

Data Path : Z:\HPCHEM1\BNA G\DATA\BG051617\
 Data File : BG026995.D
 Acq On : 17 May 2017 2:08
 Operator : SJ/MA
 Sample : SSTDC0020
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTD02091

Quant Time: May 17 05:11:13 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG042717.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed May 17 04:07:33 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	142	0.00
2	1,4-Dioxane	0.507	0.492	3.0	147	0.00
3 S	1,4-Dioxane-d8	0.444	0.438	1.4	141	0.00
4	Benzaldehyde	0.863	0.922	-6.8	185#	0.00
5 S	Phenol-d5	1.785	1.981	-11.0	158#	0.00
6	Phenol	1.809	2.021	-11.7	158#	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	1.050	1.122	-6.9	151#	0.00
8	Bis(2-Chloroethyl)ether	1.397	1.453	-4.0	144	0.00
9 S	2-Chlorophenol-d4	1.515	1.609	-6.2	154#	0.00
10	2-Chlorophenol	1.491	1.545	-3.6	146	0.00
11	2-Methylphenol	1.432	1.559	-8.9	156#	0.00
12	2,2'-oxybis(1-Chloropropane	1.458	1.614	-10.7	156#	0.00
13 S	4-Methylphenol-d8	1.481	1.588	-7.2	154#	0.00
14	Acetophenone	2.176	2.263	-4.0	144	0.00
15 P	N-Nitroso-di-n-propylamine	1.073	1.059	1.3	142	0.00
16	4-Methylphenol	1.566	1.706	-8.9	155#	0.00
17	Hexachloroethane	0.561	0.562	-0.2	150	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	144	0.00
19 S	Nitrobenzene-d5	0.167	0.173	-3.6	152#	0.00
20	Nitrobenzene	0.323	0.343	-6.2	152#	0.00
21	Isophorone	0.619	0.633	-2.3	146	0.00
22 S	2-Nitrophenol-d4	0.182	0.193	-6.0	151#	0.00
23 C	2-Nitrophenol	0.191	0.201	-5.2	153#	0.00
24	2,4-Dimethylphenol	0.349	0.372	-6.6	156#	0.00
25	Bis(2-Chloroethoxy)methane	0.423	0.428	-1.2	144	0.00
26 S	2,4-Dichlorophenol-d3	0.289	0.313	-8.3	154#	0.00
27 C	2,4-Dichlorophenol	0.288	0.311	-8.0	154#	0.00
28	Naphthalene	1.041	1.058	-1.6	145	0.00
29 S	4-Chloroaniline-d4	0.387	0.421	-8.8	142	0.00
30	4-Chloroaniline	0.365	0.401	-9.9	146	0.00
31 C	Hexachlorobutadiene	0.146	0.144	1.4	144	0.00
32	Caprolactam	0.108	0.113	-4.6	155#	0.00
33 C	4-Chloro-3-methylphenol	0.312	0.360	-15.4	162#	0.00
34	2-Methylnaphthalene	0.773	0.778	-0.6	145	0.00
35 I	Acenaphthene-d10	1.000	1.000	0.0	142	0.00
36	1,2,4,5-Tetrachlorobenzene	0.527	0.537	-1.9	146	0.00
37	Hexachlorocyclopentadiene	0.232	0.268	-15.5	180#	0.00
38 C	2,4,6-Trichlorophenol	0.363	0.390	-7.4	151#	0.00
39	2,4,5-Trichlorophenol	0.375	0.397	-5.9	148	0.00
40	1,1'-Biphenyl	1.681	1.702	-1.2	143	0.00
41	2-Chloronaphthalene	1.264	1.299	-2.8	147	0.00
42	2-Nitroaniline	0.361	0.417	-15.5	163#	0.00
43 S	Dimethylphthalate-d6	1.610	1.602	0.5	140	0.00
44	Dimethylphthalate	1.576	1.564	0.8	141	0.00

Data Path : Z:\HPCHEM1\BNA G\DATA\BG051617\
 Data File : BG026995.D
 Acq On : 17 May 2017 2:08
 Operator : SJ/MA
 Sample : SSTDC0020
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTD02091

