

Data Path : Z:\HPCHEM1\BNA_G\Data\BG080916\
 Data File : BG023555.D
 Acq On : 9 Aug 2016 12:24
 Operator : UM/SJ
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTD02018

Quant Time: Aug 10 04:59:42 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080416.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 09 02:15:40 2016
 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	120	0.00
2	1,4-Dioxane	0.494	0.532	-7.7	113	0.00
3 S	1,4-Dioxane-d8	0.470	0.449	4.5	102	0.00
4	Benzaldehyde	1.228	1.356	-10.4	115	0.00
5 S	Phenol-d5	2.015	2.070	-2.7	115	0.00
6	Phenol	2.098	2.177	-3.8	117	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	1.291	1.328	-2.9	117	0.00
8	Bis(2-Chloroethyl)ether	1.523	1.581	-3.8	116	0.00
9 S	2-Chlorophenol-d4	1.374	1.464	-6.6	120	0.00
10	2-Chlorophenol	1.400	1.458	-4.1	117	0.00
11	2-Methylphenol	1.580	1.629	-3.1	118	0.00
12	2,2'-oxybis(1-Chloropropane	2.084	2.270	-8.9	123	0.00
13 S	4-Methylphenol-d8	1.573	1.619	-2.9	119	0.00
14	Acetophenone	2.523	2.695	-6.8	117	0.00
15 P	N-Nitroso-di-n-propylamine	1.518	1.619	-6.7	119	0.00
16	4-Methylphenol	1.771	1.878	-6.0	120	0.00
17	Hexachloroethane	0.604	0.589	2.5	113	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	124	0.00
19 S	Nitrobenzene-d5	0.158	0.162	-2.5	119	0.00
20	Nitrobenzene	0.462	0.459	0.6	115	0.00
21	Isophorone	0.850	0.862	-1.4	114	0.00
22 S	2-Nitrophenol-d4	0.181	0.179	1.1	113	0.00
23 C	2-Nitrophenol	0.195	0.197	-1.0	120	0.00
24	2,4-Dimethylphenol	0.429	0.430	-0.2	115	0.00
25	Bis(2-Chloroethoxy)methane	0.490	0.492	-0.4	116	0.00
26 S	2,4-Dichlorophenol-d3	0.342	0.358	-4.7	120	0.00
27 C	2,4-Dichlorophenol	0.344	0.354	-2.9	119	0.00
28	Naphthalene	1.015	1.046	-3.1	117	0.00
29 S	4-Chloroaniline-d4	0.401	0.459	-14.5	121	0.00
30	4-Chloroaniline	0.398	0.450	-13.1	122	0.00
31 C	Hexachlorobutadiene	0.229	0.229	0.0	117	0.00
32	Caprolactam	0.143	0.138	3.5	109	0.00
33 C	4-Chloro-3-methylphenol	0.443	0.443	0.0	116	0.00
34	2-Methylnaphthalene	0.823	0.839	-1.9	116	0.00
35 I	Acenaphthene-d10	1.000	1.000	0.0	115	0.00
36	1,2,4,5-Tetrachlorobenzene	0.614	0.648	-5.5	114	0.00
37	Hexachlorocyclopentadiene	0.340	0.327	3.8	113	0.00
38 C	2,4,6-Trichlorophenol	0.425	0.451	-6.1	113	0.00
39	2,4,5-Trichlorophenol	0.456	0.476	-4.4	115	0.00
40	1,1'-Biphenyl	1.470	1.575	-7.1	116	0.00
41	2-Chloronaphthalene	1.143	1.216	-6.4	116	0.00
42	2-Nitroaniline	0.467	0.477	-2.1	108	0.00
43 S	Dimethylphthalate-d6	1.638	1.669	-1.9	110	0.00
44	Dimethylphthalate	1.622	1.640	-1.1	108	0.00

Data Path : Z:\HPCHEM1\BNA_G\Data\BG080916\
 Data File : BG023555.D
 Acq On : 9 Aug 2016 12:24
 Operator : UM/SJ
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTD02018

