

Supplementary Information for

**Enantio- and Diastereoselective Polymerization: Asymmetric Allylic Alkylation
Catalyzed by a Planar-Chiral Cp'Ru Complex**

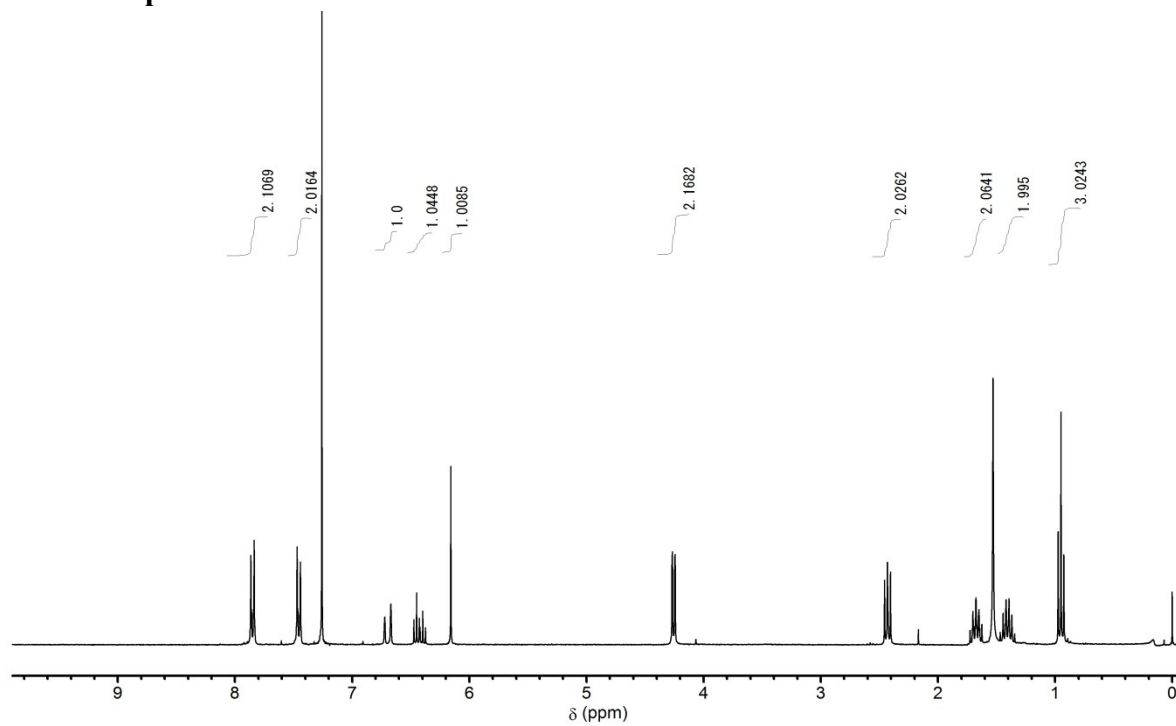
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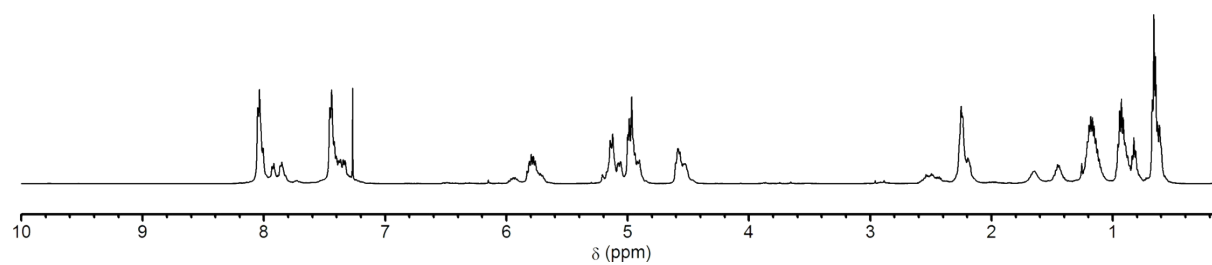
onitsuka@chem.sci.osaka-u.ac.jp

