Supporting Information for;

Synthesis and Green Phosphorescent OLED Device Performance of Cyanofluorene-Linked Phenylcarbazoles as Host Material

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Figure S1. TGA thermograms of *o*-CF-PhCz, *m*-CF-PhCz and *p*-CF-PhCz under a heating rate of 10 °C/min..



Figure S2. Cyclic voltammograms of *o*-CF-PhCz, *m*-CF-PhCz and *p*-CF-PhCz.



Figure S3. Plot of hole (A) and electron (B) mobilities of single devices with *o*-CF-PhCz, *m*-CF-PhCz and *p*-CF-PhCz against the electric filed: HOD device [PEDOT:PSS(60 nm)/host(50 nm)/Al(200 nm)], EOD single device [PEDOT:PSS(60 nm)/TSPO1(10 nm)/host(50 nm)/LiF(1.5 nm)/Al(200 nm)].



Figure S4. Current density-voltage curves of the green PhOLED devices with *o*-CF-PhCz, *m*-CF-PhCz and *p*-CF-PhCz according to the doping concentration of dopant.



Figure S5. The optimized chemical structure of *m*-CF-PhCz and *p*-CF-PhCz.



Figure S6. External quantum efficiency-luminance curves of *o*-CF-PhCz at 3, 5, and 10 % doping concentration of dopant.



Figure S7. External quantum efficiency-luminance curves of *m*-CF-PhCz at 3, 5, and 10 % doping concentration of dopant.



Figure S8. External quantum efficiency-luminance curves of *p*-CF-PhCz at 3, 5, and 10 % doping concentration of dopant.

	o-CF-PhCz	<i>m</i> -CF-PhCz	<i>p</i> -CF-PhCz
HOMO (eV)	-5.468	-5.458	-5.417
LUMO (eV)	-1.781	-1.891	-1.889
E_g (eV)	3.687	3.567	3.528
S ₁ (eV)	-3.193	-3.192	-3.182
T ₁ (eV)	-2.701	-2.656	-2.613

 Table S1. Molecular simulation results of o-CF-PhCz, m-CF-PhCz and p-CF-PhCz.