Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102318\
Data File : BF110080.D
Acq On : 23 Oct 2018 16:30
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 24 09:06:21 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 23 15:06:18 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	125	-0.01
2		1,4-Dioxane	40.000	38.891	2.8	123	-0.04
3		Pyridine	40.000	42.268	-5.7	127	-0.03
4		n-Nitrosodimethylamine	40.000	42.042	-5.1	129	-0.02
5	S	2-Fluorophenol	80.000	78.200	2.2	122	-0.02
6		Aniline	40.000	41.800	-4.5	130	-0.01
7	S	Phenol-d6	80.000	79.878	0.2	126	-0.01
8		2-Chlorophenol	40.000	40.247	-0.6	125	-0.02
9		Benzaldehyde	40.000	33.764	15.6	107	-0.01
10	C	Phenol	40.000	44.154	-10.4	138	-0.01
11		bis(2-Chloroethyl)ether	40.000	40.444	-1.1	128	-0.01
12		1,3-Dichlorobenzene	40.000	38.649	3.4	122	-0.01
13	C	1,4-Dichlorobenzene	40.000	38.844	2.9	123	-0.01
14		1,2-Dichlorobenzene	40.000	38.828	2.9	122	-0.02
15		Benzyl Alcohol	40.000	42.015	-5.0	129	-0.01
16		2,2'-oxybis(1-Chloropropane	40.000	40.886	-2.2	128	-0.01
17		2-Methylphenol	40.000	41.048	-2.6	128	-0.01
18		Hexachloroethane	40.000	40.226	-0.6	127	-0.02
19	P	n-Nitroso-di-n-propylamine	40.000	40.749	-1.9	129	0.00
20		3+4-Methylphenols	40.000	39.636	0.9	124	-0.01
21	I	Naphthalene-d8	20.000	20.000	0.0	130	-0.02
22		Acetophenone	40.000	38.376	4.1	127	-0.01
23	S	Nitrobenzene-d5	80.000	79.594	0.5	128	-0.01
24		Nitrobenzene	40.000	40.099	-0.2	129	-0.01
25		Isophorone	40.000	40.592	-1.5	132	-0.01
26	C	2-Nitrophenol	40.000	43.000	-7.5	133	-0.01
27		2,4-Dimethylphenol	40.000	39.499	1.3	128	-0.01
28		bis(2-Chloroethoxy)methane	40.000	39.244	1.9	129	-0.01
29	C	2,4-Dichlorophenol	40.000	39.914	0.2	128	-0.01
30		1,2,4-Trichlorobenzene	40.000	39.048	2.4	127	-0.01
31		Naphthalene	40.000	38.184	4.5	126	-0.02
32		Benzoic acid	40.000	43.407	-8.5	137	0.02
33		4-Chloroaniline	40.000	39.170	2.1	129	-0.01
34	C	Hexachlorobutadiene	40.000	38.819	3.0	129	-0.01
35		Caprolactam	40.000	41.664	-4.2	131	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.008	-0.0	129	0.00
37		2-Methylnaphthalene	40.000	38.334	4.2	126	-0.01
38		1-Methylnaphthalene	40.000	38.191	4.5	126	-0.01
39	I	Acenaphthene-d10	20.000	20.000	0.0	133	-0.01
40		1,2,4,5-Tetrachlorobenzene	40.000	37.527	6.2	125	-0.01
41	P	Hexachlorocyclopentadiene	40.000	40.511	-1.3	127	-0.01
42	S	2,4,6-Tribromophenol	80.000	77.501	3.1	128	0.00
43	C	2,4,6-Trichlorophenol	40.000	39.196	2.0	128	-0.01
44		2,4,5-Trichlorophenol	40.000	39.542	1.1	130	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102318\
Data File : BF110080.D
Acq On : 23 Oct 2018 16:30
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 24 09:06:21 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 23 15:06:18 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	75.311	5.9	125	-0.01