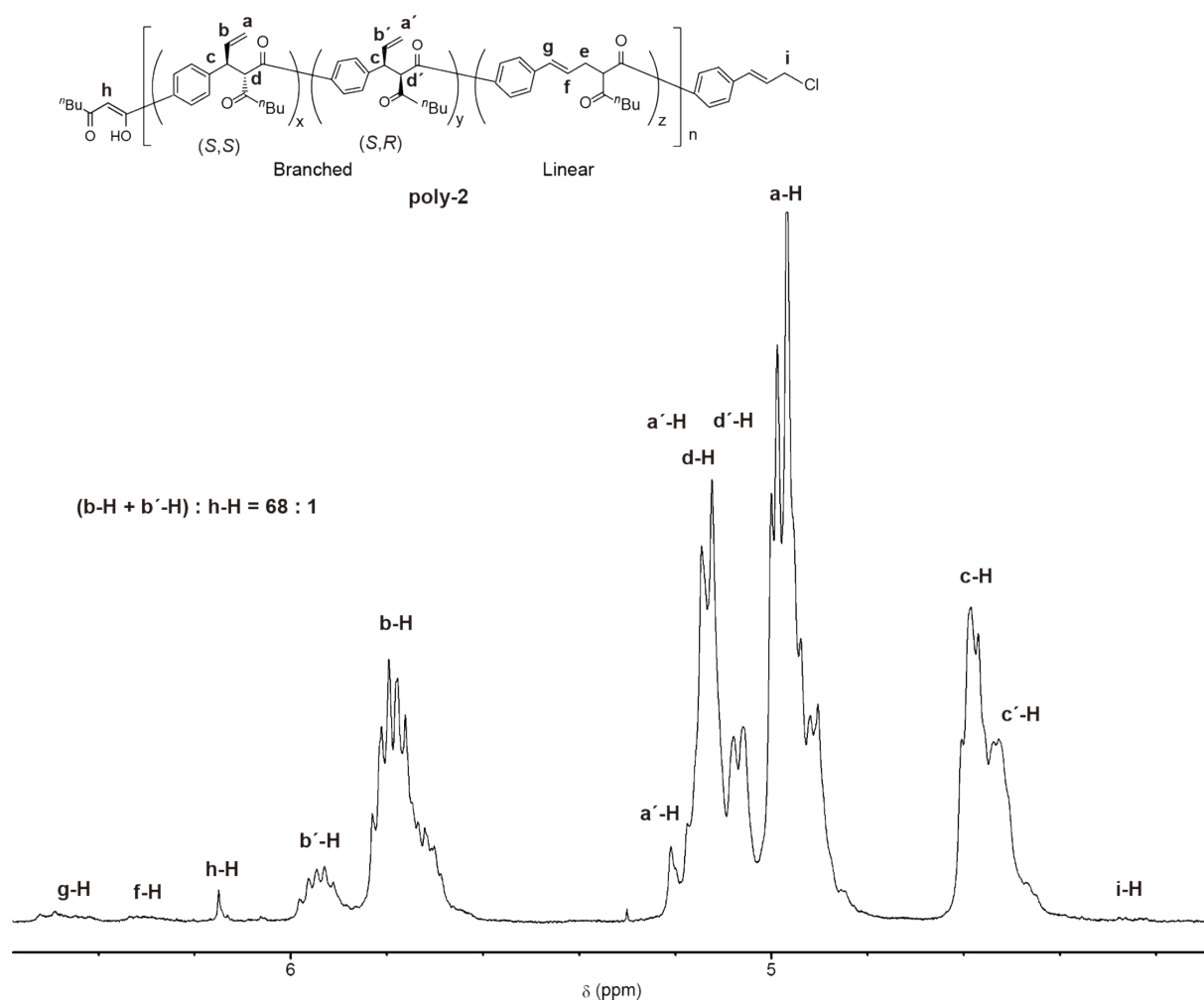
naokou@chem.sci.osaka-u.ac.jp

¹H NMR Spectrum of 2

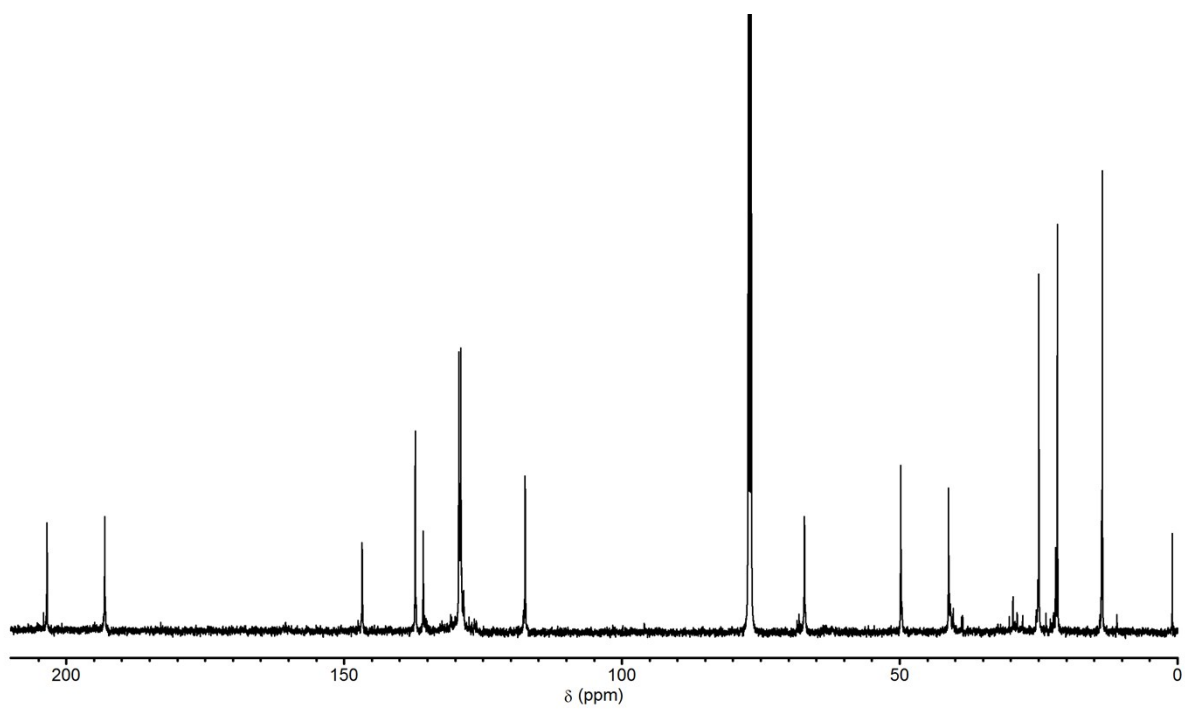


