

Supporting Information

Photovoltaic properties of 3,3'-(ethane-1,2-diylidene)-bis(indolin-2-one) based conjugated polymers

Jian Wu^a, Yingying Fu^b, Hao Huang^a, Shengxia Li^a and Zhiyuan Xie^{b,**} and Qing Zhang^{a,*}

^a*Shanghai Key Lab of Polymer and Electrical Insulation, School of Chemistry and Chemical Engineering, Shanghai Jiaotong University, Shanghai 200240, P. R. China*

^b*State Key Laboratory of Polymer Physics and Chemistry, Changchun Institute of Applied Chemistry, Chinese Academy of Science, Changchun, 130022, China.*

1. NMR spectra

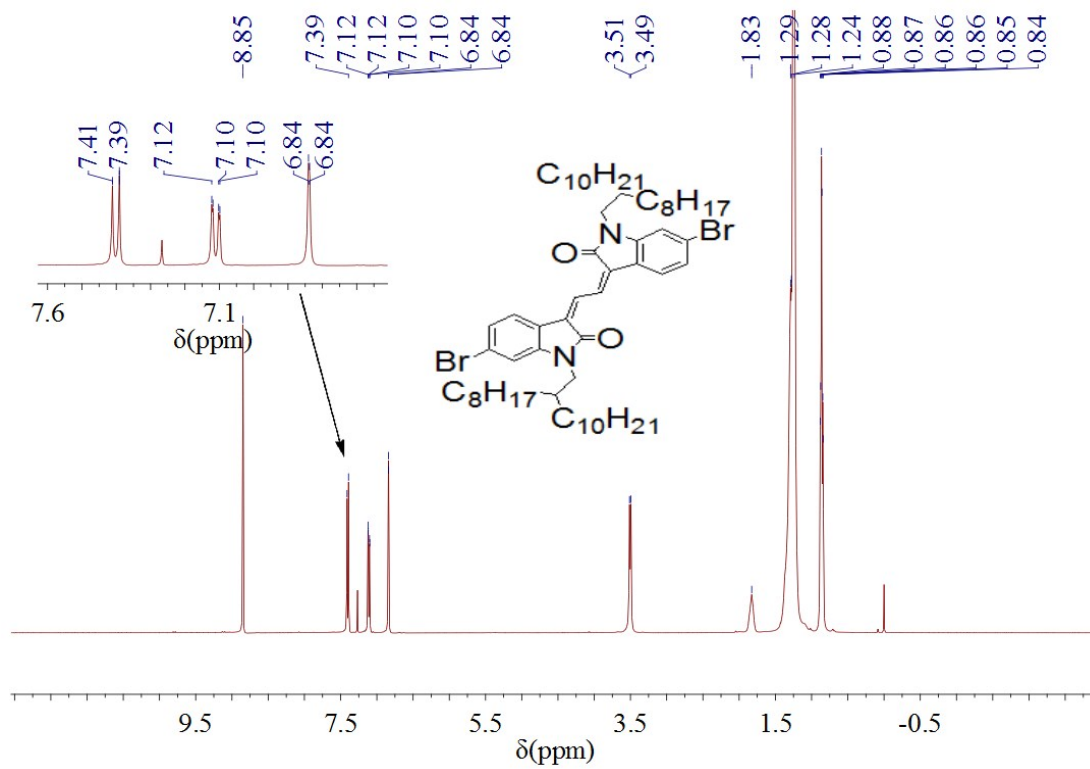


