

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081417\
 Data File : BG028323.D
 Acq On : 14 Aug 2017 22:04
 Operator : SJ/JU
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTD02068

Quant Time: Aug 15 06:40:01 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080217.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 15 06:10:38 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	131	0.00
2	1,4-Dioxane	0.451	0.499	-10.6	149	0.00
3 S	1,4-Dioxane-d8	0.429	0.383	10.7	129	0.00
4	Benzaldehyde	1.269	1.277	-0.6	123	0.00
5 S	Phenol-d5	2.197	2.075	5.6	130	0.00
6	Phenol	2.183	2.063	5.5	128	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	1.391	1.277	8.2	120	0.00
8	Bis(2-Chloroethyl)ether	1.487	1.359	8.6	121	0.00
9 S	2-Chlorophenol-d4	1.386	1.444	-4.2	133	0.00
10	2-Chlorophenol	1.350	1.384	-2.5	135	0.00
11	2-Methylphenol	1.541	1.495	3.0	128	0.00
12	2,2'-oxybis(1-Chloropropane	3.610	2.991	17.1	106	0.00
13 S	4-Methylphenol-d8	1.836	1.732	5.7	123	0.00
14	Acetophenone	2.618	2.416	7.7	118	0.00
15 P	N-Nitroso-di-n-propylamine	1.488	1.359	8.7	114	0.00
16	4-Methylphenol	1.740	1.671	4.0	125	0.00
17	Hexachloroethane	0.560	0.578	-3.2	133	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	127	0.00
19 S	Nitrobenzene-d5	0.156	0.163	-4.5	134	0.00
20	Nitrobenzene	0.453	0.443	2.2	126	0.00
21	Isophorone	0.860	0.785	8.7	116	0.00
22 S	2-Nitrophenol-d4	0.171	0.192	-12.3	142	0.00
23 C	2-Nitrophenol	0.173	0.177	-2.3	128	0.00
24	2,4-Dimethylphenol	0.430	0.429	0.2	125	0.00
25	Bis(2-Chloroethoxy)methane	0.469	0.443	5.5	119	0.00
26 S	2,4-Dichlorophenol-d3	0.355	0.383	-7.9	139	0.00
27 C	2,4-Dichlorophenol	0.323	0.348	-7.7	140	0.00
28	Naphthalene	0.991	1.016	-2.5	131	0.00
29 S	4-Chloroaniline-d4	0.395	0.455	-15.2	131	0.00
30	4-Chloroaniline	0.383	0.440	-14.9	131	0.00
31 C	Hexachlorobutadiene	0.238	0.249	-4.6	137	0.00
32	Caprolactam	0.134	0.122	9.0	117	0.00
33 C	4-Chloro-3-methylphenol	0.412	0.441	-7.0	134	0.00
34	2-Methylnaphthalene	0.796	0.816	-2.5	131	0.00
35 I	Acenaphthene-d10	1.000	1.000	0.0	126	0.00
36	1,2,4,5-Tetrachlorobenzene	0.645	0.688	-6.7	141	0.00
37	Hexachlorocyclopentadiene	0.356	0.254	28.7#	102	0.00
38 C	2,4,6-Trichlorophenol	0.402	0.474	-17.9	151#	0.00
39	2,4,5-Trichlorophenol	0.438	0.512	-16.9	150	0.00
40	1,1'-Biphenyl	1.461	1.497	-2.5	132	0.00
41	2-Chloronaphthalene	1.154	1.182	-2.4	131	0.00
42	2-Nitroaniline	0.484	0.498	-2.9	125	0.00
43 S	Dimethylphthalate-d6	1.779	1.732	2.6	122	0.00
44	Dimethylphthalate	1.638	1.616	1.3	125	0.00

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081417\
 Data File : BG028323.D
 Acq On : 14 Aug 2017 22:04
 Operator : SJ/JU
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTD02068

