Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF110317\
Data File : BF100131.D
Acq On : 3 Nov 2017 13:42
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 03 15:49:27 2017
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 03 15:46:10 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	100	0.00
2		1,4-Dioxane	40.000	39.654	0.9	98	0.00
3		Pyridine	40.000	40.226	-0.6	92	0.00
4		n-Nitrosodimethylamine	40.000	40.762	-1.9	93	0.00
5	S	2-Fluorophenol	80.000	83.466	-4.3	101	0.00
6		Aniline	40.000	39.090	2.3	96	0.00
7	S	Phenol-d6	80.000	80.309	-0.4	99	0.00
8		2-Chlorophenol	40.000	41.434	-3.6	100	0.00
9		Benzaldehyde	40.000	35.482	11.3	87	0.00
10	C	Phenol	40.000	39.547	1.1	96	0.00
11		bis(2-Chloroethyl)ether	40.000	41.081	-2.7	99	0.00
12		1,3-Dichlorobenzene	40.000	40.592	-1.5	100	0.00
13	C	1,4-Dichlorobenzene	40.000	41.004	-2.5	101	0.00
14		1,2-Dichlorobenzene	40.000	40.555	-1.4	100	0.00
15		Benzyl Alcohol	40.000	39.426	1.4	99	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	38.985	2.5	96	0.00
17		2-Methylphenol	40.000	41.117	-2.8	101	0.00
18		Hexachloroethane	40.000	40.521	-1.3	100	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.709	0.7	100	0.00
20		3+4-Methylphenols	40.000	42.661	-6.7	107	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	100	0.00
22		Acetophenone	40.000	39.676	0.8	101	0.00
23	S	Nitrobenzene-d5	80.000	77.817	2.7	98	0.00
24		Nitrobenzene	40.000	39.244	1.9	96	0.00
25		Isophorone	40.000	39.525	1.2	98	0.00
26	C	2-Nitrophenol	40.000	42.669	-6.7	101	0.00
27		2,4-Dimethylphenol	40.000	41.074	-2.7	102	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.988	0.0	100	0.00
29	C	2,4-Dichlorophenol	40.000	42.701	-6.8	105	0.00
30		1,2,4-Trichlorobenzene	40.000	40.112	-0.3	99	0.00
31		Naphthalene	40.000	39.669	0.8	99	0.00
32		Benzoic acid	40.000	39.175	2.1	99	0.00
33		4-Chloroaniline	40.000	39.236	1.9	96	0.00
34	C	Hexachlorobutadiene	40.000	40.751	-1.9	103	0.00
35		Caprolactam	40.000	39.343	1.6	96	0.00
36	C	4-Chloro-3-methylphenol	40.000	39.939	0.2	99	0.00
37		2-Methylnaphthalene	40.000	39.410	1.5	100	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	99	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	40.993	-2.5	101	0.00
40	P	Hexachlorocyclopentadiene	40.000	45.911	-14.8	125	0.00
41	S	2,4,6-Tribromophenol	80.000	75.681	5.4	93	0.00
42	C	2,4,6-Trichlorophenol	40.000	40.647	-1.6	101	0.00
43		2,4,5-Trichlorophenol	40.000	38.424	3.9	91	0.00
44	S	2-Fluorobiphenyl	80.000	79.207	1.0	96	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF110317\
Data File : BF100131.D
Acq On : 3 Nov 2017 13:42
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 03 15:49:27 2017
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 03 15:46:10 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	39.370	1.6	98	0.00
46		2-Chloronaphthalene	40.000	39.629	0.9	98	0.00
47		2-Nitroaniline	40.000	38.928	2.7	94	0.00
48		Acenaphthylene	40.000	38.752	3.1	97	0.00
49		Dimethylphthalate	40.000	36.787	8.0	91	0.00
50		2,6-Dinitrotoluene	40.000	39.409	1.5	95	0.00
51	C	Acenaphthene	40.000	38.838	2.9	97	0.00
52		3-Nitroaniline	40.000	38.738	3.2	94	0.00
53	P	2,4-Dinitrophenol	40.000	40.652	-1.6	101	0.00
54		Dibenzofuran	40.000	39.442	1.4	99	0.00
55	P	4-Nitrophenol	40.000	32.657	18.4	76	0.00
56		2,4-Dinitrotoluene	40.000	41.417	-3.5	101	0.00
57		Fluorene	40.000	38.221	4.4	96	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	39.121	2.2	94	0.00
59		Diethylphthalate	40.000	37.393	6.5	93	0.00
60		4-Chlorophenyl-phenylether	40.000	39.134	2.2	99	0.00
61		4-Nitroaniline	40.000	38.928	2.7	93	0.00
62		Azobenzene	40.000	36.761	8.1	90	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	88	0.00
64		4,6-Dinitro-2-methylphenol	40.000	41.082	-2.7	87	0.00
65	c	n-Nitrosodiphenylamine	40.000	41.701	-4.3	94	0.00
66		4-Bromophenyl-phenylether	40.000	42.610	-6.5	95	0.00
67		Hexachlorobenzene	40.000	41.338	-3.3	92	0.00
68		Atrazine	40.000	40.948	-2.4	89	0.00
69	C	Pentachlorophenol	40.000	38.827	2.9	86	0.00
70		Phenanthrene	40.000	38.287	4.3	86	0.00
71		Anthracene	40.000	38.805	3.0	88	0.00
72		Carbazole	40.000	38.012	5.0	85	0.00
73		Di-n-butylphthalate	40.000	38.632	3.4	85	0.00
74	C	Fluoranthene	40.000	34.944	12.6	78	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	72	0.00
76		Benzidine	40.000	37.268	6.8	68	0.00
77		Pyrene	40.000	42.456	-6.1	76	0.00
78	S	Terphenyl-d14	80.000	84.336	-5.4	77	0.00
79		Butylbenzylphthalate	40.000	41.338	-3.3	72	0.00
80		Benzo(a)anthracene	40.000	38.740	3.1	69	0.00
81		3,3'-Dichlorobenzidine	40.000	36.346	9.1	64	0.00
82		Chrysene	40.000	38.375	4.1	68	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	37.579	6.1	68	0.00
84	c	Di-n-octyl phthalate	40.000	37.012	7.5	65	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	37.322	6.7	70	0.00
86	I	Perylene-d12	20.000	20.000	0.0	70	0.00
87		Benzo(b)fluoranthene	40.000	39.306	1.7	72	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF110317\
Data File : BF100131.D
Acq On : 3 Nov 2017 13:42
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 03 15:49:27 2017
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 03 15:46:10 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	38.620	3.5	65	0.00
89 C	Benzo(a)pyrene	40.000	38.882	2.8	68	0.00
90	Dibenzo(a,h)anthracene	40.000	38.282	4.3	70	0.00
91	Benzo(g,h,i)perylene	40.000	38.402	4.0	69	0.00

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