Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG040919\
Data File : BG040394.D
Acq On : 9 Apr 2019 11:42
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 10 08:10:06 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG032719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 28 05:47:00 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	82	-0.02
2		1,4-Dioxane	40.000	42.006	-5.0	85	-0.03
3		Pyridine	40.000	41.181	-3.0	78	-0.02
4		n-Nitrosodimethylamine	40.000	40.751	-1.9	82	-0.02
5	S	2-Fluorophenol	80.000	81.797	-2.2	81	-0.02
6		Aniline	40.000	40.765	-1.9	80	-0.02
7	S	Phenol-d6	80.000	84.591	-5.7	84	-0.02
8		2-Chlorophenol	40.000	39.995	0.0	79	-0.02
9		Benzaldehyde	40.000	37.859	5.4	72	-0.02
10	C	Phenol	40.000	42.362	-5.9	84	-0.02
11		bis(2-Chloroethyl)ether	40.000	38.218	4.5	75	-0.02
12		1,3-Dichlorobenzene	40.000	40.710	-1.8	80	-0.02
13	C	1,4-Dichlorobenzene	40.000	40.860	-2.1	81	-0.02
14		1,2-Dichlorobenzene	40.000	40.402	-1.0	80	-0.02
15		Benzyl Alcohol	40.000	37.419	6.5	74	-0.02
16		2,2'-oxybis(1-Chloropropane	40.000	36.776	8.1	73	-0.02
17		2-Methylphenol	40.000	39.232	1.9	77	-0.02
18		Hexachloroethane	40.000	38.526	3.7	77	-0.02
19	P	n-Nitroso-di-n-propylamine	40.000	35.848	10.4	72	-0.03
20		3+4-Methylphenols	40.000	40.215	-0.5	79	-0.02
21	I	Naphthalene-d8	20.000	20.000	0.0	77	-0.02
22		Acetophenone	40.000	38.508	3.7	75	-0.02
23	S	Nitrobenzene-d5	80.000	80.944	-1.2	77	-0.02
24		Nitrobenzene	40.000	39.549	1.1	76	-0.02
25		Isophorone	40.000	36.861	7.8	71	-0.03
26	C	2-Nitrophenol	40.000	39.109	2.2	75	-0.02
27		2,4-Dimethylphenol	40.000	37.663	5.8	72	-0.02
28		bis(2-Chloroethoxy)methane	40.000	37.634	5.9	71	-0.02
29	C	2,4-Dichlorophenol	40.000	39.837	0.4	75	-0.02
30		1,2,4-Trichlorobenzene	40.000	39.000	2.5	75	-0.02
31		Naphthalene	40.000	40.582	-1.5	77	-0.02
32		Benzoic acid	40.000	30.732	23.2	63	-0.04
33		4-Chloroaniline	40.000	40.819	-2.0	77	-0.02
34	C	Hexachlorobutadiene	40.000	37.326	6.7	72	-0.02
35		Caprolactam	40.000	43.042	-7.6	81	-0.03
36	C	4-Chloro-3-methylphenol	40.000	41.156	-2.9	78	-0.02
37		2-Methylnaphthalene	40.000	38.882	2.8	74	-0.02
38		1-Methylnaphthalene	40.000	39.352	1.6	75	-0.02
39	I	Acenaphthene-d10	20.000	20.000	0.0	78	-0.02
40		1,2,4,5-Tetrachlorobenzene	40.000	39.177	2.1	74	-0.02
41	P	Hexachlorocyclopentadiene	40.000	37.835	5.4	74	-0.02
42	S	2,4,6-Tribromophenol	80.000	93.868	-17.3	89	-0.02
43	C	2,4,6-Trichlorophenol	40.000	40.093	-0.2	76	-0.02
44		2,4,5-Trichlorophenol	40.000	41.460	-3.7	80	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG040919\
Data File : BG040394.D
Acq On : 9 Apr 2019 11:42
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 10 08:10:06 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG032719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 28 05:47:00 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	80.334	-0.4	75	-0.02
