Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF110716\
Data File : BF091050.D
Acq On : 7 Nov 2016 12:43
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 07 17:42:21 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 04 11:03:12 2016
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	103	0.01
2		1,4-Dioxane	40.000	37.922	5.2	99	0.00
3		Pyridine	40.000	46.896	-17.2	114	0.01
4		n-Nitrosodimethylamine	40.000	39.975	0.1	97	0.01
5	S	2-Fluorophenol	80.000	84.173	-5.2	102	0.01
6		Aniline	40.000	39.935	0.2	93	0.01
7	S	Phenol-d6	80.000	76.198	4.8	92	0.01
8		2-Chlorophenol	40.000	39.673	0.8	99	0.00
9		Benzaldehyde	40.000	37.910	5.2	94	0.01
10	C	Phenol	40.000	40.448	-1.1	104	0.01
11		bis(2-Chloroethyl)ether	40.000	41.323	-3.3	104	0.01
12		1,3-Dichlorobenzene	40.000	41.984	-5.0	103	0.01
13	C	1,4-Dichlorobenzene	40.000	38.900	2.8	100	0.01
14		1,2-Dichlorobenzene	40.000	41.648	-4.1	100	0.01
15		Benzyl Alcohol	40.000	41.424	-3.6	102	0.01
16		2,2'-oxybis(1-Chloropropane	40.000	35.177	12.1	91	0.00
17		2-Methylphenol	40.000	38.119	4.7	90	0.00
18		Hexachloroethane	40.000	40.002	-0.0	98	0.01
19	P	n-Nitroso-di-n-propylamine	40.000	40.963	-2.4	107	0.01
20		3+4-Methylphenols	40.000	42.464	-6.2	99	0.01
21	I	Naphthalene-d8	20.000	20.000	0.0	101	0.01
22		Acetophenone	40.000	41.260	-3.1	103	0.01
23	S	Nitrobenzene-d5	80.000	79.959	0.1	103	0.00
24		Nitrobenzene	40.000	42.140	-5.4	99	0.01
25		Isophorone	40.000	38.115	4.7	99	0.01
26	C	2-Nitrophenol	40.000	44.130	-10.3	110	0.01
27		2,4-Dimethylphenol	40.000	40.810	-2.0	96	0.01
28		bis(2-Chloroethoxy)methane	40.000	41.286	-3.2	94	0.01
29	C	2,4-Dichlorophenol	40.000	40.239	-0.6	98	0.00
30		1,2,4-Trichlorobenzene	40.000	41.765	-4.4	101	0.01
31		Naphthalene	40.000	41.764	-4.4	101	0.01
32		Benzoic acid	40.000	38.839	2.9	98	0.00
33		4-Chloroaniline	40.000	41.739	-4.3	104	0.01
34	C	Hexachlorobutadiene	40.000	38.680	3.3	99	0.00
35		Caprolactam	40.000	45.283	-13.2	114	0.01
36	C	4-Chloro-3-methylphenol	40.000	38.200	4.5	91	0.01
37		2-Methylnaphthalene	40.000	38.059	4.9	98	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	93	0.01
39		1,2,4,5-Tetrachlorobenzene	40.000	46.004	-15.0	95	0.01
40	P	Hexachlorocyclopentadiene	40.000	40.559	-1.4	97	0.01
41	S	2,4,6-Tribromophenol	80.000	99.668	-24.6	121	0.02
42	C	2,4,6-Trichlorophenol	40.000	43.483	-8.7	93	0.01
43		2,4,5-Trichlorophenol	40.000	44.995	-12.5	98	0.01
44	S	2-Fluorobiphenyl	80.000	91.620	-14.5	107	0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF110716\
Data File : BF091050.D
Acq On : 7 Nov 2016 12:43
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 07 17:42:21 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 04 11:03:12 2016
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	41.486	-3.7	92	0.01
46		2-Chloronaphthalene	40.000	42.174	-5.4	92	0.01
47		2-Nitroaniline	40.000	38.110	4.7	79	0.01
48		Acenaphthylene	40.000	38.895	2.8	89	0.01
49		Dimethylphthalate	40.000	41.228	-3.1	96	0.02
50		2,6-Dinitrotoluene	40.000	48.107	-20.3	117	0.01
51	C	Acenaphthene	40.000	37.960	5.1	92	0.01
52		3-Nitroaniline	40.000	42.572	-6.4	95	0.02
53	P	2,4-Dinitrophenol	40.000	44.449	-11.1	113	0.01
54		Dibenzofuran	40.000	40.892	-2.2	97	0.01
55	P	4-Nitrophenol	40.000	39.679	0.8	92	0.01
56		2,4-Dinitrotoluene	40.000	40.714	-1.8	84	0.01
57		Fluorene	40.000	38.738	3.2	88	0.01
58		2,3,4,6-Tetrachlorophenol	40.000	43.978	-9.9	89	0.01
59		Diethylphthalate	40.000	37.894	5.3	85	0.01
60		4-Chlorophenyl-phenylether	40.000	45.758	-14.4	97	0.01
61		4-Nitroaniline	40.000	42.556	-6.4	92	0.02
62		Azobenzene	40.000	34.729	13.2	81	0.01
63	I	Phenanthrene-d10	20.000	20.000	0.0	100	0.02
64		4,6-Dinitro-2-methylphenol	40.000	41.937	-4.8	105	0.01
65	c	n-Nitrosodiphenylamine	40.000	40.655	-1.6	106	0.02
66		4-Bromophenyl-phenylether	40.000	39.859	0.4	108	0.02
67		Hexachlorobenzene	40.000	41.142	-2.9	97	0.01
68		Atrazine	40.000	34.242	14.4	76	0.01
69	C	Pentachlorophenol	40.000	43.294	-8.2	104	0.02
70		Phenanthrene	40.000	38.984	2.5	91	0.01
71		Anthracene	40.000	40.627	-1.6	108	0.02
72		Carbazole	40.000	40.065	-0.2	107	0.02
73		Di-n-butylphthalate	40.000	36.840	7.9	86	0.01
74	C	Fluoranthene	40.000	41.646	-4.1	110	0.02
75	I	Chrysene-d12	20.000	20.000	0.0	95	0.01
76		Benzidine	40.000	48.747	-21.9	106	0.02
77		Pyrene	40.000	40.845	-2.1	106	0.01
78	S	Terphenyl-d14	80.000	91.835	-14.8	104	0.02
79		Butylbenzylphthalate	40.000	40.991	-2.5	97	0.01
80		Benzo(a)anthracene	40.000	43.001	-7.5	103	0.01
81		3,3'-Dichlorobenzidine	40.000	47.833	-19.6	114	0.02
82		Chrysene	40.000	47.551	-18.9	109	0.02
83		Bis(2-ethylhexyl)phthalate	40.000	39.479	1.3	91	0.02
84	c	Di-n-octyl phthalate	40.000	38.068	4.8	85	0.01
85		Indeno(1,2,3-cd)pyrene	40.000	38.989	2.5	97	0.03
86	I	Perylene-d12	20.000	20.000	0.0	100	0.02
87		Benzo(b)fluoranthene	40.000	47.430	-18.6	107	0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF110716\
Data File : BF091050.D
Acq On : 7 Nov 2016 12:43
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 07 17:42:21 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 04 11:03:12 2016
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	38.580	3.6	96	0.02
89 C	Benzo(a)pyrene	40.000	41.104	-2.8	95	0.02
90	Dibenzo(a,h)anthracene	40.000	40.304	-0.8	95	0.03
91	Benzo(g,h,i)perylene	40.000	40.407	-1.0	96	0.03

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