

Data Path : Z:\HPCHEM1\BNA_M\Data\BM090917\
 Data File : BM011477.D
 Acq On : 09 Sep 2017 09:07
 Operator : SJ/JU
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02009

Quant Time: Sep 11 02:51:51 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090817MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 08 17:01:48 2017
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
 Max. RRF Dev : 20% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	102	0.00
2	1,4-Dioxane	0.487	0.494	-1.4	103	0.00
3 S	1,4-Dioxane-d8	0.445	0.449	-0.9	103	0.00
4	Benzaldehyde	1.097	0.991	9.7	86	0.00
5 S	Phenol-d5	1.920	1.824	5.0	100	0.00
6	Phenol	1.906	1.822	4.4	100	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	1.294	1.263	2.4	100	0.00
8	Bis(2-Chloroethyl)ether	1.501	1.479	1.5	102	0.00
9 S	2-Chlorophenol-d4	1.447	1.451	-0.3	103	0.00
10	2-Chlorophenol	1.412	1.415	-0.2	103	0.00
11	2-Methylphenol	1.445	1.396	3.4	102	0.00
12	2,2'-oxybis(1-Chloropropane	2.433	2.425	0.3	101	0.00
13 S	4-Methylphenol-d8	1.499	1.440	3.9	100	0.00
14	Acetophenone	2.292	2.241	2.2	101	0.00
15 P	N-Nitroso-di-n-propylamine	1.230	1.218	1.0	101	0.00
16	4-Methylphenol	1.582	1.559	1.5	103	0.00
17	Hexachloroethane	0.540	0.531	1.7	102	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	103	0.00
19 S	Nitrobenzene-d5	0.152	0.150	1.3	101	0.00
20	Nitrobenzene	0.357	0.355	0.6	103	0.00
21	Isophorone	0.719	0.713	0.8	102	0.00
22 S	2-Nitrophenol-d4	0.158	0.154	2.5	102	0.00
23 C	2-Nitrophenol	0.166	0.166	0.0	103	0.00
24	2,4-Dimethylphenol	0.355	0.350	1.4	101	0.00
25	Bis(2-Chloroethoxy)methane	0.469	0.461	1.7	102	0.00
26 S	2,4-Dichlorophenol-d3	0.318	0.321	-0.9	103	0.00
27 C	2,4-Dichlorophenol	0.306	0.310	-1.3	103	0.00
28	Naphthalene	1.037	1.028	0.9	102	0.00
29 S	4-Chloroaniline-d4	0.351	0.355	-1.1	118	0.00
30	4-Chloroaniline	0.333	0.339	-1.8	117	0.00
31 C	Hexachlorobutadiene	0.174	0.180	-3.4	108	0.00
32	Caprolactam	0.096	0.098	-2.1	108	0.00
33 C	4-Chloro-3-methylphenol	0.325	0.327	-0.6	104	0.00
34	2-Methylnaphthalene	0.770	0.767	0.4	102	0.00
35	1-Methylnaphthalene	0.734	0.736	-0.3	103	0.00
36 I	Acenaphthene-d10	1.000	1.000	0.0	108	0.00
37	1,2,4,5-Tetrachlorobenzene	0.593	0.598	-0.8	107	0.00
38	Hexachlorocyclopentadiene	0.333	0.286	14.1	98	0.00
39 C	2,4,6-Trichlorophenol	0.404	0.403	0.2	106	0.00
40	2,4,5-Trichlorophenol	0.410	0.410	0.0	106	0.00
41	1,1'-Biphenyl	1.664	1.640	1.4	104	0.00
42	2-Chloronaphthalene	1.252	1.238	1.1	106	0.00
43	2-Nitroaniline	0.389	0.376	3.3	101	0.00
44 S	Dimethylphthalate-d6	1.691	1.701	-0.6	106	0.00

Data Path : Z:\HPCHEM1\BNA_M\Data\BM090917\
 Data File : BM011477.D
 Acq On : 09 Sep 2017 09:07
 Operator : SJ/JU
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02009

