

Supporting Information to Accompany

Fluorescent Carboxylic and Phosphonic Acids: Comparative Photophysics from Solution to Organic Nanoparticles

Contribution from

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Experimental Section

1. Infrared spectra in CCl_4 solution

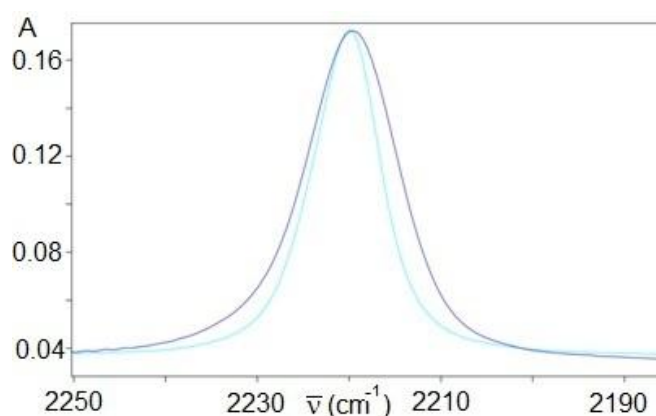


Fig. S1 Infrared absorption band of the CN unit, centered at 2220 cm^{-1} for **fPOH** (—) and **fPOEt** (---) in CCl_4 solution ($10^{-4}\text{ mol.L}^{-1}$).

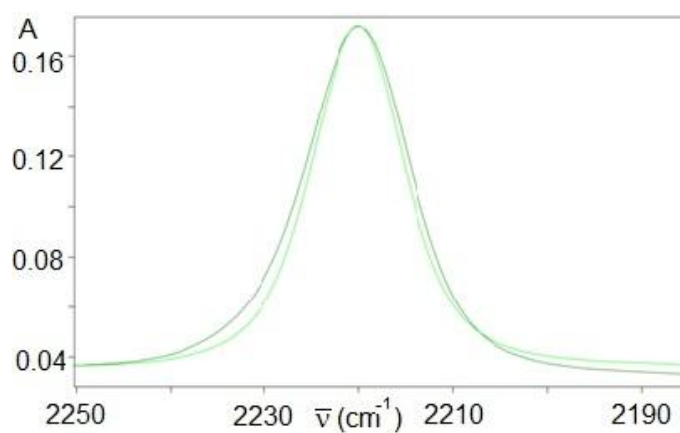


Fig. S2 Infrared absorption band of the CN unit, centered at 2220 cm^{-1} for **fPOH** (—) and **fPOEt** (—) in CHCl_3 solution ($10^{-4} \text{ mol.L}^{-1}$).

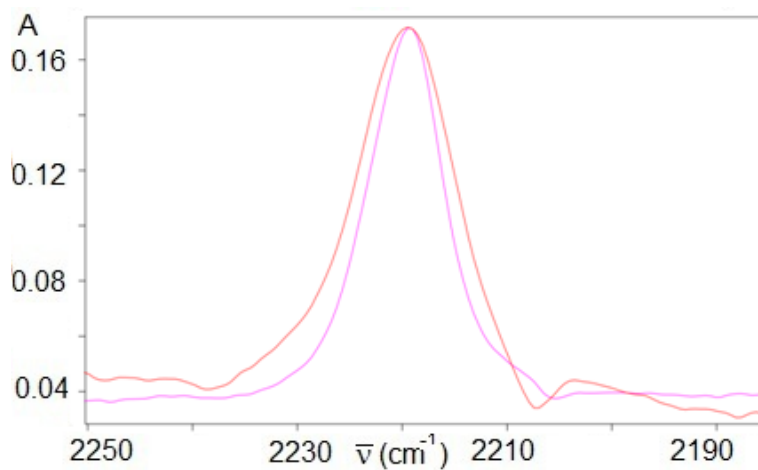


Fig. S3 Infrared absorption band of the CN unit, centered at 2220 cm^{-1} for **fPOH** (—) and **fPOEt** (—) in toluene solution ($10^{-4} \text{ mol.L}^{-1}$).

