

Data Path : Z:\HPCHEM1\BNA_G\Data\BG121916\
 Data File : BG025318.D
 Acq On : 19 Dec 2016 14:56
 Operator : UM/SJ
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTD02023

Quant Time: Dec 20 02:23:09 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 20 01:37:17 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	111	0.00
2	1,4-Dioxane	0.504	0.420	16.7	95	0.00
3 S	1,4-Dioxane-d8	0.387	0.394	-1.8	115	0.00
4	Benzaldehyde	1.299	1.416	-9.0	116	0.00
5 S	Phenol-d5	1.725	1.830	-6.1	118	0.00
6	Phenol	1.771	1.899	-7.2	116	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	1.067	1.141	-6.9	118	0.00
8	Bis(2-Chloroethyl)ether	1.296	1.360	-4.9	120	0.00
9 S	2-Chlorophenol-d4	1.207	1.275	-5.6	116	0.00
10	2-Chlorophenol	1.213	1.278	-5.4	116	0.00
11	2-Methylphenol	1.304	1.397	-7.1	119	0.00
12	2,2'-oxybis(1-Chloropropane	2.204	2.426	-10.1	120	0.00
13 S	4-Methylphenol-d8	1.442	1.522	-5.5	120	0.00
14	Acetophenone	2.194	2.348	-7.0	120	0.00
15 P	N-Nitroso-di-n-propylamine	1.267	1.421	-12.2	122	0.00
16	4-Methylphenol	1.451	1.487	-2.5	121	0.00
17	Hexachloroethane	0.537	0.596	-11.0	120	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	112	0.00
19 S	Nitrobenzene-d5	0.142	0.152	-7.0	120	0.00
20	Nitrobenzene	0.435	0.489	-12.4	130	0.00
21	Isophorone	0.824	0.895	-8.6	120	0.00
22 S	2-Nitrophenol-d4	0.177	0.180	-1.7	117	0.00
23 C	2-Nitrophenol	0.179	0.188	-5.0	118	0.00
24	2,4-Dimethylphenol	0.410	0.450	-9.8	120	0.00
25	Bis(2-Chloroethoxy)methane	0.420	0.456	-8.6	119	0.00
26 S	2,4-Dichlorophenol-d3	0.340	0.372	-9.4	122	0.00
27 C	2,4-Dichlorophenol	0.330	0.358	-8.5	120	0.00
28	Naphthalene	0.930	0.955	-2.7	115	0.00
29 S	4-Chloroaniline-d4	0.334	0.410	-22.8#	137	0.00
30	4-Chloroaniline	0.341	0.407	-19.4	131	0.00
31 C	Hexachlorobutadiene	0.283	0.312	-10.2	122	0.00
32	Caprolactam	0.137	0.144	-5.1	120	0.00
33 C	4-Chloro-3-methylphenol	0.406	0.425	-4.7	116	0.00
34	2-Methylnaphthalene	0.734	0.760	-3.5	112	0.00
35 I	Acenaphthene-d10	1.000	1.000	0.0	106	0.00
36	1,2,4,5-Tetrachlorobenzene	0.603	0.704	-16.7	127	0.00
37	Hexachlorocyclopentadiene	0.354	0.297	16.1	97	0.00
38 C	2,4,6-Trichlorophenol	0.379	0.416	-9.8	118	0.00
39	2,4,5-Trichlorophenol	0.415	0.449	-8.2	113	0.00
40	1,1'-Biphenyl	1.213	1.344	-10.8	116	0.00
41	2-Chloronaphthalene	0.965	1.055	-9.3	119	0.00
42	2-Nitroaniline	0.378	0.433	-14.6	117	0.00
43 S	Dimethylphthalate-d6	1.376	1.481	-7.6	110	0.00
44	Dimethylphthalate	1.352	1.447	-7.0	112	0.00

Data Path : Z:\HPCHEM1\BNA_G\Data\BG121916\
 Data File : BG025318.D
 Acq On : 19 Dec 2016 14:56
 Operator : UM/SJ
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTD02023

