

## Supporting Information

for

### **“Scorpion” shape mono s(carboxy)Porphyrin-(Bodipy)<sub>2</sub>: a novel triazine bridged triad, synthesis, characterization and its dye sensitized solar cell (DSSC) applications**

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**Table S1.** Coordinates of gas phase geometry optimized structure of **3 PorCOOH-(BDP)<sub>2</sub>** calculated by DFT at the B3LYP / 6-31G(d) level.  
E = - 4921.639274 Hartree / particle.

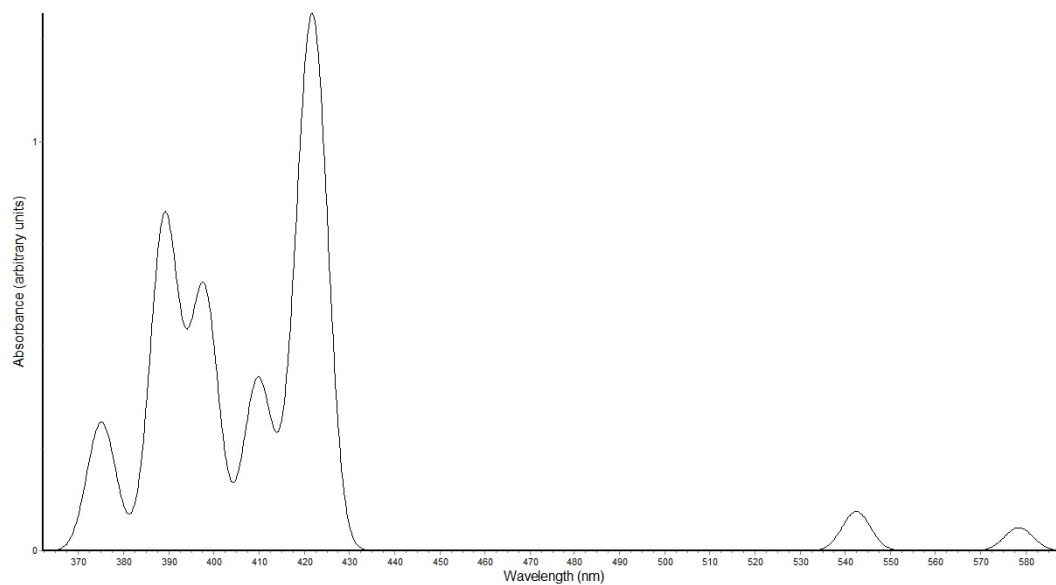
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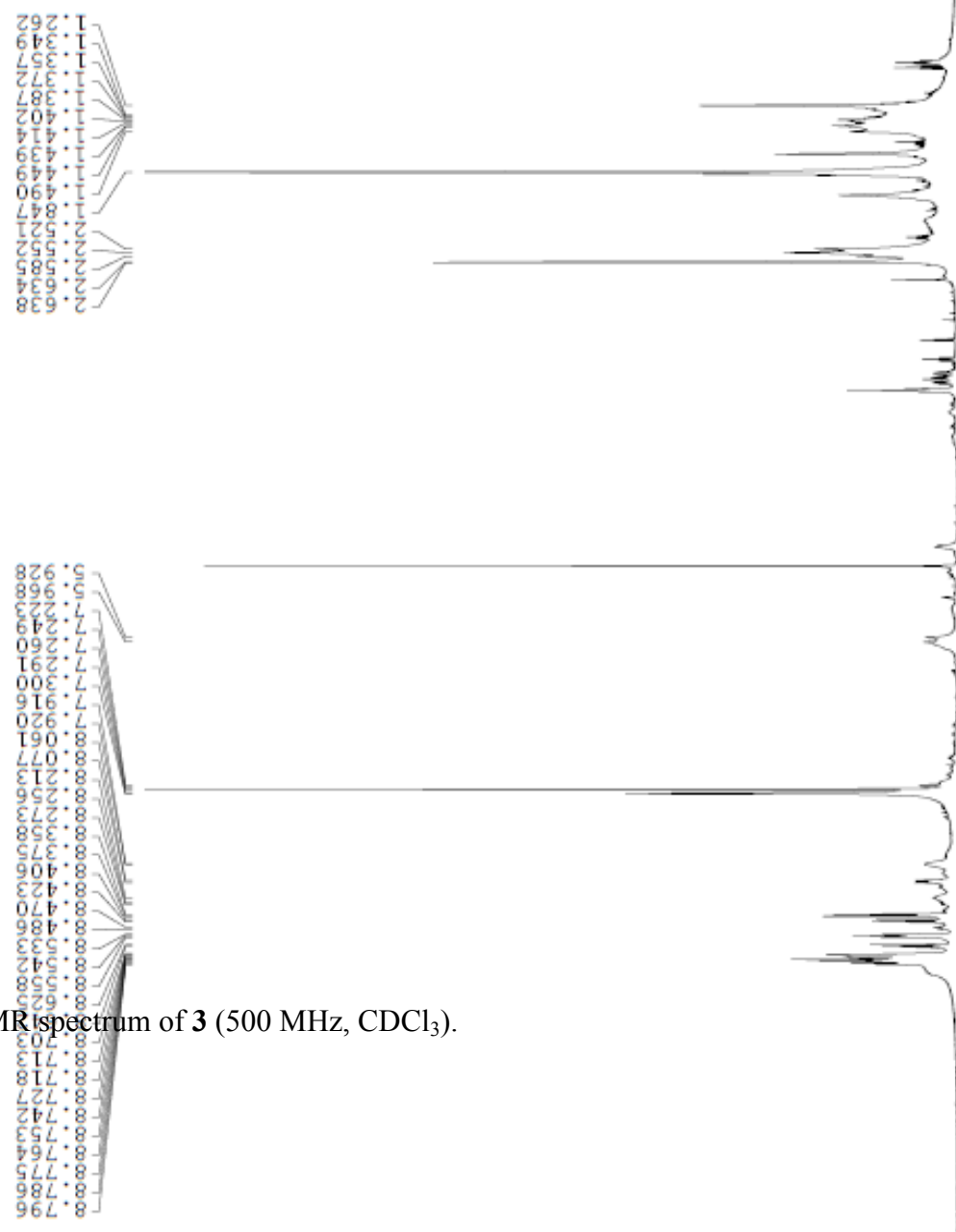
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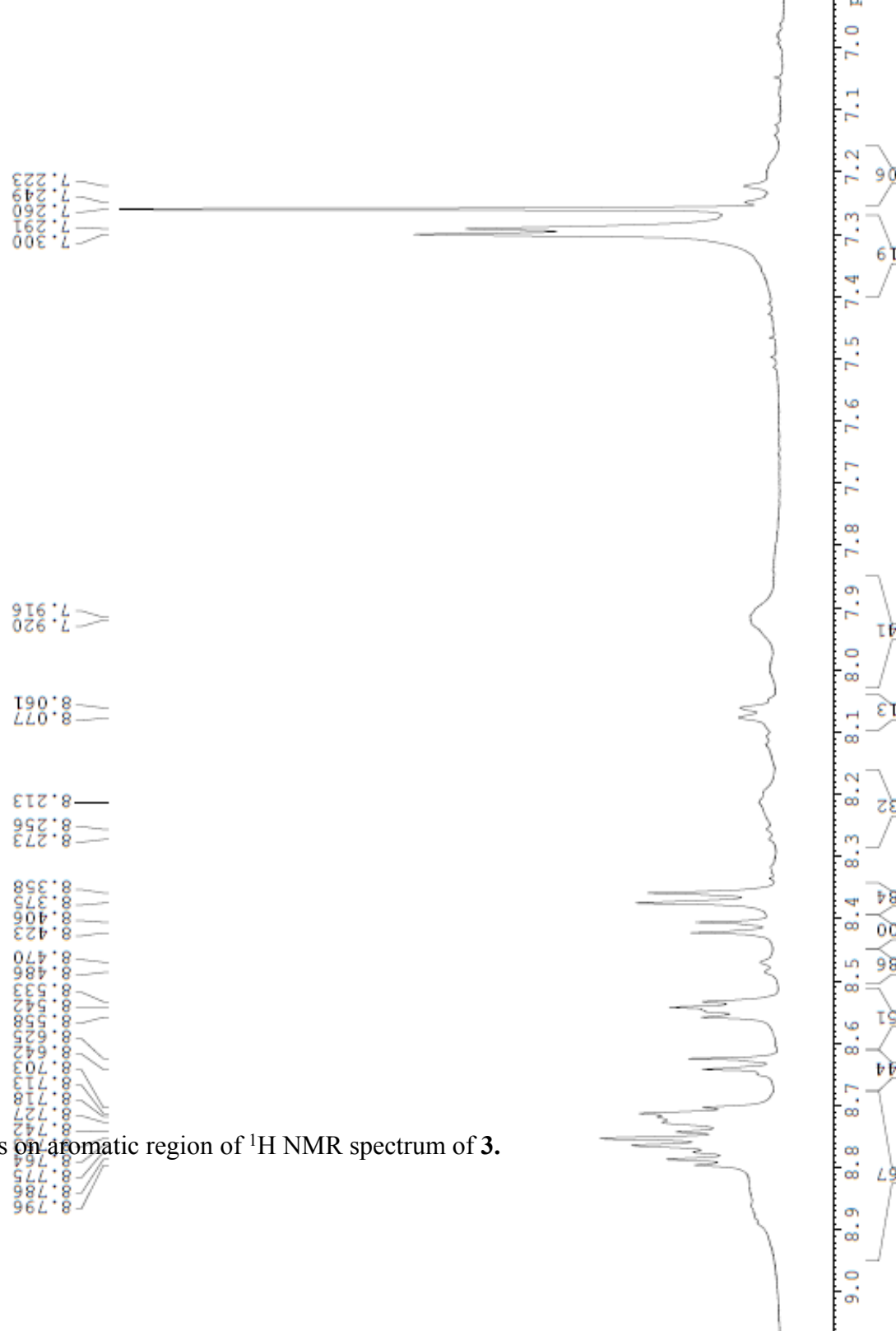
**Figure S1.** Theoretically calculated UV-vis absorption spectra of **3 PorCOOH-(BDP)<sub>2</sub>**.