Fig. S1. ^1H NMR spectrum of compound **M**

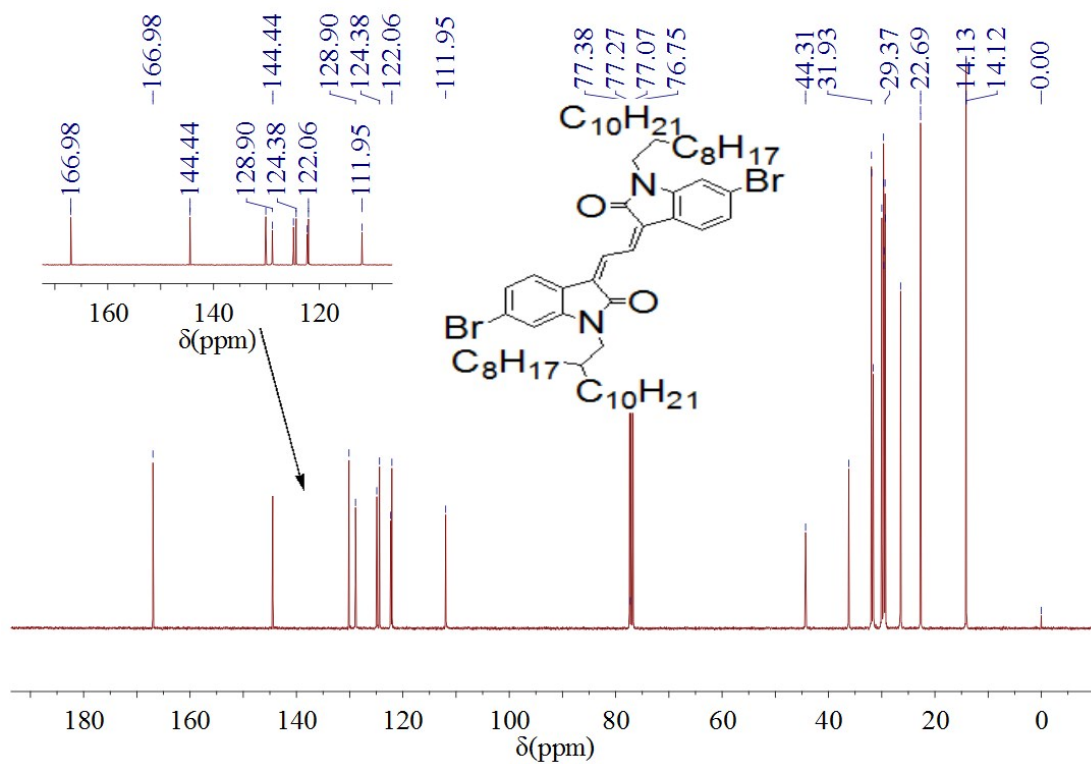


Fig. S2. ^{13}C NMR spectrum of compound **M**

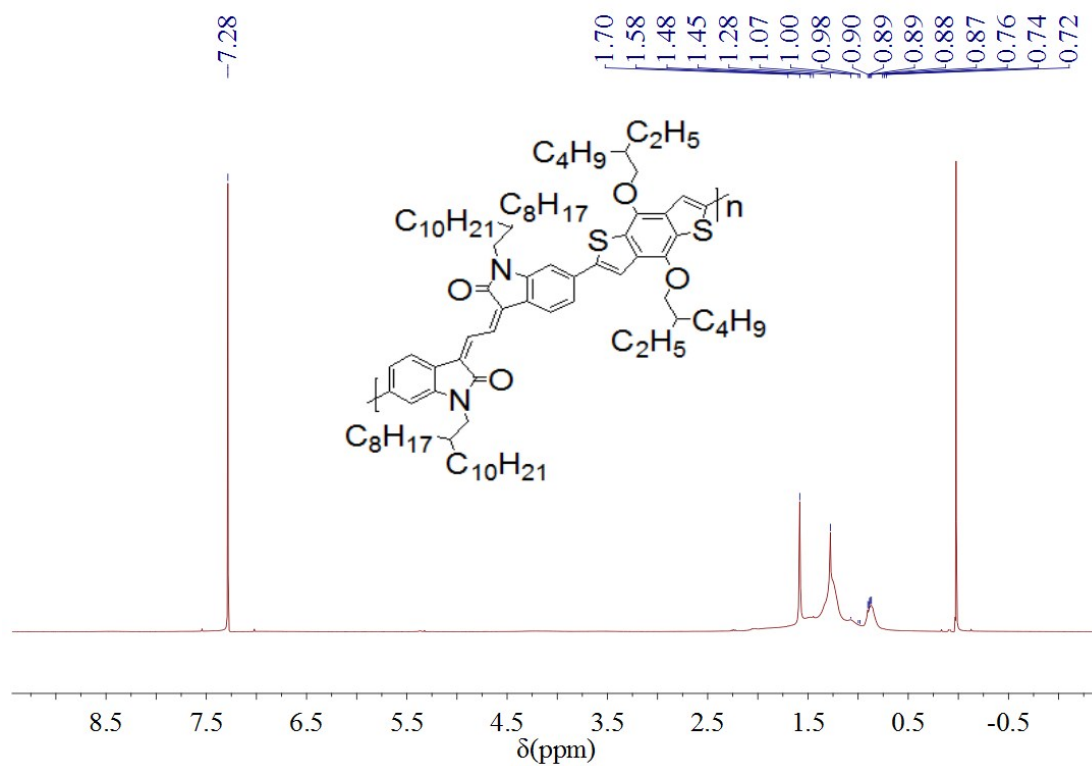


Fig. S3. ^1H NMR spectrum of **PEBI-BDTO**

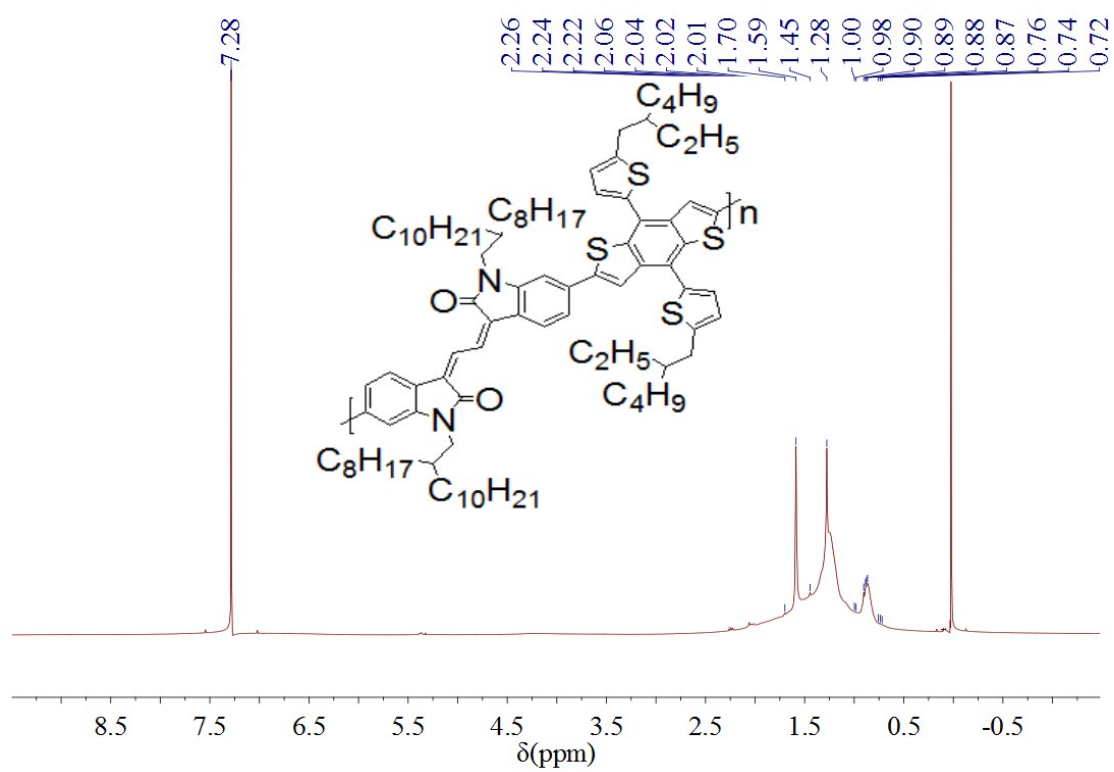


Fig. S4. ^1H NMR spectrum of PEBI-BDT