Quant Time: Sep 11 02:51:51 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090817MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 08 17:01:48 2017
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
 Max. RRF Dev : 20% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	Dimethylphthalate	1.619	1.618	0.1	106	0.00
46	2,6-Dinitrotoluene	0.286	0.295	-3.1	109	0.00
47 S	Acenaphthylene-d8	2.045	2.044	0.0	105	0.00
48	Acenaphthylene	2.058	2.054	0.2	105	0.00
49	3-Nitroaniline	0.353	0.335	5.1	105	0.00
50 C	Acenaphthene	1.391	1.376	1.1	106	0.00
51	2,4-Dinitrophenol	0.177	0.155	12.4	108	0.00
52 S	4-Nitrophenol-d4	0.318	0.298	6.3	106	0.00
53	4-Nitrophenol	0.204	0.194	4.9	108	0.00
54	Dibenzofuran	1.945	1.951	-0.3	106	0.00
55	2,4-Dinitrotoluene	0.420	0.431	-2.6	108	0.00
56	2,3,4,6-Tetrachlorophenol	0.373	0.381	-2.1	109	0.00
57	Diethylphthalate	1.685	1.691	-0.4	106	0.00
58 S	Fluorene-d10	1.465	1.474	-0.6	107	0.00
59	Fluorene	1.630	1.645	-0.9	106	0.00
60	4-Chlorophenyl-phenylether	0.764	0.774	-1.3	108	0.00
61	4-Nitroaniline	0.378	0.358	5.3	103	0.00
62 I	Phenanthrene-d10	1.000	1.000	0.0	109	0.00
63 S	4,6-Dinitro-2-methylphenol-	0.116	0.106	8.6	110	0.00
64	4,6-Dinitro-2-methylphenol	0.119	0.109	8.4	108	0.00
65	N-Nitrosodiphenylamine	0.609	0.609	0.0	108	0.00
66	4-Bromophenyl-phenylether	0.204	0.208	-2.0	112	0.00
67	Hexachlorobenzene	0.225	0.231	-2.7	113	0.00
68	Atrazine	0.211	0.206	2.4	109	0.00
69 C	Pentachlorophenol	0.144	0.135	6.2	112	0.00
70	Phenanthrene	1.115	1.118	-0.3	107	0.00
71 S	Anthracene-d10	0.985	0.986	-0.1	108	0.00
72	Anthracene	1.144	1.163	-1.7	107	0.00
73	1,2,3,4-Tetrachlorobenzene	0.267	0.260	2.6	106	0.00
74	Pentachlorobenzene	0.270	0.274	-1.5	110	0.00
75	Carbazole	0.976	1.037	-6.2	108	0.00
76	Di-n-butylphthalate	1.282	1.331	-3.8	106	0.00
77 C	Fluoranthene	1.147	1.300	-13.3	109	0.00
78 I	Chrysene-d12	1.000	1.000	0.0	112	0.00
79 S	Pyrene-d10	0.936	0.934	0.2	110	0.00
80	Pyrene	1.167	1.193	-2.2	108	0.00
81	Butylbenzylphthalate	0.537	0.535	0.4	111	0.00
82	3,3'-Dichlorobenzidine	0.363	0.369	-1.7	112	0.00
83	Benzo(a)anthracene	1.101	1.129	-2.5	110	0.00
84	Bis(2-ethylhexyl)phthalate	0.824	0.828	-0.5	108	0.00
85	Chrysene	0.998	1.019	-2.1	111	0.00
86 I	Perylene-d12	1.000	1.000	0.0	115	0.00
87	Di-n-octyl phthalate	1.419	1.476	-4.0	108	0.00

Data Path : Z:\HPCHEM1\BNA_M\Data\BM090917\
Data File : BM011477.D
Acq On : 09 Sep 2017 09:07
Operator : SJ/JU
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTD02009

Quant Time: Sep 11 02:51:51 2017
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090817MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Sep 08 17:01:48 2017
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
Max. RRF Dev : 20% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.203	1.180	1.9	113	0.00
89	Benzo(k)fluoranthene	1.156	1.125	2.7	110	0.00
90 S	Benzo(a)pyrene-d12	0.953	0.952	0.1	115	0.00
91 C	Benzo(a)pyrene	1.136	1.133	0.3	113	0.00
92	Indeno(1,2,3-cd)pyrene	1.249	1.290	-3.3	116	0.00
93	Dibenzo(a,h)anthracene	1.059	1.092	-3.1	117	0.00
94	Benzo(g,h,i)perylene	1.012	1.057	-4.4	118	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0