Supporting Information

Phosphonate Substituted 4,4'-Bis(*N*-carbazolyl)biphenyl with Dominant Electron Injection/Transport Ability for Tuning the Single-Layer Device Performance of Self-Host Phosphorescent Dendrimer

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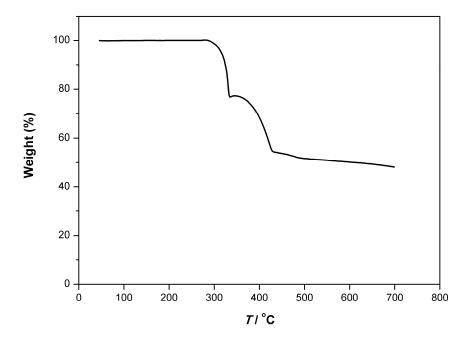


Fig. S1 TGA spectrum for PCBP, at a heating rate of 10 °C/min under nitrogen atmosphere.

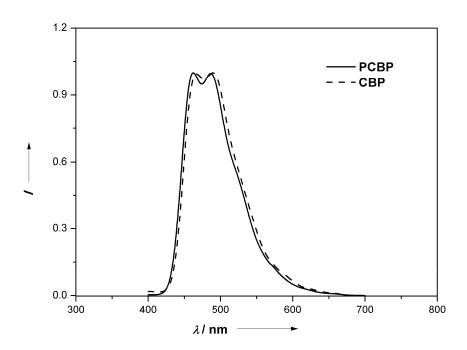


Fig. S2 Normalized phosphorenscence spectra of PCBP and CBP in 10⁻³ M toluene solution at 77K.

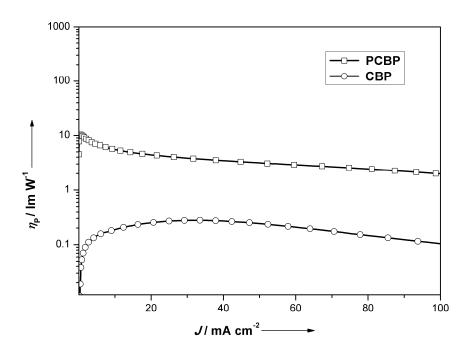


Fig. S3 The power efficiency-current density characteristics of the devices.

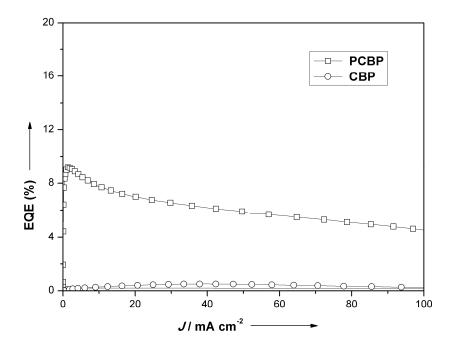


Fig. S4 The EQE-current density characteristics of the devices.

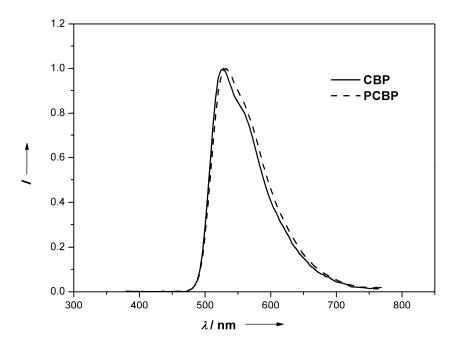


Fig. S5 Normalized EL spectra of the devices at a driving voltage of 15 V.

Table S1. Crystallographic Parameters and Refinement Results for PCBP.

Compound	PCBP
Empirical formula	C ₅₂ H ₆₀ N ₂ O ₁₂ P ₄ • 2CH ₂ Cl ₂
Formula weight	1198.75
Temperature	185(2) K
Wavelength	0.71073 Å
Crystal system	Monoclinic
Space group	<i>P</i> 2 ₁ /c
Unit cell dimensions	$a = 13.204(2) \text{ Å}, \qquad \alpha = 90^{\circ}$
	$b = 10.4398(15) \text{ Å}$ $\beta = 99.421(2)^{\circ}$
	$c = 21.992(3) \text{ Å} \qquad \gamma = 90^{\circ}$
Volume, Z	2990.6(8)ų, 2
Theta range for data collection	1.56 to 26.07°
Limiting indices	$-11 \le h \le 16, -12 \le k \le 12,$
	$-27 \le 1 \le 27$
Density (calculated)	1.331 Mg/m^3
Absorption coefficient	0.364 mm ⁻¹
F(000)	1252
Crystal size (mm³)	$0.42\times0.35\times0.10$
Reflections collected	18740
Independent reflections	5887 [R(int) = 0.0317]
Max. and min. transmission	0.9645 and 0.8622
Data / restraints / parameters	5887/0/383
Goodness-of-fit on F ²	1.008
Final R indices [I≥2σ (I)]	R1 = 0.0625, $wR2 = 0.1601$
R indices (all data)	R1 = 0.0851, $wR2 = 0.1772$
Largest diff. peak and hole	0.800 and -0.775 e/Å ⁻³