

## Supplementary Information

### Investigation on the Air Stability of P2-Layered Transition Metal Oxides by Nb Doping in Sodium Ion Batteries

Yanyan Chen <sup>1,2,†</sup>, Qin hao Shi <sup>2,†</sup>, Shengyu Zhao <sup>2</sup>, Wuliang Feng <sup>2,\*</sup>, Yang Liu <sup>2</sup>, Xinxin Yang <sup>2</sup>, Zhenwei Wang <sup>1,\*</sup> and Yufeng Zhao <sup>2,\*</sup>

<sup>1</sup> School of Chemical and Environmental Engineering, Shanghai Institute of Technology, Shanghai 201418, China

<sup>2</sup> Institute for Sustainable Energy & College of Sciences, Shanghai University, Shanghai 200444, China

\* Correspondence: wuliangfeng@shu.edu.cn (W.F.); wangzhenwei@sit.edu.cn (Z.W.); yufengzhao@shu.edu.cn (Y.Z.)

† These authors contributed equally to this work.

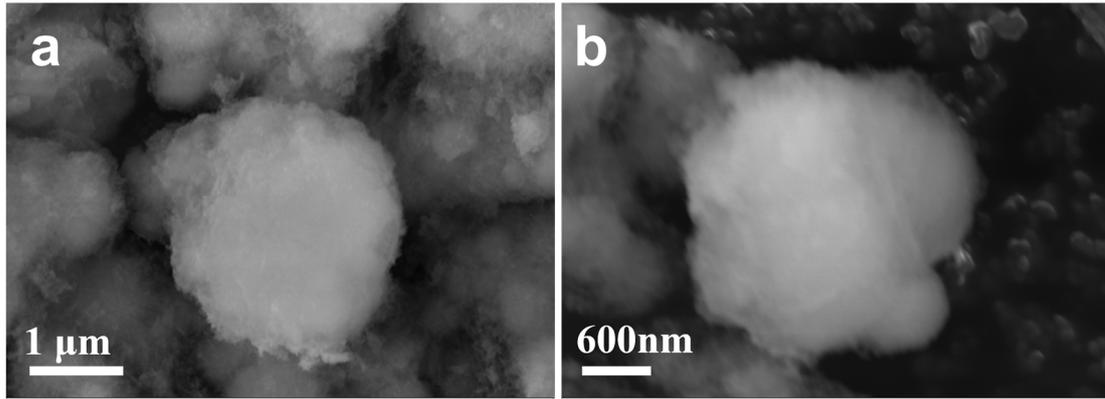


Fig. S1(a) The precursor SEM of P2-Na<sub>0.67</sub>MN and P2-Na<sub>0.67</sub>MNNb. (b) The precursor SEM of P2-Na<sub>0.67</sub>MN and P2-Na<sub>0.67</sub>MNNb.

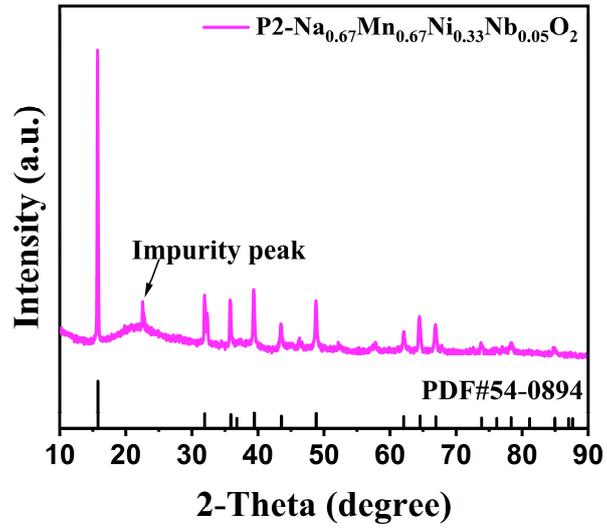


Fig. S2 The XRD patterns of P2-Na<sub>0.67</sub>Mn<sub>0.67</sub>Ni<sub>0.33</sub>Nb<sub>0.05</sub>O<sub>2</sub>.

