Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG011217\
Data File : BG025480.D
Acq On : 12 Jan 2017 13:35
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jan 13 00:57:51 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 10 11:26:22 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	117	0.00
2		1,4-Dioxane	40.000	44.749	-11.9	130	0.00
3		Pyridine	40.000	44.459	-11.1	119	0.00
4		n-Nitrosodimethylamine	40.000	46.966	-17.4	126	0.00
5	S	2-Fluorophenol	80.000	88.214	-10.3	126	0.00
6		Aniline	40.000	43.775	-9.4	122	0.00
7	S	Phenol-d6	80.000	88.543	-10.7	120	0.00
8		2-Chlorophenol	40.000	42.605	-6.5	119	0.00
9		Benzaldehyde	40.000	38.210	4.5	110	0.00
10	C	Phenol	40.000	43.634	-9.1	120	0.00
11		bis(2-Chloroethyl)ether	40.000	43.820	-9.6	124	0.00
12		1,3-Dichlorobenzene	40.000	43.253	-8.1	120	0.00
13	C	1,4-Dichlorobenzene	40.000	41.984	-5.0	118	0.00
14		1,2-Dichlorobenzene	40.000	41.767	-4.4	119	0.00
15		Benzyl Alcohol	40.000	42.924	-7.3	113	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	44.992	-12.5	125	0.00
17		2-Methylphenol	40.000	43.458	-8.6	120	0.00
18		Hexachloroethane	40.000	44.021	-10.1	118	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	44.539	-11.3	124	0.00
20		3+4-Methylphenols	40.000	43.795	-9.5	120	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	126	0.00
22		Acetophenone	40.000	40.340	-0.9	117	0.00
23	S	Nitrobenzene-d5	80.000	81.946	-2.4	119	0.00
24		Nitrobenzene	40.000	41.829	-4.6	122	0.00
25		Isophorone	40.000	41.628	-4.1	120	0.00
26	C	2-Nitrophenol	40.000	41.486	-3.7	119	0.00
27		2,4-Dimethylphenol	40.000	40.847	-2.1	118	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.125	2.2	120	0.00
29	C	2,4-Dichlorophenol	40.000	42.197	-5.5	123	0.00
30		1,2,4-Trichlorobenzene	40.000	39.787	0.5	119	0.00
31		Naphthalene	40.000	40.446	-1.1	118	0.00
32		Benzoic acid	40.000	39.669	0.8	113	0.01
33		4-Chloroaniline	40.000	42.502	-6.3	120	0.00
34	C	Hexachlorobutadiene	40.000	40.201	-0.5	119	0.00
35		Caprolactam	40.000	40.757	-1.9	119	0.01
36	C	4-Chloro-3-methylphenol	40.000	40.220	-0.5	116	0.00
37		2-Methylnaphthalene	40.000	41.244	-3.1	120	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	121	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	41.833	-4.6	120	0.00
40	P	Hexachlorocyclopentadiene	40.000	41.326	-3.3	123	0.00
41	S	2,4,6-Tribromophenol	80.000	80.467	-0.6	113	0.00
42	C	2,4,6-Trichlorophenol	40.000	40.773	-1.9	117	0.00
43		2,4,5-Trichlorophenol	40.000	40.568	-1.4	117	0.00
44	S	2-Fluorobiphenyl	80.000	80.854	-1.1	118	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG011217\
Data File : BG025480.D
Acq On : 12 Jan 2017 13:35
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jan 13 00:57:51 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 10 11:26:22 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	41.296	-3.2	119	0.00
46		2-Chloronaphthalene	40.000	40.394	-1.0	119	0.00
47		2-Nitroaniline	40.000	44.338	-10.8	124	0.00
48		Acenaphthylene	40.000	40.523	-1.3	118	0.00
49		Dimethylphthalate	40.000	40.917	-2.3	117	0.00
50		2,6-Dinitrotoluene	40.000	41.315	-3.3	118	0.00
51	C	Acenaphthene	40.000	40.290	-0.7	116	0.00
52		3-Nitroaniline	40.000	40.508	-1.3	113	0.00
53	P	2,4-Dinitrophenol	40.000	40.116	-0.3	115	0.00
54		Dibenzofuran	40.000	41.400	-3.5	119	0.00
55	P	4-Nitrophenol	40.000	41.235	-3.1	113	0.00
56		2,4-Dinitrotoluene	40.000	41.114	-2.8	116	0.00
57		Fluorene	40.000	41.212	-3.0	120	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	39.733	0.7	115	0.00
59		Diethylphthalate	40.000	40.448	-1.1	117	0.00
60		4-Chlorophenyl-phenylether	40.000	41.747	-4.4	121	0.00
61		4-Nitroaniline	40.000	41.886	-4.7	109	0.00
62		Azobenzene	40.000	42.397	-6.0	120	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	117	0.00
64		4,6-Dinitro-2-methylphenol	40.000	44.752	-11.9	114	0.00
65	c	n-Nitrosodiphenylamine	40.000	42.105	-5.3	117	0.00
66		4-Bromophenyl-phenylether	40.000	43.309	-8.3	119	0.00
67		Hexachlorobenzene	40.000	40.941	-2.4	115	0.00
68		Atrazine	40.000	33.137	17.2	92	0.00
69	C	Pentachlorophenol	40.000	38.247	4.4	101	0.00
70		Phenanthrene	40.000	41.893	-4.7	117	0.00
71		Anthracene	40.000	41.710	-4.3	118	0.00
72		Carbazole	40.000	42.153	-5.4	118	0.00
73		Di-n-butylphthalate	40.000	41.399	-3.5	116	0.00
74	C	Fluoranthene	40.000	42.271	-5.7	119	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	133	0.00
76		Benzidine	40.000	35.331	11.7	110	0.00
77		Pyrene	40.000	39.871	0.3	121	0.00
78	S	Terphenyl-d14	80.000	76.335	4.6	118	0.00
79		Butylbenzylphthalate	40.000	40.466	-1.2	128	0.00
80		Benzo(a)anthracene	40.000	40.797	-2.0	126	0.00
81		3,3'-Dichlorobenzidine	40.000	42.598	-6.5	131	0.00
82		Chrysene	40.000	40.848	-2.1	128	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	42.123	-5.3	133	0.00
84	c	Di-n-octyl phthalate	40.000	44.122	-10.3	137	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	44.406	-11.0	137	0.00
86	I	Perylene-d12	20.000	20.000	0.0	139	0.00
87		Benzo(b)fluoranthene	40.000	41.667	-4.2	138	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG011217\
Data File : BG025480.D
Acq On : 12 Jan 2017 13:35
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jan 13 00:57:51 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 10 11:26:22 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	41.145	-2.9	134	0.00
89 C	Benzo(a)pyrene	40.000	41.593	-4.0	137	0.00
90	Dibenzo(a,h)anthracene	40.000	42.027	-5.1	137	0.01
91	Benzo(g,h,i)perylene	40.000	41.929	-4.8	139	0.01

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