

Supporting information for

**High-melting-point crystals of Poly(l-lactic acid)(PLLA): the Most Efficient
Nucleating Agent to Enhance the Crystallization of PLLA**

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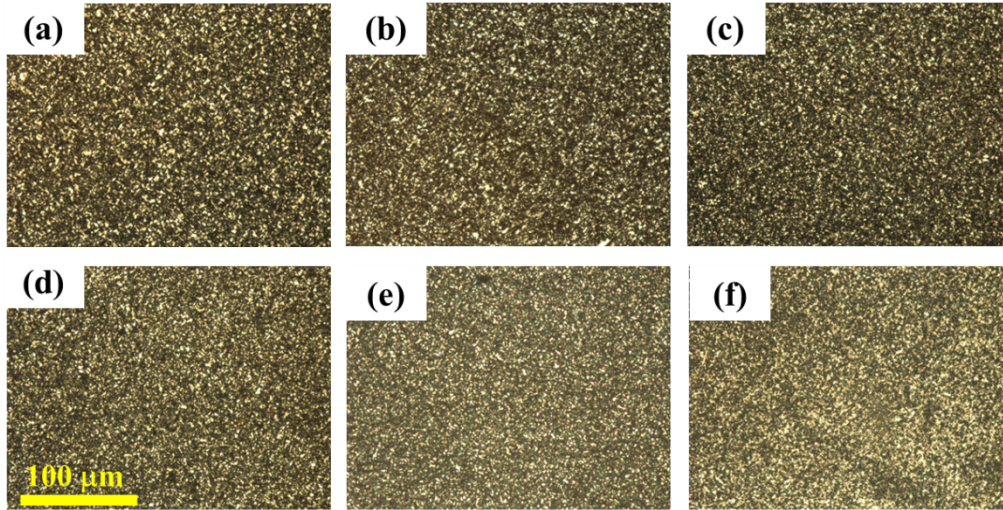


Fig. S1. Selected POM micrographs after completely crystallization at T_c of 140 °C for *h*PLLA-0.1 (a) *h*PLLA-0.2 (b) *h*PLLA-0.5 (c) *h*PLLA-1.0 (d) *h*PLLA-2.0 (e) and *h*PLLA-5.0 (f). The scale bar in the bottom micrograph of panel (d) represents 100 μm and applies to all the micrographs.

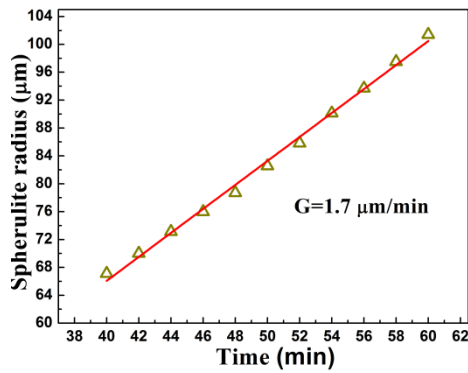


Fig. S2. The spherulite radius as a function of time for neat *l*PLLA at $T_c = 140$ °C, the sample was first melting at 180 °C for 3 min and then cooled to 140 °C at 100 °C/min

Based on the DSC data, the isothermal crystallization kinetics of PLLA in the blends crystallized at 140 °C were analyzed by the Avrami equation as follows:

$$1 - X_t = \exp[-k(T)t^n] \quad (1)$$

where n is the Avrami index and $K(T)$ is the overall rate constant. The linear form of

eq 1 can be stated as:

$$\ln[-\ln(1 - X_t)] = \ln k(T) + n \ln t \quad (2)$$

$k(T)$ and n can be estimated from the linear fitting of $\ln[-\ln(1 - X_t)]$ vs $\ln t$. To ensure the accuracy of the Avrami analysis, the data in a limited conversion range (3 – 30%) was employed for fitting. The obtained data were shown in Fig.S3 and Table.S1.

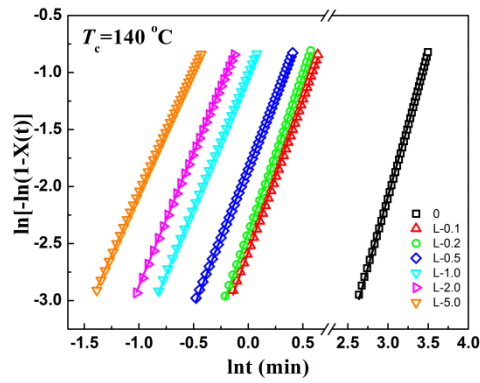


Fig. S3. Avrami plots of $\ln[-\ln(1 - X_t)]$ versus $\ln t$ for neat PLLA and the blends at T_c of 140 °C

Table S1 Values of Avrami parameters for neat PLLA and its blends at T_c of 140 °C

Sample Code	n	$\ln k(T)$ (min^{-n})
<i>l</i> PLLA	2.8	-9.57
<i>h</i> PLLA-0.1	2.7	-2.57
<i>h</i> PLLA -0.2	2.8	-2.40
<i>h</i> PLLA -0.5	2.4	-1.83
<i>h</i> PLLA -1.0	2.3	-1.02
<i>h</i> PLLA -2.0	2.3	-0.54
<i>h</i> PLLA -5.0	2.2	0.96