

Supplementary Materials: Well-Shaped Sulfonic Organosilica Nanotubes with High Activity for Hydrolysis of Cellobiose

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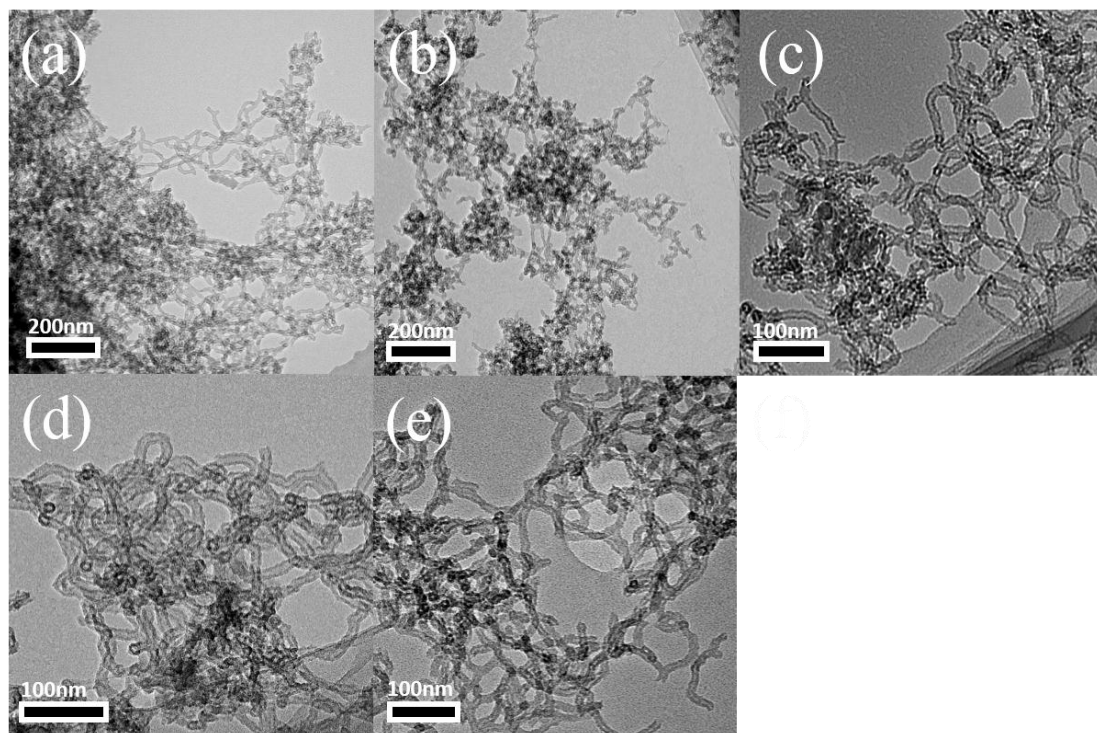


Figure S1. TEM images of mercaptopropyl organosilica nanotubes with different sulfur contents. (a) SH_{0.2}-Et-SNT; (b) SH_{0.3}-Et-SNT; (c) SH_{0.1}-Ph-SNT; (d) SH_{0.2}-Ph-SNT and (e) SH_{0.3}-Ph-SNT.

