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Supporting information

beta-manganese dioxide nanoflowers self-assembled by thin nanoplates with enhanced supercapacitive performance Li-Li Yu ^{a*}, Jun-Jie Zhu ^{a,c}, Jing-Tai Zhao ^{b,c*}

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Fig. S1 Nitrogen adsorption-desorption isotherm plots and corresponding pore size distributions (insets) of samples S_{RT} -350



Fig. S2 SEM images of samples S_{RT} - H_2SO_4 (a) and S_{RT} - HNO_3 (b)



Fig. S3 Charge/discharge cycling life tests of S_{RT} -350 at the current density of 2 A g⁻¹

E and P calculation

Energy density (E) and power density (P) of the as-prepared porous nanoflowers were calculated from the following equations:

$$E = \frac{1}{2}C(\Delta V)^2$$
$$P = \frac{E}{t}$$

Where E (W h Kg⁻¹), C (F g⁻¹), ΔV (V), and P (W kg⁻¹) are energy density specific capacitance, potential window of discharge time, and power density, respectively.