Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG101118\
Data File : BG037276.D
Acq On : 12 Oct 2018 4:28
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 12 05:03:58 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 11 13:59:58 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	128	0.00
2		1,4-Dioxane	40.000	37.569	6.1	125	0.00
3		Pyridine	40.000	40.272	-0.7	124	0.00
4		n-Nitrosodimethylamine	40.000	39.755	0.6	128	0.00
5	S	2-Fluorophenol	80.000	79.646	0.4	125	0.00
6		Aniline	40.000	39.233	1.9	124	0.00
7	S	Phenol-d6	80.000	79.430	0.7	123	0.00
8		2-Chlorophenol	40.000	40.900	-2.2	125	0.00
9		Benzaldehyde	40.000	37.861	5.3	119	0.00
10	C	Phenol	40.000	39.868	0.3	124	0.00
11		bis(2-Chloroethyl)ether	40.000	39.305	1.7	126	0.00
12		1,3-Dichlorobenzene	40.000	38.412	4.0	124	0.00
13	C	1,4-Dichlorobenzene	40.000	39.141	2.1	126	0.00
14		1,2-Dichlorobenzene	40.000	38.439	3.9	124	0.00
15		Benzyl Alcohol	40.000	40.966	-2.4	126	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	38.411	4.0	122	0.00
17		2-Methylphenol	40.000	40.260	-0.6	124	0.00
18		Hexachloroethane	40.000	39.129	2.2	123	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.376	1.6	124	0.00
20		3+4-Methylphenols	40.000	41.098	-2.7	125	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	127	0.00
22		Acetophenone	40.000	38.859	2.9	123	0.00
23	S	Nitrobenzene-d5	80.000	88.352	-10.4	131	0.00
24		Nitrobenzene	40.000	43.376	-8.4	129	0.00
25		Isophorone	40.000	39.136	2.2	121	0.00
26	C	2-Nitrophenol	40.000	43.456	-8.6	141	0.00
27		2,4-Dimethylphenol	40.000	39.359	1.6	123	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.521	1.2	123	0.00
29	C	2,4-Dichlorophenol	40.000	42.233	-5.6	124	0.00
30		1,2,4-Trichlorobenzene	40.000	39.134	2.2	125	0.00
31		Naphthalene	40.000	39.040	2.4	124	0.00
32		Benzoic acid	40.000	47.099	-17.7	146	0.01
33		4-Chloroaniline	40.000	39.203	2.0	120	0.00
34	C	Hexachlorobutadiene	40.000	37.709	5.7	121	0.00
35		Caprolactam	40.000	43.384	-8.5	126	0.00
36	C	4-Chloro-3-methylphenol	40.000	41.533	-3.8	123	0.00
37		2-Methylnaphthalene	40.000	39.538	1.2	125	0.00
38		1-Methylnaphthalene	40.000	39.006	2.5	122	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	124	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	38.333	4.2	121	0.00
41	P	Hexachlorocyclopentadiene	40.000	40.804	-2.0	126	0.00
42	S	2,4,6-Tribromophenol	80.000	84.908	-6.1	124	0.00
43	C	2,4,6-Trichlorophenol	40.000	43.319	-8.3	127	0.00
44		2,4,5-Trichlorophenol	40.000	42.748	-6.9	126	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG101118\
Data File : BG037276.D
Acq On : 12 Oct 2018 4:28
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 12 05:03:58 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 11 13:59:58 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	76.265	4.7	121	0.00
46		1,1'-Biphenyl	40.000	38.996	2.5	122	0.00
47		2-Chloronaphthalene	40.000	38.718	3.2	122	0.00
48		2-Nitroaniline	40.000	42.362	-5.9	136	0.00
49		Acenaphthylene	40.000	38.872	2.8	121	0.00
50		Dimethylphthalate	40.000	39.275	1.8	120	0.00
51		2,6-Dinitrotoluene	40.000	41.696	-4.2	132	0.00
52	C	Acenaphthene	40.000	39.264	1.8	122	0.00
53		3-Nitroaniline	40.000	42.965	-7.4	129	0.00
54	P	2,4-Dinitrophenol	40.000	38.022	4.9	121	0.00
55		Dibenzofuran	40.000	38.083	4.8	119	0.00
56	P	4-Nitrophenol	40.000	42.339	-5.8	126	0.00
57		2,4-Dinitrotoluene	40.000	41.968	-4.9	134	0.00
58		Fluorene	40.000	38.025	4.9	120	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	42.443	-6.1	121	0.00
60		Diethylphthalate	40.000	39.295	1.8	119	0.00
61		4-Chlorophenyl-phenylether	40.000	38.357	4.1	120	0.00
62		4-Nitroaniline	40.000	42.492	-6.2	128	0.00
63		Azobenzene	40.000	38.154	4.6	120	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	124	0.00
65		4,6-Dinitro-2-methylphenol	40.000	39.418	1.5	126	0.00
66	c	n-Nitrosodiphenylamine	40.000	39.365	1.6	119	0.00
67		4-Bromophenyl-phenylether	40.000	38.851	2.9	118	0.00
68		Hexachlorobenzene	40.000	38.215	4.5	117	0.00
69		Atrazine	40.000	40.043	-0.1	117	0.00
70	C	Pentachlorophenol	40.000	40.925	-2.3	122	0.00
71		Phenanthrene	40.000	37.989	5.0	117	0.00
72		Anthracene	40.000	38.886	2.8	120	0.00
73		Carbazole	40.000	38.670	3.3	118	0.00
74		Di-n-butylphthalate	40.000	41.022	-2.6	118	0.00
75	C	Fluoranthene	40.000	38.037	4.9	116	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	121	0.00
77		Benzidine	40.000	35.365	11.6	96	0.00
78		Pyrene	40.000	39.026	2.4	117	0.00
79	S	Terphenyl-d14	80.000	75.133	6.1	113	0.00
80		Butylbenzylphthalate	40.000	42.375	-5.9	123	0.00
81		Benzo(a)anthracene	40.000	38.908	2.7	117	0.00
82		3,3'-Dichlorobenzidine	40.000	40.567	-1.4	117	0.00
83		Chrysene	40.000	38.691	3.3	117	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	41.776	-4.4	121	0.00
85	c	Di-n-octyl phthalate	40.000	42.499	-6.2	123	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	41.166	-2.9	121	-0.01
87	I	Perylene-d12	20.000	20.000	0.0	126	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG101118\
Data File : BG037276.D
Acq On : 12 Oct 2018 4:28
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 12 05:03:58 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 11 13:59:58 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	38.071	4.8	118	0.00
89	Benzo(k)fluoranthene	40.000	38.911	2.7	120	0.00
90 C	Benzo(a)pyrene	40.000	38.900	2.8	120	0.00
91	Dibenzo(a,h)anthracene	40.000	38.975	2.6	120	0.00
92	Benzo(g,h,i)perylene	40.000	39.059	2.4	120	0.00

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