Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE022315\
Data File : BE089656.D
Acq On : 24 Feb 2015 00:47
Operator : TP/JJ
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Quant Time: Feb 24 04:03:36 2015
Quant Method : Z:\HPCHEM1\BNA E\METHODS\SIMPAH-BE022315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 24 03:46:57 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	120	0.00
2		1,4-Dioxane	1.000	1.011	-1.1	97	0.00
3		n-Nitrosodimethylamine	1.000	1.044	-4.4	106	0.00
4	I	Naphthalene-d8	5.000	5.000	0.0	120	0.00
5	S	Nitrobenzene-d5	1.000	0.957	4.3	97	0.00
6		Naphthalene	1.000	1.003	-0.3	104	0.00
7		2-Methylnaphthalene	1.000	1.029	-2.9	104	0.00
8	I	Acenaphthene-d10	5.000	5.000	0.0	119	0.00
9	S	2-Fluorobiphenyl	1.000	0.975	2.5	100	0.00
10		Acenaphthylene	1.000	1.038	-3.8	105	-0.01
11		Acenaphthene	1.000	1.012	-1.2	105	0.00
12		Fluorene	1.000	1.032	-3.2	104	0.00
13	I	Phenanthrene-d10	5.000	5.000	0.0	102	0.00
14		Hexachlorobenzene	1.000	1.037	-3.7	94	0.00
15		Pentachlorophenol	1.000	0.943	5.7	95	0.00
16		Phenanthrene	1.000	1.015	-1.5	90	0.00
17		Anthracene	1.000	1.032	-3.2	90	0.00
18		Fluoranthene	1.000	1.051	-5.1	102	0.00
19	I	Chrysene-d12	5.000	5.000	0.0	117	0.00
20		Pyrene	1.000	1.037	-3.7	103	0.00
21	S	Terphenyl-d14	1.000	1.085	-8.5	78	0.00
22		Benzo(a)anthracene	1.000	0.992	0.8	101	0.00
23		Chrysene	1.000	1.017	-1.7	104	0.00
24		Indeno(1,2,3-cd)pyrene	1.000	0.876	12.4	95	0.00
25	I	Perylene-d12	5.000	5.000	0.0	116	0.00
26		Benzo(b)fluoranthene	1.000	0.853	14.7	102	0.00
27		Benzo(k)fluoranthene	1.000	1.019	-1.9	100	0.00
28	C	Benzo(a)pyrene	1.000	0.883	11.7	98	0.00
29		Dibenzo(a,h)anthracene	1.000	0.880	12.0	96	0.00
30		Benzo(g,h,i)perylene	1.000	0.920	8.0	98	0.00

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