

Supporting Information

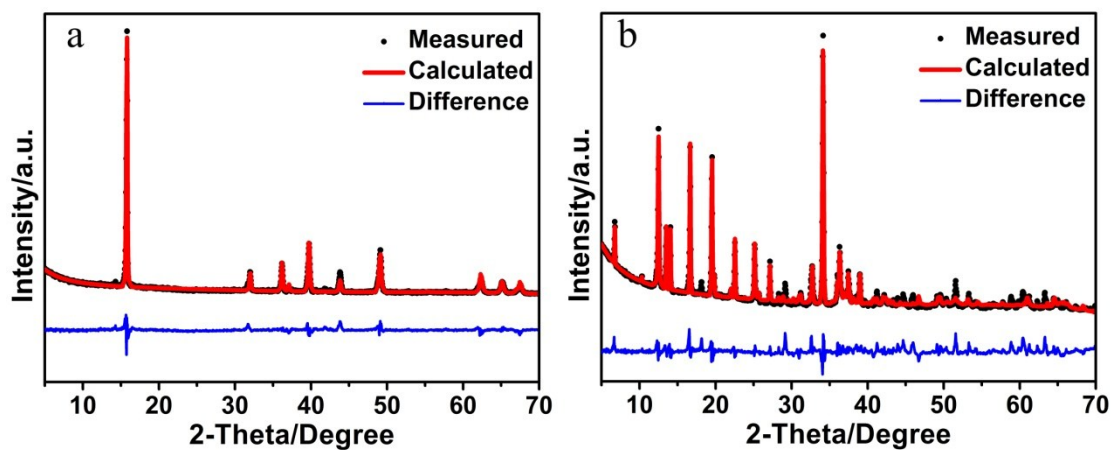


Figure S1 Rietveld refinements of (a) pure P2 -type and (b) T-type material.

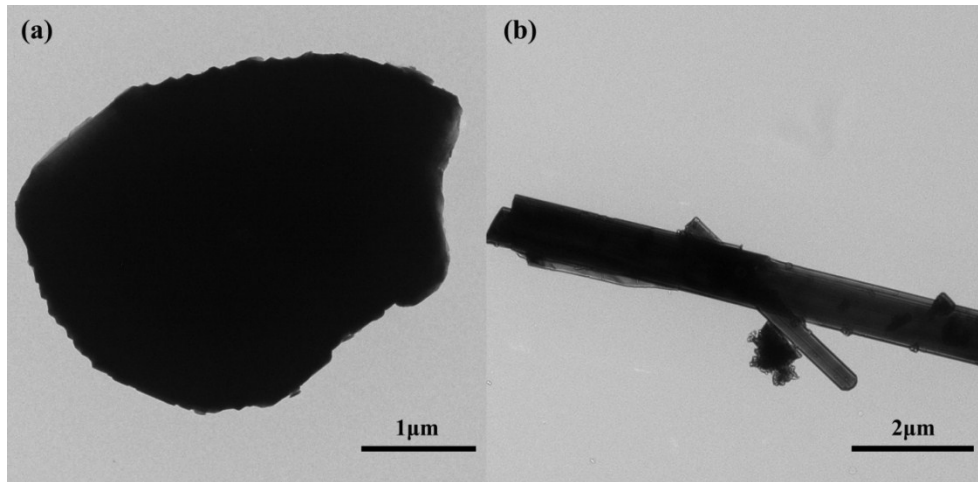


Figure S2 TEM image of (a) P2 -type and (b) T-type material.

