Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080119\
Data File : BF115909.D
Acq On : 1 Aug 2019 16:28
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 02 04:57:36 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 31 15:31:51 2019
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	125	0.00
2		1,4-Dioxane	40.000	39.332	1.7	126	0.00
3		Pyridine	40.000	39.437	1.4	125	0.02
4		n-Nitrosodimethylamine	40.000	40.924	-2.3	130	0.02
5	S	2-Fluorophenol	80.000	79.649	0.4	123	0.01
6		Aniline	40.000	40.315	-0.8	124	0.00
7	S	Phenol-d6	80.000	80.443	-0.6	126	0.00
8		2-Chlorophenol	40.000	40.248	-0.6	125	0.00
9		Benzaldehyde	40.000	38.127	4.7	118	0.00
10	C	Phenol	40.000	40.049	-0.1	124	0.01
11		bis(2-Chloroethyl)ether	40.000	39.884	0.3	124	0.00
12		1,3-Dichlorobenzene	40.000	39.235	1.9	122	0.00
13	C	1,4-Dichlorobenzene	40.000	39.627	0.9	123	0.00
14		1,2-Dichlorobenzene	40.000	39.843	0.4	121	0.00
15		Benzyl Alcohol	40.000	41.391	-3.5	126	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	39.239	1.9	121	0.00
17		2-Methylphenol	40.000	40.944	-2.4	129	0.00
18		Hexachloroethane	40.000	39.604	1.0	122	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	38.558	3.6	122	0.00
20		3+4-Methylphenols	40.000	39.497	1.3	120	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	125	0.00
22		Acetophenone	40.000	38.656	3.4	120	0.00
23	S	Nitrobenzene-d5	80.000	79.583	0.5	124	0.00
24		Nitrobenzene	40.000	39.619	1.0	123	0.00
25		Isophorone	40.000	39.445	1.4	123	0.00
26	C	2-Nitrophenol	40.000	41.395	-3.5	128	0.00
27		2,4-Dimethylphenol	40.000	39.823	0.4	123	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.319	1.7	123	0.00
29	C	2,4-Dichlorophenol	40.000	40.465	-1.2	124	0.00
30		1,2,4-Trichlorobenzene	40.000	39.753	0.6	123	0.00
31		Naphthalene	40.000	39.462	1.3	120	0.00
32		Benzoic acid	40.000	34.590	13.5	118	0.02
33		4-Chloroaniline	40.000	40.855	-2.1	126	0.00
34	C	Hexachlorobutadiene	40.000	38.833	2.9	120	0.00
35		Caprolactam	40.000	41.364	-3.4	128	0.02
36	C	4-Chloro-3-methylphenol	40.000	40.392	-1.0	126	0.00
37		2-Methylnaphthalene	40.000	38.707	3.2	119	0.00
38		1-Methylnaphthalene	40.000	39.096	2.3	120	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	123	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	39.583	1.0	120	0.00
41	P	Hexachlorocyclopentadiene	40.000	34.470	13.8	106	0.00
42	S	2,4,6-Tribromophenol	80.000	78.419	2.0	120	0.00
43	C	2,4,6-Trichlorophenol	40.000	40.183	-0.5	123	0.00
44		2,4,5-Trichlorophenol	40.000	40.294	-0.7	123	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080119\
Data File : BF115909.D
Acq On : 1 Aug 2019 16:28
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 02 04:57:36 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 31 15:31:51 2019
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	76.444	4.4	115	0.00
46		1,1'-Biphenyl	40.000	38.780	3.0	118	0.00
47		2-Chloronaphthalene	40.000	39.710	0.7	121	0.00
48		2-Nitroaniline	40.000	41.113	-2.8	126	0.00
49		Acenaphthylene	40.000	39.428	1.4	119	0.00
50		Dimethylphthalate	40.000	40.225	-0.6	123	0.00
51		2,6-Dinitrotoluene	40.000	40.338	-0.8	123	0.00
52	C	Acenaphthene	40.000	38.859	2.9	117	0.00
53		3-Nitroaniline	40.000	40.667	-1.7	125	0.00
54	P	2,4-Dinitrophenol	40.000	38.080	4.8	120	0.01
55		Dibenzofuran	40.000	39.474	1.3	118	0.00
56	P	4-Nitrophenol	40.000	38.763	3.1	118	0.01
57		2,4-Dinitrotoluene	40.000	40.056	-0.1	121	0.00
58		Fluorene	40.000	39.153	2.1	118	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	39.620	1.0	122	0.00
60		Diethylphthalate	40.000	39.266	1.8	119	0.00
61		4-Chlorophenyl-phenylether	40.000	39.145	2.1	117	0.00
62		4-Nitroaniline	40.000	40.157	-0.4	122	0.01
63		Azobenzene	40.000	39.044	2.4	118	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	121	0.00
65		4,6-Dinitro-2-methylphenol	40.000	40.737	-1.8	119	0.00
66	c	n-Nitrosodiphenylamine	40.000	39.879	0.3	120	0.00
67		4-Bromophenyl-phenylether	40.000	40.181	-0.5	120	0.00
68		Hexachlorobenzene	40.000	39.511	1.2	118	0.00
69		Atrazine	40.000	39.455	1.4	118	0.00
70	C	Pentachlorophenol	40.000	40.097	-0.2	117	0.00
71		Phenanthrene	40.000	39.662	0.8	118	0.00
72		Anthracene	40.000	39.544	1.1	118	0.00
73		Carbazole	40.000	40.196	-0.5	119	0.00
74		Di-n-butylphthalate	40.000	39.532	1.2	115	0.00
75	C	Fluoranthene	40.000	39.155	2.1	117	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	106	0.00
77		Benzidine	40.000	49.984	-25.0	124	0.00
78		Pyrene	40.000	42.939	-7.3	116	0.00
79	S	Terphenyl-d14	80.000	84.007	-5.0	113	0.00
80		Butylbenzylphthalate	40.000	42.243	-5.6	111	0.00
81		Benzo(a)anthracene	40.000	40.141	-0.4	106	0.00
82		3,3'-Dichlorobenzidine	40.000	41.528	-3.8	107	0.00
83		Chrysene	40.000	39.271	1.8	104	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	40.860	-2.1	106	0.00
85	c	Di-n-octyl phthalate	40.000	37.693	5.8	96	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	44.541	-11.4	123	0.02
87	I	Perylene-d12	20.000	20.000	0.0	103	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080119\
Data File : BF115909.D
Acq On : 1 Aug 2019 16:28
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 02 04:57:36 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 31 15:31:51 2019
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	38.566	3.6	95	0.00
89	Benzo(k)fluoranthene	40.000	39.051	2.4	104	0.00
90 C	Benzo(a)pyrene	40.000	39.564	1.1	102	0.00
91	Dibenzo(a,h)anthracene	40.000	45.392	-13.5	121	0.01
92	Benzo(g,h,i)perylene	40.000	48.062	-20.2	132	0.02

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