Novel A-(π-D-π-A)$_{1-3}$ branched fluorophores displaying high two-photon absorption

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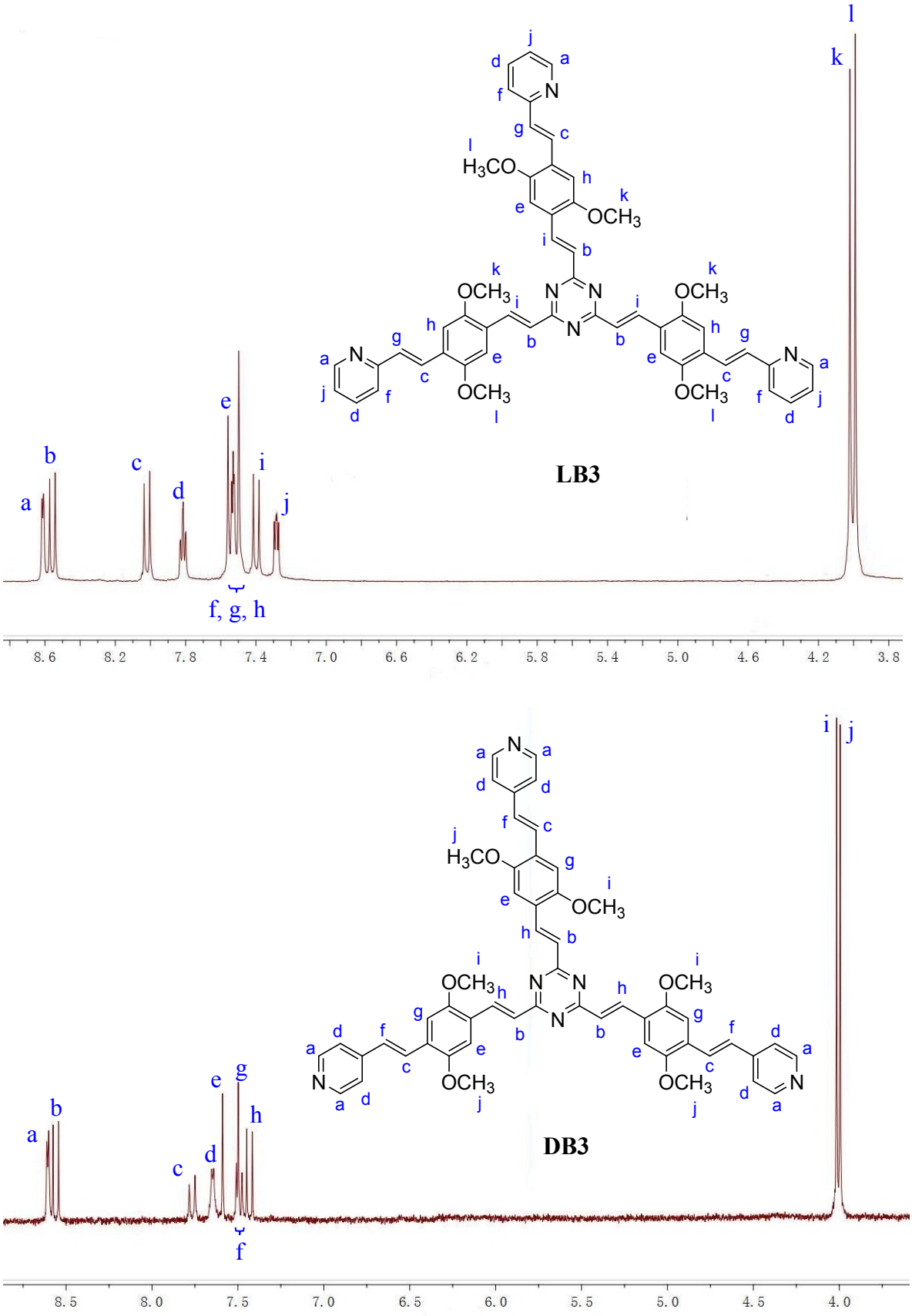


Figure S1. ^{1}H NMR spectra of **LB3** and **DB3** in DMSO-d_6.

