Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071715\
Data File : BG017947.D
Acq On : 18 Jul 2015 4:45
Operator : TP/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 20 01:48:00 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG071715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Jul 18 01:50:30 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	109	0.00
2		1,4-Dioxane	40.000	35.919	10.2	104	0.00
3		Pyridine	40.000	37.090	7.3	101	0.00
4		n-Nitrosodimethylamine	40.000	35.249	11.9	100	0.00
5	S	2-Fluorophenol	80.000	76.102	4.9	106	0.00
6		Aniline	40.000	36.902	7.7	104	0.00
7	S	Phenol-d6	80.000	75.788	5.3	107	0.00
8		2-Chlorophenol	40.000	38.227	4.4	109	0.00
9		Benzaldehyde	40.000	37.411	6.5	106	0.00
10	C	Phenol	40.000	37.276	6.8	106	0.00
11		bis(2-Chloroethyl)ether	40.000	36.889	7.8	105	0.00
12		1,3-Dichlorobenzene	40.000	37.246	6.9	106	0.00
13	C	1,4-Dichlorobenzene	40.000	37.166	7.1	106	0.00
14		1,2-Dichlorobenzene	40.000	37.291	6.8	108	0.00
15		Benzyl Alcohol	40.000	37.973	5.1	108	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	35.185	12.0	103	0.00
17		2-Methylphenol	40.000	37.188	7.0	106	0.00
18		Hexachloroethane	40.000	37.151	7.1	106	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	37.288	6.8	106	0.00
20		3+4-Methylphenols	40.000	38.114	4.7	107	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	107	0.00
22		Acetophenone	40.000	37.860	5.4	106	0.00
23	S	Nitrobenzene-d5	80.000	78.042	2.4	107	0.00
24		Nitrobenzene	40.000	38.914	2.7	105	0.00
25		Isophorone	40.000	38.258	4.4	104	0.00
26	C	2-Nitrophenol	40.000	41.865	-4.7	117	0.00
27		2,4-Dimethylphenol	40.000	38.684	3.3	108	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.262	4.3	105	0.00
29	C	2,4-Dichlorophenol	40.000	41.352	-3.4	110	0.00
30		1,2,4-Trichlorobenzene	40.000	38.644	3.4	109	0.00
31		Naphthalene	40.000	38.174	4.6	106	0.00
32		Benzoic acid	40.000	48.143	-20.4	154	0.00
33		4-Chloroaniline	40.000	38.655	3.4	105	0.00
34	C	Hexachlorobutadiene	40.000	39.642	0.9	109	0.00
35		Caprolactam	40.000	39.286	1.8	110	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.119	-0.3	109	0.00
37		2-Methylnaphthalene	40.000	38.279	4.3	107	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	105	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	39.458	1.4	107	0.00
40	P	Hexachlorocyclopentadiene	40.000	40.036	-0.1	111	0.00
41	S	2,4,6-Tribromophenol	80.000	82.863	-3.6	115	0.00
42	C	2,4,6-Trichlorophenol	40.000	41.293	-3.2	114	0.00
43		2,4,5-Trichlorophenol	40.000	42.548	-6.4	114	0.00
44	S	2-Fluorobiphenyl	80.000	77.027	3.7	105	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071715\
Data File : BG017947.D
Acq On : 18 Jul 2015 4:45
Operator : TP/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 20 01:48:00 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG071715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Jul 18 01:50:30 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	39.023	2.4	106	0.00
46		2-Chloronaphthalene	40.000	39.082	2.3	107	0.00
47		2-Nitroaniline	40.000	39.897	0.3	109	0.00
48		Acenaphthylene	40.000	39.138	2.2	105	0.00
49		Dimethylphthalate	40.000	38.662	3.3	104	0.00
50		2,6-Dinitrotoluene	40.000	40.869	-2.2	111	0.00
51	C	Acenaphthene	40.000	38.429	3.9	106	0.00
52		3-Nitroaniline	40.000	41.941	-4.9	110	0.00
53	P	2,4-Dinitrophenol	40.000	43.236	-8.1	120	0.00
54		Dibenzofuran	40.000	38.571	3.6	105	0.00
55	P	4-Nitrophenol	40.000	40.327	-0.8	110	0.00
56		2,4-Dinitrotoluene	40.000	40.828	-2.1	111	0.00
57		Fluorene	40.000	38.474	3.8	106	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	40.367	-0.9	110	0.00
59		Diethylphthalate	40.000	39.155	2.1	106	0.00
60		4-Chlorophenyl-phenylether	40.000	38.841	2.9	106	0.00
61		4-Nitroaniline	40.000	39.570	1.1	109	0.00
62		Azobenzene	40.000	37.981	5.0	103	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	106	0.00
64		4,6-Dinitro-2-methylphenol	40.000	43.106	-7.8	119	0.00
65	c	n-Nitrosodiphenylamine	40.000	39.416	1.5	105	0.00
66		4-Bromophenyl-phenylether	40.000	39.821	0.4	108	0.00
67		Hexachlorobenzene	40.000	38.659	3.4	108	0.00
68		Atrazine	40.000	40.205	-0.5	106	0.00
69	C	Pentachlorophenol	40.000	43.780	-9.5	117	0.00
70		Phenanthrene	40.000	38.215	4.5	105	0.00
71		Anthracene	40.000	39.012	2.5	104	0.00
72		Carbazole	40.000	38.956	2.6	105	0.00
73		Di-n-butylphthalate	40.000	40.030	-0.1	106	0.00
74	C	Fluoranthene	40.000	38.856	2.9	105	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	104	0.00
76		Benzidine	40.000	41.963	-4.9	111	0.00
77		Pyrene	40.000	38.836	2.9	104	0.00
78	S	Terphenyl-d14	80.000	77.499	3.1	105	0.00
79		Butylbenzylphthalate	40.000	42.173	-5.4	108	0.00
80		Benzo(a)anthracene	40.000	39.356	1.6	105	0.00
81		3,3'-Dichlorobenzidine	40.000	40.887	-2.2	107	0.00
82		Chrysene	40.000	39.276	1.8	106	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	41.723	-4.3	107	0.00
84	c	Di-n-octyl phthalate	40.000	42.908	-7.3	108	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	39.432	1.4	106	0.00
86	I	Perylene-d12	20.000	20.000	0.0	107	0.00
87		Benzo(b)fluoranthene	40.000	39.819	0.5	109	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071715\
Data File : BG017947.D
Acq On : 18 Jul 2015 4:45
Operator : TP/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 20 01:48:00 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG071715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Jul 18 01:50:30 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	37.580	6.1	101	0.00
89 C	Benzo(a)pyrene	40.000	38.894	2.8	105	0.00
90	Dibenzo(a,h)anthracene	40.000	39.009	2.5	108	0.00
91	Benzo(g,h,i)perylene	40.000	38.590	3.5	106	0.00

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