Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG093017\
Data File : BG028983.D
Acq On : 30 Sep 2017 17:27
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 01 02:45:41 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG091217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 12 16:57:11 2017
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	129	-0.08
2		1,4-Dioxane	40.000	39.587	1.0	131	-0.10
3		Pyridine	40.000	40.347	-0.9	125	-0.09
4		n-Nitrosodimethylamine	40.000	41.908	-4.8	131	-0.09
5	S	2-Fluorophenol	80.000	82.534	-3.2	127	-0.08
6		Aniline	40.000	39.606	1.0	121	-0.07
7	S	Phenol-d6	80.000	78.419	2.0	123	-0.08
8		2-Chlorophenol	40.000	41.469	-3.7	129	-0.08
9		Benzaldehyde	40.000	37.588	6.0	119	-0.08
10	C	Phenol	40.000	40.273	-0.7	125	-0.07
11		bis(2-Chloroethyl)ether	40.000	39.094	2.3	125	-0.08
12		1,3-Dichlorobenzene	40.000	41.035	-2.6	131	-0.08
13	C	1,4-Dichlorobenzene	40.000	41.527	-3.8	128	-0.08
14		1,2-Dichlorobenzene	40.000	40.643	-1.6	130	-0.08
15		Benzyl Alcohol	40.000	37.516	6.2	116	-0.07
16		2,2'-oxybis(1-Chloropropane	40.000	39.184	2.0	124	-0.07
17		2-Methylphenol	40.000	38.548	3.6	122	-0.08
18		Hexachloroethane	40.000	41.354	-3.4	128	-0.08
19	P	n-Nitroso-di-n-propylamine	40.000	37.972	5.1	119	-0.07
20		3+4-Methylphenols	40.000	39.793	0.5	124	-0.07
21	I	Naphthalene-d8	20.000	20.000	0.0	115	-0.07
22		Acetophenone	40.000	41.400	-3.5	121	-0.07
23	S	Nitrobenzene-d5	80.000	83.419	-4.3	123	-0.08
24		Nitrobenzene	40.000	41.017	-2.5	122	-0.08
25		Isophorone	40.000	39.333	1.7	116	-0.08
26	C	2-Nitrophenol	40.000	41.115	-2.8	121	-0.08
27		2,4-Dimethylphenol	40.000	40.680	-1.7	118	-0.07
28		bis(2-Chloroethoxy)methane	40.000	40.377	-0.9	118	-0.07
29	C	2,4-Dichlorophenol	40.000	42.368	-5.9	124	-0.07
30		1,2,4-Trichlorobenzene	40.000	41.556	-3.9	123	-0.07
31		Naphthalene	40.000	41.728	-4.3	123	-0.08
32		Benzoic acid	40.000	12.802	68.0#	24	-0.12
33		4-Chloroaniline	40.000	39.524	1.2	116	-0.07
34	C	Hexachlorobutadiene	40.000	41.297	-3.2	124	-0.08
35		Caprolactam	40.000	36.785	8.0	107	-0.08
36	C	4-Chloro-3-methylphenol	40.000	37.379	6.6	112	-0.06
37		2-Methylnaphthalene	40.000	40.197	-0.5	119	-0.06
38	I	Acenaphthene-d10	20.000	20.000	0.0	115	-0.06
39		1,2,4,5-Tetrachlorobenzene	40.000	43.071	-7.7	120	-0.06
40	P	Hexachlorocyclopentadiene	40.000	41.787	-4.5	122	-0.06
41	S	2,4,6-Tribromophenol	80.000	85.270	-6.6	122	-0.06
42	C	2,4,6-Trichlorophenol	40.000	40.946	-2.4	117	-0.07
43		2,4,5-Trichlorophenol	40.000	40.934	-2.3	116	-0.07
44	S	2-Fluorobiphenyl	80.000	84.691	-5.9	118	-0.06

