Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080216\
Data File : BF089540.D
Acq On : 2 Aug 2016 11:46
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 03 11:08:02 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 28 16:02:32 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	111	-0.08
2		1,4-Dioxane	40.000	40.189	-0.5	115	-0.11
3		Pyridine	40.000	51.531	-28.8#	133	-0.14
4		n-Nitrosodimethylamine	40.000	44.901	-12.3	129	-0.14
5	S	2-Fluorophenol	80.000	84.328	-5.4	113	-0.08
6		Aniline	40.000	51.068	-27.7#	132	-0.07
7	S	Phenol-d6	80.000	80.688	-0.9	107	-0.07
8		2-Chlorophenol	40.000	42.046	-5.1	107	-0.07
9		Benzaldehyde	40.000	33.430	16.4	90	-0.08
10	C	Phenol	40.000	42.880	-7.2	109	-0.06
11		bis(2-Chloroethyl)ether	40.000	40.117	-0.3	111	-0.07
12		1,3-Dichlorobenzene	40.000	42.814	-7.0	112	-0.07
13	C	1,4-Dichlorobenzene	40.000	42.094	-5.2	110	-0.07
14		1,2-Dichlorobenzene	40.000	38.894	2.8	111	-0.08
15		Benzyl Alcohol	40.000	47.828	-19.6	128	-0.07
16		2,2'-oxybis(1-Chloropropane	40.000	41.851	-4.6	110	-0.07
17		2-Methylphenol	40.000	39.507	1.2	108	-0.06
18		Hexachloroethane	40.000	42.378	-5.9	115	-0.07
19	P	n-Nitroso-di-n-propylamine	40.000	45.641	-14.1	118	-0.07
20		3+4-Methylphenols	40.000	41.730	-4.3	109	-0.06
21	I	Naphthalene-d8	20.000	20.000	0.0	111	-0.07
22		Acetophenone	40.000	43.215	-8.0	111	-0.07
23	S	Nitrobenzene-d5	80.000	85.440	-6.8	113	-0.07
24		Nitrobenzene	40.000	40.871	-2.2	112	-0.07
25		Isophorone	40.000	38.816	3.0	104	-0.07
26	C	2-Nitrophenol	40.000	39.281	1.8	108	-0.07
27		2,4-Dimethylphenol	40.000	43.829	-9.6	118	-0.06
28		bis(2-Chloroethoxy)methane	40.000	44.338	-10.8	116	-0.06
29	C	2,4-Dichlorophenol	40.000	41.590	-4.0	107	-0.06
30		1,2,4-Trichlorobenzene	40.000	41.360	-3.4	111	-0.07
31		Naphthalene	40.000	39.685	0.8	112	-0.07
32		Benzoic acid	40.000	42.984	-7.5	116	-0.05
33		4-Chloroaniline	40.000	46.111	-15.3	131	-0.07
34	C	Hexachlorobutadiene	40.000	42.130	-5.3	110	-0.07
35		Caprolactam	40.000	44.672	-11.7	120	-0.06
36	C	4-Chloro-3-methylphenol	40.000	43.780	-9.5	110	-0.06
37		2-Methylnaphthalene	40.000	41.151	-2.9	112	-0.07
38	I	Acenaphthene-d10	20.000	20.000	0.0	107	-0.07
39		1,2,4,5-Tetrachlorobenzene	40.000	37.792	5.5	107	-0.07
40	P	Hexachlorocyclopentadiene	40.000	36.714	8.2	99	-0.07
41	S	2,4,6-Tribromophenol	80.000	81.554	-1.9	105	-0.07
42	C	2,4,6-Trichlorophenol	40.000	44.824	-12.1	115	-0.07
43		2,4,5-Trichlorophenol	40.000	36.213	9.5	98	-0.07
44	S	2-Fluorobiphenyl	80.000	86.725	-8.4	112	-0.07