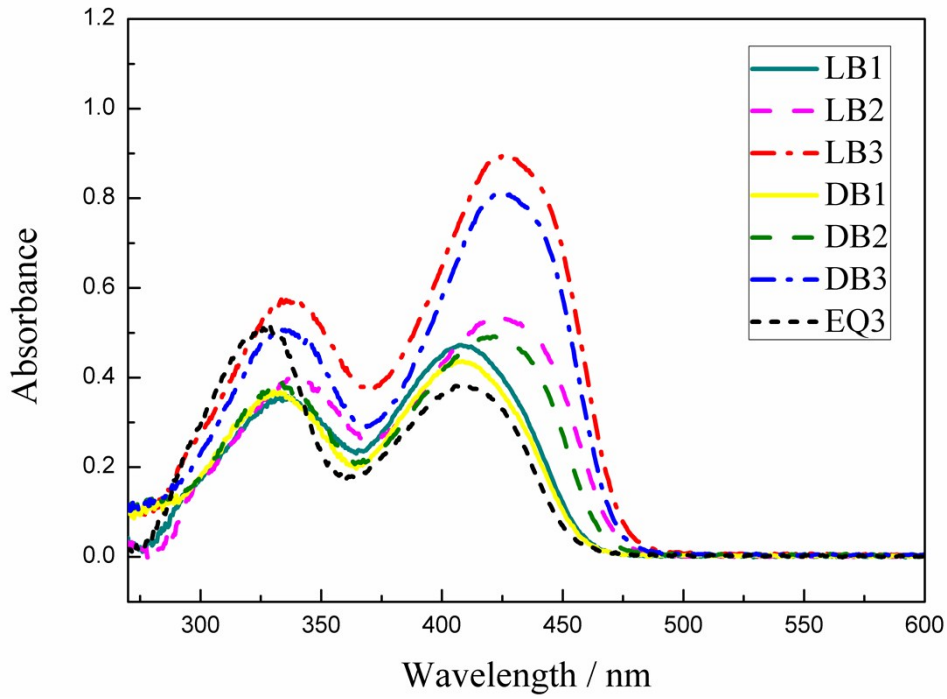
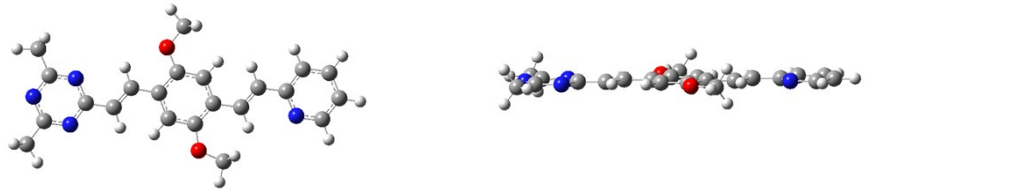


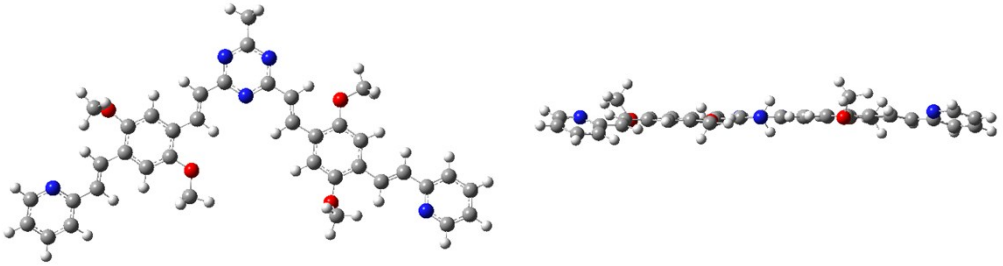
Figure S2. UV-visible absorption spectra of the target compounds (**LB1**-**LB3**, **DB1**-**DB3**, **EQ3**) in THF (c = 1×10^{-5} mol L^{-1})

