Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF120817\
Data File : BF101235.D
Acq On : 8 Dec 2017 14:53
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 08 18:21:14 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	37.175	7.1	79	0.00
3		Pyridine	40.000	38.380	4.0	76	0.00
4		n-Nitrosodimethylamine	40.000	42.696	-6.7	90	0.00
5	S	2-Fluorophenol	80.000	77.579	3.0	83	0.00
6		Aniline	40.000	39.981	0.0	86	0.00
7	S	Phenol-d6	80.000	79.557	0.6	86	0.00
8		2-Chlorophenol	40.000	39.393	1.5	85	0.00
9		Benzaldehyde	40.000	36.226	9.4	82	0.00
10	C	Phenol	40.000	39.316	1.7	85	0.00
11		bis(2-Chloroethyl)ether	40.000	38.259	4.4	83	0.00
12		1,3-Dichlorobenzene	40.000	39.184	2.0	84	0.00
13	C	1,4-Dichlorobenzene	40.000	38.629	3.4	84	0.00
14		1,2-Dichlorobenzene	40.000	39.166	2.1	86	0.00
15		Benzyl Alcohol	40.000	40.322	-0.8	85	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	37.257	6.9	79	0.00
17		2-Methylphenol	40.000	39.561	1.1	86	0.00
18		Hexachloroethane	40.000	39.721	0.7	86	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	38.863	2.8	86	0.00
20		3+4-Methylphenols	40.000	39.648	0.9	85	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	87	0.00
22		Acetophenone	40.000	39.401	1.5	84	0.00
23	S	Nitrobenzene-d5	80.000	79.072	1.2	84	0.00
24		Nitrobenzene	40.000	39.979	0.1	86	0.00
25		Isophorone	40.000	38.766	3.1	83	0.00
26	C	2-Nitrophenol	40.000	39.405	1.5	84	0.00
27		2,4-Dimethylphenol	40.000	40.197	-0.5	86	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.516	3.7	83	0.00
29	C	2,4-Dichlorophenol	40.000	39.979	0.1	83	0.00
30		1,2,4-Trichlorobenzene	40.000	39.636	0.9	85	0.00
31		Naphthalene	40.000	39.448	1.4	85	0.00
32		Benzoic acid	40.000	39.019	2.5	87	0.00
33		4-Chloroaniline	40.000	39.979	0.1	84	0.00
34	C	Hexachlorobutadiene	40.000	40.062	-0.2	86	0.00
35		Caprolactam	40.000	38.270	4.3	80	0.00
36	C	4-Chloro-3-methylphenol	40.000	39.792	0.5	85	0.00
37		2-Methylnaphthalene	40.000	39.025	2.4	83	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	84	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	40.497	-1.2	85	0.00
40	P	Hexachlorocyclopentadiene	40.000	33.537	16.2	71	0.00
41	S	2,4,6-Tribromophenol	80.000	76.795	4.0	80	0.00
42	C	2,4,6-Trichlorophenol	40.000	39.973	0.1	82	0.00
43		2,4,5-Trichlorophenol	40.000	39.881	0.3	82	0.00
44	S	2-Fluorobiphenyl	80.000	80.734	-0.9	84	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF120817\
Data File : BF101235.D
Acq On : 8 Dec 2017 14:53
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 08 18:21:14 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	39.969	0.1	84	0.00
46		2-Chloronaphthalene	40.000	40.429	-1.1	84	0.00
47		2-Nitroaniline	40.000	39.113	2.2	81	0.00
48		Acenaphthylene	40.000	39.167	2.1	82	0.00
49		Dimethylphthalate	40.000	38.791	3.0	81	0.00
50		2,6-Dinitrotoluene	40.000	39.403	1.5	83	0.00
51	C	Acenaphthene	40.000	39.445	1.4	84	0.00
52		3-Nitroaniline	40.000	39.211	2.0	81	0.00
53	P	2,4-Dinitrophenol	40.000	37.142	7.1	77	0.00
54		Dibenzofuran	40.000	39.366	1.6	82	0.00
55	P	4-Nitrophenol	40.000	36.534	8.7	73	0.00
56		2,4-Dinitrotoluene	40.000	37.674	5.8	77	0.00
57		Fluorene	40.000	39.768	0.6	83	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	37.906	5.2	78	0.00
59		Diethylphthalate	40.000	38.574	3.6	79	0.00
60		4-Chlorophenyl-phenylether	40.000	40.071	-0.2	83	0.00
61		4-Nitroaniline	40.000	38.073	4.8	79	0.00
62		Azobenzene	40.000	37.890	5.3	79	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	81	0.00
64		4,6-Dinitro-2-methylphenol	40.000	37.523	6.2	74	0.00
65	c	n-Nitrosodiphenylamine	40.000	39.450	1.4	80	0.00
66		4-Bromophenyl-phenylether	40.000	39.103	2.2	79	0.00
67		Hexachlorobenzene	40.000	38.593	3.5	79	0.00
68		Atrazine	40.000	38.801	3.0	80	0.00
69	C	Pentachlorophenol	40.000	43.177	-7.9	83	0.00
70		Phenanthrene	40.000	38.560	3.6	79	0.00
71		Anthracene	40.000	38.763	3.1	78	0.00
72		Carbazole	40.000	37.990	5.0	78	0.00
73		Di-n-butylphthalate	40.000	37.800	5.5	76	0.00
74	C	Fluoranthene	40.000	37.544	6.1	78	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	77	0.00
76		Benzidine	40.000	39.427	1.4	76	0.00
77		Pyrene	40.000	40.922	-2.3	78	0.00
78	S	Terphenyl-d14	80.000	81.686	-2.1	79	0.00
79		Butylbenzylphthalate	40.000	39.469	1.3	74	0.00
80		Benzo(a)anthracene	40.000	39.458	1.4	76	0.00
81		3,3'-Dichlorobenzidine	40.000	39.696	0.8	75	0.00
82		Chrysene	40.000	39.330	1.7	76	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	38.889	2.8	73	0.00
84	c	Di-n-octyl phthalate	40.000	37.947	5.1	71	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	34.932	12.7	68	0.00
86	I	Perylene-d12	20.000	20.000	0.0	70	0.00
87		Benzo(b)fluoranthene	40.000	42.300	-5.7	75	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF120817\
Data File : BF101235.D
Acq On : 8 Dec 2017 14:53
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 08 18:21:14 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	38.980	2.6	66	0.00
89 C	Benzo(a)pyrene	40.000	39.773	0.6	69	0.00
90	Dibenzo(a,h)anthracene	40.000	37.795	5.5	67	0.00
91	Benzo(g,h,i)perylene	40.000	38.444	3.9	69	0.00

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