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080216\
Data File : BF089540.D
Acq On : 2 Aug 2016 11:46
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 03 11:08:02 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 28 16:02:32 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	37.563	6.1	108	-0.07
46		2-Chloronaphthalene	40.000	41.900	-4.7	115	-0.07
47		2-Nitroaniline	40.000	46.474	-16.2	124	-0.07
48		Acenaphthylene	40.000	42.759	-6.9	110	-0.07
49		Dimethylphthalate	40.000	40.161	-0.4	106	-0.07
50		2,6-Dinitrotoluene	40.000	39.851	0.4	99	-0.07
51	C	Acenaphthene	40.000	42.890	-7.2	112	-0.07
52		3-Nitroaniline	40.000	49.887	-24.7	132	-0.07
53	P	2,4-Dinitrophenol	40.000	45.528	-13.8	139	-0.07
54		Dibenzofuran	40.000	41.612	-4.0	112	-0.07
55	P	4-Nitrophenol	40.000	36.341	9.1	101	-0.06
56		2,4-Dinitrotoluene	40.000	40.378	-0.9	104	-0.06
57		Fluorene	40.000	38.438	3.9	101	-0.08
58		2,3,4,6-Tetrachlorophenol	40.000	43.773	-9.4	106	-0.07
59		Diethylphthalate	40.000	36.556	8.6	104	-0.07
60		4-Chlorophenyl-phenylether	40.000	37.566	6.1	105	-0.08
61		4-Nitroaniline	40.000	45.050	-12.6	117	-0.06
62		Azobenzene	40.000	40.346	-0.9	115	-0.07
63	I	Phenanthrene-d10	20.000	20.000	0.0	112	-0.08
64		4,6-Dinitro-2-methylphenol	40.000	39.827	0.4	106	-0.07
65	c	n-Nitrosodiphenylamine	40.000	40.004	-0.0	113	-0.07
66		4-Bromophenyl-phenylether	40.000	38.150	4.6	99	-0.07
67		Hexachlorobenzene	40.000	39.329	1.7	113	-0.07
68		Atrazine	40.000	39.548	1.1	101	-0.07
69	C	Pentachlorophenol	40.000	38.415	4.0	113	-0.07
70		Phenanthrene	40.000	39.501	1.2	105	-0.07
71		Anthracene	40.000	39.594	1.0	115	-0.07
72		Carbazole	40.000	40.019	-0.0	112	-0.08
73		Di-n-butylphthalate	40.000	34.957	12.6	101	-0.07
74	C	Fluoranthene	40.000	39.816	0.5	104	-0.08
75	I	Chrysene-d12	20.000	20.000	0.0	107	-0.07
76		Benzidine	40.000	38.788	3.0	103	-0.08
77		Pyrene	40.000	41.749	-4.4	114	-0.07
78	S	Terphenyl-d14	80.000	89.231	-11.5	124	-0.07
79		Butylbenzylphthalate	40.000	45.379	-13.4	111	-0.07
80		Benzo(a)anthracene	40.000	42.460	-6.2	113	-0.07
81		3,3'-Dichlorobenzidine	40.000	40.071	-0.2	106	-0.08
82		Chrysene	40.000	42.768	-6.9	120	-0.07
83		Bis(2-ethylhexyl)phthalate	40.000	39.150	2.1	105	-0.07
84	c	Di-n-octyl phthalate	40.000	40.680	-1.7	110	-0.07
85		Indeno(1,2,3-cd)pyrene	40.000	38.416	4.0	108	-0.13
86	I	Perylene-d12	20.000	20.000	0.0	111	-0.08
87		Benzo(b)fluoranthene	40.000	45.560	-13.9	138	-0.07

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080216\
Data File : BF089540.D
Acq On : 2 Aug 2016 11:46
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 03 11:08:02 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 28 16:02:32 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	38.485	3.8	97	-0.07
89 C	Benzo(a)pyrene	40.000	39.858	0.4	109	-0.08
90	Dibenzo(a,h)anthracene	40.000	40.794	-2.0	112	-0.11
91	Benzo(g,h,i)perylene	40.000	42.705	-6.8	115	-0.13

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