Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM050415\
Data File : BM001205.D
Acq On : 05 May 2015 04:00
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: May 05 05:05:34 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM050115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon May 04 16:48:07 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	129	0.00
2		1,4-Dioxane	1.000	0.872	12.8	107	0.00
3		n-Nitrosodimethylamine	1.000	0.933	6.7	114	0.00
4	S	2-Fluorophenol	1.000	1.094	-9.4	139	-0.01
5	S	Phenol-d6	1.000	1.159	-15.9	148	0.02
6	I	Naphthalene-d8	5.000	5.000	0.0	134	0.00
7	S	Nitrobenzene-d5	1.000	1.064	-6.4	143	-0.01
8		Nitrobenzene	1.000	0.986	1.4	133	-0.01
9		Naphthalene	1.000	1.000	0.0	131	0.00
10		Hexachlorobutadiene	1.000	0.998	0.2	134	0.00
11		2-Methylnaphthalene	1.000	1.042	-4.2	139	-0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	143	-0.02
13	S	2,4,6-Tribromophenol	1.000	1.106	-10.6	192	0.00
14	S	2-Fluorobiphenyl	1.000	0.996	0.4	139	-0.01
15		Acenaphthylene	1.000	1.080	-8.0	152	-0.02
16		Acenaphthene	1.000	1.060	-6.0	153	0.00
17		Fluorene	1.000	1.130	-13.0	155	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	152	0.00
19		Hexachlorobenzene	1.000	0.952	4.8	148	0.00
20		Pentachlorophenol	1.000	1.130	-13.0	222	0.00
21		Phenanthrene	1.000	0.979	2.1	150	0.00
22		Anthracene	1.000	1.030	-3.0	154	-0.02
23		Fluoranthene	1.000	1.023	-2.3	156	-0.01
24	I	Chrysene-d12	5.000	5.000	0.0	148	0.00
25		Pyrene	1.000	1.050	-5.0	156	-0.01
26	S	Terphenyl-d14	1.000	1.052	-5.2	154	-0.01
27		Benzo(a)anthracene	1.000	1.045	-4.5	156	0.00
28		Chrysene	1.000	1.017	-1.7	146	0.00
29		Indeno(1,2,3-cd)pyrene	1.000	0.890	11.0	128	-0.01
30	I	Perylene-d12	5.000	5.000	0.0	137	-0.01
31		Benzo(b)fluoranthene	1.000	1.104	-10.4	140	-0.01
32		Benzo(k)fluoranthene	1.000	1.003	-0.3	146	-0.01
33	C	Benzo(a)pyrene	1.000	1.058	-5.8	144	-0.01
34		Dibenzo(a,h)anthracene	1.000	0.960	4.0	129	-0.02
35		Benzo(g,h,i)perylene	1.000	0.932	6.8	125	-0.01

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