Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM083117\
Data File : BM011376.D
Acq On : 31 Aug 2017 17:42
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Sep 01 05:01:31 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM083117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 31 18:39:59 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	96	0.00
2		1,4-Dioxane	40.000	37.488	6.3	90	0.00
3		Pyridine	40.000	39.484	1.3	93	0.00
4		n-Nitrosodimethylamine	40.000	38.392	4.0	92	0.00
5	S	2-Fluorophenol	80.000	76.412	4.5	92	0.00
6		Aniline	40.000	38.352	4.1	91	0.00
7	S	Phenol-d6	80.000	77.691	2.9	91	0.00
8		2-Chlorophenol	40.000	38.395	4.0	91	0.00
9		Benzaldehyde	40.000	35.990	10.0	91	0.00
10	C	Phenol	40.000	38.507	3.7	91	0.00
11		bis(2-Chloroethyl)ether	40.000	38.193	4.5	91	0.00
12		1,3-Dichlorobenzene	40.000	38.167	4.6	92	0.00
13	C	1,4-Dichlorobenzene	40.000	38.362	4.1	92	0.00
14		1,2-Dichlorobenzene	40.000	38.087	4.8	92	0.00
15		Benzyl Alcohol	40.000	38.903	2.7	91	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	38.672	3.3	92	-0.01
17		2-Methylphenol	40.000	38.534	3.7	91	0.00
18		Hexachloroethane	40.000	38.017	5.0	90	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	38.353	4.1	91	0.00
20		3+4-Methylphenols	40.000	38.864	2.8	91	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	98	0.00
22		Acetophenone	40.000	37.339	6.7	91	0.00
23	S	Nitrobenzene-d5	80.000	79.987	0.0	94	0.00
24		Nitrobenzene	40.000	39.481	1.3	92	0.00
25		Isophorone	40.000	37.488	6.3	91	0.00
26	C	2-Nitrophenol	40.000	37.706	5.7	97	0.00
27		2,4-Dimethylphenol	40.000	37.579	6.1	92	0.00
28		bis(2-Chloroethoxy)methane	40.000	37.654	5.9	92	0.00
29	C	2,4-Dichlorophenol	40.000	39.077	2.3	93	0.00
30		1,2,4-Trichlorobenzene	40.000	37.955	5.1	93	0.00
31		Naphthalene	40.000	37.684	5.8	92	0.00
32		Benzoic acid	40.000	38.027	4.9	97	0.00
33		4-Chloroaniline	40.000	37.929	5.2	92	0.00
34	C	Hexachlorobutadiene	40.000	37.881	5.3	93	0.00
35		Caprolactam	40.000	37.706	5.7	92	0.00
36	C	4-Chloro-3-methylphenol	40.000	38.178	4.6	92	0.00
37		2-Methylnaphthalene	40.000	38.021	4.9	93	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	99	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	38.462	3.8	93	0.00
40	P	Hexachlorocyclopentadiene	40.000	37.410	6.5	91	0.00
41	S	2,4,6-Tribromophenol	80.000	80.190	-0.2	96	0.00
42	C	2,4,6-Trichlorophenol	40.000	39.426	1.4	95	0.00
43		2,4,5-Trichlorophenol	40.000	39.264	1.8	94	0.00
44	S	2-Fluorobiphenyl	80.000	77.037	3.7	93	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM083117\
Data File : BM011376.D
Acq On : 31 Aug 2017 17:42
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Sep 01 05:01:31 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM083117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 31 18:39:59 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	37.715	5.7	93	0.00
46		2-Chloronaphthalene	40.000	38.043	4.9	93	0.00
47		2-Nitroaniline	40.000	38.414	4.0	98	0.00
48		Acenaphthylene	40.000	37.904	5.2	92	0.00
49		Dimethylphthalate	40.000	37.679	5.8	93	0.00
50		2,6-Dinitrotoluene	40.000	39.654	0.9	96	0.00
51	C	Acenaphthene	40.000	38.149	4.6	94	0.00
52		3-Nitroaniline	40.000	39.721	0.7	96	0.00
53	P	2,4-Dinitrophenol	40.000	39.336	1.7	100	0.00
54		Dibenzofuran	40.000	38.061	4.8	94	0.00
55	P	4-Nitrophenol	40.000	39.994	0.0	97	0.00
56		2,4-Dinitrotoluene	40.000	38.297	4.3	97	0.00
57		Fluorene	40.000	38.224	4.4	94	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	40.370	-0.9	97	0.00
59		Diethylphthalate	40.000	37.576	6.1	93	0.00
60		4-Chlorophenyl-phenylether	40.000	37.984	5.0	94	0.00
61		4-Nitroaniline	40.000	40.161	-0.4	97	0.00
62		Azobenzene	40.000	38.141	4.6	94	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	100	0.00
64		4,6-Dinitro-2-methylphenol	40.000	39.547	1.1	101	0.00
65	c	n-Nitrosodiphenylamine	40.000	37.589	6.0	94	0.00
66		4-Bromophenyl-phenylether	40.000	38.248	4.4	95	0.00
67		Hexachlorobenzene	40.000	37.936	5.2	95	0.00
68		Atrazine	40.000	38.131	4.7	95	0.00
69	C	Pentachlorophenol	40.000	40.493	-1.2	97	0.00
70		Phenanthrene	40.000	38.502	3.7	96	0.00
71		Anthracene	40.000	38.564	3.6	95	0.00
72		Carbazole	40.000	38.511	3.7	96	0.00
73		Di-n-butylphthalate	40.000	39.670	0.8	94	0.00
74	C	Fluoranthene	40.000	39.535	1.2	97	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	104	0.00
76		Benzidine	40.000	46.548	-16.4	124	0.00
77		Pyrene	40.000	38.374	4.1	97	0.00
78	S	Terphenyl-d14	80.000	76.614	4.2	98	0.00
79		Butylbenzylphthalate	40.000	38.680	3.3	96	0.00
80		Benzo(a)anthracene	40.000	38.460	3.8	99	0.00
81		3,3'-Dichlorobenzidine	40.000	40.911	-2.3	101	0.00
82		Chrysene	40.000	38.561	3.6	99	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	39.029	2.4	97	0.00
84	c	Di-n-octyl phthalate	40.000	40.964	-2.4	98	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	41.877	-4.7	109	0.00
86	I	Perylene-d12	20.000	20.000	0.0	110	0.00
87		Benzo(b)fluoranthene	40.000	38.254	4.4	105	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM083117\
Data File : BM011376.D
Acq On : 31 Aug 2017 17:42
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Sep 01 05:01:31 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM083117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 31 18:39:59 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	37.111	7.2	100	0.00
89 C	Benzo(a)pyrene	40.000	38.479	3.8	104	0.00
90	Dibenzo(a,h)anthracene	40.000	39.981	0.0	109	0.00
91	Benzo(g,h,i)perylene	40.000	39.614	1.0	109	0.00

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