Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM012819\
Data File : BM018663.D
Acq On : 28 Jan 2019 14:15
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Jan 28 15:13:40 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM010919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 18 22:01:30 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	68	-0.01
2		1,4-Dioxane	40.000	39.212	2.0	67	-0.01
3		Pyridine	40.000	39.348	1.6	63	0.00
4		n-Nitrosodimethylamine	40.000	39.277	1.8	63	-0.01
5	S	2-Fluorophenol	80.000	81.358	-1.7	68	-0.01
6		Aniline	40.000	40.352	-0.9	65	-0.02
7	S	Phenol-d6	80.000	77.720	2.9	64	-0.01
8		2-Chlorophenol	40.000	38.953	2.6	65	-0.01
9		Benzaldehyde	40.000	32.487	18.8	59	-0.01
10	C	Phenol	40.000	38.223	4.4	64	-0.01
11		bis(2-Chloroethyl)ether	40.000	37.998	5.0	63	-0.01
12		1,3-Dichlorobenzene	40.000	38.620	3.5	66	-0.01
13	C	1,4-Dichlorobenzene	40.000	38.899	2.8	66	-0.01
14		1,2-Dichlorobenzene	40.000	38.549	3.6	65	-0.01
15		Benzyl Alcohol	40.000	38.615	3.5	63	-0.01
16		2,2'-oxybis(1-Chloropropane	40.000	37.663	5.8	64	-0.01
17		2-Methylphenol	40.000	38.100	4.7	63	-0.01
18		Hexachloroethane	40.000	38.460	3.8	66	-0.01
19	P	n-Nitroso-di-n-propylamine	40.000	36.108	9.7	60	-0.01
20		3+4-Methylphenols	40.000	37.940	5.2	62	-0.01
21	I	Naphthalene-d8	20.000	20.000	0.0	63	-0.01
22		Acetophenone	40.000	38.155	4.6	60	0.00
23	S	Nitrobenzene-d5	80.000	78.375	2.0	61	-0.01
24		Nitrobenzene	40.000	38.996	2.5	61	-0.01
25		Isophorone	40.000	37.943	5.1	58	-0.02
26	C	2-Nitrophenol	40.000	43.279	-8.2	64	-0.01
27		2,4-Dimethylphenol	40.000	38.754	3.1	60	-0.01
28		bis(2-Chloroethoxy)methane	40.000	37.994	5.0	59	-0.01
29	C	2,4-Dichlorophenol	40.000	40.474	-1.2	61	-0.01
30		1,2,4-Trichlorobenzene	40.000	39.093	2.3	62	-0.01
31		Naphthalene	40.000	38.522	3.7	61	-0.01
32		Benzoic acid	40.000	44.110	-10.3	73	-0.04
33		4-Chloroaniline	40.000	42.289	-5.7	63	-0.01
34	C	Hexachlorobutadiene	40.000	38.794	3.0	62	-0.01
35		Caprolactam	40.000	37.134	7.2	56	-0.03
36	C	4-Chloro-3-methylphenol	40.000	38.273	4.3	57	-0.01
37		2-Methylnaphthalene	40.000	37.375	6.6	59	-0.01
38		1-Methylnaphthalene	40.000	37.031	7.4	58	-0.01
39	I	Acenaphthene-d10	20.000	20.000	0.0	59	-0.01
40		1,2,4,5-Tetrachlorobenzene	40.000	39.669	0.8	58	-0.01
41	P	Hexachlorocyclopentadiene	40.000	56.373	-40.9#	79	-0.02
42	S	2,4,6-Tribromophenol	80.000	80.294	-0.4	56	-0.01
43	C	2,4,6-Trichlorophenol	40.000	41.311	-3.3	58	-0.01
44		2,4,5-Trichlorophenol	40.000	41.288	-3.2	58	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM012819\
Data File : BM018663.D
Acq On : 28 Jan 2019 14:15
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Jan 28 15:13:40 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM010919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 18 22:01:30 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	78.041	2.4	57	-0.01