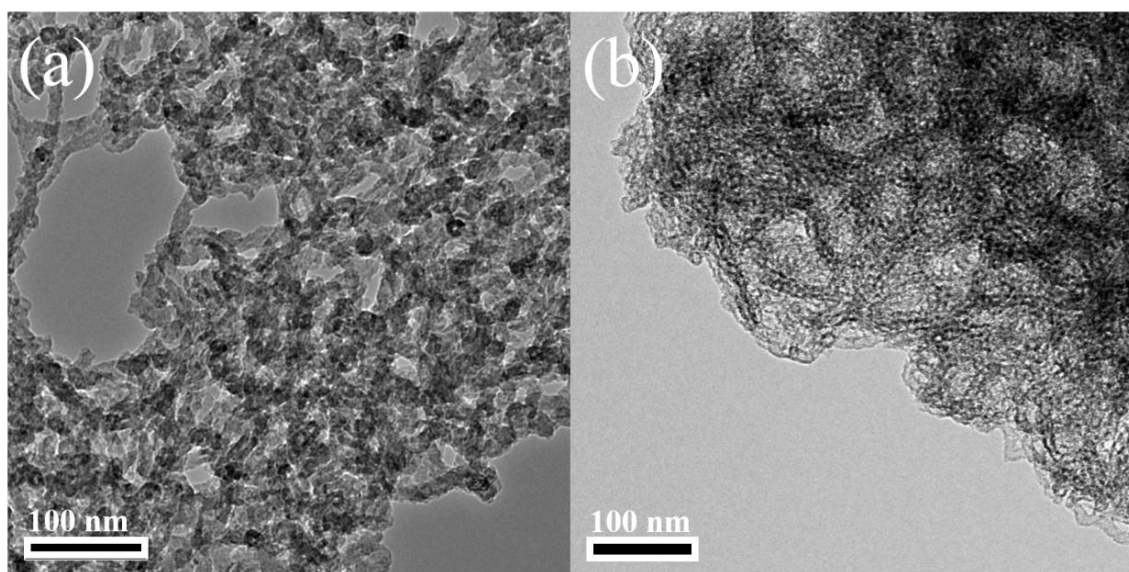


Figure S2. TEM images of (a) SH_{0.5}-Ph-SNT and (b) SH-SNT.

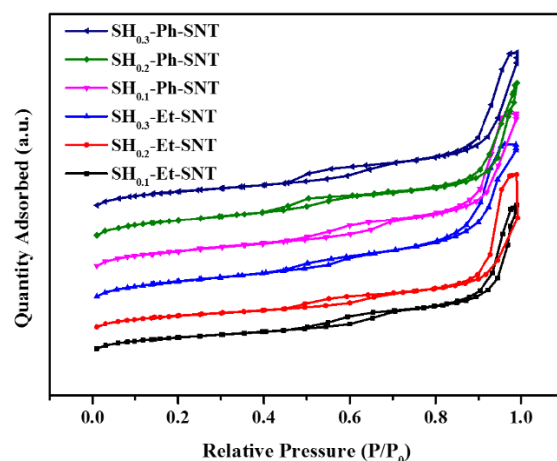


Figure S3. Nitrogen adsorption-desorption isotherms of SH_{0.1}-Et-SNT, SH_{0.2}-Et-SNT, SH_{0.3}-Et-SNT, SH_{0.1}-Ph-SNT, SH_{0.2}-Ph-SNT and SH_{0.3}-Ph-SNT.

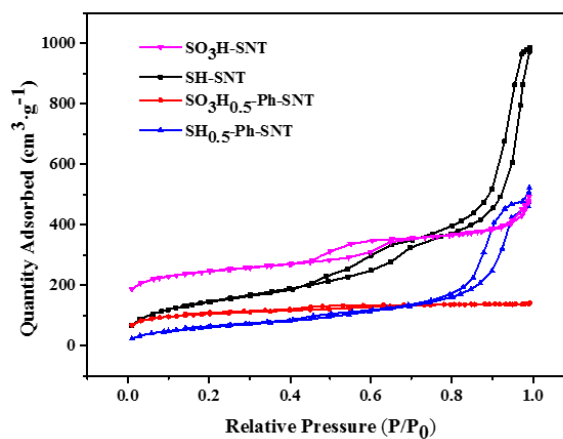


Figure S4. Nitrogen adsorption isotherms of SH-SNT, SO₃H-SNT, SH_{0.5}-Ph-SNT and SO₃H_{0.5}-Ph-SNT.

Table S1. Physicochemical properties of SH_x-Et-SNT and SH_x-Ph-SNT.