After mixed with Nb<sub>2</sub>O<sub>5</sub>, Na<sub>0.67</sub>Mn<sub>0.67</sub>Ni<sub>0.33</sub>Nb<sub>0.03</sub>O<sub>2</sub> had no impurity peak, but Na<sub>0.67</sub>Mn<sub>0.67</sub>Ni<sub>0.33</sub>Nb<sub>0.05</sub>O<sub>2</sub> had impurity peak generation, indicating redundant niobium had not been into the bulk phase. When the doping amount of Nb was 0.05, the impurity peak in 22.5° can be labeled as Na<sub>3</sub>NbO<sub>4</sub>.

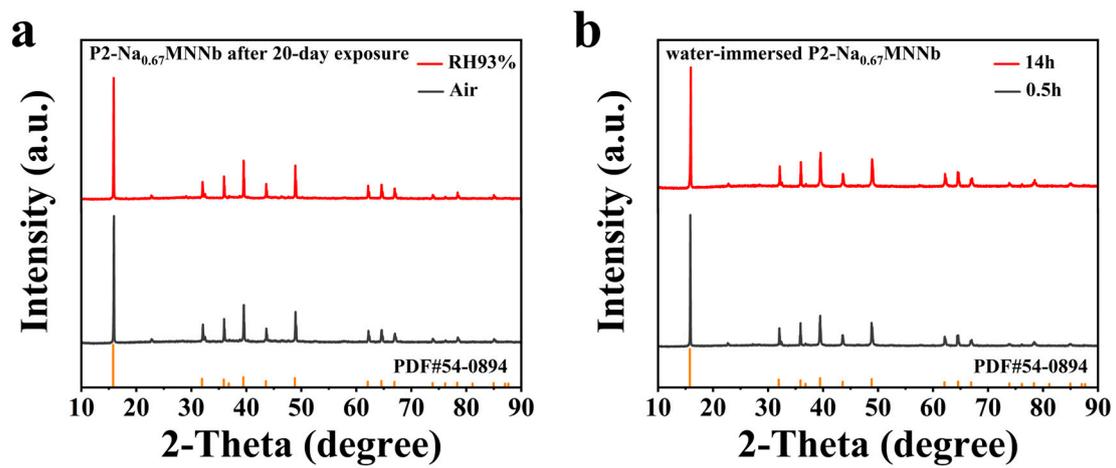


Fig. S3 (a) The XRD patterns of P2-Na<sub>0.67</sub>MNNb sample exposed in different atmosphere after 20-day exposure. (b) The XRD patterns of water-immersed P2-Na<sub>0.67</sub>MNNb sample.

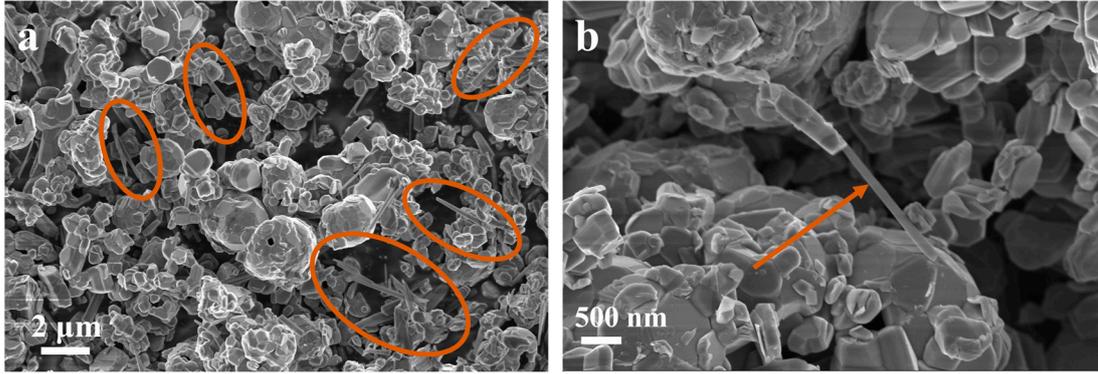


Fig. S4 (a) The SEM of P2- $\text{Na}_{0.67}\text{MN}$  after 20 days of air exposure. (b) The SEM of P2- $\text{Na}_{0.67}\text{MN}$  after 20 days of air exposure.

