

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG050917\
 Data File : BG026933.D
 Acq On : 10 May 2017 4:51
 Operator : SJ/MA
 Sample : SSTDC0020EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTD02082

Quant Time: May 10 05:25:39 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG042717.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed May 10 01:49:01 2017
 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	124	0.00
2	1,4-Dioxane	0.507	0.410	19.1	107	0.00
3 S	1,4-Dioxane-d8	0.444	0.374	15.8	105	0.00
4	Benzaldehyde	0.863	0.824	4.5	144	0.00
5 S	Phenol-d5	1.785	1.911	-7.1	133	0.00
6	Phenol	1.809	1.979	-9.4	135	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	1.050	1.136	-8.2	134	0.00
8	Bis(2-Chloroethyl)ether	1.397	1.403	-0.4	121	0.00
9 S	2-Chlorophenol-d4	1.515	1.544	-1.9	129	0.00
10	2-Chlorophenol	1.491	1.502	-0.7	124	0.00
11	2-Methylphenol	1.432	1.515	-5.8	133	0.00
12	2,2'-oxybis(1-Chloropropane	1.458	1.744	-19.6	147	0.00
13 S	4-Methylphenol-d8	1.481	1.584	-7.0	134	0.00
14	Acetophenone	2.176	2.250	-3.4	125	0.00
15 P	N-Nitroso-di-n-propylamine	1.073	1.065	0.7	125	0.00
16	4-Methylphenol	1.566	1.641	-4.8	130	0.00
17	Hexachloroethane	0.561	0.561	0.0	131	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	127	0.00
19 S	Nitrobenzene-d5	0.167	0.168	-0.6	130	0.00
20	Nitrobenzene	0.323	0.339	-5.0	133	0.00
21	Isophorone	0.619	0.624	-0.8	127	0.00
22 S	2-Nitrophenol-d4	0.182	0.184	-1.1	127	0.00
23 C	2-Nitrophenol	0.191	0.189	1.0	127	0.00
24	2,4-Dimethylphenol	0.349	0.359	-2.9	133	0.00
25	Bis(2-Chloroethoxy)methane	0.423	0.418	1.2	125	0.00
26 S	2,4-Dichlorophenol-d3	0.289	0.300	-3.8	131	0.00
27 C	2,4-Dichlorophenol	0.288	0.292	-1.4	128	0.00
28	Naphthalene	1.041	1.037	0.4	126	0.00
29 S	4-Chloroaniline-d4	0.387	0.426	-10.1	127	0.00
30	4-Chloroaniline	0.365	0.401	-9.9	129	0.00
31 C	Hexachlorobutadiene	0.146	0.142	2.7	125	0.00
32	Caprolactam	0.108	0.108	0.0	131	0.00
33 C	4-Chloro-3-methylphenol	0.312	0.340	-9.0	135	0.00
34	2-Methylnaphthalene	0.773	0.752	2.7	124	0.00
35 I	Acenaphthene-d10	1.000	1.000	0.0	121	0.00
36	1,2,4,5-Tetrachlorobenzene	0.527	0.522	0.9	122	0.00
37	Hexachlorocyclopentadiene	0.232	0.119	48.7#	68	0.00
38 C	2,4,6-Trichlorophenol	0.363	0.377	-3.9	125	0.00
39	2,4,5-Trichlorophenol	0.375	0.392	-4.5	125	0.00
40	1,1'-Biphenyl	1.681	1.674	0.4	121	0.00
41	2-Chloronaphthalene	1.264	1.272	-0.6	123	0.00
42	2-Nitroaniline	0.361	0.418	-15.8	140	0.00
43 S	Dimethylphthalate-d6	1.610	1.540	4.3	115	0.00
44	Dimethylphthalate	1.576	1.504	4.6	116	0.00

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG050917\
 Data File : BG026933.D
 Acq On : 10 May 2017 4:51
 Operator : SJ/MA
 Sample : SSTDC0020EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTD02082

