Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM010216\
Data File : BM003915.D
Acq On : 31 Dec 2015 15:40
Operator : SJ/UM
Sample : SSTDICV001
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM010216

Quant Time: Jan 04 13:45:45 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM010216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 04 13:33:21 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	173	0.00
2		1,4-Dioxane	1.000	0.959	4.1	166	0.00
3		n-Nitrosodimethylamine	1.000	1.020	-2.0	177	-0.02
4	S	2-Fluorophenol	1.000	0.967	3.3	178	0.00
5	S	Phenol-d6	1.000	1.020	-2.0	181	0.00
6		bis(2-Chloroethyl)ether	1.000	1.008	-0.8	172	-0.02
7	I	Naphthalene-d8	5.000	5.000	0.0	172	0.00
8	S	Nitrobenzene-d5	1.000	1.040	-4.0	181	-0.01
9		Nitrobenzene	1.000	1.055	-5.5	185	-0.01
10		Naphthalene	1.000	0.976	2.4	169	-0.02
11		Hexachlorobutadiene	1.000	0.969	3.1	171	0.00
12		2-Methylnaphthalene	1.000	0.991	0.9	171	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	173	0.00
14	S	2,4,6-Tribromophenol	1.000	0.990	1.0	186	0.00
15	S	2-Fluorobiphenyl	1.000	0.974	2.6	170	0.00
16		Acenaphthylene	1.000	1.054	-5.4	181	0.00
17		Acenaphthene	1.000	0.944	5.6	173	0.00
18		Fluorene	1.000	0.999	0.1	170	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	170	0.00
20		4-Bromophenyl-phenylether	1.000	0.959	4.1	173	0.00
21		Hexachlorobenzene	1.000	0.938	6.2	176	0.00
22		Pentachlorophenol	1.000	1.169	-16.9	221	0.00
23		Phenanthrene	1.000	1.059	-5.9	176	0.00
24		Anthracene	1.000	0.995	0.5	182	0.00
25		Fluoranthene	1.000	1.042	-4.2	184	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	182	0.02
27		Benzidine	1.000	1.155	-15.5	213	0.00
28		Pyrene	1.000	0.989	1.1	179	0.00
29	S	Terphenyl-d14	1.000	1.003	-0.3	184	0.00
30		Benzo(a)anthracene	1.000	0.965	3.5	180	0.02
31		3,3'-Dichlorobenzidine	1.000	1.137	-13.7	229	0.00
32		Chrysene	1.000	1.075	-7.5	193	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	1.130	-13.0	257	0.00
34		Indeno(1,2,3-cd)pyrene	1.000	0.999	0.1	181	0.01
35	I	Perylene-d12	5.000	5.000	0.0	188	0.00
36		Benzo(b)fluoranthene	1.000	0.941	5.9	189	0.00
37		Benzo(k)fluoranthene	1.000	1.006	-0.6	193	0.01
38	C	Benzo(a)pyrene	1.000	0.979	2.1	195	0.00
39		Dibenzo(a,h)anthracene	1.000	0.958	4.2	180	0.00
40		Benzo(g,h,i)perylene	1.000	0.934	6.6	177	0.01

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

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