Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122315\
Data File : BM003707.D
Acq On : 22 Dec 2015 14:31
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Dec 22 15:52:04 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM120915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 09 18:52:47 2015
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	56	-0.03
2		1,4-Dioxane	40.000	40.570	-1.4	55	-0.03
3		Pyridine	40.000	41.184	-3.0	55	-0.04
4		n-Nitrosodimethylamine	40.000	44.918	-12.3	60	-0.03
5	S	2-Fluorophenol	80.000	84.943	-6.2	57	-0.03
6		Aniline	40.000	39.430	1.4	53	-0.03
7	S	Phenol-d6	80.000	79.727	0.3	54	-0.02
8		2-Chlorophenol	40.000	40.246	-0.6	54	-0.03
9		Benzaldehyde	40.000	39.455	1.4	50	-0.03
10	C	Phenol	40.000	39.990	0.0	54	-0.02
11		bis(2-Chloroethyl)ether	40.000	38.865	2.8	53	-0.03
12		1,3-Dichlorobenzene	40.000	40.970	-2.4	56	-0.03
13	C	1,4-Dichlorobenzene	40.000	40.471	-1.2	55	-0.03
14		1,2-Dichlorobenzene	40.000	40.194	-0.5	55	-0.03
15		Benzyl Alcohol	40.000	39.378	1.6	52	-0.03
16		2,2'-oxybis(1-Chloropropane	40.000	41.446	-3.6	57	-0.02
17		2-Methylphenol	40.000	39.536	1.2	53	-0.02
18		Hexachloroethane	40.000	41.278	-3.2	55	-0.04
19	P	n-Nitroso-di-n-propylamine	40.000	38.697	3.3	52	-0.03
20		3+4-Methylphenols	40.000	38.354	4.1	52	-0.02
21	I	Naphthalene-d8	20.000	20.000	0.0	51	-0.04
22		Acetophenone	40.000	40.955	-2.4	50	-0.03
23	S	Nitrobenzene-d5	80.000	88.190	-10.2	53	-0.03
24		Nitrobenzene	40.000	43.174	-7.9	53	-0.03
25		Isophorone	40.000	40.059	-0.1	48	-0.04
26	C	2-Nitrophenol	40.000	40.451	-1.1	50	-0.03
27		2,4-Dimethylphenol	40.000	40.982	-2.5	50	-0.02
28		bis(2-Chloroethoxy)methane	40.000	40.117	-0.3	50	-0.03
29	C	2,4-Dichlorophenol	40.000	40.711	-1.8	49	-0.03
30		1,2,4-Trichlorobenzene	40.000	41.076	-2.7	51	-0.03
31		Naphthalene	40.000	41.227	-3.1	52	-0.03
32		Benzoic acid	40.000	37.666	5.8	44	-0.03
33		4-Chloroaniline	40.000	39.677	0.8	49	-0.03
34	C	Hexachlorobutadiene	40.000	41.580	-3.9	52	-0.04
35		Caprolactam	40.000	37.565	6.1	48	-0.04
36	C	4-Chloro-3-methylphenol	40.000	39.748	0.6	49	-0.02
37		2-Methylnaphthalene	40.000	39.353	1.6	49	-0.03
38	I	Acenaphthene-d10	20.000	20.000	0.0	48	-0.03
39		1,2,4,5-Tetrachlorobenzene	40.000	40.470	-1.2	47	-0.03
40	P	Hexachlorocyclopentadiene	40.000	37.474	6.3	43	-0.03
41	S	2,4,6-Tribromophenol	80.000	84.149	-5.2	49	-0.02
42	C	2,4,6-Trichlorophenol	40.000	41.432	-3.6	47	-0.02
43		2,4,5-Trichlorophenol	40.000	41.688	-4.2	48	-0.02
44	S	2-Fluorobiphenyl	80.000	81.572	-2.0	48	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122315\
Data File : BM003707.D
Acq On : 22 Dec 2015 14:31
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Dec 22 15:52:04 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM120915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 09 18:52:47 2015
