



Supplementary Materials

# Electrochemical Properties of an Sn-Doped LATP Ceramic Electrolyte and Its Derived Sandwich-Structured Composite Solid Electrolyte

Aihong Xu <sup>1,2</sup>, Ruoming Wang <sup>1,2</sup>, Mengqin Yao <sup>1,2</sup>, Jianxin Cao <sup>1,2</sup>, Mengjun Li <sup>1,2</sup>, Chunliang Yang <sup>1,2</sup>, Fei Liu <sup>1,2,\*</sup> and Jun Ma <sup>1,2,\*</sup>

<sup>1</sup> Department of Chemical Engineering, School of Chemistry and Chemical Engineering, Guizhou University, Guiyang 550025, China; 15851959055@163.com (A.X.); w1285459785@126.com (R.W.); mqyao@gzu.edu.cn (M.Y.); jxcao@gzu.edu.cn (J.C.); x1285459785@126.com (M.L.); clyang@gzu.edu.cn (C.Y.)

<sup>2</sup> Guizhou Key Laboratory for Green Chemical and Clean Energy Technology, Guiyang 550025, China

\* Correspondence: ce.feiliu@gzu.edu.cn (F.L.); jma3@gzu.edu.cn (J.M.)

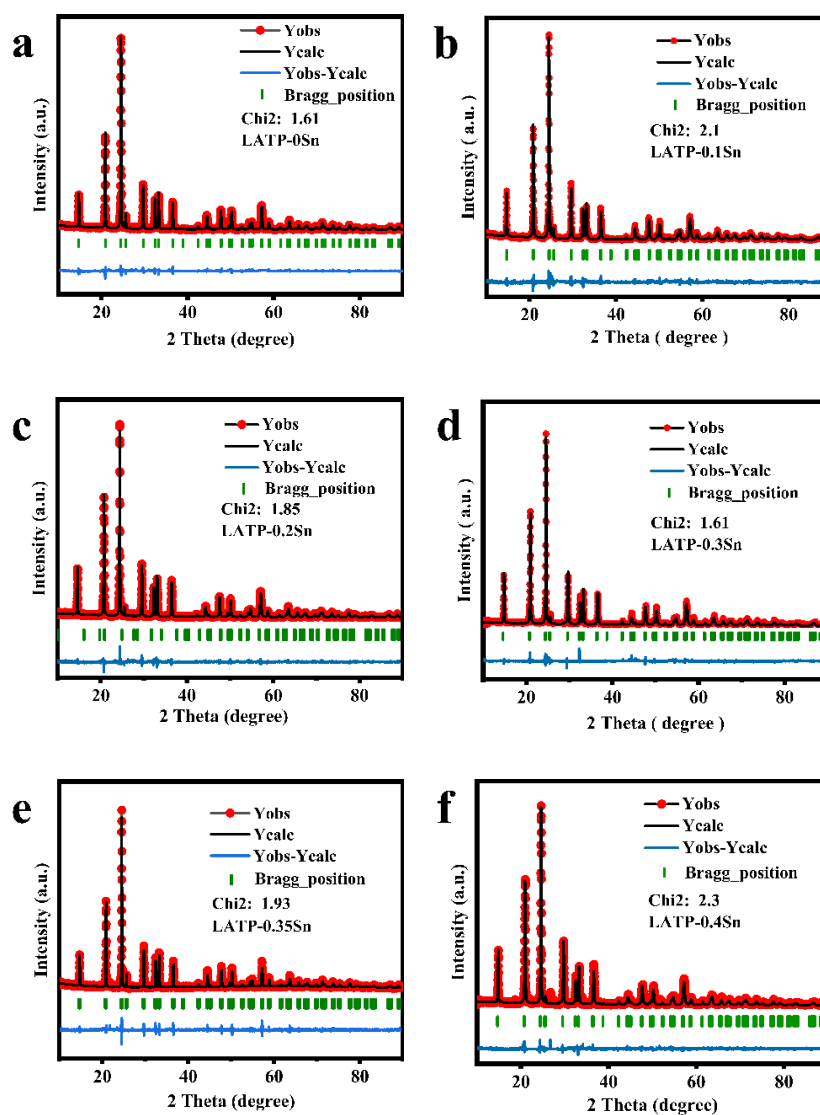


Figure S1. Rietveld refinement of XRD patterns of LATP-xSn ( $x = 0-0.4$ ) solid electrolytes.

