

Data Path : Z:\HPCHEM1\BNA G\DATA\BG063016\
 Data File : BG022729.D
 Acq On : 30 Jun 2016 16:21
 Operator : UM/SJ
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTD02007

Quant Time: Jun 30 17:06:10 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG063016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jun 30 16:28:44 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	98	0.00
2	1,4-Dioxane	0.419	0.406	3.1	107	0.00
3 S	1,4-Dioxane-d8	0.358	0.286	20.1#	80	0.00
4	Benzaldehyde	1.338	1.276	4.6	85	0.00
5 S	Phenol-d5	2.104	1.878	10.7	87	0.00
6	Phenol	2.155	1.982	8.0	89	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	1.186	1.098	7.4	87	0.00
8	Bis(2-Chloroethyl)ether	1.473	1.424	3.3	93	0.00
9 S	2-Chlorophenol-d4	1.405	1.202	14.4	84	0.00
10	2-Chlorophenol	1.429	1.287	9.9	88	0.00
11	2-Methylphenol	1.619	1.469	9.3	91	0.00
12	2,2'-oxybis(1-Chloropropane	1.657	1.526	7.9	88	0.00
13 S	4-Methylphenol-d8	1.634	1.467	10.2	87	0.00
14	Acetophenone	2.639	2.403	8.9	89	0.00
15 P	N-Nitroso-di-n-propylamine	1.586	1.404	11.5	89	0.00
16	4-Methylphenol	1.800	1.589	11.7	88	0.00
17	Hexachloroethane	0.609	0.542	11.0	83	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	99	0.00
19 S	Nitrobenzene-d5	0.159	0.140	11.9	88	0.00
20	Nitrobenzene	0.490	0.449	8.4	87	0.00
21	Isophorone	0.875	0.794	9.3	89	0.00
22 S	2-Nitrophenol-d4	0.192	0.167	13.0	88	0.00
23 C	2-Nitrophenol	0.204	0.188	7.8	89	0.00
24	2,4-Dimethylphenol	0.489	0.445	9.0	92	0.00
25	Bis(2-Chloroethoxy)methane	0.516	0.460	10.9	90	0.00
26 S	2,4-Dichlorophenol-d3	0.402	0.375	6.7	95	0.00
27 C	2,4-Dichlorophenol	0.404	0.363	10.1	91	0.00
28	Naphthalene	1.023	0.931	9.0	92	0.00
29 S	4-Chloroaniline-d4	0.341	0.303	11.1	96	0.00
30	4-Chloroaniline	0.348	0.308	11.5	94	0.00
31 C	Hexachlorobutadiene	0.311	0.276	11.3	87	0.00
32	Caprolactam	0.135	0.116	14.1	86	0.00
33 C	4-Chloro-3-methylphenol	0.511	0.448	12.3	91	0.00
34	2-Methylnaphthalene	0.846	0.766	9.5	90	0.00
35 I	Acenaphthene-d10	1.000	1.000	0.0	97	0.00
36	1,2,4,5-Tetrachlorobenzene	0.748	0.684	8.6	89	0.00
37	Hexachlorocyclopentadiene	0.451	0.382	15.3	82	0.00
38 C	2,4,6-Trichlorophenol	0.516	0.484	6.2	92	0.00
39	2,4,5-Trichlorophenol	0.543	0.501	7.7	87	0.00
40	1,1'-Biphenyl	1.521	1.423	6.4	91	0.00
41	2-Chloronaphthalene	1.232	1.142	7.3	89	0.00
42	2-Nitroaniline	0.476	0.421	11.6	86	0.00
43 S	Dimethylphthalate-d6	1.625	1.535	5.5	90	0.00
44	Dimethylphthalate	1.659	1.572	5.2	91	0.00

Data Path : Z:\HPCHEM1\BNA G\DATA\BG063016\
 Data File : BG022729.D
 Acq On : 30 Jun 2016 16:21
 Operator : UM/SJ
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTD02007

