Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA G\Data\BG111017\
Data File : BG030909.D
Acq On : 10 Nov 2017 15:40
Operator : SJ/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG111017

Quant Time: Nov 10 16:15:42 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG111017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 10 15:38:25 2017
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	97	0.00
2		1,4-Dioxane	40.000	42.121	-5.3	103	-0.02
3		Pyridine	40.000	42.322	-5.8	98	-0.02
4		n-Nitrosodimethylamine	40.000	41.434	-3.6	98	-0.02
5	S	2-Fluorophenol	80.000	81.574	-2.0	97	0.00
6		Aniline	40.000	40.704	-1.8	96	0.00
7	S	Phenol-d6	80.000	83.241	-4.1	97	0.00
8		2-Chlorophenol	40.000	41.329	-3.3	98	0.00
9		Benzaldehyde	40.000	40.297	-0.7	95	-0.02
10	C	Phenol	40.000	40.833	-2.1	96	0.00
11		bis(2-Chloroethyl)ether	40.000	41.306	-3.3	98	0.00
12		1,3-Dichlorobenzene	40.000	40.618	-1.5	98	0.00
13	C	1,4-Dichlorobenzene	40.000	40.695	-1.7	98	0.00
14		1,2-Dichlorobenzene	40.000	40.753	-1.9	97	0.00
15		Benzyl Alcohol	40.000	41.436	-3.6	96	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	41.060	-2.7	98	0.00
17		2-Methylphenol	40.000	40.748	-1.9	95	0.00
18		Hexachloroethane	40.000	40.518	-1.3	95	-0.02
19	P	n-Nitroso-di-n-propylamine	40.000	41.532	-3.8	96	0.00
20		3+4-Methylphenols	40.000	41.420	-3.6	96	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	97	0.00
22		Acetophenone	40.000	40.232	-0.6	94	0.00
23	S	Nitrobenzene-d5	80.000	81.020	-1.3	96	0.00
24		Nitrobenzene	40.000	40.559	-1.4	95	0.00
25		Isophorone	40.000	40.438	-1.1	96	0.00
26	C	2-Nitrophenol	40.000	41.750	-4.4	99	0.00
27		2,4-Dimethylphenol	40.000	41.578	-3.9	98	0.00
28		bis(2-Chloroethoxy)methane	40.000	41.291	-3.2	98	0.00
29	C	2,4-Dichlorophenol	40.000	41.829	-4.6	96	0.00
30		1,2,4-Trichlorobenzene	40.000	40.180	-0.4	95	0.00
31		Naphthalene	40.000	40.127	-0.3	97	0.00
32		Benzoic acid	40.000	39.354	1.6	97	0.00
33		4-Chloroaniline	40.000	41.019	-2.5	95	0.00
34	C	Hexachlorobutadiene	40.000	40.872	-2.2	99	0.00
35		Caprolactam	40.000	37.732	5.7	93	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.044	-0.1	94	0.00
37		2-Methylnaphthalene	40.000	40.490	-1.2	97	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	97	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	40.802	-2.0	97	0.00
40	P	Hexachlorocyclopentadiene	40.000	39.643	0.9	103	0.00
41	S	2,4,6-Tribromophenol	80.000	77.905	2.6	93	0.00
42	C	2,4,6-Trichlorophenol	40.000	40.923	-2.3	95	0.00
43		2,4,5-Trichlorophenol	40.000	40.047	-0.1	95	0.00
44	S	2-Fluorobiphenyl	80.000	81.352	-1.7	98	0.00

Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA G\Data\BG111017\
Data File : BG030909.D
Acq On : 10 Nov 2017 15:40
Operator : SJ/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG111017

Quant Time: Nov 10 16:15:42 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG111017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 10 15:38:25 2017
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	40.352	-0.9	97	0.00
46		2-Chloronaphthalene	40.000	40.382	-1.0	96	0.00
47		2-Nitroaniline	40.000	40.480	-1.2	96	0.00
48		Acenaphthylene	40.000	40.712	-1.8	97	0.00
49		Dimethylphthalate	40.000	39.514	1.2	95	0.00
50		2,6-Dinitrotoluene	40.000	40.470	-1.2	94	0.00
51	C	Acenaphthene	40.000	40.034	-0.1	96	0.00
52		3-Nitroaniline	40.000	39.717	0.7	93	0.00
53	P	2,4-Dinitrophenol	40.000	37.946	5.1	91	0.00
54		Dibenzofuran	40.000	39.774	0.6	94	0.00
55	P	4-Nitrophenol	40.000	40.474	-1.2	96	0.01
56		2,4-Dinitrotoluene	40.000	40.651	-1.6	96	0.00
57		Fluorene	40.000	39.235	1.9	95	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	40.156	-0.4	96	0.00
59		Diethylphthalate	40.000	39.725	0.7	97	0.00
60		4-Chlorophenyl-phenylether	40.000	39.965	0.1	96	0.00
61		4-Nitroaniline	40.000	39.574	1.1	95	0.00
62		Azobenzene	40.000	40.109	-0.3	96	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	95	0.00
64		4,6-Dinitro-2-methylphenol	40.000	42.365	-5.9	99	0.00
65	c	n-Nitrosodiphenylamine	40.000	40.362	-0.9	95	0.00
66		4-Bromophenyl-phenylether	40.000	40.632	-1.6	94	0.00
67		Hexachlorobenzene	40.000	40.351	-0.9	94	0.00
68		Atrazine	40.000	39.927	0.2	96	0.00
69	C	Pentachlorophenol	40.000	40.630	-1.6	96	0.00
70		Phenanthrene	40.000	40.179	-0.4	95	0.00
71		Anthracene	40.000	40.665	-1.7	97	0.00
72		Carbazole	40.000	40.187	-0.5	95	0.00
73		Di-n-butylphthalate	40.000	40.225	-0.6	96	0.00
74	C	Fluoranthene	40.000	40.495	-1.2	97	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	98	0.00
76		Benzidine	40.000	40.639	-1.6	95	0.00
77		Pyrene	40.000	41.309	-3.3	99	0.00
78	S	Terphenyl-d14	80.000	82.307	-2.9	99	0.00
79		Butylbenzylphthalate	40.000	40.927	-2.3	99	0.00
80		Benzo(a)anthracene	40.000	40.286	-0.7	98	0.00
81		3,3'-Dichlorobenzidine	40.000	40.659	-1.6	98	0.00
82		Chrysene	40.000	40.819	-2.0	99	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	40.385	-1.0	100	0.00
84	c	Di-n-octyl phthalate	40.000	40.891	-2.2	100	-0.02
85		Indeno(1,2,3-cd)pyrene	40.000	40.673	-1.7	98	0.00
86	I	Perylene-d12	20.000	20.000	0.0	99	0.00
87		Benzo(b)fluoranthene	40.000	40.506	-1.3	100	0.00

Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA G\Data\BG111017\
Data File : BG030909.D
Acq On : 10 Nov 2017 15:40
Operator : SJ/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ICVBG111017

Quant Time: Nov 10 16:15:42 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG111017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 10 15:38:25 2017
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	40.236	-0.6	99	0.00
89 C	Benzo(a)pyrene	40.000	40.200	-0.5	100	0.00
90	Dibenzo(a,h)anthracene	40.000	40.692	-1.7	101	0.00
91	Benzo(g,h,i)perylene	40.000	40.232	-0.6	100	0.00

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