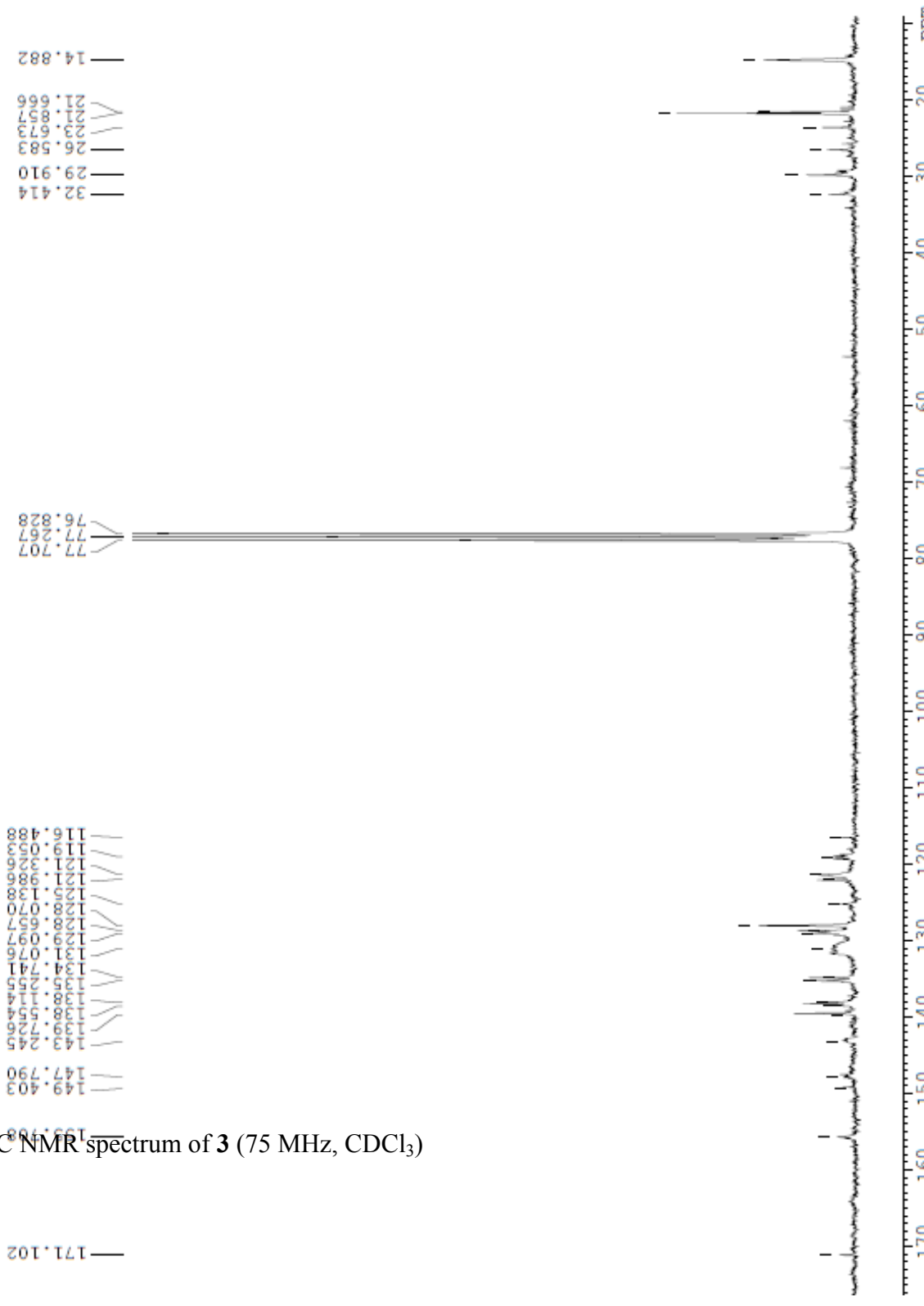




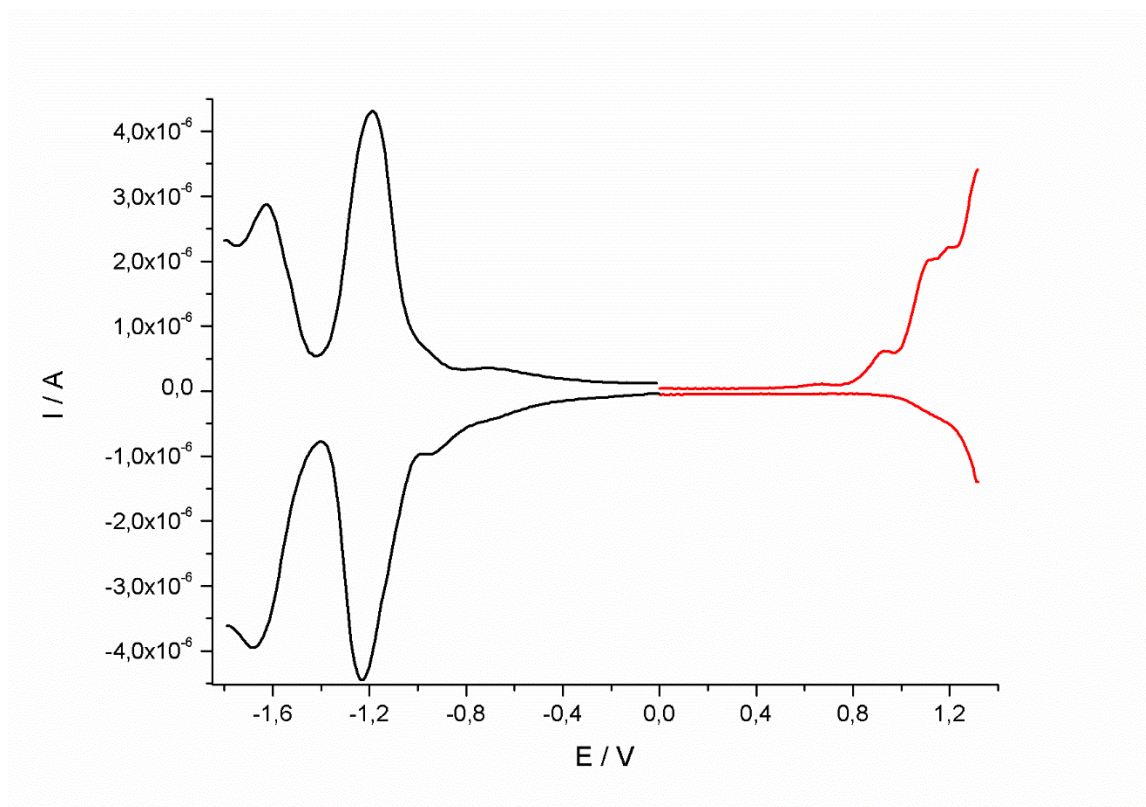
**Figure S2.**  $^1\text{H}$  NMR spectrum of **3** (500 MHz,  $\text{CDCl}_3$ ).

