Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG100618\
Data File : BG037207.D
Acq On : 6 Oct 2018 2:51
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 06 05:33:25 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Sep 24 10:41:23 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	74	-0.01
2		1,4-Dioxane	40.000	0.084	99.8#	0	-0.04
3		Pyridine	40.000	0.000	100.0#	0	-3.93#
4		n-Nitrosodimethylamine	40.000	0.112	99.7#	0	0.00
5	S	2-Fluorophenol	80.000	132.091	-65.1#	119	-0.02
6		Aniline	40.000	0.079	99.8#	0	0.19
7	S	Phenol-d6	80.000	116.662	-45.8#	105	-0.02
8		2-Chlorophenol	40.000	0.000	100.0#	0	-7.61#
9		Benzaldehyde	40.000	0.085	99.8#	0	0.00
10	C	Phenol	40.000	0.187	99.5#	0	0.33
11		bis(2-Chloroethyl)ether	40.000	0.108	99.7#	0	0.09
12		1,3-Dichlorobenzene	40.000	0.000	100.0#	0	-7.94#
13	C	1,4-Dichlorobenzene	40.000	0.000	100.0#	0	-8.09#
14		1,2-Dichlorobenzene	40.000	0.000	100.0#	0	-8.40#
15		Benzyl Alcohol	40.000	1.559	96.1#	3	0.06
16		2,2'-oxybis(1-Chloropropane	40.000	0.000	100.0#	0	-8.58#
17		2-Methylphenol	40.000	0.000	100.0#	0	-8.49#
18		Hexachloroethane	40.000	0.000	100.0#	0	-9.13#
19	P	n-Nitroso-di-n-propylamine	40.000	0.000	100.0#	0	-8.85#
20		3+4-Methylphenols	40.000	0.000	100.0#	0	-8.82#
21	I	Naphthalene-d8	20.000	20.000	0.0	64	-0.02
22		Acetophenone	40.000	0.000	100.0#	0	-8.87#
23	S	Nitrobenzene-d5	80.000	78.289	2.1	62	-0.02
24		Nitrobenzene	40.000	0.121	99.7#	0	-0.05
25		Isophorone	40.000	0.128	99.7#	0	-0.11
26	C	2-Nitrophenol	40.000	0.000	100.0#	0	-9.96#
27		2,4-Dimethylphenol	40.000	0.000	100.0#	0	-10.02#
28		bis(2-Chloroethoxy)methane	40.000	0.000	100.0#	0	-10.26#
29	C	2,4-Dichlorophenol	40.000	0.000	100.0#	0	-10.50#
30		1,2,4-Trichlorobenzene	40.000	0.000	100.0#	0	-10.71#
31		Naphthalene	40.000	0.000	100.0#	0	-10.91#
32		Benzoic acid	40.000	0.000	100.0#	0	-10.16#
33		4-Chloroaniline	40.000	0.000	100.0#	0	-11.01#
34	C	Hexachlorobutadiene	40.000	0.000	100.0#	0	-11.19#
35		Caprolactam	40.000	0.000	100.0#	0	-11.78#
36	C	4-Chloro-3-methylphenol	40.000	0.000	100.0#	0	-12.13#
37		2-Methylnaphthalene	40.000	0.000	100.0#	0	-12.51#
38	I	Acenaphthene-d10	20.000	20.000	0.0	64	-0.02
39		1,2,4,5-Tetrachlorobenzene	40.000	0.000	100.0#	0	-12.87#
40	P	Hexachlorocyclopentadiene	40.000	0.000	100.0#	0	-12.85#
41	S	2,4,6-Tribromophenol	80.000	116.089	-45.1#	90	-0.02
42	C	2,4,6-Trichlorophenol	40.000	0.000	100.0#	0	-13.11#
43		2,4,5-Trichlorophenol	40.000	0.000	100.0#	0	-13.18#
44	S	2-Fluorobiphenyl	80.000	73.517	8.1	58	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG100618\
Data File : BG037207.D
Acq On : 6 Oct 2018 2:51
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 06 05:33:25 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Sep 24 10:41:23 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	0.013	100.0#	0	0.00