¹H NMR Spectrum of poly-2

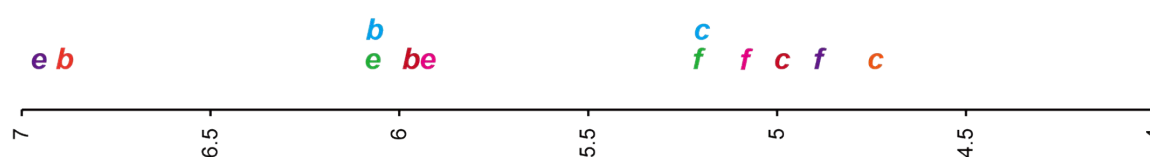
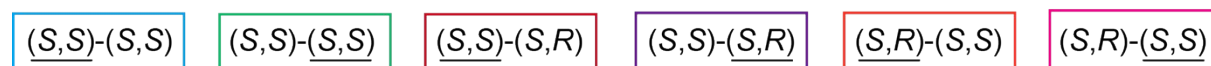
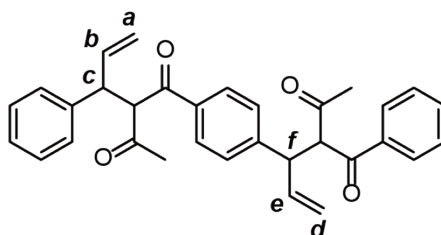




