

Data Path : Z:\HPCHEM1\BNA_G\Data\BG112215\
 Data File : BG019802.D
 Acq On : 23 Nov 2015 4:25
 Operator : UM/NP
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTD02012

Quant Time: Nov 23 06:08:10 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG111615.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 23 04:11:25 2015
 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	108	0.00
2	1,4-Dioxane	0.534	0.466	12.7	121	0.00
3 S	1,4-Dioxane-d8	0.468	0.388	17.1	99	0.00
4	Benzaldehyde	1.493	1.563	-4.7	101	0.00
5 S	Phenol-d5	2.049	2.101	-2.5	118	0.00
6	Phenol	2.212	2.275	-2.8	119	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	1.342	1.218	9.2	99	0.00
8	Bis(2-Chloroethyl)ether	1.540	1.582	-2.7	112	0.00
9 S	2-Chlorophenol-d4	1.314	1.379	-4.9	120	0.00
10	2-Chlorophenol	1.284	1.330	-3.6	121	0.00
11	2-Methylphenol	1.629	1.733	-6.4	119	0.00
12	2,2'-oxybis(1-Chloropropane	1.345	1.305	3.0	106	0.00
13 S	4-Methylphenol-d8	1.696	1.800	-6.1	121	0.00
14	Acetophenone	2.803	2.895	-3.3	116	0.00
15 P	N-Nitroso-di-n-propylamine	1.894	1.784	5.8	110	0.00
16	4-Methylphenol	1.835	1.878	-2.3	118	0.00
17	Hexachloroethane	0.610	0.646	-5.9	123	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	112	0.00
19 S	Nitrobenzene-d5	0.164	0.170	-3.7	123	0.00
20	Nitrobenzene	0.591	0.574	2.9	114	0.00
21	Isophorone	1.141	1.102	3.4	115	0.00
22 S	2-Nitrophenol-d4	0.185	0.192	-3.8	115	0.00
23 C	2-Nitrophenol	0.200	0.207	-3.5	124	0.00
24	2,4-Dimethylphenol	0.533	0.570	-6.9	128	0.00
25	Bis(2-Chloroethoxy)methane	0.565	0.540	4.4	113	0.00
26 S	2,4-Dichlorophenol-d3	0.400	0.448	-12.0	133	0.00
27 C	2,4-Dichlorophenol	0.381	0.430	-12.9	134	0.00
28	Naphthalene	1.080	1.080	0.0	118	0.00
29 S	4-Chloroaniline-d4	0.379	0.401	-5.8	111	0.00
30	4-Chloroaniline	0.395	0.410	-3.8	108	0.00
31 C	Hexachlorobutadiene	0.343	0.392	-14.3	133	0.00
32	Caprolactam	0.155	0.155	0.0	114	0.00
33 C	4-Chloro-3-methylphenol	0.531	0.559	-5.3	126	0.00
34	2-Methylnaphthalene	0.862	0.871	-1.0	116	0.00
35	1-Methylnaphthalene	0.858	0.880	-2.6	123	0.00
36 I	Acenaphthene-d10	1.000	1.000	0.0	119	0.00
37	1,2,4,5-Tetrachlorobenzene	0.802	0.878	-9.5	134	0.00
38	Hexachlorocyclopentadiene	0.467	0.286	38.8#	76	0.00
39 C	2,4,6-Trichlorophenol	0.489	0.547	-11.9	138	0.00
40	2,4,5-Trichlorophenol	0.512	0.573	-11.9	139	0.00
41	1,1'-Biphenyl	1.488	1.496	-0.5	124	0.00
42	2-Chloronaphthalene	1.156	1.166	-0.9	122	0.00
43	2-Nitroaniline	0.489	0.464	5.1	114	0.00
44 S	Dimethylphthalate-d6	1.809	1.791	1.0	121	0.00

Data Path : Z:\HPCHEM1\BNA_G\Data\BG112215\
 Data File : BG019802.D
 Acq On : 23 Nov 2015 4:25
 Operator : UM/NP
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTD02012

