

Data Path : Z:\HPCHEM1\BNA G\DATA\BG120616\
 Data File : BG024987.D
 Acq On : 6 Dec 2016 19:44
 Operator : UM/SJ
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTD02016

Quant Time: Dec 07 03:17:48 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG113016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 07 02:35:14 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	103	0.00
2	1,4-Dioxane	0.504	0.466	7.5	98	0.00
3 S	1,4-Dioxane-d8	0.387	0.382	1.3	103	0.00
4	Benzaldehyde	1.299	1.282	1.3	97	0.00
5 S	Phenol-d5	1.725	1.579	8.5	94	0.00
6	Phenol	1.771	1.750	1.2	99	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	1.067	1.029	3.6	98	0.00
8	Bis(2-Chloroethyl)ether	1.296	1.248	3.7	101	0.00
9 S	2-Chlorophenol-d4	1.207	1.166	3.4	98	0.00
10	2-Chlorophenol	1.213	1.163	4.1	98	0.00
11	2-Methylphenol	1.304	1.220	6.4	96	0.00
12	2,2'-oxybis(1-Chloropropane	2.204	2.191	0.6	100	0.00
13 S	4-Methylphenol-d8	1.442	1.372	4.9	100	0.00
14	Acetophenone	2.194	2.101	4.2	100	0.00
15 P	N-Nitroso-di-n-propylamine	1.267	1.231	2.8	98	0.00
16	4-Methylphenol	1.451	1.448	0.2	109	0.00
17	Hexachloroethane	0.537	0.501	6.7	94	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	103	0.00
19 S	Nitrobenzene-d5	0.142	0.133	6.3	97	0.00
20	Nitrobenzene	0.435	0.433	0.5	106	0.00
21	Isophorone	0.824	0.789	4.2	98	0.00
22 S	2-Nitrophenol-d4	0.177	0.163	7.9	98	0.00
23 C	2-Nitrophenol	0.179	0.165	7.8	96	0.00
24	2,4-Dimethylphenol	0.410	0.402	2.0	98	0.00
25	Bis(2-Chloroethoxy)methane	0.420	0.427	-1.7	103	0.00
26 S	2,4-Dichlorophenol-d3	0.340	0.331	2.6	100	0.00
27 C	2,4-Dichlorophenol	0.330	0.329	0.3	101	0.00
28	Naphthalene	0.930	0.887	4.6	99	0.00
29 S	4-Chloroaniline-d4	0.334	0.374	-12.0	115	0.00
30	4-Chloroaniline	0.341	0.374	-9.7	111	0.00
31 C	Hexachlorobutadiene	0.283	0.281	0.7	101	0.00
32	Caprolactam	0.137	0.128	6.6	99	0.00
33 C	4-Chloro-3-methylphenol	0.406	0.396	2.5	100	0.00
34	2-Methylnaphthalene	0.734	0.716	2.5	97	0.00
35 I	Acenaphthene-d10	1.000	1.000	0.0	102	0.00
36	1,2,4,5-Tetrachlorobenzene	0.603	0.608	-0.8	106	0.00
37	Hexachlorocyclopentadiene	0.354	0.334	5.6	105	0.00
38 C	2,4,6-Trichlorophenol	0.379	0.371	2.1	102	0.00
39	2,4,5-Trichlorophenol	0.415	0.426	-2.7	103	0.00
40	1,1'-Biphenyl	1.213	1.250	-3.1	104	0.00
41	2-Chloronaphthalene	0.965	0.960	0.5	105	0.00
42	2-Nitroaniline	0.378	0.420	-11.1	109	0.00
43 S	Dimethylphthalate-d6	1.376	1.416	-2.9	102	0.00
44	Dimethylphthalate	1.352	1.366	-1.0	102	0.00

Data Path : Z:\HPCHEM1\BNA G\DATA\BG120616\
 Data File : BG024987.D
 Acq On : 6 Dec 2016 19:44
 Operator : UM/SJ
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTD02016

