Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG120715\
Data File : BG019979.D
Acq On : 7 Dec 2015 12:11
Operator : UM/NP
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 08 00:04:38 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 02 18:50:01 2015
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	-0.02
2		1,4-Dioxane	40.000	44.069	-10.2	112	0.00
3		Pyridine	40.000	43.196	-8.0	108	-0.02
4		n-Nitrosodimethylamine	40.000	38.448	3.9	92	-0.01
5	S	2-Fluorophenol	80.000	87.609	-9.5	111	-0.01
6		Aniline	40.000	41.432	-3.6	103	-0.02
7	S	Phenol-d6	80.000	86.038	-7.5	108	-0.01
8		2-Chlorophenol	40.000	42.194	-5.5	106	-0.02
9		Benzaldehyde	40.000	37.285	6.8	95	-0.02
10	C	Phenol	40.000	43.300	-8.2	107	0.00
11		bis(2-Chloroethyl)ether	40.000	42.597	-6.5	108	-0.02
12		1,3-Dichlorobenzene	40.000	41.363	-3.4	105	-0.02
13	C	1,4-Dichlorobenzene	40.000	41.444	-3.6	106	-0.02
14		1,2-Dichlorobenzene	40.000	40.708	-1.8	103	-0.02
15		Benzyl Alcohol	40.000	38.863	2.8	99	-0.02
16		2,2'-oxybis(1-Chloropropane	40.000	41.180	-2.9	105	-0.02
17		2-Methylphenol	40.000	41.810	-4.5	104	-0.02
18		Hexachloroethane	40.000	41.084	-2.7	102	-0.02
19	P	n-Nitroso-di-n-propylamine	40.000	38.063	4.8	97	-0.02
20		3+4-Methylphenols	40.000	41.059	-2.6	100	-0.01
21	I	Naphthalene-d8	20.000	20.000	0.0	97	-0.02
22		Acetophenone	40.000	41.448	-3.6	99	-0.02
23	S	Nitrobenzene-d5	80.000	86.723	-8.4	103	-0.02
24		Nitrobenzene	40.000	43.370	-8.4	102	-0.02
25		Isophorone	40.000	39.866	0.3	96	-0.02
26	C	2-Nitrophenol	40.000	48.482	-21.2#	114	-0.01
27		2,4-Dimethylphenol	40.000	41.285	-3.2	100	-0.02
28		bis(2-Chloroethoxy)methane	40.000	42.685	-6.7	103	-0.02
29	C	2,4-Dichlorophenol	40.000	43.247	-8.1	102	-0.02
30		1,2,4-Trichlorobenzene	40.000	41.658	-4.1	99	-0.02
31		Naphthalene	40.000	41.978	-4.9	100	-0.02
32		Benzoic acid	40.000	48.616	-21.5	129	-0.01
33		4-Chloroaniline	40.000	40.826	-2.1	98	-0.02
34	C	Hexachlorobutadiene	40.000	39.622	0.9	95	-0.02
35		Caprolactam	40.000	37.940	5.2	94	-0.03
36	C	4-Chloro-3-methylphenol	40.000	39.543	1.1	96	-0.01
37		2-Methylnaphthalene	40.000	39.852	0.4	96	-0.02
38	I	Acenaphthene-d10	20.000	20.000	0.0	89	-0.02
39		1,2,4,5-Tetrachlorobenzene	40.000	40.882	-2.2	92	-0.02
40	P	Hexachlorocyclopentadiene	40.000	35.599	11.0	78	-0.02
41	S	2,4,6-Tribromophenol	80.000	77.001	3.7	87	-0.02
42	C	2,4,6-Trichlorophenol	40.000	40.553	-1.4	93	-0.02
43		2,4,5-Trichlorophenol	40.000	40.621	-1.6	94	-0.02
44	S	2-Fluorobiphenyl	80.000	81.990	-2.5	93	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG120715\
Data File : BG019979.D
Acq On : 7 Dec 2015 12:11
Operator : UM/NP
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 08 00:04:38 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 02 18:50:01 2015