46		1,1'-Biphenyl	40.000	36.927	7.7	125	-0.01
47		2-Chloronaphthalene	40.000	37.670	5.8	126	-0.01
48		2-Nitroaniline	40.000	40.705	-1.8	132	0.00
49		Acenaphthylene	40.000	37.660	5.9	126	-0.01
50		Dimethylphthalate	40.000	38.521	3.7	130	0.00
51		2,6-Dinitrotoluene	40.000	41.017	-2.5	131	0.00
52	C	Acenaphthene	40.000	38.411	4.0	129	-0.01
53		3-Nitroaniline	40.000	40.435	-1.1	130	0.00
54	P	2,4-Dinitrophenol	40.000	47.929	-19.8	170	0.00
55		Dibenzofuran	40.000	37.168	7.1	124	-0.01
56	P	4-Nitrophenol	40.000	42.760	-6.9	132	0.00
57		2,4-Dinitrotoluene	40.000	42.743	-6.9	135	0.00
58		Fluorene	40.000	37.281	6.8	126	-0.01
59		2,3,4,6-Tetrachlorophenol	40.000	39.796	0.5	130	-0.01
60		Diethylphthalate	40.000	38.279	4.3	128	-0.01
61		4-Chlorophenyl-phenylether	40.000	37.091	7.3	126	-0.01
62		4-Nitroaniline	40.000	40.818	-2.0	133	0.00
63		Azobenzene	40.000	38.229	4.4	129	-0.01
64	I	Phenanthrene-d10	20.000	20.000	0.0	132	-0.01
65		4,6-Dinitro-2-methylphenol	40.000	44.071	-10.2	141	-0.01
66	c	n-Nitrosodiphenylamine	40.000	38.154	4.6	128	-0.01
67		4-Bromophenyl-phenylether	40.000	38.485	3.8	128	-0.01
68		Hexachlorobenzene	40.000	38.977	2.6	130	-0.02
69		Atrazine	40.000	38.430	3.9	126	0.00
70	C	Pentachlorophenol	40.000	43.608	-9.0	130	-0.01
71		Phenanthrene	40.000	36.975	7.6	125	-0.01
72		Anthracene	40.000	37.298	6.8	125	-0.01
73		Carbazole	40.000	37.928	5.2	128	0.00
74		Di-n-butylphthalate	40.000	37.901	5.2	126	-0.01
75	C	Fluoranthene	40.000	37.574	6.1	127	-0.01
76	I	Chrysene-d12	20.000	20.000	0.0	130	0.00
77		Benzidine	40.000	38.422	3.9	124	-0.01
78		Pyrene	40.000	38.801	3.0	127	-0.01
79	S	Terphenyl-d14	80.000	73.963	7.5	124	-0.01
80		Butylbenzylphthalate	40.000	40.694	-1.7	130	-0.01
81		Benzo(a)anthracene	40.000	39.121	2.2	127	0.00
82		3,3'-Dichlorobenzidine	40.000	37.026	7.4	118	0.00
83		Chrysene	40.000	38.010	5.0	124	-0.01
84		Bis(2-ethylhexyl)phthalate	40.000	39.489	1.3	126	-0.02
85	c	Di-n-octyl phthalate	40.000	40.124	-0.3	129	-0.01
86		Indeno(1,2,3-cd)pyrene	40.000	38.754	3.1	139	-0.02
87	I	Perylene-d12	20.000	20.000	0.0	136	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102318\
Data File : BF110080.D
Acq On : 23 Oct 2018 16:30
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 24 09:06:21 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 23 15:06:18 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	41.176	-2.9	137	-0.01
89	Benzo(k)fluoranthene	40.000	36.013	10.0	121	-0.01
90 C	Benzo(a)pyrene	40.000	39.289	1.8	133	-0.02
91	Dibenzo(a,h)anthracene	40.000	39.498	1.3	137	-0.02
92	Benzo(g,h,i)perylene	40.000	39.804	0.5	139	-0.02

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