

Evaluation of thiocarbonyl and thioester moieties as thiol protecting groups for controlled radical polymerization

Morgane Le Neindre,^a Benoît Magny^b and Renaud Nicolay^{*a}

^a Matière Molle et Chimie (ESPCI-CNRS, UMR 7167), ESPCI ParisTech, 10 rue Vauquelin, 75005 Paris, France

^b Centre de Recherche de l'Oise, Arkema France, Parc Technologique ALATA – BP 22, 60550 Verneuil en Halatte, France.

* Corresponding author: email renaud.nicolay@espci.fr

Supporting Information

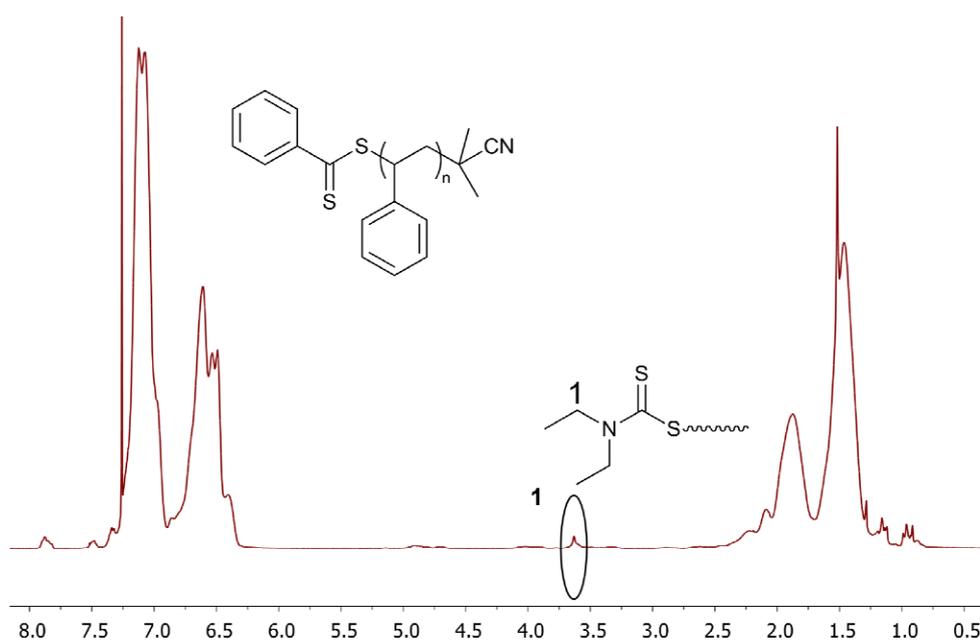


Figure S1 ¹H NMR spectrum of P1

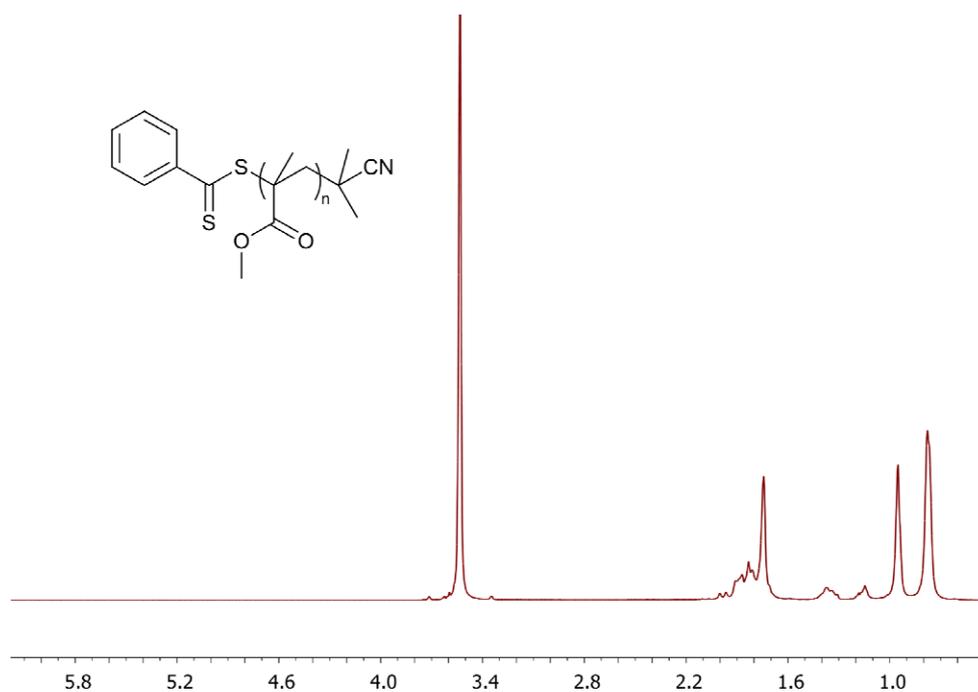


Figure S2 ¹H NMR spectrum of P2

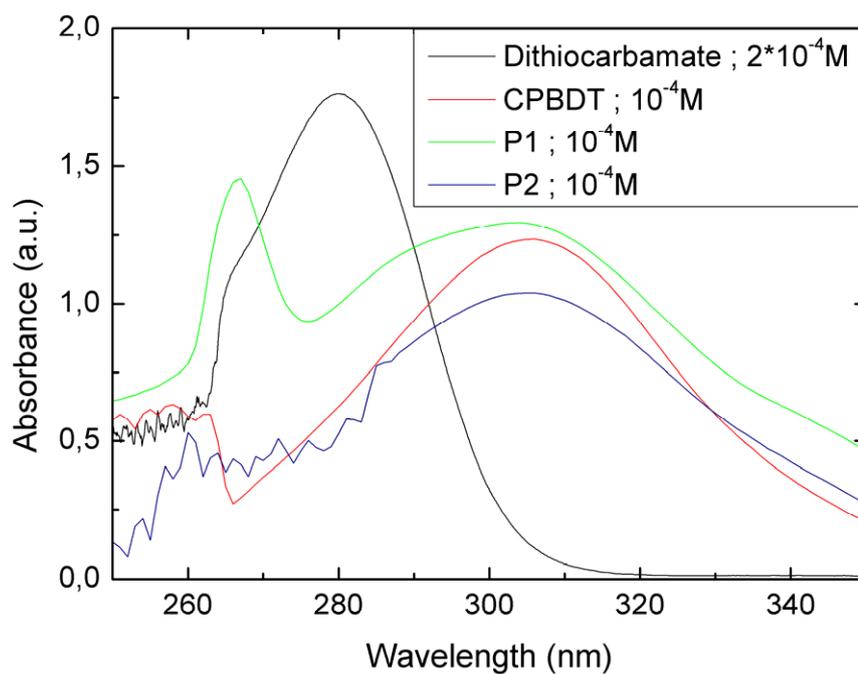


Figure S3 UV/Vis spectra of the dithiocarbamate model molecule (2×10^{-4} M in DMF), CPBDT (10^{-4} M in DMF), P1 (10^{-4} M in DMF) and P2 (10^{-4} M in DMF).

