Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061219\
Data File : BG041208.D
Acq On : 12 Jun 2019 16:17
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 13 07:41:29 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sun Jun 09 23:08:34 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	71	0.00
2		1,4-Dioxane	40.000	37.825	5.4	67	0.00
3		Pyridine	40.000	37.259	6.9	66	0.00
4		n-Nitrosodimethylamine	40.000	41.528	-3.8	73	0.00
5	S	2-Fluorophenol	80.000	78.405	2.0	69	0.00
6		Aniline	40.000	39.744	0.6	71	0.00
7	S	Phenol-d6	80.000	80.566	-0.7	72	0.00
8		2-Chlorophenol	40.000	40.942	-2.4	73	0.00
9		Benzaldehyde	40.000	40.933	-2.3	72	0.00
10	C	Phenol	40.000	40.418	-1.0	72	0.00
11		bis(2-Chloroethyl)ether	40.000	37.829	5.4	66	0.00
12		1,3-Dichlorobenzene	40.000	40.120	-0.3	71	0.00
13	C	1,4-Dichlorobenzene	40.000	40.416	-1.0	70	0.00
14		1,2-Dichlorobenzene	40.000	40.904	-2.3	72	0.00
15		Benzyl Alcohol	40.000	39.315	1.7	70	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	36.998	7.5	66	-0.01
17		2-Methylphenol	40.000	38.619	3.5	68	0.00
18		Hexachloroethane	40.000	40.745	-1.9	73	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	37.719	5.7	67	0.00
20		3+4-Methylphenols	40.000	38.835	2.9	70	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	72	0.00
22		Acetophenone	40.000	39.503	1.2	70	0.00
23	S	Nitrobenzene-d5	80.000	81.780	-2.2	73	0.00
24		Nitrobenzene	40.000	41.188	-3.0	73	0.00
25		Isophorone	40.000	37.992	5.0	67	0.00
26	C	2-Nitrophenol	40.000	43.758	-9.4	77	0.00
27		2,4-Dimethylphenol	40.000	39.264	1.8	69	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.269	4.3	68	0.00
29	C	2,4-Dichlorophenol	40.000	39.629	0.9	69	0.00
30		1,2,4-Trichlorobenzene	40.000	40.540	-1.3	71	0.00
31		Naphthalene	40.000	40.321	-0.8	71	0.00
32		Benzoic acid	40.000	38.728	3.2	70	0.00
33		4-Chloroaniline	40.000	40.438	-1.1	71	0.00
34	C	Hexachlorobutadiene	40.000	41.076	-2.7	75	0.00
35		Caprolactam	40.000	40.762	-1.9	71	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.297	-0.7	71	0.00
37		2-Methylnaphthalene	40.000	40.519	-1.3	72	0.00
38		1-Methylnaphthalene	40.000	41.395	-3.5	73	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	74	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	40.444	-1.1	73	0.00
41	P	Hexachlorocyclopentadiene	40.000	38.490	3.8	74	0.00
42	S	2,4,6-Tribromophenol	80.000	82.096	-2.6	74	0.00
43	C	2,4,6-Trichlorophenol	40.000	40.557	-1.4	73	0.00
44		2,4,5-Trichlorophenol	40.000	39.376	1.6	72	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061219\
Data File : BG041208.D
Acq On : 12 Jun 2019 16:17
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 13 07:41:29 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sun Jun 09 23:08:34 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	80.746	-0.9	73	0.00
46		1,1'-Biphenyl	40.000	39.973	0.1	71	0.00
47		2-Chloronaphthalene	40.000	41.037	-2.6	74	0.00
48		2-Nitroaniline	40.000	41.810	-4.5	75	0.00
49		Acenaphthylene	40.000	40.458	-1.1	71	0.00
50		Dimethylphthalate	40.000	39.146	2.1	70	0.00
51		2,6-Dinitrotoluene	40.000	40.016	-0.0	72	0.00
52	C	Acenaphthene	40.000	39.556	1.1	71	0.00
53		3-Nitroaniline	40.000	40.497	-1.2	71	0.00
54	P	2,4-Dinitrophenol	40.000	40.299	-0.7	77	0.00
55		Dibenzofuran	40.000	40.661	-1.7	72	0.00
56	P	4-Nitrophenol	40.000	39.142	2.1	69	0.00
57		2,4-Dinitrotoluene	40.000	40.760	-1.9	72	0.00
58		Fluorene	40.000	40.239	-0.6	72	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	40.688	-1.7	73	0.00
60		Diethylphthalate	40.000	39.404	1.5	70	0.00
61		4-Chlorophenyl-phenylether	40.000	40.090	-0.2	73	0.00
62		4-Nitroaniline	40.000	40.073	-0.2	72	0.00
63		Azobenzene	40.000	39.573	1.1	70	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	72	0.00
65		4,6-Dinitro-2-methylphenol	40.000	42.782	-7.0	83	0.00
66	c	n-Nitrosodiphenylamine	40.000	40.205	-0.5	71	0.00
67		4-Bromophenyl-phenylether	40.000	39.775	0.6	70	0.00
68		Hexachlorobenzene	40.000	39.682	0.8	71	0.00
69		Atrazine	40.000	40.025	-0.1	65	0.00
70	C	Pentachlorophenol	40.000	40.187	-0.5	71	0.00
71		Phenanthrene	40.000	41.301	-3.3	72	0.00
72		Anthracene	40.000	41.540	-3.8	73	0.00
73		Carbazole	40.000	42.103	-5.3	74	0.00
74		Di-n-butylphthalate	40.000	41.289	-3.2	71	0.00
75	C	Fluoranthene	40.000	42.494	-6.2	74	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	72	0.00
77		Benzidine	40.000	34.321	14.2	58	0.00
78		Pyrene	40.000	43.052	-7.6	75	0.00
79	S	Terphenyl-d14	80.000	88.446	-10.6	76	0.00
80		Butylbenzylphthalate	40.000	42.502	-6.3	74	0.00
81		Benzo(a)anthracene	40.000	41.880	-4.7	74	0.00
82		3,3'-Dichlorobenzidine	40.000	40.478	-1.2	73	0.00
83		Chrysene	40.000	42.099	-5.2	73	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	41.498	-3.7	73	0.00
85	c	Di-n-octyl phthalate	40.000	41.818	-4.5	73	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	32.268	19.3	58	0.00
87	I	Perylene-d12	20.000	20.000	0.0	71	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061219\
Data File : BG041208.D
Acq On : 12 Jun 2019 16:17
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 13 07:41:29 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sun Jun 09 23:08:34 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	40.322	-0.8	70	0.00
89	Benzo(k)fluoranthene	40.000	43.207	-8.0	74	0.00
90 C	Benzo(a)pyrene	40.000	40.479	-1.2	70	0.00
91	Dibenzo(a,h)anthracene	40.000	34.147	14.6	59	0.00
92	Benzo(g,h,i)perylene	40.000	31.249	21.9	55	0.00

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