Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG071718\
Data File : BG035681.D
Acq On : 17 Jul 2018 13:59
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 18 03:18:55 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	122	-0.02
2		1,4-Dioxane	40.000	34.009	15.0	112	-0.01
3		Pyridine	40.000	37.077	7.3	110	-0.02
4		n-Nitrosodimethylamine	40.000	35.365	11.6	109	-0.02
5	S	2-Fluorophenol	80.000	79.985	0.0	121	0.00
6		Aniline	40.000	37.547	6.1	115	-0.01
7	S	Phenol-d6	80.000	78.659	1.7	119	0.00
8		2-Chlorophenol	40.000	39.477	1.3	120	-0.01
9		Benzaldehyde	40.000	35.557	11.1	107	-0.02
10	C	Phenol	40.000	39.695	0.8	119	0.00
11		bis(2-Chloroethyl)ether	40.000	38.309	4.2	117	-0.02
12		1,3-Dichlorobenzene	40.000	40.010	-0.0	124	-0.02
13	C	1,4-Dichlorobenzene	40.000	39.215	2.0	121	-0.02
14		1,2-Dichlorobenzene	40.000	38.435	3.9	119	-0.01
15		Benzyl Alcohol	40.000	37.263	6.8	113	-0.01
16		2,2'-oxybis(1-Chloropropane	40.000	42.772	-6.9	131	-0.02
17		2-Methylphenol	40.000	39.728	0.7	119	-0.01
18		Hexachloroethane	40.000	38.902	2.7	122	-0.02
19	P	n-Nitroso-di-n-propylamine	40.000	35.929	10.2	112	-0.01
20		3+4-Methylphenols	40.000	40.542	-1.4	118	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	115	-0.02
22		Acetophenone	40.000	38.490	3.8	115	-0.02
23	S	Nitrobenzene-d5	80.000	77.328	3.3	113	-0.02
24		Nitrobenzene	40.000	37.530	6.2	111	-0.02
25		Isophorone	40.000	37.763	5.6	111	-0.01
26	C	2-Nitrophenol	40.000	45.284	-13.2	129	-0.02
27		2,4-Dimethylphenol	40.000	39.998	0.0	116	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.319	1.7	115	-0.02
29	C	2,4-Dichlorophenol	40.000	42.293	-5.7	119	-0.01
30		1,2,4-Trichlorobenzene	40.000	41.548	-3.9	123	-0.02
31		Naphthalene	40.000	39.742	0.6	120	-0.02
32		Benzoic acid	40.000	39.792	0.5	116	0.00
33		4-Chloroaniline	40.000	39.801	0.5	119	-0.01
34	C	Hexachlorobutadiene	40.000	40.682	-1.7	121	-0.02
35		Caprolactam	40.000	37.530	6.2	114	-0.02
36	C	4-Chloro-3-methylphenol	40.000	39.397	1.5	113	0.00
37		2-Methylnaphthalene	40.000	40.014	-0.0	118	-0.02
38	I	Acenaphthene-d10	20.000	20.000	0.0	110	-0.02
39		1,2,4,5-Tetrachlorobenzene	40.000	41.233	-3.1	121	-0.01
40	P	Hexachlorocyclopentadiene	40.000	42.897	-7.2	121	-0.02
41	S	2,4,6-Tribromophenol	80.000	83.508	-4.4	120	-0.01
42	C	2,4,6-Trichlorophenol	40.000	43.332	-8.3	119	0.00
43		2,4,5-Trichlorophenol	40.000	43.655	-9.1	129	-0.01
44	S	2-Fluorobiphenyl	80.000	80.476	-0.6	118	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG071718\
Data File : BG035681.D
Acq On : 17 Jul 2018 13:59
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 18 03:18:55 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	40.546	-1.4	117	-0.02
46		2-Chloronaphthalene	40.000	40.219	-0.5	118	-0.01
47		2-Nitroaniline	40.000	39.390	1.5	112	-0.01
48		Acenaphthylene	40.000	39.578	1.1	115	-0.01
49		Dimethylphthalate	40.000	39.101	2.2	116	-0.01
50		2,6-Dinitrotoluene	40.000	41.207	-3.0	117	-0.01
51	C	Acenaphthene	40.000	38.921	2.7	115	-0.01
52		3-Nitroaniline	40.000	39.861	0.3	112	-0.01
53	P	2,4-Dinitrophenol	40.000	7.953	80.1#	5	0.01
54		Dibenzofuran	40.000	39.999	0.0	117	-0.01
55	P	4-Nitrophenol	40.000	41.715	-4.3	117	0.00
56		2,4-Dinitrotoluene	40.000	41.618	-4.0	114	-0.01
57		Fluorene	40.000	39.261	1.8	116	-0.01
58		2,3,4,6-Tetrachlorophenol	40.000	39.604	1.0	113	-0.01
59		Diethylphthalate	40.000	38.392	4.0	112	-0.01
60		4-Chlorophenyl-phenylether	40.000	39.334	1.7	116	-0.01
61		4-Nitroaniline	40.000	41.061	-2.7	114	0.00
62		Azobenzene	40.000	36.682	8.3	107	-0.01
63	I	Phenanthrene-d10	20.000	20.000	0.0	109	-0.01
64		4,6-Dinitro-2-methylphenol	40.000	21.878	45.3#	62	-0.01
65	c	n-Nitrosodiphenylamine	40.000	40.803	-2.0	115	-0.01
66		4-Bromophenyl-phenylether	40.000	40.873	-2.2	114	-0.01
67		Hexachlorobenzene	40.000	41.025	-2.6	117	-0.01
68		Atrazine	40.000	41.295	-3.2	113	-0.01
69	C	Pentachlorophenol	40.000	34.879	12.8	95	0.00
70		Phenanthrene	40.000	39.411	1.5	113	-0.01
71		Anthracene	40.000	39.241	1.9	112	-0.01
72		Carbazole	40.000	39.415	1.5	112	0.00
73		Di-n-butylphthalate	40.000	38.878	2.8	109	0.00
74	C	Fluoranthene	40.000	38.428	3.9	109	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	122	0.00
76		Benzidine	40.000	30.614	23.5	102	-0.01
77		Pyrene	40.000	34.036	14.9	110	-0.01
78	S	Terphenyl-d14	80.000	81.516	-1.9	131	0.00
79		Butylbenzylphthalate	40.000	37.926	5.2	118	-0.01
80		Benzo(a)anthracene	40.000	35.802	10.5	115	-0.01
81		3,3'-Dichlorobenzidine	40.000	37.950	5.1	118	0.00
82		Chrysene	40.000	36.563	8.6	118	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	39.073	2.3	121	-0.01
84	c	Di-n-octyl phthalate	40.000	39.143	2.1	120	-0.01
85		Indeno(1,2,3-cd)pyrene	40.000	36.152	9.6	114	-0.03
86	I	Perylene-d12	20.000	20.000	0.0	115	-0.02
87		Benzo(b)fluoranthene	40.000	38.711	3.2	117	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG071718\
Data File : BG035681.D
Acq On : 17 Jul 2018 13:59
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 18 03:18:55 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	40.072	-0.2	118	-0.02
89 C	Benzo(a)pyrene	40.000	39.187	2.0	115	-0.02
90	Dibenzo(a,h)anthracene	40.000	39.128	2.2	115	-0.04
91	Benzo(g,h,i)perylene	40.000	39.156	2.1	114	-0.03

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