Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG121215\
Data File : BG020127.D
Acq On : 12 Dec 2015 4:35
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 12 05:30:11 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	105	-0.04
2		1,4-Dioxane	40.000	44.717	-11.8	118	-0.02
3		Pyridine	40.000	42.822	-7.1	112	-0.04
4		n-Nitrosodimethylamine	40.000	36.482	8.8	91	-0.02
5	S	2-Fluorophenol	80.000	83.921	-4.9	110	-0.02
6		Aniline	40.000	40.696	-1.7	105	-0.03
7	S	Phenol-d6	80.000	83.823	-4.8	110	-0.02
8		2-Chlorophenol	40.000	41.575	-3.9	108	-0.03
9		Benzaldehyde	40.000	34.938	12.7	92	-0.03
10	C	Phenol	40.000	41.977	-4.9	108	-0.02
11		bis(2-Chloroethyl)ether	40.000	41.662	-4.2	110	-0.03
12		1,3-Dichlorobenzene	40.000	40.396	-1.0	107	-0.03
13	C	1,4-Dichlorobenzene	40.000	39.455	1.4	105	-0.03
14		1,2-Dichlorobenzene	40.000	40.561	-1.4	107	-0.04
15		Benzyl Alcohol	40.000	39.012	2.5	104	-0.03
16		2,2'-oxybis(1-Chloropropane	40.000	38.334	4.2	101	-0.04
17		2-Methylphenol	40.000	40.897	-2.2	105	-0.03
18		Hexachloroethane	40.000	39.254	1.9	101	-0.04
19	P	n-Nitroso-di-n-propylamine	40.000	38.830	2.9	102	-0.04
20		3+4-Methylphenols	40.000	41.408	-3.5	105	-0.02
21	I	Naphthalene-d8	20.000	20.000	0.0	104	-0.04
22		Acetophenone	40.000	40.952	-2.4	105	-0.04
23	S	Nitrobenzene-d5	80.000	82.844	-3.6	106	-0.03
24		Nitrobenzene	40.000	41.270	-3.2	104	-0.03
25		Isophorone	40.000	39.262	1.8	101	-0.04
26	C	2-Nitrophenol	40.000	47.909	-19.8	121	-0.03
27		2,4-Dimethylphenol	40.000	40.581	-1.5	105	-0.03
28		bis(2-Chloroethoxy)methane	40.000	41.026	-2.6	106	-0.04
29	C	2,4-Dichlorophenol	40.000	41.541	-3.9	105	-0.03
30		1,2,4-Trichlorobenzene	40.000	39.310	1.7	100	-0.04
31		Naphthalene	40.000	40.302	-0.8	103	-0.04
32		Benzoic acid	40.000	55.557	-38.9#	161	0.00
33		4-Chloroaniline	40.000	40.411	-1.0	105	-0.04
34	C	Hexachlorobutadiene	40.000	38.586	3.5	100	-0.04
35		Caprolactam	40.000	41.143	-2.9	109	-0.04
36	C	4-Chloro-3-methylphenol	40.000	40.065	-0.2	105	-0.02
37		2-Methylnaphthalene	40.000	39.282	1.8	102	-0.04
38	I	Acenaphthene-d10	20.000	20.000	0.0	103	-0.04
39		1,2,4,5-Tetrachlorobenzene	40.000	39.034	2.4	102	-0.04
40	P	Hexachlorocyclopentadiene	40.000	41.105	-2.8	103	-0.04
41	S	2,4,6-Tribromophenol	80.000	80.706	-0.9	105	-0.03
42	C	2,4,6-Trichlorophenol	40.000	39.988	0.0	105	-0.03
43		2,4,5-Trichlorophenol	40.000	39.651	0.9	105	-0.03
44	S	2-Fluorobiphenyl	80.000	76.304	4.6	100	-0.04

