

## Electronic supplementary information

### Thermally Activated Delayed Fluorescence of Co-deposited Copper(I)

#### complexes: Cost-effective Emitters for Highly Efficient Organic

#### Light-Emitting Diodes

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#### General Information.

All chemicals and reagents were used as received from commercial sources without further purification unless stated otherwise, such as 1,3,5-tris(N-phenylbenzimidazole-2-yl) benzene (**TPBi**), 8-hydroxyquinoline lithium (**Liq**), 4,5,8,9,11-hexaaza triphenylene-hexacarbonitrile (**HAT-CN**), *N,N'*-bis(1-naphthyl)-*N,N'*-diphenyl-[1,1'-biphenyl]-4,4'-diamine (**NPB**), tris(4-carbazoyl-9-ylphenyl)amine (**TCTA**). Bis(1-(biphenyl)isoquinoline)iridium(III) acetylacetonate (**Ir(bpiq)<sub>2</sub>acac**) were purchased from Yurui (Shanghai) Chemical Co. Ltd. CzBPCb and CzBPDCb were synthesized as reported literature.<sup>36</sup>

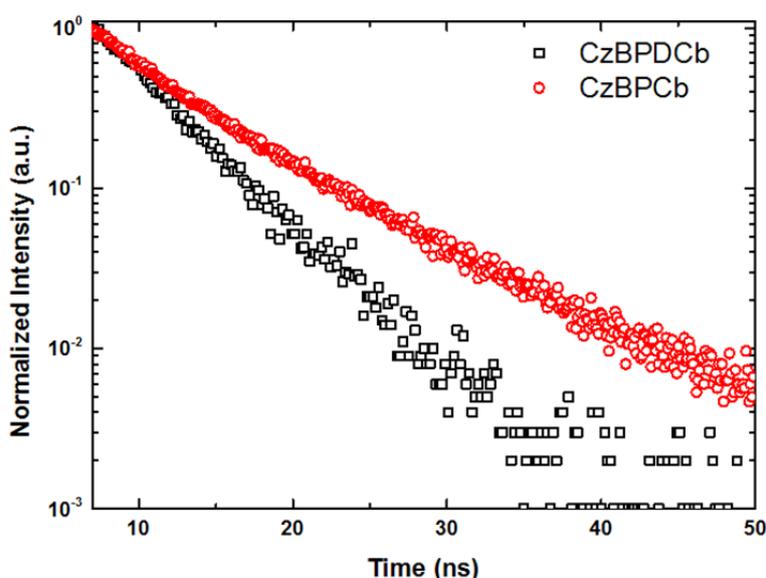


Figure S1. Transient PL decay curves of the neat films of CzBPDCb and CzBPCb.

$$\tau_p = 1/k_p, \quad \tau_d = 1/k_d \quad (1)$$

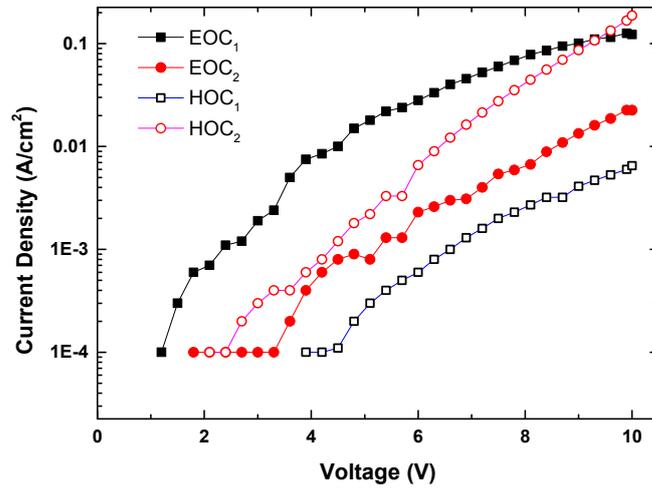
$$k_r^S = \Phi_{\text{prompt}} k_p \quad (2)$$

