

Supporting Information

A General Approach for Synthesis of Functional Metal Oxide Nanotubes and their Application in Dye-sensitized Solar Cells

Parthiban Ramasamy, Da-Hye Lim, Jungsu Kim and Jinkwon Kim*

*Department of Chemistry and GETRC, Kongju National University, 182, Shinkwondong, Kongju, 314-701,
Chungnam-do, Republic of Korea.*

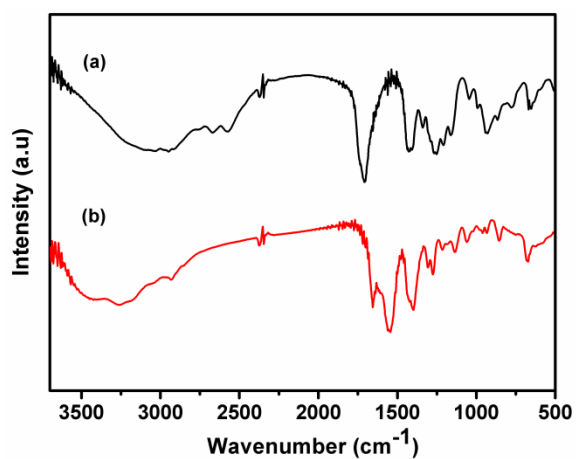


Fig. S1 FT- IR spectra of (a) MPA and (b) Se NWs functionalized with Mercaptopropionic acid.

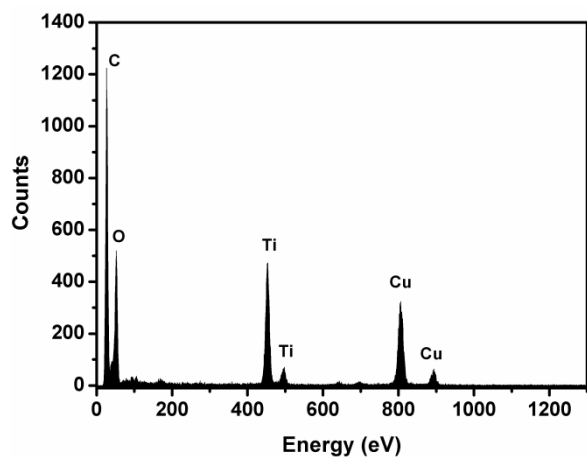


Fig. S2 EDS spectrum of TiO₂ nanotubes.

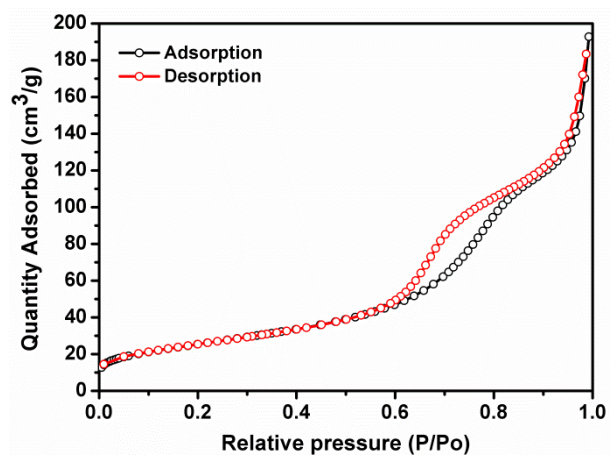


Fig. S3 Nitrogen adsorption-desorption isotherms of TiO₂ nanotubes.

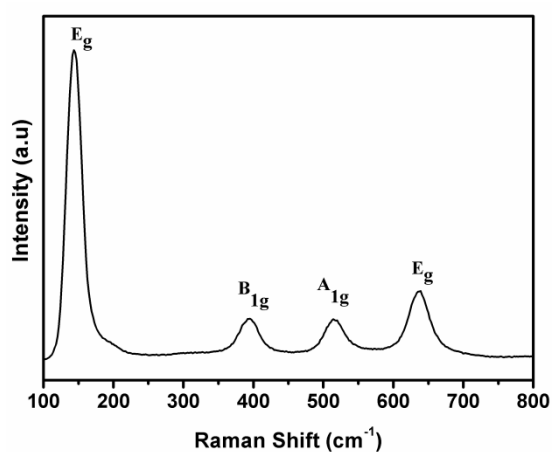


Fig. S4 Raman spectrum of TiO₂ nanotubes. The major Raman peaks observed at 144, 394, 516 and 637 cm⁻¹ can be attributed to E_g, B_{1g}, A_{1g} and E_g modes of anatase TiO₂. It is well known that the E_g peak is mainly due to symmetric stretching vibration of O-Ti-O, the B_{1g} peak is caused by symmetric bending vibration of O-Ti-O, and the A_{1g} peak is corresponds to anti-symmetric bending vibration of O-Ti-O in TiO₂.

