

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA_M\Data\BM012218\
 Data File : BM013716.D
 Acq On : 23 Jan 2018 04:21
 Operator : SJ/JU
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02034

Quant Time: Jan 23 05:48:38 2018
 Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM011018.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jan 23 00:52:14 2018
 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	122	0.00
2	1,4-Dioxane	0.554	0.431	22.2#	99	0.00
3 S	1,4-Dioxane-d8	0.506	0.415	18.0	101	0.00
4	Benzaldehyde	1.063	1.053	0.9	111	0.00
5 S	Phenol-d5	1.535	1.478	3.7	123	0.00
6	Phenol	1.540	1.500	2.6	121	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	0.930	0.885	4.8	116	0.00
8	Bis(2-Chloroethyl)ether	1.205	1.205	0.0	127	0.00
9 S	2-Chlorophenol-d4	1.251	1.247	0.3	119	0.00
10	2-Chlorophenol	1.247	1.241	0.5	124	0.00
11	2-Methylphenol	1.135	1.118	1.5	124	0.00
12	2,2'-oxybis(1-Chloropropane	1.226	1.135	7.4	114	0.00
13 S	4-Methylphenol-d8	1.180	1.172	0.7	126	0.00
14	Acetophenone	1.953	1.930	1.2	124	0.00
15 P	N-Nitroso-di-n-propylamine	0.953	0.929	2.5	124	0.00
16	4-Methylphenol	1.214	1.198	1.3	123	0.00
17	Hexachloroethane	0.607	0.565	6.9	118	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	121	0.00
19 S	Nitrobenzene-d5	0.155	0.161	-3.9	128	0.00
20	Nitrobenzene	0.401	0.397	1.0	122	0.00
21	Isophorone	0.648	0.635	2.0	121	0.00
22 S	2-Nitrophenol-d4	0.164	0.166	-1.2	125	0.00
23 C	2-Nitrophenol	0.171	0.175	-2.3	127	0.00
24	2,4-Dimethylphenol	0.360	0.361	-0.3	123	0.00
25	Bis(2-Chloroethoxy)methane	0.395	0.380	3.8	115	0.00
26 S	2,4-Dichlorophenol-d3	0.317	0.326	-2.8	124	0.00
27 C	2,4-Dichlorophenol	0.305	0.318	-4.3	127	0.00
28	Naphthalene	0.995	1.043	-4.8	130	0.00
29 S	4-Chloroaniline-d4	0.275	0.335	-21.8#	165#	0.00
30	4-Chloroaniline	0.273	0.326	-19.4	167#	0.00
31 C	Hexachlorobutadiene	0.251	0.247	1.6	122	0.00
32	Caprolactam	0.085	0.086	-1.2	127	0.00
33 C	4-Chloro-3-methylphenol	0.309	0.313	-1.3	127	0.00
34	2-Methylnaphthalene	0.742	0.731	1.5	124	0.00
35 I	Acenaphthene-d10	1.000	1.000	0.0	119	0.00
36	1,2,4,5-Tetrachlorobenzene	0.741	0.751	-1.3	121	0.00
37	Hexachlorocyclopentadiene	0.308	0.234	24.0#	103	0.00
38 C	2,4,6-Trichlorophenol	0.429	0.469	-9.3	128	0.00
39	2,4,5-Trichlorophenol	0.446	0.491	-10.1	129	0.00
40	1,1'-Biphenyl	1.665	1.716	-3.1	122	0.00
41	2-Chloronaphthalene	1.301	1.370	-5.3	123	0.00
42	2-Nitroaniline	0.359	0.375	-4.5	118	0.00
43 S	Dimethylphthalate-d6	1.648	1.642	0.4	119	0.00
44	Dimethylphthalate	1.621	1.629	-0.5	121	0.00

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA_M\Data\BM012218\
 Data File : BM013716.D
 Acq On : 23 Jan 2018 04:21
 Operator : SJ/JU
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02034

