Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091619\
Data File : BF117073.D
Acq On : 16 Sep 2019 21:05
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 17 11:45:20 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	64	0.00
2		1,4-Dioxane	40.000	38.001	5.0	64	-0.04
3		Pyridine	40.000	37.935	5.2	62	-0.02
4		n-Nitrosodimethylamine	40.000	37.474	6.3	64	-0.04
5	S	2-Fluorophenol	80.000	77.937	2.6	64	0.00
6		Aniline	40.000	38.886	2.8	65	0.00
7	S	Phenol-d6	80.000	77.773	2.8	65	0.01
8		2-Chlorophenol	40.000	38.534	3.7	64	0.00
9		Benzaldehyde	40.000	46.234	-15.6	69	0.00
10	C	Phenol	40.000	39.187	2.0	65	0.00
11		bis(2-Chloroethyl)ether	40.000	37.205	7.0	64	0.00
12		1,3-Dichlorobenzene	40.000	39.323	1.7	66	0.00
13	C	1,4-Dichlorobenzene	40.000	38.254	4.4	63	0.00
14		1,2-Dichlorobenzene	40.000	38.785	3.0	65	0.00
15		Benzyl Alcohol	40.000	37.615	6.0	62	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	36.264	9.3	61	0.00
17		2-Methylphenol	40.000	38.615	3.5	66	0.00
18		Hexachloroethane	40.000	24.173	39.6#	41	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	37.847	5.4	65	0.00
20		3+4-Methylphenols	40.000	41.147	-2.9	70	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	64	0.00
22		Acetophenone	40.000	39.263	1.8	67	0.00
23	S	Nitrobenzene-d5	80.000	77.544	3.1	65	0.00
24		Nitrobenzene	40.000	38.396	4.0	65	0.00
25		Isophorone	40.000	36.806	8.0	64	0.00
26	C	2-Nitrophenol	40.000	21.704	45.7#	36	0.00
27		2,4-Dimethylphenol	40.000	37.074	7.3	63	0.00
28		bis(2-Chloroethoxy)methane	40.000	37.158	7.1	64	0.00
29	C	2,4-Dichlorophenol	40.000	36.103	9.7	60	0.00
30		1,2,4-Trichlorobenzene	40.000	37.217	7.0	63	0.00
31		Naphthalene	40.000	37.614	6.0	63	0.00
32		Benzoic acid	40.000	23.030	42.4#	36	-0.03
33		4-Chloroaniline	40.000	37.248	6.9	62	0.00
34	C	Hexachlorobutadiene	40.000	37.480	6.3	63	0.00
35		Caprolactam	40.000	35.217	12.0	60	0.00
36	C	4-Chloro-3-methylphenol	40.000	35.848	10.4	61	0.00
37		2-Methylnaphthalene	40.000	36.416	9.0	62	0.00
38		1-Methylnaphthalene	40.000	36.659	8.4	62	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	57	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	41.095	-2.7	62	0.00
41	P	Hexachlorocyclopentadiene	40.000	0.000	100.0#	0	-8.97#
42	S	2,4,6-Tribromophenol	80.000	66.784	16.5	49	0.00
43	C	2,4,6-Trichlorophenol	40.000	37.149	7.1	55	0.00
44		2,4,5-Trichlorophenol	40.000	36.103	9.7	55	0.01

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091619\
Data File : BF117073.D
Acq On : 16 Sep 2019 21:05
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 17 11:45:20 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	83.526	-4.4	63	0.00
46		1,1'-Biphenyl	40.000	41.146	-2.9	62	0.00
47		2-Chloronaphthalene	40.000	39.712	0.7	60	0.00
48		2-Nitroaniline	40.000	43.452	-8.6	66	0.00
49		Acenaphthylene	40.000	38.249	4.4	57	0.00
50		Dimethylphthalate	40.000	39.843	0.4	60	0.00
51		2,6-Dinitrotoluene	40.000	32.519	18.7	49	0.00
52	C	Acenaphthene	40.000	37.863	5.3	56	0.00
53		3-Nitroaniline	40.000	43.666	-9.2	65	0.00
54	P	2,4-Dinitrophenol	40.000	0.000	100.0#	0	-9.91#
55		Dibenzofuran	40.000	37.552	6.1	56	0.00
56	P	4-Nitrophenol	40.000	26.172	34.6#	38	0.01
57		2,4-Dinitrotoluene	40.000	29.058	27.4#	42	0.00
58		Fluorene	40.000	35.994	10.0	54	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	34.615	13.5	50	0.00
60		Diethylphthalate	40.000	38.605	3.5	58	0.00
61		4-Chlorophenyl-phenylether	40.000	38.356	4.1	56	0.00
62		4-Nitroaniline	40.000	41.533	-3.8	61	0.00
63		Azobenzene	40.000	35.611	11.0	53	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	48	0.00
65		4,6-Dinitro-2-methylphenol	40.000	10.116	74.7#	5	-0.28
66	c	n-Nitrosodiphenylamine	40.000	41.602	-4.0	54	0.00
67		4-Bromophenyl-phenylether	40.000	40.419	-1.0	52	0.00
68		Hexachlorobenzene	40.000	38.233	4.4	50	0.00
69		Atrazine	40.000	35.752	10.6	46	0.00
70	C	Pentachlorophenol	40.000	34.669	13.3	45	0.00
71		Phenanthrene	40.000	38.563	3.6	49	0.00
72		Anthracene	40.000	39.075	2.3	50	0.00
73		Carbazole	40.000	39.154	2.1	50	0.00
74		Di-n-butylphthalate	40.000	40.237	-0.6	51	0.00
75	C	Fluoranthene	40.000	40.220	-0.5	51	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	48	0.00
77		Benzidine	40.000	51.480	-28.7#	65	0.00
78		Pyrene	40.000	38.419	4.0	52	0.00
79	S	Terphenyl-d14	80.000	80.670	-0.8	54	0.00
80		Butylbenzylphthalate	40.000	39.849	0.4	53	0.00
81		Benzo(a)anthracene	40.000	37.763	5.6	49	0.00
82		3,3'-Dichlorobenzidine	40.000	43.343	-8.4	56	0.00
83		Chrysene	40.000	37.888	5.3	49	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	41.249	-3.1	55	0.00
85	c	Di-n-octyl phthalate	40.000	43.405	-8.5	56	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	30.650	23.4	44	0.00
87	I	Perylene-d12	20.000	20.000	0.0	36	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091619\
Data File : BF117073.D
Acq On : 16 Sep 2019 21:05
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 17 11:45:20 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	36.625	8.4	37	0.00
89	Benzo(k)fluoranthene	40.000	43.082	-7.7	40	0.00
90 C	Benzo(a)pyrene	40.000	37.832	5.4	37	0.00
91	Dibenzo(a,h)anthracene	40.000	40.924	-2.3	44	0.00
92	Benzo(g,h,i)perylene	40.000	40.092	-0.2	44	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 1