Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG100518\
Data File : BG037175.D
Acq On : 5 Oct 2018 5:26
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 05 08:47:03 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	99	-0.01
2		1,4-Dioxane	40.000	48.487	-21.2	112	-0.01
3		Pyridine	40.000	46.050	-15.1	109	-0.02
4		n-Nitrosodimethylamine	40.000	46.721	-16.8	113	-0.01
5	S	2-Fluorophenol	80.000	87.600	-9.5	105	-0.02
6		Aniline	40.000	40.744	-1.9	99	-0.02
7	S	Phenol-d6	80.000	87.673	-9.6	105	-0.02
8		2-Chlorophenol	40.000	43.358	-8.4	104	-0.02
9		Benzaldehyde	40.000	40.677	-1.7	99	-0.02
10	C	Phenol	40.000	45.151	-12.9	109	-0.02
11		bis(2-Chloroethyl)ether	40.000	39.322	1.7	101	-0.02
12		1,3-Dichlorobenzene	40.000	40.450	-1.1	98	-0.02
13	C	1,4-Dichlorobenzene	40.000	40.035	-0.1	99	-0.02
14		1,2-Dichlorobenzene	40.000	39.689	0.8	99	-0.02
15		Benzyl Alcohol	40.000	40.472	-1.2	98	-0.02
16		2,2'-oxybis(1-Chloropropane	40.000	42.223	-5.6	102	-0.02
17		2-Methylphenol	40.000	42.235	-5.6	104	-0.02
18		Hexachloroethane	40.000	42.449	-6.1	103	-0.02
19	P	n-Nitroso-di-n-propylamine	40.000	40.482	-1.2	99	-0.02
20		3+4-Methylphenols	40.000	43.504	-8.8	105	-0.02
21	I	Naphthalene-d8	20.000	20.000	0.0	97	-0.02
22		Acetophenone	40.000	41.763	-4.4	103	-0.02
23	S	Nitrobenzene-d5	80.000	85.393	-6.7	102	-0.02
24		Nitrobenzene	40.000	40.936	-2.3	99	-0.01
25		Isophorone	40.000	42.476	-6.2	103	-0.02
26	C	2-Nitrophenol	40.000	44.741	-11.9	105	-0.01
27		2,4-Dimethylphenol	40.000	41.905	-4.8	100	-0.02
28		bis(2-Chloroethoxy)methane	40.000	40.723	-1.8	98	-0.02
29	C	2,4-Dichlorophenol	40.000	44.259	-10.6	103	-0.02
30		1,2,4-Trichlorobenzene	40.000	38.995	2.5	94	-0.02
31		Naphthalene	40.000	40.046	-0.1	98	-0.02
32		Benzoic acid	40.000	35.957	10.1	90	-0.01
33		4-Chloroaniline	40.000	40.364	-0.9	98	-0.02
34	C	Hexachlorobutadiene	40.000	38.535	3.7	93	-0.02
35		Caprolactam	40.000	42.343	-5.9	100	-0.01
36	C	4-Chloro-3-methylphenol	40.000	47.172	-17.9	109	-0.02
37		2-Methylnaphthalene	40.000	39.514	1.2	97	-0.02
38	I	Acenaphthene-d10	20.000	20.000	0.0	98	-0.02
39		1,2,4,5-Tetrachlorobenzene	40.000	38.244	4.4	93	-0.02
40	P	Hexachlorocyclopentadiene	40.000	32.485	18.8	75	-0.02
41	S	2,4,6-Tribromophenol	80.000	79.831	0.2	95	-0.02
42	C	2,4,6-Trichlorophenol	40.000	43.228	-8.1	102	-0.02
43		2,4,5-Trichlorophenol	40.000	44.589	-11.5	103	-0.01
44	S	2-Fluorobiphenyl	80.000	79.055	1.2	96	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG100518\
Data File : BG037175.D
Acq On : 5 Oct 2018 5:26
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 05 08:47:03 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	39.824	0.4	98	-0.02
