Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF112017\
Data File : BF100725.D
Acq On : 20 Nov 2017 12:50
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 20 14:27:41 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 17 15:22:08 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	45	0.00
2		1,4-Dioxane	40.000	37.759	5.6	42	-0.02
3		Pyridine	40.000	37.125	7.2	40	-0.03
4		n-Nitrosodimethylamine	40.000	34.405	14.0	37	-0.02
5	S	2-Fluorophenol	80.000	80.506	-0.6	46	-0.01
6		Aniline	40.000	37.216	7.0	41	-0.01
7	S	Phenol-d6	80.000	78.122	2.3	44	-0.01
8		2-Chlorophenol	40.000	40.969	-2.4	46	-0.01
9		Benzaldehyde	40.000	34.177	14.6	39	0.00
10	C	Phenol	40.000	38.304	4.2	43	-0.01
11		bis(2-Chloroethyl)ether	40.000	37.889	5.3	43	-0.01
12		1,3-Dichlorobenzene	40.000	41.150	-2.9	46	0.00
13	C	1,4-Dichlorobenzene	40.000	41.327	-3.3	47	0.00
14		1,2-Dichlorobenzene	40.000	41.734	-4.3	47	-0.01
15		Benzyl Alcohol	40.000	39.578	1.1	43	-0.01
16		2,2'-oxybis(1-Chloropropane	40.000	35.524	11.2	40	0.00
17		2-Methylphenol	40.000	39.704	0.7	45	0.00
18		Hexachloroethane	40.000	41.276	-3.2	45	-0.01
19	P	n-Nitroso-di-n-propylamine	40.000	36.864	7.8	42	-0.01
20		3+4-Methylphenols	40.000	39.648	0.9	44	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	44	-0.01
22		Acetophenone	40.000	39.153	2.1	44	0.00
23	S	Nitrobenzene-d5	80.000	89.625	-12.0	47	-0.01
24		Nitrobenzene	40.000	44.051	-10.1	45	-0.01
25		Isophorone	40.000	37.079	7.3	41	-0.01
26	C	2-Nitrophenol	40.000	49.679	-24.2#	56	-0.01
27		2,4-Dimethylphenol	40.000	40.067	-0.2	43	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.498	3.8	42	0.00
29	C	2,4-Dichlorophenol	40.000	42.615	-6.5	45	-0.01
30		1,2,4-Trichlorobenzene	40.000	43.042	-7.6	47	0.00
31		Naphthalene	40.000	40.973	-2.4	45	0.00
32		Benzoic acid	40.000	44.299	-10.7	53	0.00
33		4-Chloroaniline	40.000	40.944	-2.4	44	-0.01
34	C	Hexachlorobutadiene	40.000	42.598	-6.5	47	-0.01
35		Caprolactam	40.000	38.350	4.1	42	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.319	-0.8	43	-0.01
37		2-Methylnaphthalene	40.000	41.321	-3.3	46	-0.01
38	I	Acenaphthene-d10	20.000	20.000	0.0	42	-0.01
39		1,2,4,5-Tetrachlorobenzene	40.000	45.141	-12.9	48	-0.01
40	P	Hexachlorocyclopentadiene	40.000	44.124	-10.3	44	-0.01
41	S	2,4,6-Tribromophenol	80.000	92.999	-16.2	47	-0.01
42	C	2,4,6-Trichlorophenol	40.000	44.422	-11.1	45	0.00
43		2,4,5-Trichlorophenol	40.000	46.192	-15.5	47	-0.01
44	S	2-Fluorobiphenyl	80.000	88.785	-11.0	48	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF112017\
Data File : BF100725.D
Acq On : 20 Nov 2017 12:50
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 20 14:27:41 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 17 15:22:08 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	43.230	-8.1	47	-0.01
46		2-Chloronaphthalene	40.000	43.338	-8.3	46	-0.01
47		2-Nitroaniline	40.000	43.758	-9.4	46	-0.01
48		Acenaphthylene	40.000	42.157	-5.4	45	-0.01
49		Dimethylphthalate	40.000	41.953	-4.9	45	0.00
50		2,6-Dinitrotoluene	40.000	45.887	-14.7	46	-0.01
51	C	Acenaphthene	40.000	43.327	-8.3	47	-0.01
52		3-Nitroaniline	40.000	45.405	-13.5	46	-0.01
53	P	2,4-Dinitrophenol	40.000	59.388	-48.5#	79	0.00
54		Dibenzofuran	40.000	42.264	-5.7	45	-0.01
55	P	4-Nitrophenol	40.000	41.543	-3.9	42	0.00
56		2,4-Dinitrotoluene	40.000	47.892	-19.7	47	-0.01
57		Fluorene	40.000	44.888	-12.2	47	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	43.603	-9.0	44	0.00
59		Diethylphthalate	40.000	41.537	-3.8	45	-0.01
60		4-Chlorophenyl-phenylether	40.000	45.009	-12.5	48	0.00
61		4-Nitroaniline	40.000	46.827	-17.1	47	0.00
62		Azobenzene	40.000	37.816	5.5	41	-0.01
63	I	Phenanthrene-d10	20.000	20.000	0.0	44	0.00
64		4,6-Dinitro-2-methylphenol	40.000	52.607	-31.5#	64	0.00
65	c	n-Nitrosodiphenylamine	40.000	41.077	-2.7	45	-0.01
66		4-Bromophenyl-phenylether	40.000	42.361	-5.9	46	-0.01
67		Hexachlorobenzene	40.000	42.637	-6.6	47	0.00
68		Atrazine	40.000	41.191	-3.0	44	-0.01
69	C	Pentachlorophenol	40.000	41.455	-3.6	44	-0.01
70		Phenanthrene	40.000	41.846	-4.6	48	-0.01
71		Anthracene	40.000	41.326	-3.3	47	-0.01
72		Carbazole	40.000	40.720	-1.8	46	-0.01
73		Di-n-butylphthalate	40.000	43.175	-7.9	48	-0.01
74	C	Fluoranthene	40.000	41.412	-3.5	47	-0.01
75	I	Chrysene-d12	20.000	20.000	0.0	46	0.00
76		Benzidine	40.000	37.362	6.6	40	-0.01
77		Pyrene	40.000	40.882	-2.2	47	-0.01
78	S	Terphenyl-d14	80.000	82.516	-3.1	50	0.00
79		Butylbenzylphthalate	40.000	43.425	-8.6	48	0.00
80		Benzo(a)anthracene	40.000	41.993	-5.0	48	0.00
81		3,3'-Dichlorobenzidine	40.000	42.754	-6.9	46	-0.01
82		Chrysene	40.000	40.636	-1.6	47	-0.01
83		Bis(2-ethylhexyl)phthalate	40.000	45.773	-14.4	50	-0.01
84	c	Di-n-octyl phthalate	40.000	43.951	-9.9	48	-0.01
85		Indeno(1,2,3-cd)pyrene	40.000	37.525	6.2	44	-0.01
86	I	Perylene-d12	20.000	20.000	0.0	43	-0.01
87		Benzo(b)fluoranthene	40.000	43.431	-8.6	47	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF112017\
Data File : BF100725.D
Acq On : 20 Nov 2017 12:50
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 20 14:27:41 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 17 15:22:08 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	42.375	-5.9	43	0.00
89 C	Benzo(a)pyrene	40.000	42.462	-6.2	45	-0.01
90	Dibenzo(a,h)anthracene	40.000	40.820	-2.1	45	-0.02
91	Benzo(g,h,i)perylene	40.000	39.631	0.9	44	-0.02

(#) = Out of Range SPCC's out = 0 CCC's out = 1