Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG120215\
Data File : BG019840.D
Acq On : 2 Dec 2015 18:41
Operator : UM/NP
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG120215

Quant Time: Dec 03 00:09:46 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	104	0.00
2		1,4-Dioxane	40.000	43.398	-8.5	114	0.00
3		Pyridine	40.000	44.431	-11.1	115	0.00
4		n-Nitrosodimethylamine	40.000	43.474	-8.7	107	0.00
5	S	2-Fluorophenol	80.000	85.530	-6.9	112	0.00
6		Aniline	40.000	42.900	-7.2	109	0.00
7	S	Phenol-d6	80.000	84.475	-5.6	110	0.00
8		2-Chlorophenol	40.000	41.977	-4.9	108	0.00
9		Benzaldehyde	40.000	42.755	-6.9	112	0.00
10	C	Phenol	40.000	42.449	-6.1	108	0.00
11		bis(2-Chloroethyl)ether	40.000	42.436	-6.1	111	0.00
12		1,3-Dichlorobenzene	40.000	41.281	-3.2	108	0.00
13	C	1,4-Dichlorobenzene	40.000	40.865	-2.2	108	0.00
14		1,2-Dichlorobenzene	40.000	41.762	-4.4	109	0.00
15		Benzyl Alcohol	40.000	41.739	-4.3	110	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	43.728	-9.3	114	0.00
17		2-Methylphenol	40.000	41.940	-4.8	107	0.00
18		Hexachloroethane	40.000	41.979	-4.9	107	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	42.717	-6.8	112	0.00
20		3+4-Methylphenols	40.000	42.816	-7.0	107	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	106	0.00
22		Acetophenone	40.000	42.010	-5.0	110	0.00
23	S	Nitrobenzene-d5	80.000	86.035	-7.5	112	0.00
24		Nitrobenzene	40.000	44.002	-10.0	113	0.00
25		Isophorone	40.000	42.371	-5.9	111	0.00
26	C	2-Nitrophenol	40.000	43.265	-8.2	111	0.00
27		2,4-Dimethylphenol	40.000	40.698	-1.7	108	0.00
28		bis(2-Chloroethoxy)methane	40.000	42.273	-5.7	112	0.00
29	C	2,4-Dichlorophenol	40.000	40.388	-1.0	104	0.00
30		1,2,4-Trichlorobenzene	40.000	40.710	-1.8	106	0.00
31		Naphthalene	40.000	41.247	-3.1	108	0.00
32		Benzoic acid	40.000	42.910	-7.3	122	0.00
33		4-Chloroaniline	40.000	40.339	-0.8	106	0.00
34	C	Hexachlorobutadiene	40.000	38.761	3.1	102	0.00
35		Caprolactam	40.000	41.360	-3.4	112	0.00
36	C	4-Chloro-3-methylphenol	40.000	39.961	0.1	106	0.00
37		2-Methylnaphthalene	40.000	40.374	-0.9	107	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	100	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	41.081	-2.7	103	0.00
40	P	Hexachlorocyclopentadiene	40.000	43.625	-9.1	106	0.00
41	S	2,4,6-Tribromophenol	80.000	77.848	2.7	98	0.00
42	C	2,4,6-Trichlorophenol	40.000	40.028	-0.1	102	0.00
43		2,4,5-Trichlorophenol	40.000	40.270	-0.7	103	0.00
44	S	2-Fluorobiphenyl	80.000	83.065	-3.8	105	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG120215\
Data File : BG019840.D
Acq On : 2 Dec 2015 18:41
Operator : UM/NP
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG120215

Quant Time: Dec 03 00:09:46 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	41.539	-3.8	104	0.00
46		2-Chloronaphthalene	40.000	41.876	-4.7	105	0.00
47		2-Nitroaniline	40.000	44.216	-10.5	109	0.00
48		Acenaphthylene	40.000	41.695	-4.2	105	0.00
49		Dimethylphthalate	40.000	40.504	-1.3	103	0.00
50		2,6-Dinitrotoluene	40.000	44.673	-11.7	107	0.00
51	C	Acenaphthene	40.000	41.787	-4.5	105	0.00
52		3-Nitroaniline	40.000	44.253	-10.6	109	0.00
53	P	2,4-Dinitrophenol	40.000	42.083	-5.2	119	0.00
54		Dibenzofuran	40.000	41.230	-3.1	104	0.00
55	P	4-Nitrophenol	40.000	42.939	-7.3	105	0.00
56		2,4-Dinitrotoluene	40.000	44.744	-11.9	106	0.00
57		Fluorene	40.000	40.695	-1.7	104	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	39.111	2.2	96	0.00
59		Diethylphthalate	40.000	40.376	-0.9	104	0.00
60		4-Chlorophenyl-phenylether	40.000	40.404	-1.0	102	0.00
61		4-Nitroaniline	40.000	42.904	-7.3	105	0.00
62		Azobenzene	40.000	41.864	-4.7	106	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	97	0.00
64		4,6-Dinitro-2-methylphenol	40.000	41.578	-3.9	107	0.00
65	c	n-Nitrosodiphenylamine	40.000	41.433	-3.6	102	0.00
66		4-Bromophenyl-phenylether	40.000	41.194	-3.0	100	0.00
67		Hexachlorobenzene	40.000	40.021	-0.1	98	0.00
68		Atrazine	40.000	40.546	-1.4	98	0.00
69	C	Pentachlorophenol	40.000	42.289	-5.7	102	0.00
70		Phenanthrene	40.000	41.199	-3.0	101	0.00
71		Anthracene	40.000	41.232	-3.1	101	0.00
72		Carbazole	40.000	41.843	-4.6	102	0.00
73		Di-n-butylphthalate	40.000	41.964	-4.9	102	0.00
74	C	Fluoranthene	40.000	41.383	-3.5	100	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	103	0.00
76		Benzidine	40.000	42.531	-6.3	105	0.00
77		Pyrene	40.000	40.055	-0.1	100	0.00
78	S	Terphenyl-d14	80.000	79.993	0.0	100	0.00
79		Butylbenzylphthalate	40.000	43.135	-7.8	108	0.00
80		Benzo(a)anthracene	40.000	40.612	-1.5	104	0.00
81		3,3'-Dichlorobenzidine	40.000	40.427	-1.1	102	0.00
82		Chrysene	40.000	41.157	-2.9	105	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	42.478	-6.2	109	0.00
84	c	Di-n-octyl phthalate	40.000	42.816	-7.0	107	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	41.561	-3.9	106	0.00
86	I	Perylene-d12	20.000	20.000	0.0	102	0.00
87		Benzo(b)fluoranthene	40.000	41.366	-3.4	107	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG120215\
Data File : BG019840.D
Acq On : 2 Dec 2015 18:41
Operator : UM/NP
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG120215

Quant Time: Dec 03 00:09:46 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	40.940	-2.3	107	0.00
89 C	Benzo(a)pyrene	40.000	41.510	-3.8	107	0.00
90	Dibenzo(a,h)anthracene	40.000	41.002	-2.5	107	0.00
91	Benzo(g,h,i)perylene	40.000	40.598	-1.5	105	0.00

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