Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG021916\
Data File : BG020988.D
Acq On : 20 Feb 2016 1:30
Operator : SJ/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 20 02:40:48 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	117	0.00
2		1,4-Dioxane	40.000	37.523	6.2	113	-0.02
3		Pyridine	40.000	39.194	2.0	111	-0.01
4		n-Nitrosodimethylamine	40.000	39.740	0.6	113	-0.01
5	S	2-Fluorophenol	80.000	79.032	1.2	115	-0.01
6		Aniline	40.000	39.608	1.0	115	0.00
7	S	Phenol-d6	80.000	82.162	-2.7	119	0.00
8		2-Chlorophenol	40.000	40.529	-1.3	118	-0.01
9		Benzaldehyde	40.000	37.618	6.0	109	0.00
10	C	Phenol	40.000	40.552	-1.4	118	0.00
11		bis(2-Chloroethyl)ether	40.000	39.755	0.6	114	-0.01
12		1,3-Dichlorobenzene	40.000	40.238	-0.6	118	-0.01
13	C	1,4-Dichlorobenzene	40.000	40.833	-2.1	119	-0.01
14		1,2-Dichlorobenzene	40.000	40.369	-0.9	117	-0.01
15		Benzyl Alcohol	40.000	39.594	1.0	114	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	38.439	3.9	112	-0.01
17		2-Methylphenol	40.000	40.288	-0.7	117	0.00
18		Hexachloroethane	40.000	40.426	-1.1	117	-0.01
19	P	n-Nitroso-di-n-propylamine	40.000	39.612	1.0	116	-0.01
20		3+4-Methylphenols	40.000	40.505	-1.3	119	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	119	-0.01
22		Acetophenone	40.000	39.578	1.1	116	-0.01
23	S	Nitrobenzene-d5	80.000	80.163	-0.2	117	0.00
24		Nitrobenzene	40.000	40.138	-0.3	119	0.00
25		Isophorone	40.000	39.369	1.6	115	0.00
26	C	2-Nitrophenol	40.000	44.252	-10.6	125	-0.01
27		2,4-Dimethylphenol	40.000	39.393	1.5	113	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.170	-0.4	118	-0.01
29	C	2,4-Dichlorophenol	40.000	42.227	-5.6	126	-0.01
30		1,2,4-Trichlorobenzene	40.000	41.457	-3.6	122	-0.01
31		Naphthalene	40.000	40.289	-0.7	119	-0.01
32		Benzoic acid	40.000	40.424	-1.1	122	0.00
33		4-Chloroaniline	40.000	40.523	-1.3	120	0.00
34	C	Hexachlorobutadiene	40.000	41.656	-4.1	123	-0.02
35		Caprolactam	40.000	39.854	0.4	119	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.225	-0.6	120	0.00
37		2-Methylnaphthalene	40.000	40.556	-1.4	121	-0.01
38	I	Acenaphthene-d10	20.000	20.000	0.0	120	-0.01
39		1,2,4,5-Tetrachlorobenzene	40.000	41.328	-3.3	123	-0.01
40	P	Hexachlorocyclopentadiene	40.000	35.075	12.3	102	-0.01
41	S	2,4,6-Tribromophenol	80.000	84.442	-5.6	126	0.00
42	C	2,4,6-Trichlorophenol	40.000	42.676	-6.7	124	-0.01
43		2,4,5-Trichlorophenol	40.000	41.856	-4.6	124	0.00
44	S	2-Fluorobiphenyl	80.000	81.889	-2.4	121	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG021916\
Data File : BG020988.D
Acq On : 20 Feb 2016 1:30
Operator : SJ/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 20 02:40:48 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	40.445	-1.1	120	-0.01
46		2-Chloronaphthalene	40.000	40.698	-1.7	122	-0.01
47		2-Nitroaniline	40.000	40.430	-1.1	119	-0.01
48		Acenaphthylene	40.000	40.925	-2.3	122	-0.01
49		Dimethylphthalate	40.000	40.976	-2.4	122	-0.01
50		2,6-Dinitrotoluene	40.000	40.968	-2.4	121	-0.01
51	C	Acenaphthene	40.000	40.808	-2.0	122	-0.02
52		3-Nitroaniline	40.000	40.220	-0.5	119	0.00
53	P	2,4-Dinitrophenol	40.000	42.353	-5.9	138	0.00
54		Dibenzofuran	40.000	40.418	-1.0	122	-0.01
55	P	4-Nitrophenol	40.000	41.694	-4.2	121	0.00
56		2,4-Dinitrotoluene	40.000	42.927	-7.3	125	-0.01
57		Fluorene	40.000	40.092	-0.2	121	-0.01
58		2,3,4,6-Tetrachlorophenol	40.000	41.430	-3.6	123	-0.01
59		Diethylphthalate	40.000	40.397	-1.0	122	-0.01
60		4-Chlorophenyl-phenylether	40.000	40.896	-2.2	125	-0.01
61		4-Nitroaniline	40.000	40.569	-1.4	120	0.00
62		Azobenzene	40.000	38.913	2.7	117	-0.01
63	I	Phenanthrene-d10	20.000	20.000	0.0	121	-0.01
64		4,6-Dinitro-2-methylphenol	40.000	42.507	-6.3	134	0.00
65	c	n-Nitrosodiphenylamine	40.000	41.319	-3.3	122	-0.02
66		4-Bromophenyl-phenylether	40.000	41.656	-4.1	123	-0.02
67		Hexachlorobenzene	40.000	41.481	-3.7	125	-0.01
68		Atrazine	40.000	40.297	-0.7	121	-0.01
69	C	Pentachlorophenol	40.000	41.990	-5.0	126	-0.01
70		Phenanthrene	40.000	40.671	-1.7	122	-0.01
71		Anthracene	40.000	40.363	-0.9	121	-0.01
72		Carbazole	40.000	40.538	-1.3	121	-0.01
73		Di-n-butylphthalate	40.000	40.562	-1.4	121	-0.02
74	C	Fluoranthene	40.000	40.969	-2.4	123	-0.01
75	I	Chrysene-d12	20.000	20.000	0.0	123	-0.02
76		Benzidine	40.000	33.947	15.1	101	-0.01
77		Pyrene	40.000	40.488	-1.2	124	-0.01
78	S	Terphenyl-d14	80.000	81.849	-2.3	125	-0.02
79		Butylbenzylphthalate	40.000	40.805	-2.0	123	-0.02
80		Benzo(a)anthracene	40.000	40.781	-2.0	126	-0.02
81		3,3'-Dichlorobenzidine	40.000	40.223	-0.6	122	-0.02
82		Chrysene	40.000	40.415	-1.0	124	-0.02
83		Bis(2-ethylhexyl)phthalate	40.000	40.259	-0.6	122	-0.02
84	c	Di-n-octyl phthalate	40.000	40.697	-1.7	123	-0.03
85		Indeno(1,2,3-cd)pyrene	40.000	41.362	-3.4	126	-0.05
86	I	Perylene-d12	20.000	20.000	0.0	124	-0.03
87		Benzo(b)fluoranthene	40.000	40.535	-1.3	125	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG021916\
Data File : BG020988.D
Acq On : 20 Feb 2016 1:30
Operator : SJ/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 20 02:40:48 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	40.842	-2.1	127	-0.02
89 C	Benzo(a)pyrene	40.000	40.612	-1.5	126	-0.03
90	Dibenzo(a,h)anthracene	40.000	40.394	-1.0	126	-0.05
91	Benzo(g,h,i)perylene	40.000	39.989	0.0	124	-0.06

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