Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM113015\
Data File : BM003326.D
Acq On : 30 Nov 2015 19:35
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Dec 01 02:48:43 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM112715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 27 16:36:40 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	132	-0.02
2		1,4-Dioxane	1.000	0.993	0.7	126	-0.02
3		n-Nitrosodimethylamine	1.000	1.082	-8.2	143	-0.02
4	S	2-Fluorophenol	1.000	1.055	-5.5	143	0.00
5	S	Phenol-d6	1.000	1.156	-15.6	157	-0.02
6		bis(2-Chloroethyl)ether	1.000	1.062	-6.2	144	-0.02
7	I	Naphthalene-d8	5.000	5.000	0.0	142	-0.02
8	S	Nitrobenzene-d5	1.000	1.162	-16.2	173	-0.01
9		Nitrobenzene	1.000	1.143	-14.3	172	-0.01
10		Naphthalene	1.000	1.004	-0.4	145	0.00
11		Hexachlorobutadiene	1.000	1.073	-7.3	157	-0.02
12		2-Methylnaphthalene	1.000	1.100	-10.0	158	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	160	0.00
14	S	2,4,6-Tribromophenol	1.000	1.228	-22.8	218	0.00
15	S	2-Fluorobiphenyl	1.000	0.940	6.0	163	0.00
16		Acenaphthylene	1.000	1.146	-14.6	191	0.00
17		Acenaphthene	1.000	0.906	9.4	168	0.00
18		Fluorene	1.000	1.028	-2.8	172	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	154	0.00
20		4-Bromophenyl-phenylether	1.000	0.928	7.2	173	0.00
21		Hexachlorobenzene	1.000	0.974	2.6	164	0.00
22		Pentachlorophenol	1.000	1.140	-14.0	213	0.00
23		Phenanthrene	1.000	0.865	13.5	151	0.00
24		Anthracene	1.000	1.167	-16.7	180	0.00
25		Fluoranthene	1.000	0.910	9.0	157	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	110	-0.01
27		Benzidine	1.000	2.014	-101.4#	309	0.00
28		Pyrene	1.000	1.170	-17.0	154	0.00
29	S	Terphenyl-d14	1.000	1.137	-13.7	138	0.00
30		Benzo(a)anthracene	1.000	1.085	-8.5	135	-0.01
31		3,3'-Dichlorobenzidine	1.000	1.501	-50.1#	236	0.01
32		Chrysene	1.000	1.126	-12.6	139	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	1.906	-90.6#	284	0.00
34		Indeno(1,2,3-cd)pyrene	1.000	1.124	-12.4	143	0.00
35	I	Perylene-d12	5.000	5.000	0.0	123	0.00
36		Benzo(b)fluoranthene	1.000	1.001	-0.1	120	0.00
37		Benzo(k)fluoranthene	1.000	1.114	-11.4	144	-0.01
38	C	Benzo(a)pyrene	1.000	1.112	-11.2	143	0.00
39		Dibenzo(a,h)anthracene	1.000	1.209	-20.9	150	-0.01
40		Benzo(g,h,i)perylene	1.000	1.053	-5.3	134	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM113015\
Data File : BM003326.D
Acq On : 30 Nov 2015 19:35
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC001

Quant Time: Dec 01 02:48:43 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM112715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 27 16:36:40 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