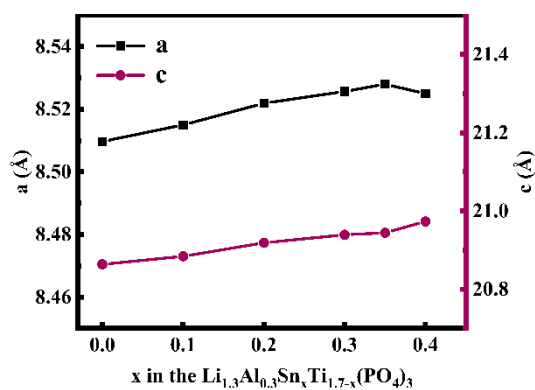
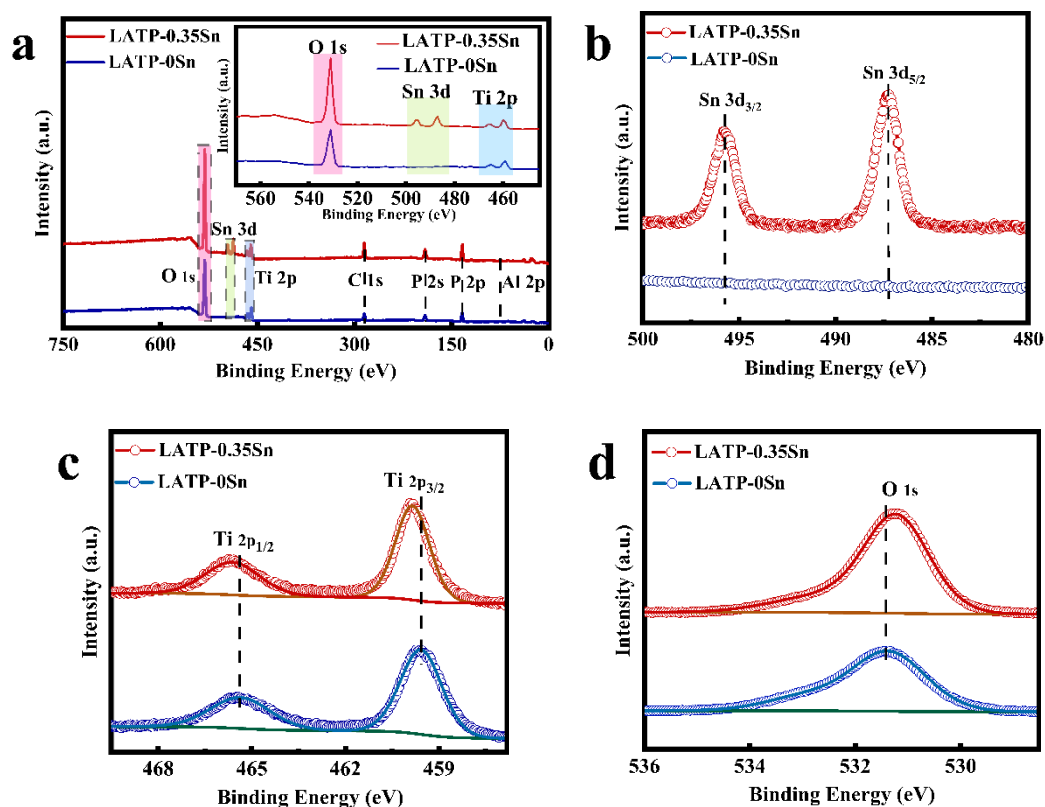
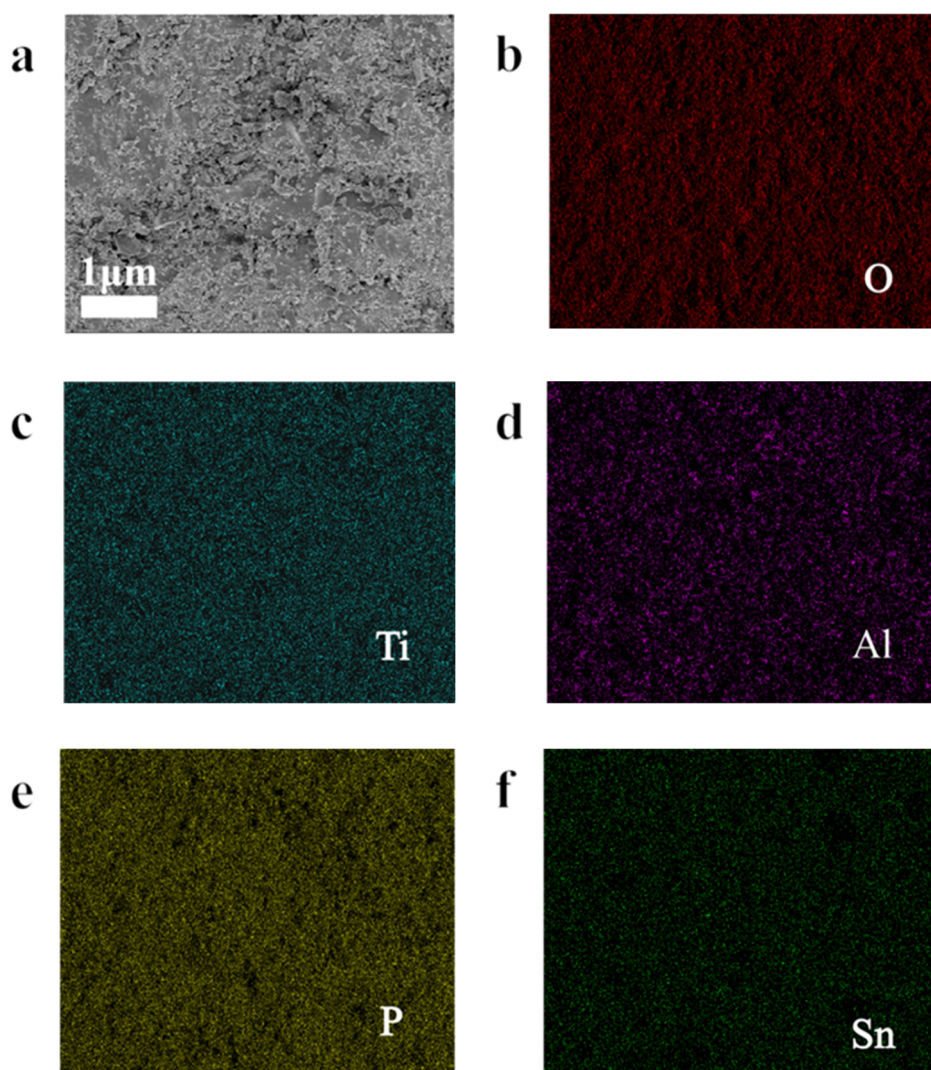


