Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051117\
Data File : BF095078.D
Acq On : 11 May 2017 14:04
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 11 16:30:01 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 03 17:24:51 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	102	-0.03
2		1,4-Dioxane	40.000	39.863	0.3	107	-0.04
3		Pyridine	40.000	37.331	6.7	98	-0.04
4		n-Nitrosodimethylamine	40.000	37.320	6.7	99	-0.04
5	S	2-Fluorophenol	80.000	82.595	-3.2	106	-0.03
6		Aniline	40.000	41.210	-3.0	101	-0.02
7	S	Phenol-d6	80.000	81.559	-1.9	104	-0.02
8		2-Chlorophenol	40.000	41.128	-2.8	106	-0.02
9		Benzaldehyde	40.000	35.602	11.0	89	-0.02
10	C	Phenol	40.000	37.299	6.8	96	-0.02
11		bis(2-Chloroethyl)ether	40.000	40.078	-0.2	102	-0.02
12		1,3-Dichlorobenzene	40.000	41.039	-2.6	113	-0.03
13	C	1,4-Dichlorobenzene	40.000	40.317	-0.8	103	-0.03
14		1,2-Dichlorobenzene	40.000	41.631	-4.1	107	-0.03
15		Benzyl Alcohol	40.000	44.701	-11.8	109	-0.02
16		2,2'-oxybis(1-Chloropropane	40.000	40.446	-1.1	106	-0.01
17		2-Methylphenol	40.000	41.891	-4.7	112	-0.02
18		Hexachloroethane	40.000	41.418	-3.5	105	-0.04
19	P	n-Nitroso-di-n-propylamine	40.000	42.330	-5.8	111	-0.02
20		3+4-Methylphenols	40.000	41.651	-4.1	107	-0.02
21	I	Naphthalene-d8	20.000	20.000	0.0	102	-0.03
22		Acetophenone	40.000	41.481	-3.7	108	-0.02
23	S	Nitrobenzene-d5	80.000	83.753	-4.7	107	-0.02
24		Nitrobenzene	40.000	42.039	-5.1	111	-0.02
25		Isophorone	40.000	42.796	-7.0	110	-0.02
26	C	2-Nitrophenol	40.000	43.548	-8.9	110	-0.02
27		2,4-Dimethylphenol	40.000	39.541	1.1	105	-0.02
28		bis(2-Chloroethoxy)methane	40.000	40.794	-2.0	106	-0.02
29	C	2,4-Dichlorophenol	40.000	37.573	6.1	100	-0.01
30		1,2,4-Trichlorobenzene	40.000	39.672	0.8	105	-0.02
31		Naphthalene	40.000	38.978	2.6	103	-0.03
32		Benzoic acid	40.000	36.599	8.5	101	0.02
33		4-Chloroaniline	40.000	42.840	-7.1	111	-0.02
34	C	Hexachlorobutadiene	40.000	39.372	1.6	104	-0.04
35		Caprolactam	40.000	41.673	-4.2	104	0.00
36	C	4-Chloro-3-methylphenol	40.000	42.182	-5.5	109	-0.01
37		2-Methylnaphthalene	40.000	40.116	-0.3	106	-0.03
38	I	Acenaphthene-d10	20.000	20.000	0.0	109	-0.03
39		1,2,4,5-Tetrachlorobenzene	40.000	37.734	5.7	98	-0.02
40	P	Hexachlorocyclopentadiene	40.000	38.982	2.5	104	-0.03
41	S	2,4,6-Tribromophenol	80.000	76.538	4.3	98	-0.03
42	C	2,4,6-Trichlorophenol	40.000	40.302	-0.8	105	-0.02
43		2,4,5-Trichlorophenol	40.000	36.500	8.8	94	0.00
44	S	2-Fluorobiphenyl	80.000	75.776	5.3	101	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051117\
Data File : BF095078.D
Acq On : 11 May 2017 14:04
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 11 16:30:01 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 03 17:24:51 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	40.520	-1.3	105	-0.03
