

A facile template method to fabricate one-dimensional $\text{Fe}_3\text{O}_4@\text{SiO}_2@\text{C}/\text{Ni}$ microtubes with efficient catalytic and adsorption performance

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Supporting Information

Preparation of $\text{Fe}_3\text{O}_4@\text{SiO}_2@\text{C}/\text{Ni}$ (700 °C)

The synthesis of $\text{Fe}_3\text{O}_4@\text{SiO}_2@\text{C}/\text{Ni}$ (700 °C) was similar with that of $\text{Fe}_3\text{O}_4@\text{SiO}_2@\text{C}/\text{Ni}$ (500 °C) just changing the carbonization temperature from 500 °C to 700 °C.

Preparation of $\text{Fe}_3\text{O}_4@\text{C}/\text{Ni}$ and $\text{Fe}_3\text{O}_4@\text{SiO}_2@\text{C}/\text{Ni}$ (200 μL)

$\text{Fe}_3\text{O}_4@\text{C}/\text{Ni}$ and $\text{Fe}_3\text{O}_4@\text{SiO}_2@\text{C}/\text{Ni}$ (200 μL) were prepared through annealing their precursor ($\text{FeOOH}@\text{PDA}-\text{Ni}^{2+}$ and $\text{FeOOH}@\text{SiO}_2@\text{PDA}-\text{Ni}^{2+}$ (200 μL)) at 500°C for 5 h in the protection of N_2 gas. The synthesis of $\text{FeOOH}@\text{PDA}-\text{Ni}^{2+}$ and $\text{FeOOH}@\text{SiO}_2@\text{PDA}-\text{Ni}^{2+}$ (200 μL) was similar with that of $\text{FeOOH}@\text{SiO}_2@\text{PDA}-\text{Ni}^{2+}$ (50 μL) just changing the amount of TEOS from 50 μL to 0 μL and 200 μL .

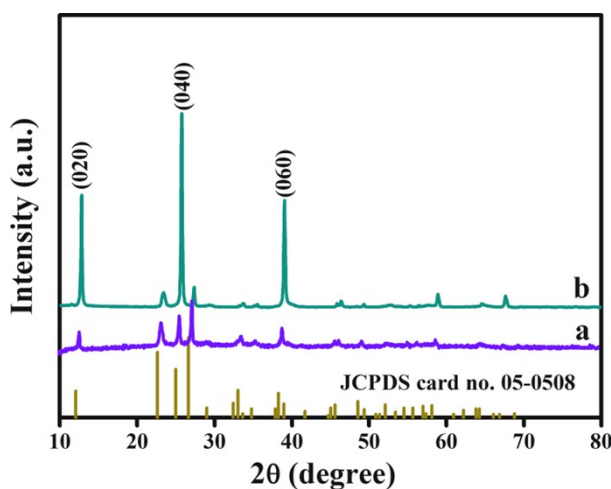


Figure S1. XRD patterns of the as-prepared MoO₃ (b) and MoO₃@FeOOH (a)

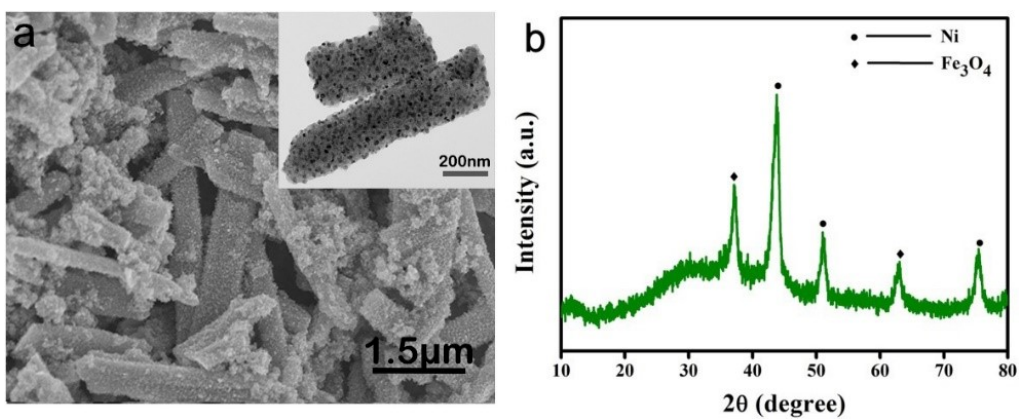


Figure S2. SEM and TEM image (a), XRD pattern of the $\text{Fe}_3\text{O}_4@\text{SiO}_2@\text{C}/\text{Ni}$ -700 (b)

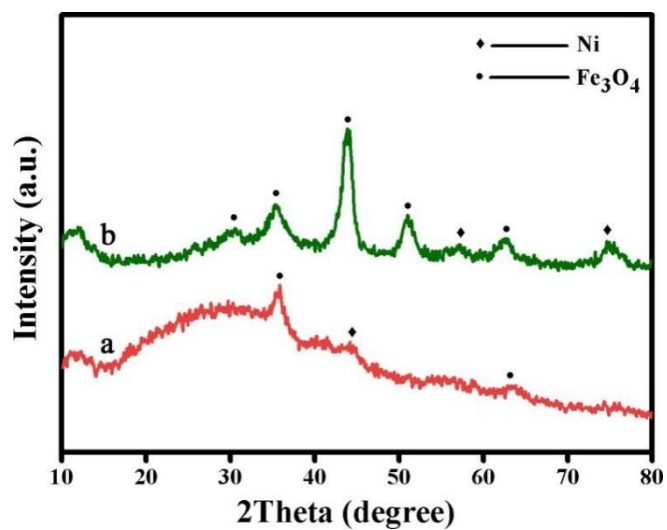


Figure S3. XRD pattern: (a) $\text{Fe}_3\text{O}_4@\text{SiO}_2@\text{C}/\text{Ni}$ microtubes (200uL TEOS); (b) $\text{Fe}_3\text{O}_4@\text{C}/\text{Ni}$ microtubes

