Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF041918\
Data File : BF104618.D
Acq On : 19 Apr 2018 9:21
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 19 14:52:17 2018
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 19 14:44:03 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	133	0.00
2		1,4-Dioxane	40.000	40.237	-0.6	132	0.00
3		Pyridine	40.000	39.873	0.3	132	0.00
4		n-Nitrosodimethylamine	40.000	37.960	5.1	125	0.00
5	S	2-Fluorophenol	80.000	74.942	6.3	126	0.00
6		Aniline	40.000	37.405	6.5	123	0.00
7	S	Phenol-d6	80.000	76.199	4.8	128	0.00
8		2-Chlorophenol	40.000	39.479	1.3	131	0.00
9		Benzaldehyde	40.000	36.663	8.3	122	0.00
10	C	Phenol	40.000	38.787	3.0	131	0.00
11		bis(2-Chloroethyl)ether	40.000	39.165	2.1	130	0.00
12		1,3-Dichlorobenzene	40.000	38.746	3.1	130	0.00
13	C	1,4-Dichlorobenzene	40.000	38.796	3.0	129	0.00
14		1,2-Dichlorobenzene	40.000	38.499	3.8	127	0.00
15		Benzyl Alcohol	40.000	36.425	8.9	122	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	36.865	7.8	123	0.00
17		2-Methylphenol	40.000	39.703	0.7	132	0.00
18		Hexachloroethane	40.000	39.723	0.7	133	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	34.388	14.0	120	0.00
20		3+4-Methylphenols	40.000	35.068	12.3	118	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	133	0.00
22		Acetophenone	40.000	35.928	10.2	121	0.00
23	S	Nitrobenzene-d5	80.000	88.981	-11.2	144	0.00
24		Nitrobenzene	40.000	41.999	-5.0	134	0.00
25		Isophorone	40.000	38.864	2.8	130	0.00
26	C	2-Nitrophenol	40.000	54.503	-36.3#	172	0.00
27		2,4-Dimethylphenol	40.000	37.195	7.0	122	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.721	3.2	130	0.00
29	C	2,4-Dichlorophenol	40.000	39.386	1.5	130	0.00
30		1,2,4-Trichlorobenzene	40.000	39.075	2.3	131	0.00
31		Naphthalene	40.000	37.601	6.0	126	0.00
32		Benzoic acid	40.000	39.468	1.3	123	0.00
33		4-Chloroaniline	40.000	40.071	-0.2	134	0.00
34	C	Hexachlorobutadiene	40.000	38.749	3.1	130	0.00
35		Caprolactam	40.000	41.436	-3.6	135	0.00
36	C	4-Chloro-3-methylphenol	40.000	39.829	0.4	132	0.00
37		2-Methylnaphthalene	40.000	38.137	4.7	128	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	130	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	40.333	-0.8	134	0.00
40	P	Hexachlorocyclopentadiene	40.000	35.409	11.5	113	0.00
41	S	2,4,6-Tribromophenol	80.000	82.389	-3.0	134	0.00
42	C	2,4,6-Trichlorophenol	40.000	40.562	-1.4	129	0.00
43		2,4,5-Trichlorophenol	40.000	40.903	-2.3	133	0.00
44	S	2-Fluorobiphenyl	80.000	72.540	9.3	120	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF041918\
Data File : BF104618.D
Acq On : 19 Apr 2018 9:21
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 19 14:52:17 2018
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 19 14:44:03 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	38.283	4.3	126	0.00
46		2-Chloronaphthalene	40.000	38.850	2.9	128	0.00
47		2-Nitroaniline	40.000	51.205	-28.0#	156	0.00
48		Acenaphthylene	40.000	37.495	6.3	122	0.00
49		Dimethylphthalate	40.000	39.574	1.1	131	0.00
50		2,6-Dinitrotoluene	40.000	47.512	-18.8	143	0.00
51	C	Acenaphthene	40.000	35.975	10.1	119	0.00
52		3-Nitroaniline	40.000	48.534	-21.3	147	0.00
53	P	2,4-Dinitrophenol	40.000	60.290	-50.7#	238	0.00
54		Dibenzofuran	40.000	36.399	9.0	119	0.00
55	P	4-Nitrophenol	40.000	44.578	-11.4	134	0.00
56		2,4-Dinitrotoluene	40.000	45.890	-14.7	143	0.00
57		Fluorene	40.000	39.333	1.7	128	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	39.855	0.4	130	0.00
59		Diethylphthalate	40.000	37.366	6.6	124	0.00
60		4-Chlorophenyl-phenylether	40.000	40.021	-0.1	130	0.00
61		4-Nitroaniline	40.000	50.816	-27.0#	151	0.00
62		Azobenzene	40.000	36.210	9.5	119	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	130	0.00
64		4,6-Dinitro-2-methylphenol	40.000	54.660	-36.6#	192	0.00
65	c	n-Nitrosodiphenylamine	40.000	37.647	5.9	124	0.00
66		4-Bromophenyl-phenylether	40.000	38.746	3.1	128	0.00
67		Hexachlorobenzene	40.000	39.268	1.8	130	0.00
68		Atrazine	40.000	38.938	2.7	129	0.00
69	C	Pentachlorophenol	40.000	37.194	7.0	115	0.00
70		Phenanthrene	40.000	36.522	8.7	121	0.00
71		Anthracene	40.000	36.740	8.1	122	0.00
72		Carbazole	40.000	37.604	6.0	126	0.00
73		Di-n-butylphthalate	40.000	35.246	11.9	117	0.00
74	C	Fluoranthene	40.000	36.875	7.8	122	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	120	0.00
76		Benzidine	40.000	42.429	-6.1	124	0.00
77		Pyrene	40.000	39.984	0.0	122	0.00
78	S	Terphenyl-d14	80.000	74.735	6.6	117	0.00
79		Butylbenzylphthalate	40.000	39.596	1.0	119	0.00
80		Benzo(a)anthracene	40.000	41.613	-4.0	129	0.00
81		3,3'-Dichlorobenzidine	40.000	39.584	1.0	116	0.00
82		Chrysene	40.000	38.411	4.0	117	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	39.435	1.4	121	0.00
84	c	Di-n-octyl phthalate	40.000	36.244	9.4	107	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	40.014	-0.0	126	0.00
86	I	Perylene-d12	20.000	20.000	0.0	112	0.00
87		Benzo(b)fluoranthene	40.000	39.939	0.2	112	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF041918\
Data File : BF104618.D
Acq On : 19 Apr 2018 9:21
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 19 14:52:17 2018
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 19 14:44:03 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	38.578	3.6	110	0.00
89 C	Benzo(a)pyrene	40.000	39.329	1.7	111	0.00
90	Dibenzo(a,h)anthracene	40.000	42.058	-5.1	122	0.00
91	Benzo(g,h,i)perylene	40.000	43.364	-8.4	130	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 1