

## Supporting information (SI)

### Oxygen Vacancies Enable Excellent Electrochemical Kinetics of Carbon Coated

#### Mesoporous SnO<sub>2</sub> Nanoparticles in Lithium Ion Battery

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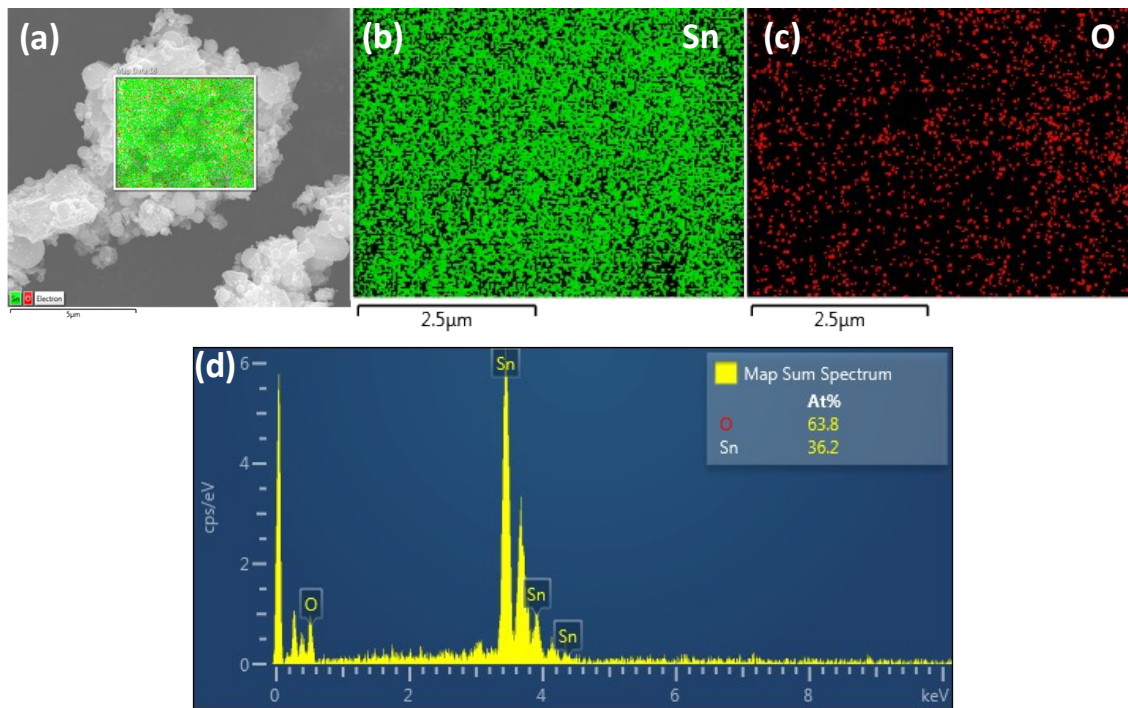
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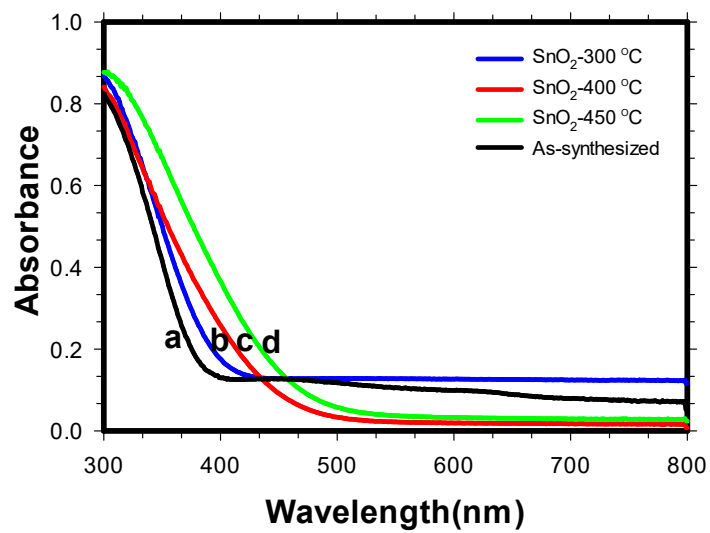