Quant Time: Dec 20 02:23:09 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 20 01:37:17 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	0.291	0.298	-2.4	106	0.00
46 S	Acenaphthylene-d8	1.507	1.660	-10.2	116	0.00
47	Acenaphthylene	1.465	1.582	-8.0	112	0.00
48	3-Nitroaniline	0.252	0.280	-11.1	113	0.00
49 C	Acenaphthene	1.004	1.120	-11.6	118	0.00
50	2,4-Dinitrophenol	0.189	0.134	29.1#	85	0.00
51 S	4-Nitrophenol-d4	0.236	0.215	8.9	95	0.00
52	4-Nitrophenol	0.284	0.269	5.3	102	0.00
53	Dibenzofuran	1.587	1.741	-9.7	115	0.00
54	2,4-Dinitrotoluene	0.438	0.456	-4.1	109	0.00
55	2,3,4,6-Tetrachlorophenol	0.439	0.472	-7.5	113	0.00
56	Diethylphthalate	1.432	1.515	-5.8	109	0.00
57 S	Fluorene-d10	1.295	1.399	-8.0	112	0.00
58	Fluorene	1.292	1.383	-7.0	111	0.00
59	4-Chlorophenyl-phenylether	0.728	0.782	-7.4	113	0.00
60	4-Nitroaniline	0.303	0.282	6.9	99	0.00
61 I	Phenanthrene-d10	1.000	1.000	0.0	106	0.00
62 S	4,6-Dinitro-2-methylphenol-	0.124	0.113	8.9	101	0.00
63	4,6-Dinitro-2-methylphenol	0.128	0.122	4.7	105	0.00
64	N-Nitrosodiphenylamine	0.515	0.530	-2.9	107	0.00
65	4-Bromophenyl-phenylether	0.226	0.248	-9.7	117	0.00
66	Hexachlorobenzene	0.254	0.277	-9.1	113	0.00
67	Atrazine	0.245	0.258	-5.3	108	0.00
68 C	Pentachlorophenol	0.153	0.142	7.2	102	0.00
69	Phenanthrene	0.955	1.004	-5.1	106	0.00
70 S	Anthracene-d10	0.844	0.872	-3.3	107	0.00
71	Anthracene	0.981	1.034	-5.4	106	0.00
72	Carbazole	0.819	0.906	-10.6	108	0.00
73	Di-n-butylphthalate	1.034	1.125	-8.8	110	0.00
74 C	Fluoranthene	1.134	1.306	-15.2	108	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	105	0.00
76 S	Pyrene-d10	0.824	0.854	-3.6	106	0.00
77	Pyrene	0.961	1.009	-5.0	105	0.00
78	Butylbenzylphthalate	0.376	0.409	-8.8	109	0.00
79	3,3'-Dichlorobenzidine	0.349	0.364	-4.3	104	0.00
80	Benzo(a)anthracene	1.039	1.102	-6.1	108	0.00
81	Bis(2-ethylhexyl)phthalate	0.518	0.575	-11.0	113	0.00
82	Chrysene	0.972	1.029	-5.9	108	0.00
83 I	Perylene-d12	1.000	1.000	0.0	105	0.00
84	Di-n-octyl phthalate	0.812	0.943	-16.1	113	0.00
85	Benzo(b)fluoranthene	0.992	1.061	-7.0	109	0.00
86	Benzo(k)fluoranthene	0.943	1.015	-7.6	110	0.00
87 S	Benzo(a)pyrene-d12	0.810	0.893	-10.2	112	0.00

Data Path : Z:\HPCHEM1\BNA_G\Data\BG121916\
Data File : BG025318.D
Acq On : 19 Dec 2016 14:56
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTD02023

Quant Time: Dec 20 02:23:09 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Dec 20 01:37:17 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	0.954	1.014	-6.3	108	0.00
89	Indeno(1,2,3-cd)pyrene	1.185	1.284	-8.4	112	0.01
90	Dibenzo(a,h)anthracene	0.998	1.072	-7.4	112	0.00
91	Benzo(g,h,i)perylene	0.989	1.053	-6.5	110	0.01

(#) = Out of Range

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