2. TGA and DSC thermograms of polymers

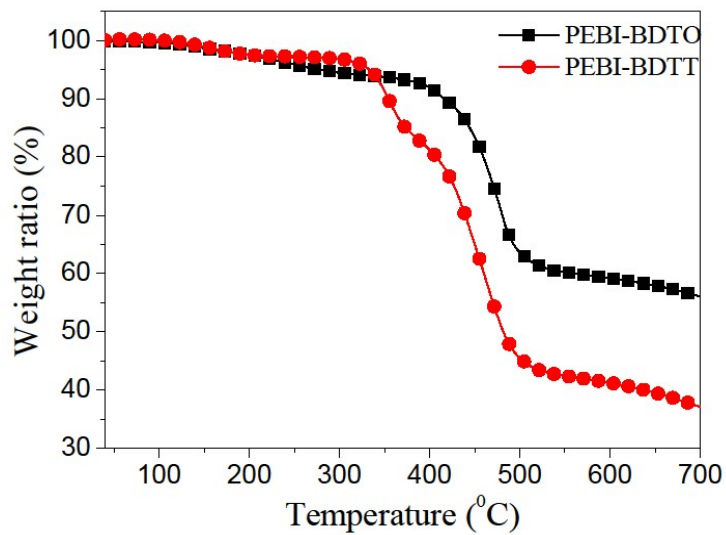


Fig. S5. TGA curves of polymers.

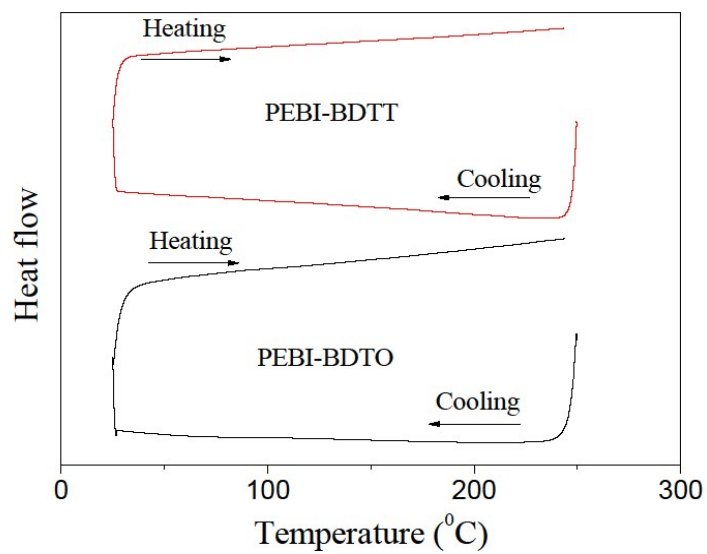


Fig. S6. DSC thermograms of polymers

The first cooling and the second heating scan of polymers with temperature ramp of 20 °C/min.