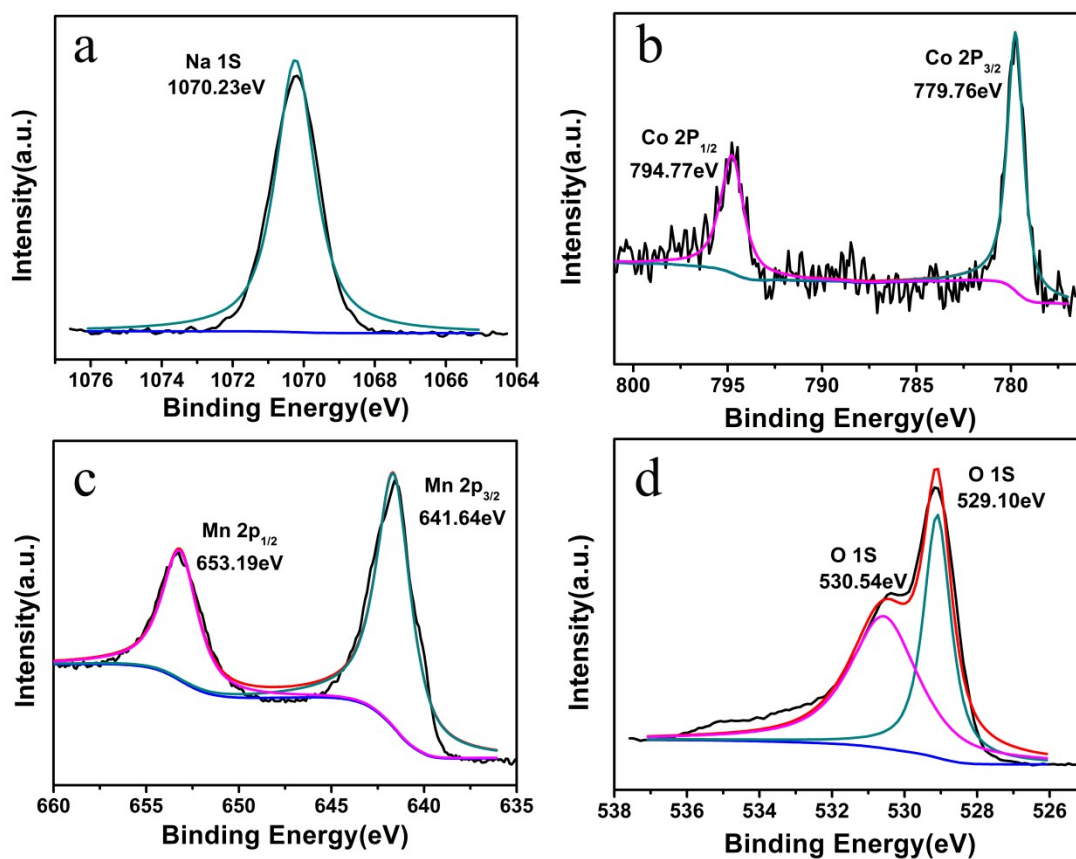


Figure S3. XPS analysis of the P2+T material: (a) XPS spectrum of Na 1s; (b)Co 2p; (c) Mn 2p; and (d) O 1s.

