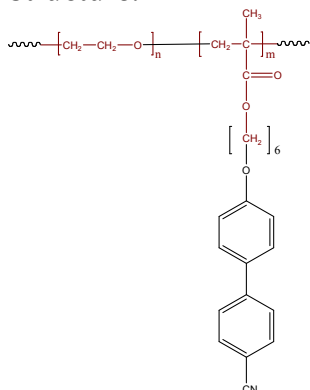


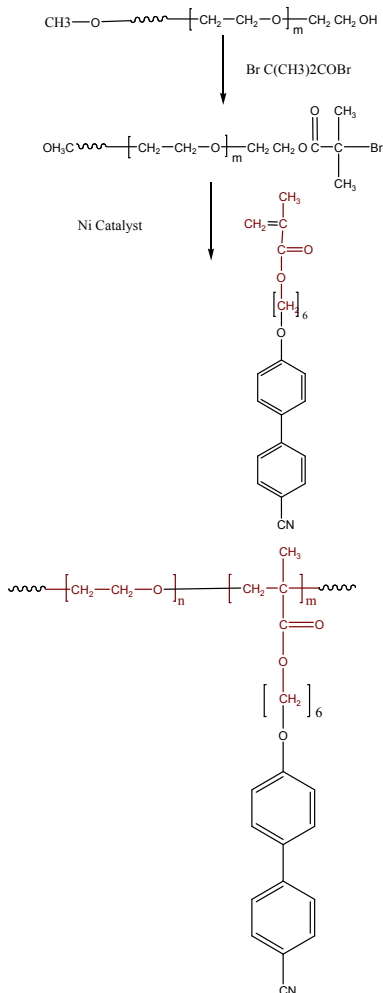
**Poly(ethylene oxide-b-6-(4'-cyanobiphenyl-4-yloxy)hexyl methacrylate**

### Structure:



Mn x 10 <sup>3</sup> PEO-b-4CNBPMA	PDI
2.0-b-9.0	1.19

The scheme of the reactions are illustrated below:



The un-reacted PEG can be removed by stirring the polymer in hot water. The obtained polymer dissolved in CHCl<sub>3</sub>/Toluene and pass through the column packed with silica. The Diblock copolymer obtained through the 2<sup>nd</sup> route where the macroinitiator of PEG bearing Br terminal group was used to initiate polymerization of 4-cyanobiphenyl-4-yloxy hexyl methacrylate. The obtained polymer solution in toluene/CHCl<sub>3</sub> was passed through a column packed with silica to remove the traces amount of Nickel catalyst. The polymer was further purified by stirring in hot water to remove unreacted PEG macroinitiator. The polymer was recovered by precipitation in cold ether/hexane mixture.

Polymer is soluble in CHCl<sub>3</sub>, THF, toluene. The polymer precipitated out from hexane.

1H NMR spectrum of P9311-EO4CNBPHMA in CDCl<sub>3</sub>. The spectrum shows peaks in the aromatic region (7.0-7.8 ppm) and aliphatic region (1.0-2.0 ppm). Integration values are shown below the peaks: 1.53, 1.00, 2.00, 4.00, 2.00, and 4.00. The x-axis is labeled 'ppm (f1)' and ranges from 10 to 0.

— Poly(ethylene oxide),  $M_n=2000$ ,  $M_w=2100$ ,  $PI=1.05$

— Block Copolymer PEO(2000)-b-4-CNBPHMA (9000). PI=1.19