

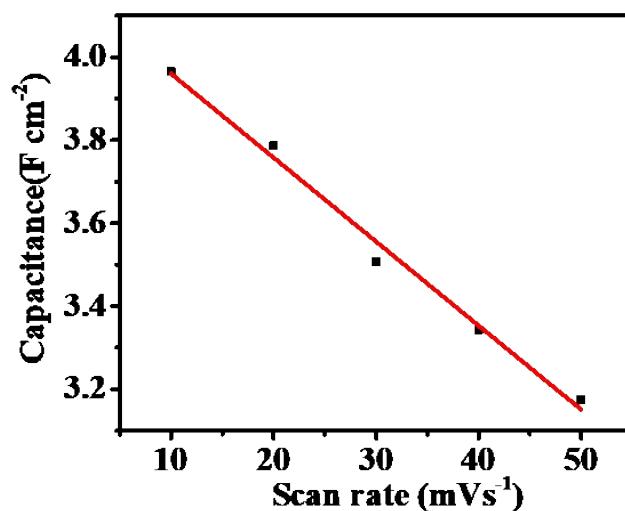
## Electronic supplementary information

### Electrochemical, top-down nanostructured pseudocapacitive electrode for enhanced specific capacitance and cycling efficiency

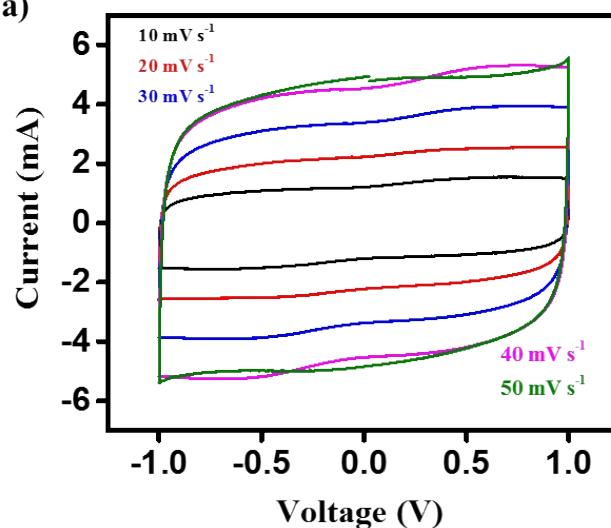
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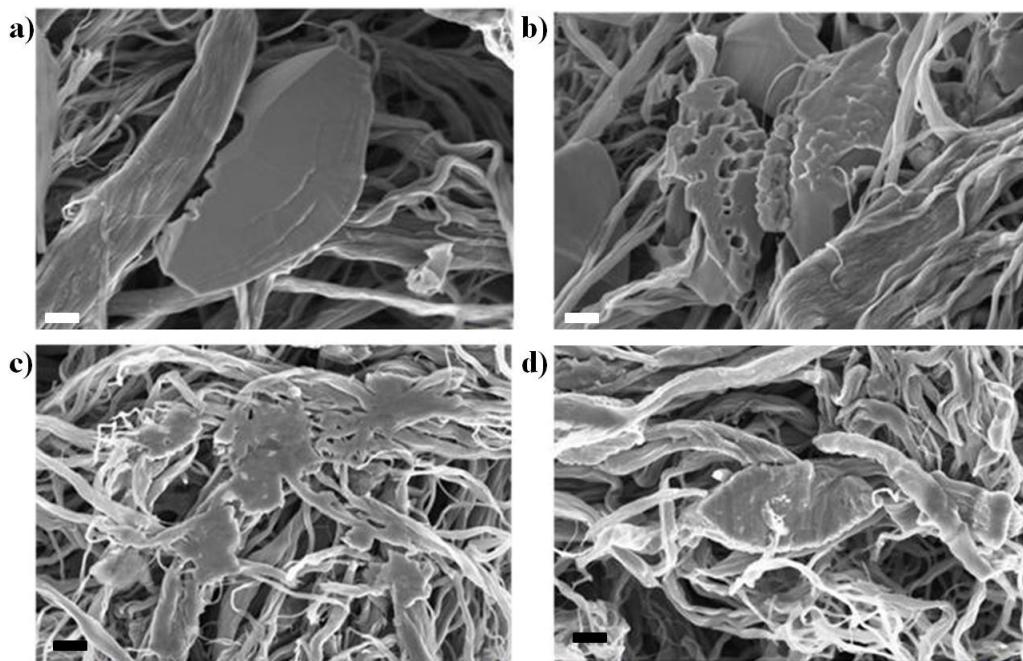


