Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF073018\
Data File : BF107801.D
Acq On : 30 Jul 2018 15:49
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 30 18:21:32 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 30 17:48:53 2018
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	0.00
2		1,4-Dioxane	40.000	39.667	0.8	101	0.00
3		Pyridine	40.000	36.668	8.3	96	0.00
4		n-Nitrosodimethylamine	40.000	34.690	13.3	94	0.00
5	S	2-Fluorophenol	80.000	70.873	11.4	85	0.00
6		Aniline	40.000	38.873	2.8	94	0.00
7	S	Phenol-d6	80.000	73.159	8.6	98	0.00
8		2-Chlorophenol	40.000	37.351	6.6	98	0.00
9		Benzaldehyde	40.000	36.736	8.2	96	0.00
10	C	Phenol	40.000	32.033	19.9	96	0.00
11		bis(2-Chloroethyl)ether	40.000	39.548	1.1	105	0.00
12		1,3-Dichlorobenzene	40.000	37.788	5.5	97	0.00
13	C	1,4-Dichlorobenzene	40.000	38.892	2.8	98	0.00
14		1,2-Dichlorobenzene	40.000	37.273	6.8	99	0.00
15		Benzyl Alcohol	40.000	39.025	2.4	96	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	36.757	8.1	96	0.00
17		2-Methylphenol	40.000	38.541	3.6	96	0.00
18		Hexachloroethane	40.000	36.338	9.2	97	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.007	2.5	98	0.00
20		3+4-Methylphenols	40.000	41.104	-2.8	97	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	103	0.00
22		Acetophenone	40.000	36.510	8.7	98	0.00
23	S	Nitrobenzene-d5	80.000	75.858	5.2	101	0.00
24		Nitrobenzene	40.000	37.347	6.6	98	0.00
25		Isophorone	40.000	37.298	6.8	95	0.00
26	C	2-Nitrophenol	40.000	39.443	1.4	97	0.00
27		2,4-Dimethylphenol	40.000	37.811	5.5	99	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.374	1.6	97	0.00
29	C	2,4-Dichlorophenol	40.000	39.453	1.4	98	0.00
30		1,2,4-Trichlorobenzene	40.000	38.637	3.4	98	0.00
31		Naphthalene	40.000	39.228	1.9	96	0.00
32		Benzoic acid	40.000	32.250	19.4	76	0.00
33		4-Chloroaniline	40.000	40.169	-0.4	102	0.00
34	C	Hexachlorobutadiene	40.000	37.427	6.4	98	0.00
35		Caprolactam	40.000	39.075	2.3	99	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.029	-0.1	98	0.00
37		2-Methylnaphthalene	40.000	40.317	-0.8	98	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	104	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	38.920	2.7	97	0.00
40	P	Hexachlorocyclopentadiene	40.000	37.031	7.4	96	0.00
41	S	2,4,6-Tribromophenol	80.000	78.049	2.4	99	0.00
42	C	2,4,6-Trichlorophenol	40.000	40.762	-1.9	98	0.00
43		2,4,5-Trichlorophenol	40.000	40.235	-0.6	97	0.00
44	S	2-Fluorobiphenyl	80.000	77.154	3.6	102	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF073018\
Data File : BF107801.D
Acq On : 30 Jul 2018 15:49
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 30 18:21:32 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 30 17:48:53 2018
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	38.736	3.2	98	0.00
46		2-Chloronaphthalene	40.000	38.703	3.2	98	0.00
47		2-Nitroaniline	40.000	39.213	2.0	98	0.00
48		Acenaphthylene	40.000	37.237	6.9	99	0.00
49		Dimethylphthalate	40.000	39.012	2.5	100	0.00
50		2,6-Dinitrotoluene	40.000	39.632	0.9	102	0.00
51	C	Acenaphthene	40.000	36.532	8.7	100	0.00
52		3-Nitroaniline	40.000	38.538	3.7	98	0.00
53	P	2,4-Dinitrophenol	40.000	37.124	7.2	95	0.00
54		Dibenzofuran	40.000	37.578	6.1	98	0.00
55	P	4-Nitrophenol	40.000	36.203	9.5	89	0.00
56		2,4-Dinitrotoluene	40.000	37.663	5.8	95	0.00
57		Fluorene	40.000	37.369	6.6	99	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	38.625	3.4	93	0.00
59		Diethylphthalate	40.000	37.739	5.7	100	0.00
60		4-Chlorophenyl-phenylether	40.000	37.693	5.8	100	0.00
61		4-Nitroaniline	40.000	39.022	2.4	101	0.00
62		Azobenzene	40.000	37.407	6.5	99	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	105	0.00
64		4,6-Dinitro-2-methylphenol	40.000	40.462	-1.2	102	0.00
65	c	n-Nitrosodiphenylamine	40.000	37.399	6.5	100	0.00
66		4-Bromophenyl-phenylether	40.000	36.095	9.8	98	0.00
67		Hexachlorobenzene	40.000	37.526	6.2	99	0.00
68		Atrazine	40.000	38.886	2.8	101	0.00
69	C	Pentachlorophenol	40.000	40.094	-0.2	99	0.00
70		Phenanthrene	40.000	37.351	6.6	98	0.00
71		Anthracene	40.000	36.668	8.3	98	0.00
72		Carbazole	40.000	38.128	4.7	100	0.00
73		Di-n-butylphthalate	40.000	36.889	7.8	98	0.00
74	C	Fluoranthene	40.000	36.803	8.0	98	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	104	0.00
76		Benzidine	40.000	39.178	2.1	98	0.00
77		Pyrene	40.000	37.465	6.3	100	0.00
78	S	Terphenyl-d14	80.000	75.326	5.8	102	0.00
79		Butylbenzylphthalate	40.000	37.953	5.1	98	0.00
80		Benzo(a)anthracene	40.000	36.991	7.5	96	0.00
81		3,3'-Dichlorobenzidine	40.000	37.192	7.0	98	0.00
82		Chrysene	40.000	37.655	5.9	101	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	37.425	6.4	99	0.00
84	c	Di-n-octyl phthalate	40.000	38.185	4.5	99	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	33.395	16.5	93	0.00
86	I	Perylene-d12	20.000	20.000	0.0	106	0.00
87		Benzo(b)fluoranthene	40.000	38.785	3.0	103	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF073018\
Data File : BF107801.D
Acq On : 30 Jul 2018 15:49
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 30 18:21:32 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 30 17:48:53 2018
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	36.778	8.1	95	0.00
89 C	Benzo(a)pyrene	40.000	36.454	8.9	99	0.00
90	Dibenzo(a,h)anthracene	40.000	34.696	13.3	96	0.00
91	Benzo(g,h,i)perylene	40.000	33.759	15.6	93	-0.01

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