$$k_{\text{nr}}^T = k_d - \Phi_{\text{prompt}} k_{\text{RISC}} \quad (3)$$

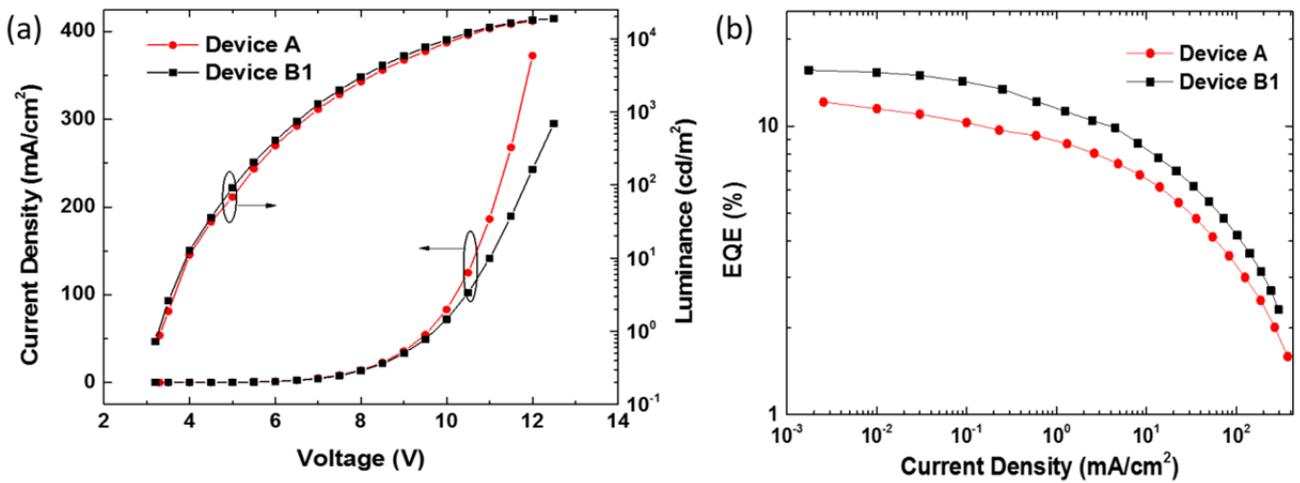
$$k_{\text{ISC}} = (1 - \Phi_{\text{prompt}}) k_p \quad (4)$$

$$k_{\text{RISC}} = \frac{k_p k_d \Phi_{\text{delayed}}}{k_{\text{ISC}} \Phi_{\text{prompt}}} \quad (5)$$

**Figure S2.** Equations for estimating the kinetic parameters



**Figure S3.** Current density-voltage (J-V) curves of EOCs and HOCs.



**Figure S4.** (a) Current density-Voltage-Luminance (J-V-L) characteristics, (b) EQE-J curves of device A and B1.

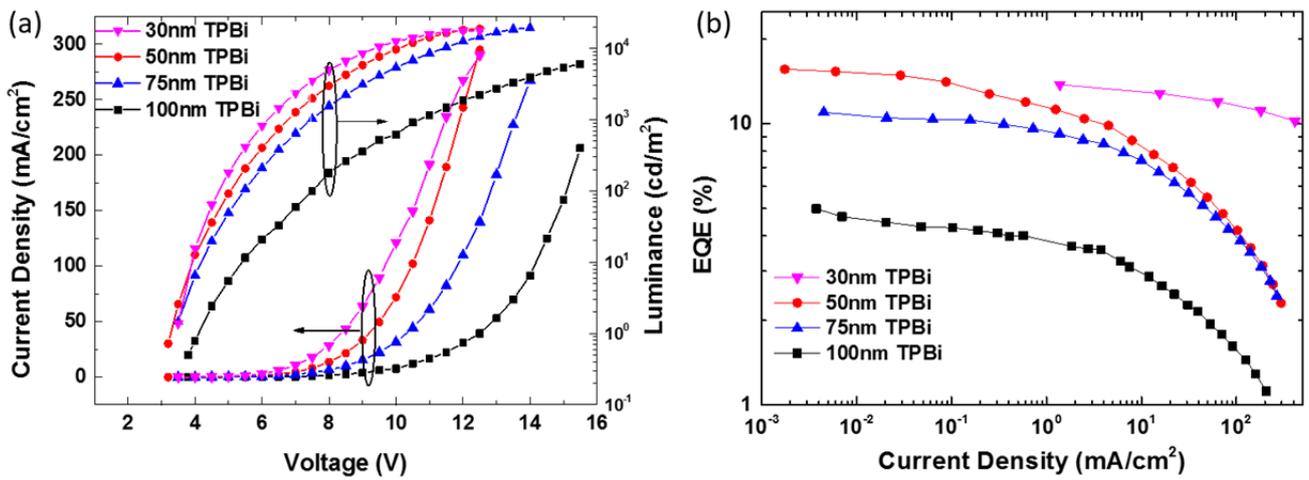


Figure S5. (a) J-V-L characteristics, (b) EQE-J curves of device B1-B4 with different thicknesses of TPBi.

