Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF100217\
Data File : BF098998.D
Acq On : 2 Oct 2017 14:29
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 03 05:17:24 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Sep 23 13:50:58 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	80	-0.02
2		1,4-Dioxane	40.000	40.772	-1.9	80	-0.06
3		Pyridine	40.000	40.493	-1.2	82	-0.05
4		n-Nitrosodimethylamine	40.000	39.880	0.3	79	-0.05
5	S	2-Fluorophenol	80.000	83.994	-5.0	83	-0.02
6		Aniline	40.000	41.264	-3.2	83	-0.02
7	S	Phenol-d6	80.000	84.678	-5.8	84	-0.02
8		2-Chlorophenol	40.000	41.653	-4.1	84	-0.02
9		Benzaldehyde	40.000	37.709	5.7	77	-0.02
10	C	Phenol	40.000	42.642	-6.6	86	-0.01
11		bis(2-Chloroethyl)ether	40.000	41.594	-4.0	84	-0.02
12		1,3-Dichlorobenzene	40.000	41.252	-3.1	83	-0.02
13	C	1,4-Dichlorobenzene	40.000	41.777	-4.4	85	-0.02
14		1,2-Dichlorobenzene	40.000	41.826	-4.6	84	-0.02
15		Benzyl Alcohol	40.000	41.134	-2.8	82	-0.02
16		2,2'-oxybis(1-Chloropropane	40.000	39.949	0.1	82	-0.02
17		2-Methylphenol	40.000	41.084	-2.7	83	-0.01
18		Hexachloroethane	40.000	41.356	-3.4	84	-0.02
19	P	n-Nitroso-di-n-propylamine	40.000	39.292	1.8	82	-0.02
20		3+4-Methylphenols	40.000	40.338	-0.8	81	-0.02
21	I	Naphthalene-d8	20.000	20.000	0.0	79	-0.02
22		Acetophenone	40.000	39.610	1.0	81	-0.02
23	S	Nitrobenzene-d5	80.000	84.591	-5.7	83	-0.02
24		Nitrobenzene	40.000	41.861	-4.7	83	-0.02
25		Isophorone	40.000	40.631	-1.6	83	-0.02
26	C	2-Nitrophenol	40.000	45.384	-13.5	87	-0.02
27		2,4-Dimethylphenol	40.000	40.823	-2.1	83	-0.02
28		bis(2-Chloroethoxy)methane	40.000	41.487	-3.7	85	-0.02
29	C	2,4-Dichlorophenol	40.000	42.051	-5.1	84	-0.02
30		1,2,4-Trichlorobenzene	40.000	41.864	-4.7	84	-0.02
31		Naphthalene	40.000	41.120	-2.8	84	-0.02
32		Benzoic acid	40.000	35.455	11.4	68	-0.01
33		4-Chloroaniline	40.000	42.128	-5.3	85	-0.02
34	C	Hexachlorobutadiene	40.000	41.666	-4.2	85	-0.02
35		Caprolactam	40.000	41.403	-3.5	85	-0.01
36	C	4-Chloro-3-methylphenol	40.000	40.863	-2.2	82	-0.01
37		2-Methylnaphthalene	40.000	41.756	-4.4	85	-0.02
38	I	Acenaphthene-d10	20.000	20.000	0.0	79	-0.02
39		1,2,4,5-Tetrachlorobenzene	40.000	42.779	-6.9	86	-0.02
40	P	Hexachlorocyclopentadiene	40.000	34.686	13.3	67	-0.02
41	S	2,4,6-Tribromophenol	80.000	85.621	-7.0	82	-0.02
42	C	2,4,6-Trichlorophenol	40.000	41.703	-4.3	83	-0.02
43		2,4,5-Trichlorophenol	40.000	42.594	-6.5	83	-0.01
44	S	2-Fluorobiphenyl	80.000	85.631	-7.0	84	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF100217\
Data File : BF098998.D
Acq On : 2 Oct 2017 14:29
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 03 05:17:24 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Sep 23 13:50:58 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	41.851	-4.6	84	-0.02