Quant Time: Aug 10 04:59:42 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080416.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 09 02:15:40 2016
 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	0.352	0.352	0.0	106	0.00
46 S	Acenaphthylene-d8	1.843	1.924	-4.4	112	0.00
47	Acenaphthylene	1.928	2.037	-5.7	113	0.00
48	3-Nitroaniline	0.331	0.363	-9.7	117	0.00
49 C	Acenaphthene	1.313	1.363	-3.8	112	0.00
50	2,4-Dinitrophenol	0.217	0.198	8.8	109	0.00
51 S	4-Nitrophenol-d4	0.280	0.267	4.6	108	0.00
52	4-Nitrophenol	0.280	0.237	15.4	91	0.00
53	Dibenzofuran	1.946	2.017	-3.6	112	0.00
54	2,4-Dinitrotoluene	0.533	0.556	-4.3	110	0.00
55	2,3,4,6-Tetrachlorophenol	0.468	0.459	1.9	106	0.00
56	Diethylphthalate	1.695	1.728	-1.9	108	0.00
57 S	Fluorene-d10	1.513	1.558	-3.0	111	0.00
58	Fluorene	1.621	1.669	-3.0	111	0.00
59	4-Chlorophenyl-phenylether	0.815	0.830	-1.8	110	0.00
60	4-Nitroaniline	0.383	0.384	-0.3	105	0.00
61 I	Phenanthrene-d10	1.000	1.000	0.0	109	0.00
62 S	4,6-Dinitro-2-methylphenol-	0.128	0.130	-1.6	108	0.00
63	4,6-Dinitro-2-methylphenol	0.135	0.136	-0.7	106	0.00
64	N-Nitrosodiphenylamine	0.559	0.599	-7.2	107	0.00
65	4-Bromophenyl-phenylether	0.229	0.235	-2.6	106	0.00
66	Hexachlorobenzene	0.244	0.251	-2.9	109	0.00
67	Atrazine	0.231	0.240	-3.9	102	0.00
68 C	Pentachlorophenol	0.135	0.128	5.2	103	0.00
69	Phenanthrene	1.036	1.073	-3.6	103	0.00
70 S	Anthracene-d10	0.901	0.941	-4.4	106	0.00
71	Anthracene	1.057	1.120	-6.0	105	0.00
72	Carbazole	0.852	0.954	-12.0	100	0.00
73	Di-n-butylphthalate	1.067	1.137	-6.6	103	0.00
74 C	Fluoranthene	1.060	1.219	-15.0	100	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	102	0.00
76 S	Pyrene-d10	1.012	1.058	-4.5	101	0.00
77	Pyrene	1.236	1.299	-5.1	99	0.00
78	Butylbenzylphthalate	0.474	0.507	-7.0	102	0.00
79	3,3'-Dichlorobenzidine	0.352	0.376	-6.8	99	0.00
80	Benzo(a)anthracene	1.126	1.176	-4.4	98	0.00
81	Bis(2-ethylhexyl)phthalate	0.631	0.686	-8.7	101	0.00
82	Chrysene	1.024	1.072	-4.7	100	0.00
83 I	Perylene-d12	1.000	1.000	0.0	104	0.00
84	Di-n-octyl phthalate	1.010	1.193	-18.1	104	0.00
85	Benzo(b)fluoranthene	1.138	1.216	-6.9	101	0.00
86	Benzo(k)fluoranthene	1.120	1.160	-3.6	102	0.00
87 S	Benzo(a)pyrene-d12	0.905	0.937	-3.5	101	0.00

Data Path : Z:\HPCHEM1\BNA_G\Data\BG080916\
Data File : BG023555.D
Acq On : 9 Aug 2016 12:24
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTD02018

Quant Time: Aug 10 04:59:42 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080416.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Aug 09 02:15:40 2016
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	1.098	1.131	-3.0	98	0.00
89	Indeno(1,2,3-cd)pyrene	1.291	1.309	-1.4	99	0.00
90	Dibenzo(a,h)anthracene	1.101	1.113	-1.1	98	0.00
91	Benzo(g,h,i)perylene	1.070	1.077	-0.7	98	0.00

(#) = Out of Range

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