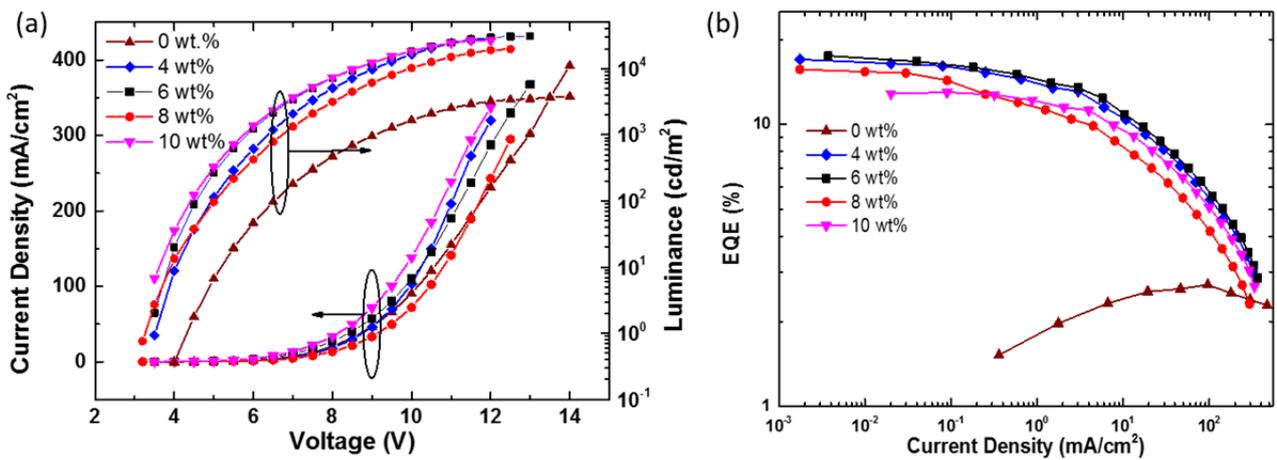


Figure S6. (a) J-V-L characteristics, (b) EQE-J curves of devices B1 and B5-B8 with different doping concentrations of Cul.

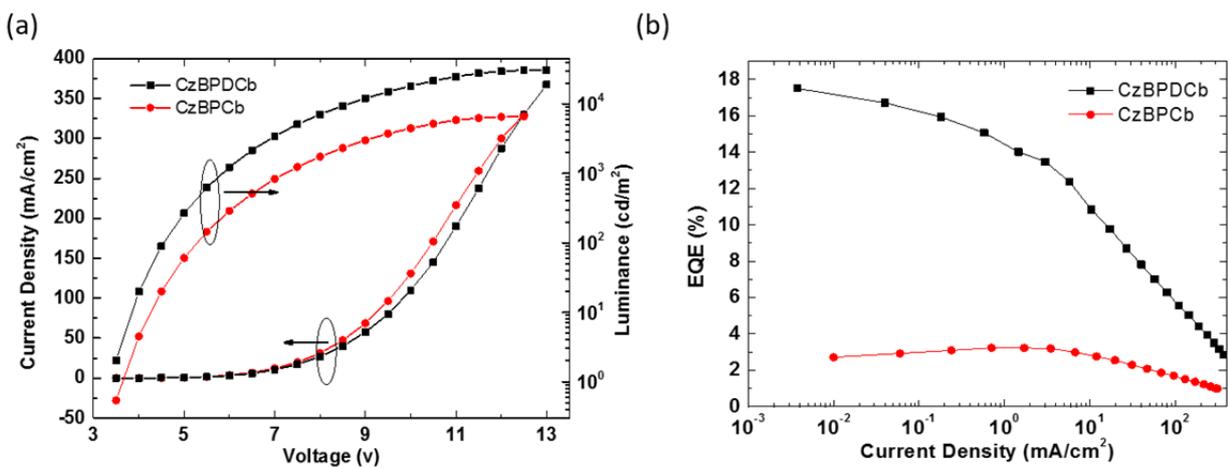


Figure S7. (a) J-V-L characteristics, (b) EQE-J curves of devices B7 and C using CzBPDCb and CzBPCb as co-deposited ligand with 6 wt% of Cul.