Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091719\
Data File : BF117094.D
Acq On : 17 Sep 2019 9:25
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 17 17:50:26 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	97	0.00
2		1,4-Dioxane	40.000	34.842	12.9	89	0.00
3		Pyridine	40.000	35.139	12.2	87	0.00
4		n-Nitrosodimethylamine	40.000	35.048	12.4	90	0.00
5	S	2-Fluorophenol	80.000	74.674	6.7	93	0.00
6		Aniline	40.000	37.324	6.7	93	0.00
7	S	Phenol-d6	80.000	75.045	6.2	95	0.01
8		2-Chlorophenol	40.000	38.449	3.9	96	0.00
9		Benzaldehyde	40.000	42.645	-6.6	97	0.00
10	C	Phenol	40.000	37.256	6.9	93	0.00
11		bis(2-Chloroethyl)ether	40.000	37.606	6.0	97	0.00
12		1,3-Dichlorobenzene	40.000	38.102	4.7	96	0.00
13	C	1,4-Dichlorobenzene	40.000	37.671	5.8	94	0.00
14		1,2-Dichlorobenzene	40.000	38.274	4.3	96	0.00
15		Benzyl Alcohol	40.000	39.533	1.2	98	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	39.001	2.5	98	0.00
17		2-Methylphenol	40.000	38.025	4.9	97	0.00
18		Hexachloroethane	40.000	38.087	4.8	96	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.390	1.5	101	0.00
20		3+4-Methylphenols	40.000	38.797	3.0	98	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	100	0.00
22		Acetophenone	40.000	37.657	5.9	100	0.00
23	S	Nitrobenzene-d5	80.000	75.640	5.4	99	0.00
24		Nitrobenzene	40.000	37.955	5.1	99	0.00
25		Isophorone	40.000	37.905	5.2	102	0.00
26	C	2-Nitrophenol	40.000	40.189	-0.5	104	0.00
27		2,4-Dimethylphenol	40.000	37.402	6.5	99	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.083	4.8	102	0.00
29	C	2,4-Dichlorophenol	40.000	39.198	2.0	101	0.00
30		1,2,4-Trichlorobenzene	40.000	37.246	6.9	98	0.00
31		Naphthalene	40.000	37.642	5.9	98	0.00
32		Benzoic acid	40.000	31.532	21.2	84	0.00
33		4-Chloroaniline	40.000	38.138	4.7	100	0.00
34	C	Hexachlorobutadiene	40.000	38.843	2.9	102	0.00
35		Caprolactam	40.000	35.916	10.2	96	0.01
36	C	4-Chloro-3-methylphenol	40.000	38.664	3.3	103	0.00
37		2-Methylnaphthalene	40.000	38.284	4.3	101	0.00
38		1-Methylnaphthalene	40.000	38.619	3.5	102	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	104	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	37.825	5.4	104	0.00
41	P	Hexachlorocyclopentadiene	40.000	42.194	-5.5	128	0.00
42	S	2,4,6-Tribromophenol	80.000	80.329	-0.4	109	0.00
43	C	2,4,6-Trichlorophenol	40.000	37.372	6.6	102	0.00
44		2,4,5-Trichlorophenol	40.000	35.789	10.5	100	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091719\
Data File : BF117094.D
Acq On : 17 Sep 2019 9:25
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 17 17:50:26 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	77.162	3.5	106	0.00
46		1,1'-Biphenyl	40.000	37.527	6.2	103	0.00
47		2-Chloronaphthalene	40.000	37.365	6.6	103	0.00
48		2-Nitroaniline	40.000	37.931	5.2	105	0.00
49		Acenaphthylene	40.000	37.510	6.2	102	0.00
50		Dimethylphthalate	40.000	38.365	4.1	105	0.00
51		2,6-Dinitrotoluene	40.000	40.216	-0.5	110	0.00
52	C	Acenaphthene	40.000	37.430	6.4	102	0.00
53		3-Nitroaniline	40.000	38.046	4.9	103	0.00
54	P	2,4-Dinitrophenol	40.000	39.103	2.2	118	0.01
55		Dibenzofuran	40.000	38.164	4.6	104	0.00
56	P	4-Nitrophenol	40.000	38.505	3.7	103	0.01
57		2,4-Dinitrotoluene	40.000	41.394	-3.5	111	0.00
58		Fluorene	40.000	38.220	4.5	104	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	38.513	3.7	102	0.00
60		Diethylphthalate	40.000	39.386	1.5	107	0.00
61		4-Chlorophenyl-phenylether	40.000	40.162	-0.4	108	0.00
62		4-Nitroaniline	40.000	37.838	5.4	102	0.00
63		Azobenzene	40.000	38.581	3.5	106	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	102	0.00
65		4,6-Dinitro-2-methylphenol	40.000	39.798	0.5	115	0.00
66	c	n-Nitrosodiphenylamine	40.000	38.246	4.4	105	0.00
67		4-Bromophenyl-phenylether	40.000	39.723	0.7	108	0.00
68		Hexachlorobenzene	40.000	39.032	2.4	109	0.00
69		Atrazine	40.000	38.639	3.4	102	0.00
70	C	Pentachlorophenol	40.000	33.518	16.2	91	0.00
71		Phenanthrene	40.000	37.376	6.6	101	0.00
72		Anthracene	40.000	38.690	3.3	104	0.00
73		Carbazole	40.000	37.143	7.1	99	0.00
74		Di-n-butylphthalate	40.000	40.999	-2.5	110	0.00
75	C	Fluoranthene	40.000	37.200	7.0	99	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	94	0.00
77		Benzidine	40.000	40.109	-0.3	98	0.00
78		Pyrene	40.000	37.372	6.6	98	0.00
79	S	Terphenyl-d14	80.000	79.202	1.0	103	0.00
80		Butylbenzylphthalate	40.000	40.939	-2.3	106	0.00
81		Benzo(a)anthracene	40.000	37.580	6.1	95	0.00
82		3,3'-Dichlorobenzidine	40.000	38.435	3.9	96	0.00
83		Chrysene	40.000	37.447	6.4	95	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	44.471	-11.2	115	0.00
85	c	Di-n-octyl phthalate	40.000	44.920	-12.3	113	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	33.232	16.9	93	0.01
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\BF091719\
Data File : BF117094.D
Acq On : 17 Sep 2019 9:25
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 17 17:50:26 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	37.485	6.3	88	0.00
89	Benzo(k)fluoranthene	40.000	40.376	-0.9	89	0.00
90 C	Benzo(a)pyrene	40.000	38.318	4.2	87	0.00
91	Dibenzo(a,h)anthracene	40.000	37.481	6.3	95	0.01
92	Benzo(g,h,i)perylene	40.000	35.167	12.1	90	0.02

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