

Supporting Information

The calculation about the loading of the Ag NPs

$$\frac{\Delta m}{m_0} = \frac{m - m_0}{m_0} = \frac{0.08\% - 0.03\%}{0.03\%} = 1.66$$

Where, m_0 is the loading of the Ag NPs onto the Ag NPs/PVDF composite, i.e. the atomic concentration of Ag in the Ag NPs/PVDF composite by XPS; m is the loading of the Ag NPs onto the Ag NPs/MWCNTs/PVDF composite, i.e. the atomic concentration of Ag in the Ag NPs/MWCNTs/PVDF composite by XPS.

Table S1. D- and G-bands shifts and Raman I_D/I_G intensity ratios of purified MWCNTs, MWCNTs/PVDF and Ag NPs/MWCNTs/PVDF composites.

| Sample | D-band [cm ⁻¹] | G-band [cm ⁻¹] | I_D/I_G |
|------------------|-------------------------------|-------------------------------|-----------|
| purified MWNTs | 1352 | 1585 | 0.88 |
| MWNTs/PVDF | 1353 | 1586 | 1.03 |
| AgNPs/MWNTs/PVDF | 1353 | 1586 | 1.31 |

Table S2. DSC parameters of the PVDF, Ag NPs/PVDF, MWCNTs/PVDF and Ag NPs/MWCNTs/PVDF composites

| Sample | ΔH_m [J/g] ^{a)} | ΔH_c [J/g] ^{b)} | T_m [°C] ^{c)} | T_c [°C] ^{d)} | X_c [%] ^{e)} |
|---------------------|-------------------------------------|-------------------------------------|-----------------------------|-----------------------------|----------------------------|
| PVDF | 50.96 | 37.35 | 162 | 139 | 48.72 |
| MWCNTs/PVDF | 25.71 | 21.00 | 162 | 139 | 24.58 |
| Ag NPs/PVDF | 45.64 | 34.81 | 162 | 139 | 43.63 |
| Ag NPs/MWCNTs /PVDF | 25.53 | 21.80 | 162 | 140 | 24.41 |

^{a)}the melting enthalpy of the first heating trace; ^{b)}the crystallization enthalpy of the cooling trace; ^{c)}the melting peak temperature; ^{d)}the crystallization temperature; ^{e)}the degree of mass crystallinity

To derive the information of sample crystallinity from the DSC results, the degree of mass crystallinity (X_c) was calculated from the melting enthalpy of the second heating trace (ΔH_m) with the following Equation 1:

$$X_c (\%) = \frac{\Delta H_m}{\Delta H_0} \quad (1)$$

where $\Delta H_0=104.6$ J/g was the heat of fusion for 100% crystalline PVDF.

Table S3. The T_i , T_{max} and Weight loss rate at T_{max} observed from **Figure S1**

| Samples | T_i [°C] ^{a)} | T_{max} [°C] ^{b)} | Weight loss rate at T_{max} (%) |
|-------------------|-----------------------------|---------------------------------|--------------------------------------|
| PVDF | 405.98 | 486.49 | 53.26% |
| MWCNTs/PVDF | 346.15 | 484.97 | 36.89% |
| Ag NPs/ PVDF | 368.27 | 478.87 | 39.78% |
| AgNPs/MWCNTs/PVDF | 359.80 | 479.63 | 36.51% |

^{a)}the initial degradation temperature; ^{b)}the temperature at the highest degradation rate

