

Data Path : Z:\HPCHEM1\BNA G\DATA\BG060116\
 Data File : BG022137.D
 Acq On : 1 Jun 2016 11:44
 Operator : UM/SJ
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTD02026

Quant Time: Jun 02 01:09:37 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG052116.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jun 01 01:07:08 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	98	0.00
2	1,4-Dioxane	0.696	0.669	3.9	97	0.00
3 S	1,4-Dioxane-d8	0.384	0.366	4.7	97	0.00
4	Benzaldehyde	0.996	0.619	37.9#	55	0.00
5 S	Phenol-d5	1.807	1.564	13.4	88	0.00
6	Phenol	1.820	1.547	15.0	83	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	0.947	0.796	15.9	83	0.00
8	Bis(2-Chloroethyl)ether	1.330	1.218	8.4	90	0.00
9 S	2-Chlorophenol-d4	1.511	1.354	10.4	88	0.00
10	2-Chlorophenol	1.489	1.337	10.2	87	0.00
11	2-Methylphenol	1.454	1.266	12.9	86	0.00
12	2,2'-oxybis(1-Chloropropane	1.589	1.287	19.0	77	0.00
13 S	4-Methylphenol-d8	1.533	1.353	11.7	86	0.00
14	Acetophenone	2.144	1.960	8.6	88	0.00
15 P	N-Nitroso-di-n-propylamine	1.107	0.995	10.1	87	0.00
16	4-Methylphenol	1.567	1.391	11.2	86	0.00
17	Hexachloroethane	0.594	0.523	12.0	86	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	98	0.00
19 S	Nitrobenzene-d5	0.148	0.133	10.1	87	0.00
20	Nitrobenzene	0.299	0.263	12.0	85	0.00
21	Isophorone	0.636	0.589	7.4	88	0.00
22 S	2-Nitrophenol-d4	0.160	0.151	5.6	90	0.00
23 C	2-Nitrophenol	0.167	0.155	7.2	88	0.00
24	2,4-Dimethylphenol	0.322	0.289	10.2	85	0.00
25	Bis(2-Chloroethoxy)methane	0.365	0.345	5.5	89	0.00
26 S	2,4-Dichlorophenol-d3	0.282	0.266	5.7	88	0.00
27 C	2,4-Dichlorophenol	0.277	0.264	4.7	92	0.00
28	Naphthalene	1.014	0.925	8.8	88	0.00
29 S	4-Chloroaniline-d4	0.307	0.336	-9.4	103	0.00
30	4-Chloroaniline	0.309	0.345	-11.7	103	0.00
31 C	Hexachlorobutadiene	0.163	0.145	11.0	85	0.00
32	Caprolactam	0.111	0.111	0.0	96	0.00
33 C	4-Chloro-3-methylphenol	0.295	0.275	6.8	89	0.00
34	2-Methylnaphthalene	0.725	0.665	8.3	88	0.00
35 I	Acenaphthene-d10	1.000	1.000	0.0	108	0.00
36	1,2,4,5-Tetrachlorobenzene	0.593	0.502	15.3	90	0.00
37	Hexachlorocyclopentadiene	0.256	0.239	6.6	115	0.00
38 C	2,4,6-Trichlorophenol	0.401	0.339	15.5	90	0.00
39	2,4,5-Trichlorophenol	0.425	0.382	10.1	99	0.00
40	1,1'-Biphenyl	1.647	1.434	12.9	92	0.00
41	2-Chloronaphthalene	1.265	1.107	12.5	93	0.00
42	2-Nitroaniline	0.285	0.232	18.6	85	0.00
43 S	Dimethylphthalate-d6	1.502	1.455	3.1	103	0.00
44	Dimethylphthalate	1.510	1.439	4.7	101	0.00

Data Path : Z:\HPCHEM1\BNA G\DATA\BG060116\
 Data File : BG022137.D
 Acq On : 1 Jun 2016 11:44
 Operator : UM/SJ
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTD02026

