Supporting information for:

## Charge Recombination Suppressing in Dye-Sensitized Solar Cells by Tuning Dielectric Constant of Triphenylamine Dyes with Altering π-Bridge from Naphthalene to Anthracene Units

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Corresponding Author: Mohammad-Reza Zamani Meymian<sup>\*</sup>, Email: r\_zamani@iust.ac.ir, Telephone Number: (+98) 21 7322 5893 . Fax: (+98) 21 7724 0497 1- FT-IR spectrum of TpAzo 1: (KBr, 298 K): aromatic C-H stretch v [3026 cm<sup>-1</sup>], in-plane bends aromatic C-H v [1150 cm<sup>-1</sup>], out-of-plane bends aromatic C-H v [725 cm<sup>-1</sup>], stretch aromatic C=C v [1550 cm<sup>-1</sup>], stretch C-N v [1277 cm<sup>-1</sup>], stretch N=N v [2300 cm<sup>-1</sup>], H-bonded stretch monomer O-H v [3450 cm<sup>-1</sup>], stretch dimer H-bonded O-H v [2950 cm<sup>-1</sup>], bend out-of-plane O-H v [937 cm<sup>-1</sup>], stretch monomer H-bonded C=O v [1600 cm<sup>-1</sup>], in-plane bend C-O-H v [1450 cm<sup>-1</sup>] and stretch C-O v [1240 cm<sup>-1</sup>].



Fig. 1S. FT-IR spectrum of TpAzo 1

2- FT-IR spectrum of TpAzo 2: (KBr, 298 K): stretch aromatic C-H v [3026 cm<sup>-1</sup>], in-plane bends aromatic C-H v [1150 cm<sup>-1</sup>], out-of-plane bends aromatic C-H v [725 cm<sup>-1</sup>], aromatic stretch C=C v [1550 cm<sup>-1</sup>], stretch C-N v [1277 cm<sup>-1</sup>], stretch N=N v [2300 cm<sup>-1</sup>], stretch monomer H-bonded O-H v [3450 cm<sup>-1</sup>], stretch dimer H-bonded O-H v [2950 cm<sup>-1</sup>], bend out-of-plane O-H v [937 cm<sup>-1</sup>], stretch monomer H-bonded C=O v [1600 cm<sup>-1</sup>], in-plane bend C-O-H v [1450 cm<sup>-1</sup>] and stretch C-O v [1240 cm<sup>-1</sup>].



Fig. 2S. FT-IR spectrum of TpAzo 2

<sup>1</sup>H&<sup>13</sup>C-NMR and Mass spectra **TpAzo 1**: <sup>1</sup>H NMR (CDCl<sub>3</sub>): δ [*ppm*]: 8.31 (dddt, *J* = 1.8, 1.7, 0.5, 0.4 Hz)), 8.29-8.34 (2H, 8.33 (tq, *J* = 1.7, 0.4 Hz), 8.07 (1H, ddt, *J* = 7.1, 1.7, 0.4 Hz), 8.01 (ddd, *J* = 7.1, 1.7, 0.4 Hz)), 7.93-8.03 (2H, 7.96 (ddt, *J* = 8.8, 1.8, 0.4 Hz), 7.73 (2H, ddd, *J* = 8.6, 2.7, 0.4 Hz), 7.61 (dddd, *J* = 8.3, 7.2, 1.7, 0.5 Hz)), 7.56-7.66 (5H, 7.62 (ddd, *J* = 8.8, 1.8, 0.4 Hz), 7.51 (2H, ddd, *J* = 8.6, 1.5, 0.4 Hz), 7.23-7.35 (6H, 7.26 (dddd, *J* = 8.3, 1.3, 1.2, 0.5 Hz), 7.32 (tt, *J* = 7.2, 1.2 Hz)). <sup>13</sup>C NMR (CDCl<sub>3</sub>): δ [*ppm*]: 167.1, 149.5, 147.9, 147.9, 147.3, 132.7, 131.8, 128.9, 128.6, 128.2, 127.8, 127.3, 127.1, 125.0, 124.5, 124.2, 119.2, 115.0, 114.4. Mass [*m*/*z*]. C<sub>29</sub>H<sub>21</sub>N<sub>3</sub>O<sub>2</sub> [M<sup>+</sup>], Found. 451; Calcd. 452.





Fig. 3S. <sup>1</sup>H&<sup>13</sup>C-NMR and Mass spectra TpAzo 1:

<sup>1</sup>H&<sup>13</sup>C-NMR and Mass spectra **TpAzo 2**: <sup>1</sup>H NMR (CDCl<sub>3</sub>): δ [*ppm*]: 8.59 (1H, dd, J = 1.7, 0.6, 0.4 Hz), 8.23-8.30 (2H, 8.25 (ddt, J = 5.9, 1.7, 0.5 Hz), 8.28 (dddd, J = 1.7, 1.5, 0.5, 0.4 Hz)), 7.96 (1H, ddd, J = 5.9, 1.7, 0.4 Hz), 7.83-7.92 (2H, 7.88 (dddd, J = 8.2, 2.2, 1.7, 0.4 Hz), 7.85 (ddq, J = 1.7, 1.5, 0.5 Hz)), 7.78 (ddd, J = 10.1, 2.2, 0.4 Hz)), 7.61 (dddd, J = 8.3, 7.2, 1.6, 0.5 Hz), 7.71 (ddd, J = 10.1, 8.2, 0.5 Hz), 7.56-7.82 (8H, 7.79 (ddd, J = 7.6, 2.6, 0.4 Hz), 7.30 (2H, ddd, J = 7.6, 1.7, 0.4 Hz), 7.30-7.42 (6H, 7.38 (dtd, J = 8.3, 1.2, 0.5 Hz), 7.33 (tt, J = 7.2, 1.2 Hz)). <sup>13</sup>C NMR (CDCl<sub>3</sub>), δ [*ppm*]: 169.3, 126.8, 126.8, 129.8, 129.6, 129.6, 125.9, 129.6, 129.6, 124.7, 130.3, 128.3, 125.7, 125.7, 125.9, 125.1, 130.8, 125.8, 125.5, 125.7, 125.7, 125.9, 125.1, 120.2, 127.7, 133.2, 130.4,



132.9, 123.0, 124.5, 145.9, 148.1, 149.7, 152.1. Mass [m/z]. C<sub>31</sub>H<sub>23</sub>N<sub>3</sub>O<sub>2</sub> (M<sup>+</sup>), Found: 475; Calcd: 476.

Fig. 4S. <sup>1</sup>H-NMR, <sup>13</sup>C NMR and Mass spectroscopy analysis of TpAzo 2 dye