Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF111517\
Data File : BF100578.D
Acq On : 15 Nov 2017 10:12
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 15 14:04:59 2017
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 14 14:17:16 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	73	0.00
2		1,4-Dioxane	40.000	39.914	0.2	73	0.00
3		Pyridine	40.000	39.432	1.4	70	0.00
4		n-Nitrosodimethylamine	40.000	39.904	0.2	71	0.00
5	S	2-Fluorophenol	80.000	79.649	0.4	73	0.00
6		Aniline	40.000	39.534	1.2	71	0.00
7	S	Phenol-d6	80.000	80.385	-0.5	73	0.00
8		2-Chlorophenol	40.000	41.280	-3.2	76	0.00
9		Benzaldehyde	40.000	37.083	7.3	68	0.00
10	C	Phenol	40.000	40.951	-2.4	76	0.00
11		bis(2-Chloroethyl)ether	40.000	39.356	1.6	72	0.00
12		1,3-Dichlorobenzene	40.000	40.284	-0.7	74	0.00
13	C	1,4-Dichlorobenzene	40.000	40.251	-0.6	74	0.00
14		1,2-Dichlorobenzene	40.000	40.145	-0.4	74	0.00
15		Benzyl Alcohol	40.000	40.230	-0.6	72	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	38.404	4.0	70	0.00
17		2-Methylphenol	40.000	40.689	-1.7	74	0.00
18		Hexachloroethane	40.000	40.357	-0.9	72	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	36.874	7.8	68	0.00
20		3+4-Methylphenols	40.000	38.626	3.4	70	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	71	0.00
22		Acetophenone	40.000	37.489	6.3	69	0.00
23	S	Nitrobenzene-d5	80.000	91.261	-14.1	78	0.00
24		Nitrobenzene	40.000	45.744	-14.4	76	0.00
25		Isophorone	40.000	38.784	3.0	69	0.00
26	C	2-Nitrophenol	40.000	51.338	-28.3#	95	0.00
27		2,4-Dimethylphenol	40.000	40.998	-2.5	72	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.664	3.3	69	0.00
29	C	2,4-Dichlorophenol	40.000	41.801	-4.5	72	0.00
30		1,2,4-Trichlorobenzene	40.000	40.554	-1.4	73	0.00
31		Naphthalene	40.000	39.618	1.0	72	0.00
32		Benzoic acid	40.000	39.861	0.3	76	0.00
33		4-Chloroaniline	40.000	41.129	-2.8	73	0.00
34	C	Hexachlorobutadiene	40.000	39.513	1.2	71	0.00
35		Caprolactam	40.000	39.900	0.3	71	0.00
36	C	4-Chloro-3-methylphenol	40.000	39.927	0.2	69	0.00
37		2-Methylnaphthalene	40.000	38.562	3.6	70	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	67	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	42.093	-5.2	71	0.00
40	P	Hexachlorocyclopentadiene	40.000	45.909	-14.8	72	0.00
41	S	2,4,6-Tribromophenol	80.000	86.317	-7.9	70	0.00
42	C	2,4,6-Trichlorophenol	40.000	43.532	-8.8	70	0.00
43		2,4,5-Trichlorophenol	40.000	44.800	-12.0	73	0.00
44	S	2-Fluorobiphenyl	80.000	79.127	1.1	69	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF111517\
Data File : BF100578.D
Acq On : 15 Nov 2017 10:12
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 15 14:04:59 2017
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 14 14:17:16 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	40.304	-0.8	70	0.00
46		2-Chloronaphthalene	40.000	41.589	-4.0	70	0.00
47		2-Nitroaniline	40.000	47.724	-19.3	80	0.00
48		Acenaphthylene	40.000	41.036	-2.6	70	0.00
49		Dimethylphthalate	40.000	40.542	-1.4	70	0.00
50		2,6-Dinitrotoluene	40.000	45.954	-14.9	74	0.00
51	C	Acenaphthene	40.000	41.342	-3.4	71	0.00
52		3-Nitroaniline	40.000	47.248	-18.1	77	0.00
53	P	2,4-Dinitrophenol	40.000	60.983	-52.5#	131	0.00
54		Dibenzofuran	40.000	40.419	-1.0	69	0.00
55	P	4-Nitrophenol	40.000	45.214	-13.0	73	0.00
56		2,4-Dinitrotoluene	40.000	48.067	-20.2	76	0.00
57		Fluorene	40.000	41.335	-3.3	69	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	41.778	-4.4	67	0.00
59		Diethylphthalate	40.000	38.859	2.9	66	0.00
60		4-Chlorophenyl-phenylether	40.000	41.211	-3.0	69	0.00
61		4-Nitroaniline	40.000	48.280	-20.7	77	0.00
62		Azobenzene	40.000	38.648	3.4	66	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	66	0.00
64		4,6-Dinitro-2-methylphenol	40.000	54.862	-37.2#	102	0.00
65	c	n-Nitrosodiphenylamine	40.000	39.810	0.5	67	0.00
66		4-Bromophenyl-phenylether	40.000	40.305	-0.8	67	0.00
67		Hexachlorobenzene	40.000	40.948	-2.4	68	0.00
68		Atrazine	40.000	39.699	0.8	65	0.00
69	C	Pentachlorophenol	40.000	41.060	-2.7	67	0.00
70		Phenanthrene	40.000	39.755	0.6	69	0.00
71		Anthracene	40.000	39.869	0.3	69	0.00
72		Carbazole	40.000	40.530	-1.3	70	0.00
73		Di-n-butylphthalate	40.000	40.912	-2.3	69	0.00
74	C	Fluoranthene	40.000	39.988	0.0	69	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	68	0.00
76		Benzidine	40.000	50.799	-27.0#	80	0.00
77		Pyrene	40.000	40.308	-0.8	68	0.00
78	S	Terphenyl-d14	80.000	76.319	4.6	68	0.00
79		Butylbenzylphthalate	40.000	42.818	-7.0	71	0.00
80		Benzo(a)anthracene	40.000	41.605	-4.0	71	0.00
81		3,3'-Dichlorobenzidine	40.000	43.140	-7.9	69	0.00
82		Chrysene	40.000	39.803	0.5	68	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	43.048	-7.6	69	0.00
84	c	Di-n-octyl phthalate	40.000	41.181	-3.0	66	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	40.878	-2.2	71	0.00
86	I	Perylene-d12	20.000	20.000	0.0	66	0.00
87		Benzo(b)fluoranthene	40.000	42.045	-5.1	70	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF111517\
Data File : BF100578.D
Acq On : 15 Nov 2017 10:12
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 15 14:04:59 2017
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 14 14:17:16 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	41.018	-2.5	64	0.00
89 C	Benzo(a)pyrene	40.000	41.663	-4.2	67	0.00
90	Dibenzo(a,h)anthracene	40.000	41.821	-4.6	70	0.00
91	Benzo(g,h,i)perylene	40.000	43.854	-9.6	74	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 1