Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122819\
Data File : BF118357.D
Acq On : 28 Dec 2019 11:13
Operator : JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 30 12:42:57 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF122419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Dec 24 15:58:52 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	128	0.00
2		1,4-Dioxane	40.000	36.551	8.6	122	0.02
3		Pyridine	40.000	37.350	6.6	123	0.02
4		n-Nitrosodimethylamine	40.000	37.134	7.2	120	0.03
5	S	2-Fluorophenol	80.000	74.728	6.6	123	0.01
6		Aniline	40.000	37.327	6.7	123	0.00
7	S	Phenol-d6	80.000	74.689	6.6	125	0.01
8		2-Chlorophenol	40.000	38.034	4.9	126	0.00
9		Benzaldehyde	40.000	34.756	13.1	120	0.00
10	C	Phenol	40.000	37.015	7.5	123	0.01
11		bis(2-Chloroethyl)ether	40.000	38.272	4.3	127	0.00
12		1,3-Dichlorobenzene	40.000	37.677	5.8	125	0.00
13	C	1,4-Dichlorobenzene	40.000	37.358	6.6	123	0.00
14		1,2-Dichlorobenzene	40.000	37.550	6.1	124	0.00
15		Benzyl Alcohol	40.000	38.040	4.9	125	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	39.026	2.4	130	0.00
17		2-Methylphenol	40.000	37.885	5.3	125	0.00
18		Hexachloroethane	40.000	37.680	5.8	125	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	37.611	6.0	126	0.00
20		3+4-Methylphenols	40.000	37.584	6.0	122	0.01
21	I	Naphthalene-d8	20.000	20.000	0.0	123	0.00
22		Acetophenone	40.000	37.664	5.8	122	0.00
23	S	Nitrobenzene-d5	80.000	75.864	5.2	124	0.00
24		Nitrobenzene	40.000	37.421	6.4	122	0.00
25		Isophorone	40.000	37.842	5.4	125	0.00
26	C	2-Nitrophenol	40.000	39.387	1.5	128	0.00
27		2,4-Dimethylphenol	40.000	37.383	6.5	123	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.760	3.1	127	0.00
29	C	2,4-Dichlorophenol	40.000	38.110	4.7	125	0.00
30		1,2,4-Trichlorobenzene	40.000	38.072	4.8	125	0.00
31		Naphthalene	40.000	37.889	5.3	123	0.00
32		Benzoic acid	40.000	34.739	13.2	115	0.02
33		4-Chloroaniline	40.000	37.959	5.1	124	0.00
34	C	Hexachlorobutadiene	40.000	38.403	4.0	125	0.00
35		Caprolactam	40.000	37.468	6.3	120	0.02
36	C	4-Chloro-3-methylphenol	40.000	37.702	5.7	124	0.00
37		2-Methylnaphthalene	40.000	37.770	5.6	122	0.00
38		1-Methylnaphthalene	40.000	37.769	5.6	124	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	123	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	38.150	4.6	124	0.00
41	P	Hexachlorocyclopentadiene	40.000	42.114	-5.3	152	0.00
42	S	2,4,6-Tribromophenol	80.000	74.679	6.7	119	0.01
43	C	2,4,6-Trichlorophenol	40.000	37.962	5.1	123	0.01
44		2,4,5-Trichlorophenol	40.000	37.275	6.8	120	0.01

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122819\
Data File : BF118357.D
Acq On : 28 Dec 2019 11:13
Operator : JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 30 12:42:57 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF122419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Dec 24 15:58:52 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	75.577	5.5	120	0.00
46		1,1'-Biphenyl	40.000	38.188	4.5	123	0.00
47		2-Chloronaphthalene	40.000	38.118	4.7	121	0.00
48		2-Nitroaniline	40.000	38.335	4.2	124	0.00
49		Acenaphthylene	40.000	37.387	6.5	119	0.00
50		Dimethylphthalate	40.000	37.783	5.5	123	0.00
51		2,6-Dinitrotoluene	40.000	38.096	4.8	123	0.01
52	C	Acenaphthene	40.000	37.764	5.6	121	0.01
53		3-Nitroaniline	40.000	37.351	6.6	118	0.00
54	P	2,4-Dinitrophenol	40.000	34.106	14.7	112	0.01
55		Dibenzofuran	40.000	37.115	7.2	118	0.00
56	P	4-Nitrophenol	40.000	32.881	17.8	105	0.02
57		2,4-Dinitrotoluene	40.000	37.565	6.1	121	0.00
58		Fluorene	40.000	37.216	7.0	118	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	36.237	9.4	115	0.01
60		Diethylphthalate	40.000	37.676	5.8	121	0.00
61		4-Chlorophenyl-phenylether	40.000	37.608	6.0	119	0.00
62		4-Nitroaniline	40.000	37.526	6.2	118	0.01
63		Azobenzene	40.000	37.241	6.9	120	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	120	0.00
65		4,6-Dinitro-2-methylphenol	40.000	38.083	4.8	118	0.01
66	c	n-Nitrosodiphenylamine	40.000	38.135	4.7	120	0.00
67		4-Bromophenyl-phenylether	40.000	37.986	5.0	120	0.00
68		Hexachlorobenzene	40.000	38.082	4.8	121	0.01
69		Atrazine	40.000	39.033	2.4	118	0.00
70	C	Pentachlorophenol	40.000	33.903	15.2	105	0.01
71		Phenanthrene	40.000	37.192	7.0	117	0.01
72		Anthracene	40.000	37.505	6.2	118	0.01
73		Carbazole	40.000	37.684	5.8	119	0.00
74		Di-n-butylphthalate	40.000	38.926	2.7	121	0.00
75	C	Fluoranthene	40.000	37.170	7.1	115	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	114	0.00
77		Benzidine	40.000	44.173	-10.4	121	0.00
78		Pyrene	40.000	37.706	5.7	117	0.00
79	S	Terphenyl-d14	80.000	74.870	6.4	114	0.00
80		Butylbenzylphthalate	40.000	40.580	-1.4	123	0.00
81		Benzo(a)anthracene	40.000	37.937	5.2	115	0.00
82		3,3'-Dichlorobenzidine	40.000	39.103	2.2	113	0.00
83		Chrysene	40.000	36.840	7.9	111	0.01
84		Bis(2-ethylhexyl)phthalate	40.000	41.396	-3.5	125	0.00
85	c	Di-n-octyl phthalate	40.000	40.678	-1.7	123	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	32.687	18.3	106	0.04
87	I	Perylene-d12	20.000	20.000	0.0	106	0.01

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122819\
Data File : BF118357.D
Acq On : 28 Dec 2019 11:13
Operator : JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 30 12:42:57 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF122419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Dec 24 15:58:52 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	37.925	5.2	108	0.01
89	Benzo(k)fluoranthene	40.000	40.027	-0.1	108	0.01
90 C	Benzo(a)pyrene	40.000	38.167	4.6	107	0.02
91	Dibenzo(a,h)anthracene	40.000	37.364	6.6	107	0.02
92	Benzo(g,h,i)perylene	40.000	36.367	9.1	104	0.04

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