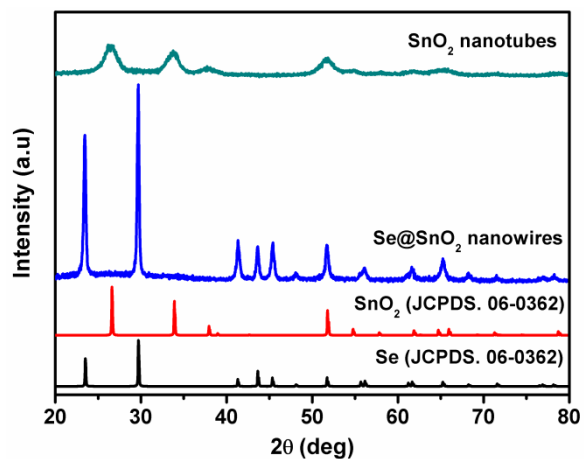


Fig. S5. XRD patterns of Se@SnO₂ core-shell nanowires and SnO₂ nanotubes.

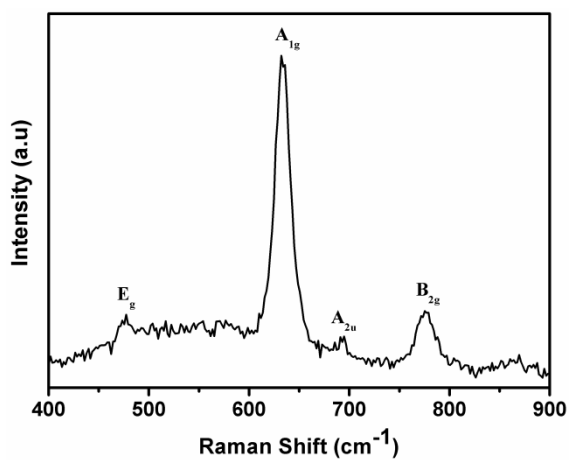


Fig. S6 Room-temperature Raman spectrum of SnO₂ nanotubes. The Raman scattering peaks at 475, 632, and 775 cm⁻¹ correspond to the E_g, A_{1g}, and B_{2g} vibrational modes, respectively. Thus, these Raman peaks further confirmed the tetragonal rutile structure of the as-synthesized SnO₂ nanotubes. The Raman peak at 694 cm⁻¹ can be assigned to the IR-active LO A_{2u} mode of SnO₂.

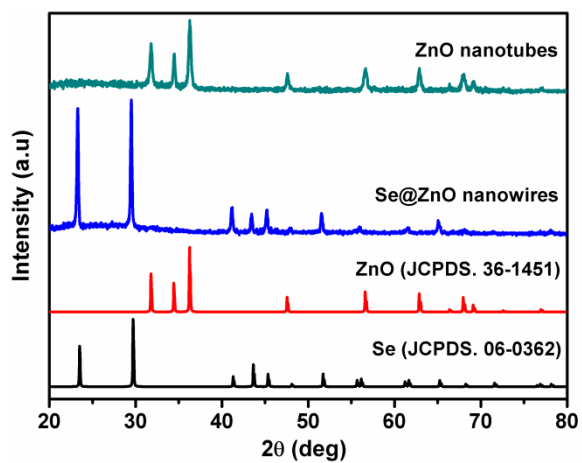


Fig. S7 XRD patterns of Se@ZnO core-shell nanowires and ZnO nanotubes.

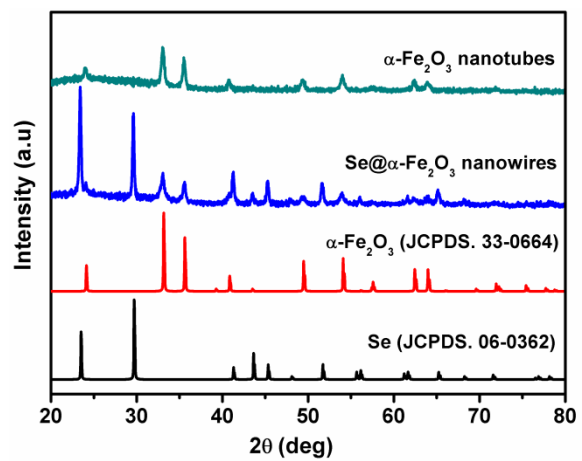


Fig. S8 XRD patterns of Se@Fe₂O₃ core-shell nanowires and Fe₂O₃ nanotubes.

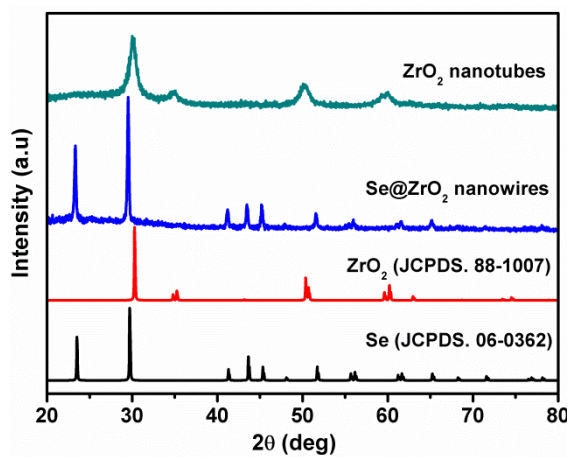


Fig. S9 XRD patterns of Se@ZrO₂ core-shell nanowires and ZrO₂ nanotubes.