Evaluate Continuing Calibration Report

Data Path : S:\HPCHEM1\BNA_M\DATA\BM022916\
Data File : BM004492.D
Acq On : 01 Mar 2016 05:42
Operator : SJ/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Mar 01 08:01:47 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BM022916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 01 00:52:28 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	94	0.00
2		1,4-Dioxane	1.000	0.943	5.7	90	0.00
3		n-Nitrosodimethylamine	1.000	0.858	14.2	86	0.00
4	S	2-Fluorophenol	1.000	0.878	12.2	88	0.00
5	S	Phenol-d6	1.000	0.914	8.6	91	0.02
6		bis(2-Chloroethyl)ether	1.000	0.933	6.7	91	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	95	0.00
8	S	Nitrobenzene-d5	1.000	0.894	10.6	90	0.00
9		Nitrobenzene	1.000	0.892	10.8	89	0.00
10		Naphthalene	1.000	0.937	6.3	93	0.00
11		Hexachlorobutadiene	1.000	0.972	2.8	94	-0.01
12		2-Methylnaphthalene	1.000	0.966	3.4	96	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	100	0.00
14	S	2,4,6-Tribromophenol	1.000	0.785	21.5	88	0.00
15	S	2-Fluorobiphenyl	1.000	0.927	7.3	97	0.00
16		Acenaphthylene	1.000	0.939	6.1	99	0.00
17		Acenaphthene	1.000	0.889	11.1	99	0.00
18		Fluorene	1.000	0.939	6.1	99	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	104	0.00
20		4-Bromophenyl-phenylether	1.000	0.803	19.7	99	0.00
21		Hexachlorobenzene	1.000	0.776	22.4	97	0.00
22		Pentachlorophenol	1.000	0.923	7.7	123	0.00
23		Phenanthrene	1.000	0.806	19.4	99	0.00
24		Anthracene	1.000	0.842	15.8	100	0.00
25		Fluoranthene	1.000	0.831	16.9	101	-0.02
26	I	Chrysene-d12	5.000	5.000	0.0	98	0.00
27		Benzidine	1.000	0.798	20.2	89	0.00
28		Pyrene	1.000	0.902	9.8	101	0.00
29	S	Terphenyl-d14	1.000	0.865	13.5	97	0.00
30		Benzo(a)anthracene	1.000	0.896	10.4	98	0.00
31		3,3'-Dichlorobenzidine	1.000	0.864	13.6	99	0.00
32		Chrysene	1.000	0.758	24.2	86	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	0.921	7.9	86	0.00
34		Indeno(1,2,3-cd)pyrene	1.000	0.933	6.7	102	-0.01
35	I	Perylene-d12	5.000	5.000	0.0	103	0.00
36		Benzo(b)fluoranthene	1.000	0.909	9.1	94	0.00
37		Benzo(k)fluoranthene	1.000	0.954	4.6	103	0.00
38	C	Benzo(a)pyrene	1.000	0.975	2.5	103	0.00
39		Dibenzo(a,h)anthracene	1.000	0.954	4.6	102	0.00
40		Benzo(g,h,i)perylene	1.000	0.966	3.4	103	-0.01

Evaluate Continuing Calibration Report

Data Path : S:\HPCHEM1\BNA_M\DATA\BM022916\
Data File : BM004492.D
Acq On : 01 Mar 2016 05:42
Operator : SJ/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Mar 01 08:01:47 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BM022916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 01 00:52:28 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