Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051215\
Data File : BF079180.D
Acq On : 13 May 2015 3:25
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 13 04:58:27 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 13 04:45:03 2015
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	99	0.00
2		1,4-Dioxane	40.000	40.964	-2.4	102	0.00
3		Pyridine	40.000	41.844	-4.6	110	0.00
4		n-Nitrosodimethylamine	40.000	36.313	9.2	93	0.00
5	S	2-Fluorophenol	80.000	80.654	-0.8	103	0.00
6		Aniline	40.000	39.757	0.6	100	0.00
7	S	Phenol-d6	80.000	81.056	-1.3	108	0.00
8		2-Chlorophenol	40.000	41.801	-4.5	112	0.00
9		Benzaldehyde	40.000	36.945	7.6	95	0.00
10	C	Phenol	40.000	42.108	-5.3	107	0.00
11		bis(2-Chloroethyl)ether	40.000	39.166	2.1	106	0.00
12		1,3-Dichlorobenzene	40.000	40.387	-1.0	108	0.00
13	C	1,4-Dichlorobenzene	40.000	39.708	0.7	105	0.00
14		1,2-Dichlorobenzene	40.000	39.775	0.6	104	0.00
15		Benzyl Alcohol	40.000	37.648	5.9	99	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	39.861	0.3	106	0.00
17		2-Methylphenol	40.000	40.217	-0.5	106	0.00
18		Hexachloroethane	40.000	38.754	3.1	99	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	38.856	2.9	97	0.00
20		3+4-Methylphenols	40.000	41.276	-3.2	106	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	97	0.00
22		Acetophenone	40.000	41.663	-4.2	108	0.00
23	S	Nitrobenzene-d5	80.000	81.386	-1.7	103	0.00
24		Nitrobenzene	40.000	40.499	-1.2	106	0.00
25		Isophorone	40.000	39.792	0.5	100	0.00
26	C	2-Nitrophenol	40.000	45.709	-14.3	110	0.00
27		2,4-Dimethylphenol	40.000	43.332	-8.3	107	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.843	0.4	98	0.00
29	C	2,4-Dichlorophenol	40.000	42.955	-7.4	108	0.00
30		1,2,4-Trichlorobenzene	40.000	40.563	-1.4	103	0.00
31		Naphthalene	40.000	41.217	-3.0	106	0.00
32		Benzoic acid	40.000	41.449	-3.6	102	0.00
33		4-Chloroaniline	40.000	40.980	-2.4	106	0.00
34	C	Hexachlorobutadiene	40.000	40.606	-1.5	105	0.00
35		Caprolactam	40.000	42.407	-6.0	105	0.00
36	C	4-Chloro-3-methylphenol	40.000	41.859	-4.6	99	0.00
37		2-Methylnaphthalene	40.000	40.943	-2.4	103	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	97	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	41.434	-3.6	108	0.00
40	P	Hexachlorocyclopentadiene	40.000	9.543	76.1#	24	0.00
41	S	2,4,6-Tribromophenol	80.000	86.317	-7.9	105	0.00
42	C	2,4,6-Trichlorophenol	40.000	43.326	-8.3	108	0.00
43		2,4,5-Trichlorophenol	40.000	43.488	-8.7	109	0.00
44	S	2-Fluorobiphenyl	80.000	76.917	3.9	98	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051215\
Data File : BF079180.D
Acq On : 13 May 2015 3:25
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 13 04:58:27 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 13 04:45:03 2015
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	40.810	-2.0	107	0.00
46		2-Chloronaphthalene	40.000	40.770	-1.9	108	0.00
47		2-Nitroaniline	40.000	44.000	-10.0	109	0.00
48		Acenaphthylene	40.000	39.949	0.1	104	0.00
49		Dimethylphthalate	40.000	40.640	-1.6	108	0.00
50		2,6-Dinitrotoluene	40.000	42.211	-5.5	106	0.00
51	C	Acenaphthene	40.000	38.035	4.9	97	0.00
52		3-Nitroaniline	40.000	46.710	-16.8	118	0.00
53	P	2,4-Dinitrophenol	40.000	22.927	42.7#	43	0.00
54		Dibenzofuran	40.000	40.253	-0.6	103	0.00
55	P	4-Nitrophenol	40.000	43.366	-8.4	111	0.00
56		2,4-Dinitrotoluene	40.000	42.286	-5.7	102	0.00
57		Fluorene	40.000	38.928	2.7	102	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	39.744	0.6	98	0.00
59		Diethylphthalate	40.000	38.451	3.9	99	0.00
60		4-Chlorophenyl-phenylether	40.000	40.645	-1.6	104	0.00
61		4-Nitroaniline	40.000	45.980	-14.9	113	0.00
62		Azobenzene	40.000	39.103	2.2	98	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	100	0.00
64		4,6-Dinitro-2-methylphenol	40.000	24.553	38.6#	52	0.00
65	c	n-Nitrosodiphenylamine	40.000	41.178	-2.9	105	0.00
66		4-Bromophenyl-phenylether	40.000	39.705	0.7	106	0.00
67		Hexachlorobenzene	40.000	38.546	3.6	104	0.00
68		Atrazine	40.000	38.671	3.3	103	0.00
69	C	Pentachlorophenol	40.000	38.172	4.6	100	0.00
70		Phenanthrene	40.000	39.781	0.5	104	0.00
71		Anthracene	40.000	39.505	1.2	104	0.00
72		Carbazole	40.000	38.873	2.8	101	0.00
73		Di-n-butylphthalate	40.000	39.980	0.1	100	0.00
74	C	Fluoranthene	40.000	41.377	-3.4	108	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	94	0.00
76		Benzidine	40.000	52.096	-30.2#	120	0.00
77		Pyrene	40.000	42.471	-6.2	107	0.00
78	S	Terphenyl-d14	80.000	80.243	-0.3	101	0.00
79		Butylbenzylphthalate	40.000	44.669	-11.7	104	0.00
80		Benzo(a)anthracene	40.000	40.955	-2.4	102	0.00
81		3,3'-Dichlorobenzidine	40.000	47.253	-18.1	113	0.00
82		Chrysene	40.000	38.224	4.4	98	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	40.149	-0.4	93	0.00
84	c	Di-n-octyl phthalate	40.000	42.373	-5.9	105	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	37.712	5.7	94	0.00
86	I	Perylene-d12	20.000	20.000	0.0	99	0.00
87		Benzo(b)fluoranthene	40.000	41.541	-3.9	106	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051215\
Data File : BF079180.D
Acq On : 13 May 2015 3:25
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 13 04:58:27 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 13 04:45:03 2015
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	40.050	-0.1	102	0.00
89 C	Benzo(a)pyrene	40.000	41.494	-3.7	106	0.00
90	Dibenzo(a,h)anthracene	40.000	36.967	7.6	94	0.00
91	Benzo(g,h,i)perylene	40.000	37.007	7.5	95	0.00

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