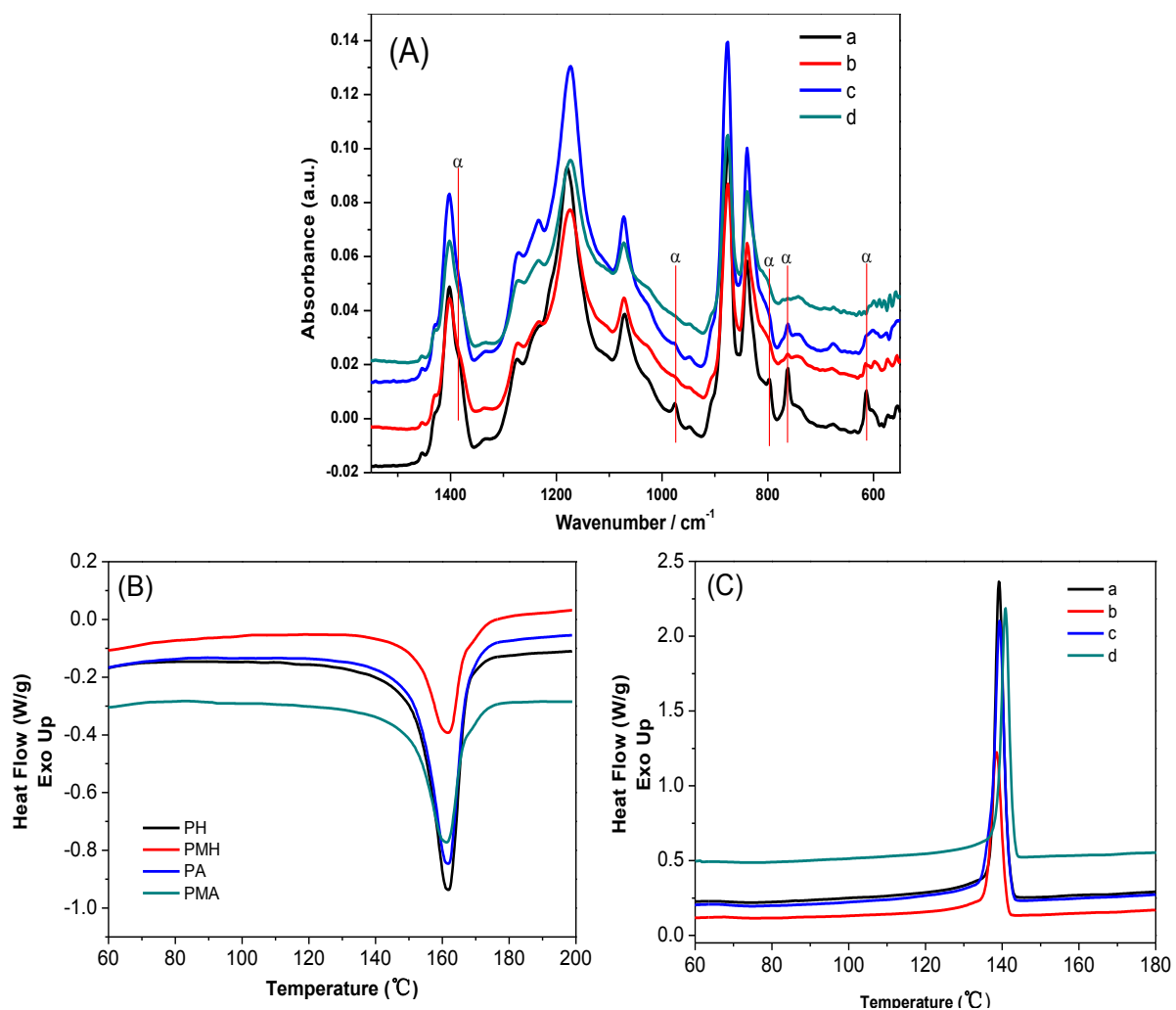


Fig. S1 FTIR spectra(A) and DSC heating (B) and cooling (C) curves of the PVDF(a), MWCNTs/PVDF(b), Ag NPs/PVDF(c) and Ag NPs/MWCNTs/PVDF(d) composites.

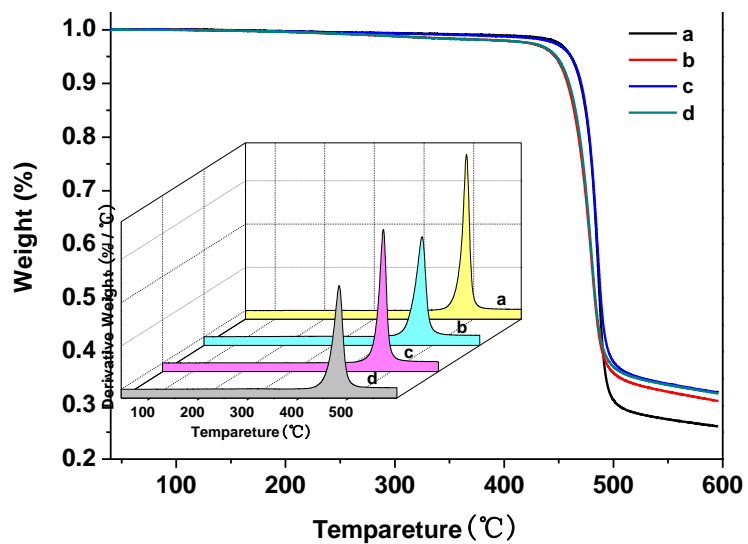


Fig. S2 TG and derivative TG (inset) curves of the PVDF(a), Ag NPs/ PVDF(b), MWCNTs/PVDF(c) and Ag NPs/ MWCNTs/PVDF(d) composites.