Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021519\
Data File : BG039524.D
Acq On : 15 Feb 2019 15:25
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 16 03:16:26 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG012919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 29 15:41:51 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	107	-0.01
2		1,4-Dioxane	40.000	38.801	3.0	106	-0.02
3		Pyridine	40.000	40.918	-2.3	105	-0.01
4		n-Nitrosodimethylamine	40.000	36.333	9.2	99	0.00
5	S	2-Fluorophenol	80.000	82.410	-3.0	112	0.00
6		Aniline	40.000	40.100	-0.3	110	-0.01
7	S	Phenol-d6	80.000	85.654	-7.1	115	0.00
8		2-Chlorophenol	40.000	42.686	-6.7	115	-0.01
9		Benzaldehyde	40.000	41.492	-3.7	111	-0.02
10	C	Phenol	40.000	43.097	-7.7	119	0.00
11		bis(2-Chloroethyl)ether	40.000	39.658	0.9	108	-0.02
12		1,3-Dichlorobenzene	40.000	39.539	1.2	109	-0.01
13	C	1,4-Dichlorobenzene	40.000	39.858	0.4	109	-0.01
14		1,2-Dichlorobenzene	40.000	39.398	1.5	109	-0.02
15		Benzyl Alcohol	40.000	41.344	-3.4	110	-0.01
16		2,2'-oxybis(1-Chloropropane	40.000	33.558	16.1	94	-0.02
17		2-Methylphenol	40.000	42.410	-6.0	116	0.00
18		Hexachloroethane	40.000	39.979	0.1	110	-0.01
19	P	n-Nitroso-di-n-propylamine	40.000	38.562	3.6	103	-0.02
20		3+4-Methylphenols	40.000	43.592	-9.0	116	-0.01
21	I	Naphthalene-d8	20.000	20.000	0.0	109	-0.02
22		Acetophenone	40.000	38.850	2.9	111	-0.02
23	S	Nitrobenzene-d5	80.000	94.681	-18.4	133	-0.02
24		Nitrobenzene	40.000	44.411	-11.0	126	-0.01
25		Isophorone	40.000	37.222	6.9	105	-0.02
26	C	2-Nitrophenol	40.000	53.823	-34.6#	173	-0.02
27		2,4-Dimethylphenol	40.000	41.268	-3.2	117	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.618	1.0	110	-0.02
29	C	2,4-Dichlorophenol	40.000	45.409	-13.5	127	0.00
30		1,2,4-Trichlorobenzene	40.000	41.425	-3.6	115	-0.01
31		Naphthalene	40.000	38.551	3.6	111	-0.02
32		Benzoic acid	40.000	48.426	-21.1	155	0.00
33		4-Chloroaniline	40.000	41.169	-2.9	116	-0.01
34	C	Hexachlorobutadiene	40.000	40.662	-1.7	114	-0.02
35		Caprolactam	40.000	41.850	-4.6	111	-0.02
36	C	4-Chloro-3-methylphenol	40.000	43.488	-8.7	119	0.00
37		2-Methylnaphthalene	40.000	38.631	3.4	111	-0.01
38		1-Methylnaphthalene	40.000	38.816	3.0	111	-0.02
39	I	Acenaphthene-d10	20.000	20.000	0.0	111	-0.01
40		1,2,4,5-Tetrachlorobenzene	40.000	40.890	-2.2	117	-0.01
41	P	Hexachlorocyclopentadiene	40.000	51.427	-28.6#	146	-0.02
42	S	2,4,6-Tribromophenol	80.000	90.293	-12.9	127	0.00
43	C	2,4,6-Trichlorophenol	40.000	45.779	-14.4	126	-0.01
44		2,4,5-Trichlorophenol	40.000	46.560	-16.4	125	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021519\
Data File : BG039524.D
Acq On : 15 Feb 2019 15:25
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 16 03:16:26 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG012919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 29 15:41:51 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	78.127	2.3	114	-0.02
