Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM022816\
Data File : BM004460.D
Acq On : 28 Feb 2016 20:56
Operator : SJ/UM
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM022816

Quant Time: Feb 29 06:36:15 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM022816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Feb 29 05:57:18 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	86	0.00
2		1,4-Dioxane	1.000	1.084	-8.4	104	0.00
3		n-Nitrosodimethylamine	1.000	1.013	-1.3	101	0.00
4	S	2-Fluorophenol	1.000	0.988	1.2	97	0.00
5	S	Phenol-d6	1.000	0.914	8.6	97	0.00
6		bis(2-Chloroethyl)ether	1.000	1.033	-3.3	100	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	91	0.00
8	S	Nitrobenzene-d5	1.000	0.935	6.5	97	0.00
9		Nitrobenzene	1.000	0.966	3.4	98	0.00
10		Hexachlorobutadiene	1.000	0.983	1.7	98	0.00
11		2-Methylnaphthalene	1.000	1.022	-2.2	103	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	95	0.00
13	S	2,4,6-Tribromophenol	1.000	0.793	20.7	89	0.00
14	S	2-Fluorobiphenyl	1.000	0.937	6.3	99	0.00
15	I	Phenanthrene-d10	5.000	5.000	0.0	94	0.00
16		Hexachlorobenzene	1.000	1.040	-4.0	104	0.00
17		Pentachlorophenol	1.000	0.803	19.7	94	0.00
18		Fluoranthene	1.000	0.891	10.9	100	0.00
19	I	Chrysene-d12	5.000	5.000	0.0	93	0.00
20		Benzidine	1.000	1.165	-16.5	158	0.00
21		Pyrene	1.000	0.940	6.0	99	0.00
22	S	Terphenyl-d14	1.000	0.968	3.2	103	0.00
23		Benzo(a)anthracene	1.000	0.888	11.2	105	0.00
24		3,3'-Dichlorobenzidine	1.000	1.047	-4.7	101	0.00
25		Chrysene	1.000	0.845	15.5	99	0.00
26		Bis(2-ethylhexyl)phthalate	1.000	0.809	19.1	86	0.00
27		Indeno(1,2,3-cd)pyrene	1.000	0.949	5.1	102	-0.01
28	I	Perylene-d12	5.000	5.000	0.0	91	-0.01
29		Benzo(b)fluoranthene	1.000	0.951	4.9	96	0.00
30		Benzo(k)fluoranthene	1.000	0.891	10.9	100	0.00
31	C	Benzo(a)pyrene	1.000	0.991	0.9	102	0.00
32		Dibenzo(a,h)anthracene	1.000	0.955	4.5	102	0.00
33		Benzo(g,h,i)perylene	1.000	0.960	4.0	103	-0.01

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