Quant Time: Aug 15 06:40:01 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080217.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 15 06:10:38 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	0.338	0.371	-9.8	134	0.00
46 S	Acenaphthylene-d8	2.006	2.104	-4.9	134	0.00
47	Acenaphthylene	1.917	1.976	-3.1	131	0.00
48	3-Nitroaniline	0.305	0.312	-2.3	125	0.00
49 C	Acenaphthene	1.293	1.344	-3.9	130	0.00
50	2,4-Dinitrophenol	0.191	0.184	3.7	143	0.00
51 S	4-Nitrophenol-d4	0.283	0.258	8.8	118	0.00
52	4-Nitrophenol	0.293	0.254	13.3	109	0.00
53	Dibenzofuran	1.886	1.952	-3.5	131	0.00
54	2,4-Dinitrotoluene	0.512	0.535	-4.5	131	0.00
55	2,3,4,6-Tetrachlorophenol	0.408	0.451	-10.5	141	0.00
56	Diethylphthalate	1.921	1.815	5.5	121	0.00
57 S	Fluorene-d10	1.596	1.637	-2.6	131	0.00
58	Fluorene	1.669	1.693	-1.4	130	0.00
59	4-Chlorophenyl-phenylether	0.896	0.921	-2.8	134	0.00
60	4-Nitroaniline	0.361	0.336	6.9	120	0.00
61 I	Phenanthrene-d10	1.000	1.000	0.0	116	0.00
62 S	4,6-Dinitro-2-methylphenol-	0.130	0.133	-2.3	134	0.00
63	4,6-Dinitro-2-methylphenol	0.132	0.135	-2.3	132	0.00
64	N-Nitrosodiphenylamine	0.526	0.585	-11.2	130	0.00
65	4-Bromophenyl-phenylether	0.215	0.238	-10.7	130	0.00
66	Hexachlorobenzene	0.229	0.246	-7.4	129	0.00
67	Atrazine	0.229	0.227	0.9	116	0.00
68 C	Pentachlorophenol	0.116	0.115	0.9	136	0.00
69	Phenanthrene	1.007	1.039	-3.2	121	0.00
70 S	Anthracene-d10	0.959	1.012	-5.5	127	0.00
71	Anthracene	1.029	1.090	-5.9	126	0.00
72	Carbazole	0.895	0.922	-3.0	123	0.00
73	Di-n-butylphthalate	1.114	1.136	-2.0	118	0.00
74 C	Fluoranthene	1.223	1.217	0.5	115	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	118	0.00
76 S	Pyrene-d10	1.026	1.009	1.7	113	0.00
77	Pyrene	1.193	1.187	0.5	116	0.00
78	Butylbenzylphthalate	0.447	0.465	-4.0	123	0.00
79	3,3'-Dichlorobenzidine	0.386	0.423	-9.6	126	0.00
80	Benzo(a)anthracene	1.156	1.179	-2.0	121	0.00
81	Bis(2-ethylhexyl)phthalate	0.636	0.641	-0.8	121	0.00
82	Chrysene	1.041	1.082	-3.9	122	0.00
83 I	Perylene-d12	1.000	1.000	0.0	113	0.00
84	Di-n-octyl phthalate	1.199	1.230	-2.6	120	0.00
85	Benzo(b)fluoranthene	1.197	1.229	-2.7	119	0.00
86	Benzo(k)fluoranthene	1.169	1.290	-10.4	123	0.00
87 S	Benzo(a)pyrene-d12	0.936	1.026	-9.6	125	0.00

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081417\
Data File : BG028323.D
Acq On : 14 Aug 2017 22:04
Operator : SJ/JU
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTD02068

Quant Time: Aug 15 06:40:01 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080217.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Aug 15 06:10:38 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	1.159	1.229	-6.0	121	0.00
89	Indeno(1,2,3-cd)pyrene	1.357	1.188	12.5	103	0.00
90	Dibenzo(a,h)anthracene	1.137	0.994	12.6	102	0.00
91	Benzo(g,h,i)perylene	1.116	0.852	23.7#	88	0.00

(#) = Out of Range

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