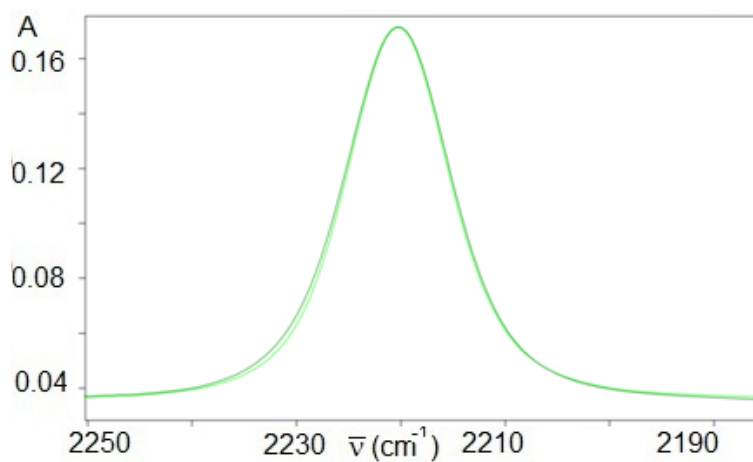


Fig. S4 Infrared absorption band of the CN unit, centered at 2220 cm^{-1} for **fCO₂H** (—) and **fOtBu** (—) in CHCl_3 solution ($10^{-4} \text{ mol.L}^{-1}$).