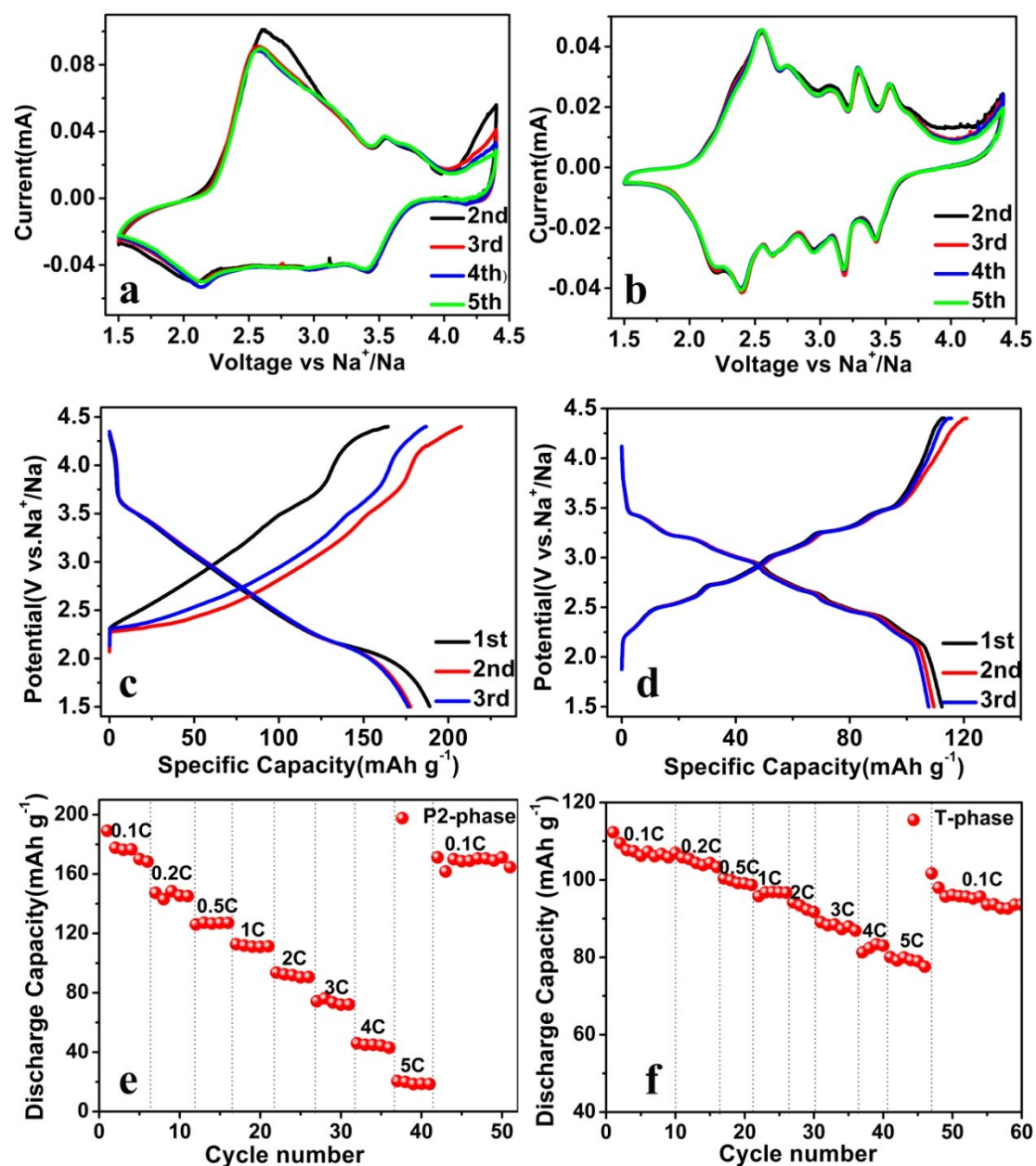


Figure S4.(a-b) Cyclic voltammogram of the P2- and T-phase electrode at a scan rate of 0.1 mVs⁻¹; (c-d) Typical charge/discharge voltage profiles of the P2- and T-phase electrode at a current rate of 0.04 C; (e-f) Rate performance of P2- and T-phase, electrode at various current rates.

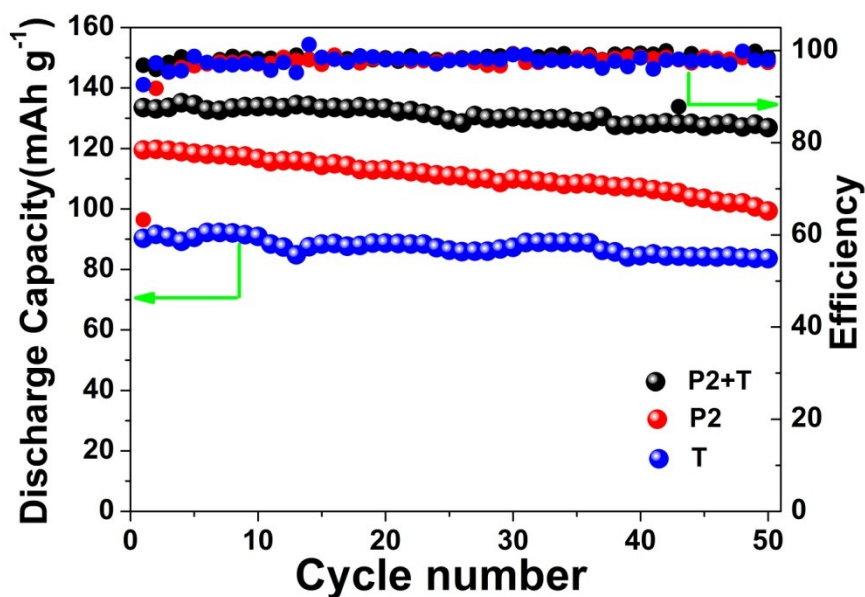


Figure S5. Specific discharge capacities and coulomb efficiencies versus cycle number for 50 cycles. Potential range: 2–4.4V.

