Supplementary Information

Fluorene-based hyperbranched copolymers with

spiro[3.3]heptane-2,6-dispirofluorene as the conjugation-uninterrupted

branching point and their application in WPLEDs

Yuling Wu,^{ab} Jie Li,^{ab} Wenqing Liang,^{ab} Junli Yang,^{ab} Jing Sun,^{ab} Hua Wang,^{ab} Xuguang Liu,^c Bingshe Xu,^{*ab} Wei Huang,^d

^a Key Laboratory of Interface Science and Engineering in Advanced Materials, Taiyuan University of Technology, Taiyuan, 030024, China.

^b Research Center of Advanced Materials Science and Technology, Taiyuan University of Technology, Taiyuan, 030024, China.

^c College of Chemistry and Chemical Engineering, Taiyuan University of Technology, Taiyuan, 030024, China.

^d Key Laboratory for Organic Electronics & Information Displays (KLOEID) and Institute of Advanced Materials, Nanjing University of Posts and Telecommunications (NUPT), Nanjing, 210046, China.

* Corresponding author. E-mail: Bingshe Xu: xubs@tyut.edu.cn



Figure S1. ¹H NMR spectrum of **TBrSDF**.



Figure S2. ¹³C NMR spectrum of **TBrSDF**.



Figure S3. ¹H NMR spectra of the copolymers.

Copolymer -	n _{TBrSDF} ^a		n _{TBrSDF} ^b	
	feed	found	feed	found
P1	100:0	100:0	0	0
P2	75.75:1	80.39:1	0.01	0.0124
P3	15.75:1	15.88:1	0.05	0.063
P4	8.25:1	8.04:1	0.10	0.1243
P5	4.50:1	4.96:1	0.20	0.2016

Table 1. Polymerization results of the copolymers

^a The integral ratio between the aromatic ring of fluorene and the spiro[3.3]heptane of **TBrSDF**.
^b The mole ratio of **TBrSDF** in copolymers.

The actual contents of spiro[3.3]heptane-2,6-dispirofluorene (**SDF**) were calculated by comparing the peak integral intensities of the proton signials of the spiro[3.3]heptane of **SDF** (δ 3.0-3.5) and the aromatic ring of fluorene (δ 7.4-8.0) of the copolymer and were close to the feed ratios.