46		1,1'-Biphenyl	40.000	38.724	3.2	57	-0.01
47		2-Chloronaphthalene	40.000	38.985	2.5	57	-0.01
48		2-Nitroaniline	40.000	41.414	-3.5	57	-0.01
49		Acenaphthylene	40.000	38.903	2.7	57	-0.01
50		Dimethylphthalate	40.000	37.461	6.3	54	-0.01
51		2,6-Dinitrotoluene	40.000	39.719	0.7	56	-0.01
52	C	Acenaphthene	40.000	38.173	4.6	56	-0.01
53		3-Nitroaniline	40.000	41.328	-3.3	57	-0.01
54	P	2,4-Dinitrophenol	40.000	37.026	7.4	56	-0.01
55		Dibenzofuran	40.000	37.981	5.0	56	-0.01
56	P	4-Nitrophenol	40.000	38.709	3.2	55	-0.01
57		2,4-Dinitrotoluene	40.000	38.062	4.8	55	-0.01
58		Fluorene	40.000	37.965	5.1	55	-0.01
59		2,3,4,6-Tetrachlorophenol	40.000	39.257	1.9	55	-0.01
60		Diethylphthalate	40.000	37.365	6.6	54	-0.01
61		4-Chlorophenyl-phenylether	40.000	37.303	6.7	55	-0.01
62		4-Nitroaniline	40.000	36.890	7.8	52	-0.01
63		Azobenzene	40.000	38.296	4.3	55	-0.01
64	I	Phenanthrene-d10	20.000	20.000	0.0	56	-0.01
65		4,6-Dinitro-2-methylphenol	40.000	39.511	1.2	58	-0.01
66	c	n-Nitrosodiphenylamine	40.000	39.399	1.5	55	-0.01
67		4-Bromophenyl-phenylether	40.000	38.834	2.9	55	0.00
68		Hexachlorobenzene	40.000	38.676	3.3	55	-0.01
69		Atrazine	40.000	37.957	5.1	51	-0.01
70	C	Pentachlorophenol	40.000	48.681	-21.7#	68	0.00
71		Phenanthrene	40.000	38.361	4.1	54	-0.01
72		Anthracene	40.000	39.188	2.0	54	0.00
73		Carbazole	40.000	39.108	2.2	54	-0.01
74		Di-n-butylphthalate	40.000	40.947	-2.4	54	-0.01
75	C	Fluoranthene	40.000	38.525	3.7	54	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	56	-0.01
77		Benzidine	40.000	38.645	3.4	46	-0.01
78		Pyrene	40.000	39.504	1.2	53	-0.01
79	S	Terphenyl-d14	80.000	80.800	-1.0	54	-0.01
80		Butylbenzylphthalate	40.000	40.940	-2.3	56	-0.01
81		Benzo(a)anthracene	40.000	38.660	3.4	53	-0.01
82		3,3'-Dichlorobenzidine	40.000	40.766	-1.9	54	-0.01
83		Chrysene	40.000	38.692	3.3	53	-0.01
84		Bis(2-ethylhexyl)phthalate	40.000	40.921	-2.3	56	0.00
85	c	Di-n-octyl phthalate	40.000	41.283	-3.2	56	-0.01
86		Indeno(1,2,3-cd)pyrene	40.000	39.551	1.1	54	-0.02
87	I	Perylene-d12	20.000	20.000	0.0	56	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM012819\
Data File : BM018663.D
Acq On : 28 Jan 2019 14:15
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Jan 28 15:13:40 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM010919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 18 22:01:30 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	39.269	1.8	55	-0.01
89	Benzo(k)fluoranthene	40.000	38.102	4.7	52	-0.01
90 C	Benzo(a)pyrene	40.000	38.939	2.7	54	-0.01
91	Dibenzo(a,h)anthracene	40.000	39.286	1.8	55	-0.02
92	Benzo(g,h,i)perylene	40.000	39.531	1.2	55	-0.02

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