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG121215\
Data File : BG020127.D
Acq On : 12 Dec 2015 4:35
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 12 05:30:11 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	38.213	4.5	99	-0.04
46		2-Chloronaphthalene	40.000	37.881	5.3	98	-0.03
47		2-Nitroaniline	40.000	41.661	-4.2	106	-0.03
48		Acenaphthylene	40.000	37.701	5.7	98	-0.04
49		Dimethylphthalate	40.000	37.928	5.2	100	-0.04
50		2,6-Dinitrotoluene	40.000	43.238	-8.1	107	-0.03
51	C	Acenaphthene	40.000	36.836	7.9	95	-0.04
52		3-Nitroaniline	40.000	44.046	-10.1	112	-0.03
53	P	2,4-Dinitrophenol	40.000	51.941	-29.9#	160	-0.03
54		Dibenzofuran	40.000	37.934	5.2	99	-0.03
55	P	4-Nitrophenol	40.000	42.403	-6.0	107	-0.02
56		2,4-Dinitrotoluene	40.000	44.979	-12.4	110	-0.03
57		Fluorene	40.000	36.874	7.8	97	-0.04
58		2,3,4,6-Tetrachlorophenol	40.000	40.331	-0.8	102	-0.03
59		Diethylphthalate	40.000	36.442	8.9	97	-0.04
60		4-Chlorophenyl-phenylether	40.000	37.130	7.2	97	-0.04
61		4-Nitroaniline	40.000	41.469	-3.7	104	-0.03
62		Azobenzene	40.000	35.485	11.3	93	-0.04
63	I	Phenanthrene-d10	20.000	20.000	0.0	98	-0.04
64		4,6-Dinitro-2-methylphenol	40.000	49.204	-23.0	131	-0.03
65	c	n-Nitrosodiphenylamine	40.000	39.675	0.8	98	-0.04
66		4-Bromophenyl-phenylether	40.000	38.442	3.9	94	-0.04
67		Hexachlorobenzene	40.000	39.987	0.0	99	-0.04
68		Atrazine	40.000	38.728	3.2	94	-0.04
69	C	Pentachlorophenol	40.000	44.824	-12.1	109	-0.04
70		Phenanthrene	40.000	38.932	2.7	96	-0.04
71		Anthracene	40.000	39.431	1.4	97	-0.04
72		Carbazole	40.000	39.296	1.8	96	-0.03
73		Di-n-butylphthalate	40.000	38.404	4.0	94	-0.05
74	C	Fluoranthene	40.000	39.286	1.8	96	-0.04
75	I	Chrysene-d12	20.000	20.000	0.0	94	-0.05
76		Benzidine	40.000	36.175	9.6	82	-0.04
77		Pyrene	40.000	41.627	-4.1	96	-0.04
78	S	Terphenyl-d14	80.000	80.677	-0.8	92	-0.04
79		Butylbenzylphthalate	40.000	38.852	2.9	89	-0.05
80		Benzo(a)anthracene	40.000	38.832	2.9	91	-0.05
81		3,3'-Dichlorobenzidine	40.000	37.442	6.4	86	-0.05
82		Chrysene	40.000	38.641	3.4	90	-0.05
83		Bis(2-ethylhexyl)phthalate	40.000	38.010	5.0	89	-0.07
84	c	Di-n-octyl phthalate	40.000	39.856	0.4	91	-0.10
85		Indeno(1,2,3-cd)pyrene	40.000	41.978	-4.9	98	-0.15
86	I	Perylene-d12	20.000	20.000	0.0	96	-0.09
87		Benzo(b)fluoranthene	40.000	38.005	5.0	92	-0.08

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG121215\
Data File : BG020127.D
Acq On : 12 Dec 2015 4:35
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 12 05:30:11 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	38.932	2.7	95	-0.08
89 C	Benzo(a)pyrene	40.000	39.218	2.0	95	-0.08
90	Dibenzo(a,h)anthracene	40.000	40.044	-0.1	98	-0.15
91	Benzo(g,h,i)perylene	40.000	40.337	-0.8	97	-0.16

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