Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM110818\
Data File : BM017551.D
Acq On : 07 Nov 2018 09:30
Operator : MJ/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Nov 07 13:36:55 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM101718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 29 15:08:36 2018
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	111	-0.02
2		1,4-Dioxane	40.000	36.790	8.0	105	-0.01
3		Pyridine	40.000	41.254	-3.1	109	-0.01
4		n-Nitrosodimethylamine	40.000	43.399	-8.5	122	-0.01
5	S	2-Fluorophenol	80.000	80.383	-0.5	109	-0.02
6		Aniline	40.000	41.713	-4.3	112	-0.02
7	S	Phenol-d6	80.000	83.907	-4.9	112	-0.02
8		2-Chlorophenol	40.000	41.506	-3.8	111	-0.02
9		Benzaldehyde	40.000	37.403	6.5	105	-0.02
10	C	Phenol	40.000	41.057	-2.6	110	-0.02
11		bis(2-Chloroethyl)ether	40.000	41.233	-3.1	112	-0.02
12		1,3-Dichlorobenzene	40.000	39.609	1.0	110	-0.02
13	C	1,4-Dichlorobenzene	40.000	39.704	0.7	110	-0.02
14		1,2-Dichlorobenzene	40.000	40.032	-0.1	110	-0.02
15		Benzyl Alcohol	40.000	47.957	-19.9	125	-0.02
16		2,2'-oxybis(1-Chloropropane	40.000	44.125	-10.3	122	-0.02
17		2-Methylphenol	40.000	44.018	-10.0	117	-0.02
18		Hexachloroethane	40.000	41.922	-4.8	114	-0.03
19	P	n-Nitroso-di-n-propylamine	40.000	49.182	-23.0	127	-0.02
20		3+4-Methylphenols	40.000	45.357	-13.4	120	-0.02
21	I	Naphthalene-d8	20.000	20.000	0.0	119	-0.02
22		Acetophenone	40.000	39.945	0.1	116	-0.02
23	S	Nitrobenzene-d5	80.000	88.543	-10.7	122	-0.02
24		Nitrobenzene	40.000	42.164	-5.4	117	-0.02
25		Isophorone	40.000	45.660	-14.1	127	-0.02
26	C	2-Nitrophenol	40.000	45.024	-12.6	139	-0.02
27		2,4-Dimethylphenol	40.000	42.692	-6.7	121	-0.02
28		bis(2-Chloroethoxy)methane	40.000	42.993	-7.5	122	-0.02
29	C	2,4-Dichlorophenol	40.000	44.237	-10.6	124	-0.02
30		1,2,4-Trichlorobenzene	40.000	40.107	-0.3	119	-0.02
31		Naphthalene	40.000	39.928	0.2	118	-0.02
32		Benzoic acid	40.000	48.995	-22.5	148	-0.01
33		4-Chloroaniline	40.000	43.121	-7.8	122	-0.02
34	C	Hexachlorobutadiene	40.000	40.004	-0.0	120	-0.03
35		Caprolactam	40.000	47.056	-17.6	136	-0.02
36	C	4-Chloro-3-methylphenol	40.000	45.913	-14.8	129	-0.02
37		2-Methylnaphthalene	40.000	42.979	-7.4	123	-0.03
38		1-Methylnaphthalene	40.000	43.659	-9.1	125	-0.02
39	I	Acenaphthene-d10	20.000	20.000	0.0	130	-0.02
40		1,2,4,5-Tetrachlorobenzene	40.000	39.342	1.6	127	-0.02
41	P	Hexachlorocyclopentadiene	40.000	43.313	-8.3	134	-0.03
42	S	2,4,6-Tribromophenol	80.000	92.871	-16.1	150	-0.02
43	C	2,4,6-Trichlorophenol	40.000	44.008	-10.0	132	-0.02
44		2,4,5-Trichlorophenol	40.000	43.665	-9.2	133	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM110818\
Data File : BM017551.D
Acq On : 07 Nov 2018 09:30
Operator : MJ/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Nov 07 13:36:55 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM101718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 29 15:08:36 2018
