Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM111615\
Data File : BM003139.D
Acq On : 16 Nov 2015 12:42
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Nov 17 13:49:52 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM110415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 17 12:50:51 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	5.000	5.000	0.0	106	0.00
2		1,4-Dioxane	1.000	0.955	4.5	99	-0.02
3		n-Nitrosodimethylamine	1.000	0.935	6.5	103	-0.02
4	S	2-Fluorophenol	1.000	0.885	11.5	98	-0.02
5	S	Phenol-d6	1.000	0.911	8.9	102	-0.02
6		bis(2-Chloroethyl)ether	1.000	0.974	2.6	106	-0.02
7	I	Naphthalene-d8	5.000	5.000	0.0	111	-0.01
8	S	Nitrobenzene-d5	1.000	0.947	5.3	111	-0.02
9		Nitrobenzene	1.000	0.950	5.0	112	-0.02
10		Naphthalene	1.000	0.974	2.6	110	-0.01
11		Hexachlorobutadiene	1.000	0.975	2.5	109	-0.01
12		2-Methylnaphthalene	1.000	0.982	1.8	111	-0.02
13	I	Acenaphthene-d10	5.000	5.000	0.0	106	0.00
14	S	2,4,6-Tribromophenol	1.000	0.963	3.7	104	0.00
15	S	2-Fluorobiphenyl	1.000	1.039	-3.9	112	-0.02
16		Acenaphthylene	1.000	0.958	4.2	106	0.00
17		Acenaphthene	1.000	1.047	-4.7	121	-0.03
18		Fluorene	1.000	1.189	-18.9	124	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	105	-0.03
20		4-Bromophenyl-phenylether	1.000	1.818	-81.8#	131	0.00
21		Hexachlorobenzene	1.000	1.303	-30.3#	99	0.00
22		Pentachlorophenol	1.000	1.222	-22.2	106	0.00
23		Phenanthrene	1.000	1.801	-80.1#	123	0.00
24		Anthracene	1.000	1.001	-0.1	105	0.00
25		Fluoranthene	1.000	1.249	-24.9	112	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	112	-0.01
27		Benzidine	1.000	1.012	-1.2	100	0.00
28		Pyrene	1.000	0.956	4.4	110	-0.02
29	S	Terphenyl-d14	1.000	0.923	7.7	100	0.00
30		Benzo(a)anthracene	1.000	0.933	6.7	105	-0.01
31		3,3'-Dichlorobenzidine	1.000	0.933	6.7	90	-0.01
32		Chrysene	1.000	0.870	13.0	86	-0.01
33		Bis(2-ethylhexyl)phthalate	1.000	0.719	28.1#	77	-0.01
34		Indeno(1,2,3-cd)pyrene	1.000	0.907	9.3	89	-0.02
35	I	Perylene-d12	5.000	5.000	0.0	91	-0.01
36		Benzo(b)fluoranthene	1.000	0.909	9.1	84	-0.01
37		Benzo(k)fluoranthene	1.000	0.954	4.6	88	-0.01
38	C	Benzo(a)pyrene	1.000	0.899	10.1	90	-0.01
39		Dibenzo(a,h)anthracene	1.000	0.944	5.6	90	-0.01
40		Benzo(g,h,i)perylene	1.000	0.910	9.0	89	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM111615\
Data File : BM003139.D
Acq On : 16 Nov 2015 12:42
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC001

Quant Time: Nov 17 13:49:52 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM110415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 17 12:50:51 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)

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