Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA B\Data\BB093016\
Data File : BB058572.D
Acq On : 30 Sep 2016 18:12
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_B
LabSampleId : SSTDCCC040

Quant Time: Sep 30 19:02:01 2016
Quant Method : Z:\HPCHEM1\BNA B\METHOD\8270-BB093016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 30 14:16:00 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	88	0.00
2		1,4-Dioxane	40.000	46.972	-17.4	105	0.00
3		Pyridine	40.000	47.552	-18.9	104	0.00
4		n-Nitrosodimethylamine	40.000	47.731	-19.3	105	0.00
5	S	2-Fluorophenol	80.000	94.814	-18.5	104	0.00
6		Aniline	40.000	46.739	-16.8	103	-0.02
7	S	Phenol-d6	80.000	94.498	-18.1	106	-0.02
8		2-Chlorophenol	40.000	47.062	-17.7	105	0.00
9		Benzaldehyde	40.000	38.526	3.7	84	0.00
10	C	Phenol	40.000	47.365	-18.4	103	-0.02
11		bis(2-Chloroethyl)ether	40.000	47.674	-19.2	105	-0.02
12		1,3-Dichlorobenzene	40.000	46.177	-15.4	101	0.00
13	C	1,4-Dichlorobenzene	40.000	46.117	-15.3	102	0.00
14		1,2-Dichlorobenzene	40.000	47.243	-18.1	103	0.00
15		Benzyl Alcohol	40.000	47.588	-19.0	105	-0.02
16		2,2'-oxybis(1-Chloropropane	40.000	49.635	-24.1	106	0.00
17		2-Methylphenol	40.000	46.787	-17.0	102	-0.02
18		Hexachloroethane	40.000	47.795	-19.5	105	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	48.002	-20.0	108	-0.04
20		3+4-Methylphenols	40.000	46.552	-16.4	104	-0.04
21	I	Naphthalene-d8	20.000	20.000	0.0	92	0.00
22		Acetophenone	40.000	44.498	-11.2	101	-0.02
23	S	Nitrobenzene-d5	80.000	91.888	-14.9	105	-0.02
24		Nitrobenzene	40.000	45.716	-14.3	103	-0.02
25		Isophorone	40.000	44.816	-12.0	105	-0.04
26	C	2-Nitrophenol	40.000	45.495	-13.7	102	0.00
27		2,4-Dimethylphenol	40.000	45.505	-13.8	106	-0.02
28		bis(2-Chloroethoxy)methane	40.000	47.233	-18.1	107	0.00
29	C	2,4-Dichlorophenol	40.000	44.897	-12.2	103	0.00
30		1,2,4-Trichlorobenzene	40.000	44.956	-12.4	103	-0.02
31		Naphthalene	40.000	44.492	-11.2	100	-0.02
32		Benzoic acid	40.000	49.095	-22.7	115	-0.09
33		4-Chloroaniline	40.000	44.856	-12.1	103	-0.02
34	C	Hexachlorobutadiene	40.000	44.118	-10.3	100	0.00
35		Caprolactam	40.000	45.832	-14.6	109	-0.11
36	C	4-Chloro-3-methylphenol	40.000	45.155	-12.9	103	-0.02
37		2-Methylnaphthalene	40.000	45.090	-12.7	99	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	104	-0.02
39		1,2,4,5-Tetrachlorobenzene	40.000	41.215	-3.0	105	0.00
40	P	Hexachlorocyclopentadiene	40.000	40.926	-2.3	100	0.00
41	S	2,4,6-Tribromophenol	80.000	82.887	-3.6	100	-0.02
42	C	2,4,6-Trichlorophenol	40.000	40.120	-0.3	102	0.00
43		2,4,5-Trichlorophenol	40.000	38.993	2.5	102	-0.02
44	S	2-Fluorobiphenyl	80.000	80.614	-0.8	94	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA B\Data\BB093016\
Data File : BB058572.D
Acq On : 30 Sep 2016 18:12
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_B
LabSampleId : SSTDCCC040

