Evaluate Continuing Calibration Report

Data Path : S:\HPCHEM1\BNA_M\DATA\BM022916\
Data File : BM004483.D
Acq On : 29 Feb 2016 22:24
Operator : SJ/UM
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM022916

Quant Time: Mar 01 05:58:08 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	96	0.00
2		1,4-Dioxane	1.000	0.977	2.3	94	0.00
3		n-Nitrosodimethylamine	1.000	0.943	5.7	96	0.00
4	S	2-Fluorophenol	1.000	0.964	3.6	98	0.00
5	S	Phenol-d6	1.000	0.989	1.1	100	0.00
6		bis(2-Chloroethyl)ether	1.000	1.003	-0.3	99	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	99	0.00
8	S	Nitrobenzene-d5	1.000	0.967	3.3	101	0.00
9		Nitrobenzene	1.000	0.963	3.7	99	0.00
10		Naphthalene	1.000	0.975	2.5	100	0.00
11		Hexachlorobutadiene	1.000	0.990	1.0	100	-0.01
12		2-Methylnaphthalene	1.000	1.000	0.0	103	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	98	0.00
14	S	2,4,6-Tribromophenol	1.000	0.946	5.4	104	0.00
15	S	2-Fluorobiphenyl	1.000	1.014	-1.4	105	0.00
16		Acenaphthylene	1.000	0.961	3.9	99	0.00
17		Acenaphthene	1.000	1.013	-1.3	111	0.00
18		Fluorene	1.000	1.042	-4.2	109	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	87	0.00
20		4-Bromophenyl-phenylether	1.000	1.205	-20.5	122	0.00
21		Hexachlorobenzene	1.000	1.226	-22.6	129	0.00
22		Pentachlorophenol	1.000	1.085	-8.5	120	0.00
23		Phenanthrene	1.000	1.174	-17.4	121	0.00
24		Anthracene	1.000	0.996	0.4	99	0.02
25		Fluoranthene	1.000	1.054	-5.4	107	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	87	0.01
27		Benzidine	1.000	1.552	-55.2#	206	0.00
28		Pyrene	1.000	1.074	-7.4	107	0.00
29	S	Terphenyl-d14	1.000	1.064	-6.4	106	0.00
30		Benzo(a)anthracene	1.000	0.934	6.6	91	0.01
31		3,3'-Dichlorobenzidine	1.000	1.489	-48.9#	163	0.00
32		Chrysene	1.000	0.931	6.9	90	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	0.817	18.3	68	-0.01
34		Indeno(1,2,3-cd)pyrene	1.000	1.043	-4.3	101	-0.01
35	I	Perylene-d12	5.000	5.000	0.0	99	-0.01
36		Benzo(b)fluoranthene	1.000	0.986	1.4	98	0.00
37		Benzo(k)fluoranthene	1.000	0.937	6.3	97	0.00
38	C	Benzo(a)pyrene	1.000	1.006	-0.6	102	0.00
39		Dibenzo(a,h)anthracene	1.000	0.977	2.3	100	0.00
40		Benzo(g,h,i)perylene	1.000	0.986	1.4	101	-0.01

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

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