46		1,1'-Biphenyl	40.000	39.426	1.4	75	-0.02
47		2-Chloronaphthalene	40.000	40.485	-1.2	77	-0.02
48		2-Nitroaniline	40.000	42.190	-5.5	79	-0.02
49		Acenaphthylene	40.000	41.955	-4.9	79	-0.02
50		Dimethylphthalate	40.000	40.951	-2.4	78	-0.02
51		2,6-Dinitrotoluene	40.000	43.949	-9.9	84	-0.02
52	C	Acenaphthene	40.000	40.814	-2.0	78	-0.02
53		3-Nitroaniline	40.000	46.143	-15.4	88	-0.02
54	P	2,4-Dinitrophenol	40.000	36.436	8.9	74	-0.02
55		Dibenzofuran	40.000	42.597	-6.5	81	-0.02
56	P	4-Nitrophenol	40.000	45.598	-14.0	88	-0.02
57		2,4-Dinitrotoluene	40.000	47.069	-17.7	89	-0.02
58		Fluorene	40.000	43.806	-9.5	84	-0.02
59		2,3,4,6-Tetrachlorophenol	40.000	44.521	-11.3	84	-0.02
60		Diethylphthalate	40.000	42.150	-5.4	80	-0.02
61		4-Chlorophenyl-phenylether	40.000	41.246	-3.1	78	-0.02
62		4-Nitroaniline	40.000	51.559	-28.9#	99	-0.02
63		Azobenzene	40.000	42.575	-6.4	81	-0.02
64	I	Phenanthrene-d10	20.000	20.000	0.0	95	-0.02
65		4,6-Dinitro-2-methylphenol	40.000	40.463	-1.2	99	-0.02
66	c	n-Nitrosodiphenylamine	40.000	36.662	8.3	85	-0.02
67		4-Bromophenyl-phenylether	40.000	35.281	11.8	82	-0.02
68		Hexachlorobenzene	40.000	38.770	3.1	91	-0.02
69		Atrazine	40.000	40.100	-0.3	92	-0.02
70	C	Pentachlorophenol	40.000	35.540	11.2	85	-0.02
71		Phenanthrene	40.000	41.255	-3.1	95	-0.02
72		Anthracene	40.000	40.843	-2.1	94	-0.02
73		Carbazole	40.000	45.645	-14.1	105	-0.02
74		Di-n-butylphthalate	40.000	42.100	-5.3	97	-0.02
75	C	Fluoranthene	40.000	48.188	-20.5#	111	-0.02
76	I	Chrysene-d12	20.000	20.000	0.0	125	-0.03
77		Benzidine	40.000	39.874	0.3	120	-0.02
78		Pyrene	40.000	37.018	7.5	112	-0.02
79	S	Terphenyl-d14	80.000	70.798	11.5	108	-0.02
80		Butylbenzylphthalate	40.000	37.452	6.4	114	-0.02
81		Benzo(a)anthracene	40.000	39.341	1.6	120	-0.03
82		3,3'-Dichlorobenzidine	40.000	40.293	-0.7	121	-0.03
83		Chrysene	40.000	39.789	0.5	121	-0.03
84		Bis(2-ethylhexyl)phthalate	40.000	36.184	9.5	111	-0.03
85	c	Di-n-octyl phthalate	40.000	36.488	8.8	110	-0.04
86		Indeno(1,2,3-cd)pyrene	40.000	40.268	-0.7	124	-0.08
87	I	Perylene-d12	20.000	20.000	0.0	126	-0.05

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG040919\
Data File : BG040394.D
Acq On : 9 Apr 2019 11:42
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 10 08:10:06 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG032719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 28 05:47:00 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	38.837	2.9	119	-0.05
89	Benzo(k)fluoranthene	40.000	39.469	1.3	123	-0.05
90 C	Benzo(a)pyrene	40.000	39.232	1.9	121	-0.05
91	Dibenzo(a,h)anthracene	40.000	40.392	-1.0	125	-0.07
92	Benzo(g,h,i)perylene	40.000	40.155	-0.4	124	-0.09

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