Quant Time: Sep 30 19:02:01 2016
Quant Method : Z:\HPCHEM1\BNA B\METHOD\8270-BB093016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 30 14:16:00 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	40.828	-2.1	99	-0.02
46		2-Chloronaphthalene	40.000	39.065	2.3	99	-0.02
47		2-Nitroaniline	40.000	40.235	-0.6	100	-0.02
48		Acenaphthylene	40.000	41.890	-4.7	108	0.00
49		Dimethylphthalate	40.000	41.765	-4.4	106	-0.04
50		2,6-Dinitrotoluene	40.000	41.876	-4.7	107	-0.02
51	C	Acenaphthene	40.000	39.430	1.4	97	-0.02
52		3-Nitroaniline	40.000	42.532	-6.3	111	-0.02
53	P	2,4-Dinitrophenol	40.000	40.444	-1.1	104	-0.04
54		Dibenzofuran	40.000	40.735	-1.8	103	0.00
55	P	4-Nitrophenol	40.000	40.125	-0.3	98	-0.04
56		2,4-Dinitrotoluene	40.000	39.793	0.5	104	-0.04
57		Fluorene	40.000	42.640	-6.6	110	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	38.457	3.9	100	0.00
59		Diethylphthalate	40.000	41.548	-3.9	108	-0.04
60		4-Chlorophenyl-phenylether	40.000	41.516	-3.8	102	0.00
61		4-Nitroaniline	40.000	42.532	-6.3	111	-0.02
62		Azobenzene	40.000	41.668	-4.2	111	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	93	0.00
64		4,6-Dinitro-2-methylphenol	40.000	42.127	-5.3	95	-0.04
65	c	n-Nitrosodiphenylamine	40.000	46.548	-16.4	110	-0.02
66		4-Bromophenyl-phenylether	40.000	41.364	-3.4	91	-0.02
67		Hexachlorobenzene	40.000	44.006	-10.0	108	0.00
68		Atrazine	40.000	40.414	-1.0	90	-0.02
69	C	Pentachlorophenol	40.000	43.047	-7.6	95	-0.02
70		Phenanthrene	40.000	44.492	-11.2	108	-0.02
71		Anthracene	40.000	43.200	-8.0	99	-0.02
72		Carbazole	40.000	41.085	-2.7	94	-0.02
73		Di-n-butylphthalate	40.000	40.790	-2.0	100	-0.02
74	C	Fluoranthene	40.000	41.643	-4.1	92	-0.02
75	I	Chrysene-d12	20.000	20.000	0.0	93	-0.02
76		Benzidine	40.000	38.782	3.0	92	0.00
77		Pyrene	40.000	40.830	-2.1	101	-0.02
78	S	Terphenyl-d14	80.000	77.869	2.7	91	-0.02
79		Butylbenzylphthalate	40.000	40.381	-1.0	89	-0.02
80		Benzo(a)anthracene	40.000	43.816	-9.5	103	-0.02
81		3,3'-Dichlorobenzidine	40.000	41.967	-4.9	101	-0.02
82		Chrysene	40.000	44.727	-11.8	104	-0.02
83		Bis(2-ethylhexyl)phthalate	40.000	42.058	-5.1	96	0.00
84	c	Di-n-octyl phthalate	40.000	42.320	-5.8	98	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	40.492	-1.2	95	-0.10
86	I	Perylene-d12	20.000	20.000	0.0	91	-0.02
87		Benzo(b)fluoranthene	40.000	40.914	-2.3	93	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA B\Data\BB093016\
Data File : BB058572.D
Acq On : 30 Sep 2016 18:12
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_B
LabSampleId : SSTDCCC040

Quant Time: Sep 30 19:02:01 2016
Quant Method : Z:\HPCHEM1\BNA B\METHOD\8270-BB093016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 30 14:16:00 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	48.560	-21.4	106	-0.04
89 C	Benzo(a)pyrene	40.000	45.022	-12.6	99	-0.02
90	Dibenzo(a,h)anthracene	40.000	43.306	-8.3	95	-0.06
91	Benzo(g,h,i)perylene	40.000	42.511	-6.3	94	-0.06

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