Supporting Information

Self-assembly of Mesoporous $ZnCo_2O_4$ Nanomaterials: Density Functional Theory Calculation and Flexible All-Solid-State Energy Storage

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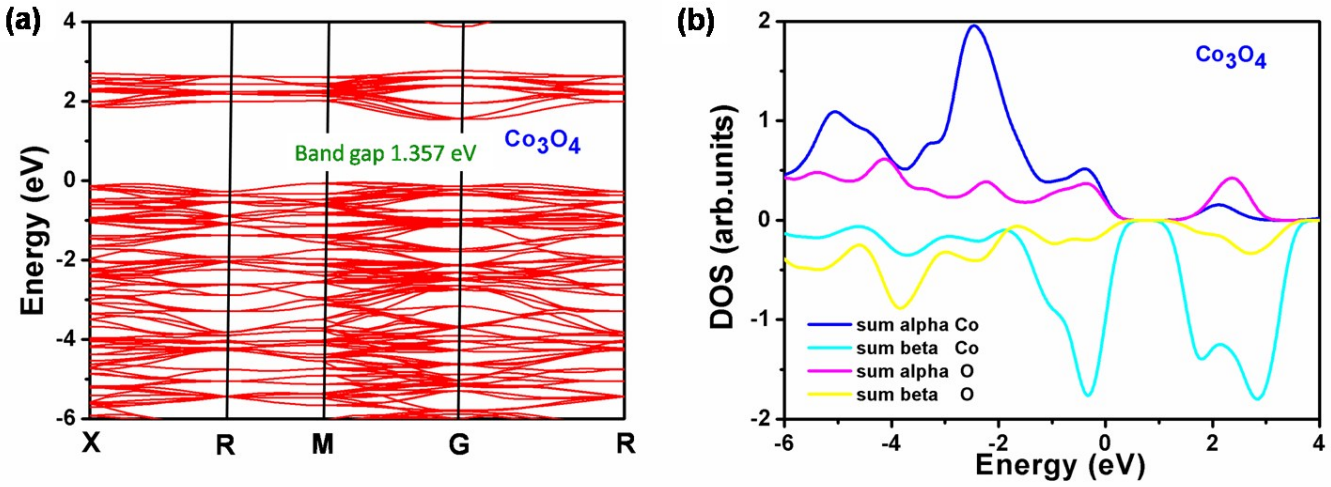

Figure S1. Calculated band structure (a) and total densities of states (b) of spinel Co_3O_4 by the CASTEP.

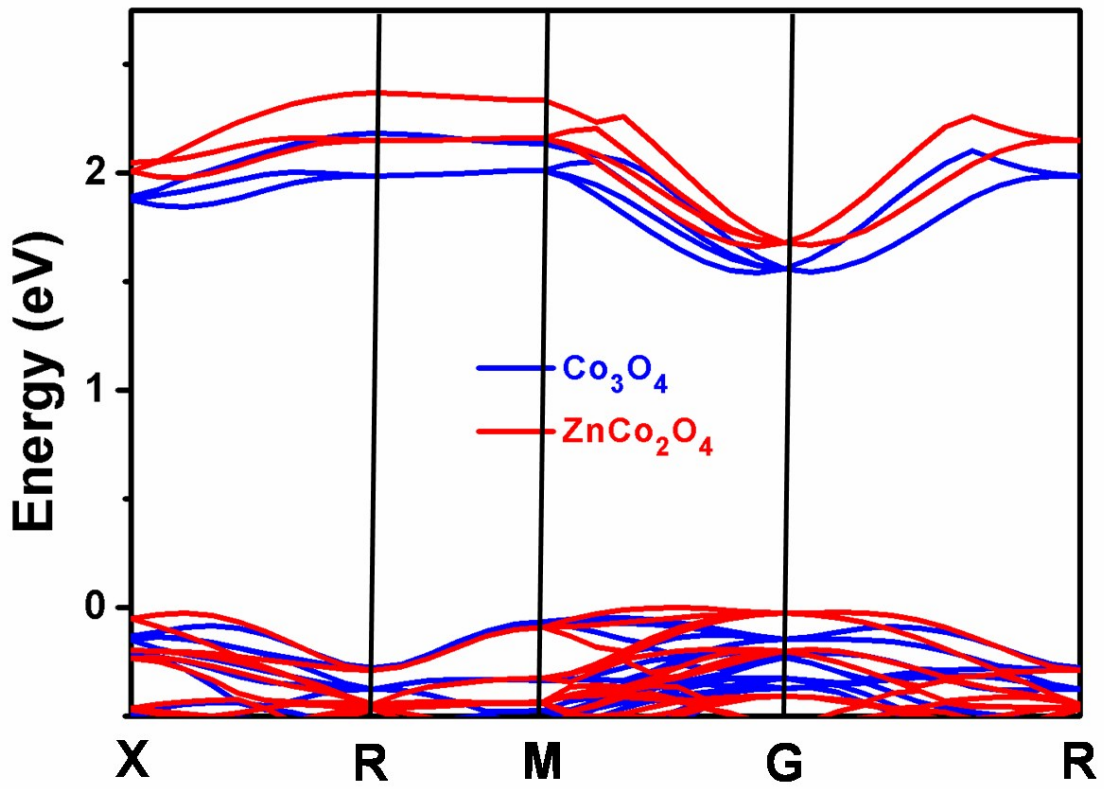

Figure S2. Calculated band structures of spinel Co_3O_4 and ZnCo2O4 near the Fermi level.