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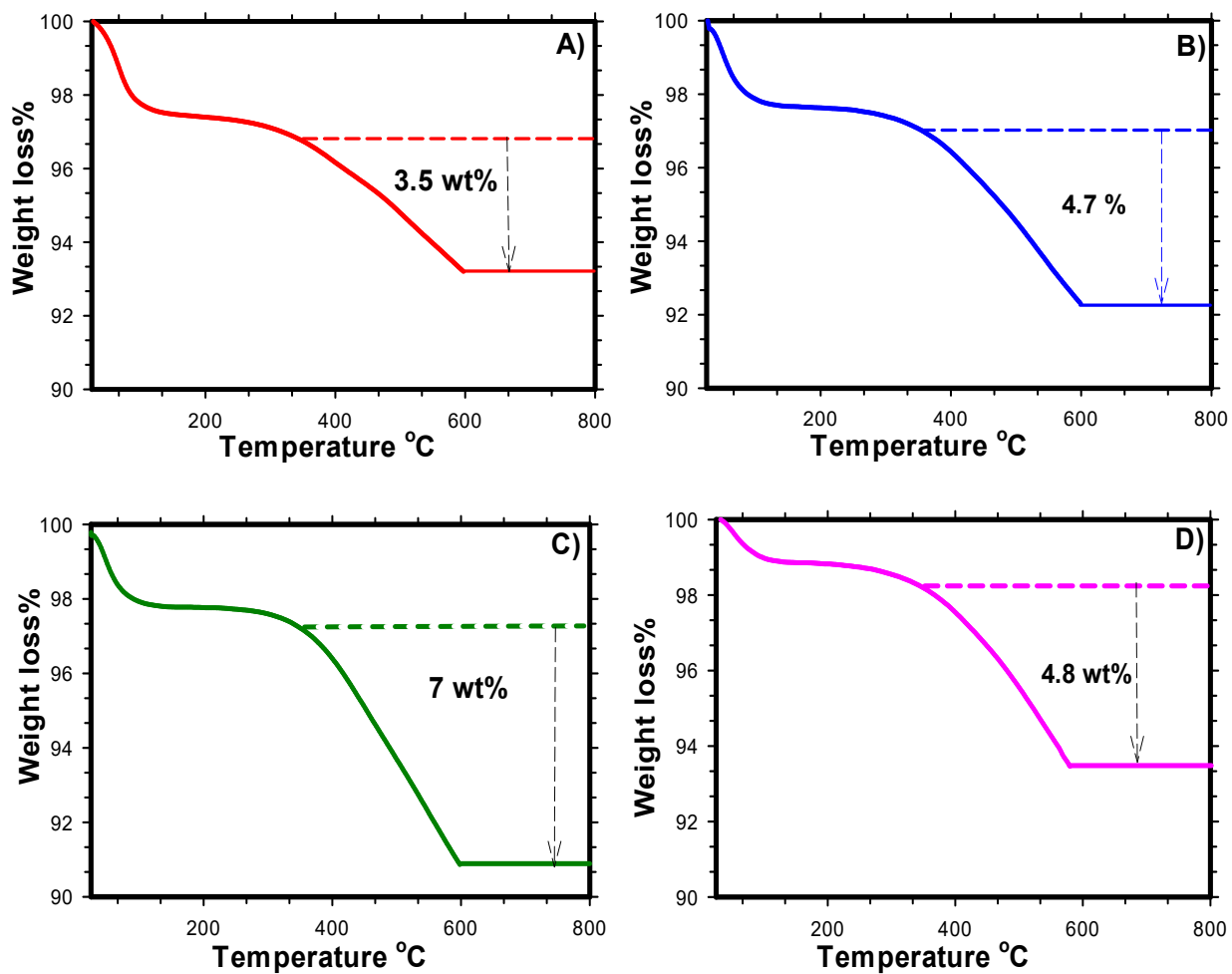
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**Fig. S1** (a) SEM image of H-SnO<sub>2</sub> particles (calcined at 400 °C), (b-c) elemental mapping of Sn and O (d) EDS analysis of H-SnO<sub>2</sub>



