Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG020520\
Data File : BG044301.D
Acq On : 5 Feb 2020 12:49
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 06 01:08:34 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jan 30 16:05:09 2020
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	80	0.00
2		1,4-Dioxane	40.000	40.669	-1.7	79	0.00
3		Pyridine	40.000	44.488	-11.2	79	0.00
4		n-Nitrosodimethylamine	40.000	40.998	-2.5	81	0.00
5	S	2-Fluorophenol	80.000	83.514	-4.4	81	0.00
6		Aniline	40.000	41.847	-4.6	80	0.00
7	S	Phenol-d6	80.000	85.988	-7.5	83	0.00
8		2-Chlorophenol	40.000	41.724	-4.3	80	0.00
9		Benzaldehyde	40.000	39.883	0.3	81	0.00
10	C	Phenol	40.000	42.836	-7.1	82	0.00
11		bis(2-Chloroethyl)ether	40.000	40.300	-0.7	78	0.00
12		1,3-Dichlorobenzene	40.000	40.722	-1.8	80	0.00
13	C	1,4-Dichlorobenzene	40.000	40.916	-2.3	79	0.00
14		1,2-Dichlorobenzene	40.000	41.038	-2.6	79	0.00
15		Benzyl Alcohol	40.000	42.450	-6.1	81	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	40.845	-2.1	79	0.00
17		2-Methylphenol	40.000	42.241	-5.6	82	0.00
18		Hexachloroethane	40.000	40.437	-1.1	79	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	41.539	-3.8	80	0.00
20		3+4-Methylphenols	40.000	43.554	-8.9	83	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	80	0.00
22		Acetophenone	40.000	39.780	0.5	80	0.00
23	S	Nitrobenzene-d5	80.000	80.830	-1.0	81	0.00
24		Nitrobenzene	40.000	40.199	-0.5	81	0.00
25		Isophorone	40.000	40.268	-0.7	81	0.00
26	C	2-Nitrophenol	40.000	41.723	-4.3	82	0.00
27		2,4-Dimethylphenol	40.000	40.805	-2.0	81	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.796	0.5	79	0.00
29	C	2,4-Dichlorophenol	40.000	41.835	-4.6	84	0.00
30		1,2,4-Trichlorobenzene	40.000	39.714	0.7	80	0.00
31		Naphthalene	40.000	40.525	-1.3	81	0.00
32		Benzoic acid	40.000	39.845	0.4	92	-0.01
33		4-Chloroaniline	40.000	42.211	-5.5	85	0.00
34	C	Hexachlorobutadiene	40.000	38.266	4.3	76	0.00
35		Caprolactam	40.000	45.432	-13.6	92	0.00
36	C	4-Chloro-3-methylphenol	40.000	44.167	-10.4	90	0.00
37		2-Methylnaphthalene	40.000	41.299	-3.2	83	0.00
38		1-Methylnaphthalene	40.000	41.425	-3.6	83	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	87	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	37.711	5.7	81	0.00
41	P	Hexachlorocyclopentadiene	40.000	39.128	2.2	82	0.00
42	S	2,4,6-Tribromophenol	80.000	83.781	-4.7	93	0.00
43	C	2,4,6-Trichlorophenol	40.000	41.303	-3.3	89	0.00
44		2,4,5-Trichlorophenol	40.000	42.128	-5.3	90	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG020520\
Data File : BG044301.D
Acq On : 5 Feb 2020 12:49
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 06 01:08:34 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jan 30 16:05:09 2020
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	78.609	1.7	85	0.00
46		1,1'-Biphenyl	40.000	39.266	1.8	85	0.00
47		2-Chloronaphthalene	40.000	39.340	1.6	86	0.00
48		2-Nitroaniline	40.000	42.176	-5.4	93	0.00
49		Acenaphthylene	40.000	40.400	-1.0	88	0.00
50		Dimethylphthalate	40.000	40.420	-1.1	89	0.00
51		2,6-Dinitrotoluene	40.000	40.607	-1.5	90	0.00
52	C	Acenaphthene	40.000	40.359	-0.9	89	0.00
53		3-Nitroaniline	40.000	41.842	-4.6	92	0.00
54	P	2,4-Dinitrophenol	40.000	38.833	2.9	87	0.00
55		Dibenzofuran	40.000	40.896	-2.2	89	0.00
56	P	4-Nitrophenol	40.000	41.330	-3.3	89	0.00
57		2,4-Dinitrotoluene	40.000	41.353	-3.4	92	0.00
58		Fluorene	40.000	41.298	-3.2	91	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	42.200	-5.5	92	0.00
60		Diethylphthalate	40.000	41.099	-2.7	91	0.00
61		4-Chlorophenyl-phenylether	40.000	39.990	0.0	88	0.00
62		4-Nitroaniline	40.000	41.905	-4.8	94	0.00
63		Azobenzene	40.000	40.661	-1.7	89	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	92	0.00
65		4,6-Dinitro-2-methylphenol	40.000	39.953	0.1	90	0.00
66	c	n-Nitrosodiphenylamine	40.000	40.380	-1.0	92	0.00
67		4-Bromophenyl-phenylether	40.000	39.645	0.9	91	0.00
68		Hexachlorobenzene	40.000	39.464	1.3	91	0.00
69		Atrazine	40.000	39.694	0.8	89	0.00
70	C	Pentachlorophenol	40.000	40.232	-0.6	91	0.00
71		Phenanthrene	40.000	40.725	-1.8	93	0.00
72		Anthracene	40.000	40.479	-1.2	93	0.00
73		Carbazole	40.000	40.823	-2.1	93	0.00
74		Di-n-butylphthalate	40.000	41.078	-2.7	92	0.00
75	C	Fluoranthene	40.000	40.824	-2.1	93	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	93	0.00
77		Benzidine	40.000	37.247	6.9	76	0.00
78		Pyrene	40.000	40.655	-1.6	93	0.00
79	S	Terphenyl-d14	80.000	75.902	5.1	93	0.00
80		Butylbenzylphthalate	40.000	41.562	-3.9	94	0.00
81		Benzo(a)anthracene	40.000	41.117	-2.8	95	0.00
82		3,3'-Dichlorobenzidine	40.000	42.422	-6.1	95	0.00
83		Chrysene	40.000	40.984	-2.5	94	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	41.201	-3.0	94	0.00
85	c	Di-n-octyl phthalate	40.000	41.791	-4.5	95	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	40.285	-0.7	94	0.00
87	I	Perylene-d12	20.000	20.000	0.0	93	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG020520\
Data File : BG044301.D
Acq On : 5 Feb 2020 12:49
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 06 01:08:34 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jan 30 16:05:09 2020
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	40.613	-1.5	94	0.00
89	Benzo(k)fluoranthene	40.000	40.270	-0.7	92	0.00
90 C	Benzo(a)pyrene	40.000	40.627	-1.6	93	0.00
91	Dibenzo(a,h)anthracene	40.000	40.240	-0.6	94	0.00
92	Benzo(g,h,i)perylene	40.000	40.115	-0.3	94	0.00

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