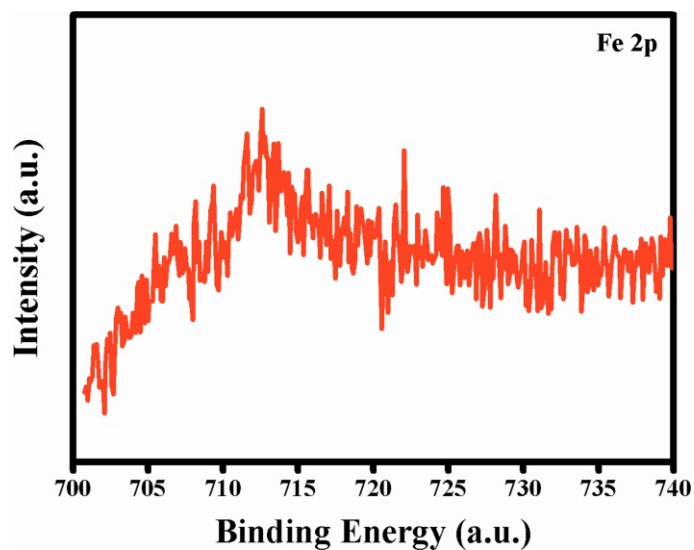


Figure S4. High-resolution of Fe 2p

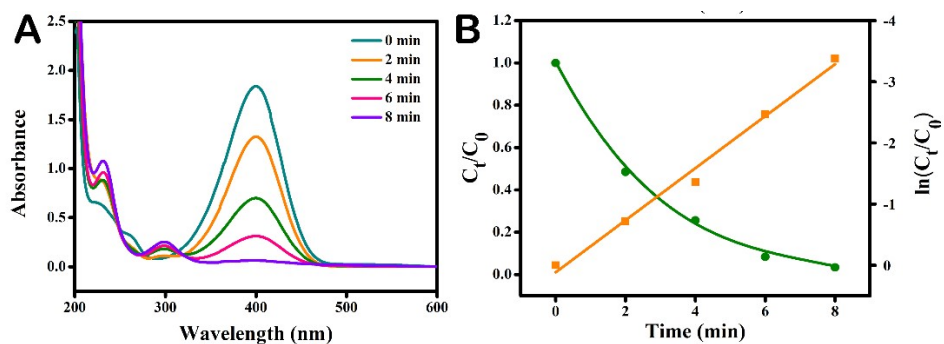


Figure S5. (A): successive reduction of 4-NP using $\text{Fe}_3\text{O}_4@\text{C}/\text{Ni}$ as catalyst; (B): C_t/C_0 and $\ln(C_t/C_0)$ versus reaction time for the reduction of 4-NP over $\text{Fe}_3\text{O}_4@\text{C}/\text{Ni}$.

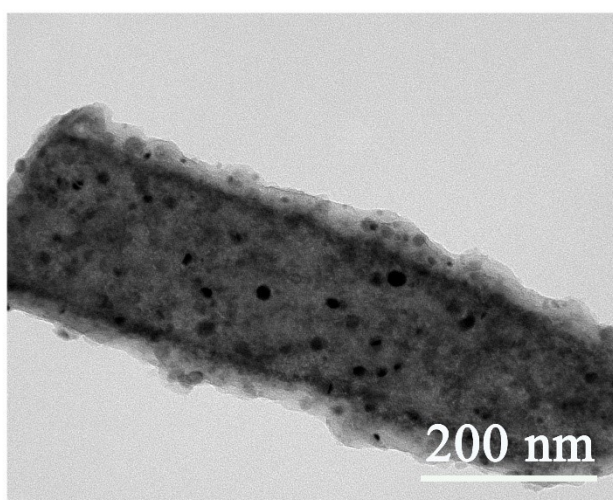


Figure S6. TEM image of $\text{Fe}_3\text{O}_4@\text{SiO}_2@\text{C}/\text{Ni}$ after five cycle

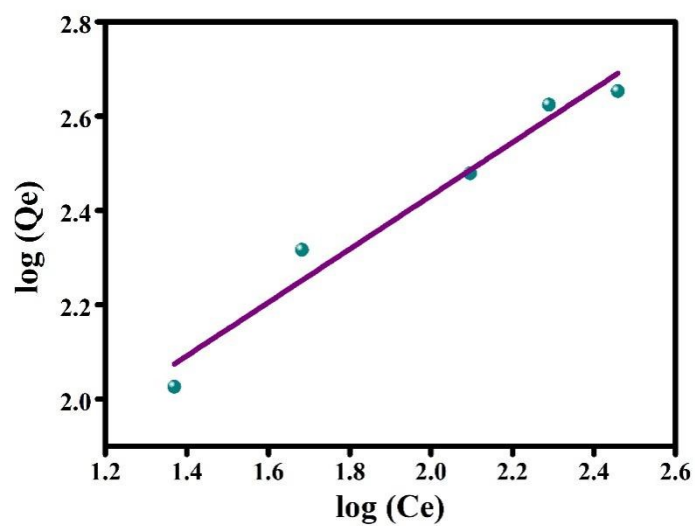


Figure S7. Linear fitting of adsorption isotherms plots based on Freundlich model