Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA G\Data\BG021116\
Data File : BG020874.D
Acq On : 12 Feb 2016 2:38
Operator : SJ/UM
Sample : SSTDCCC040
Misc : LOO 20PPM
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 12 04:51:44 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	98	0.00
2		1,4-Dioxane	40.000	39.102	2.2	98	0.00
3		Pyridine	40.000	41.259	-3.1	98	0.00
4		n-Nitrosodimethylamine	40.000	41.674	-4.2	100	0.00
5	S	2-Fluorophenol	80.000	80.524	-0.7	98	0.00
6		Aniline	40.000	40.484	-1.2	98	0.00
7	S	Phenol-d6	80.000	80.827	-1.0	98	0.00
8		2-Chlorophenol	40.000	40.254	-0.6	99	0.00
9		Benzaldehyde	40.000	40.203	-0.5	98	0.00
10	C	Phenol	40.000	39.831	0.4	98	0.00
11		bis(2-Chloroethyl)ether	40.000	39.782	0.5	96	0.00
12		1,3-Dichlorobenzene	40.000	40.030	-0.1	99	0.00
13	C	1,4-Dichlorobenzene	40.000	40.349	-0.9	98	0.00
14		1,2-Dichlorobenzene	40.000	39.883	0.3	97	0.00
15		Benzyl Alcohol	40.000	40.155	-0.4	97	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	38.591	3.5	94	0.00
17		2-Methylphenol	40.000	40.299	-0.7	99	0.00
18		Hexachloroethane	40.000	40.542	-1.4	99	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.785	0.5	98	0.00
20		3+4-Methylphenols	40.000	40.519	-1.3	100	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	98	0.00
22		Acetophenone	40.000	40.814	-2.0	98	0.00
23	S	Nitrobenzene-d5	80.000	81.562	-2.0	98	0.00
24		Nitrobenzene	40.000	40.386	-1.0	99	0.00
25		Isophorone	40.000	41.052	-2.6	99	0.00
26	C	2-Nitrophenol	40.000	41.278	-3.2	96	0.00
27		2,4-Dimethylphenol	40.000	41.154	-2.9	97	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.164	-0.4	97	0.00
29	C	2,4-Dichlorophenol	40.000	41.480	-3.7	101	0.00
30		1,2,4-Trichlorobenzene	40.000	40.821	-2.1	99	0.00
31		Naphthalene	40.000	40.767	-1.9	99	0.00
32		Benzoic acid	40.000	38.237	4.4	95	0.00
33		4-Chloroaniline	40.000	41.610	-4.0	101	0.00
34	C	Hexachlorobutadiene	40.000	40.960	-2.4	100	0.00
35		Caprolactam	40.000	43.058	-7.6	106	0.00
36	C	4-Chloro-3-methylphenol	40.000	41.618	-4.0	102	0.00
37		2-Methylnaphthalene	40.000	41.160	-2.9	101	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	103	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	40.164	-0.4	103	0.00
40	P	Hexachlorocyclopentadiene	40.000	38.299	4.3	95	0.00
41	S	2,4,6-Tribromophenol	80.000	81.189	-1.5	104	0.00
42	C	2,4,6-Trichlorophenol	40.000	41.296	-3.2	103	0.00
43		2,4,5-Trichlorophenol	40.000	40.795	-2.0	104	0.00
44	S	2-Fluorobiphenyl	80.000	80.371	-0.5	102	0.00

Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA G\Data\BG021116\
Data File : BG020874.D
Acq On : 12 Feb 2016 2:38
Operator : SJ/UM
Sample : SSTDCCC040
Misc : LOO 20PPM
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 12 04:51:44 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	40.217	-0.5	103	0.00
46		2-Chloronaphthalene	40.000	40.128	-0.3	103	0.00
47		2-Nitroaniline	40.000	40.109	-0.3	102	0.00
48		Acenaphthylene	40.000	40.791	-2.0	105	0.00
49		Dimethylphthalate	40.000	40.938	-2.3	104	0.00
50		2,6-Dinitrotoluene	40.000	41.446	-3.6	105	0.00
51	C	Acenaphthene	40.000	40.810	-2.0	105	0.00
52		3-Nitroaniline	40.000	41.754	-4.4	106	0.00
53	P	2,4-Dinitrophenol	40.000	35.850	10.4	96	0.00
54		Dibenzofuran	40.000	40.842	-2.1	106	0.00
55	P	4-Nitrophenol	40.000	41.837	-4.6	104	0.00
56		2,4-Dinitrotoluene	40.000	42.036	-5.1	105	0.00
57		Fluorene	40.000	41.132	-2.8	106	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	41.330	-3.3	106	0.00
59		Diethylphthalate	40.000	41.376	-3.4	107	0.00
60		4-Chlorophenyl-phenylether	40.000	40.327	-0.8	106	0.00
61		4-Nitroaniline	40.000	41.758	-4.4	106	0.00
62		Azobenzene	40.000	40.455	-1.1	104	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	105	0.00
64		4,6-Dinitro-2-methylphenol	40.000	37.839	5.4	101	0.00
65	c	n-Nitrosodiphenylamine	40.000	41.209	-3.0	106	0.00
66		4-Bromophenyl-phenylether	40.000	41.076	-2.7	106	0.00
67		Hexachlorobenzene	40.000	40.421	-1.1	105	0.00
68		Atrazine	40.000	40.679	-1.7	106	0.00
69	C	Pentachlorophenol	40.000	38.902	2.7	101	0.00
70		Phenanthrene	40.000	40.998	-2.5	107	0.00
71		Anthracene	40.000	40.722	-1.8	106	0.00
72		Carbazole	40.000	40.649	-1.6	106	0.00
73		Di-n-butylphthalate	40.000	40.989	-2.5	106	0.00
74	C	Fluoranthene	40.000	40.646	-1.6	106	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	105	0.00
76		Benzidine	40.000	39.250	1.9	100	0.00
77		Pyrene	40.000	41.045	-2.6	108	0.00
78	S	Terphenyl-d14	80.000	82.231	-2.8	108	0.00
79		Butylbenzylphthalate	40.000	40.661	-1.7	105	0.00
80		Benzo(a)anthracene	40.000	40.780	-2.0	108	0.00
81		3,3'-Dichlorobenzidine	40.000	40.910	-2.3	107	0.00
82		Chrysene	40.000	41.032	-2.6	108	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	40.926	-2.3	107	0.00
84	c	Di-n-octyl phthalate	40.000	41.204	-3.0	106	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	41.942	-4.9	109	-0.01
86	I	Perylene-d12	20.000	20.000	0.0	107	0.00
87		Benzo(b)fluoranthene	40.000	41.176	-2.9	109	0.00

Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA G\Data\BG021116\
Data File : BG020874.D
Acq On : 12 Feb 2016 2:38
Operator : SJ/UM
Sample : SSTDCCC040
Misc : LOO 20PPM
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 12 04:51:44 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	40.947	-2.4	109	0.00
89 C	Benzo(a)pyrene	40.000	41.154	-2.9	109	-0.01
90	Dibenzo(a,h)anthracene	40.000	41.171	-2.9	110	-0.01
91	Benzo(g,h,i)perylene	40.000	40.735	-1.8	108	-0.03

(#) = Out of Range SPCC's out = 0 CCC's out = 0