Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF073119\
Data File : BF115854.D
Acq On : 31 Jul 2019 7:59
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 31 18:20:24 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	109	0.00
2		1,4-Dioxane	40.000	39.854	0.4	111	0.00
3		Pyridine	40.000	39.388	1.5	109	0.00
4		n-Nitrosodimethylamine	40.000	41.106	-2.8	114	0.00
5	S	2-Fluorophenol	80.000	81.561	-2.0	110	0.00
6		Aniline	40.000	41.091	-2.7	110	0.00
7	S	Phenol-d6	80.000	81.807	-2.3	111	0.00
8		2-Chlorophenol	40.000	40.516	-1.3	109	0.00
9		Benzaldehyde	40.000	40.054	-0.1	108	0.00
10	C	Phenol	40.000	41.074	-2.7	111	0.00
11		bis(2-Chloroethyl)ether	40.000	40.645	-1.6	111	0.00
12		1,3-Dichlorobenzene	40.000	40.300	-0.7	109	0.00
13	C	1,4-Dichlorobenzene	40.000	40.375	-0.9	109	0.00
14		1,2-Dichlorobenzene	40.000	41.137	-2.8	109	0.00
15		Benzyl Alcohol	40.000	41.703	-4.3	110	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	40.392	-1.0	108	0.00
17		2-Methylphenol	40.000	40.804	-2.0	113	0.00
18		Hexachloroethane	40.000	40.424	-1.1	109	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.511	1.2	109	0.00
20		3+4-Methylphenols	40.000	41.255	-3.1	110	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	110	0.00
22		Acetophenone	40.000	40.258	-0.6	109	0.00
23	S	Nitrobenzene-d5	80.000	80.482	-0.6	109	0.00
24		Nitrobenzene	40.000	40.940	-2.3	111	0.00
25		Isophorone	40.000	39.295	1.8	107	0.00
26	C	2-Nitrophenol	40.000	40.158	-0.4	108	0.00
27		2,4-Dimethylphenol	40.000	40.620	-1.5	110	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.638	0.9	108	0.00
29	C	2,4-Dichlorophenol	40.000	40.624	-1.6	108	0.00
30		1,2,4-Trichlorobenzene	40.000	40.659	-1.6	109	0.00
31		Naphthalene	40.000	40.503	-1.3	108	0.00
32		Benzoic acid	40.000	36.390	9.0	109	0.00
33		4-Chloroaniline	40.000	40.536	-1.3	109	0.00
34	C	Hexachlorobutadiene	40.000	39.889	0.3	108	0.00
35		Caprolactam	40.000	40.012	-0.0	108	0.00
36	C	4-Chloro-3-methylphenol	40.000	39.935	0.2	109	0.00
37		2-Methylnaphthalene	40.000	39.747	0.6	106	0.00
38		1-Methylnaphthalene	40.000	39.858	0.4	107	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	107	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	40.662	-1.7	107	0.00
41	P	Hexachlorocyclopentadiene	40.000	39.478	1.3	107	0.00
42	S	2,4,6-Tribromophenol	80.000	79.648	0.4	105	0.00
43	C	2,4,6-Trichlorophenol	40.000	40.167	-0.4	106	0.00
44		2,4,5-Trichlorophenol	40.000	40.293	-0.7	106	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF073119\
Data File : BF115854.D
Acq On : 31 Jul 2019 7:59
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 31 18:20:24 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	80.820	-1.0	105	0.00
46		1,1'-Biphenyl	40.000	40.477	-1.2	106	0.00
47		2-Chloronaphthalene	40.000	40.864	-2.2	108	0.00
48		2-Nitroaniline	40.000	40.524	-1.3	108	0.00
49		Acenaphthylene	40.000	40.714	-1.8	106	0.00
50		Dimethylphthalate	40.000	39.235	1.9	104	0.00
51		2,6-Dinitrotoluene	40.000	39.770	0.6	105	0.00
52	C	Acenaphthene	40.000	40.148	-0.4	105	0.00
53		3-Nitroaniline	40.000	40.076	-0.2	106	0.00
54	P	2,4-Dinitrophenol	40.000	36.436	8.9	98	0.00
55		Dibenzofuran	40.000	40.811	-2.0	106	0.00
56	P	4-Nitrophenol	40.000	40.468	-1.2	106	0.00
57		2,4-Dinitrotoluene	40.000	40.205	-0.5	104	0.00
58		Fluorene	40.000	40.034	-0.1	104	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	40.887	-2.2	108	0.00
60		Diethylphthalate	40.000	39.282	1.8	103	0.00
61		4-Chlorophenyl-phenylether	40.000	40.334	-0.8	104	0.00
62		4-Nitroaniline	40.000	40.405	-1.0	106	0.00
63		Azobenzene	40.000	39.551	1.1	104	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	105	0.00
65		4,6-Dinitro-2-methylphenol	40.000	39.546	1.1	100	0.00
66	c	n-Nitrosodiphenylamine	40.000	40.516	-1.3	105	0.00
67		4-Bromophenyl-phenylether	40.000	39.978	0.1	103	0.00
68		Hexachlorobenzene	40.000	40.220	-0.5	103	0.00
69		Atrazine	40.000	39.006	2.5	101	0.00
70	C	Pentachlorophenol	40.000	39.887	0.3	101	0.00
71		Phenanthrene	40.000	40.267	-0.7	103	0.00
72		Anthracene	40.000	39.901	0.2	103	0.00
73		Carbazole	40.000	40.701	-1.8	104	0.00
74		Di-n-butylphthalate	40.000	38.944	2.6	97	0.00
75	C	Fluoranthene	40.000	39.802	0.5	103	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	102	0.00
77		Benzidine	40.000	42.558	-6.4	101	0.00
78		Pyrene	40.000	39.543	1.1	103	0.00
79	S	Terphenyl-d14	80.000	78.285	2.1	102	0.00
80		Butylbenzylphthalate	40.000	38.604	3.5	97	0.00
81		Benzo(a)anthracene	40.000	40.282	-0.7	103	0.00
82		3,3'-Dichlorobenzidine	40.000	41.309	-3.3	103	0.00
83		Chrysene	40.000	39.951	0.1	101	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	38.238	4.4	96	0.00
85	c	Di-n-octyl phthalate	40.000	38.280	4.3	94	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	36.811	8.0	98	0.00
87	I	Perylene-d12	20.000	20.000	0.0	98	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF073119\
Data File : BF115854.D
Acq On : 31 Jul 2019 7:59
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 31 18:20:24 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	39.741	0.6	93	0.00
89	Benzo(k)fluoranthene	40.000	41.754	-4.4	106	0.00
90 C	Benzo(a)pyrene	40.000	40.101	-0.3	98	0.00
91	Dibenzo(a,h)anthracene	40.000	38.565	3.6	97	0.00
92	Benzo(g,h,i)perylene	40.000	38.172	4.6	99	0.00

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