Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122116\
Data File : BG025364.D
Acq On : 21 Dec 2016 19:25
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 22 02:43:29 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 21 18:42:01 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	101	0.00
2		1,4-Dioxane	40.000	40.811	-2.0	102	0.00
3		Pyridine	40.000	43.524	-8.8	101	0.00
4		n-Nitrosodimethylamine	40.000	43.152	-7.9	100	0.00
5	S	2-Fluorophenol	80.000	87.224	-9.0	107	0.00
6		Aniline	40.000	43.880	-9.7	105	0.00
7	S	Phenol-d6	80.000	88.460	-10.6	104	0.00
8		2-Chlorophenol	40.000	42.765	-6.9	103	0.00
9		Benzaldehyde	40.000	40.634	-1.6	101	0.00
10	C	Phenol	40.000	42.867	-7.2	102	0.00
11		bis(2-Chloroethyl)ether	40.000	42.288	-5.7	103	0.00
12		1,3-Dichlorobenzene	40.000	43.555	-8.9	105	0.00
13	C	1,4-Dichlorobenzene	40.000	41.656	-4.1	101	0.00
14		1,2-Dichlorobenzene	40.000	41.260	-3.1	101	0.00
15		Benzyl Alcohol	40.000	44.625	-11.6	102	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	41.410	-3.5	100	0.00
17		2-Methylphenol	40.000	44.805	-12.0	107	0.00
18		Hexachloroethane	40.000	43.470	-8.7	101	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	42.064	-5.2	101	0.00
20		3+4-Methylphenols	40.000	43.625	-9.1	103	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	105	0.00
22		Acetophenone	40.000	39.720	0.7	96	0.00
23	S	Nitrobenzene-d5	80.000	83.666	-4.6	100	0.00
24		Nitrobenzene	40.000	40.795	-2.0	99	0.00
25		Isophorone	40.000	41.865	-4.7	100	0.00
26	C	2-Nitrophenol	40.000	41.960	-4.9	100	0.00
27		2,4-Dimethylphenol	40.000	40.827	-2.1	98	0.00
28		bis(2-Chloroethoxy)methane	40.000	41.406	-3.5	105	0.00
29	C	2,4-Dichlorophenol	40.000	42.612	-6.5	103	0.00
30		1,2,4-Trichlorobenzene	40.000	41.387	-3.5	103	0.00
31		Naphthalene	40.000	41.454	-3.6	101	0.00
32		Benzoic acid	40.000	42.283	-5.7	100	0.00
33		4-Chloroaniline	40.000	43.192	-8.0	101	0.00
34	C	Hexachlorobutadiene	40.000	41.400	-3.5	102	0.00
35		Caprolactam	40.000	41.963	-4.9	102	0.00
36	C	4-Chloro-3-methylphenol	40.000	42.048	-5.1	100	0.00
37		2-Methylnaphthalene	40.000	41.704	-4.3	101	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	101	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	43.120	-7.8	103	0.00
40	P	Hexachlorocyclopentadiene	40.000	40.457	-1.1	99	0.00
41	S	2,4,6-Tribromophenol	80.000	81.298	-1.6	96	0.00
42	C	2,4,6-Trichlorophenol	40.000	42.095	-5.2	100	0.00
43		2,4,5-Trichlorophenol	40.000	39.307	1.7	94	0.00
44	S	2-Fluorobiphenyl	80.000	82.351	-2.9	101	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122116\
Data File : BG025364.D
Acq On : 21 Dec 2016 19:25
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 22 02:43:29 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 21 18:42:01 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	41.787	-4.5	101	0.00
46		2-Chloronaphthalene	40.000	41.036	-2.6	100	0.00
47		2-Nitroaniline	40.000	43.358	-8.4	101	0.00
48		Acenaphthylene	40.000	41.110	-2.8	100	0.00
49		Dimethylphthalate	40.000	41.187	-3.0	98	0.00
50		2,6-Dinitrotoluene	40.000	42.147	-5.4	101	0.00
51	C	Acenaphthene	40.000	40.646	-1.6	98	0.00
52		3-Nitroaniline	40.000	42.258	-5.6	99	0.00
53	P	2,4-Dinitrophenol	40.000	43.311	-8.3	105	0.00
54		Dibenzofuran	40.000	41.522	-3.8	99	0.00
55	P	4-Nitrophenol	40.000	42.783	-7.0	98	0.00
56		2,4-Dinitrotoluene	40.000	42.608	-6.5	101	0.00
57		Fluorene	40.000	41.290	-3.2	100	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	41.152	-2.9	99	0.00
59		Diethylphthalate	40.000	41.106	-2.8	99	0.00
60		4-Chlorophenyl-phenylether	40.000	40.310	-0.8	97	0.00
61		4-Nitroaniline	40.000	45.128	-12.8	98	0.00
62		Azobenzene	40.000	41.790	-4.5	99	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	98	0.00
64		4,6-Dinitro-2-methylphenol	40.000	44.670	-11.7	95	0.00
65	c	n-Nitrosodiphenylamine	40.000	42.412	-6.0	99	0.00
66		4-Bromophenyl-phenylether	40.000	41.790	-4.5	96	0.00
67		Hexachlorobenzene	40.000	41.046	-2.6	97	0.00
68		Atrazine	40.000	40.668	-1.7	94	0.00
69	C	Pentachlorophenol	40.000	42.314	-5.8	94	0.00
70		Phenanthrene	40.000	42.020	-5.1	98	0.00
71		Anthracene	40.000	42.319	-5.8	100	0.00
72		Carbazole	40.000	42.490	-6.2	99	0.00
73		Di-n-butylphthalate	40.000	42.294	-5.7	99	0.00
74	C	Fluoranthene	40.000	42.841	-7.1	101	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	106	0.00
76		Benzidine	40.000	40.123	-0.3	100	0.00
77		Pyrene	40.000	42.321	-5.8	103	0.00
78	S	Terphenyl-d14	80.000	81.999	-2.5	101	0.00
79		Butylbenzylphthalate	40.000	41.729	-4.3	105	0.00
80		Benzo(a)anthracene	40.000	42.223	-5.6	104	0.00
81		3,3'-Dichlorobenzidine	40.000	41.704	-4.3	102	0.00
82		Chrysene	40.000	41.995	-5.0	105	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	41.958	-4.9	106	0.00
84	c	Di-n-octyl phthalate	40.000	42.814	-7.0	106	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	42.785	-7.0	105	-0.01
86	I	Perylene-d12	20.000	20.000	0.0	106	-0.01
87		Benzo(b)fluoranthene	40.000	41.790	-4.5	106	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122116\
Data File : BG025364.D
Acq On : 21 Dec 2016 19:25
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 22 02:43:29 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 21 18:42:01 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	41.483	-3.7	103	0.00
89 C	Benzo(a)pyrene	40.000	41.629	-4.1	105	0.00
90	Dibenzo(a,h)anthracene	40.000	42.119	-5.3	105	0.00
91	Benzo(g,h,i)perylene	40.000	41.515	-3.8	105	0.00

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