Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG121518\
Data File : BG038600.D
Acq On : 15 Dec 2018 12:13
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 16 03:51:44 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 12 15:26:27 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	96	0.00
2		1,4-Dioxane	40.000	40.404	-1.0	96	0.00
3		Pyridine	40.000	39.679	0.8	80	0.00
4		n-Nitrosodimethylamine	40.000	45.343	-13.4	100	0.00
5	S	2-Fluorophenol	80.000	84.831	-6.0	102	0.00
6		Aniline	40.000	42.821	-7.1	102	0.00
7	S	Phenol-d6	80.000	85.890	-7.4	104	0.00
8		2-Chlorophenol	40.000	42.775	-6.9	102	0.00
9		Benzaldehyde	40.000	41.171	-2.9	107	0.00
10	C	Phenol	40.000	42.796	-7.0	103	0.00
11		bis(2-Chloroethyl)ether	40.000	43.246	-8.1	106	0.00
12		1,3-Dichlorobenzene	40.000	42.868	-7.2	104	0.00
13	C	1,4-Dichlorobenzene	40.000	42.048	-5.1	103	0.00
14		1,2-Dichlorobenzene	40.000	42.597	-6.5	105	0.00
15		Benzyl Alcohol	40.000	44.935	-12.3	103	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	42.642	-6.6	104	0.00
17		2-Methylphenol	40.000	43.493	-8.7	102	0.00
18		Hexachloroethane	40.000	42.624	-6.6	104	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	42.997	-7.5	103	0.00
20		3+4-Methylphenols	40.000	45.047	-12.6	106	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	104	0.00
22		Acetophenone	40.000	40.692	-1.7	105	0.00
23	S	Nitrobenzene-d5	80.000	82.255	-2.8	107	0.00
24		Nitrobenzene	40.000	40.413	-1.0	105	0.00
25		Isophorone	40.000	38.241	4.4	85	0.00
26	C	2-Nitrophenol	40.000	39.944	0.1	103	0.00
27		2,4-Dimethylphenol	40.000	40.184	-0.5	104	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.502	-1.3	104	0.00
29	C	2,4-Dichlorophenol	40.000	42.852	-7.1	107	0.00
30		1,2,4-Trichlorobenzene	40.000	40.426	-1.1	105	0.00
31		Naphthalene	40.000	39.538	1.2	103	0.00
32		Benzoic acid	40.000	38.674	3.3	102	0.00
33		4-Chloroaniline	40.000	41.802	-4.5	105	0.00
34	C	Hexachlorobutadiene	40.000	41.780	-4.5	106	0.00
35		Caprolactam	40.000	38.453	3.9	104	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.128	-0.3	105	0.00
37		2-Methylnaphthalene	40.000	40.211	-0.5	105	0.00
38		1-Methylnaphthalene	40.000	39.689	0.8	104	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	106	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	40.220	-0.5	106	0.00
41	P	Hexachlorocyclopentadiene	40.000	38.208	4.5	99	0.00
42	S	2,4,6-Tribromophenol	80.000	81.374	-1.7	106	0.00
43	C	2,4,6-Trichlorophenol	40.000	39.652	0.9	99	0.00
44		2,4,5-Trichlorophenol	40.000	40.511	-1.3	102	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG121518\
Data File : BG038600.D
Acq On : 15 Dec 2018 12:13
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 16 03:51:44 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 12 15:26:27 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	S	2-Fluorobiphenyl	80.000	78.462	1.9	104	0.00
46		1,1'-Biphenyl	40.000	40.010	-0.0	105	0.00
47		2-Chloronaphthalene	40.000	39.235	1.9	104	0.00
48		2-Nitroaniline	40.000	41.602	-4.0	107	0.00
49		Acenaphthylene	40.000	38.908	2.7	102	0.00
50		Dimethylphthalate	40.000	39.834	0.4	104	0.00
51		2,6-Dinitrotoluene	40.000	39.763	0.6	105	0.00
52	C	Acenaphthene	40.000	39.344	1.6	104	0.00
53		3-Nitroaniline	40.000	40.559	-1.4	105	0.00
54	P	2,4-Dinitrophenol	40.000	33.039	17.4	87	0.00
55		Dibenzofuran	40.000	38.698	3.3	104	0.00
56	P	4-Nitrophenol	40.000	38.312	4.2	96	0.00
57		2,4-Dinitrotoluene	40.000	40.221	-0.6	103	0.00
58		Fluorene	40.000	39.601	1.0	104	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	37.750	5.6	96	0.00
60		Diethylphthalate	40.000	38.756	3.1	104	0.00
61		4-Chlorophenyl-phenylether	40.000	39.680	0.8	104	0.00
62		4-Nitroaniline	40.000	40.962	-2.4	103	0.00
63		Azobenzene	40.000	39.306	1.7	103	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	104	0.00
65		4,6-Dinitro-2-methylphenol	40.000	39.057	2.4	100	0.00
66	c	n-Nitrosodiphenylamine	40.000	41.005	-2.5	106	0.00
67		4-Bromophenyl-phenylether	40.000	40.220	-0.5	103	0.00
68		Hexachlorobenzene	40.000	40.347	-0.9	103	0.00
69		Atrazine	40.000	41.066	-2.7	103	0.00
70	C	Pentachlorophenol	40.000	40.186	-0.5	101	0.00
71		Phenanthrene	40.000	39.572	1.1	103	0.00
72		Anthracene	40.000	40.023	-0.1	103	0.00
73		Carbazole	40.000	40.246	-0.6	104	0.00
74		Di-n-butylphthalate	40.000	40.512	-1.3	103	0.00
75	C	Fluoranthene	40.000	39.664	0.8	105	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	100	0.00
77		Benzidine	40.000	35.453	11.4	73	0.00
78		Pyrene	40.000	40.300	-0.7	102	0.00
79	S	Terphenyl-d14	80.000	81.669	-2.1	102	0.00
80		Butylbenzylphthalate	40.000	41.508	-3.8	101	0.00
81		Benzo(a)anthracene	40.000	41.681	-4.2	102	0.00
82		3,3'-Dichlorobenzidine	40.000	42.119	-5.3	100	0.00
83		Chrysene	40.000	40.338	-0.8	100	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	41.944	-4.9	100	0.00
85	c	Di-n-octyl phthalate	40.000	42.206	-5.5	100	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	41.222	-3.1	99	-0.01
87	I	Perylene-d12	20.000	20.000	0.0	100	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG121518\
Data File : BG038600.D
Acq On : 15 Dec 2018 12:13
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 16 03:51:44 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 12 15:26:27 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(b)fluoranthene	40.000	40.882	-2.2	100	0.00
89	Benzo(k)fluoranthene	40.000	39.925	0.2	98	0.00
90 C	Benzo(a)pyrene	40.000	40.391	-1.0	99	0.00
91	Dibenzo(a,h)anthracene	40.000	40.594	-1.5	99	0.00
92	Benzo(g,h,i)perylene	40.000	39.882	0.3	98	0.00

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