Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM110415\
Data File : BM003048.D
Acq On : 04 Nov 2015 18:00
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : SSTDCCC001

Quant Time: Nov 05 08:04:04 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM110415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Nov 05 00:20:46 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	126	0.02
2		1,4-Dioxane	1.000	0.984	1.6	122	0.00
3		n-Nitrosodimethylamine	1.000	0.974	2.6	128	0.00
4	S	2-Fluorophenol	1.000	0.981	1.9	129	0.00
5	S	Phenol-d6	1.000	0.984	1.6	132	0.00
6		bis(2-Chloroethyl)ether	1.000	0.961	3.9	125	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	125	0.00
8	S	Nitrobenzene-d5	1.000	0.990	1.0	131	0.00
9		Nitrobenzene	1.000	0.982	1.8	131	0.00
10		Naphthalene	1.000	0.974	2.6	124	0.00
11		Hexachlorobutadiene	1.000	0.975	2.5	123	0.00
12		2-Methylnaphthalene	1.000	0.955	4.5	122	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	123	0.00
14	S	2,4,6-Tribromophenol	1.000	1.022	-2.2	129	0.00
15	S	2-Fluorobiphenyl	1.000	0.951	4.9	119	0.00
16		Acenaphthylene	1.000	0.978	2.2	126	0.00
17		Acenaphthene	1.000	0.900	10.0	121	0.00
18		Fluorene	1.000	0.989	1.1	120	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	122	0.00
20		4-Bromophenyl-phenylether	1.000	1.340	-34.0#	113	0.00
21		Hexachlorobenzene	1.000	1.303	-30.3#	115	0.00
22		Pentachlorophenol	1.000	1.301	-30.1#	133	0.00
23		Phenanthrene	1.000	1.402	-40.2#	114	0.00
24		Anthracene	1.000	1.002	-0.2	123	0.00
25		Fluoranthene	1.000	1.133	-13.3	118	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	126	0.00
27		Benzidine	1.000	1.194	-19.4	133	0.00
28		Pyrene	1.000	0.919	8.1	118	0.00
29	S	Terphenyl-d14	1.000	0.948	5.2	115	0.00
30		Benzo(a)anthracene	1.000	1.010	-1.0	128	0.00
31		3,3'-Dichlorobenzidine	1.000	1.178	-17.8	135	0.00
32		Chrysene	1.000	1.036	-3.6	110	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	0.860	14.0	112	0.00
34		Indeno(1,2,3-cd)pyrene	1.000	1.179	-17.9	131	0.00
35	I	Perylene-d12	5.000	5.000	0.0	126	0.00
36		Benzo(b)fluoranthene	1.000	0.933	6.7	120	0.00
37		Benzo(k)fluoranthene	1.000	0.993	0.7	126	0.00
38	C	Benzo(a)pyrene	1.000	0.940	6.0	131	0.00
39		Dibenzo(a,h)anthracene	1.000	0.992	0.8	131	0.01
40		Benzo(g,h,i)perylene	1.000	0.972	2.8	131	0.00

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Compound	Amount	Calc.	%Dev	Area%	Dev(min)

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