Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM051815\
Data File : BM001340.D
Acq On : 18 May 2015 20:59
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: May 19 02:32:58 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue May 19 02:11:49 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	101	0.00
2		1,4-Dioxane	0.582	0.411	29.4#	66	-0.01
3		n-Nitrosodimethylamine	0.566	0.525	7.2	107	0.00
4	S	2-Fluorophenol	0.974	0.913	6.3	100	-0.02
5	S	Phenol-d6	1.099	1.065	3.1	106	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	102	0.00
7	S	Nitrobenzene-d5	0.263	0.257	2.3	106	0.00
8		Nitrobenzene	0.276	0.260	5.8	105	0.00
9		Naphthalene	1.064	1.026	3.6	102	0.00
10		Hexachlorobutadiene	0.208	0.198	4.8	104	0.00
11		2-Methylnaphthalene	0.692	0.662	4.3	104	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	101	0.00
13	S	2,4,6-Tribromophenol	0.149	0.105	29.5#	86	0.00
14	S	2-Fluorobiphenyl	1.552	1.494	3.7	103	0.00
15		Acenaphthylene	2.058	1.950	5.2	103	0.00
16		Acenaphthene	1.318	1.248	5.3	103	0.00
17		Fluorene	1.194	1.011	15.3	97	-0.03
18	I	Phenanthrene-d10	1.000	1.000	0.0	130	0.00
19		Hexachlorobenzene	0.322	0.230	28.6#	100	0.00
20		Pentachlorophenol	0.067	0.049	26.9#	134	0.00
21		Phenanthrene	0.760	0.729	4.1	136	0.00
22		Anthracene	0.741	0.687	7.3	132	0.00
23		Fluoranthene	1.193	0.840	29.6#	101	0.00
24	I	Chrysene-d12	1.000	1.000	0.0	101	-0.01
25		Pyrene	1.545	1.435	7.1	103	0.00
26	S	Terphenyl-d14	0.672	0.642	4.5	103	-0.01
27		Benzo(a)anthracene	1.357	1.253	7.7	105	-0.01
28		Chrysene	1.376	1.303	5.3	102	-0.01
29		Indeno(1,2,3-cd)pyrene	1.398	1.317	5.8	102	-0.01
30	I	Perylene-d12	1.000	1.000	0.0	101	-0.01
31		Benzo(b)fluoranthene	1.465	1.327	9.4	100	-0.01
32		Benzo(k)fluoranthene	1.394	1.361	2.4	105	-0.01
33	C	Benzo(a)pyrene	1.269	1.180	7.0	103	-0.01
34		Dibenzo(a,h)anthracene	1.193	1.132	5.1	102	-0.01
35		Benzo(g,h,i)perylene	1.309	1.233	5.8	102	-0.02

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