Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG101918\
Data File : BG037464.D
Acq On : 19 Oct 2018 23:46
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 20 00:23:26 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG101818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 18 14:06:42 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	114	0.00
2		1,4-Dioxane	40.000	38.613	3.5	113	0.00
3		Pyridine	40.000	40.431	-1.1	117	0.00
4		n-Nitrosodimethylamine	40.000	40.402	-1.0	118	0.00
5	S	2-Fluorophenol	80.000	79.006	1.2	113	0.00
6		Aniline	40.000	38.947	2.6	112	0.00
7	S	Phenol-d6	80.000	77.620	3.0	110	0.00
8		2-Chlorophenol	40.000	38.189	4.5	110	0.00
9		Benzaldehyde	40.000	37.629	5.9	108	0.00
10	C	Phenol	40.000	38.688	3.3	111	0.00
11		bis(2-Chloroethyl)ether	40.000	39.357	1.6	114	0.00
12		1,3-Dichlorobenzene	40.000	38.786	3.0	113	0.00
13	C	1,4-Dichlorobenzene	40.000	37.798	5.5	110	0.00
14		1,2-Dichlorobenzene	40.000	37.626	5.9	109	0.00
15		Benzyl Alcohol	40.000	38.989	2.5	109	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	40.011	-0.0	116	0.00
17		2-Methylphenol	40.000	39.027	2.4	110	0.00
18		Hexachloroethane	40.000	39.628	0.9	114	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	38.936	2.7	109	0.00
20		3+4-Methylphenols	40.000	38.950	2.6	111	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	112	0.00
22		Acetophenone	40.000	38.862	2.8	109	0.00
23	S	Nitrobenzene-d5	80.000	80.928	-1.2	113	0.00
24		Nitrobenzene	40.000	40.596	-1.5	112	0.00
25		Isophorone	40.000	40.438	-1.1	110	0.00
26	C	2-Nitrophenol	40.000	42.329	-5.8	115	0.00
27		2,4-Dimethylphenol	40.000	38.717	3.2	107	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.513	1.2	109	0.00
29	C	2,4-Dichlorophenol	40.000	38.828	2.9	108	0.00
30		1,2,4-Trichlorobenzene	40.000	38.546	3.6	109	0.00
31		Naphthalene	40.000	39.174	2.1	111	0.00
32		Benzoic acid	40.000	44.576	-11.4	121	0.00
33		4-Chloroaniline	40.000	39.101	2.2	108	0.00
34	C	Hexachlorobutadiene	40.000	38.922	2.7	108	0.00
35		Caprolactam	40.000	39.760	0.6	107	0.01
36	C	4-Chloro-3-methylphenol	40.000	38.688	3.3	107	0.00
37		2-Methylnaphthalene	40.000	39.055	2.4	108	0.00
38		1-Methylnaphthalene	40.000	38.756	3.1	110	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	109	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	40.149	-0.4	109	0.00
41	P	Hexachlorocyclopentadiene	40.000	42.482	-6.2	112	0.00
42	S	2,4,6-Tribromophenol	80.000	80.232	-0.3	104	0.00
43	C	2,4,6-Trichlorophenol	40.000	40.339	-0.8	106	0.00
44		2,4,5-Trichlorophenol	40.000	40.280	-0.7	106	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG101918\
Data File : BG037464.D
Acq On : 19 Oct 2018 23:46
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 20 00:23:26 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG101818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 18 14:06:42 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	S	2-Fluorobiphenyl	80.000	77.891	2.6	107	0.00
46		1,1'-Biphenyl	40.000	39.504	1.2	108	0.00
47		2-Chloronaphthalene	40.000	39.487	1.3	108	0.00
48		2-Nitroaniline	40.000	42.334	-5.8	113	0.00
49		Acenaphthylene	40.000	40.049	-0.1	108	0.00
50		Dimethylphthalate	40.000	39.637	0.9	107	0.00
51		2,6-Dinitrotoluene	40.000	40.924	-2.3	108	0.00
52	C	Acenaphthene	40.000	39.495	1.3	109	0.00
53		3-Nitroaniline	40.000	41.483	-3.7	108	0.00
54	P	2,4-Dinitrophenol	40.000	45.900	-14.7	136	0.00
55		Dibenzofuran	40.000	38.770	3.1	107	0.00
56	P	4-Nitrophenol	40.000	41.256	-3.1	108	0.00
57		2,4-Dinitrotoluene	40.000	42.591	-6.5	110	0.00
58		Fluorene	40.000	39.013	2.5	107	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	40.485	-1.2	108	0.00
60		Diethylphthalate	40.000	39.650	0.9	108	0.00
61		4-Chlorophenyl-phenylether	40.000	38.629	3.4	105	0.00
62		4-Nitroaniline	40.000	41.754	-4.4	109	0.00
63		Azobenzene	40.000	39.438	1.4	109	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	108	0.00
65		4,6-Dinitro-2-methylphenol	40.000	41.108	-2.8	120	0.00
66	c	n-Nitrosodiphenylamine	40.000	39.177	2.1	106	0.00
67		4-Bromophenyl-phenylether	40.000	39.364	1.6	107	0.00
68		Hexachlorobenzene	40.000	38.375	4.1	104	0.00
69		Atrazine	40.000	39.758	0.6	105	0.00
70	C	Pentachlorophenol	40.000	41.207	-3.0	108	0.00
71		Phenanthrene	40.000	38.707	3.2	106	0.00
72		Anthracene	40.000	39.325	1.7	107	0.00
73		Carbazole	40.000	39.862	0.3	107	0.00
74		Di-n-butylphthalate	40.000	40.867	-2.2	108	0.00
75	C	Fluoranthene	40.000	39.216	2.0	106	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	109	0.00
77		Benzidine	40.000	44.784	-12.0	118	0.00
78		Pyrene	40.000	39.460	1.3	107	0.00
79	S	Terphenyl-d14	80.000	77.345	3.3	105	0.00
80		Butylbenzylphthalate	40.000	42.407	-6.0	110	0.00
81		Benzo(a)anthracene	40.000	39.284	1.8	106	0.00
82		3,3'-Dichlorobenzidine	40.000	40.685	-1.7	106	0.00
83		Chrysene	40.000	39.508	1.2	107	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	42.184	-5.5	109	0.00
85	c	Di-n-octyl phthalate	40.000	42.650	-6.6	111	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	40.174	-0.4	106	0.00
87	I	Perylene-d12	20.000	20.000	0.0	110	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG101918\
Data File : BG037464.D
Acq On : 19 Oct 2018 23:46
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 20 00:23:26 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG101818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 18 14:06:42 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(b)fluoranthene	40.000	38.678	3.3	105	0.00
89	Benzo(k)fluoranthene	40.000	40.075	-0.2	109	0.00
90 C	Benzo(a)pyrene	40.000	40.011	-0.0	108	0.00
91	Dibenzo(a,h)anthracene	40.000	38.916	2.7	104	0.00
92	Benzo(g,h,i)perylene	40.000	38.967	2.6	106	0.00

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