Quant Time: May 10 05:25:39 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG042717.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed May 10 01:49:01 2017
 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	0.333	0.336	-0.9	122	0.00
46 S	Acenaphthylene-d8	2.022	2.054	-1.6	122	0.00
47	Acenaphthylene	2.060	2.079	-0.9	122	0.00
48	3-Nitroaniline	0.350	0.376	-7.4	133	0.00
49 C	Acenaphthene	1.433	1.425	0.6	121	0.00
50	2,4-Dinitrophenol	0.170	0.130	23.5#	84	0.00
51 S	4-Nitrophenol-d4	0.290	0.248	14.5	108	0.00
52	4-Nitrophenol	0.170	0.151	11.2	109	0.00
53	Dibenzofuran	1.907	1.848	3.1	117	0.00
54	2,4-Dinitrotoluene	0.476	0.465	2.3	117	0.00
55	2,3,4,6-Tetrachlorophenol	0.295	0.276	6.4	111	0.00
56	Diethylphthalate	1.649	1.587	3.8	116	0.00
57 S	Fluorene-d10	1.405	1.345	4.3	117	0.00
58	Fluorene	1.557	1.489	4.4	116	0.00
59	4-Chlorophenyl-phenylether	0.679	0.631	7.1	114	0.00
60	4-Nitroaniline	0.386	0.382	1.0	130	0.00
61 I	Phenanthrene-d10	1.000	1.000	0.0	110	0.00
62 S	4,6-Dinitro-2-methylphenol-	0.121	0.094	22.3#	90	0.00
63	4,6-Dinitro-2-methylphenol	0.126	0.097	23.0#	88	0.00
64	N-Nitrosodiphenylamine	0.657	0.682	-3.8	114	0.00
65	4-Bromophenyl-phenylether	0.200	0.201	-0.5	110	0.00
66	Hexachlorobenzene	0.215	0.206	4.2	108	0.00
67	Atrazine	0.205	0.214	-4.4	112	0.00
68 C	Pentachlorophenol	0.098	0.091	7.1	120	0.00
69	Phenanthrene	1.127	1.124	0.3	109	0.00
70 S	Anthracene-d10	0.971	0.972	-0.1	110	0.00
71	Anthracene	1.156	1.163	-0.6	109	0.00
72	Carbazole	0.977	1.070	-9.5	110	0.00
73	Di-n-butylphthalate	1.272	1.370	-7.7	114	0.00
74 C	Fluoranthene	1.046	1.136	-8.6	108	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	109	0.00
76 S	Pyrene-d10	1.056	1.042	1.3	107	0.00
77	Pyrene	1.342	1.331	0.8	107	0.00
78	Butylbenzylphthalate	0.623	0.723	-16.1	125	0.00
79	3,3'-Dichlorobenzidine	0.385	0.383	0.5	106	0.00
80	Benzo(a)anthracene	1.170	1.184	-1.2	111	0.00
81	Bis(2-ethylhexyl)phthalate	0.889	1.034	-16.3	124	0.00
82	Chrysene	1.087	1.088	-0.1	110	0.00
83 I	Perylene-d12	1.000	1.000	0.0	114	0.00
84	Di-n-octyl phthalate	1.374	1.653	-20.3#	127	0.00
85	Benzo(b)fluoranthene	1.143	1.131	1.0	116	0.00
86	Benzo(k)fluoranthene	1.123	1.119	0.4	113	0.00
87 S	Benzo(a)pyrene-d12	0.946	0.954	-0.8	116	0.00

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG050917\
Data File : BG026933.D
Acq On : 10 May 2017 4:51
Operator : SJ/MA
Sample : SSTDCCC020EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTD02082

Quant Time: May 10 05:25:39 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG042717.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed May 10 01:49:01 2017
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	1.125	1.118	0.6	114	0.00
89	Indeno(1,2,3-cd)pyrene	1.333	1.218	8.6	106	0.01
90	Dibenzo(a,h)anthracene	1.130	1.057	6.5	108	0.00
91	Benzo(g,h,i)perylene	1.106	0.872	21.2#	92	0.00

(#) = Out of Range

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