**Figure S3.** Focus on aromatic region of  $^1\text{H}$  NMR spectrum of **3**.

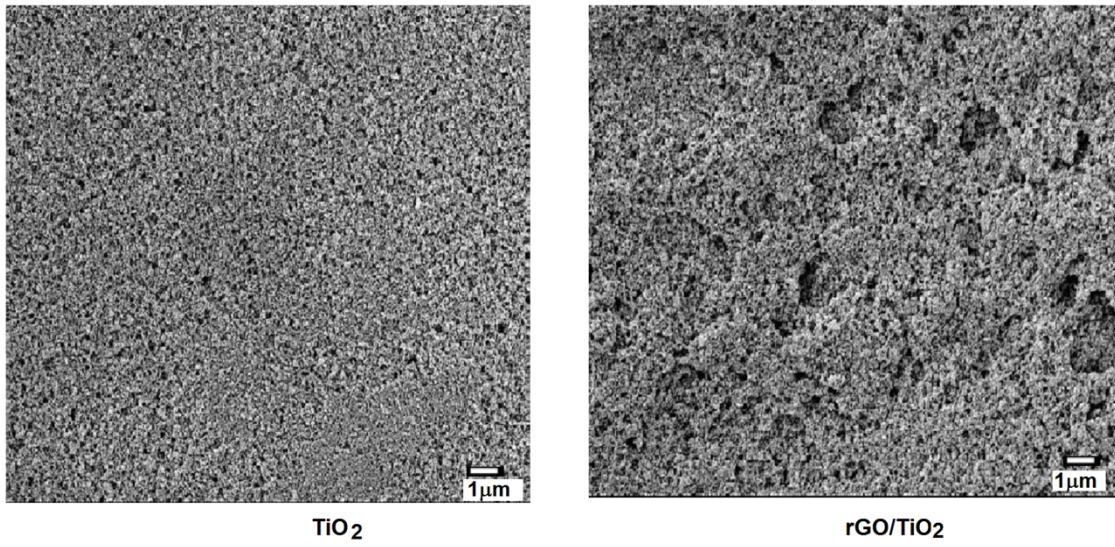




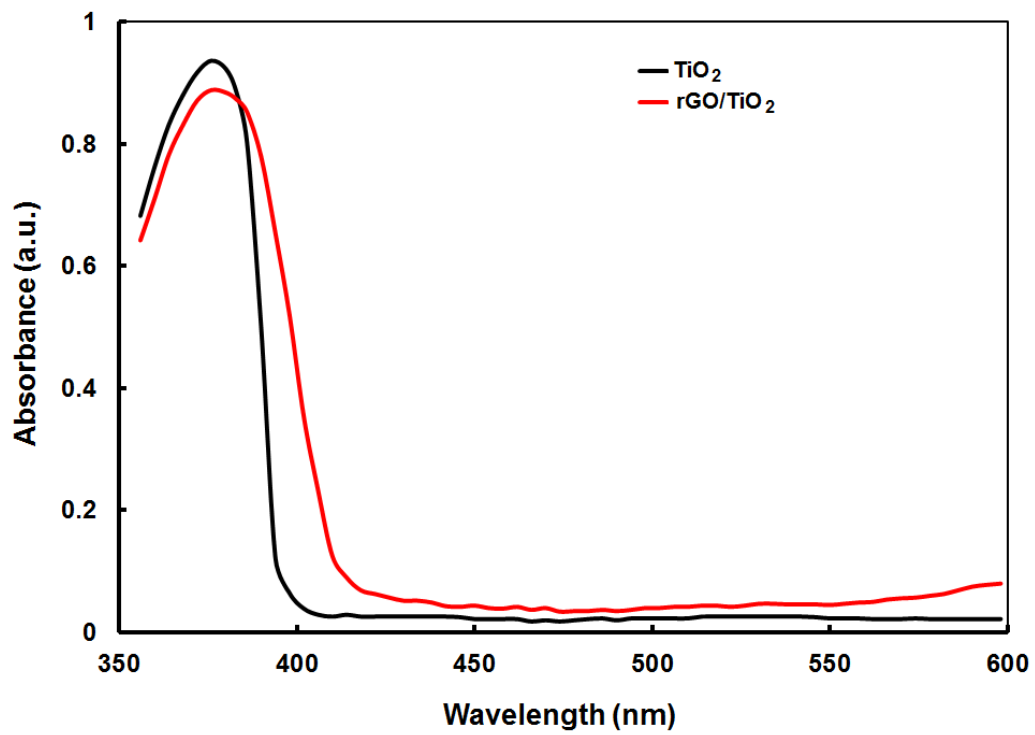
**Figure S4.**  $^{13}\text{C}$  NMR spectrum of **3** (75 MHz,  $\text{CDCl}_3$ )



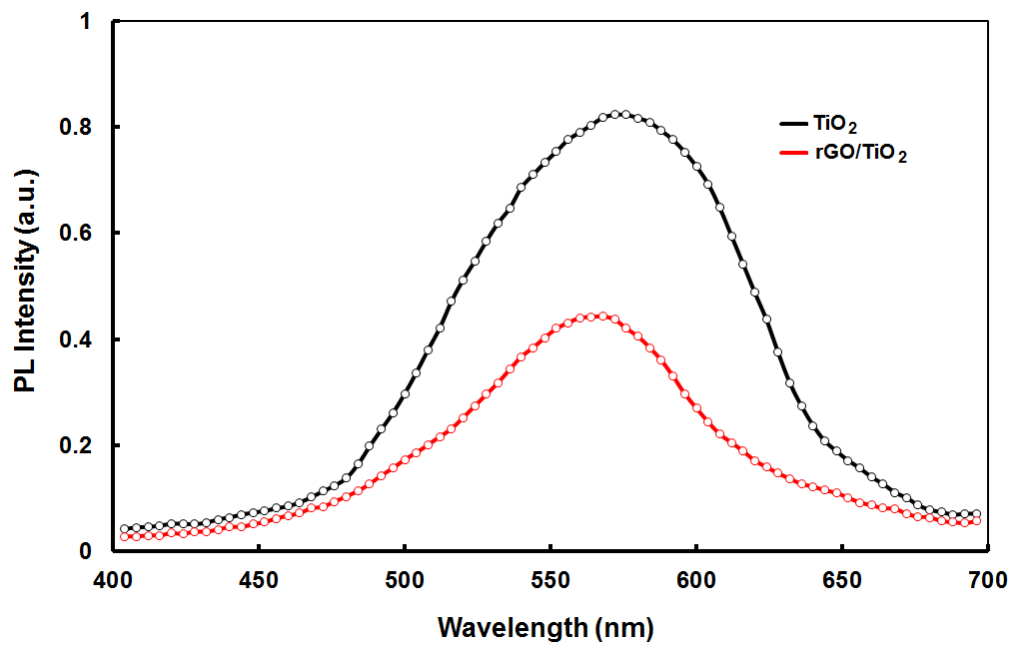
**Figure S5.** Square-wave voltammogram of **3**. The ferrocene/ferrocenium (Fc/Fc<sup>+</sup>) redox couple wave (not shown) is found to be at 0.56 V. Voltage is reported vs SCE.