¹³C NMR Spectrum of poly-2

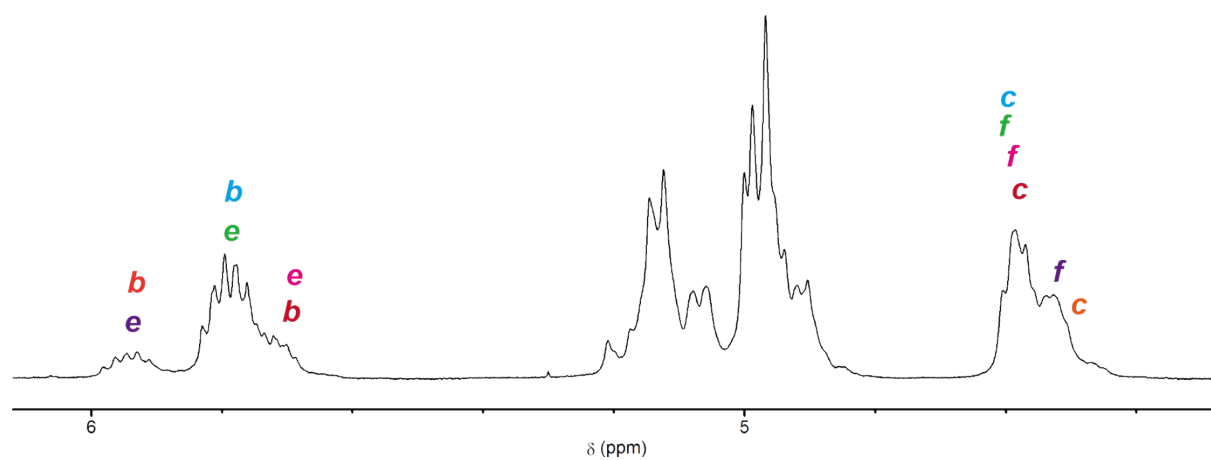


Simulation of ^1H NMR spectrum of **6 Using DFT (B3LYP-GIAO/6-311+G(2df,2p)) Calculation**

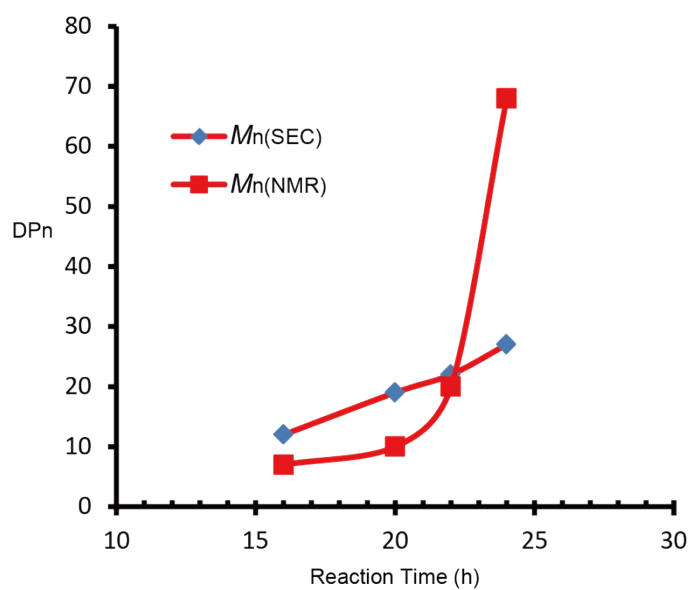


	6 -(<i>S,S</i>),(<i>S,S</i>)	6 -(<i>S,S</i>),(<i>S,R</i>)	6 -(<i>S,R</i>),(<i>S,S</i>)
b	6.08	6.07	6.91
c	5.19	4.99	4.74
e	5.97	6.94	5.96
f	5.19	4.89	5.08

