Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091319\
Data File : BF117026.D
Acq On : 13 Sep 2019 17:02
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 14 00:22:06 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 13 16:23:17 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	102	0.00
2		1,4-Dioxane	40.000	38.993	2.5	105	0.00
3		Pyridine	40.000	38.171	4.6	100	0.00
4		n-Nitrosodimethylamine	40.000	37.363	6.6	102	0.00
5	S	2-Fluorophenol	80.000	77.613	3.0	102	0.00
6		Aniline	40.000	38.842	2.9	103	0.00
7	S	Phenol-d6	80.000	77.085	3.6	103	0.00
8		2-Chlorophenol	40.000	38.105	4.7	101	0.00
9		Benzaldehyde	40.000	42.323	-5.8	102	0.00
10	C	Phenol	40.000	38.746	3.1	103	0.00
11		bis(2-Chloroethyl)ether	40.000	37.302	6.7	102	0.00
12		1,3-Dichlorobenzene	40.000	38.640	3.4	103	0.00
13	C	1,4-Dichlorobenzene	40.000	38.672	3.3	102	0.00
14		1,2-Dichlorobenzene	40.000	38.717	3.2	103	0.00
15		Benzyl Alcohol	40.000	38.217	4.5	100	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	37.769	5.6	101	0.00
17		2-Methylphenol	40.000	38.068	4.8	103	0.00
18		Hexachloroethane	40.000	38.640	3.4	103	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	37.713	5.7	103	0.00
20		3+4-Methylphenols	40.000	37.994	5.0	102	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	102	0.00
22		Acetophenone	40.000	38.178	4.6	104	0.00
23	S	Nitrobenzene-d5	80.000	77.623	3.0	103	0.00
24		Nitrobenzene	40.000	38.513	3.7	103	0.00
25		Isophorone	40.000	37.054	7.4	102	0.00
26	C	2-Nitrophenol	40.000	39.113	2.2	103	0.00
27		2,4-Dimethylphenol	40.000	37.963	5.1	102	0.00
28		bis(2-Chloroethoxy)methane	40.000	37.651	5.9	103	0.00
29	C	2,4-Dichlorophenol	40.000	37.916	5.2	100	0.00
30		1,2,4-Trichlorobenzene	40.000	37.515	6.2	101	0.00
31		Naphthalene	40.000	37.813	5.5	101	0.00
32		Benzoic acid	40.000	36.825	7.9	104	0.00
33		4-Chloroaniline	40.000	38.021	4.9	101	0.00
34	C	Hexachlorobutadiene	40.000	37.763	5.6	101	0.00
35		Caprolactam	40.000	36.740	8.1	100	0.00
36	C	4-Chloro-3-methylphenol	40.000	37.392	6.5	101	0.00
37		2-Methylnaphthalene	40.000	37.783	5.5	102	0.00
38		1-Methylnaphthalene	40.000	37.474	6.3	101	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	100	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	38.408	4.0	101	0.00
41	P	Hexachlorocyclopentadiene	40.000	36.981	7.5	105	0.00
42	S	2,4,6-Tribromophenol	80.000	75.873	5.2	99	0.00
43	C	2,4,6-Trichlorophenol	40.000	37.490	6.3	98	0.00
44		2,4,5-Trichlorophenol	40.000	37.408	6.5	100	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091319\
Data File : BF117026.D
Acq On : 13 Sep 2019 17:02
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 14 00:22:06 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 13 16:23:17 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	76.544	4.3	101	0.00
46		1,1'-Biphenyl	40.000	38.049	4.9	101	0.00
47		2-Chloronaphthalene	40.000	37.821	5.4	100	0.00
48		2-Nitroaniline	40.000	38.740	3.1	103	0.00
49		Acenaphthylene	40.000	37.933	5.2	99	0.00
50		Dimethylphthalate	40.000	36.882	7.8	97	0.00
51		2,6-Dinitrotoluene	40.000	39.360	1.6	104	0.00
52	C	Acenaphthene	40.000	38.086	4.8	100	0.00
53		3-Nitroaniline	40.000	38.387	4.0	100	0.00
54	P	2,4-Dinitrophenol	40.000	37.496	6.3	108	0.00
55		Dibenzofuran	40.000	38.126	4.7	100	0.00
56	P	4-Nitrophenol	40.000	38.485	3.8	99	0.00
57		2,4-Dinitrotoluene	40.000	40.447	-1.1	104	0.00
58		Fluorene	40.000	37.611	6.0	99	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	36.430	8.9	93	0.00
60		Diethylphthalate	40.000	36.696	8.3	96	0.00
61		4-Chlorophenyl-phenylether	40.000	37.336	6.7	97	0.00
62		4-Nitroaniline	40.000	38.157	4.6	99	0.00
63		Azobenzene	40.000	37.265	6.8	98	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	96	0.00
65		4,6-Dinitro-2-methylphenol	40.000	38.376	4.1	103	0.00
66	c	n-Nitrosodiphenylamine	40.000	37.904	5.2	98	0.00
67		4-Bromophenyl-phenylether	40.000	37.281	6.8	95	0.00
68		Hexachlorobenzene	40.000	37.443	6.4	98	0.00
69		Atrazine	40.000	38.044	4.9	95	0.00
70	C	Pentachlorophenol	40.000	38.270	4.3	97	0.00
71		Phenanthrene	40.000	38.190	4.5	97	0.00
72		Anthracene	40.000	38.530	3.7	97	0.00
73		Carbazole	40.000	37.929	5.2	95	0.00
74		Di-n-butylphthalate	40.000	37.611	6.0	95	0.00
75	C	Fluoranthene	40.000	37.949	5.1	95	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	92	0.00
77		Benzidine	40.000	39.098	2.3	93	0.00
78		Pyrene	40.000	37.194	7.0	96	0.00
79	S	Terphenyl-d14	80.000	74.442	6.9	95	0.00
80		Butylbenzylphthalate	40.000	35.600	11.0	90	0.00
81		Benzo(a)anthracene	40.000	37.520	6.2	93	0.00
82		3,3'-Dichlorobenzidine	40.000	36.954	7.6	90	0.00
83		Chrysene	40.000	37.238	6.9	93	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	34.950	12.6	89	0.00
85	c	Di-n-octyl phthalate	40.000	35.711	10.7	88	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	34.985	12.5	96	-0.02
87	I	Perylene-d12	20.000	20.000	0.0	86	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091319\
Data File : BF117026.D
Acq On : 13 Sep 2019 17:02
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 14 00:22:06 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 13 16:23:17 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	37.999	5.0	90	0.00
89	Benzo(k)fluoranthene	40.000	39.889	0.3	88	0.00
90 C	Benzo(a)pyrene	40.000	38.157	4.6	88	0.00
91	Dibenzo(a,h)anthracene	40.000	37.120	7.2	94	-0.01
92	Benzo(g,h,i)perylene	40.000	37.012	7.5	96	0.00

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