46		2-Chloronaphthalene	40.000	39.576	1.1	97	-0.02
47		2-Nitroaniline	40.000	44.590	-11.5	104	-0.02
48		Acenaphthylene	40.000	40.443	-1.1	99	-0.02
49		Dimethylphthalate	40.000	39.816	0.5	97	-0.02
50		2,6-Dinitrotoluene	40.000	40.916	-2.3	99	-0.01
51	C	Acenaphthene	40.000	39.637	0.9	97	-0.02
52		3-Nitroaniline	40.000	40.937	-2.3	97	-0.01
53	P	2,4-Dinitrophenol	40.000	39.068	2.3	99	-0.01
54		Dibenzofuran	40.000	40.110	-0.3	97	-0.02
55	P	4-Nitrophenol	40.000	49.333	-23.3	115	-0.02
56		2,4-Dinitrotoluene	40.000	40.944	-2.4	98	-0.01
57		Fluorene	40.000	39.929	0.2	98	-0.02
58		2,3,4,6-Tetrachlorophenol	40.000	42.423	-6.1	99	-0.01
59		Diethylphthalate	40.000	40.450	-1.1	99	-0.01
60		4-Chlorophenyl-phenylether	40.000	38.278	4.3	93	-0.02
61		4-Nitroaniline	40.000	39.366	1.6	94	-0.02
62		Azobenzene	40.000	42.600	-6.5	104	-0.01
63	I	Phenanthrene-d10	20.000	20.000	0.0	97	-0.02
64		4,6-Dinitro-2-methylphenol	40.000	41.221	-3.1	99	-0.01
65	c	n-Nitrosodiphenylamine	40.000	40.983	-2.5	98	-0.01
66		4-Bromophenyl-phenylether	40.000	38.529	3.7	91	-0.02
67		Hexachlorobenzene	40.000	38.647	3.4	92	-0.02
68		Atrazine	40.000	38.037	4.9	89	-0.02
69	C	Pentachlorophenol	40.000	37.580	6.1	87	-0.02
70		Phenanthrene	40.000	39.917	0.2	96	-0.02
71		Anthracene	40.000	40.776	-1.9	98	-0.02
72		Carbazole	40.000	40.797	-2.0	97	-0.02
73		Di-n-butylphthalate	40.000	41.764	-4.4	99	-0.02
74	C	Fluoranthene	40.000	39.631	0.9	95	-0.02
75	I	Chrysene-d12	20.000	20.000	0.0	99	-0.02
76		Benzidine	40.000	48.178	-20.4	120	-0.02
77		Pyrene	40.000	38.750	3.1	95	-0.02
78	S	Terphenyl-d14	80.000	76.443	4.4	95	-0.02
79		Butylbenzylphthalate	40.000	42.067	-5.2	104	-0.02
80		Benzo(a)anthracene	40.000	38.785	3.0	98	-0.02
81		3,3'-Dichlorobenzidine	40.000	39.991	0.0	96	-0.02
82		Chrysene	40.000	39.127	2.2	98	-0.02
83		Bis(2-ethylhexyl)phthalate	40.000	41.720	-4.3	105	-0.02
84	c	Di-n-octyl phthalate	40.000	42.594	-6.5	105	-0.03
85		Indeno(1,2,3-cd)pyrene	40.000	40.595	-1.5	98	-0.05
86	I	Perylene-d12	20.000	20.000	0.0	100	-0.03
87		Benzo(b)fluoranthene	40.000	38.362	4.1	94	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG100518\
Data File : BG037175.D
Acq On : 5 Oct 2018 5:26
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 05 08:47:03 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	40.435	-1.1	102	-0.03
89 C	Benzo(a)pyrene	40.000	39.887	0.3	99	-0.03
90	Dibenzo(a,h)anthracene	40.000	40.203	-0.5	99	-0.06
91	Benzo(g,h,i)perylene	40.000	40.410	-1.0	98	-0.06

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