46		2-Chloronaphthalene	40.000	0.000	100.0#	0	-13.55#
47		2-Nitroaniline	40.000	0.484	98.8#	1	-0.49
48		Acenaphthylene	40.000	0.000	100.0#	0	-14.40#
49		Dimethylphthalate	40.000	0.000	100.0#	0	-14.13#
50		2,6-Dinitrotoluene	40.000	0.000	100.0#	0	-14.25#
51	C	Acenaphthene	40.000	0.035	99.9#	0	-0.08
52		3-Nitroaniline	40.000	0.000	100.0#	0	-14.58#
53	P	2,4-Dinitrophenol	40.000	0.000	100.0#	0	-14.79#
54		Dibenzofuran	40.000	0.000	100.0#	0	-15.07#
55	P	4-Nitrophenol	40.000	0.000	100.0#	0	-14.89#
56		2,4-Dinitrotoluene	40.000	0.000	100.0#	0	-15.03#
57		Fluorene	40.000	0.718	98.2#	1	0.42
58		2,3,4,6-Tetrachlorophenol	40.000	0.000	100.0#	0	-15.30#
59		Diethylphthalate	40.000	0.000	100.0#	0	-15.49#
60		4-Chlorophenyl-phenylether	40.000	0.000	100.0#	0	-15.71#
61		4-Nitroaniline	40.000	0.000	100.0#	0	-15.75#
62		Azobenzene	40.000	0.098	99.8#	0	0.15
63	I	Phenanthrene-d10	20.000	20.000	0.0	74	-0.02
64		4,6-Dinitro-2-methylphenol	40.000	0.136	99.7#	0	0.35
65	c	n-Nitrosodiphenylamine	40.000	0.693	98.3#	1	0.22
66		4-Bromophenyl-phenylether	40.000	0.000	100.0#	0	-16.61#
67		Hexachlorobenzene	40.000	0.000	100.0#	0	-16.72#
68		Atrazine	40.000	0.000	100.0#	0	-16.88#
69	C	Pentachlorophenol	40.000	0.000	100.0#	0	-17.07#
70		Phenanthrene	40.000	0.000	100.0#	0	-17.47#
71		Anthracene	40.000	0.000	100.0#	0	-17.55#
72		Carbazole	40.000	0.000	100.0#	0	-17.82#
73		Di-n-butylphthalate	40.000	0.006	100.0#	0	0.00
74	C	Fluoranthene	40.000	0.000	100.0#	0	-19.47#
75	I	Chrysene-d12	20.000	20.000	0.0	83	-0.02
76		Benzidine	40.000	1.839	95.4#	4	0.36
77		Pyrene	40.000	0.190	99.5#	0	0.18
78	S	Terphenyl-d14	80.000	87.637	-9.5	91	-0.02
79		Butylbenzylphthalate	40.000	0.019	100.0#	0	-0.01
80		Benzo(a)anthracene	40.000	0.089	99.8#	0	-0.01
81		3,3'-Dichlorobenzidine	40.000	0.000	100.0#	0	-21.58#
82		Chrysene	40.000	0.029	99.9#	0	-0.02
83		Bis(2-ethylhexyl)phthalate	40.000	0.034	99.9#	0	-0.02
84	c	Di-n-octyl phthalate	40.000	0.029	99.9#	0	-0.04
85		Indeno(1,2,3-cd)pyrene	40.000	0.000	100.0#	0	-28.55#
86	I	Perylene-d12	20.000	20.000	0.0	80	-0.05
87		Benzo(b)fluoranthene	40.000	0.018	100.0#	0	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG100618\
Data File : BG037207.D
Acq On : 6 Oct 2018 2:51
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 06 05:33:25 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Sep 24 10:41:23 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	0.005	100.0#	0	-0.04
89 C	Benzo(a)pyrene	40.000	0.004	100.0#	0	-0.05
90	Dibenzo(a,h)anthracene	40.000	0.000	100.0#	0	-28.62#
91	Benzo(g,h,i)perylene	40.000	0.000	100.0#	0	-29.70#

(#) = Out of Range SPCC's out = 0 CCC's out = 13