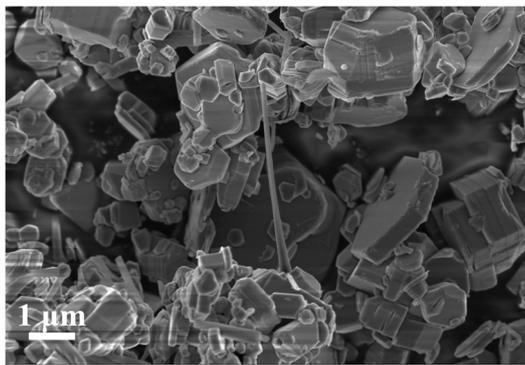


Fig. S5 The SEM of P2-Na<sub>0.67</sub>MNNb after 20 days of air exposure.

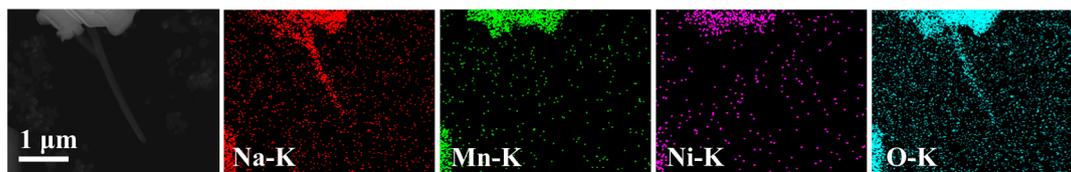


Fig. S6 SEM-EDS elemental mappings of rod-like particles in P2-Na<sub>0.67</sub>MN sample.

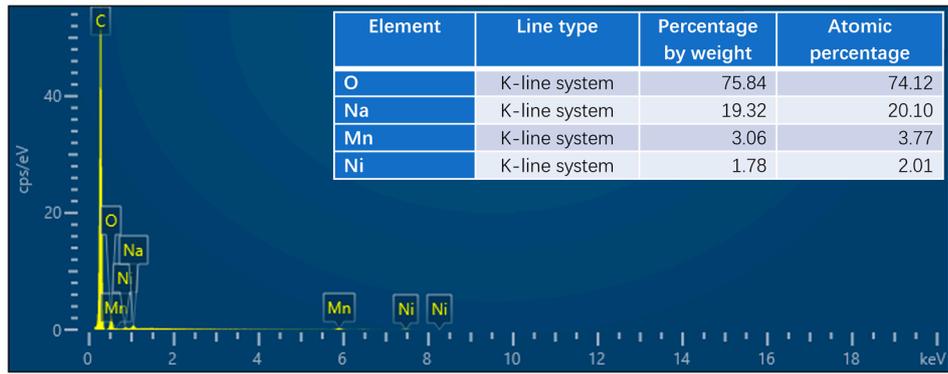


Fig. S7 The spectrum diagram of total distribution diagram for SEM-EDS of rod-like particles in P2-Na<sub>0.67</sub>Mn sample.

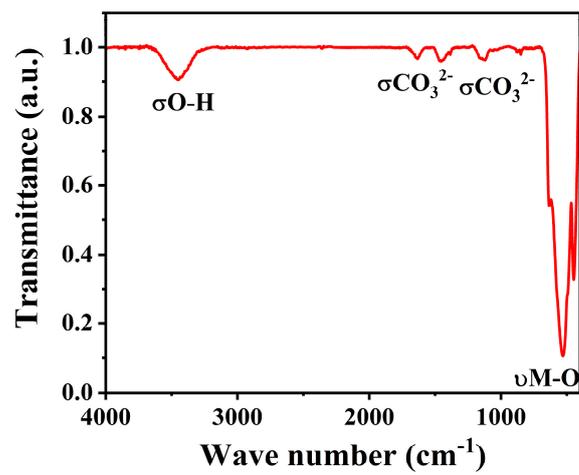


Fig. S8 The FTIR of P2-Na<sub>0.67</sub>MN after 20 days of air exposure.