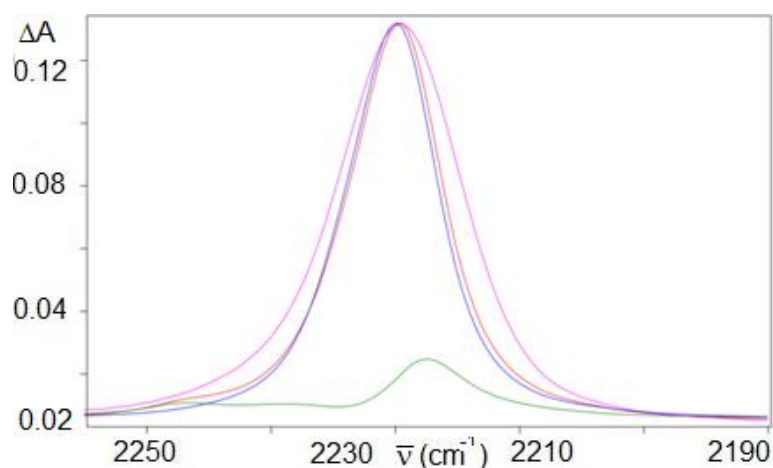
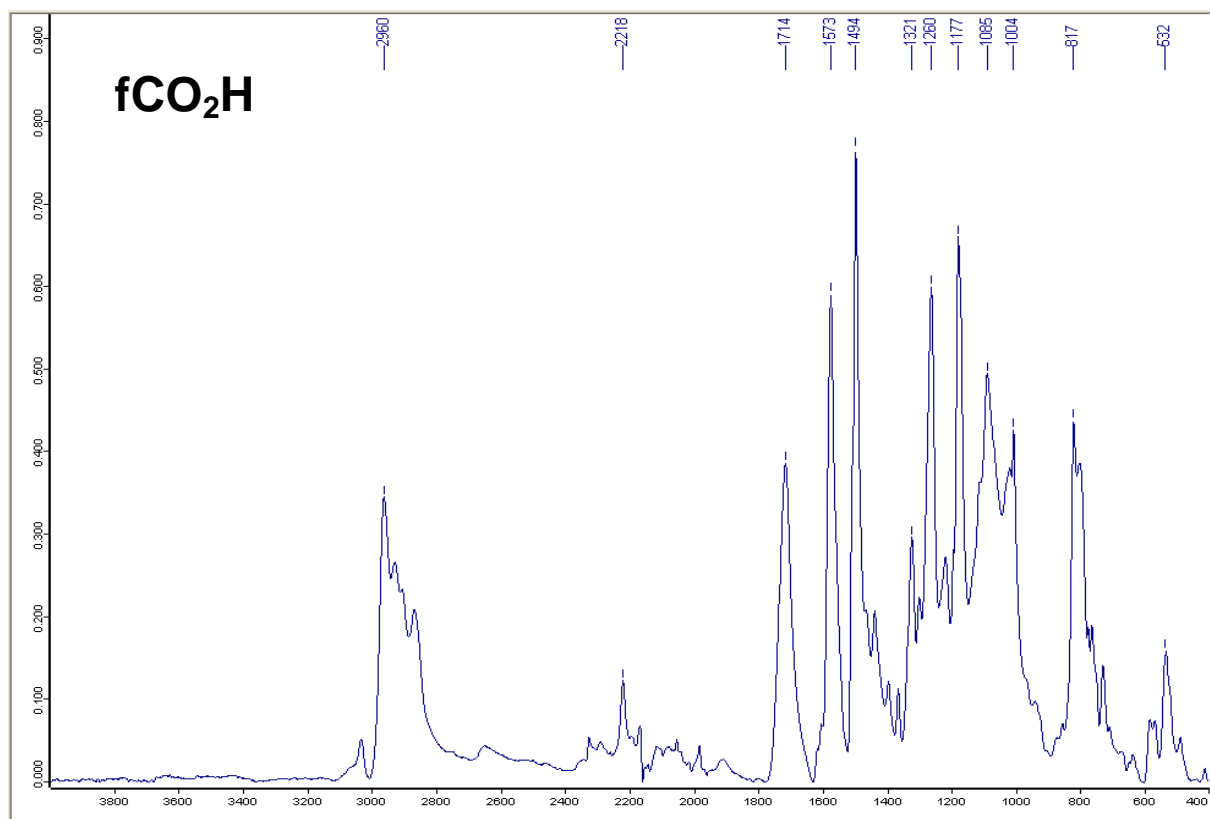