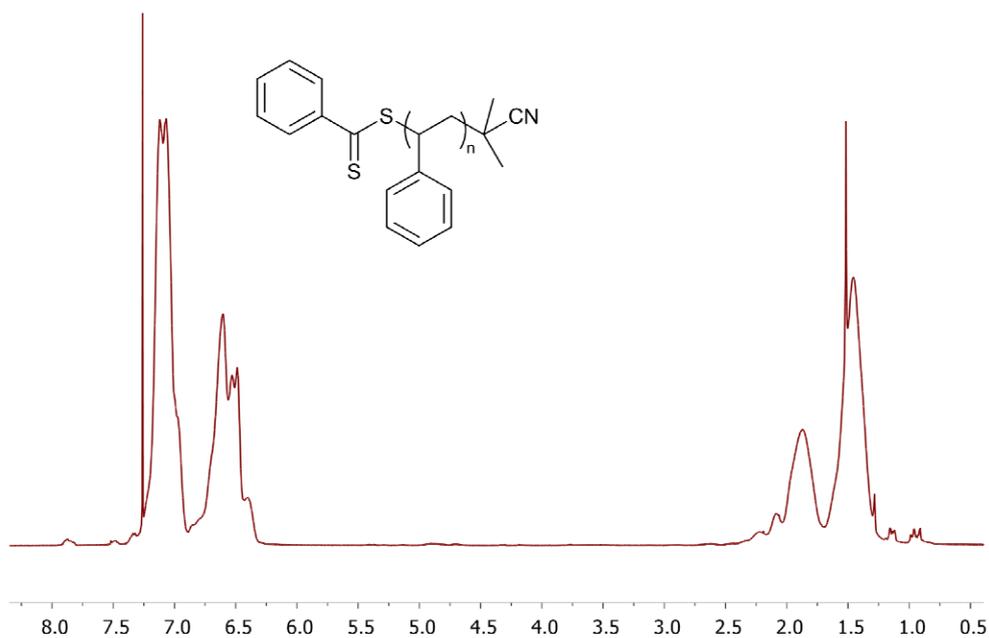


Figure S4 ¹H NMR spectrum of P4

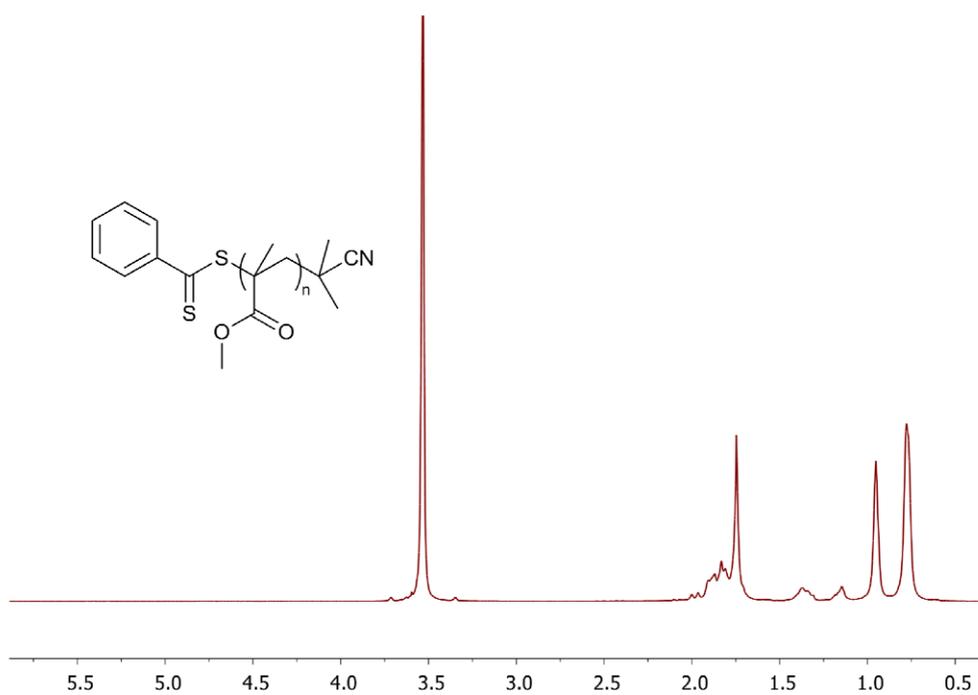


Figure S5 ¹H NMR spectrum of P5

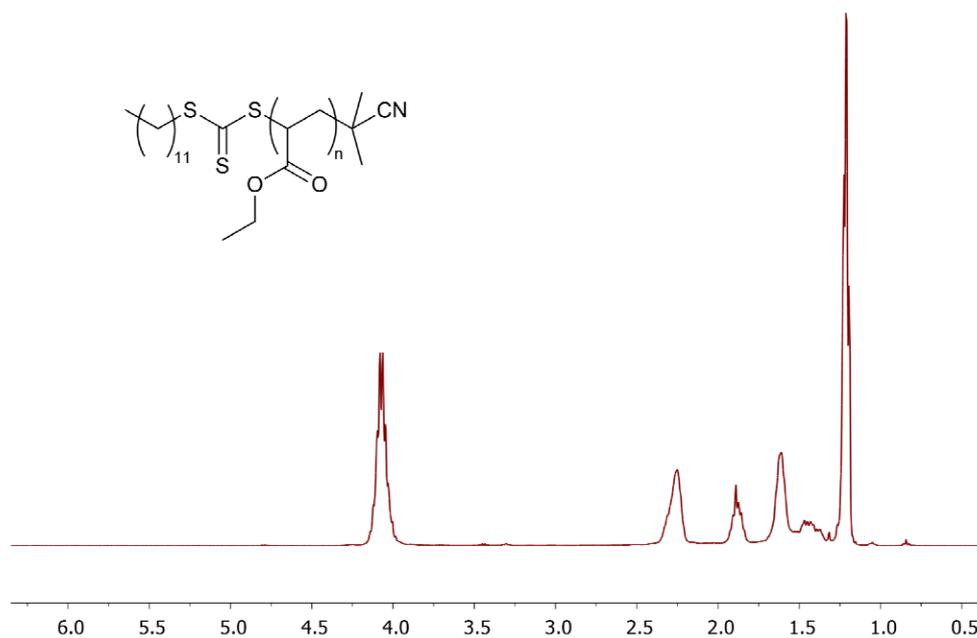


Figure S6 ¹H NMR spectrum of P6

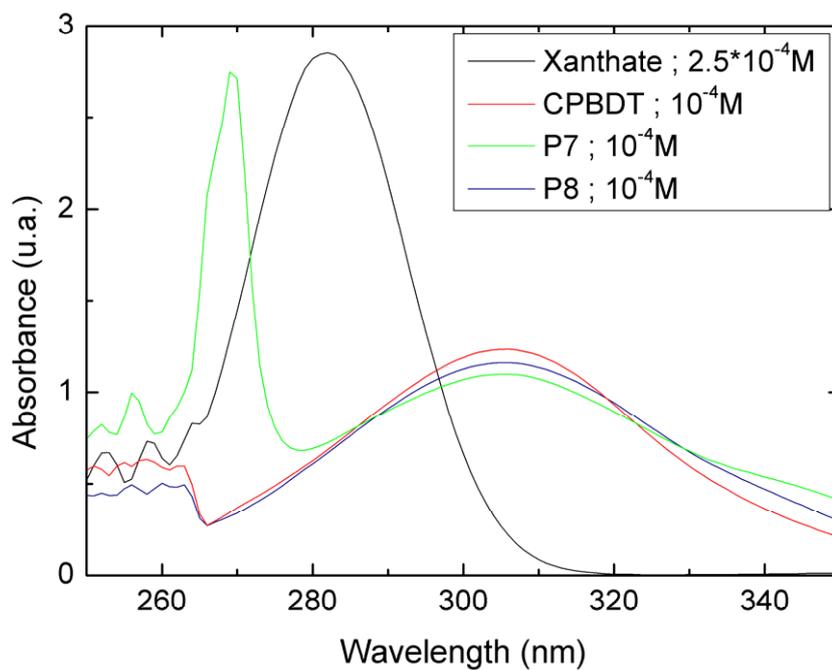