46		2-Chloronaphthalene	40.000	41.855	-4.6	84	-0.02
47		2-Nitroaniline	40.000	42.330	-5.8	82	-0.01
48		Acenaphthylene	40.000	41.700	-4.3	84	-0.02
49		Dimethylphthalate	40.000	41.871	-4.7	85	-0.02
50		2,6-Dinitrotoluene	40.000	42.907	-7.3	83	-0.01
51	C	Acenaphthene	40.000	41.780	-4.5	84	-0.02
52		3-Nitroaniline	40.000	43.516	-8.8	83	-0.01
53	P	2,4-Dinitrophenol	40.000	39.660	0.9	80	0.00
54		Dibenzofuran	40.000	41.954	-4.9	84	-0.02
55	P	4-Nitrophenol	40.000	40.209	-0.5	77	0.00
56		2,4-Dinitrotoluene	40.000	44.618	-11.5	85	-0.01
57		Fluorene	40.000	42.032	-5.1	85	-0.02
58		2,3,4,6-Tetrachlorophenol	40.000	41.455	-3.6	82	-0.02
59		Diethylphthalate	40.000	41.246	-3.1	83	-0.02
60		4-Chlorophenyl-phenylether	40.000	41.175	-2.9	83	-0.02
61		4-Nitroaniline	40.000	43.575	-8.9	84	-0.01
62		Azobenzene	40.000	40.589	-1.5	82	-0.02
63	I	Phenanthrene-d10	20.000	20.000	0.0	77	-0.01
64		4,6-Dinitro-2-methylphenol	40.000	44.110	-10.3	84	-0.01
65	c	n-Nitrosodiphenylamine	40.000	41.643	-4.1	83	-0.01
66		4-Bromophenyl-phenylether	40.000	42.325	-5.8	84	-0.02
67		Hexachlorobenzene	40.000	42.753	-6.9	85	-0.02
68		Atrazine	40.000	41.897	-4.7	82	-0.01
69	C	Pentachlorophenol	40.000	38.246	4.4	72	-0.01
70		Phenanthrene	40.000	41.525	-3.8	84	-0.02
71		Anthracene	40.000	41.672	-4.2	83	-0.02
72		Carbazole	40.000	41.449	-3.6	82	-0.01
73		Di-n-butylphthalate	40.000	41.590	-4.0	82	-0.02
74	C	Fluoranthene	40.000	40.952	-2.4	82	-0.02
75	I	Chrysene-d12	20.000	20.000	0.0	76	-0.01
76		Benzidine	40.000	41.209	-3.0	79	-0.02
77		Pyrene	40.000	42.727	-6.8	82	-0.02
78	S	Terphenyl-d14	80.000	86.284	-7.9	83	-0.02
79		Butylbenzylphthalate	40.000	44.057	-10.1	82	-0.02
80		Benzo(a)anthracene	40.000	41.794	-4.5	81	-0.01
81		3,3'-Dichlorobenzidine	40.000	40.662	-1.7	77	-0.02
82		Chrysene	40.000	41.602	-4.0	81	-0.01
83		Bis(2-ethylhexyl)phthalate	40.000	43.554	-8.9	82	-0.02
84	c	Di-n-octyl phthalate	40.000	43.617	-9.0	85	-0.02
85		Indeno(1,2,3-cd)pyrene	40.000	45.016	-12.5	94	-0.02
86	I	Perylene-d12	20.000	20.000	0.0	80	-0.01
87		Benzo(b)fluoranthene	40.000	43.319	-8.3	85	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF100217\
Data File : BF098998.D
Acq On : 2 Oct 2017 14:29
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 03 05:17:24 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Sep 23 13:50:58 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	39.120	2.2	79	-0.01
89 C	Benzo(a)pyrene	40.000	41.716	-4.3	84	-0.02
90	Dibenzo(a,h)anthracene	40.000	44.899	-12.2	93	-0.02
91	Benzo(g,h,i)perylene	40.000	45.941	-14.9	95	-0.02

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