Figure S2. Lattice parameters of LATP-xSn ( $x = 0-0.4$ ) solid electrolytes.

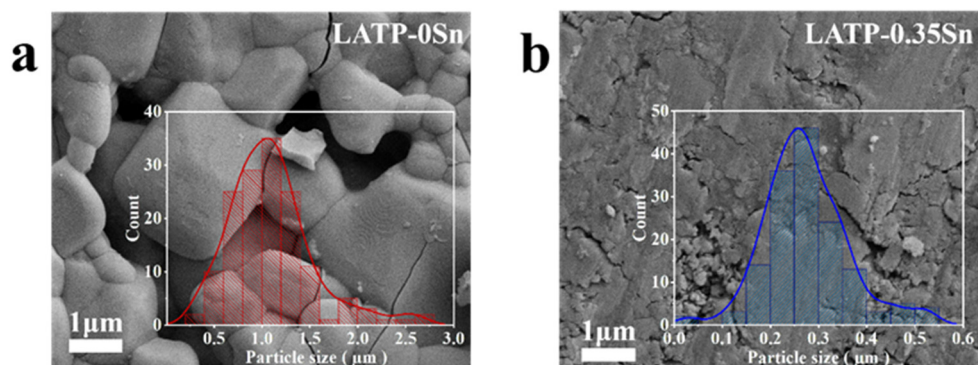
**Table S1.** The deviation degrees of  $\text{Sn}^{4+}$  at  $\text{P}^{5+}$  site,  $\text{Al}^{3+}$  site and  $\text{Ti}^{4+}$  site respectively under different coordination environments.

Ions	$\text{Sn}^{4+}$	$\text{P}^{5+}$	$\text{Al}^{3+}$	$\text{Ti}^{4+}$
Crystal ionic radii with tetrahedral coordination	69 pm	31 pm	/	/
Crystal ionic radii with octahedral coordination	83 pm	/	67.5 pm	74.5 pm
Electronegativity with tetrahedral coordination	1.95	3.003	/	/
Electronegativity with octahedral coordination	1.706	/	1.513	1.730
Deviation degree	/	1.58	0.357	0.128

**Figure S3.** XPS spectra of LATP-0Sn and LATP-0.35Sn solid electrolytes (a) survey spectra; (b) Sn 3d spectra; (c) Ti 2p spectra; (d) O 1s spectra.