Figure S7 UV/Vis spectra of the xanthate model molecule (2.5×10^{-4} M in DMF), CPBDT (10^{-4} M in DMF), P7 (10^{-4} M in DMF) and P8 (10^{-4} M in DMF).

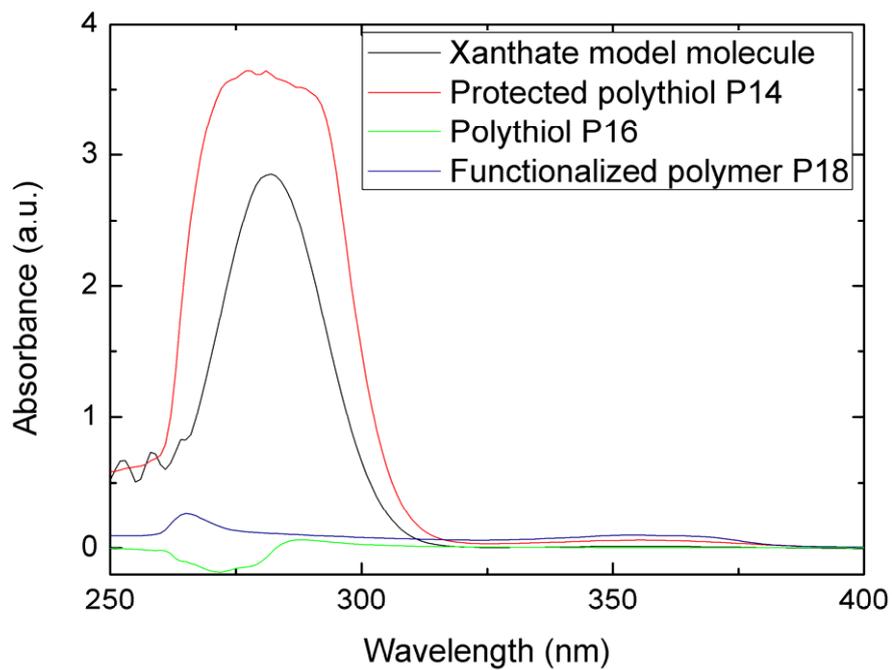


Figure S8 UV/Vis spectra of the xanthate model molecule (2.5×10^{-4} M in DMF), **P14** (10^{-4} M in DMF), **P17** (10^{-4} M in DMF) and **P18** (10^{-4} M in DMF).