Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM060815\
Data File : BM001604.D
Acq On : 09 Jun 2015 03:36
Operator : TP/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM060815

Quant Time: Jun 09 18:10:11 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM060815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 09 17:36:38 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	92	0.00
2		1,4-Dioxane	1.000	0.980	2.0	95	0.00
3		n-Nitrosodimethylamine	1.000	1.050	-5.0	98	0.00
4	S	2-Fluorophenol	1.000	1.023	-2.3	96	-0.02
5	S	Phenol-d6	1.000	1.007	-0.7	96	-0.02
6	I	Naphthalene-d8	5.000	5.000	0.0	94	0.00
7	S	Nitrobenzene-d5	1.000	0.986	1.4	95	0.00
8		Nitrobenzene	1.000	0.986	1.4	95	0.00
9		Naphthalene	1.000	1.028	-2.8	98	0.00
10		Hexachlorobutadiene	1.000	1.011	-1.1	96	0.00
11		2-Methylnaphthalene	1.000	1.027	-2.7	99	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	96	0.00
13	S	2,4,6-Tribromophenol	1.000	1.006	-0.6	104	0.00
14	S	2-Fluorobiphenyl	1.000	1.028	-2.8	99	0.00
15		Acenaphthylene	1.000	1.012	-1.2	99	0.00
16		Acenaphthene	1.000	1.036	-3.6	101	0.00
17		Fluorene	1.000	1.071	-7.1	105	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	91	0.03
19		4-Bromophenyl-phenylether	1.000	1.098	-9.8	100	0.00
20		Hexachlorobenzene	1.000	1.134	-13.4	106	0.00
21		Pentachlorophenol	1.000	1.129	-12.9	104	0.00
22		Phenanthrene	1.000	1.119	-11.9	93	0.00
23		Anthracene	1.000	1.187	-18.7	112	0.00
24		Fluoranthene	1.000	1.076	-7.6	103	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	96	0.00
26		Pyrene	1.000	1.026	-2.6	104	0.00
27	S	Terphenyl-d14	1.000	1.043	-4.3	104	0.00
28		Benzo(a)anthracene	1.000	0.998	0.2	101	0.00
29		Chrysene	1.000	1.050	-5.0	102	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.892	10.8	96	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.967	3.3	98	0.00
32	I	Perylene-d12	5.000	5.000	0.0	101	0.00
33		Benzo(b)fluoranthene	1.000	0.970	3.0	102	0.00
34		Benzo(k)fluoranthene	1.000	0.970	3.0	99	0.00
35	C	Benzo(a)pyrene	1.000	0.942	5.8	98	0.00
36		Dibenzo(a,h)anthracene	1.000	0.922	7.8	97	0.00
37		Benzo(g,h,i)perylene	1.000	0.940	6.0	99	0.00

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