Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF112918\
Data File : BF111087.D
Acq On : 29 Nov 2018 13:08
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 29 13:41:44 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF112618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Nov 29 10:35:31 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	109	0.00
2		1,4-Dioxane	40.000	40.136	-0.3	110	0.01
3		Pyridine	40.000	35.263	11.8	94	0.01
4		n-Nitrosodimethylamine	40.000	39.040	2.4	108	0.02
5	S	2-Fluorophenol	80.000	81.472	-1.8	107	0.00
6		Aniline	40.000	41.526	-3.8	110	0.00
7	S	Phenol-d6	80.000	80.737	-0.9	111	0.00
8		2-Chlorophenol	40.000	41.222	-3.1	112	0.00
9		Benzaldehyde	40.000	39.854	0.4	109	0.00
10	C	Phenol	40.000	40.569	-1.4	108	0.00
11		bis(2-Chloroethyl)ether	40.000	39.978	0.1	111	0.00
12		1,3-Dichlorobenzene	40.000	40.819	-2.0	112	0.00
13	C	1,4-Dichlorobenzene	40.000	40.430	-1.1	111	0.00
14		1,2-Dichlorobenzene	40.000	40.379	-0.9	111	0.00
15		Benzyl Alcohol	40.000	40.043	-0.1	107	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	40.518	-1.3	112	0.00
17		2-Methylphenol	40.000	39.509	1.2	108	0.00
18		Hexachloroethane	40.000	41.001	-2.5	112	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.889	0.3	109	0.00
20		3+4-Methylphenols	40.000	38.767	3.1	106	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	110	0.00
22		Acetophenone	40.000	40.826	-2.1	111	0.00
23	S	Nitrobenzene-d5	80.000	80.944	-1.2	110	0.00
24		Nitrobenzene	40.000	40.239	-0.6	111	0.00
25		Isophorone	40.000	39.882	0.3	108	0.00
26	C	2-Nitrophenol	40.000	41.622	-4.1	110	0.00
27		2,4-Dimethylphenol	40.000	40.907	-2.3	108	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.728	-1.8	111	0.00
29	C	2,4-Dichlorophenol	40.000	41.226	-3.1	108	0.00
30		1,2,4-Trichlorobenzene	40.000	40.523	-1.3	109	0.00
31		Naphthalene	40.000	40.075	-0.2	110	0.00
32		Benzoic acid	40.000	37.146	7.1	98	0.00
33		4-Chloroaniline	40.000	41.872	-4.7	112	0.00
34	C	Hexachlorobutadiene	40.000	40.705	-1.8	110	0.00
35		Caprolactam	40.000	38.321	4.2	103	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.093	-0.2	106	0.00
37		2-Methylnaphthalene	40.000	39.692	0.8	108	0.00
38		1-Methylnaphthalene	40.000	38.811	3.0	107	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	106	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	40.847	-2.1	107	0.00
41	P	Hexachlorocyclopentadiene	40.000	51.948	-29.9#	171	0.00
42	S	2,4,6-Tribromophenol	80.000	80.691	-0.9	105	0.00
43	C	2,4,6-Trichlorophenol	40.000	40.738	-1.8	104	0.00
44		2,4,5-Trichlorophenol	40.000	40.264	-0.7	103	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF112918\
Data File : BF111087.D
Acq On : 29 Nov 2018 13:08
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 29 13:41:44 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF112618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Nov 29 10:35:31 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	79.118	1.1	107	0.00
46		1,1'-Biphenyl	40.000	40.336	-0.8	106	0.00
47		2-Chloronaphthalene	40.000	40.303	-0.8	106	0.00
48		2-Nitroaniline	40.000	40.818	-2.0	107	0.00
49		Acenaphthylene	40.000	39.225	1.9	104	0.00
50		Dimethylphthalate	40.000	38.619	3.5	103	0.00
51		2,6-Dinitrotoluene	40.000	39.415	1.5	103	0.00
52	C	Acenaphthene	40.000	39.868	0.3	106	0.00
53		3-Nitroaniline	40.000	39.998	0.0	104	0.00
54	P	2,4-Dinitrophenol	40.000	40.937	-2.3	102	0.00
55		Dibenzofuran	40.000	38.748	3.1	104	0.00
56	P	4-Nitrophenol	40.000	35.344	11.6	90	0.00
57		2,4-Dinitrotoluene	40.000	39.599	1.0	103	0.00
58		Fluorene	40.000	39.090	2.3	104	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	37.146	7.1	98	0.00
60		Diethylphthalate	40.000	39.004	2.5	105	0.00
61		4-Chlorophenyl-phenylether	40.000	39.687	0.8	105	0.00
62		4-Nitroaniline	40.000	40.977	-2.4	106	0.00
63		Azobenzene	40.000	38.661	3.3	103	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	102	0.00
65		4,6-Dinitro-2-methylphenol	40.000	39.373	1.6	97	0.00
66	c	n-Nitrosodiphenylamine	40.000	41.648	-4.1	105	0.00
67		4-Bromophenyl-phenylether	40.000	41.143	-2.9	103	0.00
68		Hexachlorobenzene	40.000	41.254	-3.1	104	0.00
69		Atrazine	40.000	38.818	3.0	97	0.00
70	C	Pentachlorophenol	40.000	36.768	8.1	87	0.00
71		Phenanthrene	40.000	40.342	-0.9	103	0.00
72		Anthracene	40.000	39.948	0.1	101	0.00
73		Carbazole	40.000	40.374	-0.9	102	0.00
74		Di-n-butylphthalate	40.000	39.529	1.2	100	0.00
75	C	Fluoranthene	40.000	38.683	3.3	99	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	90	0.00
77		Benzidine	40.000	45.784	-14.5	114	0.00
78		Pyrene	40.000	44.115	-10.3	99	0.00
79	S	Terphenyl-d14	80.000	86.025	-7.5	98	0.00
80		Butylbenzylphthalate	40.000	43.253	-8.1	95	0.00
81		Benzo(a)anthracene	40.000	40.191	-0.5	89	0.00
82		3,3'-Dichlorobenzidine	40.000	40.583	-1.5	89	0.00
83		Chrysene	40.000	39.504	1.2	88	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	42.047	-5.1	93	0.00
85	c	Di-n-octyl phthalate	40.000	38.289	4.3	84	-0.01
86		Indeno(1,2,3-cd)pyrene	40.000	42.179	-5.4	99	0.00
87	I	Perylene-d12	20.000	20.000	0.0	90	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF112918\
Data File : BF111087.D
Acq On : 29 Nov 2018 13:08
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 29 13:41:44 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF112618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Nov 29 10:35:31 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	38.775	3.1	85	0.00
89	Benzo(k)fluoranthene	40.000	40.358	-0.9	88	0.00
90 C	Benzo(a)pyrene	40.000	39.236	1.9	88	0.00
91	Dibenzo(a,h)anthracene	40.000	44.223	-10.6	99	0.00
92	Benzo(g,h,i)perylene	40.000	45.707	-14.3	101	0.00

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