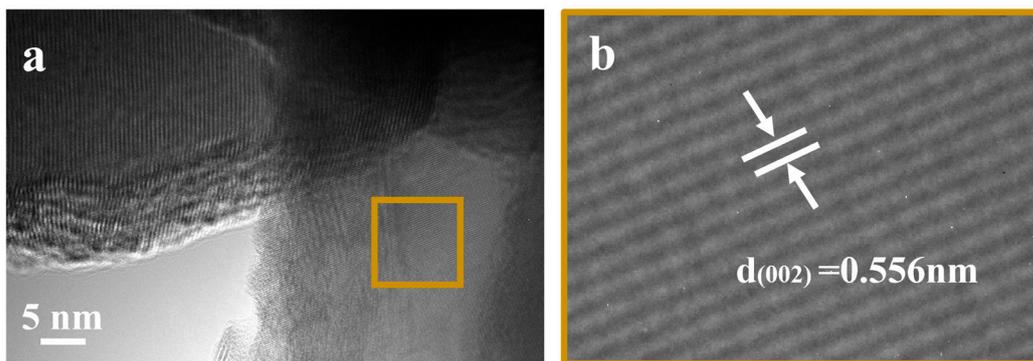


Fig. S9 (a) The TEM images of P2-Na<sub>0.67</sub>MN. (b) The TEM images of P2-Na<sub>0.67</sub>MN.

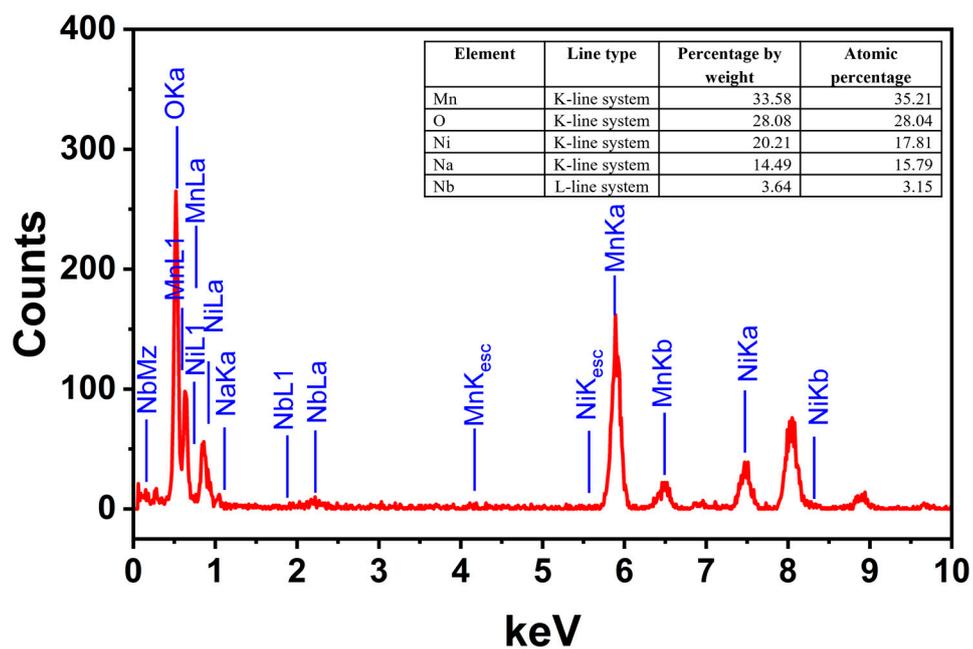


Fig. S10 The spectrum diagram of total distribution diagram for HADDF-EDS of P2-

Na<sub>0.67</sub>MNNb.