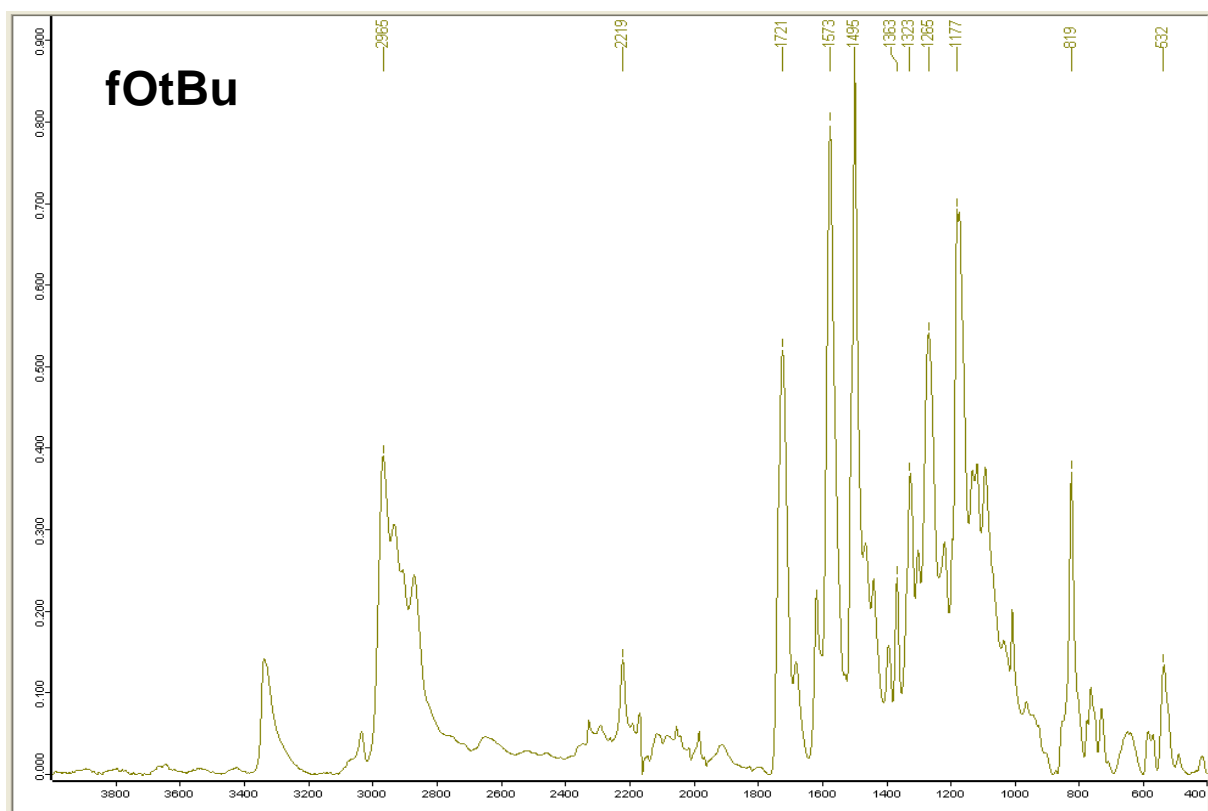
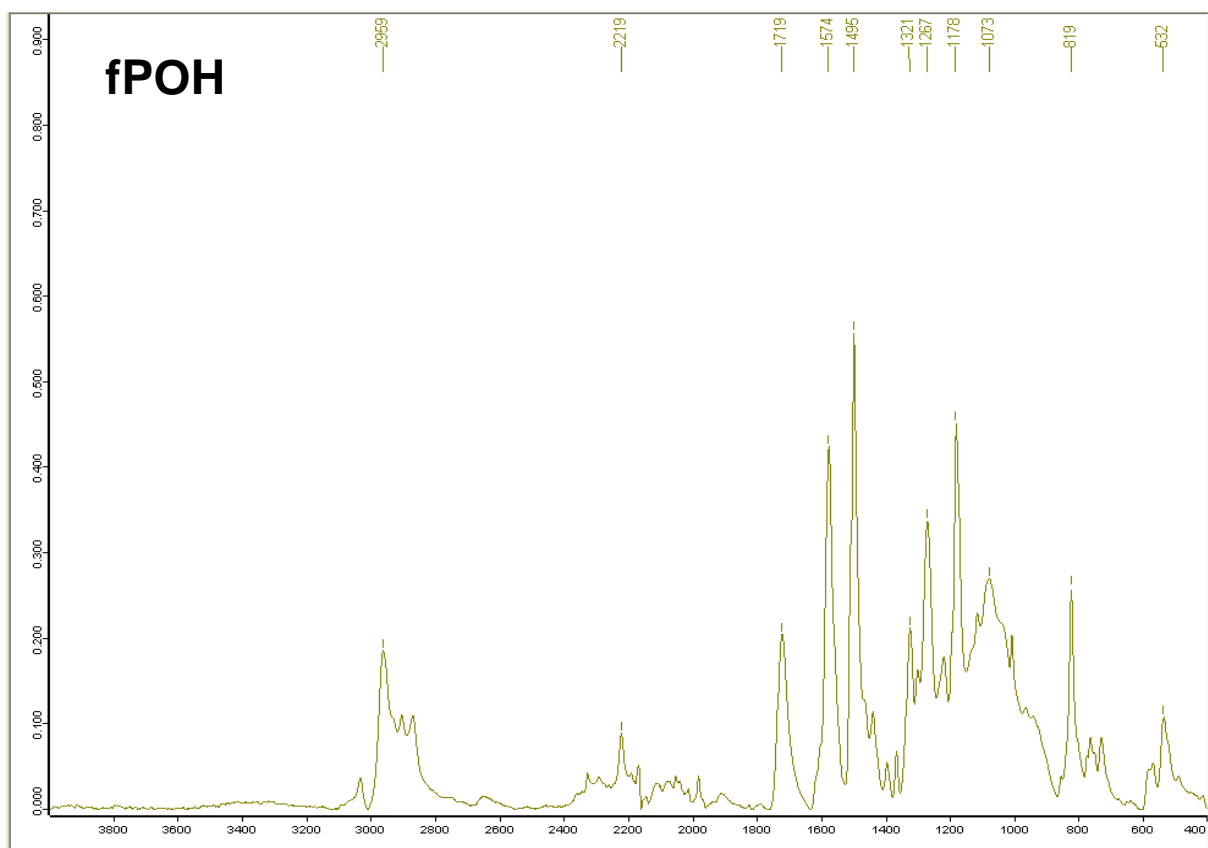


