Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF110918\
Data File : BF110615.D
Acq On : 9 Nov 2018 9:29
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 09 15:48:17 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 09 15:13:13 2018
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	40.256	-0.6	104	0.00
3		Pyridine	40.000	42.438	-6.1	103	0.00
4		n-Nitrosodimethylamine	40.000	41.539	-3.8	103	0.00
5	S	2-Fluorophenol	80.000	81.602	-2.0	103	0.00
6		Aniline	40.000	39.595	1.0	104	0.00
7	S	Phenol-d6	80.000	80.094	-0.1	104	0.00
8		2-Chlorophenol	40.000	39.809	0.5	103	0.00
9		Benzaldehyde	40.000	39.750	0.6	101	0.00
10	C	Phenol	40.000	39.746	0.6	102	0.00
11		bis(2-Chloroethyl)ether	40.000	39.444	1.4	103	0.00
12		1,3-Dichlorobenzene	40.000	39.717	0.7	102	0.00
13	C	1,4-Dichlorobenzene	40.000	39.095	2.3	102	0.00
14		1,2-Dichlorobenzene	40.000	39.692	0.8	104	0.00
15		Benzyl Alcohol	40.000	40.192	-0.5	103	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	39.490	1.3	102	0.00
17		2-Methylphenol	40.000	39.485	1.3	103	0.00
18		Hexachloroethane	40.000	39.116	2.2	101	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.579	1.1	104	0.00
20		3+4-Methylphenols	40.000	39.685	0.8	104	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	102	0.00
22		Acetophenone	40.000	40.116	-0.3	102	0.00
23	S	Nitrobenzene-d5	80.000	81.398	-1.7	103	0.00
24		Nitrobenzene	40.000	40.218	-0.5	102	0.00
25		Isophorone	40.000	39.946	0.1	103	0.00
26	C	2-Nitrophenol	40.000	41.764	-4.4	103	0.00
27		2,4-Dimethylphenol	40.000	40.610	-1.5	103	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.687	0.8	101	0.00
29	C	2,4-Dichlorophenol	40.000	40.948	-2.4	102	0.00
30		1,2,4-Trichlorobenzene	40.000	40.188	-0.5	101	0.00
31		Naphthalene	40.000	39.693	0.8	102	0.00
32		Benzoic acid	40.000	43.596	-9.0	103	0.00
33		4-Chloroaniline	40.000	38.988	2.5	103	0.00
34	C	Hexachlorobutadiene	40.000	39.348	1.6	101	0.00
35		Caprolactam	40.000	40.800	-2.0	103	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.542	-1.4	102	0.00
37		2-Methylnaphthalene	40.000	39.809	0.5	102	0.00
38		1-Methylnaphthalene	40.000	39.919	0.2	102	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	102	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	40.074	-0.2	101	0.00
41	P	Hexachlorocyclopentadiene	40.000	36.220	9.5	92	0.00
42	S	2,4,6-Tribromophenol	80.000	77.602	3.0	98	0.00
43	C	2,4,6-Trichlorophenol	40.000	40.019	-0.0	99	0.00
44		2,4,5-Trichlorophenol	40.000	40.034	-0.1	100	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF110918\
Data File : BF110615.D
Acq On : 9 Nov 2018 9:29
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 09 15:48:17 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 09 15:13:13 2018
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	S	2-Fluorobiphenyl	80.000	77.695	2.9	100	0.00
46		1,1'-Biphenyl	40.000	39.470	1.3	101	0.00
47		2-Chloronaphthalene	40.000	39.636	0.9	102	0.00
48		2-Nitroaniline	40.000	40.038	-0.1	100	0.00
49		Acenaphthylene	40.000	39.401	1.5	100	0.00
50		Dimethylphthalate	40.000	38.258	4.4	98	0.00
51		2,6-Dinitrotoluene	40.000	39.196	2.0	99	0.00
52	C	Acenaphthene	40.000	38.901	2.7	100	0.00
53		3-Nitroaniline	40.000	40.204	-0.5	102	0.00
54	P	2,4-Dinitrophenol	40.000	37.214	7.0	93	0.00
55		Dibenzofuran	40.000	38.809	3.0	99	0.00
56	P	4-Nitrophenol	40.000	39.999	0.0	97	0.00
57		2,4-Dinitrotoluene	40.000	40.002	-0.0	100	0.00
58		Fluorene	40.000	38.659	3.4	100	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	40.050	-0.1	101	0.00
60		Diethylphthalate	40.000	38.170	4.6	99	0.00
61		4-Chlorophenyl-phenylether	40.000	38.859	2.9	99	0.00
62		4-Nitroaniline	40.000	39.307	1.7	99	0.00
63		Azobenzene	40.000	38.998	2.5	99	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	98	0.00
65		4,6-Dinitro-2-methylphenol	40.000	38.595	3.5	92	0.00
66	c	n-Nitrosodiphenylamine	40.000	40.256	-0.6	100	0.00
67		4-Bromophenyl-phenylether	40.000	39.319	1.7	98	0.00
68		Hexachlorobenzene	40.000	40.072	-0.2	100	0.00
69		Atrazine	40.000	37.484	6.3	93	0.00
70	C	Pentachlorophenol	40.000	40.901	-2.3	97	0.00
71		Phenanthrene	40.000	39.533	1.2	100	0.00
72		Anthracene	40.000	39.567	1.1	99	0.00
73		Carbazole	40.000	40.182	-0.5	100	0.00
74		Di-n-butylphthalate	40.000	38.990	2.5	96	0.00
75	C	Fluoranthene	40.000	39.685	0.8	100	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	102	0.00
77		Benzidine	40.000	37.034	7.4	98	0.00
78		Pyrene	40.000	38.820	2.9	100	0.00
79	S	Terphenyl-d14	80.000	75.807	5.2	100	0.00
80		Butylbenzylphthalate	40.000	39.185	2.0	98	0.00
81		Benzo(a)anthracene	40.000	39.664	0.8	102	0.00
82		3,3'-Dichlorobenzidine	40.000	39.133	2.2	102	0.00
83		Chrysene	40.000	39.562	1.1	105	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	39.280	1.8	101	0.00
85	c	Di-n-octyl phthalate	40.000	40.335	-0.8	107	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	36.412	9.0	96	0.00
87	I	Perylene-d12	20.000	20.000	0.0	98	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF110918\
Data File : BF110615.D
Acq On : 9 Nov 2018 9:29
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 09 15:48:17 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 09 15:13:13 2018
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(b)fluoranthene	40.000	41.151	-2.9	98	0.00
89	Benzo(k)fluoranthene	40.000	42.507	-6.3	108	0.00
90 C	Benzo(a)pyrene	40.000	40.289	-0.7	99	0.00
91	Dibenzo(a,h)anthracene	40.000	39.737	0.7	95	-0.01
92	Benzo(g,h,i)perylene	40.000	40.054	-0.1	95	0.00

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