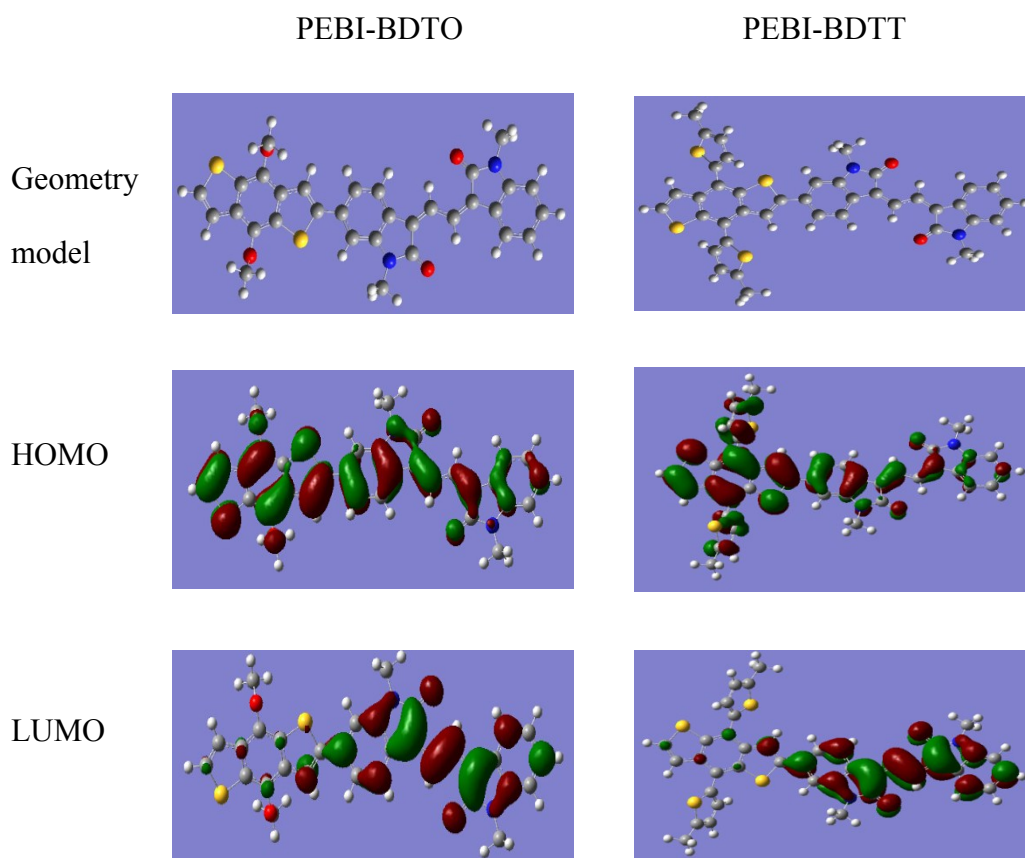


Fig. S7. The geometry, the LUMO and HOMO orbital of model **PEBI-BDTO** (left) and **PEBI-BDTT** (right) by DFT calculations with the B3LYP/6-31G** basis set.