Quant Time: Jan 23 05:48:38 2018
 Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM011018.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jan 23 00:52:14 2018
 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	2,6-Dinitrotoluene	0.314	0.326	-3.8	117	0.00
46 S	Acenaphthylene-d8	2.102	2.174	-3.4	123	0.00
47	Acenaphthylene	1.923	1.995	-3.7	121	0.00
48	3-Nitroaniline	0.247	0.282	-14.2	145	0.00
49 C	Acenaphthene	1.332	1.356	-1.8	121	0.00
50	2,4-Dinitrophenol	0.180	0.135	25.0#	103	0.00
51 S	4-Nitrophenol-d4	0.268	0.250	6.7	123	0.00
52	4-Nitrophenol	0.272	0.254	6.6	110	0.00
53	Dibenzofuran	2.004	2.039	-1.7	119	0.00
54	2,4-Dinitrotoluene	0.458	0.462	-0.9	116	0.00
55	2,3,4,6-Tetrachlorophenol	0.420	0.425	-1.2	123	0.00
56	Diethylphthalate	1.631	1.565	4.0	117	0.00
57 S	Fluorene-d10	1.450	1.480	-2.1	120	0.00
58	Fluorene	1.641	1.645	-0.2	119	0.00
59	4-Chlorophenyl-phenylether	0.887	0.848	4.4	113	0.00
60	4-Nitroaniline	0.289	0.289	0.0	114	0.00
61 I	Phenanthrene-d10	1.000	1.000	0.0	112	0.00
62 S	4,6-Dinitro-2-methylphenol-	0.120	0.111	7.5	116	0.00
63	4,6-Dinitro-2-methylphenol	0.127	0.116	8.7	111	0.00
64	N-Nitrosodiphenylamine	0.588	0.612	-4.1	120	0.00
65	4-Bromophenyl-phenylether	0.236	0.242	-2.5	117	0.00
66	Hexachlorobenzene	0.254	0.258	-1.6	116	0.00
67	Atrazine	0.235	0.232	1.3	117	0.00
68 C	Pentachlorophenol	0.135	0.129	4.4	123	0.00
69	Phenanthrene	1.112	1.152	-3.6	118	0.00
70 S	Anthracene-d10	0.974	1.011	-3.8	117	0.00
71	Anthracene	1.147	1.179	-2.8	118	0.00
72	Carbazole	0.947	0.963	-1.7	117	0.00
73	Di-n-butylphthalate	1.113	1.086	2.4	111	0.00
74 C	Fluoranthene	1.327	1.349	-1.7	116	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	110	0.00
76 S	Pyrene-d10	0.931	0.982	-5.5	117	0.00
77	Pyrene	1.216	1.273	-4.7	117	0.00
78	Butylbenzylphthalate	0.428	0.444	-3.7	115	0.00
79	3,3'-Dichlorobenzidine	0.331	0.368	-11.2	129	0.00
80	Benzo(a)anthracene	1.209	1.251	-3.5	116	0.00
81	Bis(2-ethylhexyl)phthalate	0.643	0.617	4.0	106	0.00
82	Chrysene	1.109	1.144	-3.2	114	0.00
83 I	Perylene-d12	1.000	1.000	0.0	114	0.00
84	Di-n-octyl phthalate	1.205	1.048	13.0	107	0.00
85	Benzo(b)fluoranthene	1.228	1.272	-3.6	116	0.00
86	Benzo(k)fluoranthene	1.187	1.255	-5.7	118	0.00
87 S	Benzo(a)pyrene-d12	0.919	0.963	-4.8	119	0.00

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA_M\Data\BM012218\
Data File : BM013716.D
Acq On : 23 Jan 2018 04:21
Operator : SJ/JU
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTD02034

Quant Time: Jan 23 05:48:38 2018
Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM011018.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jan 23 00:52:14 2018
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 C	Benzo(a)pyrene	1.160	1.197	-3.2	116	0.00
89	Indeno(1,2,3-cd)pyrene	1.370	1.384	-1.0	116	0.00
90	Dibenzo(a,h)anthracene	1.157	1.177	-1.7	116	0.00
91	Benzo(g,h,i)perylene	1.127	1.151	-2.1	116	0.00

(#) = Out of Range

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