46		2-Chloronaphthalene	40.000	39.779	0.6	106	-0.02
47		2-Nitroaniline	40.000	40.135	-0.3	107	-0.02
48		Acenaphthylene	40.000	36.664	8.3	98	-0.03
49		Dimethylphthalate	40.000	35.472	11.3	94	-0.03
50		2,6-Dinitrotoluene	40.000	37.899	5.3	99	-0.02
51	C	Acenaphthene	40.000	38.758	3.1	103	-0.04
52		3-Nitroaniline	40.000	41.703	-4.3	108	0.00
53	P	2,4-Dinitrophenol	40.000	37.197	7.0	101	0.00
54		Dibenzofuran	40.000	39.776	0.6	108	-0.03
55	P	4-Nitrophenol	40.000	34.795	13.0	87	0.02
56		2,4-Dinitrotoluene	40.000	40.995	-2.5	106	-0.01
57		Fluorene	40.000	38.235	4.4	105	-0.03
58		2,3,4,6-Tetrachlorophenol	40.000	40.628	-1.6	109	-0.02
59		Diethylphthalate	40.000	39.431	1.4	108	-0.03
60		4-Chlorophenyl-phenylether	40.000	39.532	1.2	105	-0.03
61		4-Nitroaniline	40.000	39.862	0.3	107	-0.01
62		Azobenzene	40.000	40.885	-2.2	106	-0.03
63	I	Phenanthrene-d10	20.000	20.000	0.0	103	-0.02
64		4,6-Dinitro-2-methylphenol	40.000	42.710	-6.8	108	-0.01
65	c	n-Nitrosodiphenylamine	40.000	40.842	-2.1	106	-0.02
66		4-Bromophenyl-phenylether	40.000	39.911	0.2	101	-0.04
67		Hexachlorobenzene	40.000	40.161	-0.4	104	-0.02
68		Atrazine	40.000	41.017	-2.5	111	-0.02
69	C	Pentachlorophenol	40.000	37.667	5.8	109	-0.02
70		Phenanthrene	40.000	39.848	0.4	105	-0.02
71		Anthracene	40.000	40.798	-2.0	107	-0.02
72		Carbazole	40.000	41.492	-3.7	110	-0.02
73		Di-n-butylphthalate	40.000	43.219	-8.0	110	-0.03
74	C	Fluoranthene	40.000	39.013	2.5	100	-0.02
75	I	Chrysene-d12	20.000	20.000	0.0	104	-0.02
76		Benzidine	40.000	38.270	4.3	97	-0.02
77		Pyrene	40.000	41.355	-3.4	106	-0.02
78	S	Terphenyl-d14	80.000	81.378	-1.7	106	-0.02
79		Butylbenzylphthalate	40.000	43.319	-8.3	110	-0.02
80		Benzo(a)anthracene	40.000	39.032	2.4	101	-0.02
81		3,3'-Dichlorobenzidine	40.000	38.571	3.6	96	-0.02
82		Chrysene	40.000	40.042	-0.1	103	-0.02
83		Bis(2-ethylhexyl)phthalate	40.000	42.222	-5.6	107	-0.03
84	c	Di-n-octyl phthalate	40.000	40.958	-2.4	103	-0.03
85		Indeno(1,2,3-cd)pyrene	40.000	36.972	7.6	98	-0.02
86	I	Perylene-d12	20.000	20.000	0.0	98	-0.01
87		Benzo(b)fluoranthene	40.000	42.463	-6.2	96	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051117\
Data File : BF095078.D
Acq On : 11 May 2017 14:04
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 11 16:30:01 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 03 17:24:51 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	40.927	-2.3	98	-0.02
89 C	Benzo(a)pyrene	40.000	39.881	0.3	94	-0.02
90	Dibenzo(a,h)anthracene	40.000	40.718	-1.8	98	-0.02
91	Benzo(g,h,i)perylene	40.000	40.750	-1.9	101	-0.02

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