Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG041720\
Data File : BG045042.D
Acq On : 17 Apr 2020 21:47
Operator : CG/JU
Sample : SSTDCCC40
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC40

Quant Time: Apr 17 22:29:38 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG041520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Apr 17 11:08:29 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	97	0.00
2		1,4-Dioxane	40.000	34.756	13.1	84	0.00
3		Pyridine	40.000	35.047	12.4	88	0.00
4		n-Nitrosodimethylamine	40.000	41.512	-3.8	105	0.00
5	S	2-Fluorophenol	80.000	77.853	2.7	97	0.00
6		Aniline	40.000	41.238	-3.1	101	0.00
7	S	Phenol-d6	80.000	85.593	-7.0	105	0.00
8		2-Chlorophenol	40.000	42.198	-5.5	104	0.00
9		Benzaldehyde	40.000	38.211	4.5	94	0.00
10	C	Phenol	40.000	42.611	-6.5	104	0.00
11		bis(2-Chloroethyl)ether	40.000	40.751	-1.9	100	0.00
12		1,3-Dichlorobenzene	40.000	39.796	0.5	99	0.00
13	C	1,4-Dichlorobenzene	40.000	40.501	-1.3	100	0.00
14		1,2-Dichlorobenzene	40.000	41.035	-2.6	100	0.00
15		Benzyl Alcohol	40.000	43.916	-9.8	108	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	40.257	-0.6	98	0.00
17		2-Methylphenol	40.000	42.166	-5.4	105	0.00
18		Hexachloroethane	40.000	41.094	-2.7	100	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	43.094	-7.7	105	0.00
20		3+4-Methylphenols	40.000	43.067	-7.7	105	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	102	0.00
22		Acetophenone	40.000	39.265	1.8	104	0.00
23	S	Nitrobenzene-d5	80.000	80.003	-0.0	107	0.00
24		Nitrobenzene	40.000	39.799	0.5	106	0.00
25		Isophorone	40.000	40.238	-0.6	104	0.00
26	C	2-Nitrophenol	40.000	42.403	-6.0	112	0.00
27		2,4-Dimethylphenol	40.000	38.013	5.0	101	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.603	1.0	103	0.00
29	C	2,4-Dichlorophenol	40.000	41.667	-4.2	109	0.00
30		1,2,4-Trichlorobenzene	40.000	40.595	-1.5	108	0.00
31		Naphthalene	40.000	40.024	-0.1	104	0.00
32		Benzoic acid	40.000	45.994	-15.0	131	0.01
33		4-Chloroaniline	40.000	39.836	0.4	105	0.00
34	C	Hexachlorobutadiene	40.000	41.414	-3.5	110	0.00
35		Caprolactam	40.000	37.897	5.3	99	0.00
36	C	4-Chloro-3-methylphenol	40.000	41.625	-4.1	111	0.00
37		2-Methylnaphthalene	40.000	41.249	-3.1	108	0.00
38		1-Methylnaphthalene	40.000	41.318	-3.3	108	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	107	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	40.690	-1.7	111	0.00
41	P	Hexachlorocyclopentadiene	40.000	36.529	8.7	97	0.00
42	S	2,4,6-Tribromophenol	80.000	82.794	-3.5	113	0.00
43	C	2,4,6-Trichlorophenol	40.000	41.218	-3.0	111	0.00
44		2,4,5-Trichlorophenol	40.000	40.773	-1.9	111	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG041720\
Data File : BG045042.D
Acq On : 17 Apr 2020 21:47
Operator : CG/JU
Sample : SSTDCCC40
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC40

Quant Time: Apr 17 22:29:38 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG041520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Apr 17 11:08:29 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)
45	S	2-Fluorobiphenyl	80.000	81.052	-1.3	109	0.00
46		1,1'-Biphenyl	40.000	39.868	0.3	108	0.00
47		2-Chloronaphthalene	40.000	39.864	0.3	109	0.00
48		2-Nitroaniline	40.000	40.066	-0.2	111	0.00
49		Acenaphthylene	40.000	39.592	1.0	108	0.00
50		Dimethylphthalate	40.000	39.280	1.8	107	0.00
51		2,6-Dinitrotoluene	40.000	40.130	-0.3	110	0.00
52	C	Acenaphthene	40.000	40.365	-0.9	110	0.00
53		3-Nitroaniline	40.000	38.455	3.9	105	0.00
54	P	2,4-Dinitrophenol	40.000	43.961	-9.9	123	0.00
55		Dibenzofuran	40.000	39.846	0.4	109	0.00
56	P	4-Nitrophenol	40.000	39.429	1.4	108	0.00
57		2,4-Dinitrotoluene	40.000	39.423	1.4	110	0.00
58		Fluorene	40.000	40.296	-0.7	110	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	42.178	-5.4	116	0.00
60		Diethylphthalate	40.000	38.653	3.4	106	0.00
61		4-Chlorophenyl-phenylether	40.000	41.127	-2.8	113	0.00
62		4-Nitroaniline	40.000	37.720	5.7	104	0.00
63		Azobenzene	40.000	39.809	0.5	110	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	105	0.00
65		4,6-Dinitro-2-methylphenol	40.000	42.063	-5.2	116	0.00
66	c	n-Nitrosodiphenylamine	40.000	40.469	-1.2	108	0.00
67		4-Bromophenyl-phenylether	40.000	41.894	-4.7	112	0.00
68		Hexachlorobenzene	40.000	40.841	-2.1	111	0.00
69		Atrazine	40.000	37.845	5.4	99	0.00
70	C	Pentachlorophenol	40.000	42.644	-6.6	112	0.00
71		Phenanthrene	40.000	39.644	0.9	106	0.00
72		Anthracene	40.000	38.995	2.5	104	0.00
73		Carbazole	40.000	37.081	7.3	103	0.00
74		Di-n-butylphthalate	40.000	39.509	1.2	102	0.00
75	C	Fluoranthene	40.000	39.673	0.8	104	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	102	0.00
77		Benzidine	40.000	31.474	21.3	79	0.00
78		Pyrene	40.000	38.432	3.9	102	0.00
79	S	Terphenyl-d14	80.000	83.372	-4.2	104	0.00
80		Butylbenzylphthalate	40.000	39.739	0.7	103	0.00
81		Benzo(a)anthracene	40.000	40.523	-1.3	105	0.00
82		3,3'-Dichlorobenzidine	40.000	41.422	-3.6	107	0.00
83		Chrysene	40.000	40.553	-1.4	106	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	39.602	1.0	103	0.00
85	c	Di-n-octyl phthalate	40.000	40.080	-0.2	104	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	41.070	-2.7	109	0.00
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\BG041720\
Data File : BG045042.D
Acq On : 17 Apr 2020 21:47
Operator : CG/JU
Sample : SSTDCCC40
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC40

Quant Time: Apr 17 22:29:38 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG041520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Apr 17 11:08:29 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)
88	Benzo(b)fluoranthene	40.000	42.201	-5.5	109	0.00
89	Benzo(k)fluoranthene	40.000	41.515	-3.8	105	0.00
90 C	Benzo(a)pyrene	40.000	41.656	-4.1	107	0.00
91	Dibenzo(a,h)anthracene	40.000	41.695	-4.2	108	0.00
92	Benzo(g,h,i)perylene	40.000	41.737	-4.3	108	0.00

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