Table SI. Chemical shifts in ppm (^{1}H-NMR) for compounds 2-8c in $CDCl_3$ or DMSO-d_6.

Compound	H-4	H-5	H-7	H-13	H-15	H-16
2[a]	8.53	7.62	7.63	-	-	-
6[a]	8.11	7.63	7.70-7.64	7.91	7.70-7.64	-
3[b]	8.77	8.13	7.91	-	-	-
7[b]	8.53	8.10	7.90	8.08	7.92	7.52
4a[b]	8.93	8.47	8.03	-	-	-
4c[a]	8.89	8.1	8.01	-	-	-
5[a]	-	7.80	7.77	-	-	-
8a[b]	8.54	8.33	8.05	8.22	8.01	7.69
8b[b]	8.98	8.79	8.60	8.85	8.60	8.22
8c[b]	8.98	8.78	8.58	8.85	8.58	8.21

NMR spectra were recorded in a Bruker advance TM-400 spectrometer (400 MHz ^{1}H NMR).
[a]$CDCl_3$; [b]DMSO-d_6.