Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF062817\
Data File : BF096243.D
Acq On : 28 Jun 2017 12:15
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 29 01:28:00 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF062717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 28 15:41:49 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	100	0.00
2		1,4-Dioxane	40.000	40.838	-2.1	103	0.00
3		Pyridine	40.000	41.040	-2.6	101	-0.02
4		n-Nitrosodimethylamine	40.000	40.514	-1.3	99	0.02
5	S	2-Fluorophenol	80.000	80.801	-1.0	102	0.02
6		Aniline	40.000	40.602	-1.5	101	0.01
7	S	Phenol-d6	80.000	82.107	-2.6	105	0.02
8		2-Chlorophenol	40.000	40.632	-1.6	101	0.01
9		Benzaldehyde	40.000	41.755	-4.4	101	0.00
10	C	Phenol	40.000	41.419	-3.5	102	0.02
11		bis(2-Chloroethyl)ether	40.000	40.290	-0.7	101	0.02
12		1,3-Dichlorobenzene	40.000	39.710	0.7	101	0.01
13	C	1,4-Dichlorobenzene	40.000	40.059	-0.1	101	0.01
14		1,2-Dichlorobenzene	40.000	39.697	0.8	100	0.01
15		Benzyl Alcohol	40.000	42.728	-6.8	105	0.01
16		2,2'-oxybis(1-Chloropropane	40.000	39.310	1.7	98	0.00
17		2-Methylphenol	40.000	41.059	-2.6	103	0.02
18		Hexachloroethane	40.000	40.775	-1.9	100	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.761	0.6	99	0.02
20		3+4-Methylphenols	40.000	40.509	-1.3	103	0.02
21	I	Naphthalene-d8	20.000	20.000	0.0	100	0.01
22		Acetophenone	40.000	39.462	1.3	100	0.00
23	S	Nitrobenzene-d5	80.000	88.752	-10.9	107	-0.01
24		Nitrobenzene	40.000	44.614	-11.5	104	0.02
25		Isophorone	40.000	38.967	2.6	99	0.02
26	C	2-Nitrophenol	40.000	47.168	-17.9	122	0.00
27		2,4-Dimethylphenol	40.000	39.818	0.5	101	0.01
28		bis(2-Chloroethoxy)methane	40.000	38.859	2.9	100	0.01
29	C	2,4-Dichlorophenol	40.000	41.438	-3.6	101	0.01
30		1,2,4-Trichlorobenzene	40.000	39.117	2.2	99	0.01
31		Naphthalene	40.000	39.177	2.1	101	0.01
32		Benzoic acid	40.000	39.247	1.9	101	-0.05
33		4-Chloroaniline	40.000	39.624	0.9	100	0.01
34	C	Hexachlorobutadiene	40.000	39.327	1.7	99	0.00
35		Caprolactam	40.000	41.781	-4.5	101	0.07
36	C	4-Chloro-3-methylphenol	40.000	40.154	-0.4	99	0.02
37		2-Methylnaphthalene	40.000	38.863	2.8	100	0.01
38	I	Acenaphthene-d10	20.000	20.000	0.0	99	0.01
39		1,2,4,5-Tetrachlorobenzene	40.000	39.407	1.5	102	0.01
40	P	Hexachlorocyclopentadiene	40.000	48.001	-20.0	112	0.00
41	S	2,4,6-Tribromophenol	80.000	104.585	-30.7#	128	0.01
42	C	2,4,6-Trichlorophenol	40.000	40.559	-1.4	100	0.01
43		2,4,5-Trichlorophenol	40.000	42.754	-6.9	101	0.02
44	S	2-Fluorobiphenyl	80.000	84.224	-5.3	99	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF062817\
Data File : BF096243.D
Acq On : 28 Jun 2017 12:15
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 29 01:28:00 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF062717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 28 15:41:49 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	40.226	-0.6	100	0.00
46		2-Chloronaphthalene	40.000	38.324	4.2	99	0.01
47		2-Nitroaniline	40.000	46.184	-15.5	111	0.00
48		Acenaphthylene	40.000	38.856	2.9	100	0.01
49		Dimethylphthalate	40.000	38.734	3.2	99	0.02
50		2,6-Dinitrotoluene	40.000	43.480	-8.7	106	-0.01
51	C	Acenaphthene	40.000	38.870	2.8	101	0.02
52		3-Nitroaniline	40.000	43.486	-8.7	107	0.00
53	P	2,4-Dinitrophenol	40.000	52.266	-30.7#	149	0.00
54		Dibenzofuran	40.000	39.424	1.4	98	0.00
55	P	4-Nitrophenol	40.000	43.691	-9.2	104	0.00
56		2,4-Dinitrotoluene	40.000	44.553	-11.4	105	0.00
57		Fluorene	40.000	42.719	-6.8	101	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	42.531	-6.3	102	0.01
59		Diethylphthalate	40.000	37.972	5.1	95	0.02
60		4-Chlorophenyl-phenylether	40.000	41.520	-3.8	101	0.00
61		4-Nitroaniline	40.000	43.658	-9.1	108	-0.02
62		Azobenzene	40.000	38.892	2.8	99	0.02
63	I	Phenanthrene-d10	20.000	20.000	0.0	101	0.01
64		4,6-Dinitro-2-methylphenol	40.000	47.660	-19.1	132	0.00
65	c	n-Nitrosodiphenylamine	40.000	38.169	4.6	99	0.01
66		4-Bromophenyl-phenylether	40.000	39.246	1.9	100	0.00
67		Hexachlorobenzene	40.000	42.310	-5.8	107	0.01
68		Atrazine	40.000	40.041	-0.1	102	0.02
69	C	Pentachlorophenol	40.000	46.573	-16.4	108	0.01
70		Phenanthrene	40.000	41.819	-4.5	100	0.00
71		Anthracene	40.000	39.931	0.2	101	0.00
72		Carbazole	40.000	37.854	5.4	99	0.02
73		Di-n-butylphthalate	40.000	39.630	0.9	99	0.00
74	C	Fluoranthene	40.000	38.216	4.5	100	0.02
75	I	Chrysene-d12	20.000	20.000	0.0	109	0.02
76		Benzidine	40.000	38.688	3.3	99	0.00
77		Pyrene	40.000	36.097	9.8	98	0.02
78	S	Terphenyl-d14	80.000	76.421	4.5	107	0.01
79		Butylbenzylphthalate	40.000	39.626	0.9	102	0.00
80		Benzo(a)anthracene	40.000	35.113	12.2	94	0.02
81		3,3'-Dichlorobenzidine	40.000	39.916	0.2	100	0.02
82		Chrysene	40.000	35.743	10.6	96	0.02
83		Bis(2-ethylhexyl)phthalate	40.000	39.246	1.9	98	0.00
84	c	Di-n-octyl phthalate	40.000	38.104	4.7	98	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	38.655	3.4	105	0.04
86	I	Perylene-d12	20.000	20.000	0.0	100	0.02
87		Benzo(b)fluoranthene	40.000	36.372	9.1	96	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF062817\
Data File : BF096243.D
Acq On : 28 Jun 2017 12:15
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 29 01:28:00 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF062717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 28 15:41:49 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	36.145	9.6	89	0.02
89 C	Benzo(a)pyrene	40.000	36.621	8.4	92	0.02
90	Dibenzo(a,h)anthracene	40.000	40.842	-2.1	105	0.04
91	Benzo(g,h,i)perylene	40.000	40.634	-1.6	107	0.05

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