

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA G\Data\BG021116\
 Data File : BG020872.D
 Acq On : 12 Feb 2016 1:20
 Operator : SJ/UM
 Sample : SSTDCCC040
 Misc : L00 20PPM
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Feb 12 04:51:03 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG021116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 11 15:26:53 2016
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	95	0.00
2	1,4-Dioxane	40.000	39.297	1.8	96	0.00
3	Pyridine	40.000	41.082	-2.7	94	0.00
4	n-Nitrosodimethylamine	40.000	41.065	-2.7	95	0.00
5 S	2-Fluorophenol	80.000	81.563	-2.0	96	0.00
6	Aniline	40.000	40.787	-2.0	96	0.00
7 S	Phenol-d6	80.000	82.912	-3.6	97	0.00
8	2-Chlorophenol	40.000	41.185	-3.0	98	0.00
9	Benzaldehyde	40.000	41.407	-3.5	97	0.00
10 C	Phenol	40.000	40.247	-0.6	95	0.00
11	bis(2-Chloroethyl)ether	40.000	40.573	-1.4	94	0.00
12	1,3-Dichlorobenzene	40.000	40.657	-1.6	97	0.00
13 C	1,4-Dichlorobenzene	40.000	41.079	-2.7	97	0.00
14	1,2-Dichlorobenzene	40.000	41.306	-3.3	97	0.00
15	Benzyl Alcohol	40.000	40.850	-2.1	96	0.00
16	2,2'-oxybis(1-Chloropropane)	40.000	39.342	1.6	93	0.00
17	2-Methylphenol	40.000	40.516	-1.3	96	0.00
18	Hexachloroethane	40.000	40.865	-2.2	96	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	40.782	-2.0	97	0.00
20	3+4-Methylphenols	40.000	41.039	-2.6	98	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	97	0.00
22	Acetophenone	40.000	40.866	-2.2	97	0.00
23 S	Nitrobenzene-d5	80.000	81.210	-1.5	97	0.00
24	Nitrobenzene	40.000	40.262	-0.7	97	0.00
25	Isophorone	40.000	40.819	-2.0	97	0.00
26 C	2-Nitrophenol	40.000	41.758	-4.4	96	0.00
27	2,4-Dimethylphenol	40.000	41.430	-3.6	97	0.00
28	bis(2-Chloroethoxy)methane	40.000	40.710	-1.8	98	0.00
29 C	2,4-Dichlorophenol	40.000	41.056	-2.6	99	0.00
30	1,2,4-Trichlorobenzene	40.000	40.809	-2.0	98	0.00
31	Naphthalene	40.000	40.752	-1.9	98	0.00
32	Benzoic acid	40.000	39.198	2.0	96	0.00
33	4-Chloroaniline	40.000	40.476	-1.2	98	0.00
34 C	Hexachlorobutadiene	40.000	40.676	-1.7	98	0.00
35	Caprolactam	40.000	42.277	-5.7	103	0.00
36 C	4-Chloro-3-methylphenol	40.000	41.084	-2.7	100	0.00
37	2-Methylnaphthalene	40.000	40.960	-2.4	100	0.00
38 I	Acenaphthene-d10	20.000	20.000	0.0	100	0.00
39	1,2,4,5-Tetrachlorobenzene	40.000	40.628	-1.6	101	0.00
40 P	Hexachlorocyclopentadiene	40.000	39.611	1.0	96	0.00
41 S	2,4,6-Tribromophenol	80.000	83.630	-4.5	105	0.00
42 C	2,4,6-Trichlorophenol	40.000	41.785	-4.5	102	0.00
43	2,4,5-Trichlorophenol	40.000	41.613	-4.0	103	0.00
44 S	2-Fluorobiphenyl	80.000	81.801	-2.3	101	0.00