In a narrow voltage of 2-4.4V, the P2+T phase material exhibits specific discharge capacities of 133.5 mAh g⁻¹ and 126.9 mAh g⁻¹ at the 1st and 50th cycle, resulting in 95% capacity retention (2nd–50th cycles). This performance is better than the other two materials. The P2 phase material shows a specific discharge capacity of 99.3 mAh g⁻¹ after 50th cycles, with a cycle efficiency of 83.1%.

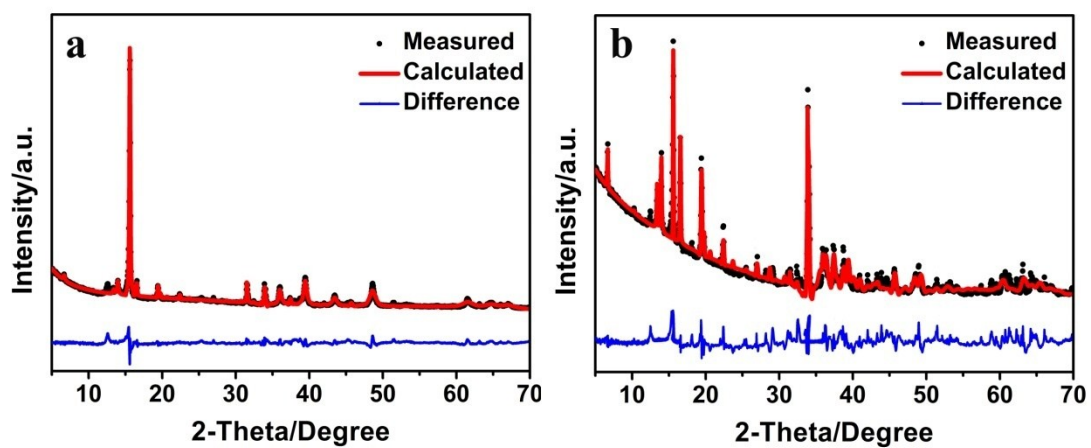


Figure S6 Rietveld refinements of P2+T-type $\text{Na}_x\text{Co}_{0.1}\text{Mn}_{0.9}\text{O}_2$ with varied P2: T ratio of 50%:50% (a), and 20%:80% (b).