1. Polymer solar cell devices performances

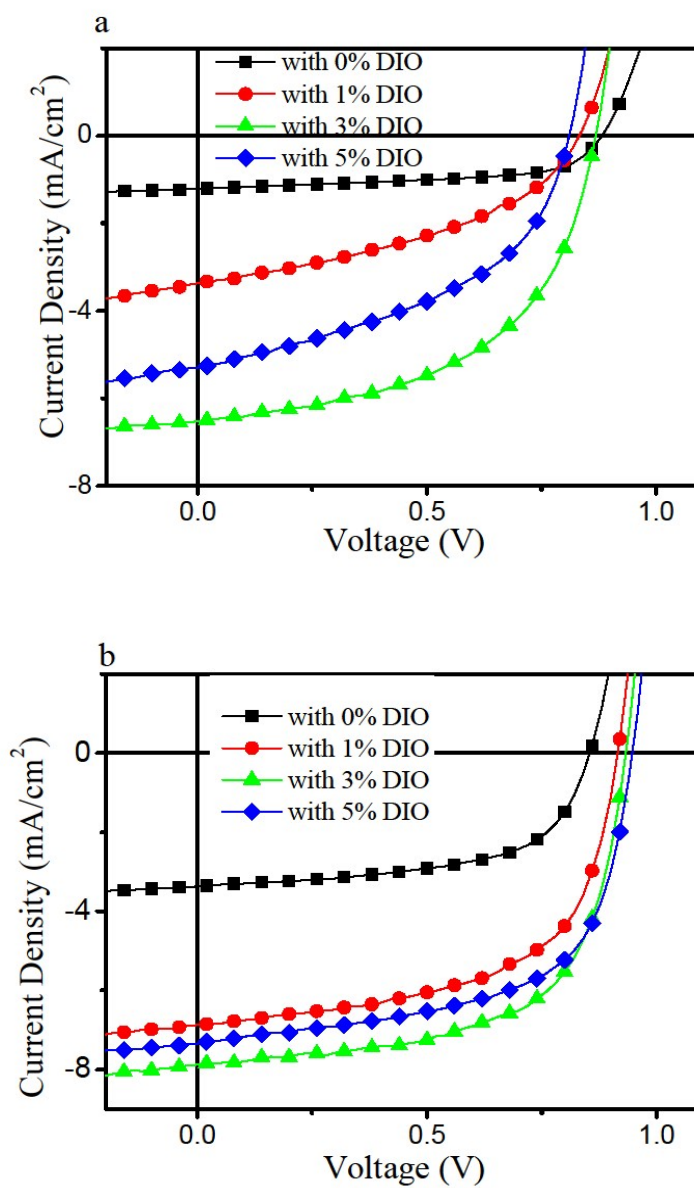


Fig. S8. $I-V$ curves of PSC devices processed with different amount of DIO (a) PEBI-BDTO:PC₇₁BM and (b) PEBI-BDTT:PC₇₁BM

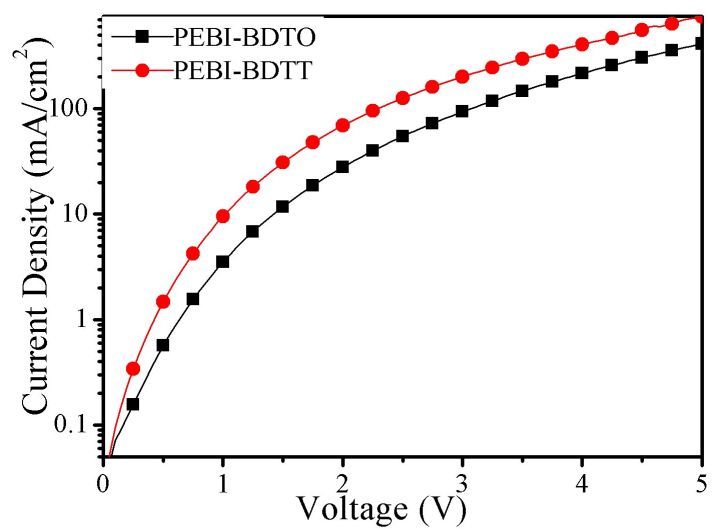


Fig. S9. *I-V* characteristics of **PEBI-BDTO** and **PEBI-BDTT** hole-only devices.

Table.S1. The solar cell performances with different amount of DIO

Active layer (w/w)		DIO(%)	J_{sc}	Voc	FF	PCE (%)
			(mA/cm ²)	(V)		
Polymer:PC ₇₁ BM	D:A					
PEBI-BDTO: PC ₇₁ BM	1:2	0	1.21	0.88	58.7	2.39±0.08
	1:2	1	3.37	0.84	41.2	2.54±0.13
	1:2	3	6.53	0.86	53.4	2.74±0.06
	1:2	5	5.48	0.82	51.7	3.00±0.12
PEBI-BDTT: PC ₇₁ BM	1:1.5	0	3.37	0.86	58.8	1.70±0.17
	1:1.5	1	6.87	0.92	58.5	3.70±0.11
	1:1.5	3	7.88	0.94	62.1	4.59±0.15
	1:1.5	5	7.34	0.94	61.4	4.23±0.07