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA G\Data\BG021116\
 Data File : BG020872.D
 Acq On : 12 Feb 2016 1:20
 Operator : SJ/UM
 Sample : SSTDCCC040
 Misc : L00 20PPM
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Feb 12 04:51:03 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG021116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 11 15:26:53 2016
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
45	1,1'-Biphenyl	40.000	40.879	-2.2	102	0.00
46	2-Chloronaphthalene	40.000	40.667	-1.7	102	0.00
47	2-Nitroaniline	40.000	41.502	-3.8	102	0.00
48	Acenaphthylene	40.000	40.985	-2.5	102	0.00
49	Dimethylphthalate	40.000	41.821	-4.6	104	0.00
50	2,6-Dinitrotoluene	40.000	41.971	-4.9	104	0.00
51 C	Acenaphthene	40.000	41.524	-3.8	104	0.00
52	3-Nitroaniline	40.000	42.088	-5.2	104	0.00
53 P	2,4-Dinitrophenol	40.000	36.701	8.2	96	0.00
54	Dibenzofuran	40.000	41.639	-4.1	105	0.00
55 P	4-Nitrophenol	40.000	42.814	-7.0	104	0.00
56	2,4-Dinitrotoluene	40.000	43.051	-7.6	105	0.00
57	Fluorene	40.000	41.963	-4.9	105	0.00
58	2,3,4,6-Tetrachlorophenol	40.000	42.070	-5.2	105	0.00
59	Diethylphthalate	40.000	42.060	-5.2	106	0.00
60	4-Chlorophenyl-phenylether	40.000	40.895	-2.2	105	0.00
61	4-Nitroaniline	40.000	42.718	-6.8	105	0.00
62	Azobenzene	40.000	40.790	-2.0	102	0.00
63 I	Phenanthrene-d10	20.000	20.000	0.0	105	0.00
64	4,6-Dinitro-2-methylphenol	40.000	38.582	3.5	104	0.00
65 c	n-Nitrosodiphenylamine	40.000	40.734	-1.8	104	0.00
66	4-Bromophenyl-phenylether	40.000	40.815	-2.0	105	0.00
67	Hexachlorobenzene	40.000	40.595	-1.5	106	0.00
68	Atrazine	40.000	41.525	-3.8	108	0.00
69 C	Pentachlorophenol	40.000	39.620	1.0	103	0.00
70	Phenanthrene	40.000	40.822	-2.1	106	0.00
71	Anthracene	40.000	40.742	-1.9	106	0.00
72	Carbazole	40.000	41.081	-2.7	107	0.00
73	Di-n-butylphthalate	40.000	41.282	-3.2	107	0.00
74 C	Fluoranthene	40.000	41.040	-2.6	107	0.00
75 I	Chrysene-d12	20.000	20.000	0.0	105	0.00
76	Benzidine	40.000	40.069	-0.2	102	0.00
77	Pyrene	40.000	40.809	-2.0	108	0.00
78 S	Terphenyl-d14	80.000	82.100	-2.6	107	0.00
79	Butylbenzylphthalate	40.000	41.070	-2.7	106	0.00
80	Benzo(a)anthracene	40.000	40.918	-2.3	108	0.00
81	3,3'-Dichlorobenzidine	40.000	40.634	-1.6	106	0.00
82	Chrysene	40.000	40.560	-1.4	107	0.00
83	Bis(2-ethylhexyl)phthalate	40.000	41.008	-2.5	107	0.00
84 c	Di-n-octyl phthalate	40.000	41.486	-3.7	107	0.00
85	Indeno(1,2,3-cd)pyrene	40.000	41.625	-4.1	108	-0.02
86 I	Perylene-d12	20.000	20.000	0.0	106	-0.01
87	Benzo(b)fluoranthene	40.000	41.210	-3.0	108	0.00

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA G\Data\BG021116\
Data File : BG020872.D
Acq On : 12 Feb 2016 1:20
Operator : SJ/UM
Sample : SSTDCCC040
Misc : L00 20PPM
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Feb 12 04:51:03 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG021116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Feb 11 15:26:53 2016
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	40.000	41.201	-3.0	109	0.00
89 C	Benzo(a)pyrene	40.000	41.433	-3.6	109	0.00
90	Dibenzo(a,h)anthracene	40.000	41.014	-2.5	109	-0.02
91	Benzo(a,h,i)perylene	40.000	41.024	-2.6	108	-0.03

(#) = Out of Range

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