Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061215\
Data File : BM001660.D
Acq On : 12 Jun 2015 03:38
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Jun 12 16:48:28 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM061215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 12 16:27:37 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	89	0.00
2		1,4-Dioxane	1.000	1.021	-2.1	90	0.00
3		n-Nitrosodimethylamine	1.000	1.013	-1.3	92	0.00
4	S	2-Fluorophenol	1.000	0.943	5.7	88	0.00
5	S	Phenol-d6	1.000	0.946	5.4	89	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	90	0.00
7	S	Nitrobenzene-d5	1.000	0.928	7.2	88	0.00
8		Nitrobenzene	1.000	0.925	7.5	88	0.00
9		Naphthalene	1.000	0.942	5.8	90	0.00
10		Hexachlorobutadiene	1.000	0.954	4.6	93	0.00
11		2-Methylnaphthalene	1.000	0.963	3.7	92	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	92	0.00
13	S	2,4,6-Tribromophenol	1.000	0.994	0.6	90	0.00
14	S	2-Fluorobiphenyl	1.000	0.955	4.5	93	0.00
15		Acenaphthylene	1.000	0.957	4.3	92	-0.01
16		Acenaphthene	1.000	0.939	6.1	92	0.00
17		Fluorene	1.000	1.012	-1.2	99	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	92	0.00
19		4-Bromophenyl-phenylether	1.000	0.967	3.3	87	0.00
20		Hexachlorobenzene	1.000	1.095	-9.5	101	0.00
21		Pentachlorophenol	1.000	0.994	0.6	108	0.00
22		Phenanthrene	1.000	0.877	12.3	86	0.00
23		Anthracene	1.000	1.061	-6.1	98	0.00
24		Fluoranthene	1.000	0.972	2.8	92	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	90	0.00
26		Pyrene	1.000	0.974	2.6	91	0.00
27	S	Terphenyl-d14	1.000	0.967	3.3	89	-0.01
28		Benzo(a)anthracene	1.000	0.878	12.2	82	0.00
29		Chrysene	1.000	0.984	1.6	92	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.992	0.8	95	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.820	18.0	81	0.00
32	I	Perylene-d12	5.000	5.000	0.0	83	0.00
33		Benzo(b)fluoranthene	1.000	0.969	3.1	86	0.00
34		Benzo(k)fluoranthene	1.000	0.940	6.0	82	0.00
35	C	Benzo(a)pyrene	1.000	0.928	7.2	82	0.00
36		Dibenzo(a,h)anthracene	1.000	0.880	12.0	80	0.00
37		Benzo(g,h,i)perylene	1.000	0.992	0.8	84	0.00

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