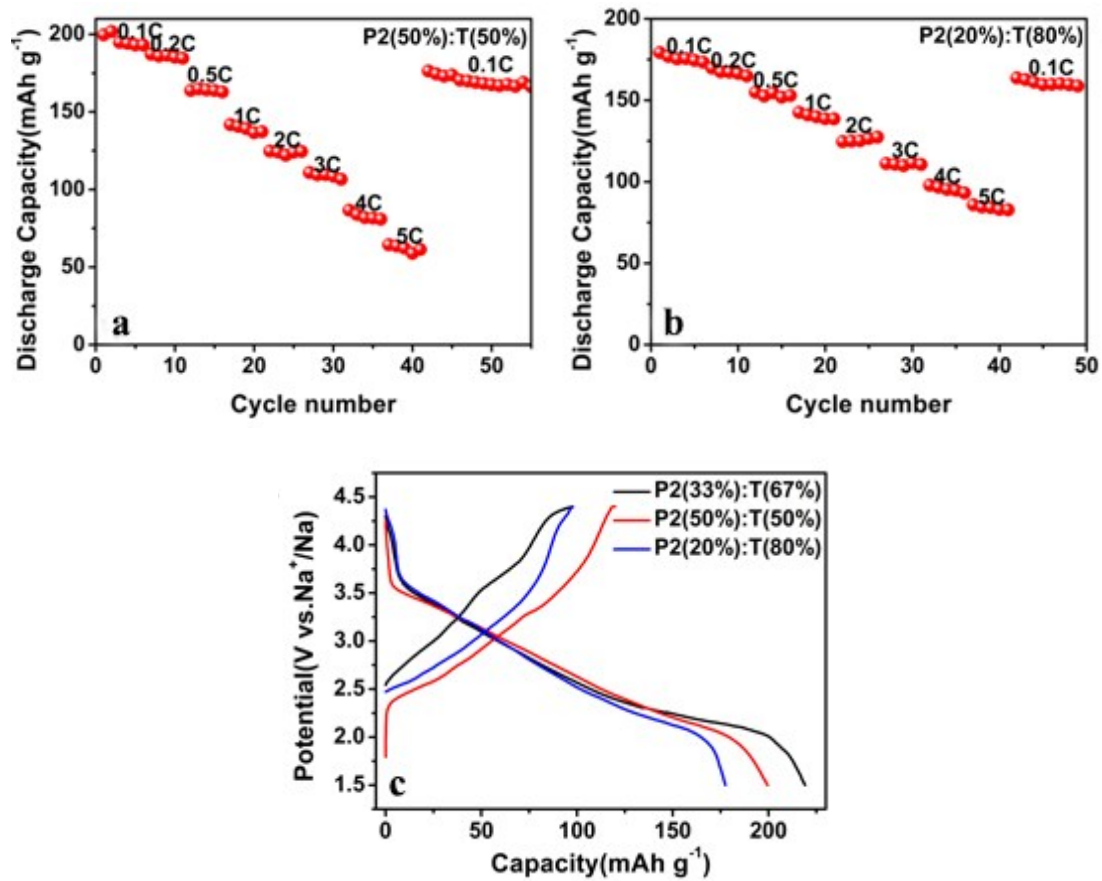


Figure S7 The C-rate performance of $\text{Na}_x\text{Co}_{0.1}\text{Mn}_{0.9}\text{O}_2$ with the ratio between P2 and T of 50%:50% (a), and 20%:80% (b); (c) Charge/discharge curves of $\text{Na}_x\text{Co}_{0.1}\text{Mn}_{0.9}\text{O}_2$ with varied P2:T ratio.

Table S1 Lattice parameters of the P2 compound sample after Rietveld refinement.

Phase	P2 (JCPDS: 27-0751)	
Space Group	P63/mmc	
	a (Å)	2.8275
	b (Å)	2.8275
	c (Å)	11.0295(9)
	Volume (Å ³)	76.3651
Cell parameters	α (°)	90
	β (°)	90
	γ (°)	120
Agreement factors	Rwp (%)	6.88
	Rp (%)	8.56

Table S2 Lattice parameters of the T compound sample after Rietveld refinement.

Phase	T (JCPDS: 27-0751)	
Space Group	Pbam	
	a (Å)	9.1133(9)
	b (Å)	21.3177(3)
	c (Å)	2.8897(9)
	Volume (Å ³)	693.0517
Cell parameters	α (°)	90
	β (°)	90
	γ (°)	90
Agreement factors	Rwp (%)	5.97
	Rp (%)	9.8

Table S3 Lattice parameters of the P2+T compound sample after Rietveld refinement.

Phase	P2 (JCPDS: 27-0751)		T (JCPDS: 27-0750)	
Phase ratio	32.39%		63.61%	
space group	P63/mmc		Pbam	
	a (Å)	2.8791(1)	a (Å)	9.1125(1)
	b (Å)	2.8791(1)	b (Å)	26.4026(8)
	c (Å)	11.3380(5)	c (Å)	2.8235(1)
	Volume (Å ³)	81.396(4)	Volume (Å ³)	679.4678(5)
Cell parameters	α (°)	90	α (°)	90
	β (°)	90	β (°)	90
	γ (°)	120	γ (°)	90
Agreement factors	Rwp(%)	5.49	Rwp(%)	5.49
	Rp(%)	7.69	Rp(%)	7.69

