Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100820\
Data File : BG046811.D
Acq On : 8 Oct 2020 11:03
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 08 13:13:41 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG092220.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 07 11:38:17 2020
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	116	0.00
2		1,4-Dioxane	40.000	34.647	13.4	107	0.00
3		Pyridine	40.000	37.221	6.9	110	0.00
4		n-Nitrosodimethylamine	40.000	36.683	8.3	111	0.00
5	S	2-Fluorophenol	80.000	76.928	3.8	116	0.00
6		Aniline	40.000	36.931	7.7	111	0.00
7	S	Phenol-d6	80.000	77.566	3.0	117	0.00
8		2-Chlorophenol	40.000	38.069	4.8	116	0.00
9		Benzaldehyde	40.000	41.072	-2.7	111	0.00
10	C	Phenol	40.000	36.777	8.1	113	0.00
11		bis(2-Chloroethyl)ether	40.000	36.623	8.4	114	0.00
12		1,3-Dichlorobenzene	40.000	38.850	2.9	117	0.00
13	C	1,4-Dichlorobenzene	40.000	39.144	2.1	118	0.00
14		1,2-Dichlorobenzene	40.000	39.459	1.4	118	0.00
15		Benzyl Alcohol	40.000	36.431	8.9	110	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	32.014	20.0	98	0.00
17		2-Methylphenol	40.000	37.329	6.7	115	0.00
18		Hexachloroethane	40.000	37.510	6.2	116	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	35.173	12.1	106	0.00
20		3+4-Methylphenols	40.000	38.167	4.6	115	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	111	0.00
22		Acetophenone	40.000	39.004	2.5	115	0.00
23	S	Nitrobenzene-d5	80.000	92.941	-16.2	133	0.00
24		Nitrobenzene	40.000	43.606	-9.0	126	0.00
25		Isophorone	40.000	36.889	7.8	108	0.00
26	C	2-Nitrophenol	40.000	51.905	-29.8#	147	0.00
27		2,4-Dimethylphenol	40.000	38.290	4.3	115	0.00
28		bis(2-Chloroethoxy)methane	40.000	36.706	8.2	109	0.00
29	C	2,4-Dichlorophenol	40.000	42.852	-7.1	124	0.00
30		1,2,4-Trichlorobenzene	40.000	41.570	-3.9	121	0.00
31		Naphthalene	40.000	40.154	-0.4	118	0.00
32		Benzoic acid	40.000	50.749	-26.9#	162	0.02
33		4-Chloroaniline	40.000	39.517	1.2	118	0.00
34	C	Hexachlorobutadiene	40.000	42.230	-5.6	123	0.00
35		Caprolactam	40.000	40.178	-0.4	118	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.155	-0.4	119	0.00
37		2-Methylnaphthalene	40.000	40.538	-1.3	121	0.00
38		1-Methylnaphthalene	40.000	41.076	-2.7	120	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	116	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	40.777	-1.9	123	0.00
41	P	Hexachlorocyclopentadiene	40.000	36.999	7.5	110	0.00
42	S	2,4,6-Tribromophenol	80.000	97.665	-22.1	145	0.00
43	C	2,4,6-Trichlorophenol	40.000	43.120	-7.8	126	0.00
44		2,4,5-Trichlorophenol	40.000	44.250	-10.6	134	0.00
45	S	2-Fluorobiphenyl	80.000	78.633	1.7	119	0.00
46		1,1'-Biphenyl	40.000	40.099	-0.2	120	0.00
47		2-Chloronaphthalene	40.000	40.251	-0.6	123	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100820\
Data File : BG046811.D
Acq On : 8 Oct 2020 11:03
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 08 13:13:41 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG092220.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 07 11:38:17 2020
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)
48		2-Nitroaniline	40.000	52.352	-30.9#	152	0.00
49		Acenaphthylene	40.000	39.956	0.1	123	0.00
50		Dimethylphthalate	40.000	39.772	0.6	120	0.00
51		2,6-Dinitrotoluene	40.000	53.364	-33.4#	155	0.00
52	C	Acenaphthene	40.000	41.026	-2.6	126	0.00
53		3-Nitroaniline	40.000	50.101	-25.3#	147	0.00
54	P	2,4-Dinitrophenol	40.000	51.559	-28.9#	161	0.00
55		Dibenzofuran	40.000	40.644	-1.6	124	0.00
56	P	4-Nitrophenol	40.000	51.000	-27.5#	151	0.00
57		2,4-Dinitrotoluene	40.000	57.857	-44.6#	167	0.00
58		Fluorene	40.000	40.443	-1.1	123	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	46.016	-15.0	139	0.00
60		Diethylphthalate	40.000	38.811	3.0	119	0.00
61		4-Chlorophenyl-phenylether	40.000	41.147	-2.9	124	0.00
62		4-Nitroaniline	40.000	49.780	-24.5	148	0.00
63		Azobenzene	40.000	36.954	7.6	113	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	119	0.00
65		4,6-Dinitro-2-methylphenol	40.000	54.105	-35.3#	182	0.00
66	c	n-Nitrosodiphenylamine	40.000	38.343	4.1	121	0.00
67		4-Bromophenyl-phenylether	40.000	40.911	-2.3	129	0.00
68		Hexachlorobenzene	40.000	40.430	-1.1	129	0.00
69		Atrazine	40.000	38.146	4.6	126	0.00
70	C	Pentachlorophenol	40.000	45.128	-12.8	138	0.00
71		Phenanthrene	40.000	39.799	0.5	127	0.00
72		Anthracene	40.000	39.822	0.4	128	0.00
73		Carbazole	40.000	40.163	-0.4	129	0.00
74		Di-n-butylphthalate	40.000	38.142	4.6	122	0.00
75	C	Fluoranthene	40.000	41.605	-4.0	134	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	136	0.00
77		Benzidine	40.000	40.984	-2.5	144	0.00
78		Pyrene	40.000	37.959	5.1	137	0.00
79	S	Terphenyl-d14	80.000	74.646	6.7	134	0.00
80		Butylbenzylphthalate	40.000	40.128	-0.3	142	0.00
81		Benzo(a)anthracene	40.000	39.311	1.7	143	0.00
82		3,3'-Dichlorobenzidine	40.000	38.548	3.6	138	0.00
83		Chrysene	40.000	39.617	1.0	144	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	38.300	4.3	137	0.00
85	c	Di-n-octyl phthalate	40.000	37.954	5.1	134	0.00
86	I	Perylene-d12	20.000	20.000	0.0	132	0.00
87		Indeno(1,2,3-cd)pyrene	40.000	40.356	-0.9	143	0.00
88		Benzo(b)fluoranthene	40.000	40.776	-1.9	143	0.00
89		Benzo(k)fluoranthene	40.000	40.109	-0.3	143	0.00
90	C	Benzo(a)pyrene	40.000	40.172	-0.4	141	0.00
91		Dibenzo(a,h)anthracene	40.000	40.324	-0.8	141	0.00
92		Benzo(g,h,i)perylene	40.000	39.685	0.8	139	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100820\
Data File : BG046811.D
Acq On : 8 Oct 2020 11:03
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Oct 08 13:13:41 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG092220.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 07 11:38:17 2020
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)

(#) = Out of Range SPCC's out = 0 CCC's out = 1