Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF121217\
Data File : BF101327.D
Acq On : 12 Dec 2017 13:58
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 13 01:18:14 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 08 18:17:31 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	88	0.00
2		1,4-Dioxane	40.000	40.335	-0.8	86	0.00
3		Pyridine	40.000	37.692	5.8	74	0.00
4		n-Nitrosodimethylamine	40.000	43.949	-9.9	92	0.00
5	S	2-Fluorophenol	80.000	85.923	-7.4	92	0.00
6		Aniline	40.000	42.251	-5.6	91	0.00
7	S	Phenol-d6	80.000	84.544	-5.7	91	0.00
8		2-Chlorophenol	40.000	42.160	-5.4	91	0.00
9		Benzaldehyde	40.000	40.917	-2.3	92	0.00
10	C	Phenol	40.000	41.964	-4.9	91	0.00
11		bis(2-Chloroethyl)ether	40.000	42.013	-5.0	90	0.00
12		1,3-Dichlorobenzene	40.000	42.160	-5.4	90	0.00
13	C	1,4-Dichlorobenzene	40.000	41.577	-3.9	90	0.00
14		1,2-Dichlorobenzene	40.000	41.600	-4.0	91	0.00
15		Benzyl Alcohol	40.000	43.528	-8.8	92	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	40.961	-2.4	87	0.00
17		2-Methylphenol	40.000	42.514	-6.3	93	0.00
18		Hexachloroethane	40.000	42.310	-5.8	92	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	41.674	-4.2	92	0.00
20		3+4-Methylphenols	40.000	41.227	-3.1	89	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	89	0.00
22		Acetophenone	40.000	41.248	-3.1	91	0.00
23	S	Nitrobenzene-d5	80.000	83.392	-4.2	91	0.00
24		Nitrobenzene	40.000	41.876	-4.7	92	0.00
25		Isophorone	40.000	41.443	-3.6	92	0.00
26	C	2-Nitrophenol	40.000	42.079	-5.2	93	0.00
27		2,4-Dimethylphenol	40.000	41.495	-3.7	91	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.605	-1.5	90	0.00
29	C	2,4-Dichlorophenol	40.000	42.439	-6.1	91	0.00
30		1,2,4-Trichlorobenzene	40.000	41.332	-3.3	92	0.00
31		Naphthalene	40.000	41.041	-2.6	91	0.00
32		Benzoic acid	40.000	40.736	-1.8	93	0.01
33		4-Chloroaniline	40.000	41.281	-3.2	89	0.00
34	C	Hexachlorobutadiene	40.000	41.944	-4.9	93	0.00
35		Caprolactam	40.000	40.125	-0.3	86	0.00
36	C	4-Chloro-3-methylphenol	40.000	42.612	-6.5	94	0.00
37		2-Methylnaphthalene	40.000	40.935	-2.3	90	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	89	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	41.176	-2.9	91	0.00
40	P	Hexachlorocyclopentadiene	40.000	38.600	3.5	90	0.00
41	S	2,4,6-Tribromophenol	80.000	78.677	1.7	86	0.00
42	C	2,4,6-Trichlorophenol	40.000	40.476	-1.2	88	0.00
43		2,4,5-Trichlorophenol	40.000	40.272	-0.7	87	0.00
44	S	2-Fluorobiphenyl	80.000	82.370	-3.0	91	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF121217\
Data File : BF101327.D
Acq On : 12 Dec 2017 13:58
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 13 01:18:14 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 08 18:17:31 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	40.403	-1.0	90	0.00
46		2-Chloronaphthalene	40.000	41.123	-2.8	91	0.00
47		2-Nitroaniline	40.000	40.937	-2.3	89	0.00
48		Acenaphthylene	40.000	40.428	-1.1	89	0.00
49		Dimethylphthalate	40.000	39.685	0.8	88	0.00
50		2,6-Dinitrotoluene	40.000	40.760	-1.9	91	0.00
51	C	Acenaphthene	40.000	40.301	-0.8	91	0.00
52		3-Nitroaniline	40.000	40.173	-0.4	88	0.00
53	P	2,4-Dinitrophenol	40.000	41.495	-3.7	92	0.00
54		Dibenzofuran	40.000	39.448	1.4	87	0.00
55	P	4-Nitrophenol	40.000	36.806	8.0	78	0.01
56		2,4-Dinitrotoluene	40.000	38.522	3.7	83	0.00
57		Fluorene	40.000	40.474	-1.2	89	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	36.060	9.8	79	0.00
59		Diethylphthalate	40.000	40.279	-0.7	88	0.00
60		4-Chlorophenyl-phenylether	40.000	41.650	-4.1	91	0.00
61		4-Nitroaniline	40.000	39.029	2.4	86	0.00
62		Azobenzene	40.000	40.250	-0.6	89	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	85	0.00
64		4,6-Dinitro-2-methylphenol	40.000	38.151	4.6	79	0.00
65	c	n-Nitrosodiphenylamine	40.000	41.172	-2.9	87	0.00
66		4-Bromophenyl-phenylether	40.000	41.632	-4.1	88	0.00
67		Hexachlorobenzene	40.000	41.340	-3.4	88	0.00
68		Atrazine	40.000	38.146	4.6	82	0.00
69	C	Pentachlorophenol	40.000	38.724	3.2	78	0.00
70		Phenanthrene	40.000	40.379	-0.9	86	0.00
71		Anthracene	40.000	40.589	-1.5	85	0.00
72		Carbazole	40.000	40.033	-0.1	85	0.00
73		Di-n-butylphthalate	40.000	40.599	-1.5	85	0.00
74	C	Fluoranthene	40.000	39.206	2.0	84	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	83	0.00
76		Benzidine	40.000	37.548	6.1	77	0.00
77		Pyrene	40.000	41.252	-3.1	84	0.00
78	S	Terphenyl-d14	80.000	82.537	-3.2	86	0.00
79		Butylbenzylphthalate	40.000	41.413	-3.5	83	0.00
80		Benzo(a)anthracene	40.000	39.876	0.3	82	0.00
81		3,3'-Dichlorobenzidine	40.000	40.729	-1.8	83	0.00
82		Chrysene	40.000	40.559	-1.4	84	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	40.391	-1.0	81	0.00
84	c	Di-n-octyl phthalate	40.000	40.932	-2.3	82	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	38.526	3.7	80	0.04
86	I	Perylene-d12	20.000	20.000	0.0	82	0.01
87		Benzo(b)fluoranthene	40.000	41.570	-3.9	86	0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF121217\
Data File : BF101327.D
Acq On : 12 Dec 2017 13:58
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 13 01:18:14 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 08 18:17:31 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	38.147	4.6	76	0.01
89 C	Benzo(a)pyrene	40.000	39.918	0.2	81	0.02
90	Dibenzo(a,h)anthracene	40.000	38.739	3.2	80	0.03
91	Benzo(g,h,i)perylene	40.000	38.835	2.9	82	0.04

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