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	40.872	-2.2	92	-0.02
46		2-Chloronaphthalene	40.000	40.913	-2.3	92	-0.02
47		2-Nitroaniline	40.000	45.474	-13.7	100	-0.02
48		Acenaphthylene	40.000	40.219	-0.5	91	-0.02
49		Dimethylphthalate	40.000	38.004	5.0	87	-0.02
50		2,6-Dinitrotoluene	40.000	44.251	-10.6	96	-0.02
51	C	Acenaphthene	40.000	39.737	0.7	89	-0.02
52		3-Nitroaniline	40.000	44.738	-11.8	98	-0.02
53	P	2,4-Dinitrophenol	40.000	33.867	15.3	80	-0.02
54		Dibenzofuran	40.000	39.512	1.2	90	-0.02
55	P	4-Nitrophenol	40.000	41.778	-4.4	92	0.00
56		2,4-Dinitrotoluene	40.000	45.137	-12.8	96	-0.01
57		Fluorene	40.000	38.498	3.8	88	-0.02
58		2,3,4,6-Tetrachlorophenol	40.000	39.123	2.2	86	-0.01
59		Diethylphthalate	40.000	36.211	9.5	84	-0.02
60		4-Chlorophenyl-phenylether	40.000	37.899	5.3	86	-0.02
61		4-Nitroaniline	40.000	41.951	-4.9	92	-0.02
62		Azobenzene	40.000	37.919	5.2	86	-0.02
63	I	Phenanthrene-d10	20.000	20.000	0.0	82	-0.02
64		4,6-Dinitro-2-methylphenol	40.000	42.402	-6.0	92	-0.02
65	c	n-Nitrosodiphenylamine	40.000	41.254	-3.1	85	-0.02
66		4-Bromophenyl-phenylether	40.000	40.627	-1.6	83	-0.02
67		Hexachlorobenzene	40.000	40.504	-1.3	84	-0.02
68		Atrazine	40.000	38.457	3.9	78	-0.02
69	C	Pentachlorophenol	40.000	45.213	-13.0	92	-0.02
70		Phenanthrene	40.000	41.025	-2.6	85	-0.02
71		Anthracene	40.000	40.899	-2.2	84	-0.02
72		Carbazole	40.000	40.960	-2.4	84	-0.02
73		Di-n-butylphthalate	40.000	41.282	-3.2	85	-0.02
74	C	Fluoranthene	40.000	40.834	-2.1	84	-0.02
75	I	Chrysene-d12	20.000	20.000	0.0	83	-0.03
76		Benzidine	40.000	36.999	7.5	75	-0.02
77		Pyrene	40.000	41.064	-2.7	84	-0.02
78	S	Terphenyl-d14	80.000	80.784	-1.0	82	-0.02
79		Butylbenzylphthalate	40.000	42.475	-6.2	87	-0.03
80		Benzo(a)anthracene	40.000	40.346	-0.9	84	-0.02
81		3,3'-Dichlorobenzidine	40.000	38.056	4.9	78	-0.03
82		Chrysene	40.000	40.437	-1.1	84	-0.03
83		Bis(2-ethylhexyl)phthalate	40.000	41.690	-4.2	87	-0.04
84	c	Di-n-octyl phthalate	40.000	43.414	-8.5	88	-0.06
85		Indeno(1,2,3-cd)pyrene	40.000	42.771	-6.9	89	-0.09
86	I	Perylene-d12	20.000	20.000	0.0	85	-0.05
87		Benzo(b)fluoranthene	40.000	40.219	-0.5	86	-0.05

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG120715\
Data File : BG019979.D
Acq On : 7 Dec 2015 12:11
Operator : UM/NP
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 08 00:04:38 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 02 18:50:01 2015
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	40.301	-0.8	87	-0.05
89 C	Benzo(a)pyrene	40.000	40.387	-1.0	87	-0.05
90	Dibenzo(a,h)anthracene	40.000	41.140	-2.9	89	-0.08
91	Benzo(g,h,i)perylene	40.000	41.087	-2.7	88	-0.09

(#) = Out of Range SPCC's out = 0 CCC's out = 1