Quant Time: Jun 02 01:09:37 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG052116.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jun 01 01:07:08 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	0.324	0.316	2.5	103	0.00
46 S	Acenaphthylene-d8	2.098	1.885	10.2	95	0.00
47	Acenaphthylene	2.137	1.903	10.9	95	0.00
48	3-Nitroaniline	0.310	0.256	17.4	95	0.00
49 C	Acenaphthene	1.476	1.314	11.0	94	0.00
50	2,4-Dinitrophenol	0.125	0.096	23.2#	105	0.00
51 S	4-Nitrophenol-d4	0.259	0.209	19.3	94	0.00
52	4-Nitrophenol	0.146	0.113	22.6#	89	0.00
53	Dibenzofuran	1.825	1.714	6.1	99	0.00
54	2,4-Dinitrotoluene	0.437	0.429	1.8	101	0.00
55	2,3,4,6-Tetrachlorophenol	0.336	0.298	11.3	93	0.00
56	Diethylphthalate	1.566	1.518	3.1	103	0.00
57 S	Fluorene-d10	1.407	1.332	5.3	100	0.00
58	Fluorene	1.541	1.476	4.2	101	0.00
59	4-Chlorophenyl-phenylether	0.694	0.663	4.5	100	0.00
60	4-Nitroaniline	0.341	0.226	33.7#	68	0.00
61 I	Phenanthrene-d10	1.000	1.000	0.0	124	0.00
62 S	4,6-Dinitro-2-methylphenol-	0.105	0.088	16.2	108	0.00
63	4,6-Dinitro-2-methylphenol	0.108	0.091	15.7	117	0.00
64	N-Nitrosodiphenylamine	0.632	0.540	14.6	102	0.00
65	4-Bromophenyl-phenylether	0.202	0.176	12.9	105	0.00
66	Hexachlorobenzene	0.235	0.209	11.1	115	0.00
67	Atrazine	0.218	0.197	9.6	108	0.00
68 C	Pentachlorophenol	0.122	0.074	39.3#	78	0.00
69	Phenanthrene	1.112	1.007	9.4	109	0.00
70 S	Anthracene-d10	0.960	0.880	8.3	112	0.00
71	Anthracene	1.127	0.997	11.5	105	0.00
72	Carbazole	0.977	0.881	9.8	111	0.00
73	Di-n-butylphthalate	1.309	1.195	8.7	110	0.00
74 C	Fluoranthene	1.137	1.114	2.0	119	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	155#	0.00
76 S	Pyrene-d10	1.128	0.937	16.9	120	0.00
77	Pyrene	1.407	1.145	18.6	117	0.00
78	Butylbenzylphthalate	0.667	0.546	18.1	120	0.00
79	3,3'-Dichlorobenzidine	0.376	0.333	11.4	129	0.00
80	Benzo(a)anthracene	1.173	1.019	13.1	128	0.00
81	Bis(2-ethylhexyl)phthalate	0.949	0.797	16.0	124	0.00
82	Chrysene	1.125	0.992	11.8	134	0.00
83 I	Perylene-d12	1.000	1.000	0.0	142	0.00
84	Di-n-octyl phthalate	1.630	1.461	10.4	123	0.00
85	Benzo(b)fluoranthene	1.170	1.112	5.0	132	0.00
86	Benzo(k)fluoranthene	1.139	1.017	10.7	129	0.00
87 S	Benzo(a)pyrene-d12	0.928	0.857	7.7	129	0.00

Data Path : Z:\HPCHEM1\BNA G\DATA\BG060116\
Data File : BG022137.D
Acq On : 1 Jun 2016 11:44
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTD02026

Quant Time: Jun 02 01:09:37 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG052116.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jun 01 01:07:08 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	1.129	1.030	8.8	129	0.00
89	Indeno(1,2,3-cd)pyrene	1.295	1.188	8.3	132	0.00
90	Dibenzo(a,h)anthracene	1.083	0.971	10.3	127	0.00
91	Benzo(a,h,i)perylene	1.037	1.002	3.4	143	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 1