Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM051615\
Data File : BM001311.D
Acq On : 15 May 2015 02:43
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: May 15 06:13:39 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIMPAH-BM051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 13 05:24:15 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	96	-0.02
2		1,4-Dioxane	1.000	0.528	47.2#	70	0.00
3		n-Nitrosodimethylamine	1.000	1.031	-3.1	113	0.00
4	S	2-Fluorophenol	1.000	0.950	5.0	96	-0.02
5	S	Phenol-d6	1.000	0.937	6.3	97	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	97	0.00
7	S	Nitrobenzene-d5	1.000	0.871	12.9	95	0.00
8		Nitrobenzene	1.000	0.851	14.9	95	0.00
9		Naphthalene	1.000	0.961	3.9	96	0.00
10		Hexachlorobutadiene	1.000	0.937	6.3	96	0.00
11		2-Methylnaphthalene	1.000	0.938	6.2	97	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	95	0.00
13	S	2,4,6-Tribromophenol	1.000	0.689	31.1#	80	0.00
14	S	2-Fluorobiphenyl	1.000	0.954	4.6	96	0.00
15		Acenaphthylene	1.000	0.934	6.6	96	0.00
16		Acenaphthene	1.000	0.933	6.7	96	0.00
17		Fluorene	1.000	0.880	12.0	95	-0.03
18	I	Phenanthrene-d10	5.000	5.000	0.0	126	0.00
19		Hexachlorobenzene	1.000	0.668	33.2#	91	0.00
20		Pentachlorophenol	1.000	0.915	8.5	131	0.00
21		Phenanthrene	1.000	1.057	-5.7	145	0.00
22		Anthracene	1.000	0.974	2.6	134	0.00
23		Fluoranthene	1.000	0.701	29.9#	98	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	99	-0.01
25		Pyrene	1.000	0.909	9.1	98	0.00
26	S	Terphenyl-d14	1.000	0.888	11.2	94	-0.01
27		Benzo(a)anthracene	1.000	0.933	6.7	104	-0.01
28		Chrysene	1.000	0.927	7.3	97	-0.01
29		Indeno(1,2,3-cd)pyrene	1.000	1.033	-3.3	109	-0.01
30	I	Perylene-d12	5.000	5.000	0.0	103	-0.02
31		Benzo(b)fluoranthene	1.000	0.937	6.3	105	-0.01
32		Benzo(k)fluoranthene	1.000	0.913	8.7	100	-0.01
33	C	Benzo(a)pyrene	1.000	0.931	6.9	105	-0.01
34		Dibenzo(a,h)anthracene	1.000	0.991	0.9	108	-0.02
35		Benzo(g,h,i)perylene	1.000	0.987	1.3	109	-0.02

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