Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM061915\
Data File : BM001748.D
Acq On : 18 Jun 2015 21:59
Operator : TP/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM061915

Quant Time: Jun 19 05:13:28 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM061915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 19 05:05:14 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	0.974	2.6	113	0.00
3		n-Nitrosodimethylamine	1.000	0.981	1.9	117	-0.02
4	S	2-Fluorophenol	1.000	0.943	5.7	113	0.00
5	S	Phenol-d6	1.000	0.929	7.1	114	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	120	0.00
7	S	Nitrobenzene-d5	1.000	0.930	7.0	114	0.00
8		Nitrobenzene	1.000	0.937	6.3	114	0.00
9		Naphthalene	1.000	0.942	5.8	114	0.00
10		Hexachlorobutadiene	1.000	0.949	5.1	117	0.00
11		2-Methylnaphthalene	1.000	0.950	5.0	115	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	122	-0.01
13	S	2,4,6-Tribromophenol	1.000	0.972	2.8	122	0.00
14	S	2-Fluorobiphenyl	1.000	0.942	5.8	115	0.00
15		Acenaphthylene	1.000	0.936	6.4	116	0.00
16		Acenaphthene	1.000	0.939	6.1	116	0.00
17		Fluorene	1.000	0.994	0.6	109	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	114	0.00
19		4-Bromophenyl-phenylether	1.000	1.013	-1.3	128	0.00
20		Hexachlorobenzene	1.000	0.929	7.1	114	0.00
21		Pentachlorophenol	1.000	0.933	6.7	115	0.00
22		Phenanthrene	1.000	0.904	9.6	109	0.00
23		Anthracene	1.000	0.910	9.0	103	0.00
24		Fluoranthene	1.000	0.963	3.7	114	-0.01
25	I	Chrysene-d12	5.000	5.000	0.0	117	0.00
26		Pyrene	1.000	0.931	6.9	115	-0.01
27	S	Terphenyl-d14	1.000	0.945	5.5	115	0.00
28		Benzo(a)anthracene	1.000	0.935	6.5	113	0.00
29		Chrysene	1.000	0.964	3.6	115	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.924	7.6	118	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.937	6.3	110	0.00
32	I	Perylene-d12	5.000	5.000	0.0	117	0.00
33		Benzo(b)fluoranthene	1.000	0.974	2.6	115	-0.01
34		Benzo(k)fluoranthene	1.000	0.904	9.6	108	0.00
35	C	Benzo(a)pyrene	1.000	0.941	5.9	112	0.00
36		Dibenzo(a,h)anthracene	1.000	0.940	6.0	110	0.00
37		Benzo(g,h,i)perylene	1.000	0.938	6.2	111	0.00

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