Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG121817\
Data File : BG031627.D
Acq On : 18 Dec 2017 10:17
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 19 07:14:38 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG121117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Dec 11 17:25:55 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	164	-0.01
2		1,4-Dioxane	40.000	32.780	18.0	136	-0.01
3		Pyridine	40.000	35.025	12.4	136	-0.01
4		n-Nitrosodimethylamine	40.000	33.700	15.7	131	-0.01
5	S	2-Fluorophenol	80.000	77.534	3.1	151	-0.01
6		Aniline	40.000	37.655	5.9	150	-0.01
7	S	Phenol-d6	80.000	77.856	2.7	153	0.00
8		2-Chlorophenol	40.000	39.774	0.6	157	-0.01
9		Benzaldehyde	40.000	35.887	10.3	144	-0.01
10	C	Phenol	40.000	38.371	4.1	149	0.00
11		bis(2-Chloroethyl)ether	40.000	37.828	5.4	148	-0.01
12		1,3-Dichlorobenzene	40.000	39.158	2.1	157	-0.01
13	C	1,4-Dichlorobenzene	40.000	38.524	3.7	156	-0.02
14		1,2-Dichlorobenzene	40.000	38.512	3.7	157	-0.01
15		Benzyl Alcohol	40.000	36.632	8.4	146	-0.01
16		2,2'-oxybis(1-Chloropropane	40.000	33.122	17.2	134	0.00
17		2-Methylphenol	40.000	38.594	3.5	151	-0.01
18		Hexachloroethane	40.000	38.913	2.7	156	-0.01
19	P	n-Nitroso-di-n-propylamine	40.000	34.983	12.5	140	-0.01
20		3+4-Methylphenols	40.000	36.995	7.5	149	-0.01
21	I	Naphthalene-d8	20.000	20.000	0.0	153	-0.01
22		Acetophenone	40.000	39.988	0.0	146	-0.01
23	S	Nitrobenzene-d5	80.000	78.519	1.9	143	-0.01
24		Nitrobenzene	40.000	39.137	2.2	142	0.00
25		Isophorone	40.000	38.464	3.8	137	-0.01
26	C	2-Nitrophenol	40.000	43.942	-9.9	161	0.00
27		2,4-Dimethylphenol	40.000	41.694	-4.2	150	-0.01
28		bis(2-Chloroethoxy)methane	40.000	40.137	-0.3	147	-0.01
29	C	2,4-Dichlorophenol	40.000	44.386	-11.0	157	-0.01
30		1,2,4-Trichlorobenzene	40.000	42.154	-5.4	160	-0.01
31		Naphthalene	40.000	38.855	2.9	147	-0.01
32		Benzoic acid	40.000	27.882	30.3#	91	0.01
33		4-Chloroaniline	40.000	40.004	-0.0	147	-0.01
34	C	Hexachlorobutadiene	40.000	42.647	-6.6	166	-0.01
35		Caprolactam	40.000	38.641	3.4	141	0.00
36	C	4-Chloro-3-methylphenol	40.000	38.765	3.1	144	0.00
37		2-Methylnaphthalene	40.000	39.041	2.4	146	-0.01
38	I	Acenaphthene-d10	20.000	20.000	0.0	150	-0.01
39		1,2,4,5-Tetrachlorobenzene	40.000	41.558	-3.9	152	0.00
40	P	Hexachlorocyclopentadiene	40.000	29.510	26.2#	95	-0.01
41	S	2,4,6-Tribromophenol	80.000	89.052	-11.3	159	0.00
42	C	2,4,6-Trichlorophenol	40.000	42.902	-7.3	150	0.00
43		2,4,5-Trichlorophenol	40.000	42.273	-5.7	147	-0.01
44	S	2-Fluorobiphenyl	80.000	79.544	0.6	146	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG121817\
Data File : BG031627.D
Acq On : 18 Dec 2017 10:17
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 19 07:14:38 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG121117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Dec 11 17:25:55 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	39.422	1.4	145	0.00