Quant Time: Dec 07 03:17:48 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG113016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 07 02:35:14 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
45	2,6-Dinitrotoluene	0.291	0.274	5.8	93	0.00
46 S	Acenaphthylene-d8	1.507	1.528	-1.4	102	0.00
47	Acenaphthylene	1.465	1.511	-3.1	103	0.00
48	3-Nitroaniline	0.252	0.291	-15.5	113	0.00
49 C	Acenaphthene	1.004	1.016	-1.2	103	0.00
50	2,4-Dinitrophenol	0.189	0.195	-3.2	118	0.00
51 S	4-Nitrophenol-d4	0.236	0.251	-6.4	107	0.00
52	4-Nitrophenol	0.284	0.305	-7.4	111	0.00
53	Dibenzofuran	1.587	1.647	-3.8	105	0.00
54	2,4-Dinitrotoluene	0.438	0.478	-9.1	110	0.00
55	2,3,4,6-Tetrachlorophenol	0.439	0.442	-0.7	102	0.00
56	Diethylphthalate	1.432	1.436	-0.3	100	0.00
57 S	Fluorene-d10	1.295	1.317	-1.7	102	0.00
58	Fluorene	1.292	1.327	-2.7	102	0.00
59	4-Chlorophenyl-phenylether	0.728	0.749	-2.9	104	0.00
60	4-Nitroaniline	0.303	0.315	-4.0	106	0.00
61 I	Phenanthrene-d10	1.000	1.000	0.0	106	0.00
62 S	4,6-Dinitro-2-methylphenol-	0.124	0.127	-2.4	115	0.00
63	4,6-Dinitro-2-methylphenol	0.128	0.133	-3.9	114	0.00
64	N-Nitrosodiphenylamine	0.515	0.519	-0.8	105	0.00
65	4-Bromophenyl-phenylether	0.226	0.228	-0.9	107	0.00
66	Hexachlorobenzene	0.254	0.260	-2.4	106	0.00
67	Atrazine	0.245	0.247	-0.8	104	0.00
68 C	Pentachlorophenol	0.153	0.141	7.8	102	0.00
69	Phenanthrene	0.955	0.990	-3.7	105	0.00
70 S	Anthracene-d10	0.844	0.878	-4.0	107	0.00
71	Anthracene	0.981	1.021	-4.1	105	0.00
72	Carbazole	0.819	0.895	-9.3	107	0.00
73	Di-n-butylphthalate	1.034	1.080	-4.4	106	0.00
74 C	Fluoranthene	1.134	1.313	-15.8	108	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	117	0.00
76 S	Pyrene-d10	0.824	0.795	3.5	110	0.00
77	Pyrene	0.961	0.936	2.6	109	0.00
78	Butylbenzylphthalate	0.376	0.383	-1.9	114	0.00
79	3,3'-Dichlorobenzidine	0.349	0.384	-10.0	123	0.00
80	Benzo(a)anthracene	1.039	1.054	-1.4	115	0.00
81	Bis(2-ethylhexyl)phthalate	0.518	0.515	0.6	113	0.00
82	Chrysene	0.972	0.998	-2.7	117	0.00
83 I	Perylene-d12	1.000	1.000	0.0	118	0.00
84	Di-n-octyl phthalate	0.812	0.845	-4.1	114	0.00
85	Benzo(b)fluoranthene	0.992	0.990	0.2	114	0.00
86	Benzo(k)fluoranthene	0.943	0.991	-5.1	121	0.00
87 S	Benzo(a)pyrene-d12	0.810	0.817	-0.9	116	0.00

Data Path : Z:\HPCHEM1\BNA G\DATA\BG120616\
Data File : BG024987.D
Acq On : 6 Dec 2016 19:44
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTD02016

Quant Time: Dec 07 03:17:48 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Dec 07 02:35:14 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	0.975	-2.2	118	0.00
89	Indeno(1,2,3-cd)pyrene	1.185	1.197	-1.0	118	0.00
90	Dibenzo(a,h)anthracene	0.998	1.015	-1.7	120	0.00
91	Benzo(a,h,i)perylene	0.989	0.991	-0.2	117	0.00

(#) = Out of Range

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