Evaluate Continuing Calibration Report

Data Path : U:\HPCHEM1\BNA G\Data\BG020218\
Data File : BG032524.D
Acq On : 2 Feb 2018 23:20
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 03 00:46:54 2018
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 29 15:32:55 2018
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	128	0.00
2		1,4-Dioxane	40.000	39.662	0.8	139	0.00
3		Pyridine	40.000	40.536	-1.3	127	0.00
4		n-Nitrosodimethylamine	40.000	41.374	-3.4	125	0.00
5	S	2-Fluorophenol	80.000	83.956	-4.9	131	0.00
6		Aniline	40.000	40.116	-0.3	123	0.00
7	S	Phenol-d6	80.000	84.433	-5.5	134	0.01
8		2-Chlorophenol	40.000	41.527	-3.8	130	0.00
9		Benzaldehyde	40.000	38.489	3.8	123	0.00
10	C	Phenol	40.000	42.141	-5.4	131	0.00
11		bis(2-Chloroethyl)ether	40.000	41.863	-4.7	133	0.00
12		1,3-Dichlorobenzene	40.000	42.789	-7.0	139	0.00
13	C	1,4-Dichlorobenzene	40.000	43.053	-7.6	139	-0.01
14		1,2-Dichlorobenzene	40.000	42.707	-6.8	135	-0.01
15		Benzyl Alcohol	40.000	41.438	-3.6	127	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	38.466	3.8	122	0.00
17		2-Methylphenol	40.000	39.966	0.1	125	0.00
18		Hexachloroethane	40.000	42.534	-6.3	139	-0.01
19	P	n-Nitroso-di-n-propylamine	40.000	41.245	-3.1	127	0.00
20		3+4-Methylphenols	40.000	42.285	-5.7	128	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	128	0.00
22		Acetophenone	40.000	38.646	3.4	125	0.00
23	S	Nitrobenzene-d5	80.000	81.855	-2.3	129	0.00
24		Nitrobenzene	40.000	40.591	-1.5	131	0.00
25		Isophorone	40.000	40.347	-0.9	130	0.00
26	C	2-Nitrophenol	40.000	40.911	-2.3	125	0.00
27		2,4-Dimethylphenol	40.000	38.109	4.7	122	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.177	2.1	126	0.00
29	C	2,4-Dichlorophenol	40.000	41.194	-3.0	132	0.00
30		1,2,4-Trichlorobenzene	40.000	39.665	0.8	131	-0.01
31		Naphthalene	40.000	40.194	-0.5	130	0.00
32		Benzoic acid	40.000	36.509	8.7	118	0.01
33		4-Chloroaniline	40.000	39.191	2.0	127	0.00
34	C	Hexachlorobutadiene	40.000	40.327	-0.8	132	-0.01
35		Caprolactam	40.000	42.348	-5.9	133	0.03
36	C	4-Chloro-3-methylphenol	40.000	40.845	-2.1	133	0.00
37		2-Methylnaphthalene	40.000	40.004	-0.0	129	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	129	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	38.836	2.9	129	-0.01
40	P	Hexachlorocyclopentadiene	40.000	36.653	8.4	120	-0.02
41	S	2,4,6-Tribromophenol	80.000	87.469	-9.3	144	-0.01
42	C	2,4,6-Trichlorophenol	40.000	39.697	0.8	133	-0.01
43		2,4,5-Trichlorophenol	40.000	40.939	-2.3	135	0.00
44	S	2-Fluorobiphenyl	80.000	76.679	4.2	128	-0.01

