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SUPPLEMENTARY INFORMATION

Extensive methodology screening of *meso*-tetrakys-(furan-2-yl)-porphyrin microwave-assisted synthesis

F. Bosca,^a S. Tagliapietra,^a C. Garino,^b G. Cravotto *^a and A. Barge*^a

- Figure S1. ¹H NMR spectrum of compound **1** (300 MHz, THF-d ₈)
- Figure S2. ¹H NMR spectrum of compound **2** (300 MHz, CDCl₃)
- Figure S3. ¹H NMR spectrum of compound **3** (300 MHz, CDCl₃)
- Figure S4. ¹H NMR spectrum of compound **5** (300 MHz, THF-d ₈)
- Figure S5. Electronic absorption and photoemission spectra of compounds 1 and 2
- Figure S6. Extinction Coefficients of compounds 1 and 2.
- Figure S7. HPLC-MS analysis of compound **4**.
- Figure S8. APCI-MS (positive molality) analysis of compound 4
- Figure S9. HPLC-MS of compound **1**.
- Figure S10. IR spectrum of compound 2.
- Figure S9. Epsilon of compounds 1 and 2.







Figure S4. ¹H NMR spectrum of compound **5** (300 MHz, THF-d $_8$)



Figure S5. Electronic absorption (left) and photoemission (right) spectra of *meso*-tetrakys-(furan-2-yl)-porphyrin (red) and Zn(II)-*meso*-tetrakys-(furan-2-yl)-porphyrin (blue) in THF solution.



Figure S6. Extinction Coefficient (left) and photoemission (right) spectra of *meso*-tetrakys-(furan-2-yl)-porphyrin (red) and Zn(II)-*meso*-tetrakys-(furan-2-yl)-porphyrin (blue) in THF solution.



Figure S7. HPLC-MS analysis of compound 4.



Figure S8. APCI-MS (positive molality) analysis of compound 4



Figure S9. HPLC-MS of compound **1**. Upper trace: ESI+ profile of m/z= 575.5 (peak of interest); middle trace: UV profile at 459.5 nm; bottom trace: Diode array max plot. Rt1 and Rt2 indicate the two regions considered in the impurity analysis.



Figure S10. IR spectrum of compound 2.