Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM072017\
Data File : BM010925.D
Acq On : 21 Jul 2017 01:05
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Jul 21 03:06:24 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM071217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 12 14:21:16 2017
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	90	-0.02
2		1,4-Dioxane	40.000	34.357	14.1	77	0.00
3		Pyridine	40.000	38.895	2.8	85	0.00
4		n-Nitrosodimethylamine	40.000	42.196	-5.5	95	0.00
5	S	2-Fluorophenol	80.000	74.753	6.6	84	0.00
6		Aniline	40.000	40.688	-1.7	91	-0.01
7	S	Phenol-d6	80.000	79.855	0.2	90	0.00
8		2-Chlorophenol	40.000	40.060	-0.2	89	-0.01
9		Benzaldehyde	40.000	38.749	3.1	87	0.00
10	C	Phenol	40.000	40.836	-2.1	92	0.00
11		bis(2-Chloroethyl)ether	40.000	39.939	0.2	91	-0.01
12		1,3-Dichlorobenzene	40.000	39.048	2.4	90	-0.01
13	C	1,4-Dichlorobenzene	40.000	38.185	4.5	88	-0.01
14		1,2-Dichlorobenzene	40.000	38.727	3.2	89	-0.01
15		Benzyl Alcohol	40.000	43.213	-8.0	98	-0.01
16		2,2'-oxybis(1-Chloropropane	40.000	43.101	-7.8	97	-0.01
17		2-Methylphenol	40.000	42.784	-7.0	95	-0.01
18		Hexachloroethane	40.000	41.259	-3.1	95	-0.02
19	P	n-Nitroso-di-n-propylamine	40.000	46.407	-16.0	107	-0.01
20		3+4-Methylphenols	40.000	44.299	-10.7	99	-0.01
21	I	Naphthalene-d8	20.000	20.000	0.0	105	-0.01
22		Acetophenone	40.000	37.639	5.9	101	-0.01
23	S	Nitrobenzene-d5	80.000	79.915	0.1	106	0.00
24		Nitrobenzene	40.000	39.245	1.9	105	0.00
25		Isophorone	40.000	41.235	-3.1	109	0.00
26	C	2-Nitrophenol	40.000	39.247	1.9	105	-0.01
27		2,4-Dimethylphenol	40.000	39.623	0.9	104	-0.01
28		bis(2-Chloroethoxy)methane	40.000	39.704	0.7	104	-0.01
29	C	2,4-Dichlorophenol	40.000	39.674	0.8	106	-0.01
30		1,2,4-Trichlorobenzene	40.000	38.463	3.8	103	-0.02
31		Naphthalene	40.000	39.897	0.3	106	-0.01
32		Benzoic acid	40.000	43.276	-8.2	110	0.02
33		4-Chloroaniline	40.000	40.907	-2.3	106	-0.01
34	C	Hexachlorobutadiene	40.000	39.718	0.7	105	-0.02
35		Caprolactam	40.000	46.680	-16.7	119	0.01
36	C	4-Chloro-3-methylphenol	40.000	45.463	-13.7	119	0.00
37		2-Methylnaphthalene	40.000	42.950	-7.4	115	-0.01
38	I	Acenaphthene-d10	20.000	20.000	0.0	127	-0.01
39		1,2,4,5-Tetrachlorobenzene	40.000	36.718	8.2	117	-0.01
40	P	Hexachlorocyclopentadiene	40.000	36.081	9.8	110	-0.01
41	S	2,4,6-Tribromophenol	80.000	92.310	-15.4	145	0.00
42	C	2,4,6-Trichlorophenol	40.000	39.480	1.3	127	0.00
43		2,4,5-Trichlorophenol	40.000	39.848	0.4	122	0.00
44	S	2-Fluorobiphenyl	80.000	76.959	3.8	124	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM072017\
Data File : BM010925.D
Acq On : 21 Jul 2017 01:05
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Jul 21 03:06:24 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM071217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 12 14:21:16 2017
