Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061215\
Data File : BM001675.D
Acq On : 12 Jun 2015 13:17
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Jun 12 18:09:23 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	105	0.00
2		1,4-Dioxane	1.000	0.975	2.5	102	0.00
3		n-Nitrosodimethylamine	1.000	0.992	0.8	106	0.00
4	S	2-Fluorophenol	1.000	1.026	-2.6	113	0.00
5	S	Phenol-d6	1.000	1.045	-4.5	116	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	107	0.00
7	S	Nitrobenzene-d5	1.000	1.059	-5.9	120	0.00
8		Nitrobenzene	1.000	1.043	-4.3	118	0.00
9		Naphthalene	1.000	0.950	5.0	108	0.00
10		Hexachlorobutadiene	1.000	0.943	5.7	109	0.00
11		2-Methylnaphthalene	1.000	0.975	2.5	111	-0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	110	0.00
13	S	2,4,6-Tribromophenol	1.000	1.243	-24.3	135	0.00
14	S	2-Fluorobiphenyl	1.000	0.950	5.0	111	0.00
15		Acenaphthylene	1.000	1.018	-1.8	117	-0.01
16		Acenaphthene	1.000	0.959	4.1	113	0.00
17		Fluorene	1.000	1.042	-4.2	122	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	109	0.00
19		4-Bromophenyl-phenylether	1.000	0.946	5.4	102	0.00
20		Hexachlorobenzene	1.000	1.071	-7.1	118	0.00
21		Pentachlorophenol	1.000	1.097	-9.7	148	0.00
22		Phenanthrene	1.000	0.859	14.1	101	0.00
23		Anthracene	1.000	1.135	-13.5	125	0.00
24		Fluoranthene	1.000	0.949	5.1	108	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	93	0.00
26		Pyrene	1.000	1.074	-7.4	104	0.00
27	S	Terphenyl-d14	1.000	1.073	-7.3	103	0.00
28		Benzo(a)anthracene	1.000	0.982	1.8	95	0.00
29		Chrysene	1.000	0.964	3.6	94	0.02
30		Bis(2-ethylhexyl)phthalate	1.000	1.360	-36.0#	145	0.02
31		Indeno(1,2,3-cd)pyrene	1.000	1.001	-0.1	102	0.01
32	I	Perylene-d12	5.000	5.000	0.0	91	0.01
33		Benzo(b)fluoranthene	1.000	0.979	2.1	95	0.01
34		Benzo(k)fluoranthene	1.000	0.937	6.3	89	0.01
35	C	Benzo(a)pyrene	1.000	1.014	-1.4	98	0.01
36		Dibenzo(a,h)anthracene	1.000	1.050	-5.0	104	0.00
37		Benzo(g,h,i)perylene	1.000	1.089	-8.9	101	0.01

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