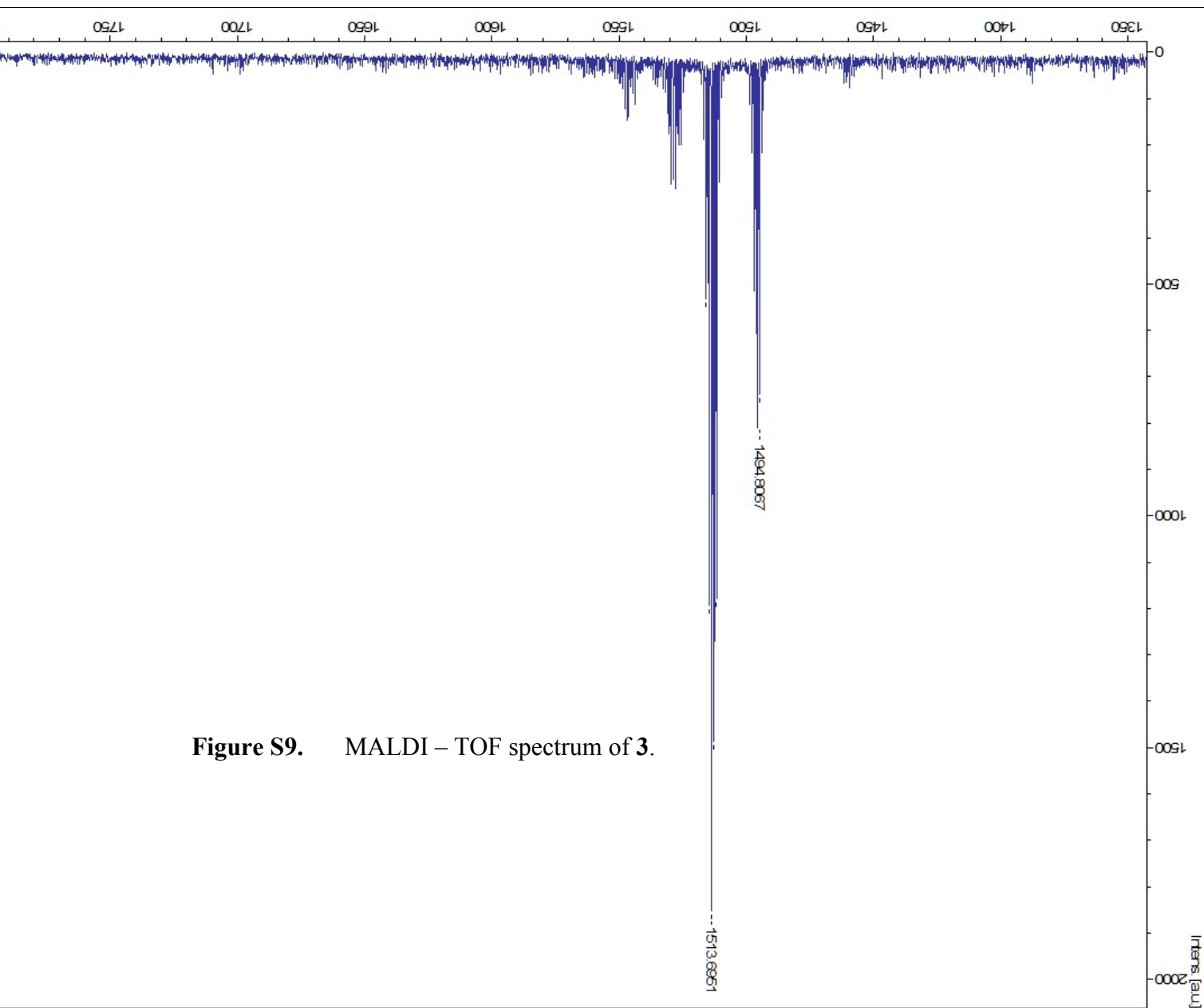
**Figure S6.** SEM images of  $\text{TiO}_2$  (left) and  $\text{rGO/TiO}_2$  (right) films.



**Figure S7.** UV-visible absorption spectra of TiO<sub>2</sub> (black line) and rGO/TiO<sub>2</sub> (red line) films



**Figure S8.** PL spectra of TiO<sub>2</sub> (black line) and rGO/TiO<sub>2</sub> (red line) films.



**Figure S9.** MALDI – TOF spectrum of **3**.