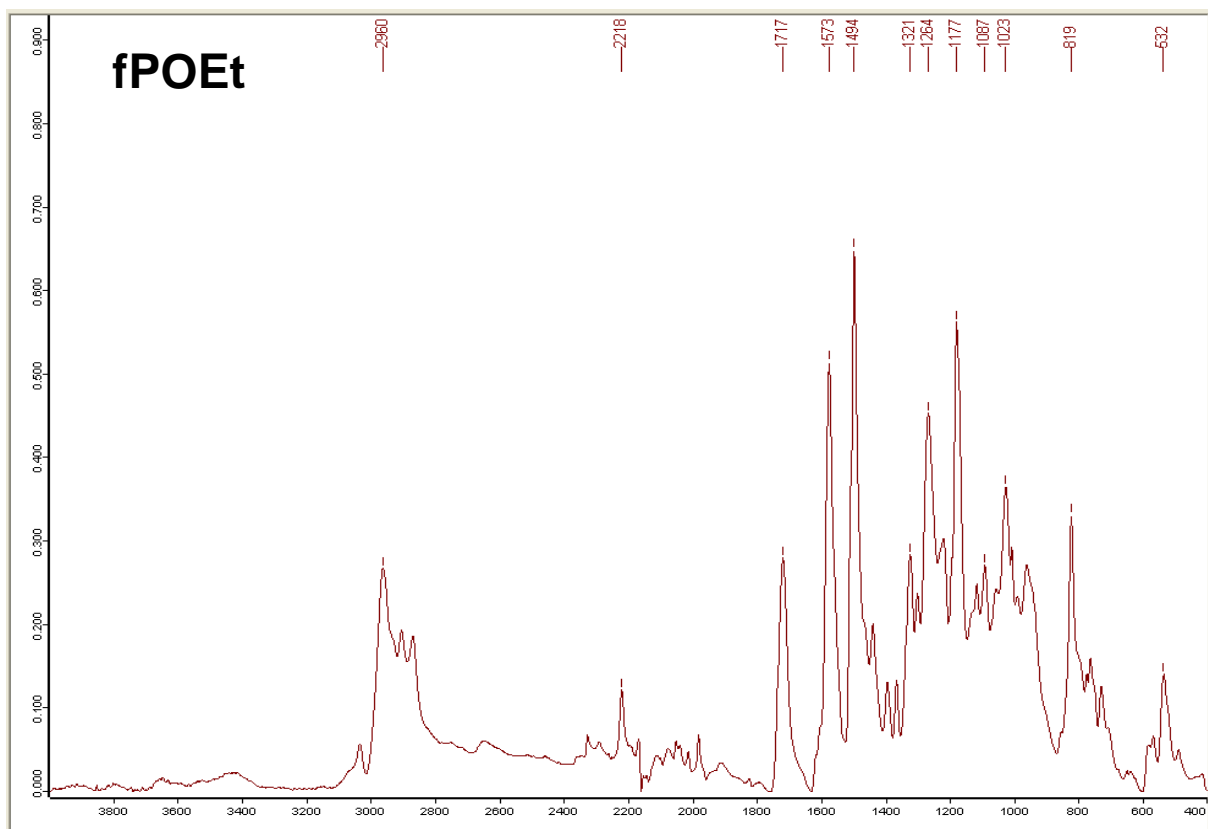
Fig. S5 Infrared absorption band of the CN unit, centered at 2220 cm^{-1} for **fPOH** (—) and **fPOEt** before (—) and (—) after adding 4-fluorophenol (4FP) in excess in CCl_4 solution ($10^{-4} \text{ mol.L}^{-1}$). The spectrum resulting from the subtraction after and before adding 4FP to the solution of fPOEt (—) is indicated and shows a neat bathochromic shift, featuring strong hydrogen bonding.

