Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG072418\
Data File : BG035858.D
Acq On : 24 Jul 2018 22:33
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 25 03:20:16 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 24 14:48:00 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	106	0.00
2		1,4-Dioxane	40.000	36.222	9.4	103	0.00
3		Pyridine	40.000	38.096	4.8	97	0.00
4		n-Nitrosodimethylamine	40.000	35.222	11.9	94	0.00
5	S	2-Fluorophenol	80.000	78.183	2.3	102	0.00
6		Aniline	40.000	37.051	7.4	98	0.00
7	S	Phenol-d6	80.000	78.435	2.0	103	0.00
8		2-Chlorophenol	40.000	40.450	-1.1	106	0.00
9		Benzaldehyde	40.000	34.863	12.8	90	0.00
10	C	Phenol	40.000	39.714	0.7	103	0.00
11		bis(2-Chloroethyl)ether	40.000	38.603	3.5	102	0.00
12		1,3-Dichlorobenzene	40.000	38.082	4.8	102	0.00
13	C	1,4-Dichlorobenzene	40.000	40.157	-0.4	107	0.00
14		1,2-Dichlorobenzene	40.000	38.959	2.6	104	0.00
15		Benzyl Alcohol	40.000	36.449	8.9	96	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	35.038	12.4	93	0.00
17		2-Methylphenol	40.000	38.151	4.6	98	0.00
18		Hexachloroethane	40.000	38.462	3.8	104	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	34.547	13.6	93	0.00
20		3+4-Methylphenols	40.000	39.172	2.1	99	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	98	0.00
22		Acetophenone	40.000	37.633	5.9	96	0.00
23	S	Nitrobenzene-d5	80.000	77.373	3.3	96	0.00
24		Nitrobenzene	40.000	37.581	6.0	95	0.00
25		Isophorone	40.000	36.760	8.1	92	0.00
26	C	2-Nitrophenol	40.000	43.377	-8.4	105	0.00
27		2,4-Dimethylphenol	40.000	38.852	2.9	96	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.393	4.0	96	0.00
29	C	2,4-Dichlorophenol	40.000	43.113	-7.8	104	0.00
30		1,2,4-Trichlorobenzene	40.000	41.348	-3.4	104	0.00
31		Naphthalene	40.000	39.080	2.3	101	0.00
32		Benzoic acid	40.000	32.777	18.1	81	0.00
33		4-Chloroaniline	40.000	39.271	1.8	100	0.00
34	C	Hexachlorobutadiene	40.000	40.892	-2.2	104	0.00
35		Caprolactam	40.000	36.557	8.6	95	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.739	-1.8	100	0.00
37		2-Methylnaphthalene	40.000	40.144	-0.4	101	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	94	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	41.067	-2.7	104	0.00
40	P	Hexachlorocyclopentadiene	40.000	37.534	6.2	91	0.00
41	S	2,4,6-Tribromophenol	80.000	86.645	-8.3	107	0.00
42	C	2,4,6-Trichlorophenol	40.000	42.631	-6.6	101	0.00
43		2,4,5-Trichlorophenol	40.000	43.302	-8.3	110	0.00
44	S	2-Fluorobiphenyl	80.000	81.323	-1.7	103	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG072418\
Data File : BG035858.D
Acq On : 24 Jul 2018 22:33
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 25 03:20:16 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 24 14:48:00 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	41.245	-3.1	103	0.00
46		2-Chloronaphthalene	40.000	40.951	-2.4	104	0.00
47		2-Nitroaniline	40.000	41.147	-2.9	101	0.00
48		Acenaphthylene	40.000	40.564	-1.4	101	0.00
49		Dimethylphthalate	40.000	39.067	2.3	100	0.00
50		2,6-Dinitrotoluene	40.000	43.547	-8.9	106	0.00
51	C	Acenaphthene	40.000	40.532	-1.3	103	0.00
52		3-Nitroaniline	40.000	43.217	-8.0	104	0.00
53	P	2,4-Dinitrophenol	40.000	36.915	7.7	95	0.00
54		Dibenzofuran	40.000	40.582	-1.5	102	0.00
55	P	4-Nitrophenol	40.000	44.353	-10.9	107	0.00
56		2,4-Dinitrotoluene	40.000	43.418	-8.5	103	0.00
57		Fluorene	40.000	40.392	-1.0	103	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	43.835	-9.6	108	0.00
59		Diethylphthalate	40.000	39.061	2.3	98	0.00
60		4-Chlorophenyl-phenylether	40.000	41.494	-3.7	106	0.00
61		4-Nitroaniline	40.000	43.040	-7.6	103	0.00
62		Azobenzene	40.000	38.163	4.6	96	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	97	0.00
64		4,6-Dinitro-2-methylphenol	40.000	39.957	0.1	100	0.00
65	c	n-Nitrosodiphenylamine	40.000	41.110	-2.8	103	0.00
66		4-Bromophenyl-phenylether	40.000	41.299	-3.2	102	0.00
67		Hexachlorobenzene	40.000	41.848	-4.6	105	0.00
68		Atrazine	40.000	38.944	2.6	94	0.00
69	C	Pentachlorophenol	40.000	38.212	4.5	92	0.00
70		Phenanthrene	40.000	40.875	-2.2	104	0.00
71		Anthracene	40.000	41.215	-3.0	104	0.00
72		Carbazole	40.000	40.796	-2.0	102	0.00
73		Di-n-butylphthalate	40.000	40.240	-0.6	100	0.00
74	C	Fluoranthene	40.000	40.746	-1.9	102	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	101	0.00
76		Benzidine	40.000	39.969	0.1	110	0.00
77		Pyrene	40.000	39.364	1.6	105	0.00
78	S	Terphenyl-d14	80.000	78.010	2.5	103	0.00
79		Butylbenzylphthalate	40.000	41.740	-4.4	107	0.00
80		Benzo(a)anthracene	40.000	40.162	-0.4	107	0.00
81		3,3'-Dichlorobenzidine	40.000	42.008	-5.0	108	0.00
82		Chrysene	40.000	40.322	-0.8	107	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	42.024	-5.1	107	0.00
84	c	Di-n-octyl phthalate	40.000	43.702	-9.3	111	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	40.342	-0.9	105	0.00
86	I	Perylene-d12	20.000	20.000	0.0	105	0.00
87		Benzo(b)fluoranthene	40.000	39.609	1.0	109	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG072418\
Data File : BG035858.D
Acq On : 24 Jul 2018 22:33
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 25 03:20:16 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 24 14:48:00 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	41.198	-3.0	110	0.00
89 C	Benzo(a)pyrene	40.000	41.145	-2.9	110	0.00
90	Dibenzo(a,h)anthracene	40.000	39.741	0.6	106	0.00
91	Benzo(g,h,i)perylene	40.000	37.870	5.3	101	0.00

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