

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110618\
 Data File : BM017530.D
 Acq On : 06 Nov 2018 15:58
 Operator : MJ/SJ
 Sample : SSTDCCC020EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02092

Quant Time: Nov 06 16:33:35 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102418MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 06 12:04:23 2018
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% 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	0.438	0.470	-7.3	109	0.00
3 S	1,4-Dioxane-d8	0.435	0.451	-3.7	106	0.00
4	Benzaldehyde	0.348	0.316	9.2	95	0.00
5 S	Phenol-d5	1.827	1.772	3.0	102	0.00
6	Phenol	1.814	1.775	2.1	103	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	1.065	1.121	-5.3	109	0.00
8	Bis(2-Chloroethyl)ether	1.374	1.390	-1.2	103	0.00
9 S	2-Chlorophenol-d4	1.416	1.454	-2.7	105	0.00
10	2-Chlorophenol	1.423	1.441	-1.3	104	0.00
11	2-Methylphenol	1.366	1.356	0.7	105	0.00
12	2,2'-oxybis(1-Chloropropane	1.831	2.069	-13.0	115	0.00
13 S	4-Methylphenol-d8	1.445	1.463	-1.2	107	0.00
14	Acetophenone	2.226	2.271	-2.0	105	0.00
15 P	N-Nitroso-di-n-propylamine	1.093	1.181	-8.1	109	0.00
16	4-Methylphenol	1.471	1.471	0.0	104	0.00
17	Hexachloroethane	0.556	0.592	-6.5	110	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	102	0.00
19 S	Nitrobenzene-d5	0.156	0.157	-0.6	105	0.00
20	Nitrobenzene	0.369	0.382	-3.5	108	0.00
21	Isophorone	0.671	0.722	-7.6	110	0.00
22 S	2-Nitrophenol-d4	0.175	0.179	-2.3	106	0.00
23 C	2-Nitrophenol	0.185	0.187	-1.1	104	0.00
24	2,4-Dimethylphenol	0.367	0.385	-4.9	108	0.00
25	Bis(2-Chloroethoxy)methane	0.426	0.446	-4.7	108	0.00
26 S	2,4-Dichlorophenol-d3	0.326	0.341	-4.6	108	0.00
27 C	2,4-Dichlorophenol	0.312	0.321	-2.9	106	0.00
28	Naphthalene	1.046	1.054	-0.8	105	0.00
29 S	4-Chloroaniline-d4	0.266	0.303	-13.9	123	0.00
30	4-Chloroaniline	0.272	0.305	-12.1	119	0.00
31 C	Hexachlorobutadiene	0.206	0.214	-3.9	110	0.00
32	Caprolactam	0.106	0.108	-1.9	108	0.00
33 C	4-Chloro-3-methylphenol	0.343	0.376	-9.6	112	0.00
34	2-Methylnaphthalene	0.769	0.789	-2.6	107	0.00
35	1-Methylnaphthalene	0.737	0.759	-3.0	106	0.00
36 I	Acenaphthene-d10	1.000	1.000	0.0	110	0.00
37	1,2,4,5-Tetrachlorobenzene	0.672	0.662	1.5	109	0.00
38	Hexachlorocyclopentadiene	0.429	0.389	9.3	103	0.00
39 C	2,4,6-Trichlorophenol	0.426	0.437	-2.6	112	0.00
40	2,4,5-Trichlorophenol	0.459	0.476	-3.7	113	0.00
41	1,1'-Biphenyl	1.655	1.625	1.8	107	0.00
42	2-Chloronaphthalene	1.293	1.277	1.2	109	0.00
43	2-Nitroaniline	0.364	0.393	-8.0	118	0.00
44 S	Dimethylphthalate-d6	1.715	1.775	-3.5	113	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110618\
 Data File : BM017530.D
 Acq On : 06 Nov 2018 15:58
 Operator : MJ/SJ
 Sample : SSTDC0020EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02092