top view front view

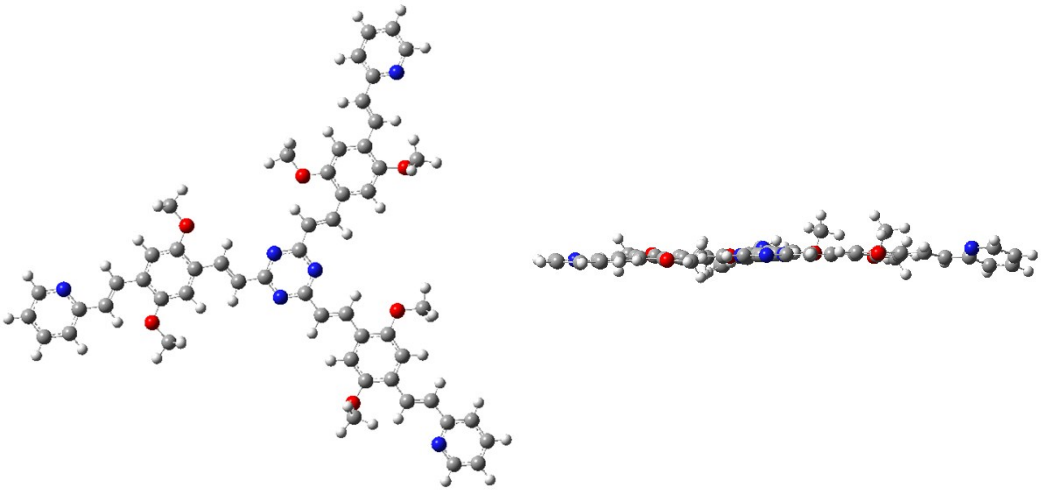
LB1



LB2

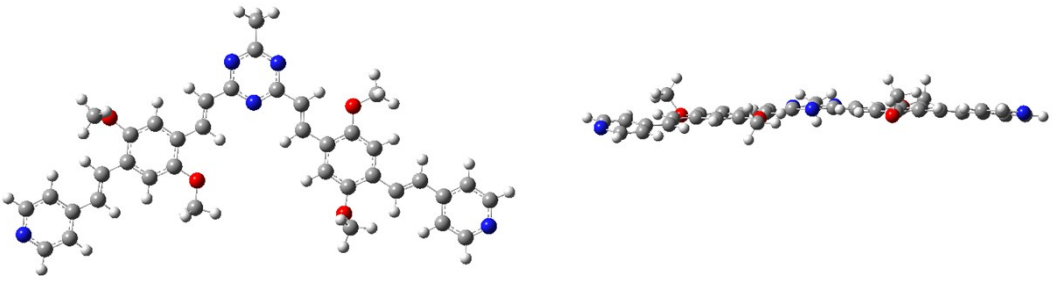


LB3

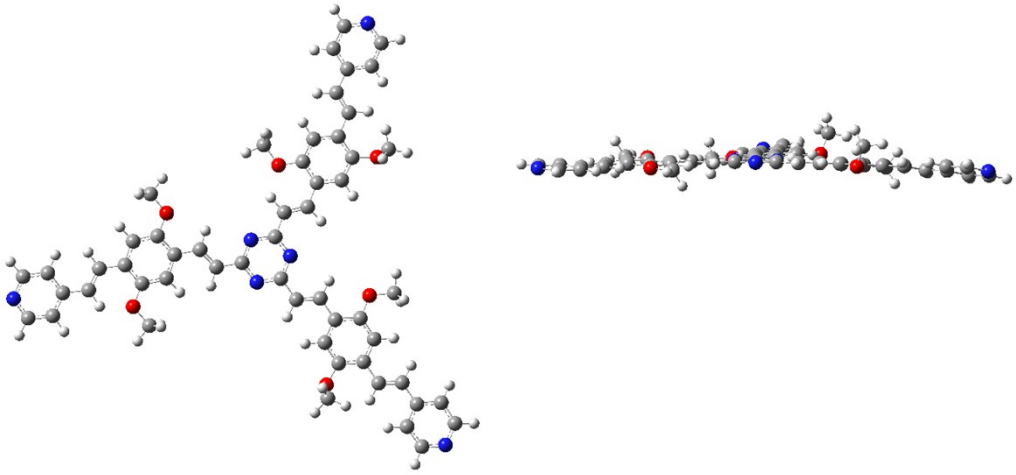


DB1

DB2



DB3



EQ3

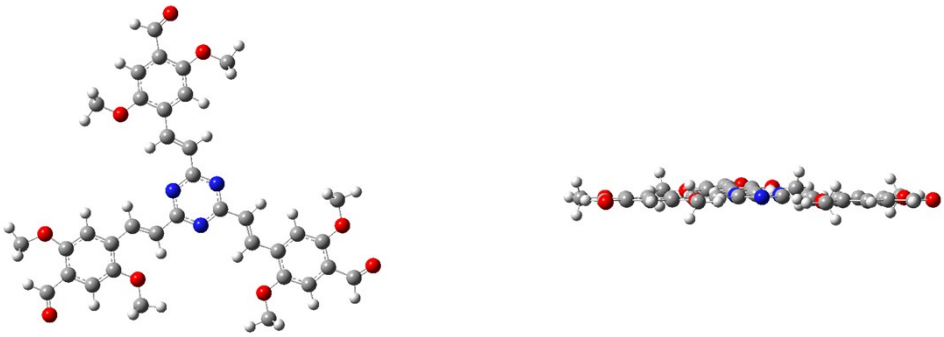
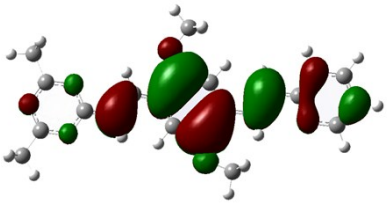