**Fig. S2:** UV-vis- absorption spectrum of (a) H-SnO<sub>2</sub> As-synthesized, H-SnO<sub>2</sub> calcined at (b) 300 °C, (c) 400 °C and (d) 450 °C.



**Fig. S3:** TGA curves of the non-conventional carbon coated H-SnO<sub>2</sub>: (a) NCC-1 (b) NCC-2 (c) NCC-3 and (d) NCC-4.

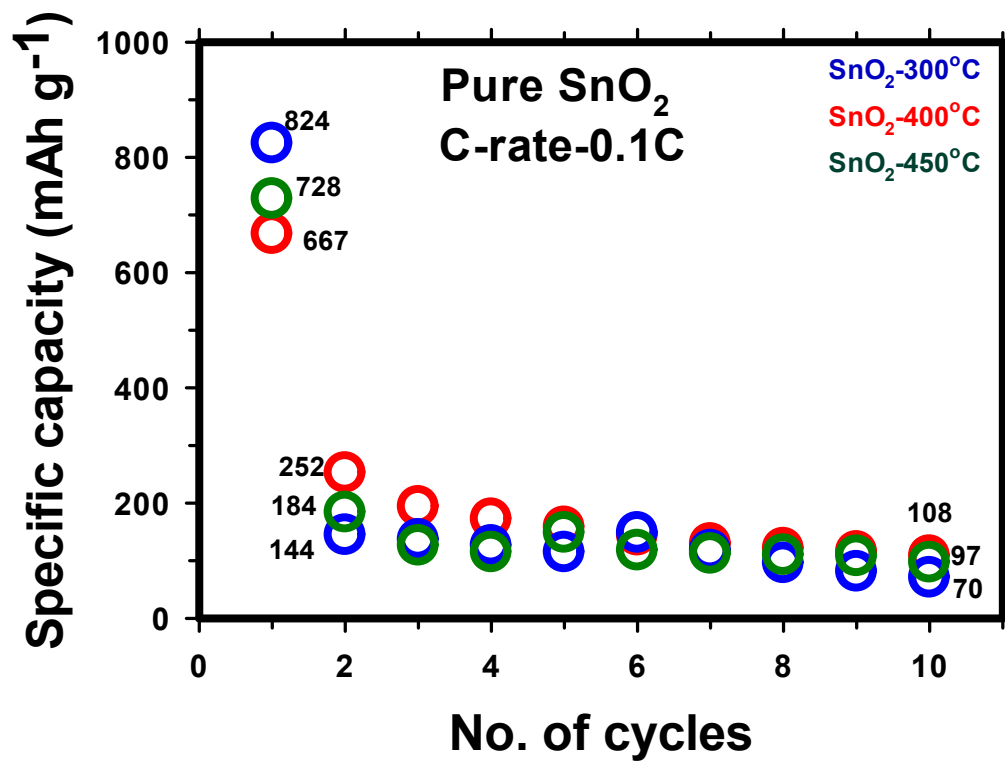
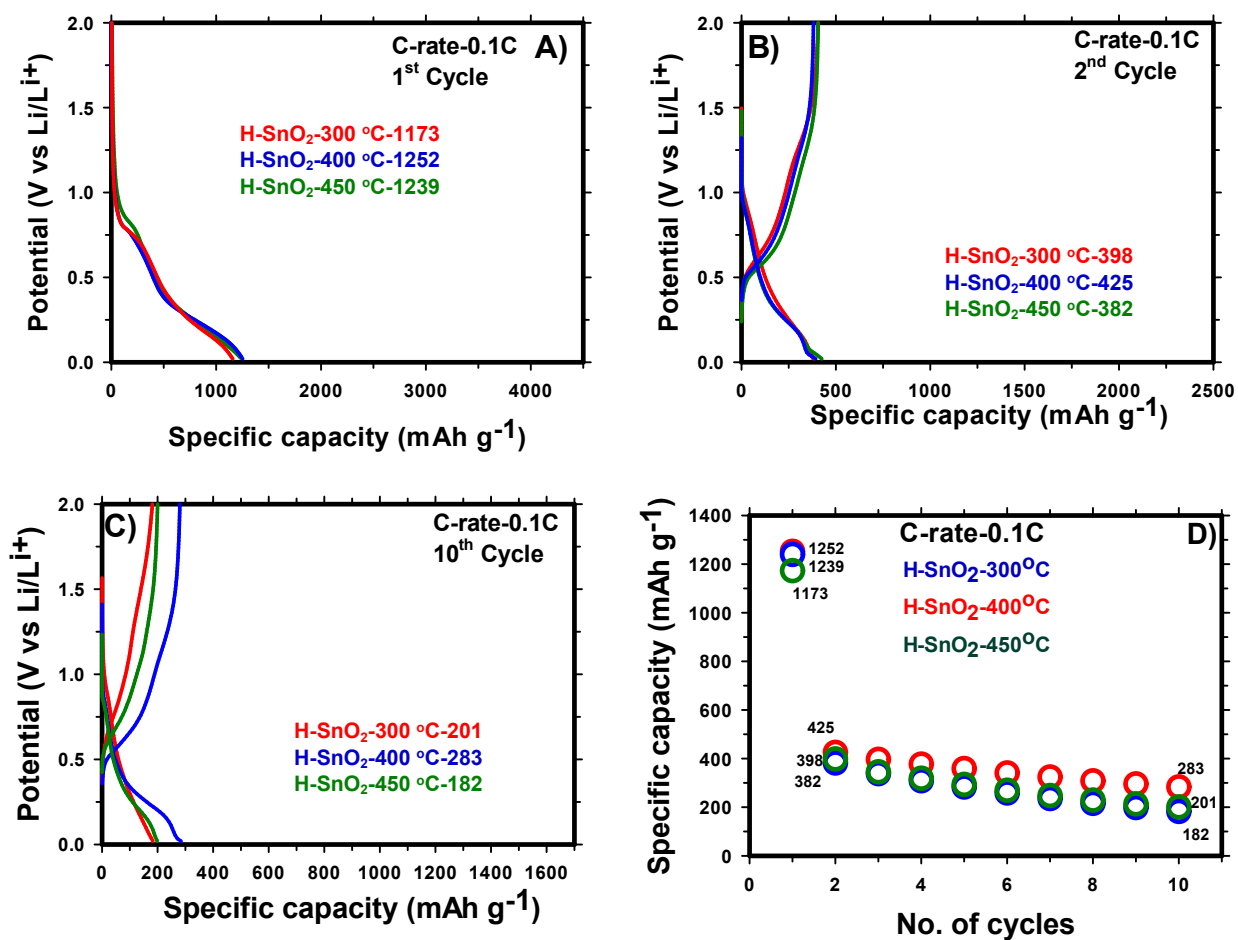
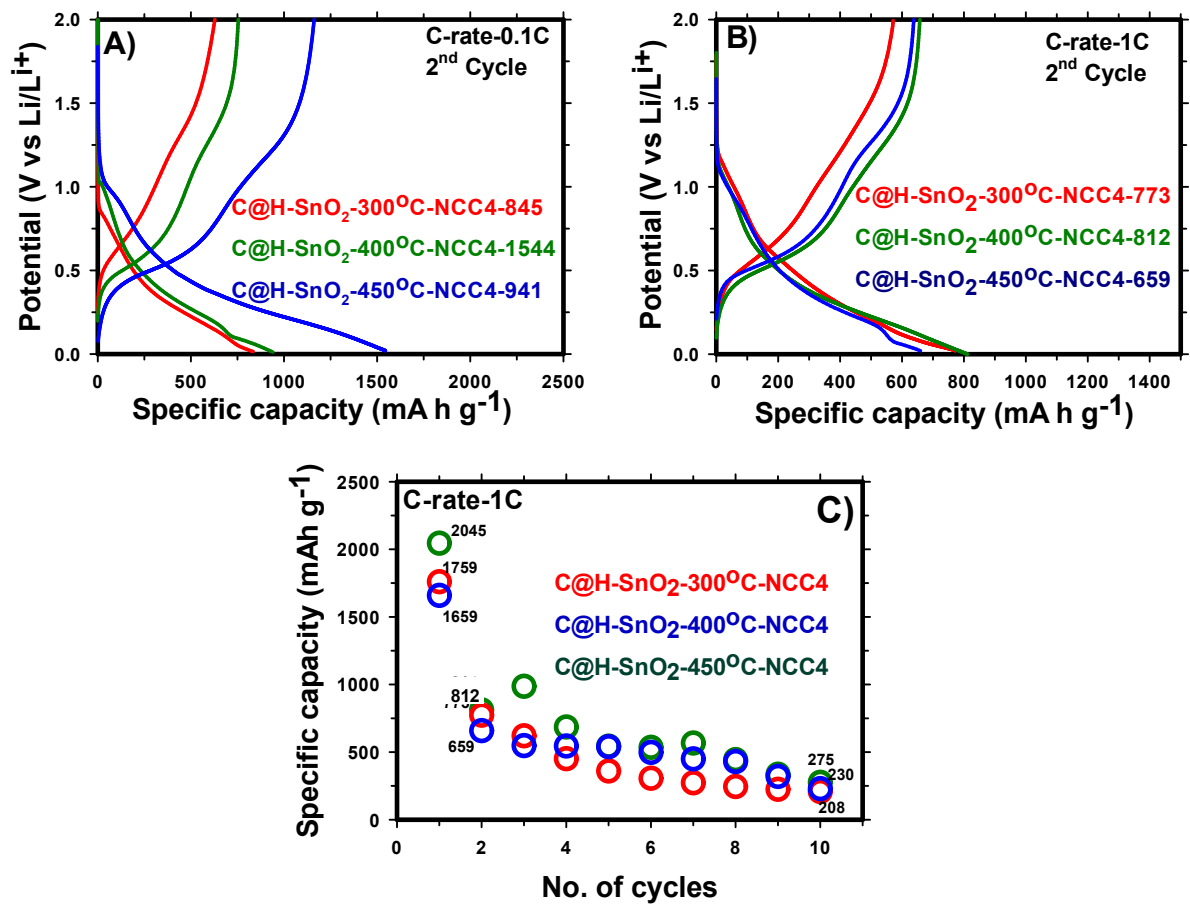


