Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG101218\
Data File : BG037311.D
Acq On : 13 Oct 2018 4:48
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 13 05:24:55 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 11 13:59:58 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	38.064	4.8	132	0.00
3		Pyridine	40.000	41.263	-3.2	132	0.00
4		n-Nitrosodimethylamine	40.000	38.426	3.9	129	0.00
5	S	2-Fluorophenol	80.000	84.114	-5.1	138	0.00
6		Aniline	40.000	38.875	2.8	128	0.00
7	S	Phenol-d6	80.000	82.180	-2.7	133	0.00
8		2-Chlorophenol	40.000	41.087	-2.7	131	0.00
9		Benzaldehyde	40.000	38.852	2.9	128	0.00
10	C	Phenol	40.000	41.485	-3.7	134	0.00
11		bis(2-Chloroethyl)ether	40.000	38.945	2.6	130	0.00
12		1,3-Dichlorobenzene	40.000	39.739	0.7	134	0.00
13	C	1,4-Dichlorobenzene	40.000	39.469	1.3	132	0.00
14		1,2-Dichlorobenzene	40.000	39.011	2.5	132	0.00
15		Benzyl Alcohol	40.000	41.009	-2.5	131	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	36.590	8.5	121	-0.01
17		2-Methylphenol	40.000	41.032	-2.6	132	0.00
18		Hexachloroethane	40.000	39.200	2.0	129	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.251	1.9	129	0.00
20		3+4-Methylphenols	40.000	41.835	-4.6	133	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	132	0.00
22		Acetophenone	40.000	38.705	3.2	127	0.00
23	S	Nitrobenzene-d5	80.000	90.707	-13.4	141	0.00
24		Nitrobenzene	40.000	44.124	-10.3	137	0.00
25		Isophorone	40.000	38.863	2.8	126	0.00
26	C	2-Nitrophenol	40.000	49.118	-22.8#	172	0.00
27		2,4-Dimethylphenol	40.000	40.124	-0.3	131	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.910	2.7	126	0.00
29	C	2,4-Dichlorophenol	40.000	43.653	-9.1	134	0.00
30		1,2,4-Trichlorobenzene	40.000	38.511	3.7	128	0.00
31		Naphthalene	40.000	38.955	2.6	129	0.00
32		Benzoic acid	40.000	51.310	-28.3#	165	0.00
33		4-Chloroaniline	40.000	39.749	0.6	127	0.00
34	C	Hexachlorobutadiene	40.000	38.050	4.9	128	-0.01
35		Caprolactam	40.000	45.329	-13.3	138	0.00
36	C	4-Chloro-3-methylphenol	40.000	43.073	-7.7	133	0.00
37		2-Methylnaphthalene	40.000	39.584	1.0	130	0.00
38		1-Methylnaphthalene	40.000	39.141	2.1	128	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	131	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	38.627	3.4	129	0.00
41	P	Hexachlorocyclopentadiene	40.000	43.092	-7.7	141	0.00
42	S	2,4,6-Tribromophenol	80.000	88.633	-10.8	137	0.00
43	C	2,4,6-Trichlorophenol	40.000	45.369	-13.4	141	0.00
44		2,4,5-Trichlorophenol	40.000	44.563	-11.4	138	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG101218\
Data File : BG037311.D
Acq On : 13 Oct 2018 4:48
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 13 05:24:55 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 11 13:59:58 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	76.572	4.3	128	0.00
46		1,1'-Biphenyl	40.000	38.591	3.5	128	0.00
47		2-Chloronaphthalene	40.000	38.977	2.6	130	0.00
48		2-Nitroaniline	40.000	44.389	-11.0	152	0.00
49		Acenaphthylene	40.000	39.124	2.2	129	0.00
50		Dimethylphthalate	40.000	39.293	1.8	126	0.00
51		2,6-Dinitrotoluene	40.000	43.912	-9.8	148	0.00
52	C	Acenaphthene	40.000	39.518	1.2	130	0.00
53		3-Nitroaniline	40.000	45.743	-14.4	145	0.00
54	P	2,4-Dinitrophenol	40.000	39.086	2.3	132	0.00
55		Dibenzofuran	40.000	38.484	3.8	128	0.00
56	P	4-Nitrophenol	40.000	43.766	-9.4	138	0.00
57		2,4-Dinitrotoluene	40.000	45.950	-14.9	157	0.00
58		Fluorene	40.000	38.525	3.7	128	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	44.015	-10.0	133	0.00
60		Diethylphthalate	40.000	38.952	2.6	124	0.00
61		4-Chlorophenyl-phenylether	40.000	38.303	4.2	126	0.00
62		4-Nitroaniline	40.000	44.087	-10.2	140	0.00
63		Azobenzene	40.000	38.010	5.0	126	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	130	0.00
65		4,6-Dinitro-2-methylphenol	40.000	46.475	-16.2	161	0.00
66	c	n-Nitrosodiphenylamine	40.000	40.089	-0.2	128	0.00
67		4-Bromophenyl-phenylether	40.000	39.267	1.8	126	0.00
68		Hexachlorobenzene	40.000	39.090	2.3	127	0.00
69		Atrazine	40.000	40.190	-0.5	124	0.00
70	C	Pentachlorophenol	40.000	42.043	-5.1	132	0.00
71		Phenanthrene	40.000	38.419	4.0	125	0.00
72		Anthracene	40.000	39.032	2.4	127	0.00
73		Carbazole	40.000	38.693	3.3	125	0.00
74		Di-n-butylphthalate	40.000	41.120	-2.8	124	0.00
75	C	Fluoranthene	40.000	38.044	4.9	122	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	127	0.00
77		Benzidine	40.000	40.326	-0.8	115	0.00
78		Pyrene	40.000	39.185	2.0	123	0.00
79	S	Terphenyl-d14	80.000	75.819	5.2	120	0.00
80		Butylbenzylphthalate	40.000	43.806	-9.5	133	0.00
81		Benzo(a)anthracene	40.000	38.894	2.8	122	0.00
82		3,3'-Dichlorobenzidine	40.000	40.893	-2.2	123	-0.01
83		Chrysene	40.000	39.133	2.2	124	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	42.946	-7.4	130	-0.01
85	c	Di-n-octyl phthalate	40.000	43.728	-9.3	132	-0.02
86		Indeno(1,2,3-cd)pyrene	40.000	40.558	-1.4	125	-0.04
87	I	Perylene-d12	20.000	20.000	0.0	129	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG101218\
Data File : BG037311.D
Acq On : 13 Oct 2018 4:48
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 13 05:24:55 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 11 13:59:58 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	39.525	1.2	126	-0.01
89	Benzo(k)fluoranthene	40.000	38.850	2.9	123	-0.01
90 C	Benzo(a)pyrene	40.000	39.502	1.2	125	-0.01
91	Dibenzo(a,h)anthracene	40.000	39.299	1.8	124	-0.02
92	Benzo(g,h,i)perylene	40.000	39.363	1.6	124	-0.04

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