Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM051915\
Data File : BM001344.D
Acq On : 19 May 2015 12:26
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: May 20 04:10:26 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 20 03:51:55 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	96	-0.02
2		1,4-Dioxane	0.582	0.623	-7.0	95	0.00
3		n-Nitrosodimethylamine	0.566	0.572	-1.1	111	0.00
4	S	2-Fluorophenol	0.974	0.962	1.2	100	-0.02
5	S	Phenol-d6	1.099	1.107	-0.7	105	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	99	0.00
7	S	Nitrobenzene-d5	0.263	0.265	-0.8	106	0.00
8		Nitrobenzene	0.276	0.274	0.7	107	0.00
9		Naphthalene	1.064	1.078	-1.3	104	0.00
10		Hexachlorobutadiene	0.208	0.210	-1.0	106	0.00
11		2-Methylnaphthalene	0.692	0.716	-3.5	109	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	103	0.00
13	S	2,4,6-Tribromophenol	0.149	0.108	27.5#	90	0.00
14	S	2-Fluorobiphenyl	1.552	1.588	-2.3	111	0.00
15		Acenaphthylene	2.058	2.030	1.4	110	0.00
16		Acenaphthene	1.318	1.302	1.2	110	0.00
17		Fluorene	1.194	1.106	7.4	108	-0.03
18	I	Phenanthrene-d10	1.000	1.000	0.0	129	0.00
19		Hexachlorobenzene	0.322	0.234	27.3#	102	0.00
20		Pentachlorophenol	0.067	0.052	22.4	140	0.00
21		Phenanthrene	0.760	0.848	-11.6	157#	0.00
22		Anthracene	0.741	0.761	-2.7	145	0.00
23		Fluoranthene	1.193	0.820	31.3#	98	0.00
24	I	Chrysene-d12	1.000	1.000	0.0	88	-0.01
25		Pyrene	1.545	1.575	-1.9	99	0.00
26	S	Terphenyl-d14	0.672	0.702	-4.5	99	-0.01
27		Benzo(a)anthracene	1.357	1.279	5.7	94	0.00
28		Chrysene	1.376	1.355	1.5	92	-0.01
29		Indeno(1,2,3-cd)pyrene	1.398	1.349	3.5	92	-0.01
30	I	Perylene-d12	1.000	1.000	0.0	88	-0.01
31		Benzo(b)fluoranthene	1.465	1.535	-4.8	100	0.00
32		Benzo(k)fluoranthene	1.394	1.235	11.4	83	-0.01
33	C	Benzo(a)pyrene	1.269	1.197	5.7	91	0.00
34		Dibenzo(a,h)anthracene	1.193	1.167	2.2	91	-0.01
35		Benzo(g,h,i)perylene	1.309	1.295	1.1	93	-0.02

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