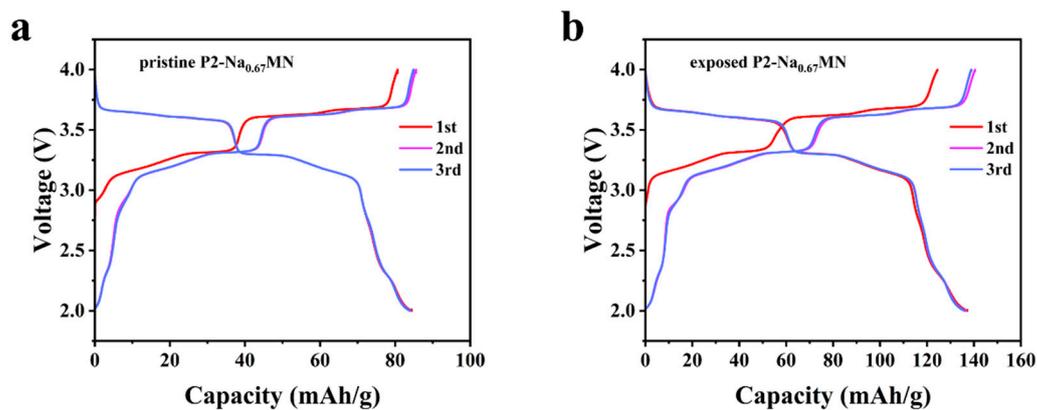


Fig. S11 Galvanostatic charge/discharge voltage profiles of (a) the pristine P2-Na<sub>0.67</sub>MN and (b) exposed P2-Na<sub>0.67</sub>MN at the first three cycles at 0.2 C in the voltage range of 2-4 V (exposed P2-Na<sub>0.67</sub>MN refers to exposing pristine P2-Na<sub>0.67</sub>MN to RH93% humid environment for 20 days).

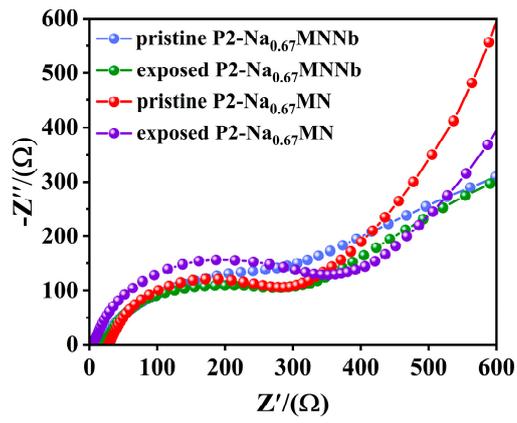


Fig. S12 Nyquist plot of the coin cells that based on pristine P2-Na<sub>0.67</sub>MN, exposed P2-Na<sub>0.67</sub>MN, pristine P2-Na<sub>0.67</sub>MNNb and exposed P2-Na<sub>0.67</sub>MNNb electrode materials.

Table. S1 The ICP date of P2-Na<sub>0.67</sub>MN and P2-Na<sub>0.67</sub>MNNb cathode materials.

ICP results (mol mL <sup>-1</sup> )	Na	Mn	Ni	Nb
Na <sub>0.67</sub> Mn <sub>0.67</sub> Ni <sub>0.33</sub> O <sub>2</sub>	0.67	0.67	0.33	
Na <sub>0.67</sub> Mn <sub>0.67</sub> Ni <sub>0.33</sub> Nb <sub>0.03</sub> O <sub>2</sub>	0.67	0.67	0.30	0.03

Table. S2 Crystallographic parameters of P2-Na<sub>0.67</sub>MNNb refined by the Rietveld method.

Atoms	Wyckoff position	X	Y	Z	Occupancy
Na1	2b	0	0	0.25000	0.30
Na2	2d	0.66670	0.333330	0.25000	0.37
Ni	2a	0	0	0	0.31
Mn	2a	0	0	0	0.67
O	4f	0.66670	0.33330	0.07950	1
Nb	2a	0	0	0	0.03

Space group: P6<sub>3</sub>/mmc (space group no. 194), a=2.88749 Å, c=11.16065 Å, Rwp=4.32%.

Table. S3 Crystallographic parameters of P2-Na<sub>0.67</sub>MN refined by the Rietveld method.

Atoms	Wyckoff position	X	Y	Z	Occupancy
Na1	2b	0	0	0.25000	0.36
Na2	2d	0.66670	0.33330	0.25000	0.31
Ni	2a	0	0	0	0.33
Mn	2a	0	0	0	0.67
O	4f	0.66670	0.33330	0.05180	1

Space group: P6<sub>3</sub>/mmc (space group no. 194), a=2.88504 Å, c=11.15540 Å, Rwp=5.62%.