Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF080919\
Data File : BF116099.D
Acq On : 9 Aug 2019 8:46
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 10 02:20:03 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 07 11:13:18 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	124	0.00
2		1,4-Dioxane	40.000	38.306	4.2	121	-0.01
3		Pyridine	40.000	37.735	5.7	118	-0.01
4		n-Nitrosodimethylamine	40.000	39.822	0.4	125	-0.01
5	S	2-Fluorophenol	80.000	82.684	-3.4	126	0.00
6		Aniline	40.000	39.651	0.9	121	0.00
7	S	Phenol-d6	80.000	83.386	-4.2	129	0.00
8		2-Chlorophenol	40.000	43.169	-7.9	132	0.00
9		Benzaldehyde	40.000	37.538	6.2	115	0.00
10	C	Phenol	40.000	40.646	-1.6	125	0.00
11		bis(2-Chloroethyl)ether	40.000	43.637	-9.1	134	0.00
12		1,3-Dichlorobenzene	40.000	40.821	-2.1	125	0.00
13	C	1,4-Dichlorobenzene	40.000	41.489	-3.7	127	0.00
14		1,2-Dichlorobenzene	40.000	40.584	-1.5	122	0.00
15		Benzyl Alcohol	40.000	41.111	-2.8	123	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	41.929	-4.8	127	0.00
17		2-Methylphenol	40.000	42.382	-6.0	132	0.00
18		Hexachloroethane	40.000	42.502	-6.3	129	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	41.967	-4.9	131	0.00
20		3+4-Methylphenols	40.000	41.594	-4.0	125	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	127	0.00
22		Acetophenone	40.000	40.425	-1.1	127	0.00
23	S	Nitrobenzene-d5	80.000	81.867	-2.3	129	0.00
24		Nitrobenzene	40.000	41.506	-3.8	131	0.00
25		Isophorone	40.000	42.637	-6.6	135	0.00
26	C	2-Nitrophenol	40.000	44.699	-11.7	140	0.00
27		2,4-Dimethylphenol	40.000	40.608	-1.5	128	0.00
28		bis(2-Chloroethoxy)methane	40.000	42.784	-7.0	135	0.00
29	C	2,4-Dichlorophenol	40.000	42.416	-6.0	131	0.00
30		1,2,4-Trichlorobenzene	40.000	41.299	-3.2	129	0.00
31		Naphthalene	40.000	41.171	-2.9	127	0.00
32		Benzoic acid	40.000	40.911	-2.3	142	0.00
33		4-Chloroaniline	40.000	40.415	-1.0	126	0.00
34	C	Hexachlorobutadiene	40.000	41.742	-4.4	131	0.00
35		Caprolactam	40.000	41.729	-4.3	131	0.00
36	C	4-Chloro-3-methylphenol	40.000	42.520	-6.3	134	0.00
37		2-Methylnaphthalene	40.000	40.865	-2.2	127	0.00
38		1-Methylnaphthalene	40.000	41.151	-2.9	128	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	125	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	41.818	-4.5	129	0.00
41	P	Hexachlorocyclopentadiene	40.000	38.848	2.9	123	0.00
42	S	2,4,6-Tribromophenol	80.000	81.661	-2.1	126	0.00
43	C	2,4,6-Trichlorophenol	40.000	40.017	-0.0	124	0.00
44		2,4,5-Trichlorophenol	40.000	41.042	-2.6	127	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF080919\
Data File : BF116099.D
Acq On : 9 Aug 2019 8:46
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 10 02:20:03 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 07 11:13:18 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	79.272	0.9	121	0.00
46		1,1'-Biphenyl	40.000	40.972	-2.4	126	0.00
47		2-Chloronaphthalene	40.000	40.326	-0.8	125	0.00
48		2-Nitroaniline	40.000	41.249	-3.1	128	0.00
49		Acenaphthylene	40.000	40.277	-0.7	123	0.00
50		Dimethylphthalate	40.000	41.738	-4.3	130	0.00
51		2,6-Dinitrotoluene	40.000	42.140	-5.4	130	0.00
52	C	Acenaphthene	40.000	40.099	-0.2	123	0.00
53		3-Nitroaniline	40.000	40.765	-1.9	127	0.00
54	P	2,4-Dinitrophenol	40.000	39.480	1.3	128	0.00
55		Dibenzofuran	40.000	38.806	3.0	118	0.00
56	P	4-Nitrophenol	40.000	34.379	14.1	106	0.00
57		2,4-Dinitrotoluene	40.000	38.779	3.1	118	0.00
58		Fluorene	40.000	40.960	-2.4	125	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	38.589	3.5	120	0.00
60		Diethylphthalate	40.000	41.321	-3.3	127	0.00
61		4-Chlorophenyl-phenylether	40.000	40.936	-2.3	124	0.00
62		4-Nitroaniline	40.000	40.770	-1.9	126	0.00
63		Azobenzene	40.000	41.675	-4.2	128	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	122	0.00
65		4,6-Dinitro-2-methylphenol	40.000	46.669	-16.7	137	0.00
66	c	n-Nitrosodiphenylamine	40.000	41.802	-4.5	126	0.00
67		4-Bromophenyl-phenylether	40.000	42.749	-6.9	128	0.00
68		Hexachlorobenzene	40.000	41.477	-3.7	124	0.00
69		Atrazine	40.000	33.074	17.3	100	0.00
70	C	Pentachlorophenol	40.000	35.364	11.6	104	0.00
71		Phenanthrene	40.000	40.738	-1.8	122	0.00
72		Anthracene	40.000	40.904	-2.3	123	0.00
73		Carbazole	40.000	41.664	-4.2	124	0.00
74		Di-n-butylphthalate	40.000	42.461	-6.2	124	0.00
75	C	Fluoranthene	40.000	40.365	-0.9	121	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	111	0.00
77		Benzidine	40.000	40.912	-2.3	106	0.00
78		Pyrene	40.000	42.319	-5.8	119	0.00
79	S	Terphenyl-d14	80.000	85.001	-6.3	120	0.00
80		Butylbenzylphthalate	40.000	45.491	-13.7	124	0.00
81		Benzo(a)anthracene	40.000	43.711	-9.3	121	0.00
82		3,3'-Dichlorobenzidine	40.000	44.309	-10.8	119	0.00
83		Chrysene	40.000	42.832	-7.1	118	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	46.773	-16.9	127	0.00
85	c	Di-n-octyl phthalate	40.000	45.238	-13.1	121	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	43.742	-9.4	126	0.00
87	I	Perylene-d12	20.000	20.000	0.0	113	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF080919\
Data File : BF116099.D
Acq On : 9 Aug 2019 8:46
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 10 02:20:03 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 07 11:13:18 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	42.082	-5.2	113	0.00
89	Benzo(k)fluoranthene	40.000	40.009	-0.0	117	0.00
90 C	Benzo(a)pyrene	40.000	41.746	-4.4	117	0.00
91	Dibenzo(a,h)anthracene	40.000	42.497	-6.2	124	0.00
92	Benzo(g,h,i)perylene	40.000	43.553	-8.9	131	0.01

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