Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF013017\
Data File : BF092641.D
Acq On : 30 Jan 2017 18:39
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 31 01:14:53 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF013017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 30 14:50:05 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	105	0.00
2		1,4-Dioxane	40.000	47.150	-17.9	127	0.01
3		Pyridine	40.000	51.607	-29.0#	134	0.00
4		n-Nitrosodimethylamine	40.000	50.639	-26.6#	133	0.01
5	S	2-Fluorophenol	80.000	83.901	-4.9	106	-0.01
6		Aniline	40.000	40.644	-1.6	110	0.00
7	S	Phenol-d6	80.000	80.861	-1.1	107	0.00
8		2-Chlorophenol	40.000	41.083	-2.7	108	0.00
9		Benzaldehyde	40.000	42.175	-5.4	109	0.00
10	C	Phenol	40.000	38.243	4.4	106	0.00
11		bis(2-Chloroethyl)ether	40.000	39.637	0.9	109	0.00
12		1,3-Dichlorobenzene	40.000	39.734	0.7	107	0.00
13	C	1,4-Dichlorobenzene	40.000	40.096	-0.2	109	0.00
14		1,2-Dichlorobenzene	40.000	42.153	-5.4	110	0.00
15		Benzyl Alcohol	40.000	39.727	0.7	108	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	39.621	0.9	106	0.00
17		2-Methylphenol	40.000	43.620	-9.0	110	0.00
18		Hexachloroethane	40.000	41.511	-3.8	109	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	37.953	5.1	110	0.00
20		3+4-Methylphenols	40.000	41.634	-4.1	107	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	110	0.00
22		Acetophenone	40.000	37.582	6.0	107	0.00
23	S	Nitrobenzene-d5	80.000	83.246	-4.1	114	0.00
24		Nitrobenzene	40.000	35.659	10.9	105	0.00
25		Isophorone	40.000	38.663	3.3	109	0.00
26	C	2-Nitrophenol	40.000	43.299	-8.2	110	0.00
27		2,4-Dimethylphenol	40.000	41.645	-4.1	109	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.431	-1.1	112	0.00
29	C	2,4-Dichlorophenol	40.000	38.778	3.1	112	0.01
30		1,2,4-Trichlorobenzene	40.000	38.403	4.0	108	0.00
31		Naphthalene	40.000	39.726	0.7	113	0.00
32		Benzoic acid	40.000	39.241	1.9	105	0.00
33		4-Chloroaniline	40.000	39.959	0.1	108	0.00
34	C	Hexachlorobutadiene	40.000	38.672	3.3	107	0.00
35		Caprolactam	40.000	41.220	-3.0	106	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.911	-2.3	111	0.00
37		2-Methylnaphthalene	40.000	36.665	8.3	101	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	109	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	41.476	-3.7	114	0.01
40	P	Hexachlorocyclopentadiene	40.000	41.572	-3.9	112	0.00
41	S	2,4,6-Tribromophenol	80.000	86.281	-7.9	108	0.00
42	C	2,4,6-Trichlorophenol	40.000	41.564	-3.9	107	0.00
43		2,4,5-Trichlorophenol	40.000	39.588	1.0	112	0.00
44	S	2-Fluorobiphenyl	80.000	74.943	6.3	96	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF013017\
Data File : BF092641.D
Acq On : 30 Jan 2017 18:39
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 31 01:14:53 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF013017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 30 14:50:05 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	41.895	-4.7	110	0.00
46		2-Chloronaphthalene	40.000	39.527	1.2	101	0.00
47		2-Nitroaniline	40.000	41.736	-4.3	121	0.01
48		Acenaphthylene	40.000	39.301	1.7	115	0.01
49		Dimethylphthalate	40.000	37.247	6.9	101	0.00
50		2,6-Dinitrotoluene	40.000	46.312	-15.8	123	0.01
51	C	Acenaphthene	40.000	39.845	0.4	115	0.00
52		3-Nitroaniline	40.000	46.995	-17.5	120	0.00
53	P	2,4-Dinitrophenol	40.000	42.885	-7.2	119	0.01
54		Dibenzofuran	40.000	39.169	2.1	114	0.00
55	P	4-Nitrophenol	40.000	39.542	1.1	109	0.00
56		2,4-Dinitrotoluene	40.000	43.860	-9.6	106	0.00
57		Fluorene	40.000	38.831	2.9	109	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	43.874	-9.7	107	0.00
59		Diethylphthalate	40.000	38.728	3.2	102	0.00
60		4-Chlorophenyl-phenylether	40.000	40.259	-0.6	113	0.01
61		4-Nitroaniline	40.000	42.251	-5.6	116	0.00
62		Azobenzene	40.000	40.539	-1.3	111	0.01
63	I	Phenanthrene-d10	20.000	20.000	0.0	107	0.00
64		4,6-Dinitro-2-methylphenol	40.000	43.854	-9.6	118	0.00
65	c	n-Nitrosodiphenylamine	40.000	38.642	3.4	102	0.00
66		4-Bromophenyl-phenylether	40.000	40.305	-0.8	111	0.01
67		Hexachlorobenzene	40.000	38.590	3.5	106	0.00
68		Atrazine	40.000	37.237	6.9	106	0.00
69	C	Pentachlorophenol	40.000	41.680	-4.2	114	0.00
70		Phenanthrene	40.000	40.272	-0.7	107	0.00
71		Anthracene	40.000	37.716	5.7	104	0.00
72		Carbazole	40.000	41.702	-4.3	105	0.00
73		Di-n-butylphthalate	40.000	39.048	2.4	105	0.00
74	C	Fluoranthene	40.000	39.241	1.9	103	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	105	-0.01
76		Benzidine	40.000	38.717	3.2	102	0.00
77		Pyrene	40.000	41.219	-3.0	102	0.00
78	S	Terphenyl-d14	80.000	78.730	1.6	108	0.00
79		Butylbenzylphthalate	40.000	40.493	-1.2	106	0.00
80		Benzo(a)anthracene	40.000	39.853	0.4	103	0.00
81		3,3'-Dichlorobenzidine	40.000	40.046	-0.1	102	0.00
82		Chrysene	40.000	44.027	-10.1	107	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	43.256	-8.1	109	0.00
84	c	Di-n-octyl phthalate	40.000	44.570	-11.4	111	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	38.207	4.5	105	0.00
86	I	Perylene-d12	20.000	20.000	0.0	105	0.00
87		Benzo(b)fluoranthene	40.000	44.565	-11.4	111	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF013017\
Data File : BF092641.D
Acq On : 30 Jan 2017 18:39
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 31 01:14:53 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF013017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 30 14:50:05 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	35.215	12.0	101	0.00
89 C	Benzo(a)pyrene	40.000	40.159	-0.4	106	0.00
90	Dibenzo(a,h)anthracene	40.000	37.631	5.9	105	0.00
91	Benzo(g,h,i)perylene	40.000	37.133	7.2	104	-0.01

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