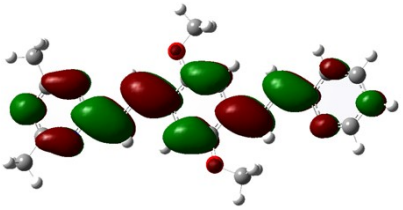
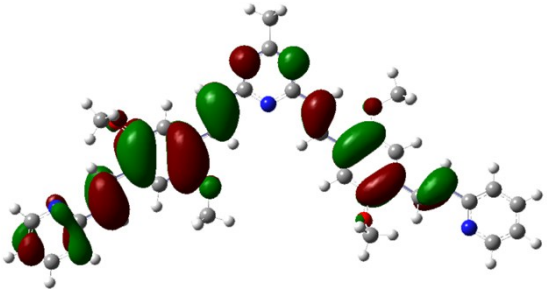
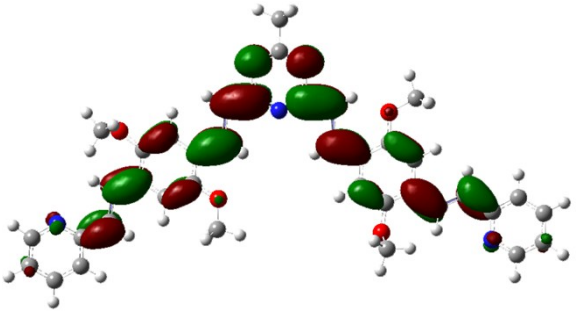
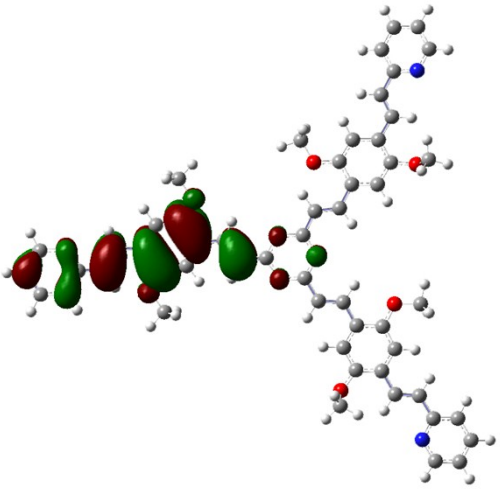
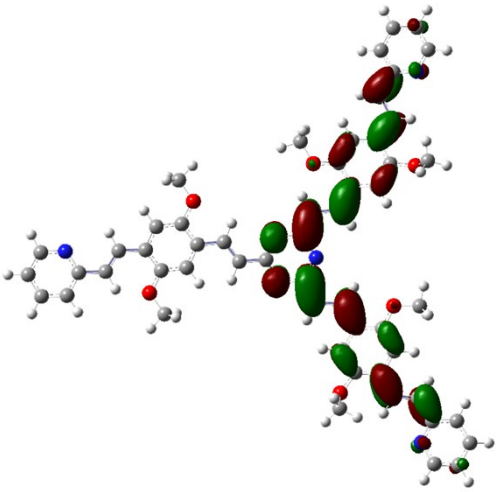
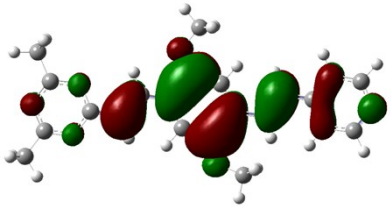
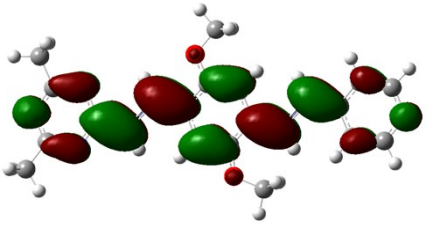