Quant Time: Nov 23 06:08:10 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG111615.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 23 04:11:25 2015
 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	1.794	1.778	0.9	120	0.00
46	2,6-Dinitrotoluene	0.359	0.383	-6.7	133	0.00
47 S	Acenaphthylene-d8	1.961	1.978	-0.9	124	0.00
48	Acenaphthylene	1.957	1.941	0.8	120	0.00
49	3-Nitroaniline	0.299	0.296	1.0	106	0.00
50 C	Acenaphthene	1.340	1.332	0.6	122	0.00
51	2,4-Dinitrophenol	0.236	0.189	19.9	103	0.00
52 S	4-Nitrophenol-d4	0.285	0.264	7.4	111	0.00
53	4-Nitrophenol	0.347	0.314	9.5	106	0.00
54	Dibenzofuran	1.964	2.021	-2.9	128	0.00
55	2,4-Dinitrotoluene	0.558	0.548	1.8	118	0.00
56	2,3,4,6-Tetrachlorophenol	0.543	0.604	-11.2	134	0.00
57	Diethylphthalate	1.792	1.723	3.9	118	0.00
58 S	Fluorene-d10	1.619	1.620	-0.1	123	0.00
59	Fluorene	1.717	1.713	0.2	124	0.00
60	4-Chlorophenyl-phenylether	0.999	1.035	-3.6	129	0.00
61	4-Nitroaniline	0.360	0.346	3.9	105	0.00
62 I	Phenanthrene-d10	1.000	1.000	0.0	110	0.00
63 S	4,6-Dinitro-2-methylphenol-	0.129	0.132	-2.3	117	0.00
64	4,6-Dinitro-2-methylphenol	0.137	0.139	-1.5	113	0.00
65	N-Nitrosodiphenylamine	0.539	0.551	-2.2	117	0.00
66	4-Bromophenyl-phenylether	0.240	0.263	-9.6	125	0.00
67	Hexachlorobenzene	0.254	0.278	-9.4	125	0.00
68	Atrazine	0.251	0.266	-6.0	118	0.00
69 C	Pentachlorophenol	0.142	0.147	-3.5	123	0.00
70	Phenanthrene	1.070	1.076	-0.6	116	0.00
71 S	Anthracene-d10	0.971	0.994	-2.4	117	0.00
72	Anthracene	1.079	1.085	-0.6	114	0.00
73	1,2,3,4-Tetrachlorobenzene	0.276	0.331	-19.9	138	0.00
74	Pentachlorobenzene	0.288	0.342	-18.8	136	0.00
75	Carbazole	0.867	0.875	-0.9	107	0.00
76	Di-n-butylphthalate	1.124	1.046	6.9	104	0.00
77 C	Fluoranthene	1.364	1.370	-0.4	104	0.00
78 I	Chrysene-d12	1.000	1.000	0.0	103	0.00
79 S	Pyrene-d10	0.911	0.914	-0.3	108	0.00
80	Pyrene	1.168	1.139	2.5	104	0.00
81	Butylbenzylphthalate	0.384	0.381	0.8	102	0.00
82	3,3'-Dichlorobenzidine	0.366	0.371	-1.4	95	0.00
83	Benzo(a)anthracene	1.212	1.240	-2.3	107	0.00
84	Bis(2-ethylhexyl)phthalate	0.563	0.560	0.5	103	0.00
85	Chrysene	1.141	1.144	-0.3	106	0.00
86 I	Perylene-d12	1.000	1.000	0.0	105	0.00
87	Di-n-octyl phthalate	0.971	0.979	-0.8	103	0.00

Data Path : Z:\HPCHEM1\BNA_G\Data\BG112215\
Data File : BG019802.D
Acq On : 23 Nov 2015 4:25
Operator : UM/NP
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTD02012

Quant Time: Nov 23 06:08:10 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG111615.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Nov 23 04:11:25 2015
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	1.232	1.212	1.6	105	0.00
89	Benzo(k)fluoranthene	1.189	1.192	-0.3	109	0.00
90 S	Benzo(a)pyrene-d12	1.044	1.068	-2.3	110	0.00
91 C	Benzo(a)pyrene	1.186	1.186	0.0	109	0.00
92	Indeno(1,2,3-cd)pyrene	1.328	1.355	-2.0	111	0.00
93	Dibenzo(a,h)anthracene	1.091	1.118	-2.5	109	0.00
94	Benzo(g,h,i)perylene	1.102	1.136	-3.1	111	0.00

(#) = Out of Range

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