Fig. S4: Cyclic stability of pure SnO<sub>2</sub> tested at 0.1C rate



**Fig. S5:** Electrochemical performance of H-SnO<sub>2</sub> calcined at 300 °C, 400 °C and 450 °C tested at 0.1C (A) 1<sup>st</sup> cycle, (B) 2<sup>nd</sup> cycle and (C) 10<sup>th</sup> cycle and (D) cyclic stability of H-SnO<sub>2</sub> calcined at 300 °C, 400 °C and 450 °C



**Fig. S6:** Charge-discharge profile of carbon coated H-SnO<sub>2</sub> calcined at 300, 400 and 450 °C carbon coated by NCC-4 tested at (A) 0.1C, (B) 1C, (C) cyclic stability tested at 1C rate

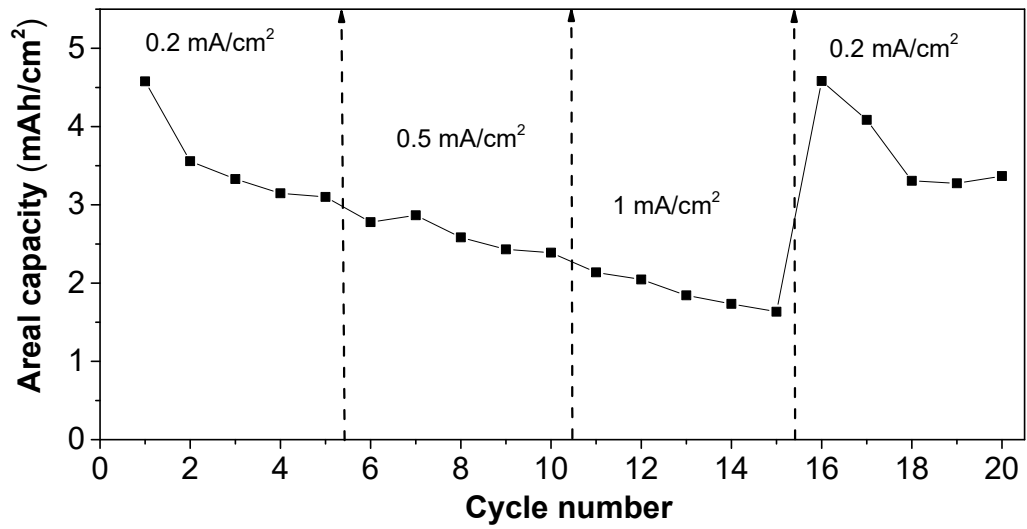


Fig. S7 Rate performance of C@H-SnO<sub>2</sub>-400 °C-NCC4



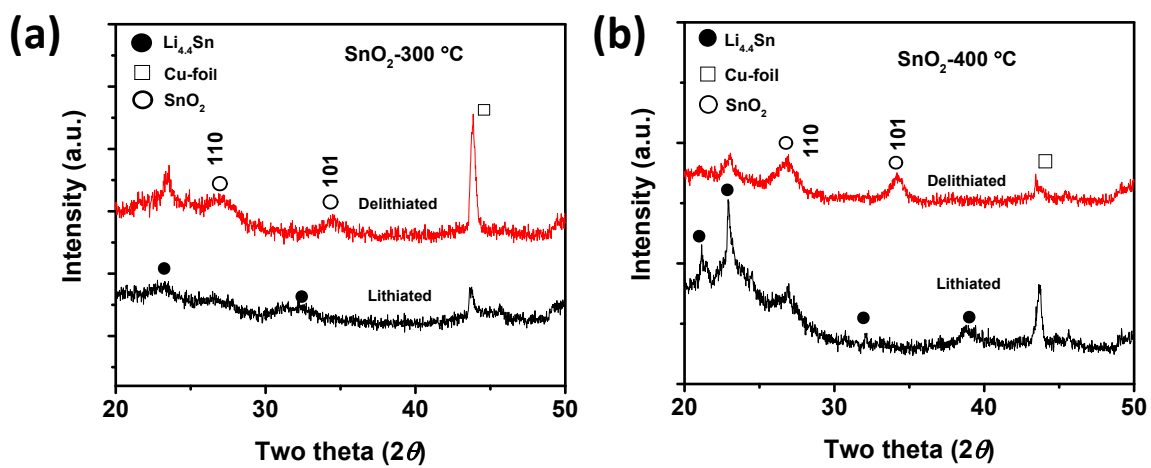


Fig. S8: XRD pattern for H-SnO<sub>2</sub> calcined at (a) 300 °C (b) 400 °C carbon coated by NCC-4

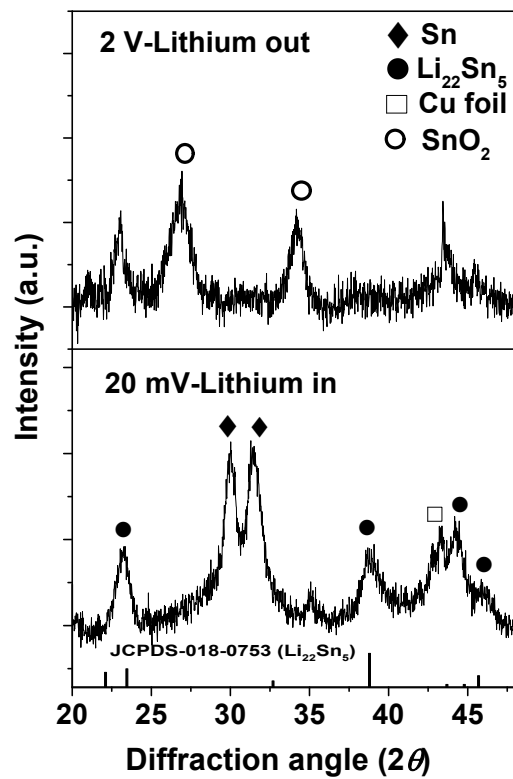


Fig. S9. Ex-situ XRD of H-SnO<sub>2</sub> after 50 cycles.

**Table S1:** XRD data of pure H-SnO<sub>2</sub> calcined at different temperatures and C@H-SnO<sub>2</sub> nanoparticles, UV-Vis data with UV-vis and TGA results

Sample	Crystallite size (nm)	Lattice Parameters		Unit cell volume (Å <sup>3</sup> )	Band gap (e.V)	Carbon content (TGA)
