Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF082019\
Data File : BF116437.D
Acq On : 20 Aug 2019 18:34
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 21 01:23:55 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 19 06:03:31 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	78	-0.01
2		1,4-Dioxane	40.000	40.742	-1.9	81	-0.02
3		Pyridine	40.000	37.036	7.4	73	-0.02
4		n-Nitrosodimethylamine	40.000	37.459	6.4	73	-0.02
5	S	2-Fluorophenol	80.000	86.911	-8.6	83	0.00
6		Aniline	40.000	39.638	0.9	76	-0.01
7	S	Phenol-d6	80.000	85.377	-6.7	83	0.00
8		2-Chlorophenol	40.000	44.368	-10.9	85	-0.01
9		Benzaldehyde	40.000	37.643	5.9	72	-0.01
10	C	Phenol	40.000	41.756	-4.4	80	0.00
11		bis(2-Chloroethyl)ether	40.000	43.255	-8.1	84	-0.01
12		1,3-Dichlorobenzene	40.000	42.372	-5.9	82	-0.01
13	C	1,4-Dichlorobenzene	40.000	43.420	-8.6	83	-0.01
14		1,2-Dichlorobenzene	40.000	43.351	-8.4	82	-0.01
15		Benzyl Alcohol	40.000	41.121	-2.8	77	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	42.672	-6.7	81	-0.01
17		2-Methylphenol	40.000	42.218	-5.5	83	0.00
18		Hexachloroethane	40.000	42.849	-7.1	82	-0.02
19	P	n-Nitroso-di-n-propylamine	40.000	45.234	-13.1	88	-0.01
20		3+4-Methylphenols	40.000	47.763	-19.4	90	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	81	-0.01
22		Acetophenone	40.000	43.437	-8.6	87	-0.01
23	S	Nitrobenzene-d5	80.000	81.799	-2.2	83	-0.01
24		Nitrobenzene	40.000	41.962	-4.9	85	-0.01
25		Isophorone	40.000	42.857	-7.1	87	-0.01
26	C	2-Nitrophenol	40.000	43.533	-8.8	87	-0.01
27		2,4-Dimethylphenol	40.000	41.516	-3.8	83	0.00
28		bis(2-Chloroethoxy)methane	40.000	43.239	-8.1	88	-0.01
29	C	2,4-Dichlorophenol	40.000	43.167	-7.9	86	-0.01
30		1,2,4-Trichlorobenzene	40.000	41.561	-3.9	83	-0.02
31		Naphthalene	40.000	42.905	-7.3	85	-0.01
32		Benzoic acid	40.000	32.553	18.6	72	0.00
33		4-Chloroaniline	40.000	41.601	-4.0	83	-0.01
34	C	Hexachlorobutadiene	40.000	42.021	-5.1	84	-0.01
35		Caprolactam	40.000	40.734	-1.8	82	-0.01
36	C	4-Chloro-3-methylphenol	40.000	43.837	-9.6	89	0.00
37		2-Methylnaphthalene	40.000	42.649	-6.6	85	-0.01
38		1-Methylnaphthalene	40.000	43.061	-7.7	86	-0.01
39	I	Acenaphthene-d10	20.000	20.000	0.0	83	-0.01
40		1,2,4,5-Tetrachlorobenzene	40.000	41.750	-4.4	86	-0.01
41	P	Hexachlorocyclopentadiene	40.000	38.166	4.6	80	-0.02
42	S	2,4,6-Tribromophenol	80.000	82.037	-2.5	84	-0.01
43	C	2,4,6-Trichlorophenol	40.000	36.304	9.2	75	-0.01
44		2,4,5-Trichlorophenol	40.000	34.732	13.2	71	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF082019\
Data File : BF116437.D
Acq On : 20 Aug 2019 18:34
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 21 01:23:55 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 19 06:03:31 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	83.530	-4.4	85	-0.01
