Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM013016\
Data File : BM004156.D
Acq On : 29 Jan 2016 21:01
Operator : SJ/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Jan 29 22:16:02 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM012216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 22 14:46:41 2016
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	74	0.00
2		1,4-Dioxane	1.000	0.966	3.4	75	0.00
3		n-Nitrosodimethylamine	1.000	0.939	6.1	76	0.00
4	S	2-Fluorophenol	1.000	0.976	2.4	80	0.00
5	S	Phenol-d6	1.000	1.000	0.0	83	0.00
6		bis(2-Chloroethyl)ether	1.000	0.918	8.2	75	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	77	0.00
8	S	Nitrobenzene-d5	1.000	0.978	2.2	84	0.00
9		Nitrobenzene	1.000	0.953	4.7	83	0.00
10		Naphthalene	1.000	0.910	9.0	76	0.00
11		Hexachlorobutadiene	1.000	0.916	8.4	77	0.00
12		2-Methylnaphthalene	1.000	0.938	6.2	79	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	82	0.00
14	S	2,4,6-Tribromophenol	1.000	1.031	-3.1	98	0.00
15	S	2-Fluorobiphenyl	1.000	0.900	10.0	80	0.00
16		Acenaphthylene	1.000	0.919	8.1	83	0.00
17		Acenaphthene	1.000	0.873	12.7	80	0.00
18		Fluorene	1.000	0.939	6.1	86	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	88	0.00
20		4-Bromophenyl-phenylether	1.000	0.894	10.6	81	0.00
21		Hexachlorobenzene	1.000	0.878	12.2	82	0.00
22		Pentachlorophenol	1.000	1.201	-20.1	143	0.00
23		Phenanthrene	1.000	0.882	11.8	85	0.00
24		Anthracene	1.000	0.922	7.8	87	0.00
25		Fluoranthene	1.000	0.965	3.5	92	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	83	0.00
27		Benzidine	1.000	1.097	-9.7	94	0.02
28		Pyrene	1.000	0.997	0.3	90	0.00
29	S	Terphenyl-d14	1.000	0.995	0.5	91	0.00
30		Benzo(a)anthracene	1.000	0.890	11.0	83	0.00
31		3,3'-Dichlorobenzidine	1.000	1.350	-35.0#	126	0.00
32		Chrysene	1.000	1.169	-16.9	107	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	1.503	-50.3#	151	0.00
34		Indeno(1,2,3-cd)pyrene	1.000	1.050	-5.0	99	0.01
35	I	Perylene-d12	5.000	5.000	0.0	96	0.00
36		Benzo(b)fluoranthene	1.000	0.875	12.5	90	0.00
37		Benzo(k)fluoranthene	1.000	0.984	1.6	107	0.00
38	C	Benzo(a)pyrene	1.000	0.949	5.1	102	0.00
39		Dibenzo(a,h)anthracene	1.000	0.930	7.0	99	0.00
40		Benzo(g,h,i)perylene	1.000	0.913	8.7	98	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM013016\
Data File : BM004156.D
Acq On : 29 Jan 2016 21:01
Operator : SJ/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC001

Quant Time: Jan 29 22:16:02 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM012216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 22 14:46:41 2016
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