		a=b	c			
H-SnO <sub>2</sub> -300 °C	4.5	4.7311	3.1867	71.33	3.03	
H-SnO <sub>2</sub> -400 °C	5.1	4.7342	3.186	71.4	2.76	
H-SnO <sub>2</sub> -450 °C	6.8	4.7381	3.186	71.54	2.51	
C@H-SnO <sub>2</sub> -NCC1	7.4	4.7461	3.189	71.8		3.5%

C@H-SnO <sub>2</sub> -NCC2	7.3	4.7459	3.1752	71.52		4.7%
C@H-SnO <sub>2</sub> -NCC3	7.1	4.708	3.189	71.4		7.0%
C@H-SnO <sub>2</sub> -NCC4	7.7	4.7581	3.175	71.9		4.8%

**Table S2.** XPS core level of Sn3d, O1s and C1s of C@H-SnO<sub>2</sub> NPs by CC1 and NCC4 condition

Sample	Crystallite size (nm)	Binding energy (eV)	Binding energy (eV) C 1s	$\Delta$ (eV) for Sn 3d (spin-orbit coupling)	Binding energy (eV) O-1s (Deconvoluted)
		Sn-3d <sub>5/2</sub> and Sn-5d <sub>3/2</sub>			
C@H-SnO <sub>2</sub> -CC	4.5	487.087	284.5	8.417	530.909
		495.310	286.199		531.587

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C@H-SnO <sub>2</sub> -NCC4	7.7	486.671	284.5	8.031	530.617
		495.210	285.437		532.179
			288.629		

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