46		1,1'-Biphenyl	40.000	42.274	-5.7	87	-0.02
47		2-Chloronaphthalene	40.000	41.493	-3.7	85	-0.01
48		2-Nitroaniline	40.000	41.041	-2.6	85	-0.01
49		Acenaphthylene	40.000	41.923	-4.8	86	-0.02
50		Dimethylphthalate	40.000	41.512	-3.8	86	-0.01
51		2,6-Dinitrotoluene	40.000	42.024	-5.1	86	0.00
52	C	Acenaphthene	40.000	41.611	-4.0	85	-0.02
53		3-Nitroaniline	40.000	40.542	-1.4	84	-0.01
54	P	2,4-Dinitrophenol	40.000	31.330	21.7	63	0.00
55		Dibenzofuran	40.000	40.582	-1.5	82	-0.01
56	P	4-Nitrophenol	40.000	66.912	-67.3#	137	0.00
57		2,4-Dinitrotoluene	40.000	39.708	0.7	81	0.00
58		Fluorene	40.000	43.799	-9.5	89	-0.01
59		2,3,4,6-Tetrachlorophenol	40.000	36.097	9.8	75	-0.01
60		Diethylphthalate	40.000	43.804	-9.5	89	-0.01
61		4-Chlorophenyl-phenylether	40.000	44.229	-10.6	89	-0.01
62		4-Nitroaniline	40.000	37.881	5.3	78	0.00
63		Azobenzene	40.000	43.540	-8.8	89	-0.02
64	I	Phenanthrene-d10	20.000	20.000	0.0	83	-0.02
65		4,6-Dinitro-2-methylphenol	40.000	42.105	-5.3	85	0.00
66	c	n-Nitrosodiphenylamine	40.000	43.046	-7.6	89	-0.01
67		4-Bromophenyl-phenylether	40.000	43.057	-7.6	88	-0.01
68		Hexachlorobenzene	40.000	41.510	-3.8	85	-0.02
69		Atrazine	40.000	36.658	8.4	76	-0.01
70	C	Pentachlorophenol	40.000	34.476	13.8	69	-0.01
71		Phenanthrene	40.000	42.120	-5.3	86	-0.01
72		Anthracene	40.000	43.209	-8.0	89	-0.01
73		Carbazole	40.000	43.372	-8.4	88	-0.01
74		Di-n-butylphthalate	40.000	46.396	-16.0	92	-0.01
75	C	Fluoranthene	40.000	43.448	-8.6	89	-0.01
76	I	Chrysene-d12	20.000	20.000	0.0	83	-0.01
77		Benzidine	40.000	40.094	-0.2	78	-0.02
78		Pyrene	40.000	41.097	-2.7	87	-0.01
79	S	Terphenyl-d14	80.000	83.118	-3.9	88	-0.01
80		Butylbenzylphthalate	40.000	44.653	-11.6	92	-0.02
81		Benzo(a)anthracene	40.000	42.755	-6.9	89	-0.01
82		3,3'-Dichlorobenzidine	40.000	43.374	-8.4	88	-0.01
83		Chrysene	40.000	43.067	-7.7	89	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	48.049	-20.1	98	-0.02
85	c	Di-n-octyl phthalate	40.000	48.241	-20.6#	97	-0.01
86		Indeno(1,2,3-cd)pyrene	40.000	42.560	-6.4	93	-0.01
87	I	Perylene-d12	20.000	20.000	0.0	90	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF082019\
Data File : BF116437.D
Acq On : 20 Aug 2019 18:34
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 21 01:23:55 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 19 06:03:31 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	41.019	-2.5	88	-0.01
89	Benzo(k)fluoranthene	40.000	41.092	-2.7	96	0.00
90 C	Benzo(a)pyrene	40.000	41.436	-3.6	93	-0.02
91	Dibenzo(a,h)anthracene	40.000	40.018	-0.0	93	-0.02
92	Benzo(g,h,i)perylene	40.000	38.617	3.5	92	-0.02

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