Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080119\
Data File : BG041865.D
Acq On : 1 Aug 2019 16:12
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Aug 02 01:32:00 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG072719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 31 08:57:26 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	93	0.00
2		1,4-Dioxane	40.000	35.747	10.6	83	0.00
3		Pyridine	40.000	37.441	6.4	85	0.00
4		n-Nitrosodimethylamine	40.000	38.491	3.8	89	0.00
5	S	2-Fluorophenol	80.000	79.331	0.8	91	0.00
6		Aniline	40.000	39.276	1.8	90	0.00
7	S	Phenol-d6	80.000	81.488	-1.9	93	0.00
8		2-Chlorophenol	40.000	40.436	-1.1	93	0.00
9		Benzaldehyde	40.000	38.127	4.7	86	0.00
10	C	Phenol	40.000	40.384	-1.0	93	0.00
11		bis(2-Chloroethyl)ether	40.000	38.199	4.5	87	0.00
12		1,3-Dichlorobenzene	40.000	39.667	0.8	92	0.00
13	C	1,4-Dichlorobenzene	40.000	39.352	1.6	92	0.00
14		1,2-Dichlorobenzene	40.000	39.244	1.9	91	0.00
15		Benzyl Alcohol	40.000	40.849	-2.1	93	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	38.732	3.2	88	0.00
17		2-Methylphenol	40.000	40.341	-0.9	93	0.00
18		Hexachloroethane	40.000	41.005	-2.5	98	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.740	0.6	91	0.00
20		3+4-Methylphenols	40.000	41.267	-3.2	94	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	94	0.00
22		Acetophenone	40.000	39.185	2.0	92	0.00
23	S	Nitrobenzene-d5	80.000	85.040	-6.3	99	0.00
24		Nitrobenzene	40.000	41.857	-4.6	98	0.00
25		Isophorone	40.000	40.230	-0.6	94	0.00
26	C	2-Nitrophenol	40.000	48.942	-22.4#	118	0.00
27		2,4-Dimethylphenol	40.000	41.138	-2.8	95	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.733	0.7	92	0.00
29	C	2,4-Dichlorophenol	40.000	43.019	-7.5	100	0.00
30		1,2,4-Trichlorobenzene	40.000	40.571	-1.4	95	0.00
31		Naphthalene	40.000	39.867	0.3	93	0.00
32		Benzoic acid	40.000	43.184	-8.0	115	0.00
33		4-Chloroaniline	40.000	40.123	-0.3	92	0.00
34	C	Hexachlorobutadiene	40.000	40.420	-1.1	96	0.00
35		Caprolactam	40.000	45.017	-12.5	104	0.00
36	C	4-Chloro-3-methylphenol	40.000	43.270	-8.2	99	0.00
37		2-Methylnaphthalene	40.000	41.059	-2.6	95	0.00
38		1-Methylnaphthalene	40.000	41.170	-2.9	96	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	98	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	39.552	1.1	97	0.00
41	P	Hexachlorocyclopentadiene	40.000	42.367	-5.9	104	0.00
42	S	2,4,6-Tribromophenol	80.000	89.974	-12.5	111	0.00
43	C	2,4,6-Trichlorophenol	40.000	42.769	-6.9	104	0.00
44		2,4,5-Trichlorophenol	40.000	43.697	-9.2	107	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080119\
Data File : BG041865.D
Acq On : 1 Aug 2019 16:12
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Aug 02 01:32:00 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG072719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 31 08:57:26 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	80.542	-0.7	96	0.00
46		1,1'-Biphenyl	40.000	39.542	1.1	96	0.00
47		2-Chloronaphthalene	40.000	39.781	0.5	97	0.00
48		2-Nitroaniline	40.000	46.287	-15.7	114	0.00
49		Acenaphthylene	40.000	40.374	-0.9	98	0.00
50		Dimethylphthalate	40.000	42.265	-5.7	102	0.00
51		2,6-Dinitrotoluene	40.000	46.301	-15.8	114	0.00
52	C	Acenaphthene	40.000	40.918	-2.3	100	0.00
53		3-Nitroaniline	40.000	44.782	-12.0	110	0.00
54	P	2,4-Dinitrophenol	40.000	50.059	-25.1#	135	0.00
55		Dibenzofuran	40.000	40.831	-2.1	99	0.00
56	P	4-Nitrophenol	40.000	44.771	-11.9	111	0.00
57		2,4-Dinitrotoluene	40.000	48.403	-21.0	120	0.00
58		Fluorene	40.000	41.190	-3.0	100	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	44.380	-11.0	109	0.00
60		Diethylphthalate	40.000	42.083	-5.2	102	0.00
61		4-Chlorophenyl-phenylether	40.000	41.133	-2.8	101	0.00
62		4-Nitroaniline	40.000	43.543	-8.9	106	0.00
63		Azobenzene	40.000	39.806	0.5	97	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	101	0.00
65		4,6-Dinitro-2-methylphenol	40.000	47.638	-19.1	135	0.00
66	c	n-Nitrosodiphenylamine	40.000	39.969	0.1	101	0.00
67		4-Bromophenyl-phenylether	40.000	40.035	-0.1	101	0.00
68		Hexachlorobenzene	40.000	39.818	0.5	101	0.00
69		Atrazine	40.000	41.290	-3.2	104	0.00
70	C	Pentachlorophenol	40.000	44.100	-10.3	112	0.00
71		Phenanthrene	40.000	40.599	-1.5	101	0.00
72		Anthracene	40.000	40.459	-1.1	101	0.00
73		Carbazole	40.000	40.594	-1.5	101	0.00
74		Di-n-butylphthalate	40.000	42.548	-6.4	104	0.00
75	C	Fluoranthene	40.000	41.026	-2.6	102	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	103	0.00
77		Benzidine	40.000	37.855	5.4	91	0.00
78		Pyrene	40.000	41.337	-3.3	101	0.00
79	S	Terphenyl-d14	80.000	76.405	4.5	101	0.00
80		Butylbenzylphthalate	40.000	43.217	-8.0	108	0.00
81		Benzo(a)anthracene	40.000	40.744	-1.9	102	0.00
82		3,3'-Dichlorobenzidine	40.000	41.261	-3.2	102	0.00
83		Chrysene	40.000	40.599	-1.5	101	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	42.971	-7.4	106	0.00
85	c	Di-n-octyl phthalate	40.000	43.969	-9.9	109	-0.01
86		Indeno(1,2,3-cd)pyrene	40.000	40.324	-0.8	103	-0.01
87	I	Perylene-d12	20.000	20.000	0.0	102	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080119\
Data File : BG041865.D
Acq On : 1 Aug 2019 16:12
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Aug 02 01:32:00 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG072719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 31 08:57:26 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	39.790	0.5	100	0.00
89	Benzo(k)fluoranthene	40.000	40.820	-2.1	104	0.00
90 C	Benzo(a)pyrene	40.000	40.318	-0.8	102	-0.01
91	Dibenzo(a,h)anthracene	40.000	40.134	-0.3	103	-0.02
92	Benzo(g,h,i)perylene	40.000	39.924	0.2	103	-0.02

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