Quant Time: May 17 05:11:13 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG042717.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed May 17 04:07:33 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
45	2,6-Dinitrotoluene	0.333	0.350	-5.1	148	0.00
46 S	Acenaphthylene-d8	2.022	2.125	-5.1	147	0.00
47	Acenaphthylene	2.060	2.118	-2.8	145	0.00
48	3-Nitroaniline	0.350	0.387	-10.6	159#	0.00
49 C	Acenaphthene	1.433	1.448	-1.0	144	0.00
50	2,4-Dinitrophenol	0.170	0.231	-35.9#	174#	0.00
51 S	4-Nitrophenol-d4	0.290	0.246	15.2	124	0.00
52	4-Nitrophenol	0.170	0.142	16.5	119	0.00
53	Dibenzofuran	1.907	1.905	0.1	141	0.00
54	2,4-Dinitrotoluene	0.476	0.487	-2.3	142	0.00
55	2,3,4,6-Tetrachlorophenol	0.295	0.288	2.4	135	0.00
56	Diethylphthalate	1.649	1.634	0.9	140	0.00
57 S	Fluorene-d10	1.405	1.408	-0.2	142	0.00
58	Fluorene	1.557	1.522	2.2	138	0.00
59	4-Chlorophenyl-phenylether	0.679	0.664	2.2	140	0.00
60	4-Nitroaniline	0.386	0.389	-0.8	154#	0.00
61 I	Phenanthrene-d10	1.000	1.000	0.0	131	0.00
62 S	4,6-Dinitro-2-methylphenol-	0.121	0.119	1.7	135	0.00
63	4,6-Dinitro-2-methylphenol	0.126	0.123	2.4	133	0.00
64	N-Nitrosodiphenylamine	0.657	0.699	-6.4	139	0.00
65	4-Bromophenyl-phenylether	0.200	0.216	-8.0	141	0.00
66	Hexachlorobenzene	0.215	0.225	-4.7	140	0.00
67	Atrazine	0.205	0.227	-10.7	141	0.00
68 C	Pentachlorophenol	0.098	0.094	4.1	147	0.00
69	Phenanthrene	1.127	1.140	-1.2	131	0.00
70 S	Anthracene-d10	0.971	1.000	-3.0	135	0.00
71	Anthracene	1.156	1.194	-3.3	133	0.00
72	Carbazole	0.977	1.089	-11.5	133	0.00
73	Di-n-butylphthalate	1.272	1.394	-9.6	138	0.00
74 C	Fluoranthene	1.046	1.185	-13.3	133	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	134	0.00
76 S	Pyrene-d10	1.056	1.053	0.3	133	0.00
77	Pyrene	1.342	1.324	1.3	131	0.00
78	Butylbenzylphthalate	0.623	0.713	-14.4	152#	0.00
79	3,3'-Dichlorobenzidine	0.385	0.445	-15.6	151#	0.00
80	Benzo(a)anthracene	1.170	1.210	-3.4	138	0.00
81	Bis(2-ethylhexyl)phthalate	0.889	1.011	-13.7	148	0.00
82	Chrysene	1.087	1.116	-2.7	138	0.00
83 I	Perylene-d12	1.000	1.000	0.0	136	0.00
84	Di-n-octyl phthalate	1.374	1.657	-20.6#	152#	0.00
85	Benzo(b)fluoranthene	1.143	1.186	-3.8	145	0.00
86	Benzo(k)fluoranthene	1.123	1.220	-8.6	147	0.00
87 S	Benzo(a)pyrene-d12	0.946	0.997	-5.4	145	0.00

Data Path : Z:\HPCHEM1\BNA G\DATA\BG051617\
Data File : BG026995.D
Acq On : 17 May 2017 2:08
Operator : SJ/MA
Sample : SSTDCCC020
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTD02091

Quant Time: May 17 05:11:13 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG042717.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed May 17 04:07:33 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	1.196	-6.3	145	0.00
89	Indeno(1,2,3-cd)pyrene	1.333	1.137	14.7	118	0.00
90	Dibenzo(a,h)anthracene	1.130	0.991	12.3	121	0.00
91	Benzo(a,h,i)perylene	1.106	0.872	21.2#	110	0.00

(#) = Out of Range

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