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	80.119	-0.1	129	-0.02
46		1,1'-Biphenyl	40.000	39.389	1.5	127	-0.02
47		2-Chloronaphthalene	40.000	39.162	2.1	125	-0.02
48		2-Nitroaniline	40.000	45.285	-13.2	150	-0.02
49		Acenaphthylene	40.000	41.726	-4.3	129	-0.02
50		Dimethylphthalate	40.000	43.469	-8.7	136	-0.02
51		2,6-Dinitrotoluene	40.000	44.661	-11.7	140	-0.02
52	C	Acenaphthene	40.000	40.250	-0.6	128	-0.02
53		3-Nitroaniline	40.000	43.634	-9.1	137	-0.02
54	P	2,4-Dinitrophenol	40.000	45.623	-14.1	161	-0.02
55		Dibenzofuran	40.000	39.853	0.4	129	-0.02
56	P	4-Nitrophenol	40.000	40.593	-1.5	132	-0.02
57		2,4-Dinitrotoluene	40.000	46.762	-16.9	145	-0.02
58		Fluorene	40.000	41.912	-4.8	134	-0.02
59		2,3,4,6-Tetrachlorophenol	40.000	45.024	-12.6	137	-0.02
60		Diethylphthalate	40.000	46.379	-15.9	143	-0.02
61		4-Chlorophenyl-phenylether	40.000	42.304	-5.8	134	-0.02
62		4-Nitroaniline	40.000	39.501	1.2	128	-0.02
63		Azobenzene	40.000	44.935	-12.3	137	-0.02
64	I	Phenanthrene-d10	20.000	20.000	0.0	134	-0.02
65		4,6-Dinitro-2-methylphenol	40.000	46.131	-15.3	164	-0.02
66	c	n-Nitrosodiphenylamine	40.000	42.250	-5.6	136	-0.02
67		4-Bromophenyl-phenylether	40.000	43.986	-10.0	141	-0.02
68		Hexachlorobenzene	40.000	41.659	-4.1	138	-0.02
69		Atrazine	40.000	42.811	-7.0	138	-0.02
70	C	Pentachlorophenol	40.000	44.504	-11.3	143	-0.02
71		Phenanthrene	40.000	40.030	-0.1	133	-0.02
72		Anthracene	40.000	41.510	-3.8	133	-0.02
73		Carbazole	40.000	42.289	-5.7	137	-0.02
74		Di-n-butylphthalate	40.000	49.180	-22.9	159	-0.02
75	C	Fluoranthene	40.000	42.544	-6.4	137	-0.02
76	I	Chrysene-d12	20.000	20.000	0.0	140	-0.02
77		Benzidine	40.000	46.913	-17.3	149	-0.02
78		Pyrene	40.000	40.834	-2.1	134	-0.02
79	S	Terphenyl-d14	80.000	85.320	-6.6	143	-0.02
80		Butylbenzylphthalate	40.000	51.111	-27.8#	195	-0.02
81		Benzo(a)anthracene	40.000	40.746	-1.9	139	-0.02
82		3,3'-Dichlorobenzidine	40.000	46.228	-15.6	155	-0.02
83		Chrysene	40.000	39.629	0.9	137	-0.02
84		Bis(2-ethylhexyl)phthalate	40.000	50.316	-25.8#	183	-0.02
85	c	Di-n-octyl phthalate	40.000	49.713	-24.3#	182	-0.02
86		Indeno(1,2,3-cd)pyrene	40.000	37.442	6.4	130	-0.04
87	I	Perylene-d12	20.000	20.000	0.0	137	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM110818\
Data File : BM017551.D
Acq On : 07 Nov 2018 09:30
Operator : MJ/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Nov 07 13:36:55 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM101718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 29 15:08:36 2018
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	42.039	-5.1	141	-0.03
89	Benzo(k)fluoranthene	40.000	40.869	-2.2	136	-0.02
90 C	Benzo(a)pyrene	40.000	41.150	-2.9	137	-0.03
91	Dibenzo(a,h)anthracene	40.000	39.060	2.3	132	-0.04
92	Benzo(g,h,i)perylene	40.000	37.177	7.1	127	-0.04

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