorobenz		0.485	0.420	0.509	0.497	0.383	0.459	11.87
83)	Benzo(a)anthracen	1.390	1.413	1.141	1.317	1.335		1.319	8.11
84)	Bis(2-ethylhexyl)	0.624	0.680	0.586	0.701	0.737		0.666	9.10
85)	Chrysene	1.358	1.360	1.092	1.277	1.286		1.274	8.57
86) I	Perylene-d12		-----ISTD-----						
87)	Di-n-octyl phthal		1.136	0.973	1.177	1.232	1.022	1.108	9.73
88)	Benzo(b)fluoranth	1.380	1.401	1.113	1.323	1.397		1.323	9.17
89)	Benzo(k)fluoranth	1.324	1.401	1.097	1.323	1.325		1.294	8.89
90) S	Benzo(a)pyrene-d1	1.075	1.127	0.911	1.126	1.166		1.081	9.29
91) C	Benzo(a)pyrene	1.284	1.278	1.026	1.231	1.263		1.216	8.93
92)	Indeno(1,2,3-cd)p	1.484	1.544	1.249	1.515	1.590		1.476	9.02
93)	Dibenzo(a,h)anthr	1.245	1.284	1.041	1.253	1.299		1.224	8.55
94)	Benzo(g,h,i)peryl	1.258	1.289	1.039	1.249	1.319		1.231	8.98

(#) = Out of Range