Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM051315\
Data File : BM001274.D
Acq On : 13 May 2015 05:04
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: May 13 06:46:32 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	94	0.00
2		1,4-Dioxane	0.582	0.644	-10.7	97	0.00
3		n-Nitrosodimethylamine	0.566	0.500	11.7	96	0.00
4	S	2-Fluorophenol	0.974	0.923	5.2	95	0.00
5	S	Phenol-d6	1.099	1.029	6.4	96	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	94	0.00
7	S	Nitrobenzene-d5	0.263	0.247	6.1	94	0.00
8		Nitrobenzene	0.276	0.254	8.0	94	0.00
9		Naphthalene	1.064	1.035	2.7	95	0.00
10		Hexachlorobutadiene	0.208	0.198	4.8	95	0.00
11		2-Methylnaphthalene	0.692	0.655	5.3	95	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	93	0.00
13	S	2,4,6-Tribromophenol	0.149	0.114	23.5	86	0.00
14	S	2-Fluorobiphenyl	1.552	1.487	4.2	94	0.00
15		Acenaphthylene	2.058	1.919	6.8	93	0.00
16		Acenaphthene	1.318	1.244	5.6	94	0.00
17		Fluorene	1.194	0.993	16.8	87	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	112	0.00
19		Hexachlorobenzene	0.322	0.254	21.1	96	0.00
20		Pentachlorophenol	0.067	0.050	25.4#	117	0.00
21		Phenanthrene	0.760	0.622	18.2	100	0.00
22		Anthracene	0.741	0.610	17.7	101	0.00
23		Fluoranthene	1.193	0.897	24.8	94	0.00
24	I	Chrysene-d12	1.000	1.000	0.0	97	-0.01
25		Pyrene	1.545	1.406	9.0	97	0.00
26	S	Terphenyl-d14	0.672	0.608	9.5	94	-0.01
27		Benzo(a)anthracene	1.357	1.206	11.1	98	0.00
28		Chrysene	1.376	1.302	5.4	98	-0.01
29		Indeno(1,2,3-cd)pyrene	1.398	1.387	0.8	103	0.00
30	I	Perylene-d12	1.000	1.000	0.0	100	-0.01
31		Benzo(b)fluoranthene	1.465	1.421	3.0	106	0.00
32		Benzo(k)fluoranthene	1.394	1.194	14.3	91	0.00
33	C	Benzo(a)pyrene	1.269	1.153	9.1	100	0.00
34		Dibenzo(a,h)anthracene	1.193	1.154	3.3	103	-0.01
35		Benzo(g,h,i)perylene	1.309	1.279	2.3	104	-0.01

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