Evaluate Continuing Calibration Report

Data Path : U:\HPCHEM1\BNA G\Data\BG020218\
Data File : BG032524.D
Acq On : 2 Feb 2018 23:20
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 03 00:46:54 2018
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 29 15:32:55 2018
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	39.437	1.4	131	0.00
46		2-Chloronaphthalene	40.000	39.578	1.1	132	-0.01
47		2-Nitroaniline	40.000	43.129	-7.8	142	0.00
48		Acenaphthylene	40.000	39.956	0.1	133	0.00
49		Dimethylphthalate	40.000	39.662	0.8	136	0.00
50		2,6-Dinitrotoluene	40.000	41.588	-4.0	141	0.00
51	C	Acenaphthene	40.000	37.533	6.2	124	0.00
52		3-Nitroaniline	40.000	42.457	-6.1	140	0.00
53	P	2,4-Dinitrophenol	40.000	17.257	56.9#	51	0.00
54		Dibenzofuran	40.000	39.611	1.0	134	0.00
55	P	4-Nitrophenol	40.000	34.286	14.3	114	0.01
56		2,4-Dinitrotoluene	40.000	42.416	-6.0	140	0.00
57		Fluorene	40.000	39.879	0.3	136	-0.01
58		2,3,4,6-Tetrachlorophenol	40.000	41.685	-4.2	136	-0.01
59		Diethylphthalate	40.000	41.060	-2.7	141	-0.01
60		4-Chlorophenyl-phenylether	40.000	39.927	0.2	134	-0.01
61		4-Nitroaniline	40.000	41.279	-3.2	142	0.00
62		Azobenzene	40.000	41.086	-2.7	141	-0.01
63	I	Phenanthrene-d10	20.000	20.000	0.0	138	0.00
64		4,6-Dinitro-2-methylphenol	40.000	25.660	35.9#	88	0.00
65	c	n-Nitrosodiphenylamine	40.000	39.136	2.2	138	0.00
66		4-Bromophenyl-phenylether	40.000	40.407	-1.0	141	-0.01
67		Hexachlorobenzene	40.000	40.092	-0.2	142	-0.01
68		Atrazine	40.000	41.363	-3.4	142	0.00
69	C	Pentachlorophenol	40.000	35.762	10.6	126	0.00
70		Phenanthrene	40.000	39.075	2.3	139	-0.01
71		Anthracene	40.000	40.470	-1.2	142	-0.01
72		Carbazole	40.000	40.249	-0.6	142	0.00
73		Di-n-butylphthalate	40.000	41.698	-4.2	146	-0.01
74	C	Fluoranthene	40.000	40.334	-0.8	146	-0.01
75	I	Chrysene-d12	20.000	20.000	0.0	149	0.00
76		Benzidine	40.000	39.045	2.4	131	0.00
77		Pyrene	40.000	38.136	4.7	145	-0.01
78	S	Terphenyl-d14	80.000	75.416	5.7	145	0.00
79		Butylbenzylphthalate	40.000	43.001	-7.5	159	0.00
80		Benzo(a)anthracene	40.000	39.986	0.0	152	-0.01
81		3,3'-Dichlorobenzidine	40.000	40.238	-0.6	151	0.00
82		Chrysene	40.000	40.134	-0.3	155	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	43.154	-7.9	160	-0.01
84	c	Di-n-octyl phthalate	40.000	43.943	-9.9	163	-0.02
85		Indeno(1,2,3-cd)pyrene	40.000	42.529	-6.3	160	0.00
86	I	Perylene-d12	20.000	20.000	0.0	156	0.00
87		Benzo(b)fluoranthene	40.000	39.281	1.8	156	0.00

Evaluate Continuing Calibration Report

Data Path : U:\HPCHEM1\BNA G\Data\BG020218\
Data File : BG032524.D
Acq On : 2 Feb 2018 23:20
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 03 00:46:54 2018
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 29 15:32:55 2018
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	40.099	-0.2	161	-0.01
89 C	Benzo(a)pyrene	40.000	40.053	-0.1	160	0.00
90	Dibenzo(a,h)anthracene	40.000	40.537	-1.3	160	0.00
91	Benzo(g,h,i)perylene	40.000	40.075	-0.2	158	0.00

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