2. Infrared spectra in the solid state

Solid state spectra were recorded in the ATR mode using a FTIR Bruker Vertex 70 spectrometer.

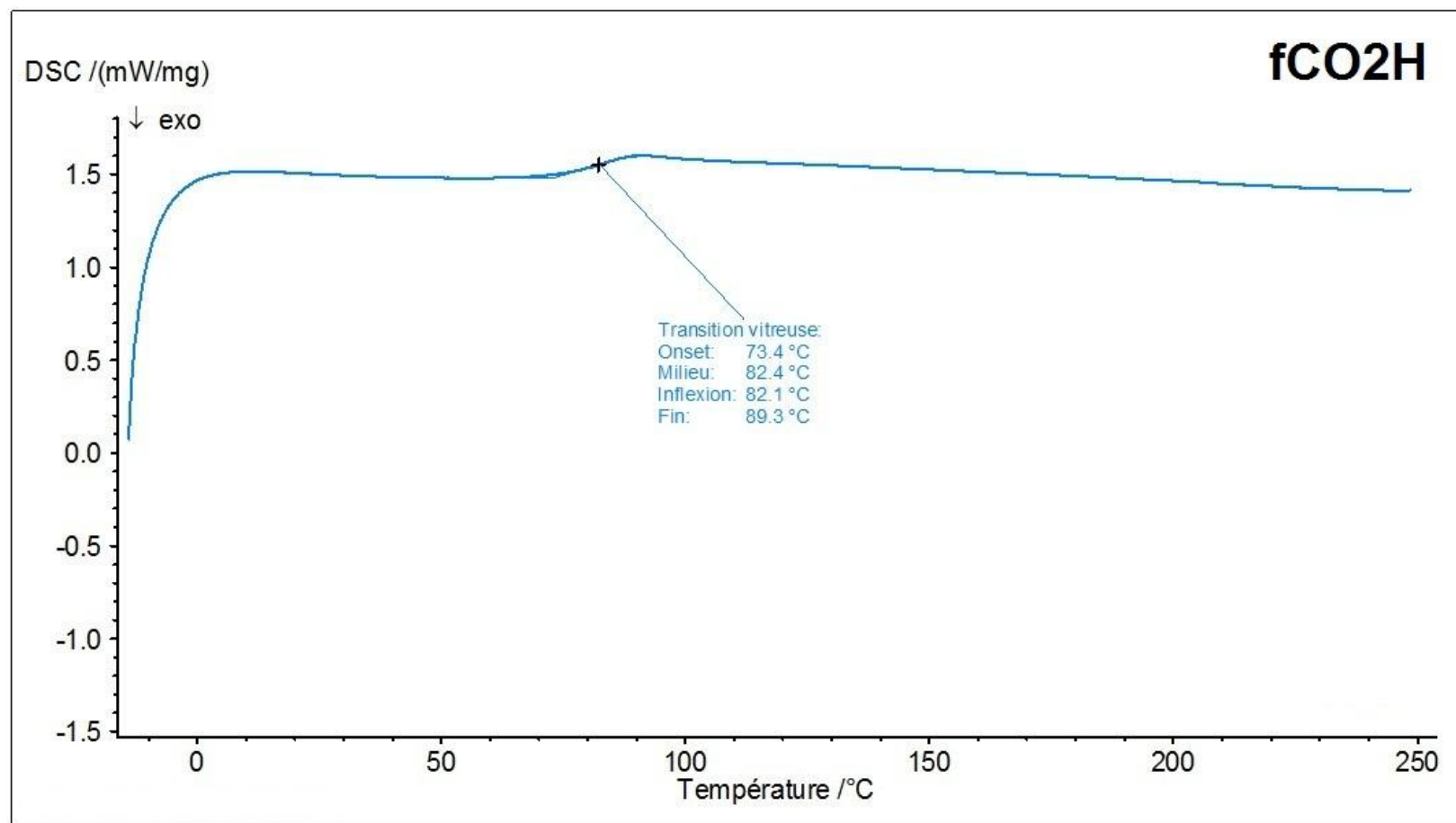


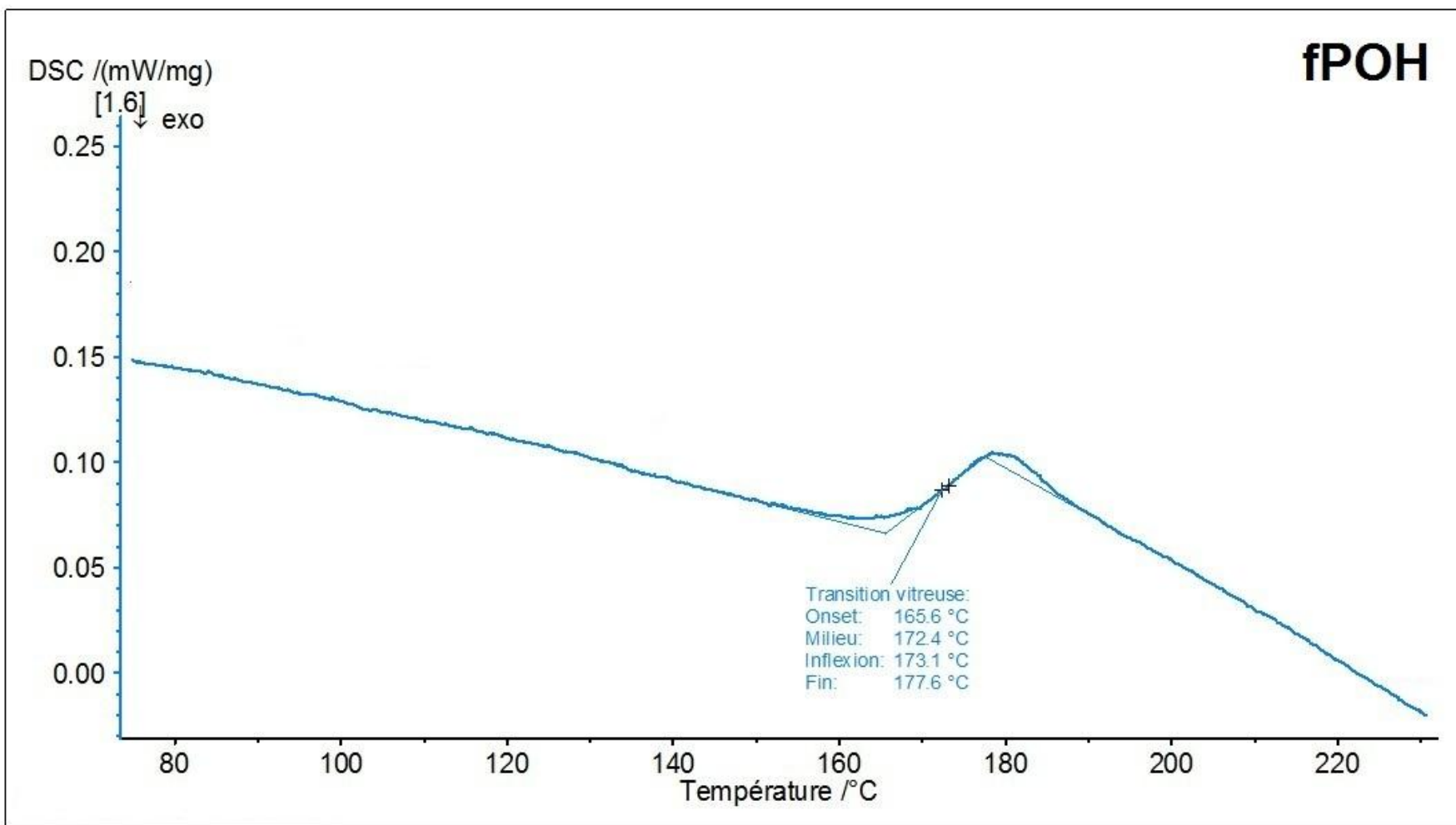


