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Supporting Information

## Self-assembled Mn<sub>3</sub>O<sub>4</sub> nano-clusters over carbon nanotube threads with enhanced supercapacitor performance

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Fig. S1 Textural properties of  $Mn_3O_4$  and  $Mn_3O_4@CNT$  (a)  $N_2$  adsorption-desorption isotherms, (b) Pore size distribution curves.

Table S1 Summary of textural properties of Mn<sub>3</sub>O<sub>4</sub> and Mn<sub>3</sub>O<sub>4</sub>@CNT nano-clusters.

Sample	Crystallite size (nm)	Specific surface area $(m^2/g)$	Total pore volume $(cm^{3/g})$	Average pore diameter (nm)
Mn <sub>3</sub> O <sub>4</sub>	35.6	44	0.17	15.6
Mn <sub>3</sub> O <sub>4</sub> @CNT	36.9	45	0.23	20.8

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Fig. S2 FE-SEM images for Mn<sub>3</sub>O<sub>4</sub> nanoclusters



Fig. S3 (a) Cyclic voltamograms of pristine Mn<sub>3</sub>O<sub>4</sub> nanoparticles electrode at different scan rates, (b) galvanostatic charge-discharge curves of pristine Mn<sub>3</sub>O<sub>4</sub> nanoparticles at different potential



Fig. S4 calculated ESR for Mn<sub>3</sub>O<sub>4</sub>@CNT at different current densities.