Figure S3. Optimized geometries of the target compounds (**LB1-LB3**, **DB1-DB3**, **EQ3**)

	HOMO	LUMO
LB1		
LB2		
LB3		
DB1		

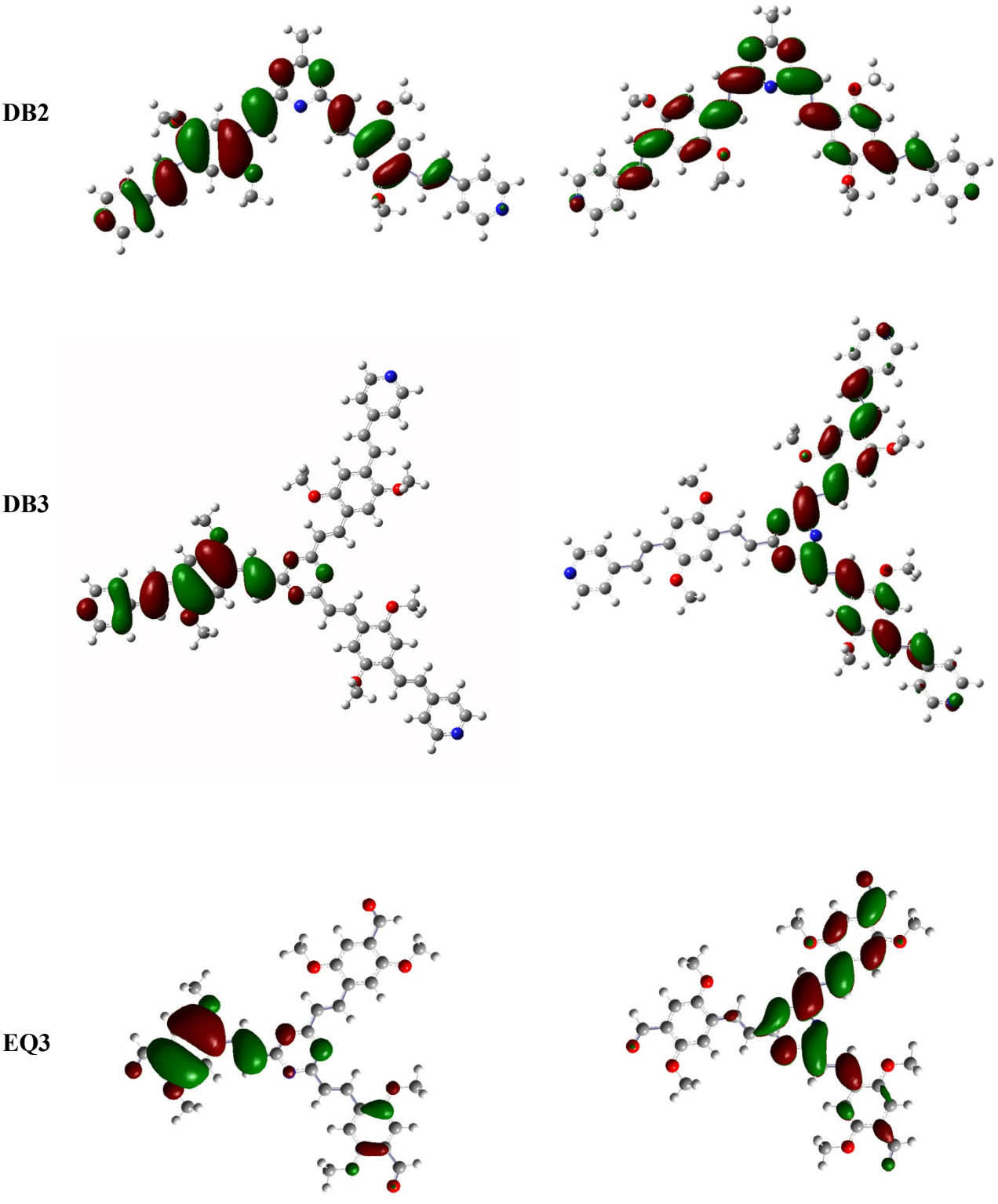


Figure S4. Frontier molecular orbitals of the target compounds (**LB1-LB3**, **DB1-DB3**, **EQ3**)