Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG042919\
Data File : BG040655.D
Acq On : 29 Apr 2019 11:02
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 29 11:36:28 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG042319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 24 05:35:33 2019
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	118	0.00
2		1,4-Dioxane	40.000	41.833	-4.6	128	0.00
3		Pyridine	40.000	42.953	-7.4	125	0.00
4		n-Nitrosodimethylamine	40.000	40.531	-1.3	125	0.00
5	S	2-Fluorophenol	80.000	84.706	-5.9	129	0.00
6		Aniline	40.000	40.584	-1.5	122	0.00
7	S	Phenol-d6	80.000	85.403	-6.8	130	0.00
8		2-Chlorophenol	40.000	40.072	-0.2	120	0.00
9		Benzaldehyde	40.000	47.158	-17.9	135	0.00
10	C	Phenol	40.000	42.391	-6.0	128	0.00
11		bis(2-Chloroethyl)ether	40.000	41.306	-3.3	125	0.00
12		1,3-Dichlorobenzene	40.000	41.758	-4.4	127	0.00
13	C	1,4-Dichlorobenzene	40.000	41.284	-3.2	126	0.00
14		1,2-Dichlorobenzene	40.000	40.634	-1.6	124	0.00
15		Benzyl Alcohol	40.000	40.240	-0.6	121	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	40.012	-0.0	122	0.00
17		2-Methylphenol	40.000	40.827	-2.1	125	0.00
18		Hexachloroethane	40.000	40.401	-1.0	124	-0.01
19	P	n-Nitroso-di-n-propylamine	40.000	39.310	1.7	122	0.00
20		3+4-Methylphenols	40.000	41.109	-2.8	124	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	122	0.00
22		Acetophenone	40.000	40.380	-1.0	123	0.00
23	S	Nitrobenzene-d5	80.000	86.201	-7.8	133	0.00
24		Nitrobenzene	40.000	43.610	-9.0	136	0.00
25		Isophorone	40.000	39.113	2.2	121	-0.01
26	C	2-Nitrophenol	40.000	44.865	-12.2	140	0.00
27		2,4-Dimethylphenol	40.000	39.505	1.2	122	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.280	1.8	120	0.00
29	C	2,4-Dichlorophenol	40.000	41.154	-2.9	125	0.00
30		1,2,4-Trichlorobenzene	40.000	41.783	-4.5	128	0.00
31		Naphthalene	40.000	40.156	-0.4	122	0.00
32		Benzoic acid	40.000	49.385	-23.5	157	0.00
33		4-Chloroaniline	40.000	39.951	0.1	125	0.00
34	C	Hexachlorobutadiene	40.000	42.060	-5.2	130	0.00
35		Caprolactam	40.000	38.885	2.8	118	0.00
36	C	4-Chloro-3-methylphenol	40.000	41.252	-3.1	130	0.00
37		2-Methylnaphthalene	40.000	40.371	-0.9	123	0.00
38		1-Methylnaphthalene	40.000	39.918	0.2	123	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	129	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	39.459	1.4	127	0.00
41	P	Hexachlorocyclopentadiene	40.000	33.966	15.1	110	0.00
42	S	2,4,6-Tribromophenol	80.000	82.193	-2.7	134	0.00
43	C	2,4,6-Trichlorophenol	40.000	40.377	-0.9	134	0.00
44		2,4,5-Trichlorophenol	40.000	41.587	-4.0	136	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG042919\
Data File : BG040655.D
Acq On : 29 Apr 2019 11:02
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 29 11:36:28 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG042319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 24 05:35:33 2019
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	78.872	1.4	126	0.00
46		1,1'-Biphenyl	40.000	38.269	4.3	123	0.00
47		2-Chloronaphthalene	40.000	39.538	1.2	128	0.00
48		2-Nitroaniline	40.000	45.391	-13.5	150	0.00
49		Acenaphthylene	40.000	38.784	3.0	124	0.00
50		Dimethylphthalate	40.000	39.302	1.7	126	0.00
51		2,6-Dinitrotoluene	40.000	43.976	-9.9	142	0.00
52	C	Acenaphthene	40.000	38.452	3.9	127	0.00
53		3-Nitroaniline	40.000	43.934	-9.8	142	0.00
54	P	2,4-Dinitrophenol	40.000	37.396	6.5	128	0.00
55		Dibenzofuran	40.000	39.125	2.2	128	0.00
56	P	4-Nitrophenol	40.000	40.879	-2.2	136	0.00
57		2,4-Dinitrotoluene	40.000	46.544	-16.4	153	0.00
58		Fluorene	40.000	38.823	2.9	126	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	41.606	-4.0	134	0.00
60		Diethylphthalate	40.000	38.436	3.9	124	0.00
61		4-Chlorophenyl-phenylether	40.000	39.545	1.1	129	0.00
62		4-Nitroaniline	40.000	42.070	-5.2	137	0.00
63		Azobenzene	40.000	38.496	3.8	124	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	130	0.00
65		4,6-Dinitro-2-methylphenol	40.000	41.873	-4.7	153	0.00
66	c	n-Nitrosodiphenylamine	40.000	37.686	5.8	124	0.00
67		4-Bromophenyl-phenylether	40.000	39.044	2.4	129	0.00
68		Hexachlorobenzene	40.000	38.322	4.2	128	0.00
69		Atrazine	40.000	37.304	6.7	119	0.00
70	C	Pentachlorophenol	40.000	40.601	-1.5	133	0.00
71		Phenanthrene	40.000	38.849	2.9	126	0.00
72		Anthracene	40.000	38.613	3.5	125	0.00
73		Carbazole	40.000	38.224	4.4	124	0.00
74		Di-n-butylphthalate	40.000	37.800	5.5	123	-0.01
75	C	Fluoranthene	40.000	38.919	2.7	127	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	116	0.00
77		Benzidine	40.000	41.391	-3.5	113	0.00
78		Pyrene	40.000	43.632	-9.1	124	0.00
79	S	Terphenyl-d14	80.000	87.377	-9.2	124	0.00
80		Butylbenzylphthalate	40.000	43.857	-9.6	127	0.00
81		Benzo(a)anthracene	40.000	43.235	-8.1	124	0.00
82		3,3'-Dichlorobenzidine	40.000	41.466	-3.7	118	0.00
83		Chrysene	40.000	43.701	-9.3	125	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	41.354	-3.4	119	0.00
85	c	Di-n-octyl phthalate	40.000	41.984	-5.0	121	-0.02
86		Indeno(1,2,3-cd)pyrene	40.000	42.470	-6.2	124	-0.01
87	I	Perylene-d12	20.000	20.000	0.0	125	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG042919\
Data File : BG040655.D
Acq On : 29 Apr 2019 11:02
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 29 11:36:28 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG042319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 24 05:35:33 2019
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	40.441	-1.1	126	0.00
89	Benzo(k)fluoranthene	40.000	40.323	-0.8	126	0.00
90 C	Benzo(a)pyrene	40.000	40.138	-0.3	125	0.00
91	Dibenzo(a,h)anthracene	40.000	39.413	1.5	123	0.00
92	Benzo(g,h,i)perylene	40.000	39.262	1.8	122	0.00

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