Quant Time: Nov 06 16:33:35 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102418MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 06 12:04:23 2018
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	Dimethylphthalate	1.662	1.718	-3.4	113	0.00
46	2,6-Dinitrotoluene	0.334	0.349	-4.5	112	0.00
47 S	Acenaphthylene-d8	2.083	2.126	-2.1	111	0.00
48	Acenaphthylene	1.954	1.982	-1.4	110	0.00
49	3-Nitroaniline	0.314	0.317	-1.0	119	0.00
50 C	Acenaphthene	1.399	1.410	-0.8	111	0.00
51	2,4-Dinitrophenol	0.222	0.214	3.6	119	0.00
52 S	4-Nitrophenol-d4	0.320	0.317	0.9	112	0.00
53	4-Nitrophenol	0.243	0.260	-7.0	117	0.00
54	Dibenzofuran	1.982	2.018	-1.8	111	0.00
55	2,4-Dinitrotoluene	0.498	0.535	-7.4	114	0.00
56	2,3,4,6-Tetrachlorophenol	0.425	0.451	-6.1	115	0.00
57	Diethylphthalate	1.661	1.810	-9.0	118	0.00
58 S	Fluorene-d10	1.490	1.529	-2.6	112	0.00
59	Fluorene	1.634	1.704	-4.3	114	0.00
60	4-Chlorophenyl-phenylether	0.836	0.880	-5.3	114	0.00
61	4-Nitroaniline	0.377	0.345	8.5	98	0.00
62 I	Phenanthrene-d10	1.000	1.000	0.0	113	0.00
63 S	4,6-Dinitro-2-methylphenol-	0.139	0.138	0.7	119	0.00
64	4,6-Dinitro-2-methylphenol	0.142	0.142	0.0	117	0.00
65	N-Nitrosodiphenylamine	0.601	0.612	-1.8	116	0.00
66	4-Bromophenyl-phenylether	0.225	0.228	-1.3	117	0.00
67	Hexachlorobenzene	0.253	0.256	-1.2	115	0.00
68	Atrazine	0.221	0.231	-4.5	120	0.00
69 C	Pentachlorophenol	0.158	0.156	1.3	116	0.00
70	Phenanthrene	1.148	1.150	-0.2	113	0.00
71 S	Anthracene-d10	0.990	1.009	-1.9	115	0.00
72	Anthracene	1.166	1.185	-1.6	114	0.00
73	1,2,3,4-Tetrachlorobenzene	0.282	0.266	5.7	109	0.00
74	Pentachlorobenzene	0.290	0.277	4.5	109	0.00
75	Carbazole	1.020	1.041	-2.1	113	0.00
76	Di-n-butylphthalate	1.192	1.353	-13.5	127	0.00
77 C	Fluoranthene	1.316	1.414	-7.4	115	0.00
78 I	Chrysene-d12	1.000	1.000	0.0	112	0.00
79 S	Pyrene-d10	0.927	0.948	-2.3	115	0.00
80	Pyrene	1.161	1.187	-2.2	114	0.00
81	Butylbenzylphthalate	0.464	0.534	-15.1	128	0.00
82	3,3'-Dichlorobenzidine	0.391	0.413	-5.6	124	0.00
83	Benzo(a)anthracene	1.224	1.247	-1.9	114	0.00
84	Bis(2-ethylhexyl)phthalate	0.678	0.829	-22.3	135	0.00
85	Chrysene	1.164	1.179	-1.3	113	0.00
86 I	Perylene-d12	1.000	1.000	0.0	102	0.00
87	Di-n-octyl phthalate	1.042	1.415	-35.8#	133	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110618\
Data File : BM017530.D
Acq On : 06 Nov 2018 15:58
Operator : MJ/SJ
Sample : SSTDCCC020EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTD02092

Quant Time: Nov 06 16:33:35 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102418MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 06 12:04:23 2018
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.171	1.242	-6.1	104	0.00
89	Benzo(k)fluoranthene	1.144	1.231	-7.6	112	0.00
90 S	Benzo(a)pyrene-d12	1.056	1.097	-3.9	106	0.00
91 C	Benzo(a)pyrene	1.144	1.188	-3.8	105	0.00
92	Indeno(1,2,3-cd)pyrene	1.392	1.301	6.5	96	0.00
93	Dibenzo(a,h)anthracene	1.162	1.100	5.3	96	0.00
94	Benzo(g,h,i)perylene	1.182	1.057	10.6	93	0.00

(#) = Out of Range

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