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG093017\
Data File : BG028983.D
Acq On : 30 Sep 2017 17:27
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 01 02:45:41 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG091217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 12 16:57:11 2017
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		1,1'-Biphenyl	40.000	41.831	-4.6	117	-0.06
46		2-Chloronaphthalene	40.000	41.816	-4.5	117	-0.06
47		2-Nitroaniline	40.000	41.907	-4.8	116	-0.06
48		Acenaphthylene	40.000	40.767	-1.9	115	-0.06
49		Dimethylphthalate	40.000	40.770	-1.9	116	-0.06
50		2,6-Dinitrotoluene	40.000	41.630	-4.1	116	-0.06
51	C	Acenaphthene	40.000	40.525	-1.3	115	-0.06
52		3-Nitroaniline	40.000	41.385	-3.5	116	-0.06
53	P	2,4-Dinitrophenol	40.000	20.246	49.4#	45	-0.06
54		Dibenzofuran	40.000	40.627	-1.6	115	-0.06
55	P	4-Nitrophenol	40.000	42.230	-5.6	113	-0.06
56		2,4-Dinitrotoluene	40.000	40.455	-1.1	111	-0.05
57		Fluorene	40.000	40.405	-1.0	114	-0.06
58		2,3,4,6-Tetrachlorophenol	40.000	43.779	-9.4	124	-0.06
59		Diethylphthalate	40.000	40.280	-0.7	117	-0.05
60		4-Chlorophenyl-phenylether	40.000	40.141	-0.4	116	-0.06
61		4-Nitroaniline	40.000	41.887	-4.7	113	-0.06
62		Azobenzene	40.000	40.176	-0.4	115	-0.06
63	I	Phenanthrene-d10	20.000	20.000	0.0	112	-0.06
64		4,6-Dinitro-2-methylphenol	40.000	31.112	22.2	84	-0.06
65	c	n-Nitrosodiphenylamine	40.000	41.041	-2.6	116	-0.06
66		4-Bromophenyl-phenylether	40.000	41.606	-4.0	117	-0.06
67		Hexachlorobenzene	40.000	40.454	-1.1	115	-0.06
68		Atrazine	40.000	41.940	-4.8	115	-0.06
69	C	Pentachlorophenol	40.000	32.861	17.8	89	-0.06
70		Phenanthrene	40.000	41.637	-4.1	117	-0.06
71		Anthracene	40.000	41.417	-3.5	115	-0.06
72		Carbazole	40.000	42.927	-7.3	117	-0.06
73		Di-n-butylphthalate	40.000	42.738	-6.8	118	-0.06
74	C	Fluoranthene	40.000	43.122	-7.8	117	-0.06
75	I	Chrysene-d12	20.000	20.000	0.0	119	-0.08
76		Benzidine	40.000	44.598	-11.5	126	-0.06
77		Pyrene	40.000	38.612	3.5	117	-0.06
78	S	Terphenyl-d14	80.000	79.323	0.8	118	-0.06
79		Butylbenzylphthalate	40.000	39.894	0.3	120	-0.06
80		Benzo(a)anthracene	40.000	40.689	-1.7	121	-0.08
81		3,3'-Dichlorobenzidine	40.000	40.435	-1.1	118	-0.08
82		Chrysene	40.000	40.777	-1.9	121	-0.08
83		Bis(2-ethylhexyl)phthalate	40.000	40.492	-1.2	119	-0.08
84	c	Di-n-octyl phthalate	40.000	41.039	-2.6	120	-0.10
85		Indeno(1,2,3-cd)pyrene	40.000	41.598	-4.0	124	-0.23
86	I	Perylene-d12	20.000	20.000	0.0	119	-0.15
87		Benzo(b)fluoranthene	40.000	41.164	-2.9	123	-0.13

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG093017\
Data File : BG028983.D
Acq On : 30 Sep 2017 17:27
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 01 02:45:41 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG091217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 12 16:57:11 2017
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(k)fluoranthene	40.000	40.814	-2.0	121	-0.13
89 C	Benzo(a)pyrene	40.000	41.522	-3.8	123	-0.15
90	Dibenzo(a,h)anthracene	40.000	41.763	-4.4	123	-0.24
91	Benzo(g,h,i)perylene	40.000	41.510	-3.8	123	-0.26

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