Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM050515\
Data File : BM001221.D
Acq On : 06 May 2015 10:39
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: May 06 13:13:27 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM050515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 06 12:46:35 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	94	0.00
2		1,4-Dioxane	1.000	0.929	7.1	84	0.00
3		n-Nitrosodimethylamine	1.000	0.892	10.8	93	0.00
4	S	2-Fluorophenol	1.000	1.011	-1.1	93	0.00
5	S	Phenol-d6	1.000	0.970	3.0	88	-0.02
6	I	Naphthalene-d8	5.000	5.000	0.0	91	0.00
7	S	Nitrobenzene-d5	1.000	0.989	1.1	87	0.00
8		Nitrobenzene	1.000	0.950	5.0	89	0.00
9		Naphthalene	1.000	1.027	-2.7	91	0.00
10		Hexachlorobutadiene	1.000	1.027	-2.7	94	0.00
11		2-Methylnaphthalene	1.000	1.006	-0.6	89	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	86	0.00
13	S	2,4,6-Tribromophenol	1.000	0.896	10.4	83	0.00
14	S	2-Fluorobiphenyl	1.000	1.046	-4.6	88	-0.01
15		Acenaphthylene	1.000	1.007	-0.7	85	0.00
16		Acenaphthene	1.000	1.031	-3.1	87	0.00
17		Fluorene	1.000	1.004	-0.4	81	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	83	0.00
19		Hexachlorobenzene	1.000	1.103	-10.3	92	0.00
20		Pentachlorophenol	1.000	1.019	-1.9	81	0.00
21		Phenanthrene	1.000	0.987	1.3	80	0.00
22		Anthracene	1.000	0.968	3.2	82	0.00
23		Fluoranthene	1.000	1.073	-7.3	86	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	92	0.00
25		Pyrene	1.000	0.954	4.6	88	0.00
26	S	Terphenyl-d14	1.000	0.950	5.0	86	0.00
27		Benzo(a)anthracene	1.000	0.991	0.9	91	0.00
28		Chrysene	1.000	1.043	-4.3	94	0.00
29		Indeno(1,2,3-cd)pyrene	1.000	1.226	-22.6	110	0.00
30	I	Perylene-d12	5.000	5.000	0.0	103	0.00
31		Benzo(b)fluoranthene	1.000	1.045	-4.5	111	0.01
32		Benzo(k)fluoranthene	1.000	0.939	6.1	90	0.01
33	C	Benzo(a)pyrene	1.000	1.013	-1.3	103	0.00
34		Dibenzo(a,h)anthracene	1.000	1.074	-7.4	109	0.00
35		Benzo(g,h,i)perylene	1.000	1.102	-10.2	112	0.00

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