Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031119\
Data File : BM019105.D
Acq On : 11 Mar 2019 22:25
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Mar 12 04:45:39 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Mar 11 17:30:46 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	25	0.00
2		1,4-Dioxane	40.000	36.197	9.5	24	0.00
3		Pyridine	40.000	42.670	-6.7	26	0.01
4		n-Nitrosodimethylamine	40.000	34.653	13.4	22	0.00
5	S	2-Fluorophenol	80.000	77.807	2.7	25	0.00
6		Aniline	40.000	40.228	-0.6	26	0.00
7	S	Phenol-d6	80.000	81.563	-2.0	26	-0.01
8		2-Chlorophenol	40.000	40.750	-1.9	26	0.00
9		Benzaldehyde	40.000	41.097	-2.7	26	0.00
10	C	Phenol	40.000	41.309	-3.3	26	0.00
11		bis(2-Chloroethyl)ether	40.000	38.420	3.9	24	0.00
12		1,3-Dichlorobenzene	40.000	39.240	1.9	26	0.00
13	C	1,4-Dichlorobenzene	40.000	39.492	1.3	26	0.00
14		1,2-Dichlorobenzene	40.000	40.458	-1.1	26	0.00
15		Benzyl Alcohol	40.000	40.357	-0.9	25	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	40.861	-2.2	27	-0.01
17		2-Methylphenol	40.000	41.346	-3.4	26	0.00
18		Hexachloroethane	40.000	39.126	2.2	25	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.930	0.2	25	0.00
20		3+4-Methylphenols	40.000	41.567	-3.9	26	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	26	0.00
22		Acetophenone	40.000	39.422	1.4	26	0.00
23	S	Nitrobenzene-d5	80.000	77.623	3.0	26	0.00
24		Nitrobenzene	40.000	40.047	-0.1	26	0.00
25		Isophorone	40.000	38.305	4.2	25	0.00
26	C	2-Nitrophenol	40.000	37.724	5.7	24	0.00
27		2,4-Dimethylphenol	40.000	39.474	1.3	26	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.864	2.8	26	0.00
29	C	2,4-Dichlorophenol	40.000	39.848	0.4	26	0.00
30		1,2,4-Trichlorobenzene	40.000	39.665	0.8	27	-0.01
31		Naphthalene	40.000	39.602	1.0	27	0.00
32		Benzoic acid	40.000	26.618	33.5#	17	-0.07
33		4-Chloroaniline	40.000	39.388	1.5	26	0.00
34	C	Hexachlorobutadiene	40.000	38.018	5.0	26	0.00
35		Caprolactam	40.000	36.858	7.9	23	-0.02
36	C	4-Chloro-3-methylphenol	40.000	39.684	0.8	26	-0.01
37		2-Methylnaphthalene	40.000	39.698	0.8	27	0.00
38		1-Methylnaphthalene	40.000	39.772	0.6	27	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	26	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	39.181	2.0	26	0.00
41	P	Hexachlorocyclopentadiene	40.000	29.049	27.4#	19	0.00
42	S	2,4,6-Tribromophenol	80.000	76.870	3.9	25	0.00
43	C	2,4,6-Trichlorophenol	40.000	38.995	2.5	25	0.00
44		2,4,5-Trichlorophenol	40.000	38.954	2.6	25	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031119\
Data File : BM019105.D
Acq On : 11 Mar 2019 22:25
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Mar 12 04:45:39 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Mar 11 17:30:46 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	78.385	2.0	26	0.00
46		1,1'-Biphenyl	40.000	39.502	1.2	27	0.00
47		2-Chloronaphthalene	40.000	40.120	-0.3	27	0.00
48		2-Nitroaniline	40.000	37.497	6.3	24	0.00
49		Acenaphthylene	40.000	40.019	-0.0	27	0.00
50		Dimethylphthalate	40.000	39.301	1.7	27	-0.01
51		2,6-Dinitrotoluene	40.000	38.659	3.4	25	0.00
52	C	Acenaphthene	40.000	39.757	0.6	27	0.00
53		3-Nitroaniline	40.000	35.924	10.2	24	0.00
54	P	2,4-Dinitrophenol	40.000	25.909	35.2#	17	0.00
55		Dibenzofuran	40.000	39.395	1.5	27	0.00
56	P	4-Nitrophenol	40.000	34.467	13.8	23	0.00
57		2,4-Dinitrotoluene	40.000	36.315	9.2	24	0.00
58		Fluorene	40.000	38.605	3.5	26	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	39.452	1.4	26	0.00
60		Diethylphthalate	40.000	39.066	2.3	26	0.00
61		4-Chlorophenyl-phenylether	40.000	38.657	3.4	26	0.00
62		4-Nitroaniline	40.000	35.290	11.8	24	-0.01
63		Azobenzene	40.000	39.583	1.0	26	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	27	0.00
65		4,6-Dinitro-2-methylphenol	40.000	30.694	23.3	21	0.00
66	c	n-Nitrosodiphenylamine	40.000	39.076	2.3	26	0.00
67		4-Bromophenyl-phenylether	40.000	38.704	3.2	26	0.00
68		Hexachlorobenzene	40.000	39.083	2.3	27	0.00
69		Atrazine	40.000	38.648	3.4	26	0.00
70	C	Pentachlorophenol	40.000	35.526	11.2	24	0.00
71		Phenanthrene	40.000	38.900	2.8	27	0.00
72		Anthracene	40.000	38.844	2.9	26	0.00
73		Carbazole	40.000	40.135	-0.3	27	0.00
74		Di-n-butylphthalate	40.000	38.708	3.2	26	0.00
75	C	Fluoranthene	40.000	40.313	-0.8	28	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	31	0.00
77		Benzidine	40.000	38.475	3.8	30	0.00
78		Pyrene	40.000	37.406	6.5	28	0.00
79	S	Terphenyl-d14	80.000	74.182	7.3	28	0.00
80		Butylbenzylphthalate	40.000	37.297	6.8	27	0.00
81		Benzo(a)anthracene	40.000	39.011	2.5	31	0.00
82		3,3'-Dichlorobenzidine	40.000	41.231	-3.1	31	0.00
83		Chrysene	40.000	39.894	0.3	31	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	38.265	4.3	28	0.00
85	c	Di-n-octyl phthalate	40.000	41.047	-2.6	32	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	57.883	-44.7#	51	-0.01
87	I	Perylene-d12	20.000	20.000	0.0	40	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031119\
Data File : BM019105.D
Acq On : 11 Mar 2019 22:25
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Mar 12 04:45:39 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Mar 11 17:30:46 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	36.576	8.6	37	0.00
89	Benzo(k)fluoranthene	40.000	38.717	3.2	38	0.00
90 C	Benzo(a)pyrene	40.000	39.292	1.8	39	0.00
91	Dibenzo(a,h)anthracene	40.000	48.852	-22.1	51	-0.02
92	Benzo(g,h,i)perylene	40.000	51.419	-28.5#	54	-0.01

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