Electronic Supplementary Material (ESI) for New Journal of Chemistry.

Table S3. The pseudo-first-order kinetic model and the pseudo-second-order kinetic model adsorption kinetic parameters of Li^+ on NFs@LMO-6:6

Absorbents	Pseudo-first-order kinetic model			Pseudo-second-order kinetic model		
	Q_1(mg·g^{-1})	k_1(min^{-1})	R^2	Q_2(mg·g^{-1})	k_2(g·mg^{-1}·min^{-1})	R^2
NF@LIS-6:6	24.26	0.0536	0.9087	25.00	0.0130	0.9977