

Data Path : Z:\HPCHEM1\BNA_G\Data\BG082016\
 Data File : BG023818.D
 Acq On : 20 Aug 2016 15:19
 Operator : UM/SJ
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTD02017

Quant Time: Aug 22 04:54:13 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG081816.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Aug 22 03:42:09 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	110	0.00
2	1,4-Dioxane	0.462	0.495	-7.1	109	0.00
3 S	1,4-Dioxane-d8	0.441	0.459	-4.1	114	0.00
4	Benzaldehyde	1.210	1.367	-13.0	109	0.00
5 S	Phenol-d5	1.875	1.968	-5.0	110	0.00
6	Phenol	1.921	2.075	-8.0	113	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	1.285	1.355	-5.4	109	0.00
8	Bis(2-Chloroethyl)ether	1.466	1.584	-8.0	113	0.00
9 S	2-Chlorophenol-d4	1.331	1.408	-5.8	113	0.00
10	2-Chlorophenol	1.358	1.441	-6.1	112	0.00
11	2-Methylphenol	1.461	1.550	-6.1	112	0.00
12	2,2'-oxybis(1-Chloropropane	2.096	2.199	-4.9	107	0.00
13 S	4-Methylphenol-d8	1.466	1.574	-7.4	114	0.00
14	Acetophenone	2.384	2.521	-5.7	110	0.00
15 P	N-Nitroso-di-n-propylamine	1.477	1.501	-1.6	107	0.00
16	4-Methylphenol	1.589	1.641	-3.3	109	0.00
17	Hexachloroethane	0.599	0.646	-7.8	114	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	109	0.00
19 S	Nitrobenzene-d5	0.162	0.163	-0.6	109	0.00
20	Nitrobenzene	0.473	0.476	-0.6	110	0.00
21	Isophorone	0.875	0.870	0.6	107	0.00
22 S	2-Nitrophenol-d4	0.184	0.191	-3.8	108	0.00
23 C	2-Nitrophenol	0.198	0.207	-4.5	111	0.00
24	2,4-Dimethylphenol	0.428	0.441	-3.0	108	0.00
25	Bis(2-Chloroethoxy)methane	0.492	0.501	-1.8	109	0.00
26 S	2,4-Dichlorophenol-d3	0.333	0.345	-3.6	110	0.00
27 C	2,4-Dichlorophenol	0.333	0.349	-4.8	113	0.00
28	Naphthalene	1.023	1.045	-2.2	108	0.00
29 S	4-Chloroaniline-d4	0.408	0.439	-7.6	109	0.00
30	4-Chloroaniline	0.400	0.429	-7.2	108	0.00
31 C	Hexachlorobutadiene	0.239	0.256	-7.1	115	0.00
32	Caprolactam	0.130	0.110	15.4	92	0.00
33 C	4-Chloro-3-methylphenol	0.411	0.405	1.5	104	0.00
34	2-Methylnaphthalene	0.779	0.791	-1.5	107	0.00
35 I	Acenaphthene-d10	1.000	1.000	0.0	103	0.00
36	1,2,4,5-Tetrachlorobenzene	0.664	0.727	-9.5	112	0.00
37	Hexachlorocyclopentadiene	0.290	0.199	31.4#	83	0.00
38 C	2,4,6-Trichlorophenol	0.437	0.456	-4.3	108	0.00
39	2,4,5-Trichlorophenol	0.452	0.461	-2.0	103	0.00
40	1,1'-Biphenyl	1.564	1.621	-3.6	104	0.00
41	2-Chloronaphthalene	1.189	1.238	-4.1	105	0.00
42	2-Nitroaniline	0.493	0.474	3.9	97	0.00
43 S	Dimethylphthalate-d6	1.658	1.648	0.6	100	0.00
44	Dimethylphthalate	1.658	1.634	1.4	99	0.00

Data Path : Z:\HPCHEM1\BNA_G\Data\BG082016\
 Data File : BG023818.D
 Acq On : 20 Aug 2016 15:19
 Operator : UM/SJ
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTD02017

