Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM060615\
Data File : BM001578.D
Acq On : 06 Jun 2015 05:06
Operator : TP/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM060615

Quant Time: Jun 08 05:41:17 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM060615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 08 05:27:25 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	101	0.00
2		1,4-Dioxane	1.000	0.935	6.5	101	0.00
3		n-Nitrosodimethylamine	1.000	0.986	1.4	102	0.00
4	S	2-Fluorophenol	1.000	0.969	3.1	102	0.00
5	S	Phenol-d6	1.000	0.968	3.2	102	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	100	0.00
7	S	Nitrobenzene-d5	1.000	0.952	4.8	100	0.00
8		Nitrobenzene	1.000	0.948	5.2	100	0.00
9		Naphthalene	1.000	0.966	3.4	100	0.00
10		Hexachlorobutadiene	1.000	0.965	3.5	101	0.00
11		2-Methylnaphthalene	1.000	0.953	4.7	99	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	96	0.00
13	S	2,4,6-Tribromophenol	1.000	0.940	6.0	103	0.00
14	S	2-Fluorobiphenyl	1.000	0.974	2.6	97	0.00
15		Acenaphthylene	1.000	0.963	3.7	96	0.00
16		Acenaphthene	1.000	0.937	6.3	94	0.00
17		Fluorene	1.000	0.982	1.8	102	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	71	0.00
19		4-Bromophenyl-phenylether	1.000	1.180	-18.0	104	0.00
20		Hexachlorobenzene	1.000	1.328	-32.8#	105	0.00
21		Pentachlorophenol	1.000	1.197	-19.7	115	0.00
22		Phenanthrene	1.000	1.021	-2.1	77	0.00
23		Anthracene	1.000	1.254	-25.4#	104	0.00
24		Fluoranthene	1.000	1.210	-21.0	95	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	101	0.00
26		Pyrene	1.000	0.935	6.5	96	0.00
27	S	Terphenyl-d14	1.000	0.916	8.4	93	0.00
28		Benzo(a)anthracene	1.000	0.916	8.4	97	0.00
29		Chrysene	1.000	0.935	6.5	98	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.873	12.7	98	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.883	11.7	96	0.01
32	I	Perylene-d12	5.000	5.000	0.0	99	0.00
33		Benzo(b)fluoranthene	1.000	0.899	10.1	98	0.00
34		Benzo(k)fluoranthene	1.000	0.992	0.8	98	0.01
35	C	Benzo(a)pyrene	1.000	0.936	6.4	98	0.01
36		Dibenzo(a,h)anthracene	1.000	0.890	11.0	96	0.01
37		Benzo(g,h,i)perylene	1.000	0.897	10.3	97	0.00

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