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	40.164	-0.4	47	-0.03
46		2-Chloronaphthalene	40.000	41.809	-4.5	49	-0.03
47		2-Nitroaniline	40.000	42.608	-6.5	50	-0.02
48		Acenaphthylene	40.000	41.591	-4.0	49	-0.03
49		Dimethylphthalate	40.000	40.533	-1.3	49	-0.02
50		2,6-Dinitrotoluene	40.000	41.327	-3.3	48	-0.03
51	C	Acenaphthene	40.000	41.348	-3.4	49	-0.03
52		3-Nitroaniline	40.000	43.440	-8.6	50	-0.02
53	P	2,4-Dinitrophenol	40.000	37.131	7.2	47	-0.02
54		Dibenzofuran	40.000	41.088	-2.7	49	-0.03
55	P	4-Nitrophenol	40.000	41.531	-3.8	51	-0.01
56		2,4-Dinitrotoluene	40.000	43.164	-7.9	50	-0.02
57		Fluorene	40.000	42.268	-5.7	51	-0.03
58		2,3,4,6-Tetrachlorophenol	40.000	42.400	-6.0	50	-0.02
59		Diethylphthalate	40.000	41.683	-4.2	50	-0.03
60		4-Chlorophenyl-phenylether	40.000	40.783	-2.0	50	-0.03
61		4-Nitroaniline	40.000	44.169	-10.4	54	-0.02
62		Azobenzene	40.000	44.319	-10.8	52	-0.03
63	I	Phenanthrene-d10	20.000	20.000	0.0	52	-0.02
64		4,6-Dinitro-2-methylphenol	40.000	40.813	-2.0	52	-0.02
65	c	n-Nitrosodiphenylamine	40.000	41.028	-2.6	50	-0.03
66		4-Bromophenyl-phenylether	40.000	40.249	-0.6	49	-0.03
67		Hexachlorobenzene	40.000	40.518	-1.3	51	-0.03
68		Atrazine	40.000	43.276	-8.2	52	-0.02
69	C	Pentachlorophenol	40.000	38.038	4.9	48	-0.02
70		Phenanthrene	40.000	41.645	-4.1	53	-0.02
71		Anthracene	40.000	42.303	-5.8	54	-0.03
72		Carbazole	40.000	44.277	-10.7	56	-0.02
73		Di-n-butylphthalate	40.000	45.424	-13.6	56	-0.03
74	C	Fluoranthene	40.000	45.289	-13.2	59	-0.02
75	I	Chrysene-d12	20.000	20.000	0.0	66	-0.02
76		Benzidine	40.000	41.839	-4.6	63	-0.02
77		Pyrene	40.000	36.519	8.7	60	-0.02
78	S	Terphenyl-d14	80.000	80.394	-0.5	66	-0.02
79		Butylbenzylphthalate	40.000	39.131	2.2	64	-0.02
80		Benzo(a)anthracene	40.000	40.563	-1.4	65	-0.02
81		3,3'-Dichlorobenzidine	40.000	45.926	-14.8	69	-0.02
82		Chrysene	40.000	40.115	-0.3	66	-0.02
83		Bis(2-ethylhexyl)phthalate	40.000	41.515	-3.8	66	-0.02
84	c	Di-n-octyl phthalate	40.000	43.963	-9.9	70	-0.03
85		Indeno(1,2,3-cd)pyrene	40.000	45.456	-13.6	69	-0.05
86	I	Perylene-d12	20.000	20.000	0.0	69	-0.04
87		Benzo(b)fluoranthene	40.000	40.803	-2.0	70	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122315\
Data File : BM003707.D
Acq On : 22 Dec 2015 14:31
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Dec 22 15:52:04 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM120915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 09 18:52:47 2015
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	39.776	0.6	67	-0.03
89 C	Benzo(a)pyrene	40.000	40.929	-2.3	68	-0.04
90	Dibenzo(a,h)anthracene	40.000	42.216	-5.5	70	-0.06
91	Benzo(g,h,i)perylene	40.000	41.063	-2.7	68	-0.05

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