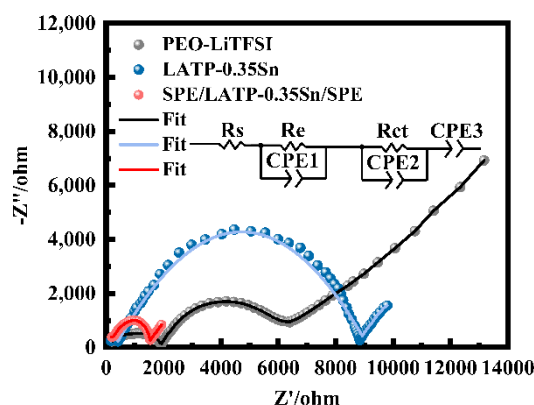
**Figure S4.** (a) SEM image of LATP-0.35Sn solid electrolyte and (b–f) the corresponding EDS elemental maps.



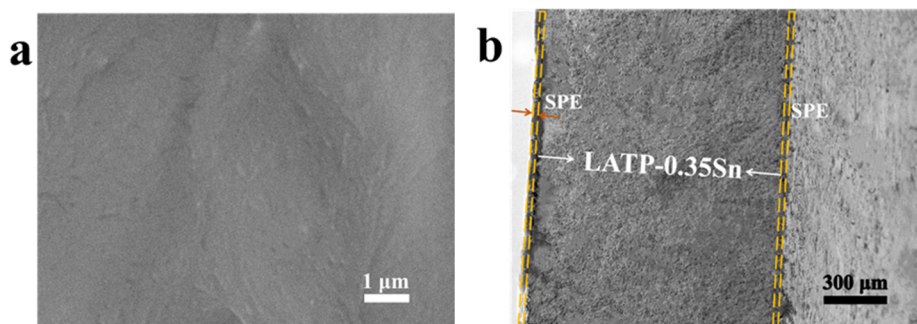
**Figure S5.** SEM images and grain size distribution curves of (a) LATP-0Sn solid electrolyte and (b) LATP-0.35Sn solid electrolyte.

**Table S2.** Performance comparison of LATP-xSn obtained with other typical solid electrolytes.

Composition	Electrolyte-type	Preparation method	Conductivity (mS/cm)	Reference
$\text{Li}_7\text{La}_3\text{Zr}_2\text{O}_{12}$	Garnet	solid-state reaction	0.38	58
$\text{Li}_{3/8}\text{Sr}_{7/16}\text{Hf}_{1/4}\text{Ta}_{3/4}\text{O}_3$	Perovskite	solid-state reaction	0.38	59
$\text{Li}_{10.42}\text{Ge}_{1.5}\text{P}_{1.5}\text{C}_{10.08}\text{O}_{11.92}$	LISICON	melt-casting	0.037	60
$\text{Li}_3\text{PO}_4$	LiPON	magnetron sputtering	0.001	61
$\beta\text{-Li}_3\text{PS}_4$	Sulfides	liquid-phase reaction	0.27	62
$\text{Li}_3\text{OCl}$	Anti-perovskite	PLD	0.2	63

**Figure S6.** EIS curves of LATP-0.35Sn and SPE/LATP-0.35Sn/SPE solid electrolytes at 25 °C.**Table S3.** Electrolyte resistance ( $R_e$ ), charge transfer resistance ( $R_{ct}$ ), total resistance ( $R_t$ ) and total ionic conductivities ( $\sigma_t$ ) of PEO-LiTFSI, LATP-0.35Sn and SPE/LATP-0.35Sn/SPE solid electrolytes at 25 °C.

Electrolytes	$R_e$ ( $\Omega$ )	$R_{ct}$ ( $\Omega$ )	$R_t$ ( $\Omega$ )	$\sigma_t$ (S/cm)
PEO-LiTFSI	1739	4408	6344	$1.0 \times 10^{-6}$
LATP-0.35Sn	413	8514	8927	$1.0 \times 10^{-5}$
SPE/LATP-0.35Sn/SPE	239	1344	1583	$5.9 \times 10^{-5}$

**Figure S7.** SEM images of SPE/LATP-0.35Sn/SPE solid electrolytes (a) surface and (b) cross-section.