Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM062215\
Data File : BM001777.D
Acq On : 22 Jun 2015 17:48
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Jun 23 04:43:58 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	128	0.00
2		1,4-Dioxane	1.000	0.983	1.7	122	0.00
3		n-Nitrosodimethylamine	1.000	1.015	-1.5	129	-0.02
4	S	2-Fluorophenol	1.000	1.031	-3.1	132	0.00
5	S	Phenol-d6	1.000	1.045	-4.5	136	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	132	-0.02
7	S	Nitrobenzene-d5	1.000	1.003	-0.3	135	-0.01
8		Nitrobenzene	1.000	1.006	-0.6	135	-0.01
9		Naphthalene	1.000	0.977	2.3	130	0.00
10		Hexachlorobutadiene	1.000	0.957	4.3	130	0.00
11		2-Methylnaphthalene	1.000	0.998	0.2	134	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	136	-0.01
13	S	2,4,6-Tribromophenol	1.000	1.176	-17.6	164	0.00
14	S	2-Fluorobiphenyl	1.000	0.973	2.7	133	0.00
15		Acenaphthylene	1.000	1.011	-1.1	140	0.00
16		Acenaphthene	1.000	0.983	1.7	136	0.00
17		Fluorene	1.000	0.954	4.6	116	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	120	0.00
19		4-Bromophenyl-phenylether	1.000	1.229	-22.9	163	0.00
20		Hexachlorobenzene	1.000	0.944	5.6	121	0.00
21		Pentachlorophenol	1.000	1.079	-7.9	139	0.00
22		Phenanthrene	1.000	1.031	-3.1	130	0.00
23		Anthracene	1.000	0.986	1.4	118	0.00
24		Fluoranthene	1.000	1.098	-9.8	137	-0.01
25	I	Chrysene-d12	5.000	5.000	0.0	133	0.00
26		Pyrene	1.000	0.985	1.5	139	-0.01
27	S	Terphenyl-d14	1.000	1.023	-2.3	142	0.00
28		Benzo(a)anthracene	1.000	0.968	3.2	133	0.00
29		Chrysene	1.000	0.949	5.1	129	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.027	-2.7	149	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.961	3.9	128	0.00
32	I	Perylene-d12	5.000	5.000	0.0	131	0.00
33		Benzo(b)fluoranthene	1.000	0.974	2.6	129	0.00
34		Benzo(k)fluoranthene	1.000	0.990	1.0	133	0.00
35	C	Benzo(a)pyrene	1.000	1.003	-0.3	134	0.01
36		Dibenzo(a,h)anthracene	1.000	0.970	3.0	127	0.00
37		Benzo(g,h,i)perylene	1.000	0.962	3.8	127	0.00

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