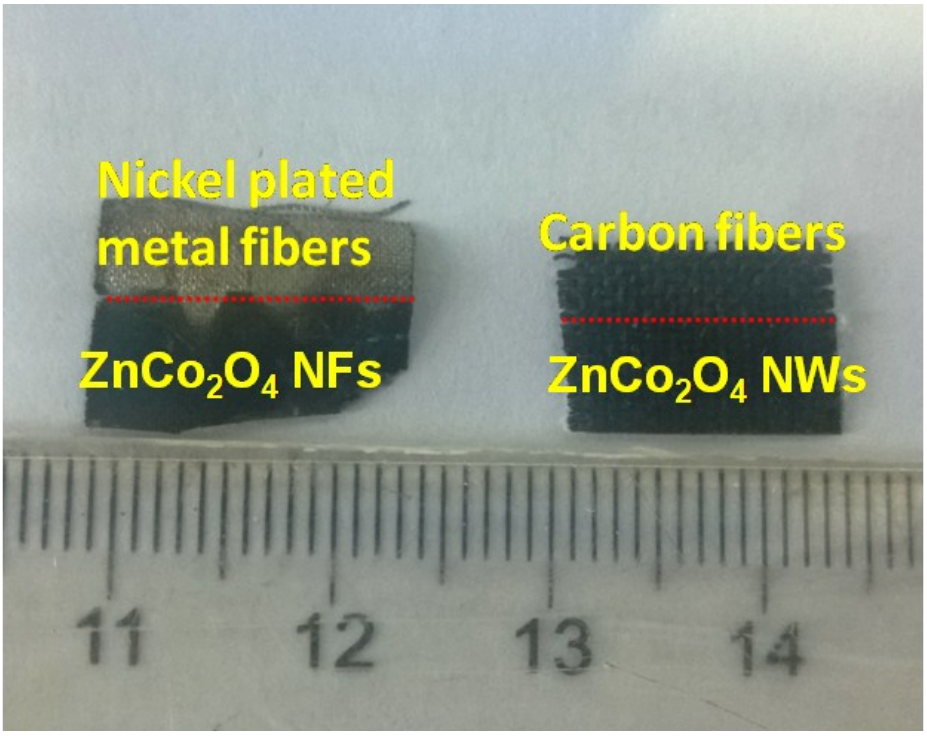

Figure S3. Optical images of the $ZnCo_2O_4$ NFs and $ZnCo_2O_4$ NWs on the substrates.

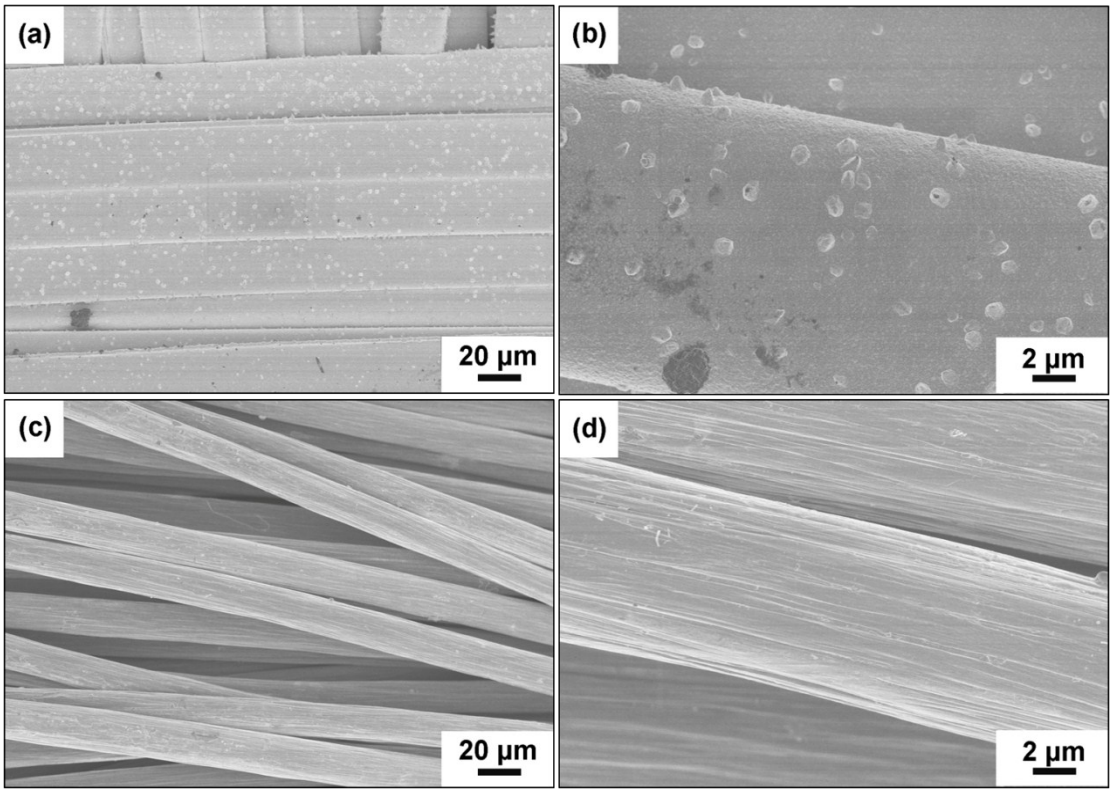

Figure S4. (a, b) Low-magnification SEM images of the nickel plated steel mesh; (c, d) low-magnification SEM images of the carbon fibers.

Figure S5. (a, b) Optical photographs of flowers in the real world; c, d) low-magnification and high-magnification SEM images of the $ZnCo_2O_4$ NFs.