

A spirobifluorene and diketopyrrolopyrrole moieties based non-fullerene acceptor for efficient and thermally stable PSCs

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Abstract:

Recently, the exploration of non-fullerene electron acceptors have gained considerable attention and we report herein the design of a new non-fullerene acceptor, $SF(DPPB)_4$, which shows a cross-shaped molecular geometry to help suppress strong intermolecular aggregation. In addition, the energy levels of $SF(DPPB)_4$ match well with those of P3HT, which not only allows achievement of efficient exciton dissociation, but also ensures high V_{oc} (1.14 V) and efficiency (5.16%) in P3HT:SF(DPPB)₄ PSCs. Moreover, the devices show excellent thermal stability. The abovementioned results demonstrate that the new design strategy of employing a high-performance non-fullerene acceptor has promising practical applications.





Fig. 1 (a) Chemical structure of $SF(DPPB)_4$; (b) ground-state geometry of $SF(DPPB)_4$ calculated by DFT method; (c) UV-vis absorption spectra of P3HT film, SF(DPPB)₄ solution and film, and P3HT:SF(DPPB)₄ (2:1, by wt.) blended film; (d) cyclic voltammogram for SF(DPPB)₄ in dichloromethane solution.



Fig. 3 (a) 2D GIWAXS images of P3HT, SF(DPPB)₄ and 2:1 P3HT:SF(DPPB)₄ thin films with or without the thermal annealing at 120 °C for 10 min; (b) intensity profiles along q_z axis of P3HT, $SF(DPPB)_4$ and 2:1 P3HT:SF(DPPB)_4 thin films; (c) corresponding intensity profiles along q_r axis.





Fig. 4 (a) Optical microscopy images showing morphological stability of P3HT:PC₆₁BM (1:0.8, by wt.) and P3HT:SF(DPPB)₄ (2:1, by wt.) blended films after thermal treatment at 150 °C for various time; (b) graph showing the stability of the PCEs for the devices based on P3HT:SF(DPPB)₄ (2:1, by wt.) and P3HT:PC₆₁BM (1:0.8, by wt.) blended films after thermal treatment at 150 °C for various time.

Table 1	The photovoltaic	performances	of the PSCs based	on P3HT:SF(DPPB)
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D:A (w/w)	V _{oc} (V)	J _{sc} (mA cm ⁻²)	FF	PCE (%)
P3HT:SF(DPPB) ₄ (2:1) ^a	1.20	1.91	0.28	0.63 (0.60) ^d
P3HT:SF(DPPB) ₄ (1:1) ^b	1.13	5.78	0.57	3.70 (3.57) ^d
P3HT:SF(DPPB) ₄ (2:1) ^b	1.14	8.29	0.55	5.16 (5.10) ^d
P3HT:SF(DPPB) ₄ (3:1) ^b	1.13	7.13	0.46	3.68 (3.55) ^d
P3HT:PC ₆₁ BM (1:0.8) ^c	0.62	8.27	0.62	3.18 (3.14) ^d

Fig. 2 (a) Energy levels of P3HT, SF(DPPB)₄ and PC₆₁BM; (b) the diagram of device structure; (c) J-V curves of PSCs with the structure of ITO/PEDOT:PSS/P3HT:SF(DPPB)₄/PFN/AI; (d) EQE spectra of P3HT:SF(DPPB)₄ based PSCs.

500

Wavelength (nm)

400

600

700

0.4 0.6 0.8 1.0 1.2 1.4

Voltage (V)

^a As-cast. ^b Annealed at 120 °C for 10 min. ^c Annealed at 170 °C for 10 min. ^d The values in the parentheses are the average PCEs from 10 devices.

Conclusions

In summary, a spirobifluorene core based non-fullerene electron acceptor, SF(DPPB)₄, was designed and synthesized. SF(DPPB)₄ possesses a cruciform molecular configuration, which assures fine phase separation in the active layer of corresponding PSCs. In addition, SF(DPPB)₄ exhibits appropriate absorption bands and matching energy levels for P3HT. Owing to the good morphology and an outstandingly high V_{oc} of 1.14 V, the PSCs based on the P3HT:SF(DPPB)_4 blended films provided the best PCE of 5.16%. To the best of our knowledge, this PCE is one of the highest values reported in the literature to date for P3HT-based fullerene-free PSCs. The P3HT:SF(DPPB)₄ based devices also show excellent thermal stability upon thermal treatment at 150 °C for up to 3 h, demonstrating that SF(DPPB)₄ is a promising non-fullerene acceptor for future practical application of PSCs.

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