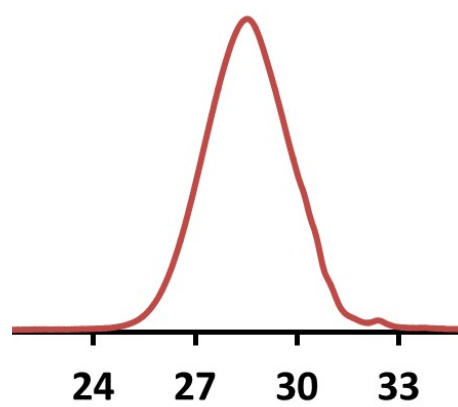
Assignment of ^1H NMR Spectrum of Poly-2 Based on DFT (B3LYP-GIAO/6-311+G(2df,2p)) Calculation



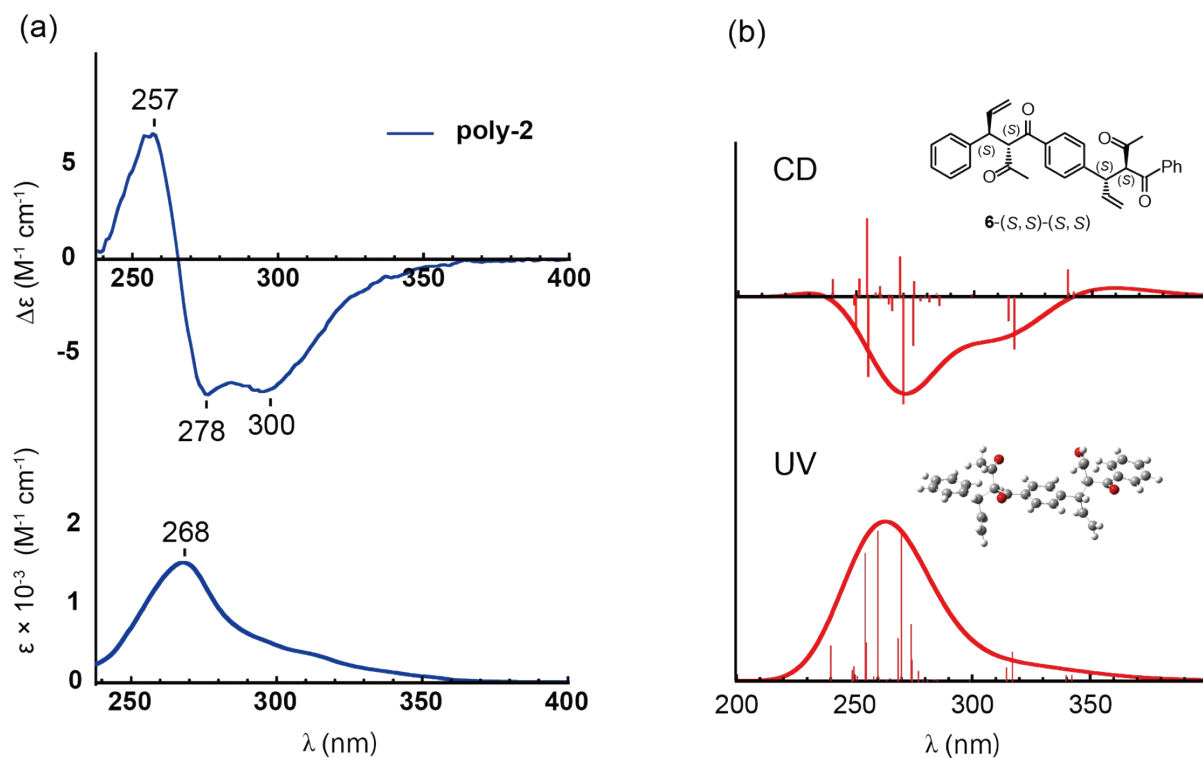
Time dependence of polymerization degree determined by ^1H NMR or SEC.



The SEC spectrum of crude product poly-2 (Table 1, entry 5).



TD-DFT (B3LYP/6-31G) simulated CD and UV spectra of 6-(*S,S*)-(*S,S*)**



The proposed coordination structure of (*S*)-7. The structures are based on simulated structures of 5-(*S,S*) and 5-(*S,R*).