46		2-Chloronaphthalene	40.000	39.152	2.1	144	-0.01
47		2-Nitroaniline	40.000	36.964	7.6	133	-0.01
48		Acenaphthylene	40.000	38.681	3.3	141	0.00
49		Dimethylphthalate	40.000	39.138	2.2	143	0.00
50		2,6-Dinitrotoluene	40.000	40.686	-1.7	147	-0.01
51	C	Acenaphthene	40.000	39.546	1.1	146	-0.01
52		3-Nitroaniline	40.000	39.280	1.8	144	-0.01
53	P	2,4-Dinitrophenol	40.000	19.717	50.7#	53	0.00
54		Dibenzofuran	40.000	38.401	4.0	143	0.00
55	P	4-Nitrophenol	40.000	34.792	13.0	127	0.00
56		2,4-Dinitrotoluene	40.000	39.764	0.6	144	0.00
57		Fluorene	40.000	38.886	2.8	143	-0.01
58		2,3,4,6-Tetrachlorophenol	40.000	41.476	-3.7	143	0.00
59		Diethylphthalate	40.000	38.594	3.5	140	0.00
60		4-Chlorophenyl-phenylether	40.000	38.701	3.2	143	-0.01
61		4-Nitroaniline	40.000	38.545	3.6	140	0.00
62		Azobenzene	40.000	34.639	13.4	125	-0.01
63	I	Phenanthrene-d10	20.000	20.000	0.0	143	-0.01
64		4,6-Dinitro-2-methylphenol	40.000	33.955	15.1	119	0.00
65	c	n-Nitrosodiphenylamine	40.000	41.433	-3.6	145	-0.01
66		4-Bromophenyl-phenylether	40.000	42.353	-5.9	149	-0.01
67		Hexachlorobenzene	40.000	42.632	-6.6	153	0.00
68		Atrazine	40.000	42.573	-6.4	146	0.00
69	C	Pentachlorophenol	40.000	32.484	18.8	112	0.00
70		Phenanthrene	40.000	39.366	1.6	141	-0.01
71		Anthracene	40.000	40.042	-0.1	141	0.00
72		Carbazole	40.000	40.442	-1.1	142	0.00
73		Di-n-butylphthalate	40.000	40.371	-0.9	140	-0.01
74	C	Fluoranthene	40.000	39.030	2.4	139	-0.01
75	I	Chrysene-d12	20.000	20.000	0.0	138	-0.01
76		Benzidine	40.000	44.723	-11.8	142	0.00
77		Pyrene	40.000	41.066	-2.7	137	0.00
78	S	Terphenyl-d14	80.000	80.200	-0.3	136	-0.01
79		Butylbenzylphthalate	40.000	43.750	-9.4	140	-0.01
80		Benzo(a)anthracene	40.000	39.789	0.5	134	0.00
81		3,3'-Dichlorobenzidine	40.000	41.587	-4.0	133	-0.02
82		Chrysene	40.000	39.460	1.3	133	-0.01
83		Bis(2-ethylhexyl)phthalate	40.000	40.571	-1.4	132	-0.02
84	c	Di-n-octyl phthalate	40.000	39.340	1.6	126	-0.02
85		Indeno(1,2,3-cd)pyrene	40.000	35.832	10.4	119	-0.05
86	I	Perylene-d12	20.000	20.000	0.0	126	-0.02
87		Benzo(b)fluoranthene	40.000	41.115	-2.8	127	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG121817\
Data File : BG031627.D
Acq On : 18 Dec 2017 10:17
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 19 07:14:38 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG121117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Dec 11 17:25:55 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	40.706	-1.8	126	-0.02
89 C	Benzo(a)pyrene	40.000	40.596	-1.5	124	-0.02
90	Dibenzo(a,h)anthracene	40.000	39.006	2.5	119	-0.04
91	Benzo(g,h,i)perylene	40.000	38.723	3.2	119	-0.04

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