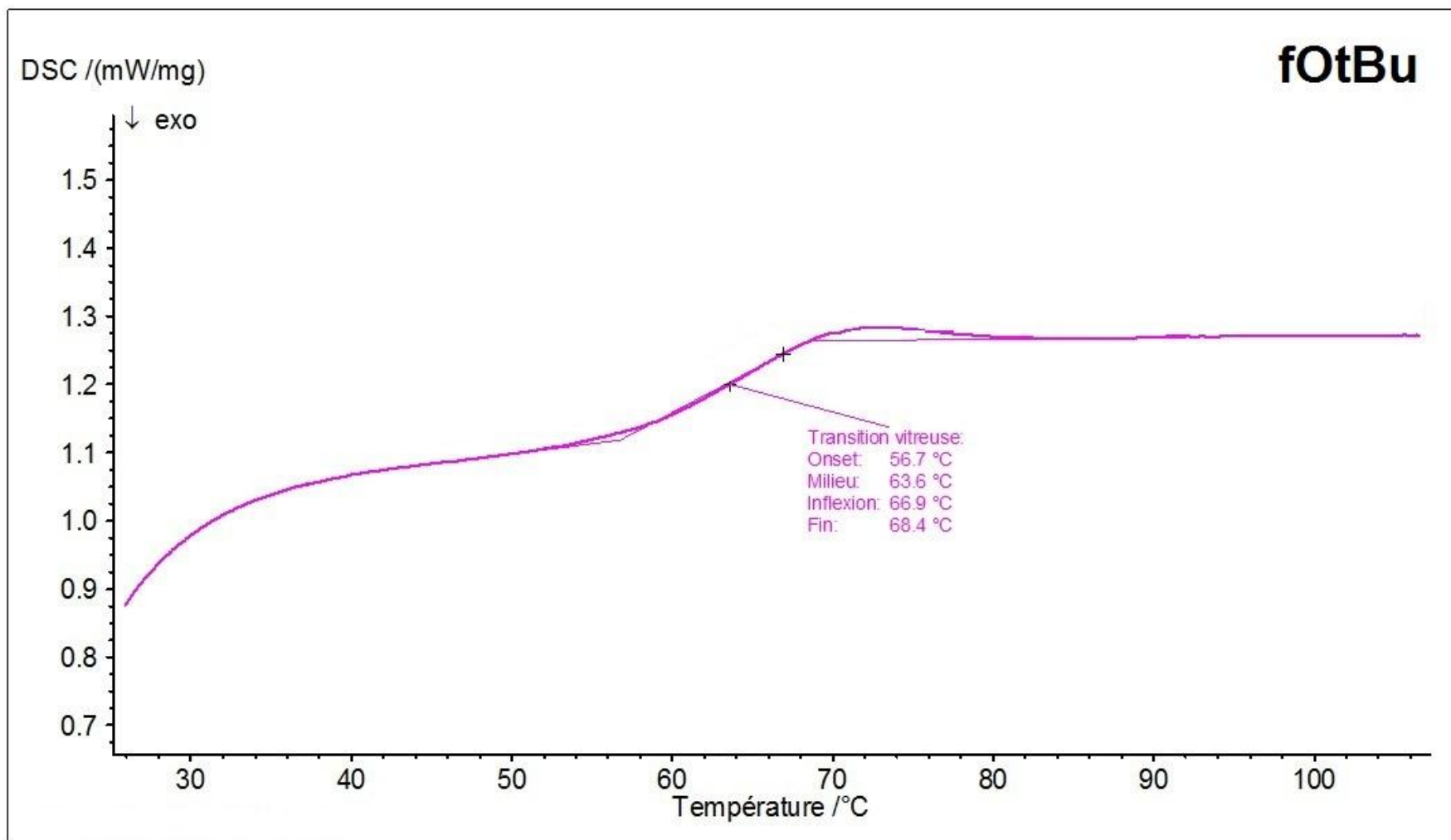


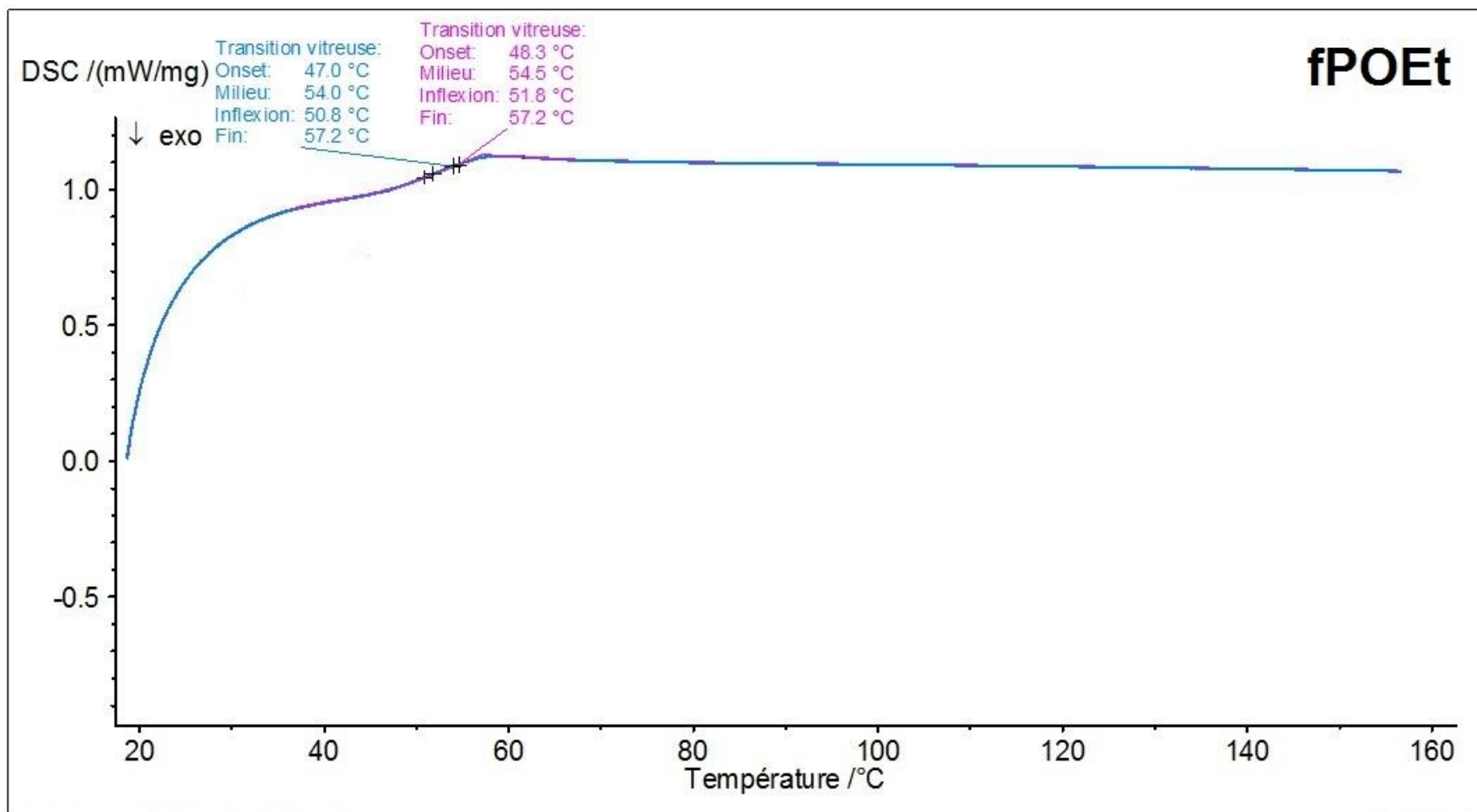
3. DSC thermal analyses

Thermal properties were measured by using differential scanning calorimetry (Netzsch- Maia DSC 200 F3) in alumina caps under a nitrogen flow at a scan rate of $30\text{ }^{\circ}\text{C}\cdot\text{min}^{-1}$ over the temperature range $[-10^{\circ}\text{C} - 250^{\circ}\text{C}]$









Thermal gradient: blue curve 20 °C.min⁻¹ ; pink curve : 30 °C.min⁻¹.