Quant Time: Jun 30 17:06:10 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG063016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jun 30 16:28:44 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	0.359	0.341	5.0	92	0.00
46 S	Acenaphthylene-d8	1.862	1.759	5.5	91	0.00
47	Acenaphthylene	1.875	1.773	5.4	90	0.00
48	3-Nitroaniline	0.281	0.272	3.2	94	0.00
49 C	Acenaphthene	1.274	1.183	7.1	90	0.00
50	2,4-Dinitrophenol	0.200	0.139	30.5#	80	0.00
51 S	4-Nitrophenol-d4	0.255	0.240	5.9	91	0.00
52	4-Nitrophenol	0.324	0.301	7.1	89	0.00
53	Dibenzofuran	1.968	1.882	4.4	93	0.00
54	2,4-Dinitrotoluene	0.519	0.494	4.8	95	0.00
55	2,3,4,6-Tetrachlorophenol	0.552	0.526	4.7	90	0.00
56	Diethylphthalate	1.686	1.592	5.6	90	0.00
57 S	Fluorene-d10	1.539	1.449	5.8	91	0.00
58	Fluorene	1.581	1.478	6.5	90	0.00
59	4-Chlorophenyl-phenylether	0.947	0.873	7.8	91	0.00
60	4-Nitroaniline	0.337	0.336	0.3	93	0.00
61 I	Phenanthrene-d10	1.000	1.000	0.0	99	0.00
62 S	4,6-Dinitro-2-methylphenol-	0.129	0.100	22.5#	82	0.00
63	4,6-Dinitro-2-methylphenol	0.136	0.111	18.4	85	0.00
64	N-Nitrosodiphenylamine	0.589	0.545	7.5	91	0.00
65	4-Bromophenyl-phenylether	0.268	0.242	9.7	91	0.00
66	Hexachlorobenzene	0.283	0.255	9.9	90	0.00
67	Atrazine	0.246	0.233	5.3	91	0.00
68 C	Pentachlorophenol	0.163	0.126	22.7	80	0.00
69	Phenanthrene	1.023	0.987	3.5	93	0.00
70 S	Anthracene-d10	0.930	0.861	7.4	91	0.00
71	Anthracene	1.051	0.996	5.2	90	0.00
72	Carbazole	0.778	0.805	-3.5	90	0.00
73	Di-n-butylphthalate	0.999	0.961	3.8	90	0.00
74 C	Fluoranthene	1.066	1.168	-9.6	90	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	101	0.00
76 S	Pyrene-d10	0.955	0.898	6.0	92	0.00
77	Pyrene	1.120	1.044	6.8	92	0.00
78	Butylbenzylphthalate	0.434	0.395	9.0	91	0.00
79	3,3'-Dichlorobenzidine	0.379	0.348	8.2	88	0.00
80	Benzo(a)anthracene	1.173	1.107	5.6	91	0.00
81	Bis(2-ethylhexyl)phthalate	0.560	0.500	10.7	90	0.00
82	Chrysene	1.068	1.014	5.1	93	0.00
83 I	Perylene-d12	1.000	1.000	0.0	100	0.00
84	Di-n-octyl phthalate	0.881	0.874	0.8	89	0.00
85	Benzo(b)fluoranthene	1.250	1.127	9.8	90	0.00
86	Benzo(k)fluoranthene	1.174	1.148	2.2	93	0.00
87 S	Benzo(a)pyrene-d12	0.950	0.870	8.4	91	0.00

Data Path : Z:\HPCHEM1\BNA G\DATA\BG063016\
Data File : BG022729.D
Acq On : 30 Jun 2016 16:21
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTD02007

Quant Time: Jun 30 17:06:10 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG063016.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 30 16:28:44 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	1.187	1.082	8.8	90	0.00
89	Indeno(1,2,3-cd)pyrene	1.255	1.079	14.0	84	0.00
90	Dibenzo(a,h)anthracene	1.038	0.843	18.8	78	0.00
91	Benzo(a,h,i)perylene	0.984	0.791	19.6	79	0.00

(#) = Out of Range

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