Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM051915\
Data File : BM001342.D
Acq On : 19 May 2015 10:54
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: May 20 03:59:56 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	78	-0.02
2		1,4-Dioxane	1.000	0.859	14.1	71	0.00
3		n-Nitrosodimethylamine	1.000	0.950	5.0	85	0.00
4	S	2-Fluorophenol	1.000	0.998	0.2	82	-0.02
5	S	Phenol-d6	1.000	0.985	1.5	83	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	78	0.00
7	S	Nitrobenzene-d5	1.000	0.921	7.9	82	0.00
8		Nitrobenzene	1.000	0.905	9.5	82	0.00
9		Naphthalene	1.000	1.011	-1.1	82	0.00
10		Hexachlorobutadiene	1.000	0.999	0.1	83	0.00
11		2-Methylnaphthalene	1.000	0.984	1.6	82	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	76	0.00
13	S	2,4,6-Tribromophenol	1.000	0.702	29.8#	65	0.00
14	S	2-Fluorobiphenyl	1.000	1.010	-1.0	81	0.00
15		Acenaphthylene	1.000	0.956	4.4	79	0.00
16		Acenaphthene	1.000	0.979	2.1	80	0.00
17		Fluorene	1.000	0.955	4.5	82	-0.03
18	I	Phenanthrene-d10	5.000	5.000	0.0	101	0.00
19		Hexachlorobenzene	1.000	0.699	30.1#	77	0.00
20		Pentachlorophenol	1.000	1.002	-0.2	124	0.00
21		Phenanthrene	1.000	1.080	-8.0	119	0.00
22		Anthracene	1.000	1.000	0.0	111	0.00
23		Fluoranthene	1.000	0.730	27.0#	82	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	81	-0.01
25		Pyrene	1.000	0.937	6.3	83	0.00
26	S	Terphenyl-d14	1.000	0.943	5.7	82	-0.01
27		Benzo(a)anthracene	1.000	0.942	5.8	86	-0.01
28		Chrysene	1.000	0.985	1.5	85	-0.01
29		Indeno(1,2,3-cd)pyrene	1.000	1.041	-4.1	91	-0.01
30	I	Perylene-d12	5.000	5.000	0.0	86	-0.01
31		Benzo(b)fluoranthene	1.000	0.922	7.8	86	-0.01
32		Benzo(k)fluoranthene	1.000	0.993	0.7	91	-0.01
33	C	Benzo(a)pyrene	1.000	0.949	5.1	89	0.00
34		Dibenzo(a,h)anthracene	1.000	0.984	1.6	90	-0.01
35		Benzo(g,h,i)perylene	1.000	1.001	-0.1	92	-0.02

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