

Supporting Information

High performance nano-sized $\text{LiMn}_{1-x}\text{Fe}_x\text{PO}_4$ cathode material for advanced lithium-ion batteries

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Table S1 Refined cell parameters of $\text{LiMn}_{1-x}\text{Fe}_x\text{PO}_4/\text{C}$ ($0 \leq x \leq 1$) materials

Iron proportion	a(Å)	b(Å)	c(Å)	Volume (Å ³)	Carbon content (%)
x = 0	6.10553	10.45085	4.74697	302.89	5.34
x = 0.05	6.10227	10.45149	4.74786	302.81	5.64
x = 0.10	6.09713	10.43212	4.74042	301.52	5.50
x = 0.15	6.08934	10.43192	4.73890	301.03	6.48
x = 0.20	6.09331	10.43535	4.74054	301.43	6.08
x = 0.25	6.08084	10.43030	4.74531	300.97	5.86
x = 0.30	6.07246	10.41302	4.73188	299.12	7.35
x = 0.40	6.06129	10.40572	4.72702	298.14	6.55
x = 0.50	6.05326	10.39118	4.71638	296.66	6.68
x = 1	5.99712	10.32512	4.69412	290.66	6.04

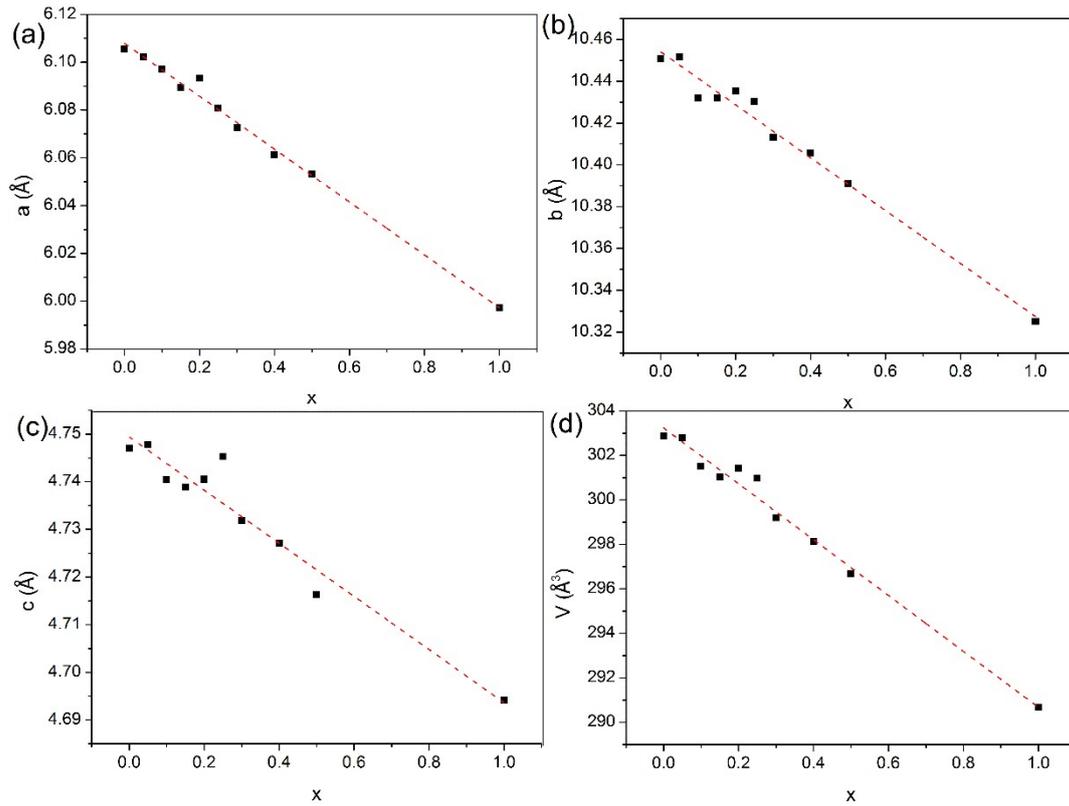


Figure S1. The crystal lattice parameters comparison of $\text{LiMn}_{1-x}\text{Fe}_x\text{PO}_4/\text{C}$ ($0 \leq x \leq 1$) materials.

Table S2. Atomic position of $\text{LiMn}_{0.7}\text{Mn}_{0.3}\text{PO}_4$ determined by Rietveld structure refinement.