Sample	BET Surface Area (m ² ·g ⁻¹) ^a	Pore Volume (cm ³ ·g ⁻¹) ^b	Pore Size (nm) ^c
SH _{0.1} -Et-SNT	799	2.6	5.5
SH _{0.2} -Et-SNT	625	2.2	5.2
SH _{0.3} -Et-SNT	490	1.4	4.6
SH _{0.1} -Ph-SNT	834	2.1	6.5
SH _{0.2} -Ph-SNT	560	1.6	4.4
SH _{0.3} -Ph-SNT	530	1.2	5.4
SH _{0.5} -Ph-SNT	365	-	-
SH-SNT	540	1.6	6.3

^a Surface area was determined using the Brunauer-Emmett-Teller (BET) model; ^b Pore volume was obtained at the relative pressure of 0.8; ^c Pore size was estimated Barrett-Joyner-Halenda (BJH) method with adsorption branch.

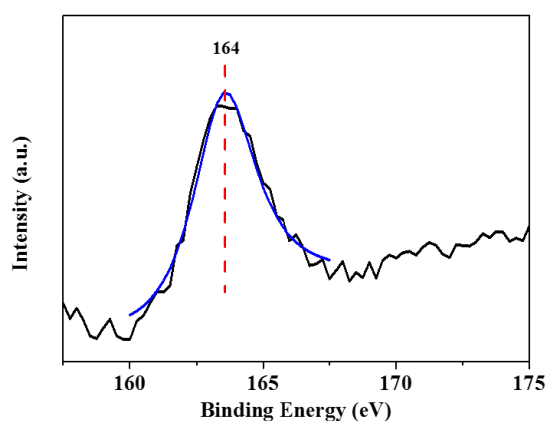


Figure S5. XPS spectrum of SH_{0.1}-Et-SNT.

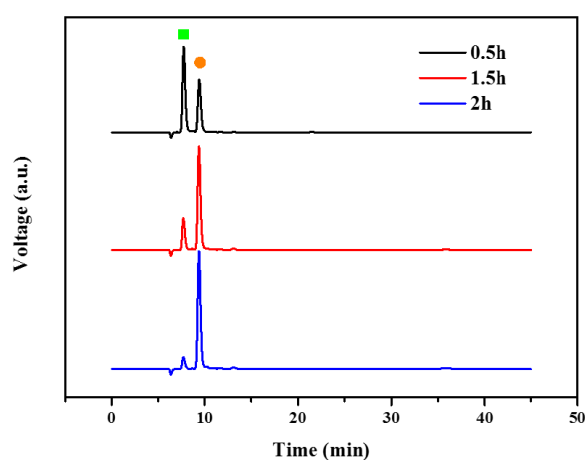


Figure S6. Liquid chromatography spectra of mixture including cellulose (green square) and glucose (orange circle) after reaction using SO₃H_{0.3}-Ph-SNT as catalyst.

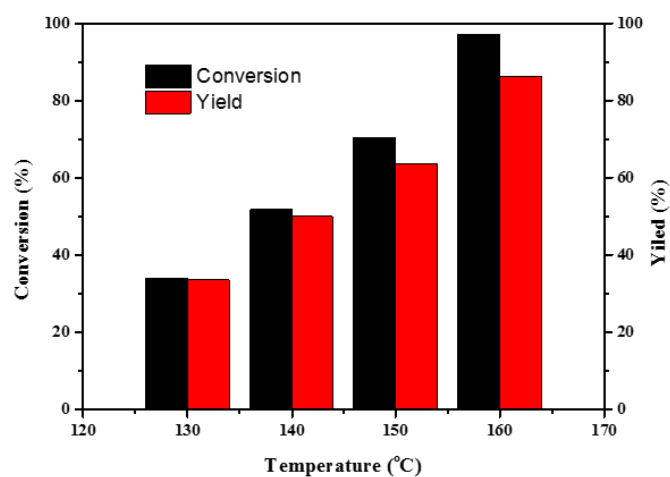


Figure S7. Temperature effect on hydrolysis of cellulose using SO₃H_{0.1}-Ph-SNT as catalyst. Hydrolysis of cellulose by using SO₃H_{0.1}-Ph-SNT as catalysts at 130 °C, 140 °C, 150 °C and 160 °C, respectively. Reaction conditions: 50 mL autoclave, 20 mL deionized water, substrates: acid active sites = 14, 0.2 g cellulose, 2.5 MPa Nitrogen, 800 rpm, 2 h.