Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG113018\
Data File : BG038252.D
Acq On : 30 Nov 2018 11:27
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Nov 30 13:14:21 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG111318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 13 16:39:28 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	77	0.00
2		1,4-Dioxane	40.000	36.074	9.8	73	0.00
3		Pyridine	40.000	37.797	5.5	72	0.00
4		n-Nitrosodimethylamine	40.000	48.997	-22.5	93	0.00
5	S	2-Fluorophenol	80.000	83.631	-4.5	81	0.00
6		Aniline	40.000	42.401	-6.0	81	0.00
7	S	Phenol-d6	80.000	85.871	-7.3	81	0.00
8		2-Chlorophenol	40.000	42.949	-7.4	82	0.00
9		Benzaldehyde	40.000	37.949	5.1	73	0.00
10	C	Phenol	40.000	42.915	-7.3	81	0.00
11		bis(2-Chloroethyl)ether	40.000	42.416	-6.0	82	0.00
12		1,3-Dichlorobenzene	40.000	41.090	-2.7	80	0.00
13	C	1,4-Dichlorobenzene	40.000	40.913	-2.3	78	0.00
14		1,2-Dichlorobenzene	40.000	40.397	-1.0	79	0.00
15		Benzyl Alcohol	40.000	47.280	-18.2	87	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	48.968	-22.4	94	0.00
17		2-Methylphenol	40.000	43.882	-9.7	83	0.00
18		Hexachloroethane	40.000	42.847	-7.1	83	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	42.704	-6.8	81	0.00
20		3+4-Methylphenols	40.000	43.993	-10.0	84	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	81	0.00
22		Acetophenone	40.000	40.445	-1.1	81	0.00
23	S	Nitrobenzene-d5	80.000	85.678	-7.1	85	0.00
24		Nitrobenzene	40.000	43.275	-8.2	88	0.00
25		Isophorone	40.000	41.280	-3.2	83	0.00
26	C	2-Nitrophenol	40.000	42.655	-6.6	83	0.00
27		2,4-Dimethylphenol	40.000	42.024	-5.1	83	0.00
28		bis(2-Chloroethoxy)methane	40.000	41.543	-3.9	83	0.00
29	C	2,4-Dichlorophenol	40.000	41.992	-5.0	82	0.00
30		1,2,4-Trichlorobenzene	40.000	41.280	-3.2	83	0.00
31		Naphthalene	40.000	40.761	-1.9	82	0.00
32		Benzoic acid	40.000	32.283	19.3	63	0.00
33		4-Chloroaniline	40.000	40.575	-1.4	81	0.00
34	C	Hexachlorobutadiene	40.000	41.425	-3.6	83	0.00
35		Caprolactam	40.000	39.318	1.7	77	0.01
36	C	4-Chloro-3-methylphenol	40.000	42.634	-6.6	83	0.00
37		2-Methylnaphthalene	40.000	40.440	-1.1	82	0.00
38		1-Methylnaphthalene	40.000	40.576	-1.4	82	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	80	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	42.020	-5.1	83	0.00
41	P	Hexachlorocyclopentadiene	40.000	35.468	11.3	68	0.00
42	S	2,4,6-Tribromophenol	80.000	81.566	-2.0	79	0.00
43	C	2,4,6-Trichlorophenol	40.000	41.203	-3.0	80	0.00
44		2,4,5-Trichlorophenol	40.000	39.544	1.1	77	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG113018\
Data File : BG038252.D
Acq On : 30 Nov 2018 11:27
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Nov 30 13:14:21 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG111318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 13 16:39:28 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	82.719	-3.4	83	0.00
46		1,1'-Biphenyl	40.000	41.308	-3.3	82	0.00
47		2-Chloronaphthalene	40.000	40.787	-2.0	81	0.00
48		2-Nitroaniline	40.000	45.181	-13.0	87	0.00
49		Acenaphthylene	40.000	41.108	-2.8	81	0.00
50		Dimethylphthalate	40.000	40.815	-2.0	81	0.00
51		2,6-Dinitrotoluene	40.000	41.870	-4.7	81	0.00
52	C	Acenaphthene	40.000	40.815	-2.0	80	0.00
53		3-Nitroaniline	40.000	41.624	-4.1	82	0.00
54	P	2,4-Dinitrophenol	40.000	16.937	57.7#	20	0.00
55		Dibenzofuran	40.000	40.547	-1.4	81	0.00
56	P	4-Nitrophenol	40.000	34.152	14.6	66	0.00
57		2,4-Dinitrotoluene	40.000	42.190	-5.5	81	0.00
58		Fluorene	40.000	40.574	-1.4	81	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	39.723	0.7	76	0.00
60		Diethylphthalate	40.000	40.537	-1.3	81	0.00
61		4-Chlorophenyl-phenylether	40.000	41.113	-2.8	82	0.00
62		4-Nitroaniline	40.000	41.680	-4.2	80	0.00
63		Azobenzene	40.000	42.070	-5.2	84	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	78	0.00
65		4,6-Dinitro-2-methylphenol	40.000	20.477	48.8#	38	0.00
66	c	n-Nitrosodiphenylamine	40.000	41.848	-4.6	80	0.00
67		4-Bromophenyl-phenylether	40.000	43.111	-7.8	82	0.00
68		Hexachlorobenzene	40.000	43.409	-8.5	84	0.00
69		Atrazine	40.000	41.660	-4.1	79	0.00
70	C	Pentachlorophenol	40.000	37.222	6.9	68	0.00
71		Phenanthrene	40.000	41.327	-3.3	80	0.00
72		Anthracene	40.000	41.390	-3.5	80	0.00
73		Carbazole	40.000	41.584	-4.0	79	0.00
74		Di-n-butylphthalate	40.000	42.976	-7.4	82	0.00
75	C	Fluoranthene	40.000	41.519	-3.8	80	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	80	0.00
77		Benzidine	40.000	39.853	0.4	71	0.00
78		Pyrene	40.000	40.957	-2.4	80	0.00
79	S	Terphenyl-d14	80.000	83.961	-5.0	82	0.00
80		Butylbenzylphthalate	40.000	41.714	-4.3	80	0.00
81		Benzo(a)anthracene	40.000	41.450	-3.6	82	0.00
82		3,3'-Dichlorobenzidine	40.000	40.946	-2.4	77	0.00
83		Chrysene	40.000	41.315	-3.3	81	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	42.497	-6.2	83	0.00
85	c	Di-n-octyl phthalate	40.000	43.511	-8.8	82	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	37.569	6.1	72	-0.01
87	I	Perylene-d12	20.000	20.000	0.0	80	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG113018\
Data File : BG038252.D
Acq On : 30 Nov 2018 11:27
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Nov 30 13:14:21 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG111318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 13 16:39:28 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	42.150	-5.4	82	0.00
89	Benzo(k)fluoranthene	40.000	40.844	-2.1	81	0.00
90 C	Benzo(a)pyrene	40.000	42.338	-5.8	82	0.00
91	Dibenzo(a,h)anthracene	40.000	36.963	7.6	72	-0.02
92	Benzo(g,h,i)perylene	40.000	35.772	10.6	69	0.00

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