Quant Time: Aug 22 04:54:13 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG081816.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Aug 22 03:42:09 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	0.354	0.351	0.8	101	0.00
46 S	Acenaphthylene-d8	1.878	1.951	-3.9	104	0.00
47	Acenaphthylene	1.956	2.032	-3.9	103	0.00
48	3-Nitroaniline	0.333	0.342	-2.7	101	0.00
49 C	Acenaphthene	1.310	1.360	-3.8	106	0.00
50	2,4-Dinitrophenol	0.188	0.166	11.7	105	0.00
51 S	4-Nitrophenol-d4	0.268	0.241	10.1	94	0.00
52	4-Nitrophenol	0.269	0.292	-8.6	113	0.00
53	Dibenzofuran	1.905	1.954	-2.6	102	0.00
54	2,4-Dinitrotoluene	0.517	0.511	1.2	101	0.00
55	2,3,4,6-Tetrachlorophenol	0.417	0.424	-1.7	102	0.00
56	Diethylphthalate	1.694	1.652	2.5	97	0.00
57 S	Fluorene-d10	1.448	1.462	-1.0	102	0.00
58	Fluorene	1.543	1.578	-2.3	101	0.00
59	4-Chlorophenyl-phenylether	0.780	0.807	-3.5	106	0.00
60	4-Nitroaniline	0.370	0.352	4.9	94	0.00
61 I	Phenanthrene-d10	1.000	1.000	0.0	101	0.00
62 S	4,6-Dinitro-2-methylphenol-	0.127	0.132	-3.9	105	0.00
63	4,6-Dinitro-2-methylphenol	0.132	0.137	-3.8	105	0.00
64	N-Nitrosodiphenylamine	0.592	0.605	-2.2	98	0.00
65	4-Bromophenyl-phenylether	0.230	0.235	-2.2	99	0.00
66	Hexachlorobenzene	0.248	0.255	-2.8	101	0.00
67	Atrazine	0.243	0.254	-4.5	99	0.00
68 C	Pentachlorophenol	0.108	0.113	-4.6	118	0.00
69	Phenanthrene	1.059	1.093	-3.2	98	0.00
70 S	Anthracene-d10	0.922	0.941	-2.1	97	0.00
71	Anthracene	1.089	1.136	-4.3	100	0.00
72	Carbazole	0.888	0.988	-11.3	97	0.00
73	Di-n-butylphthalate	1.202	1.215	-1.1	95	0.00
74 C	Fluoranthene	1.128	1.316	-16.7	98	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	99	0.00
76 S	Pyrene-d10	0.960	1.014	-5.6	100	0.00
77	Pyrene	1.182	1.246	-5.4	97	0.00
78	Butylbenzylphthalate	0.535	0.530	0.9	95	0.00
79	3,3'-Dichlorobenzidine	0.397	0.421	-6.0	95	0.00
80	Benzo(a)anthracene	1.156	1.203	-4.1	99	0.00
81	Bis(2-ethylhexyl)phthalate	0.719	0.718	0.1	95	0.00
82	Chrysene	1.041	1.101	-5.8	100	0.00
83 I	Perylene-d12	1.000	1.000	0.0	102	0.00
84	Di-n-octyl phthalate	1.150	1.256	-9.2	97	0.00
85	Benzo(b)fluoranthene	1.163	1.222	-5.1	103	0.00
86	Benzo(k)fluoranthene	1.123	1.153	-2.7	101	0.00
87 S	Benzo(a)pyrene-d12	0.916	0.952	-3.9	104	0.00

Data Path : Z:\HPCHEM1\BNA_G\Data\BG082016\
Data File : BG023818.D
Acq On : 20 Aug 2016 15:19
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTD02017

Quant Time: Aug 22 04:54:13 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG081816.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 22 03:42:09 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	1.116	1.152	-3.2	102	0.00
89	Indeno(1,2,3-cd)pyrene	1.312	1.342	-2.3	101	0.00
90	Dibenzo(a,h)anthracene	1.125	1.149	-2.1	101	0.00
91	Benzo(g,h,i)perylene	1.085	1.114	-2.7	102	0.00

(#) = Out of Range

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