Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG071618\
Data File : BG035662.D
Acq On : 16 Jul 2018 10:07
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 16 11:15:04 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 12 14:43:32 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	104	0.00
2		1,4-Dioxane	40.000	37.489	6.3	105	0.00
3		Pyridine	40.000	41.332	-3.3	104	0.00
4		n-Nitrosodimethylamine	40.000	39.251	1.9	103	0.00
5	S	2-Fluorophenol	80.000	79.985	0.0	104	0.00
6		Aniline	40.000	40.156	-0.4	105	0.00
7	S	Phenol-d6	80.000	82.979	-3.7	107	0.00
8		2-Chlorophenol	40.000	42.372	-5.9	110	0.00
9		Benzaldehyde	40.000	40.017	-0.0	102	0.00
10	C	Phenol	40.000	42.359	-5.9	109	0.00
11		bis(2-Chloroethyl)ether	40.000	41.282	-3.2	108	0.00
12		1,3-Dichlorobenzene	40.000	40.125	-0.3	106	0.00
13	C	1,4-Dichlorobenzene	40.000	41.901	-4.8	110	0.00
14		1,2-Dichlorobenzene	40.000	41.441	-3.6	110	0.00
15		Benzyl Alcohol	40.000	39.777	0.6	103	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	39.622	0.9	104	0.00
17		2-Methylphenol	40.000	41.393	-3.5	106	0.00
18		Hexachloroethane	40.000	39.753	0.6	106	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	41.152	-2.9	109	0.00
20		3+4-Methylphenols	40.000	42.635	-6.6	106	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	105	0.00
22		Acetophenone	40.000	38.892	2.8	106	0.00
23	S	Nitrobenzene-d5	80.000	78.747	1.6	105	0.00
24		Nitrobenzene	40.000	39.423	1.4	107	0.00
25		Isophorone	40.000	39.122	2.2	106	0.00
26	C	2-Nitrophenol	40.000	42.753	-6.9	112	0.00
27		2,4-Dimethylphenol	40.000	40.833	-2.1	108	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.340	1.6	106	0.00
29	C	2,4-Dichlorophenol	40.000	42.341	-5.9	110	0.00
30		1,2,4-Trichlorobenzene	40.000	40.488	-1.2	110	0.00
31		Naphthalene	40.000	39.110	2.2	109	0.00
32		Benzoic acid	40.000	33.633	15.9	90	0.00
33		4-Chloroaniline	40.000	40.858	-2.1	112	0.00
34	C	Hexachlorobutadiene	40.000	41.153	-2.9	112	0.00
35		Caprolactam	40.000	37.349	6.6	104	-0.01
36	C	4-Chloro-3-methylphenol	40.000	40.849	-2.1	108	0.00
37		2-Methylnaphthalene	40.000	40.525	-1.3	109	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	104	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	38.867	2.8	109	0.00
40	P	Hexachlorocyclopentadiene	40.000	39.118	2.2	105	0.00
41	S	2,4,6-Tribromophenol	80.000	81.189	-1.5	111	0.00
42	C	2,4,6-Trichlorophenol	40.000	41.842	-4.6	109	0.00
43		2,4,5-Trichlorophenol	40.000	41.771	-4.4	117	0.00
44	S	2-Fluorobiphenyl	80.000	79.939	0.1	111	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG071618\
Data File : BG035662.D
Acq On : 16 Jul 2018 10:07
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 16 11:15:04 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 12 14:43:32 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	39.570	1.1	109	0.00
46		2-Chloronaphthalene	40.000	40.031	-0.1	112	0.00
47		2-Nitroaniline	40.000	41.752	-4.4	113	0.00
48		Acenaphthylene	40.000	39.376	1.6	109	0.00
49		Dimethylphthalate	40.000	38.794	3.0	109	0.00
50		2,6-Dinitrotoluene	40.000	42.862	-7.2	115	0.00
51	C	Acenaphthene	40.000	39.611	1.0	111	0.00
52		3-Nitroaniline	40.000	41.093	-2.7	109	0.00
53	P	2,4-Dinitrophenol	40.000	36.952	7.6	105	0.00
54		Dibenzofuran	40.000	39.750	0.6	110	0.00
55	P	4-Nitrophenol	40.000	36.193	9.5	97	0.00
56		2,4-Dinitrotoluene	40.000	42.069	-5.2	110	0.00
57		Fluorene	40.000	39.524	1.2	111	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	39.650	0.9	107	0.00
59		Diethylphthalate	40.000	38.576	3.6	107	0.00
60		4-Chlorophenyl-phenylether	40.000	40.241	-0.6	113	0.00
61		4-Nitroaniline	40.000	40.331	-0.8	107	0.00
62		Azobenzene	40.000	39.122	2.2	109	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	104	0.00
64		4,6-Dinitro-2-methylphenol	40.000	40.854	-2.1	110	0.00
65	c	n-Nitrosodiphenylamine	40.000	40.651	-1.6	109	0.00
66		4-Bromophenyl-phenylether	40.000	40.269	-0.7	107	0.00
67		Hexachlorobenzene	40.000	40.270	-0.7	109	0.00
68		Atrazine	40.000	38.593	3.5	100	0.00
69	C	Pentachlorophenol	40.000	37.338	6.7	97	0.00
70		Phenanthrene	40.000	39.244	1.9	107	0.00
71		Anthracene	40.000	40.161	-0.4	108	0.00
72		Carbazole	40.000	38.836	2.9	104	0.00
73		Di-n-butylphthalate	40.000	38.700	3.2	103	0.00
74	C	Fluoranthene	40.000	38.038	4.9	102	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	97	0.00
76		Benzidine	40.000	33.322	16.7	88	0.00
77		Pyrene	40.000	40.722	-1.8	104	0.00
78	S	Terphenyl-d14	80.000	80.671	-0.8	102	0.00
79		Butylbenzylphthalate	40.000	41.121	-2.8	101	0.00
80		Benzo(a)anthracene	40.000	39.544	1.1	101	0.00
81		3,3'-Dichlorobenzidine	40.000	41.074	-2.7	102	0.00
82		Chrysene	40.000	39.745	0.6	102	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	41.366	-3.4	101	0.00
84	c	Di-n-octyl phthalate	40.000	41.720	-4.3	102	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	41.148	-2.9	103	0.00
86	I	Perylene-d12	20.000	20.000	0.0	98	0.00
87		Benzo(b)fluoranthene	40.000	39.731	0.7	102	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG071618\
Data File : BG035662.D
Acq On : 16 Jul 2018 10:07
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 16 11:15:04 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 12 14:43:32 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	39.969	0.1	100	0.00
89 C	Benzo(a)pyrene	40.000	40.289	-0.7	101	0.00
90	Dibenzo(a,h)anthracene	40.000	41.445	-3.6	104	-0.01
91	Benzo(g,h,i)perylene	40.000	40.697	-1.7	101	0.00

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