Table S4 ICP-AES results of the P2+T compound sample

Theoretical chemical formula	Measured atomic ratio		
	Na	Mn	Co
Na_xMn_{0.9}Co_{0.1}O₂	0.48	0.89	0.09

Table S5 Electrochemical performance comparison between the previously reported $\text{Na}_x\text{Mn}_y\text{M}_{1-y}\text{O}_2$ cathodes and P2+T-type $\text{Na}_x\text{Co}_{0.1}\text{Mn}_{0.9}\text{O}_2$ composite cathode

Cathode materials	Rate capability	Cycling stability	First discharge capacity	Ref.
P2-Type	80 mAh g ⁻¹	~120 mAh g ⁻¹ after	180 mAh g ⁻¹ (12	21
$\text{Na}_{0.78}\text{Ni}_{0.23}\text{Mn}_{0.69}\text{O}_2$	(605 mA g ⁻¹)	20 cycles	mA g ⁻¹)	
		(12 mA g ⁻¹)		
$\text{Na}_{0.67}\text{Mn}_{0.67}\text{Ni}_{0.33}\text{O}_2$	35 mAh g ⁻¹	~120 mAh g ⁻¹	150 mAh g ⁻¹ (17	23
	(340 mA g ⁻¹)	after 100 cycles	(17 mA g ⁻¹)	
		(17 mA g ⁻¹)		
P2-type	58 mAh g ⁻¹	~150 mAh g ⁻¹ after	190 mAh g ⁻¹ (13	25
$\text{Na}_x\text{Fe}_{1/2}\text{Mn}_{1/2}\text{O}_2$	(1040 mA g ⁻¹)	30 cycles	mA g ⁻¹)	
		(13 mA g ⁻¹)		
P2/O3 composite	69 mAh g ⁻¹	~125 mAh g ⁻¹	200 mAh g ⁻¹ (10	26
$\text{Na}_{0.66}\text{Li}_{0.18}\text{Mn}_{0.71}\text{Ni}_{0.2}$	(500 mA g ⁻¹)	after 150 cycles	mA g ⁻¹)	
$\text{Co}_{0.08}\text{O}_{2+}$		(50 mA g ⁻¹)		
P2/O3	Na ₁₋ 78 mAh g ⁻¹	~134 mAh g ⁻¹	140 mAh g ⁻¹ (15	27
$\text{Li}_x\text{Ni}_{0.5}\text{Mn}_{0.5}\text{O}_{2+d}$	(150 mA g ⁻¹)	after 20 cycles	mA g ⁻¹)	
		(15 mA g ⁻¹)		
P2/O3	110 mAh g ⁻¹	~130 mAh g ⁻¹	154 mAh g ⁻¹ (18	30
$\text{Na}_x\text{Mn}_y\text{Ni}_z\text{Fe}_{0.1}\text{Mg}_{0.1}$	(900 mA g ⁻¹)	after 50 cycles	mA g ⁻¹)	
O_2		(180 mA g ⁻¹)		
P2- $\text{Na}_x\text{Co}_y\text{Mn}_{1-y}\text{O}_2$ (y	120 mAh g ⁻¹	~117 mAh g ⁻¹	188 mAh g ⁻¹ (20	39
= 0, 0.1)	(400 mA g ⁻¹)	after 150 cycles	mA g ⁻¹)	
		(50 mA g ⁻¹)		
P2+T	117 mAh g ⁻¹	~145 mAh g ⁻¹	219 mAh g ⁻¹ (18	This
$\text{Na}_x\text{Mn}_{0.9}\text{Co}_{0.1}\text{O}_2$	(900 mA g ⁻¹)	after 50 cycles	mA g ⁻¹)	work
		(180 mA g ⁻¹)		

Table S6 Na⁺ Diffusion Coefficient(D) of P2, T and P2+T materials Calculated from the CV and EIS.

		P2	T	P2+T
D for CV	Average slope	0.00568	0.01436	0.01671
	$D(\times 10^{-12}cm^2/s)$	2.156	76.55	434.24
D for EIS	Slope	13.07	32.83	8.23
	$D(\times 10^{-12}cm^2/s)$	1.827	41.67	127.07