Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010320\
Data File : BG043881.D
Acq On : 3 Jan 2020 10:15
Operator : JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jan 03 10:51:50 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Dec 31 13:32:23 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	91	0.00
2		1,4-Dioxane	40.000	39.959	0.1	90	0.00
3		Pyridine	40.000	41.885	-4.7	93	0.00
4		n-Nitrosodimethylamine	40.000	38.046	4.9	87	0.00
5	S	2-Fluorophenol	80.000	82.020	-2.5	90	0.00
6		Aniline	40.000	40.256	-0.6	91	0.00
7	S	Phenol-d6	80.000	81.841	-2.3	91	0.00
8		2-Chlorophenol	40.000	40.509	-1.3	91	0.00
9		Benzaldehyde	40.000	36.703	8.2	88	0.00
10	C	Phenol	40.000	40.713	-1.8	92	0.00
11		bis(2-Chloroethyl)ether	40.000	40.349	-0.9	91	0.00
12		1,3-Dichlorobenzene	40.000	40.355	-0.9	91	0.00
13	C	1,4-Dichlorobenzene	40.000	40.874	-2.2	91	0.00
14		1,2-Dichlorobenzene	40.000	40.516	-1.3	92	0.00
15		Benzyl Alcohol	40.000	39.679	0.8	90	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	40.029	-0.1	92	0.00
17		2-Methylphenol	40.000	39.958	0.1	91	0.00
18		Hexachloroethane	40.000	40.206	-0.5	91	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.998	0.0	91	0.00
20		3+4-Methylphenols	40.000	40.734	-1.8	92	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	92	0.00
22		Acetophenone	40.000	39.894	0.3	92	0.00
23	S	Nitrobenzene-d5	80.000	76.131	4.8	90	0.00
24		Nitrobenzene	40.000	39.271	1.8	91	0.00
25		Isophorone	40.000	39.803	0.5	92	0.00
26	C	2-Nitrophenol	40.000	37.649	5.9	89	0.00
27		2,4-Dimethylphenol	40.000	38.979	2.6	92	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.200	-0.5	92	0.00
29	C	2,4-Dichlorophenol	40.000	39.681	0.8	92	0.00
30		1,2,4-Trichlorobenzene	40.000	40.181	-0.5	93	0.00
31		Naphthalene	40.000	40.891	-2.2	92	0.00
32		Benzoic acid	40.000	33.490	16.3	89	0.00
33		4-Chloroaniline	40.000	39.536	1.2	90	0.00
34	C	Hexachlorobutadiene	40.000	40.107	-0.3	94	0.00
35		Caprolactam	40.000	39.280	1.8	91	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.511	-1.3	94	0.00
37		2-Methylnaphthalene	40.000	40.846	-2.1	94	0.00
38		1-Methylnaphthalene	40.000	41.125	-2.8	95	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	94	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	39.549	1.1	94	0.00
41	P	Hexachlorocyclopentadiene	40.000	39.371	1.6	94	0.00
42	S	2,4,6-Tribromophenol	80.000	83.392	-4.2	97	0.00
43	C	2,4,6-Trichlorophenol	40.000	39.616	1.0	94	0.00
44		2,4,5-Trichlorophenol	40.000	39.757	0.6	94	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010320\
Data File : BG043881.D
Acq On : 3 Jan 2020 10:15
Operator : JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jan 03 10:51:50 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Dec 31 13:32:23 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	84.367	-5.5	94	0.00
46		1,1'-Biphenyl	40.000	40.891	-2.2	94	0.00
47		2-Chloronaphthalene	40.000	40.137	-0.3	94	0.00
48		2-Nitroaniline	40.000	39.395	1.5	93	0.00
49		Acenaphthylene	40.000	41.113	-2.8	94	0.00
50		Dimethylphthalate	40.000	41.536	-3.8	95	0.00
51		2,6-Dinitrotoluene	40.000	39.981	0.0	95	0.00
52	C	Acenaphthene	40.000	40.586	-1.5	96	0.00
53		3-Nitroaniline	40.000	41.135	-2.8	96	0.00
54	P	2,4-Dinitrophenol	40.000	37.102	7.2	91	0.00
55		Dibenzofuran	40.000	41.431	-3.6	95	0.00
56	P	4-Nitrophenol	40.000	42.360	-5.9	97	0.00
57		2,4-Dinitrotoluene	40.000	41.349	-3.4	96	0.00
58		Fluorene	40.000	41.845	-4.6	97	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	40.785	-2.0	95	0.00
60		Diethylphthalate	40.000	41.940	-4.8	96	0.00
61		4-Chlorophenyl-phenylether	40.000	40.894	-2.2	97	0.00
62		4-Nitroaniline	40.000	42.733	-6.8	99	0.00
63		Azobenzene	40.000	41.955	-4.9	96	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	98	0.00
65		4,6-Dinitro-2-methylphenol	40.000	37.169	7.1	96	0.00
66	c	n-Nitrosodiphenylamine	40.000	39.216	2.0	98	0.00
67		4-Bromophenyl-phenylether	40.000	38.790	3.0	98	0.00
68		Hexachlorobenzene	40.000	39.080	2.3	99	0.00
69		Atrazine	40.000	40.708	-1.8	96	0.00
70	C	Pentachlorophenol	40.000	39.813	0.5	95	0.00
71		Phenanthrene	40.000	40.763	-1.9	98	0.00
72		Anthracene	40.000	41.178	-2.9	98	0.00
73		Carbazole	40.000	41.613	-4.0	99	0.00
74		Di-n-butylphthalate	40.000	40.871	-2.2	99	0.00
75	C	Fluoranthene	40.000	40.720	-1.8	100	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	98	0.00
77		Benzidine	40.000	47.149	-17.9	103	0.00
78		Pyrene	40.000	40.780	-2.0	101	0.00
79	S	Terphenyl-d14	80.000	82.414	-3.0	100	0.00
80		Butylbenzylphthalate	40.000	40.270	-0.7	98	0.00
81		Benzo(a)anthracene	40.000	40.695	-1.7	99	0.00
82		3,3'-Dichlorobenzidine	40.000	41.054	-2.6	99	0.00
83		Chrysene	40.000	41.343	-3.4	100	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	40.384	-1.0	98	0.00
85	c	Di-n-octyl phthalate	40.000	40.954	-2.4	98	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	39.891	0.3	99	-0.01
87	I	Perylene-d12	20.000	20.000	0.0	99	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010320\
Data File : BG043881.D
Acq On : 3 Jan 2020 10:15
Operator : JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jan 03 10:51:50 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Dec 31 13:32:23 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	39.824	0.4	100	0.00
89	Benzo(k)fluoranthene	40.000	40.475	-1.2	99	0.00
90 C	Benzo(a)pyrene	40.000	40.008	-0.0	99	0.00
91	Dibenzo(a,h)anthracene	40.000	39.803	0.5	100	0.00
92	Benzo(g,h,i)perylene	40.000	39.701	0.7	100	0.00

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