Label	x	y	z	Symmetry multiplicity
Li	0	0	0	4
Mn	0.28243	0.25	0.97211	4
Fe	0.27907	0.25	0.96776	4
P	0.09389	0.25	0.41314	4
O1	0.095852	0.25	0.73001	4
O2	0.45568	0.25	0.21398	4
O3	0.16138	0.05160	0.27820	8

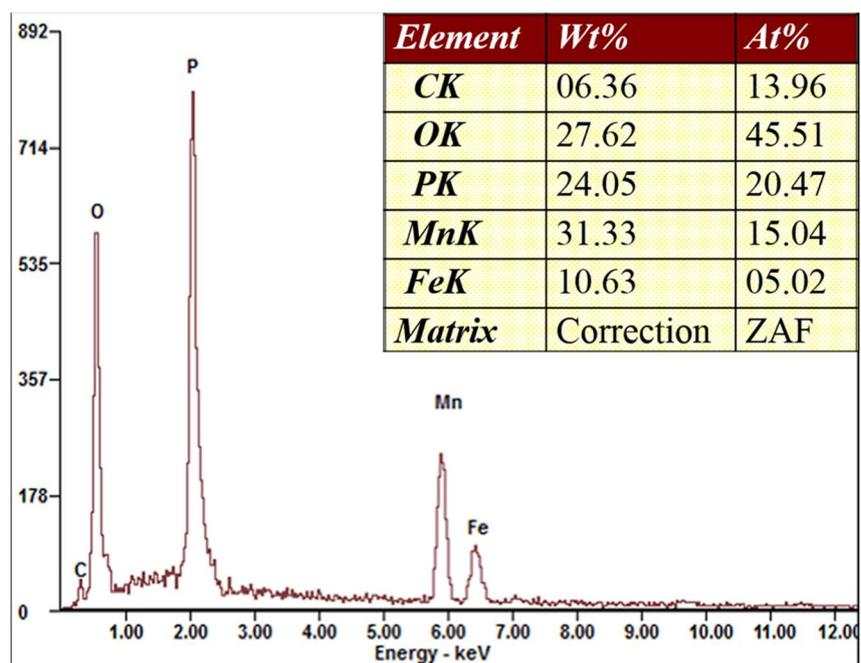


Figure S2. The SEM-EDX spectrum of LiMn_{0.75}Fe_{0.25}PO₄/C material.

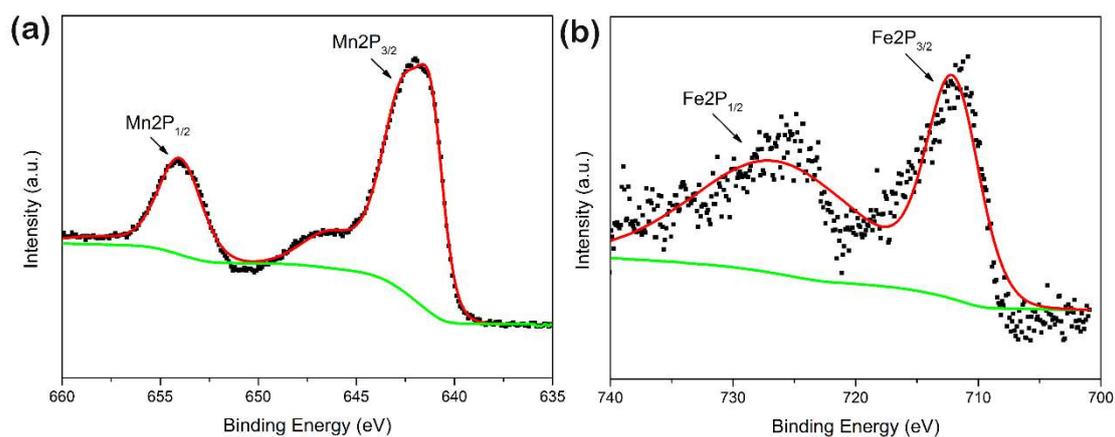


Figure S3. XPS spectra of Mn and Fe element in LiMn_{0.7}Fe_{0.3}PO₄/C material.

Table S3 Electronic conductivities of Selected LiMn_{1-x}Fe_xPO₄/C materials

Material	Conductivity (S cm ⁻¹)
LiMn _{0.7} Fe _{0.3} PO ₄ /C	5.8×10^{-5}
LiMn _{0.75} Fe _{0.25} PO ₄ /C	3.1×10^{-5}
LiMnPO ₄ /C	2.3×10^{-5}
LiFePO ₄ /C	6.8×10^{-5}