Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA F\Data\BF110517\
Data File : BF100216.D
Acq On : 5 Nov 2017 11:41
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 06 06:10:18 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	90	0.00
2		1,4-Dioxane	40.000	37.938	5.2	84	-0.02
3		Pyridine	40.000	36.581	8.5	75	-0.01
4		n-Nitrosodimethylamine	40.000	38.099	4.8	78	-0.03
5	S	2-Fluorophenol	80.000	82.771	-3.5	90	0.00
6		Aniline	40.000	39.827	0.4	88	0.00
7	S	Phenol-d6	80.000	80.859	-1.1	90	0.00
8		2-Chlorophenol	40.000	40.743	-1.9	88	0.00
9		Benzaldehyde	40.000	35.712	10.7	79	0.00
10	C	Phenol	40.000	38.197	4.5	83	0.00
11		bis(2-Chloroethyl)ether	40.000	39.243	1.9	85	0.00
12		1,3-Dichlorobenzene	40.000	39.376	1.6	87	0.00
13	C	1,4-Dichlorobenzene	40.000	40.467	-1.2	89	0.00
14		1,2-Dichlorobenzene	40.000	39.917	0.2	88	0.00
15		Benzyl Alcohol	40.000	42.672	-6.7	96	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	37.281	6.8	82	-0.01
17		2-Methylphenol	40.000	41.178	-2.9	90	0.00
18		Hexachloroethane	40.000	38.221	4.4	85	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	41.214	-3.0	93	0.00
20		3+4-Methylphenols	40.000	45.828	-14.6	103	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	92	0.00
22		Acetophenone	40.000	40.598	-1.5	95	0.00
23	S	Nitrobenzene-d5	80.000	75.477	5.7	87	-0.01
24		Nitrobenzene	40.000	38.595	3.5	86	0.00
25		Isophorone	40.000	38.312	4.2	87	0.00
26	C	2-Nitrophenol	40.000	39.029	2.4	85	0.00
27		2,4-Dimethylphenol	40.000	40.392	-1.0	92	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.625	3.4	89	0.00
29	C	2,4-Dichlorophenol	40.000	41.256	-3.1	93	0.00
30		1,2,4-Trichlorobenzene	40.000	38.322	4.2	87	0.00
31		Naphthalene	40.000	39.237	1.9	90	0.00
32		Benzoic acid	40.000	41.660	-4.1	96	0.00
33		4-Chloroaniline	40.000	40.488	-1.2	91	0.00
34	C	Hexachlorobutadiene	40.000	38.624	3.4	89	-0.01
35		Caprolactam	40.000	41.711	-4.3	93	-0.02
36	C	4-Chloro-3-methylphenol	40.000	40.515	-1.3	92	0.00
37		2-Methylnaphthalene	40.000	39.734	0.7	92	-0.01
38	I	Acenaphthene-d10	20.000	20.000	0.0	93	-0.01
39		1,2,4,5-Tetrachlorobenzene	40.000	39.426	1.4	92	0.00
40	P	Hexachlorocyclopentadiene	40.000	49.888	-24.7	134	-0.01
41	S	2,4,6-Tribromophenol	80.000	77.942	2.6	91	-0.01
42	C	2,4,6-Trichlorophenol	40.000	38.172	4.6	89	0.00
43		2,4,5-Trichlorophenol	40.000	34.144	14.6	76	0.00
44	S	2-Fluorobiphenyl	80.000	75.776	5.3	87	-0.01

Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA F\Data\BF110517\
Data File : BF100216.D
Acq On : 5 Nov 2017 11:41
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 06 06:10:18 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	38.584	3.5	91	-0.01
46		2-Chloronaphthalene	40.000	38.376	4.1	89	-0.01
47		2-Nitroaniline	40.000	39.465	1.3	90	0.00
48		Acenaphthylene	40.000	38.910	2.7	92	-0.01
49		Dimethylphthalate	40.000	36.815	8.0	86	0.00
50		2,6-Dinitrotoluene	40.000	38.568	3.6	88	-0.01
51	C	Acenaphthene	40.000	38.575	3.6	91	-0.01
52		3-Nitroaniline	40.000	40.409	-1.0	93	0.00
53	P	2,4-Dinitrophenol	40.000	35.342	11.6	81	0.01
54		Dibenzofuran	40.000	40.318	-0.8	96	0.00
55	P	4-Nitrophenol	40.000	38.483	3.8	84	0.00
56		2,4-Dinitrotoluene	40.000	44.770	-11.9	103	0.00
57		Fluorene	40.000	39.459	1.4	93	-0.01
58		2,3,4,6-Tetrachlorophenol	40.000	40.188	-0.5	91	0.00
59		Diethylphthalate	40.000	37.554	6.1	88	-0.01
60		4-Chlorophenyl-phenylether	40.000	39.616	1.0	94	-0.01
61		4-Nitroaniline	40.000	40.787	-2.0	92	0.00
62		Azobenzene	40.000	37.160	7.1	85	-0.01
63	I	Phenanthrene-d10	20.000	20.000	0.0	90	0.00
64		4,6-Dinitro-2-methylphenol	40.000	36.549	8.6	79	0.00
65	c	n-Nitrosodiphenylamine	40.000	40.048	-0.1	92	-0.01
66		4-Bromophenyl-phenylether	40.000	40.370	-0.9	92	0.00
67		Hexachlorobenzene	40.000	39.882	0.3	90	0.00
68		Atrazine	40.000	38.561	3.6	85	-0.01
69	C	Pentachlorophenol	40.000	34.435	13.9	78	0.00
70		Phenanthrene	40.000	39.436	1.4	90	-0.01
71		Anthracene	40.000	40.086	-0.2	92	-0.01
72		Carbazole	40.000	39.796	0.5	90	0.00
73		Di-n-butylphthalate	40.000	37.756	5.6	85	-0.01
74	C	Fluoranthene	40.000	37.324	6.7	85	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	71	-0.01
76		Benzidine	40.000	30.254	24.4	55	-0.01
77		Pyrene	40.000	46.659	-16.6	83	-0.01
78	S	Terphenyl-d14	80.000	91.731	-14.7	83	-0.01
79		Butylbenzylphthalate	40.000	42.361	-5.9	73	-0.02
80		Benzo(a)anthracene	40.000	37.848	5.4	67	-0.01
81		3,3'-Dichlorobenzidine	40.000	35.259	11.9	62	-0.01
82		Chrysene	40.000	38.857	2.9	69	-0.01
83		Bis(2-ethylhexyl)phthalate	40.000	37.754	5.6	68	-0.01
84	c	Di-n-octyl phthalate	40.000	32.870	17.8	57	-0.02
85		Indeno(1,2,3-cd)pyrene	40.000	38.612	3.5	72	-0.02
86	I	Perylene-d12	20.000	20.000	0.0	61	0.00
87		Benzo(b)fluoranthene	40.000	35.394	11.5	56	-0.01

Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA F\Data\BF110517\
Data File : BF100216.D
Acq On : 5 Nov 2017 11:41
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 06 06:10:18 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	43.268	-8.2	62	-0.01
89 C	Benzo(a)pyrene	40.000	39.362	1.6	59	-0.01
90	Dibenzo(a,h)anthracene	40.000	44.916	-12.3	70	-0.02
91	Benzo(g,h,i)perylene	40.000	46.371	-15.9	72	-0.02

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