Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080416\
Data File : BF089607.D
Acq On : 5 Aug 2016 1:52
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 05 05:48:21 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF080416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 05 01:39:23 2016
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	98	0.00
2		1,4-Dioxane	40.000	35.842	10.4	103	0.00
3		Pyridine	40.000	42.136	-5.3	107	0.00
4		n-Nitrosodimethylamine	40.000	38.005	5.0	93	0.00
5	S	2-Fluorophenol	80.000	80.771	-1.0	101	0.00
6		Aniline	40.000	41.606	-4.0	101	0.00
7	S	Phenol-d6	80.000	80.710	-0.9	101	0.00
8		2-Chlorophenol	40.000	40.638	-1.6	102	0.00
9		Benzaldehyde	40.000	46.277	-15.7	106	0.00
10	C	Phenol	40.000	41.516	-3.8	100	0.00
11		bis(2-Chloroethyl)ether	40.000	39.690	0.8	102	0.00
12		1,3-Dichlorobenzene	40.000	39.222	1.9	101	0.00
13	C	1,4-Dichlorobenzene	40.000	38.354	4.1	101	0.00
14		1,2-Dichlorobenzene	40.000	38.211	4.5	103	0.00
15		Benzyl Alcohol	40.000	40.374	-0.9	97	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	38.329	4.2	100	0.00
17		2-Methylphenol	40.000	40.513	-1.3	102	-0.01
18		Hexachloroethane	40.000	38.069	4.8	100	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.941	0.1	100	0.00
20		3+4-Methylphenols	40.000	41.418	-3.5	101	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	100	0.00
22		Acetophenone	40.000	43.074	-7.7	101	0.00
23	S	Nitrobenzene-d5	80.000	80.978	-1.2	101	0.00
24		Nitrobenzene	40.000	41.799	-4.5	102	0.00
25		Isophorone	40.000	37.791	5.5	100	0.00
26	C	2-Nitrophenol	40.000	39.942	0.1	101	0.00
27		2,4-Dimethylphenol	40.000	41.383	-3.5	103	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.204	-0.5	100	0.00
29	C	2,4-Dichlorophenol	40.000	39.057	2.4	99	0.00
30		1,2,4-Trichlorobenzene	40.000	39.032	2.4	99	0.00
31		Naphthalene	40.000	38.931	2.7	103	0.00
32		Benzoic acid	40.000	36.983	7.5	93	0.00
33		4-Chloroaniline	40.000	42.399	-6.0	102	0.00
34	C	Hexachlorobutadiene	40.000	38.260	4.4	101	0.00
35		Caprolactam	40.000	39.872	0.3	98	0.00
36	C	4-Chloro-3-methylphenol	40.000	41.146	-2.9	100	0.00
37		2-Methylnaphthalene	40.000	38.282	4.3	100	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	103	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	38.826	2.9	100	0.00
40	P	Hexachlorocyclopentadiene	40.000	41.600	-4.0	101	0.00
41	S	2,4,6-Tribromophenol	80.000	79.680	0.4	97	0.00
42	C	2,4,6-Trichlorophenol	40.000	40.191	-0.5	100	0.00
43		2,4,5-Trichlorophenol	40.000	39.601	1.0	101	0.00
44	S	2-Fluorobiphenyl	80.000	74.685	6.6	101	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080416\
Data File : BF089607.D
Acq On : 5 Aug 2016 1:52
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 05 05:48:21 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF080416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 05 01:39:23 2016
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	38.533	3.7	100	0.00
46		2-Chloronaphthalene	40.000	38.974	2.6	102	0.00
47		2-Nitroaniline	40.000	38.992	2.5	100	0.00
48		Acenaphthylene	40.000	38.006	5.0	101	0.00
49		Dimethylphthalate	40.000	38.195	4.5	100	0.00
50		2,6-Dinitrotoluene	40.000	40.873	-2.2	98	0.00
51	C	Acenaphthene	40.000	38.220	4.5	100	0.00
52		3-Nitroaniline	40.000	37.787	5.5	95	0.00
53	P	2,4-Dinitrophenol	40.000	37.765	5.6	96	0.00
54		Dibenzofuran	40.000	38.572	3.6	101	0.00
55	P	4-Nitrophenol	40.000	32.343	19.1	81	0.01
56		2,4-Dinitrotoluene	40.000	38.812	3.0	98	0.00
57		Fluorene	40.000	37.537	6.2	99	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	38.852	2.9	100	0.00
59		Diethylphthalate	40.000	38.122	4.7	99	0.00
60		4-Chlorophenyl-phenylether	40.000	38.799	3.0	99	0.00
61		4-Nitroaniline	40.000	40.377	-0.9	98	0.00
62		Azobenzene	40.000	39.829	0.4	99	-0.01
63	I	Phenanthrene-d10	20.000	20.000	0.0	98	-0.01
64		4,6-Dinitro-2-methylphenol	40.000	38.323	4.2	96	0.00
65	c	n-Nitrosodiphenylamine	40.000	39.195	2.0	100	0.00
66		4-Bromophenyl-phenylether	40.000	40.725	-1.8	99	0.00
67		Hexachlorobenzene	40.000	37.139	7.2	98	0.00
68		Atrazine	40.000	43.370	-8.4	100	0.00
69	C	Pentachlorophenol	40.000	42.613	-6.5	94	0.00
70		Phenanthrene	40.000	38.350	4.1	98	0.00
71		Anthracene	40.000	37.399	6.5	98	0.00
72		Carbazole	40.000	38.917	2.7	96	-0.01
73		Di-n-butylphthalate	40.000	38.007	5.0	94	-0.01
74	C	Fluoranthene	40.000	38.570	3.6	96	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	94	0.00
76		Benzidine	40.000	44.133	-10.3	97	-0.01
77		Pyrene	40.000	36.918	7.7	93	-0.01
78	S	Terphenyl-d14	80.000	73.192	8.5	94	0.00
79		Butylbenzylphthalate	40.000	39.245	1.9	91	-0.01
80		Benzo(a)anthracene	40.000	38.666	3.3	95	0.00
81		3,3'-Dichlorobenzidine	40.000	40.316	-0.8	92	0.00
82		Chrysene	40.000	37.973	5.1	92	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	38.043	4.9	91	0.00
84	c	Di-n-octyl phthalate	40.000	37.832	5.4	86	-0.01
85		Indeno(1,2,3-cd)pyrene	40.000	37.991	5.0	99	0.00
86	I	Perylene-d12	20.000	20.000	0.0	93	0.00
87		Benzo(b)fluoranthene	40.000	40.110	-0.3	95	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080416\
Data File : BF089607.D
Acq On : 5 Aug 2016 1:52
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 05 05:48:21 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF080416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 05 01:39:23 2016
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	38.710	3.2	92	0.00
89 C	Benzo(a)pyrene	40.000	38.024	4.9	92	0.00
90	Dibenzo(a,h)anthracene	40.000	39.972	0.1	100	0.00
91	Benzo(g,h,i)perylene	40.000	40.513	-1.3	101	0.00

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