Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM092019\
Data File : BM022732.D
Acq On : 20 Sep 2019 19:06
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Sep 21 02:12:26 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM092019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 20 16:34:01 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	99	0.00
2		1,4-Dioxane	40.000	38.957	2.6	99	0.00
3		Pyridine	40.000	41.395	-3.5	98	0.00
4		n-Nitrosodimethylamine	40.000	41.595	-4.0	106	0.00
5	S	2-Fluorophenol	80.000	79.559	0.6	100	0.00
6		Aniline	40.000	40.616	-1.5	103	0.00
7	S	Phenol-d6	80.000	80.637	-0.8	102	0.00
8		2-Chlorophenol	40.000	40.032	-0.1	101	0.00
9		Benzaldehyde	40.000	44.650	-11.6	97	0.00
10	C	Phenol	40.000	40.416	-1.0	102	0.00
11		bis(2-Chloroethyl)ether	40.000	40.723	-1.8	103	0.00
12		1,3-Dichlorobenzene	40.000	39.373	1.6	100	0.00
13	C	1,4-Dichlorobenzene	40.000	39.337	1.7	100	0.00
14		1,2-Dichlorobenzene	40.000	39.547	1.1	100	0.00
15		Benzyl Alcohol	40.000	40.380	-1.0	102	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	41.377	-3.4	105	0.00
17		2-Methylphenol	40.000	40.499	-1.2	102	0.00
18		Hexachloroethane	40.000	40.077	-0.2	102	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	41.830	-4.6	105	0.00
20		3+4-Methylphenols	40.000	40.131	-0.3	102	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	104	0.00
22		Acetophenone	40.000	43.178	-7.9	104	0.00
23	S	Nitrobenzene-d5	80.000	79.613	0.5	104	0.00
24		Nitrobenzene	40.000	39.739	0.7	104	0.00
25		Isophorone	40.000	39.478	1.3	104	0.00
26	C	2-Nitrophenol	40.000	37.230	6.9	99	0.00
27		2,4-Dimethylphenol	40.000	39.104	2.2	103	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.770	0.6	105	0.00
29	C	2,4-Dichlorophenol	40.000	38.963	2.6	103	0.00
30		1,2,4-Trichlorobenzene	40.000	38.448	3.9	101	0.00
31		Naphthalene	40.000	39.063	2.3	103	0.00
32		Benzoic acid	40.000	42.736	-6.8	102	0.00
33		4-Chloroaniline	40.000	39.515	1.2	104	0.00
34	C	Hexachlorobutadiene	40.000	38.310	4.2	102	0.00
35		Caprolactam	40.000	38.941	2.6	103	0.00
36	C	4-Chloro-3-methylphenol	40.000	39.387	1.5	104	0.00
37		2-Methylnaphthalene	40.000	39.010	2.5	103	0.00
38		1-Methylnaphthalene	40.000	39.141	2.1	103	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	105	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	43.868	-9.7	103	0.00
41	P	Hexachlorocyclopentadiene	40.000	37.580	6.1	102	0.00
42	S	2,4,6-Tribromophenol	80.000	77.534	3.1	103	0.00
43	C	2,4,6-Trichlorophenol	40.000	38.309	4.2	102	0.00
44		2,4,5-Trichlorophenol	40.000	38.472	3.8	103	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM092019\
Data File : BM022732.D
Acq On : 20 Sep 2019 19:06
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Sep 21 02:12:26 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM092019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 20 16:34:01 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	78.491	1.9	103	0.00
46		1,1'-Biphenyl	40.000	43.151	-7.9	104	0.00
47		2-Chloronaphthalene	40.000	38.939	2.7	104	0.00
48		2-Nitroaniline	40.000	40.302	-0.8	106	0.00
49		Acenaphthylene	40.000	38.927	2.7	104	0.00
50		Dimethylphthalate	40.000	38.730	3.2	104	0.00
51		2,6-Dinitrotoluene	40.000	38.636	3.4	103	0.00
52	C	Acenaphthene	40.000	39.202	2.0	104	0.00
53		3-Nitroaniline	40.000	39.442	1.4	104	0.00
54	P	2,4-Dinitrophenol	40.000	35.858	10.4	99	0.00
55		Dibenzofuran	40.000	38.819	3.0	104	0.00
56	P	4-Nitrophenol	40.000	39.345	1.6	103	0.00
57		2,4-Dinitrotoluene	40.000	39.191	2.0	104	0.00
58		Fluorene	40.000	39.419	1.5	105	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	38.002	5.0	102	0.00
60		Diethylphthalate	40.000	38.955	2.6	105	0.00
61		4-Chlorophenyl-phenylether	40.000	39.224	1.9	104	0.00
62		4-Nitroaniline	40.000	39.837	0.4	105	0.00
63		Azobenzene	40.000	40.357	-0.9	106	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	103	0.00
65		4,6-Dinitro-2-methylphenol	40.000	36.942	7.6	100	0.00
66	c	n-Nitrosodiphenylamine	40.000	39.871	0.3	104	0.00
67		4-Bromophenyl-phenylether	40.000	38.876	2.8	102	0.00
68		Hexachlorobenzene	40.000	39.489	1.3	104	0.00
69		Atrazine	40.000	39.487	1.3	104	0.00
70	C	Pentachlorophenol	40.000	38.858	2.9	102	0.00
71		Phenanthrene	40.000	39.402	1.5	103	0.00
72		Anthracene	40.000	39.707	0.7	103	0.00
73		Carbazole	40.000	39.715	0.7	103	0.00
74		Di-n-butylphthalate	40.000	39.815	0.5	104	0.00
75	C	Fluoranthene	40.000	39.046	2.4	102	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	102	0.00
77		Benzidine	40.000	40.394	-1.0	104	0.00
78		Pyrene	40.000	39.850	0.4	103	0.00
79	S	Terphenyl-d14	80.000	82.645	-3.3	104	0.00
80		Butylbenzylphthalate	40.000	38.887	2.8	102	0.00
81		Benzo(a)anthracene	40.000	39.680	0.8	103	0.00
82		3,3'-Dichlorobenzidine	40.000	38.998	2.5	99	0.00
83		Chrysene	40.000	39.711	0.7	103	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	39.418	1.5	102	-0.01
85	c	Di-n-octyl phthalate	40.000	38.161	4.6	101	-0.01
86		Indeno(1,2,3-cd)pyrene	40.000	38.639	3.4	101	-0.01
87	I	Perylene-d12	20.000	20.000	0.0	101	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM092019\
Data File : BM022732.D
Acq On : 20 Sep 2019 19:06
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Sep 21 02:12:26 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM092019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 20 16:34:01 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	40.531	-1.3	103	-0.01
89	Benzo(k)fluoranthene	40.000	39.043	2.4	98	0.00
90 C	Benzo(a)pyrene	40.000	39.730	0.7	101	-0.01
91	Dibenzo(a,h)anthracene	40.000	39.685	0.8	100	-0.01
92	Benzo(g,h,i)perylene	40.000	39.248	1.9	100	-0.01

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