Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG041218\
Data File : BG033774.D
Acq On : 12 Apr 2018 16:08
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 13 02:52:51 2018
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG031918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 10 16:51:36 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	130	0.00
2		1,4-Dioxane	40.000	39.383	1.5	129	0.00
3		Pyridine	40.000	42.200	-5.5	124	0.00
4		n-Nitrosodimethylamine	40.000	43.308	-8.3	139	0.00
5	S	2-Fluorophenol	80.000	83.683	-4.6	125	0.00
6		Aniline	40.000	42.363	-5.9	127	0.00
7	S	Phenol-d6	80.000	83.859	-4.8	132	0.00
8		2-Chlorophenol	40.000	40.793	-2.0	122	0.00
9		Benzaldehyde	40.000	33.299	16.8	110	0.00
10	C	Phenol	40.000	41.819	-4.5	128	0.00
11		bis(2-Chloroethyl)ether	40.000	40.411	-1.0	120	0.00
12		1,3-Dichlorobenzene	40.000	39.847	0.4	127	0.00
13	C	1,4-Dichlorobenzene	40.000	37.641	5.9	122	0.00
14		1,2-Dichlorobenzene	40.000	39.903	0.2	123	0.00
15		Benzyl Alcohol	40.000	42.334	-5.8	128	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	42.135	-5.3	134	0.00
17		2-Methylphenol	40.000	40.945	-2.4	129	0.00
18		Hexachloroethane	40.000	40.690	-1.7	131	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	40.484	-1.2	120	0.00
20		3+4-Methylphenols	40.000	41.803	-4.5	126	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	125	0.00
22		Acetophenone	40.000	40.304	-0.8	117	0.00
23	S	Nitrobenzene-d5	80.000	80.498	-0.6	122	0.00
24		Nitrobenzene	40.000	39.934	0.2	120	0.00
25		Isophorone	40.000	40.138	-0.3	121	0.00
26	C	2-Nitrophenol	40.000	41.864	-4.7	121	0.00
27		2,4-Dimethylphenol	40.000	42.276	-5.7	134	0.00
28		bis(2-Chloroethoxy)methane	40.000	37.586	6.0	113	0.00
29	C	2,4-Dichlorophenol	40.000	39.413	1.5	117	0.00
30		1,2,4-Trichlorobenzene	40.000	39.058	2.4	120	0.00
31		Naphthalene	40.000	38.503	3.7	118	0.00
32		Benzoic acid	40.000	33.153	17.1	116	0.01
33		4-Chloroaniline	40.000	37.653	5.9	112	0.00
34	C	Hexachlorobutadiene	40.000	37.395	6.5	119	0.00
35		Caprolactam	40.000	40.591	-1.5	122	0.01
36	C	4-Chloro-3-methylphenol	40.000	39.452	1.4	115	0.00
37		2-Methylnaphthalene	40.000	38.118	4.7	120	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	118	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	38.372	4.1	113	0.00
40	P	Hexachlorocyclopentadiene	40.000	42.172	-5.4	121	0.00
41	S	2,4,6-Tribromophenol	80.000	73.354	8.3	105	0.00
42	C	2,4,6-Trichlorophenol	40.000	39.606	1.0	111	0.00
43		2,4,5-Trichlorophenol	40.000	40.334	-0.8	109	0.00
44	S	2-Fluorobiphenyl	80.000	76.456	4.4	111	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG041218\
Data File : BG033774.D
Acq On : 12 Apr 2018 16:08
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 13 02:52:51 2018
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG031918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 10 16:51:36 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	39.322	1.7	114	0.00
46		2-Chloronaphthalene	40.000	39.451	1.4	114	0.00
47		2-Nitroaniline	40.000	43.107	-7.8	121	0.00
48		Acenaphthylene	40.000	39.005	2.5	114	0.00
49		Dimethylphthalate	40.000	38.911	2.7	113	0.00
50		2,6-Dinitrotoluene	40.000	40.878	-2.2	114	0.00
51	C	Acenaphthene	40.000	40.625	-1.6	116	0.00
52		3-Nitroaniline	40.000	39.818	0.5	114	0.00
53	P	2,4-Dinitrophenol	40.000	42.690	-6.7	129	0.00
54		Dibenzofuran	40.000	38.197	4.5	112	0.00
55	P	4-Nitrophenol	40.000	37.329	6.7	105	0.00
56		2,4-Dinitrotoluene	40.000	39.162	2.1	113	0.00
57		Fluorene	40.000	38.437	3.9	114	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	37.346	6.6	105	0.00
59		Diethylphthalate	40.000	39.140	2.1	115	0.00
60		4-Chlorophenyl-phenylether	40.000	37.561	6.1	112	0.00
61		4-Nitroaniline	40.000	40.665	-1.7	117	0.00
62		Azobenzene	40.000	39.930	0.2	116	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	120	0.00
64		4,6-Dinitro-2-methylphenol	40.000	38.977	2.6	110	0.00
65	c	n-Nitrosodiphenylamine	40.000	38.725	3.2	114	0.00
66		4-Bromophenyl-phenylether	40.000	36.695	8.3	107	0.00
67		Hexachlorobenzene	40.000	37.100	7.2	110	0.00
68		Atrazine	40.000	37.967	5.1	111	0.00
69	C	Pentachlorophenol	40.000	32.627	18.4	96	0.00
70		Phenanthrene	40.000	37.813	5.5	112	0.00
71		Anthracene	40.000	38.527	3.7	114	0.00
72		Carbazole	40.000	39.575	1.1	116	0.00
73		Di-n-butylphthalate	40.000	40.510	-1.3	118	0.00
74	C	Fluoranthene	40.000	37.793	5.5	112	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	122	0.00
76		Benzidine	40.000	35.490	11.3	98	0.00
77		Pyrene	40.000	38.001	5.0	114	0.00
78	S	Terphenyl-d14	80.000	75.198	6.0	114	0.00
79		Butylbenzylphthalate	40.000	40.650	-1.6	121	0.00
80		Benzo(a)anthracene	40.000	38.077	4.8	115	0.00
81		3,3'-Dichlorobenzidine	40.000	40.318	-0.8	116	0.00
82		Chrysene	40.000	37.751	5.6	114	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	41.474	-3.7	123	0.00
84	c	Di-n-octyl phthalate	40.000	43.476	-8.7	127	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	41.745	-4.4	124	0.02
86	I	Perylene-d12	20.000	20.000	0.0	127	0.00
87		Benzo(b)fluoranthene	40.000	36.739	8.2	115	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG041218\
Data File : BG033774.D
Acq On : 12 Apr 2018 16:08
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 13 02:52:51 2018
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG031918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 10 16:51:36 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	38.796	3.0	121	0.00
89 C	Benzo(a)pyrene	40.000	38.430	3.9	120	0.00
90	Dibenzo(a,h)anthracene	40.000	40.075	-0.2	121	0.00
91	Benzo(g,h,i)perylene	40.000	39.742	0.6	123	0.02

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