Supporting Information (SI)

## Pentafluorobenzene boronic acid with strong Lewis acidity for

## the modification of PEO-based polymer electrolytes

Liwei Su,\*a Yijuan Zhu, a Xingyi Zhan, a Kang Yu, a Ke Gu, a Hao Wu, a Lianbang

Wang\*a, Yuanhao Wang,b\* Xiaoxiang Wang<sup>c</sup>

<sup>a</sup> State Key Laboratory Breeding Base of Green Chemistry-Synthesis Technology, College of Chemical Engineering, Zhejiang University of Technology, Hangzhou 310014, P. R. China
<sup>b</sup> Hoffmann Institute of Advanced Materials, Shenzhen Polytechnic, Shenzhen 518055, P. R. China
<sup>c</sup> Anteotech, Queensland 4113, Australia
\*Corresponding author.
E-mail: suliwei@zjut.edu.cn (L. Su); wanglb99@zjut.edu.cn (L. Wang);





Fig. S1. Diagram of the preparation process of PEO-based SSEs

Electrolytes	T <sub>g</sub> (°C)	$T_m(^{\circ}C)$	$\Delta H_m$ (J g <sup>-1</sup> )	<i>χ<sub>c</sub></i> (%)
Pristine	-50.44	48.60	46.86	23.08
0.5%PFBBA	-57.41	47.76	44.53	21.93
1.0%PFBBA	-59.76	46.55	39.99	19.70
1.5%PFBBA	-58.90	46.96	33.39	16.45

Table S1. Thermodynamic values and crystallinity of electrolytes

Table S2. Ionic conductivity (S cm<sup>-1</sup>) of the four electrolytes at different temperatures

Electrolytes	30°C	40°C	50°C	60°C
Pristine	8.01×10 <sup>-6</sup>	1.99×10 <sup>-5</sup>	5.26×10-5	1.92×10 <sup>-4</sup>
0.5%PFBBA	2.03×10 <sup>-5</sup>	4.15×10 <sup>-5</sup>	8.85×10 <sup>-5</sup>	2.37×10-4
1.0%PFBBA	3.09×10 <sup>-5</sup>	6.68×10 <sup>-5</sup>	1.56×10-4	4.58×10 <sup>-4</sup>
1.5%PFBBA	1.20×10 <sup>-5</sup>	2.46×10 <sup>-5</sup>	5.32×10 <sup>-5</sup>	1.28×10 <sup>-4</sup>

Table S3. FT-IR characteristic peaks and the vibration types of functional groups

Wavenumber/cm <sup>-1</sup>	Chemical bond	Vibration types
1400-1300	-CH <sub>2</sub>	rocking vibration
990-910	-CH <sub>2</sub>	stretching vibration
1668, 1671	-OLi	/
1358, 962	B-OH and -O-	hydrogen bond interaction
1348	B-O	/
1096, 1056	C-O-C	symmetric tensile vibration



Fig. S2. Arrhenius diagram of electrolytes: (a) Pristine and (b) 1.0% PFBBA



Fig. S3. DC polarization curves and impedance of symmetric Li//Li battery before and after polarization at 30°C: (a) Pristine, (b) 0.5% PFBBA, (c) 1.0% PFBBA, and (d) 1.5% PFBBA



Fig. S4. DC polarization curves and impedance of symmetric Li//Li battery before and after polarization at 60°C: (a) Pristine, (b) 0.5% PFBBA, (c) 1.0% PFBBA, and (d) 1.5% PFBBA



Fig. S5. SEM images of the electrolyte surface exfoliated from the (a, c) Li/Pristine/Li and (b, d) Li/1.0%PFBBA/Li batteries after cycling for 100 h at the current density of 0.05 mA cm<sup>-2</sup>



Fig. S6. The equivalent circuit used to fit the EIS plots in Fig. 5, S3, and S4.

Cycles	Pristine $R_{int}(\Omega)$	1.0% PFBBA R <sub>int</sub> (Ω)
0	1949	1590
1	2821	1712
5	5877	1861
10	3174	1797

Table S4. The interface impedance of Li//Li batteries before and after polarization



Fig. S7. The charge/discharge curves at different rates of the Li/Pristine/LFP and Li/1.0%PFBBA/LFP full batteries at 30°C



Fig. S8. Cycling performance and Coulombic efficiency of the Li//LiCoO<sub>2</sub> batteries: (a) the Pristine electrolyte and (b) the 1.0%PFBBA electrolyte