46		1,1'-Biphenyl	40.000	38.809	3.0	113	-0.02
47		2-Chloronaphthalene	40.000	40.389	-1.0	117	-0.01
48		2-Nitroaniline	40.000	47.598	-19.0	152	-0.01
49		Acenaphthylene	40.000	39.408	1.5	114	-0.01
50		Dimethylphthalate	40.000	39.160	2.1	113	-0.02
51		2,6-Dinitrotoluene	40.000	48.136	-20.3	146	-0.02
52	C	Acenaphthene	40.000	38.641	3.4	109	-0.01
53		3-Nitroaniline	40.000	46.264	-15.7	141	0.00
54	P	2,4-Dinitrophenol	40.000	45.073	-12.7	137	-0.01
55		Dibenzofuran	40.000	39.839	0.4	116	-0.02
56	P	4-Nitrophenol	40.000	41.532	-3.8	124	0.00
57		2,4-Dinitrotoluene	40.000	51.271	-28.2#	164	-0.01
58		Fluorene	40.000	39.124	2.2	113	-0.02
59		2,3,4,6-Tetrachlorophenol	40.000	44.312	-10.8	121	-0.01
60		Diethylphthalate	40.000	37.787	5.5	110	-0.02
61		4-Chlorophenyl-phenylether	40.000	41.342	-3.4	121	-0.02
62		4-Nitroaniline	40.000	45.763	-14.4	136	-0.01
63		Azobenzene	40.000	36.399	9.0	106	-0.02
64	I	Phenanthrene-d10	20.000	20.000	0.0	109	-0.01
65		4,6-Dinitro-2-methylphenol	40.000	49.610	-24.0	156	-0.01
66	c	n-Nitrosodiphenylamine	40.000	39.342	1.6	111	-0.01
67		4-Bromophenyl-phenylether	40.000	41.652	-4.1	118	-0.02
68		Hexachlorobenzene	40.000	42.138	-5.3	120	-0.01
69		Atrazine	40.000	37.406	6.5	104	-0.01
70	C	Pentachlorophenol	40.000	37.109	7.2	101	-0.01
71		Phenanthrene	40.000	38.847	2.9	111	-0.01
72		Anthracene	40.000	39.153	2.1	112	-0.02
73		Carbazole	40.000	38.523	3.7	111	-0.01
74		Di-n-butylphthalate	40.000	38.206	4.5	109	-0.02
75	C	Fluoranthene	40.000	39.547	1.1	115	-0.02
76	I	Chrysene-d12	20.000	20.000	0.0	116	-0.01
77		Benzidine	40.000	34.915	12.7	107	-0.01
78		Pyrene	40.000	38.447	3.9	116	-0.01
79	S	Terphenyl-d14	80.000	75.702	5.4	115	-0.02
80		Butylbenzylphthalate	40.000	39.803	0.5	117	-0.02
81		Benzo(a)anthracene	40.000	39.023	2.4	118	-0.02
82		3,3'-Dichlorobenzidine	40.000	40.385	-1.0	120	-0.02
83		Chrysene	40.000	39.039	2.4	117	-0.02
84		Bis(2-ethylhexyl)phthalate	40.000	39.006	2.5	115	-0.04
85	c	Di-n-octyl phthalate	40.000	40.076	-0.2	118	-0.05
86		Indeno(1,2,3-cd)pyrene	40.000	40.334	-0.8	119	-0.05
87	I	Perylene-d12	20.000	20.000	0.0	119	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021519\
Data File : BG039524.D
Acq On : 15 Feb 2019 15:25
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 16 03:16:26 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG012919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 29 15:41:51 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	38.616	3.5	118	-0.03
89	Benzo(k)fluoranthene	40.000	37.765	5.6	118	-0.03
90 C	Benzo(a)pyrene	40.000	38.705	3.2	119	-0.03
91	Dibenzo(a,h)anthracene	40.000	38.513	3.7	119	-0.06
92	Benzo(g,h,i)perylene	40.000	37.213	7.0	114	-0.05

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