**Fig S1.** a)  
capacitance with



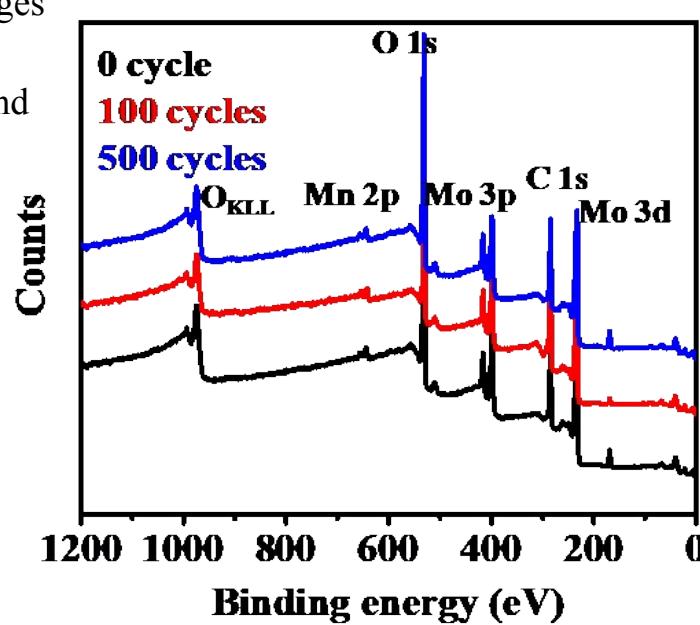
Variation of areal  
scan rate.

**Fig S2.** Cycling voltamograms of the CNT bucky paper at different scan rate indicating no pseudocapacitive component.

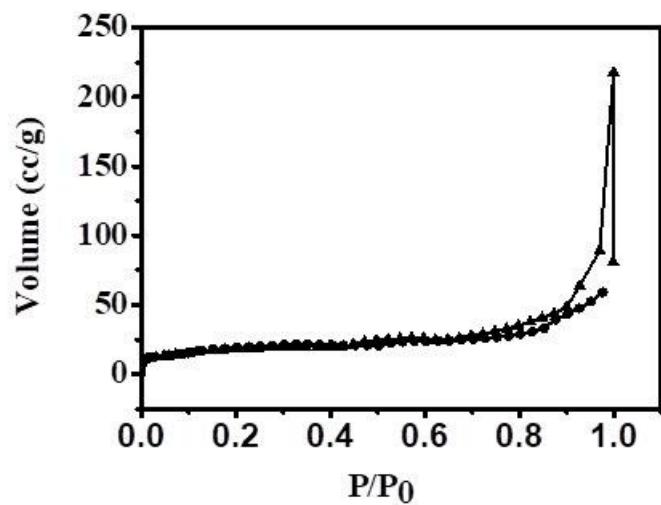


**Fig S3.** SEM images  
CNT/MnMoO<sub>4</sub>  
(a, b) 100 cycle and  
cycles.

of  
electrode after  
(c, d) 500



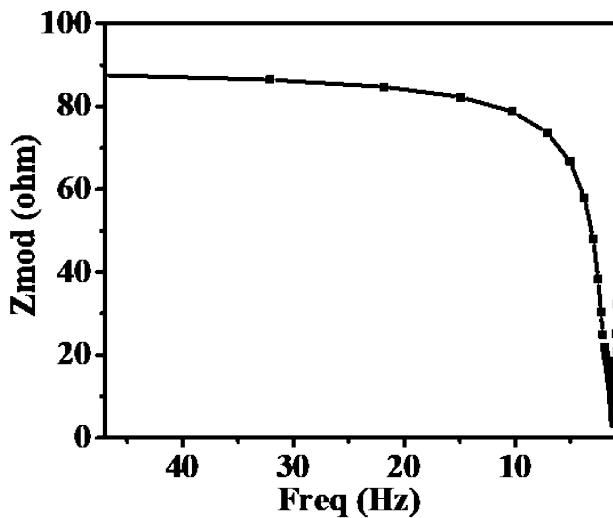
**Fig S4.** Survey spectrum of CNT/MnMoO<sub>4</sub> electrode at different stages of electrochemical cycling.