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		1,1'-Biphenyl	40.000	38.549	3.6	123	-0.01
46		2-Chloronaphthalene	40.000	38.479	3.8	124	-0.01
47		2-Nitroaniline	40.000	46.121	-15.3	141	0.00
48		Acenaphthylene	40.000	41.121	-2.8	129	-0.01
49		Dimethylphthalate	40.000	41.212	-3.0	133	-0.01
50		2,6-Dinitrotoluene	40.000	44.328	-10.8	139	0.00
51	C	Acenaphthene	40.000	41.721	-4.3	134	0.00
52		3-Nitroaniline	40.000	43.020	-7.6	135	0.00
53	P	2,4-Dinitrophenol	40.000	34.215	14.5	107	0.00
54		Dibenzofuran	40.000	41.960	-4.9	134	-0.01
55	P	4-Nitrophenol	40.000	43.279	-8.2	134	0.00
56		2,4-Dinitrotoluene	40.000	46.566	-16.4	143	0.00
57		Fluorene	40.000	44.637	-11.6	143	-0.01
58		2,3,4,6-Tetrachlorophenol	40.000	43.153	-7.9	135	0.00
59		Diethylphthalate	40.000	45.339	-13.3	144	0.00
60		4-Chlorophenyl-phenylether	40.000	43.087	-7.7	141	-0.01
61		4-Nitroaniline	40.000	44.979	-12.4	137	0.00
62		Azobenzene	40.000	47.730	-19.3	150	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	150	0.00
64		4,6-Dinitro-2-methylphenol	40.000	37.046	7.4	138	0.00
65	c	n-Nitrosodiphenylamine	40.000	38.429	3.9	144	-0.01
66		4-Bromophenyl-phenylether	40.000	39.917	0.2	148	-0.01
67		Hexachlorobenzene	40.000	38.436	3.9	146	-0.01
68		Atrazine	40.000	41.034	-2.6	151	0.00
69	C	Pentachlorophenol	40.000	39.364	1.6	143	-0.01
70		Phenanthrene	40.000	40.346	-0.9	152	-0.01
71		Anthracene	40.000	41.198	-3.0	154	-0.01
72		Carbazole	40.000	41.371	-3.4	155	0.00
73		Di-n-butylphthalate	40.000	44.034	-10.1	160	0.00
74	C	Fluoranthene	40.000	41.792	-4.5	156	-0.01
75	I	Chrysene-d12	20.000	20.000	0.0	140	0.00
76		Benzidine	40.000	40.260	-0.6	130	-0.01
77		Pyrene	40.000	44.453	-11.1	155	-0.01
78	S	Terphenyl-d14	80.000	88.040	-10.1	157	0.00
79		Butylbenzylphthalate	40.000	46.788	-17.0	159	0.00
80		Benzo(a)anthracene	40.000	40.560	-1.4	143	0.00
81		3,3'-Dichlorobenzidine	40.000	39.387	1.5	134	0.00
82		Chrysene	40.000	40.251	-0.6	141	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	44.668	-11.7	148	0.00
84	c	Di-n-octyl phthalate	40.000	40.938	-2.3	138	-0.01
85		Indeno(1,2,3-cd)pyrene	40.000	30.557	23.6	108	-0.02
86	I	Perylene-d12	20.000	20.000	0.0	115	-0.01
87		Benzo(b)fluoranthene	40.000	42.110	-5.3	121	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM072017\
Data File : BM010925.D
Acq On : 21 Jul 2017 01:05
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Jul 21 03:06:24 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM071217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 12 14:21:16 2017
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(k)fluoranthene	40.000	42.744	-6.9	123	-0.01
89 C	Benzo(a)pyrene	40.000	41.438	-3.6	119	-0.01
90	Dibenzo(a,h)anthracene	40.000	37.272	6.8	107	-0.02
91	Benzo(g,h,i)perylene	40.000	36.998	7.5	106	-0.02

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