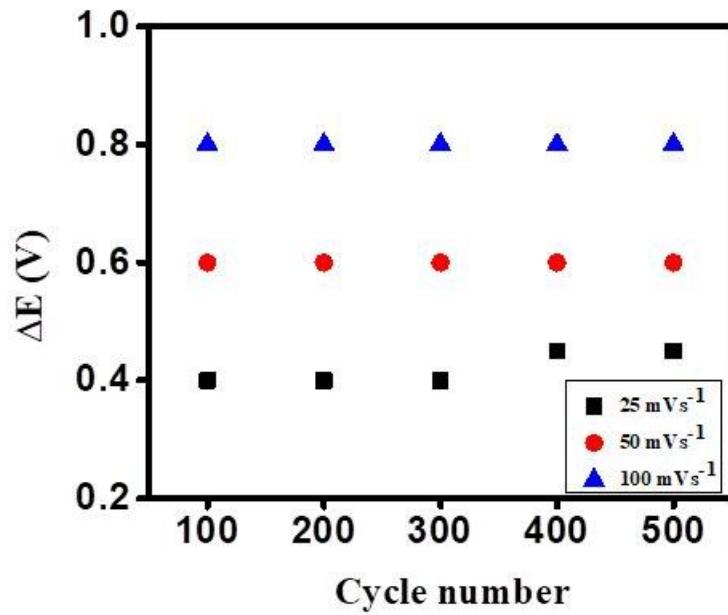
**Fig S5.** Nitrogen adsorption isotherm for CNT bucky paper.

**Table T1:** Comparison of surface area parameters for CNT bucky paper and CNT/MnMoO<sub>4</sub> electrode at different cycles.

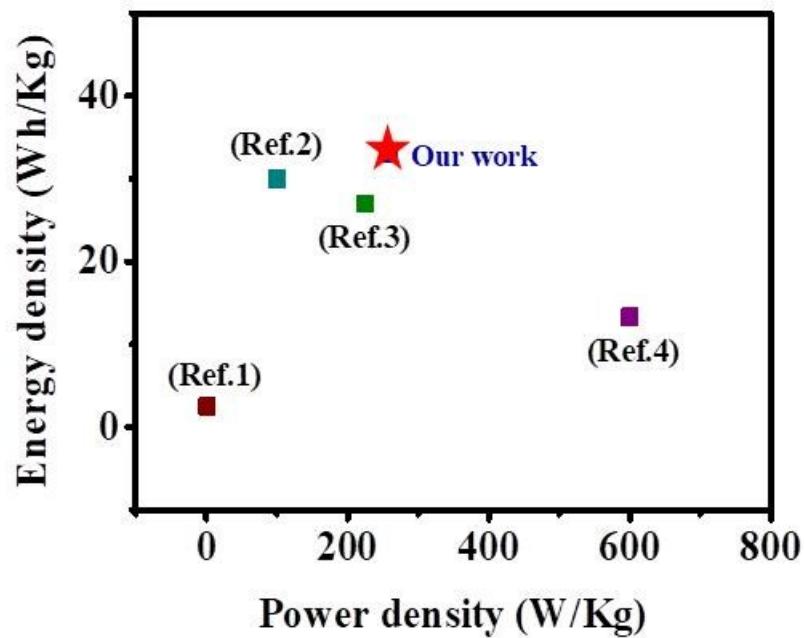
Materials	BET Specific surface area (m <sup>2</sup> /g)	Pore volume (cc/g)	% change in SSA	External surface area from <i>V-t</i> plots (cc/g.A)
CNT- bucky paper	714	0.33	+ 0 %	3.5
CNT/MnMoO <sub>4</sub> (0 cycles)	1086	10.37	+ 52%	2.2
CNT/MnMoO <sub>4</sub> (100 cycles)	1337	13.62	+ 87%	1.8
CNT/MnMoO <sub>4</sub> (500 cycles)	1726	13.74	+ 142%	2.9



**Fig S6.** Bode plot indicating the dependence of phase angle on frequency of CNT/MnMoO<sub>4</sub> electrode.



**Fig S7.** Evidence of mass diffusion during electrochemical cycling of CNT/MnMoO<sub>4</sub> electrode.



**Fig S8.** Ragone plot for the comparison of energy and power density with other reports based on MnO<sub>2</sub>-CNT hybrid electrode.

## **References:**

1. Y. Zhao, M. P. Li, S. Liu, and M. F. Islam, *ACS Applied Materials & Interfaces*, 2017, **9**, **28**, 23810.
2. E. M. Jin, J. G. Lim, and S. M. Jeong, *Journal of Industrial and Engineering Chemistry*, 2017, **54**, 421.
3. D. Gueon and J. H. Moon, *ACS Sustainable Chemistry & Engineering*, 2017, **5**, **3**, 2445.
4. L. Li, Z. A. Hu, N. An, Y. Y. Yang, Z. M. Li and H. Y. Wu, *J. Phys. Chem. C*, 2014, **118**, **40**, 22865.