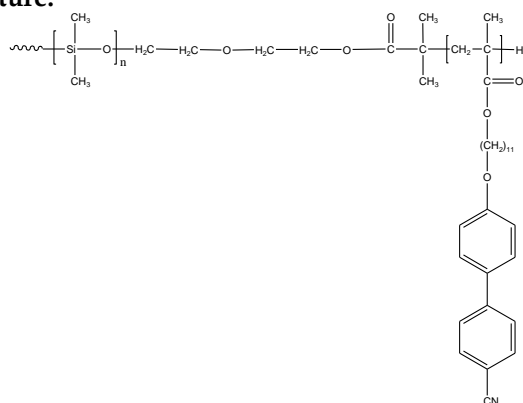


**Sample Name:** Poly(dimethylsiloxane-b-11-(4'-cyanobiphenyl-4-yloxy)undecylmethacrylate)

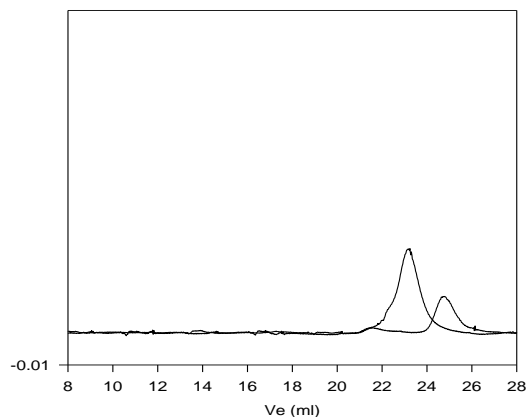
**Sample #:** P9917-DMS4CNBP11CMA

**Structure:**



**SEC profile of the block copolymer**

**P9917-DMS4CNBP11CMA**



Size exclusion chromatography of the product:

— Polydimethylsiloxane,  $M_n=5000$ ,  $M_w=5400$ ,  $PI=1.05$   
 — Block Copolymer PDMS(5000)-b-4CNBP11CMA(20,000),  $PI=1.3$

**Composition:**

$M_n \times 10^3$ DMS-b-4CNBP-11LCMA	$M_w/M_n$ (PDI)
5-b-20	1.3

**Synthesis Procedure:**

Poly(dimethylsiloxane-b-6-(4'-cyanobiphenyl-4-yloxy)undecylmethacrylate) is prepared by ATR-polymerization of 6-(4'-cyanobiphenyl-4-yloxy)undecylmethacrylate with bromo-terminated polydimethylsiloxane as macro-initiator.

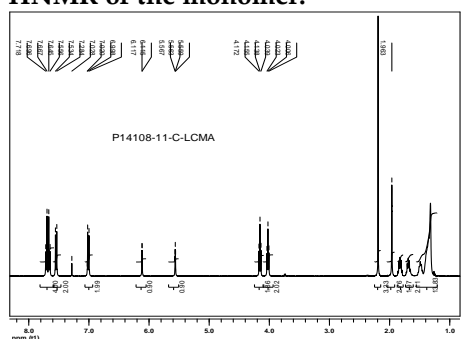
**Characterization:**

An aliquot of the bromo-terminated poly(dimethyl siloxane) was analyzed by size exclusion chromatography (SEC) to obtain the molecular weight and polydispersity index (PDI). The final block copolymer composition was calculated from  $^1H$ -NMR spectroscopy by comparing the peak area of the dimethyl siloxane protons near 0 ppm with the 4CNBPMA protons at about 3.6 ppm. Block copolymer PDI is determined by SEC.

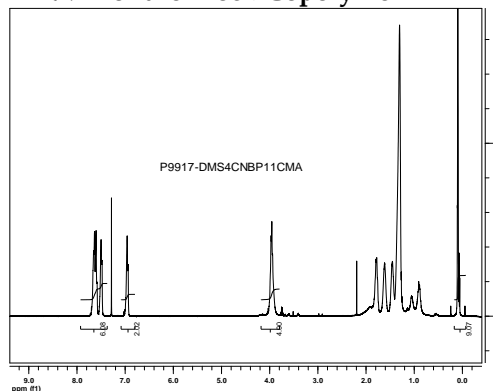
**Solubility:**

Poly(dimethylsiloxane-b-6-(4'-cyanobiphenyl-4-yloxy)undecylmethacrylate) is soluble in THF, chloroform and toluene. It is precipitated in methanol.

**$^1H$ NMR of the monomer:**



**$^1H$  NMR of the Block Copolymer**



## Thermal analysis of the P9917-DMS4CNBP11CMA

Thermal analysis of the samples was carried out on a TA Q100 differential scanning calorimeter at a heating rate of 10°C/min. The midpoint of the slope change of the heat flow plot of the second heating scan was considered as the glass transition temperature ( $T_g$ ).

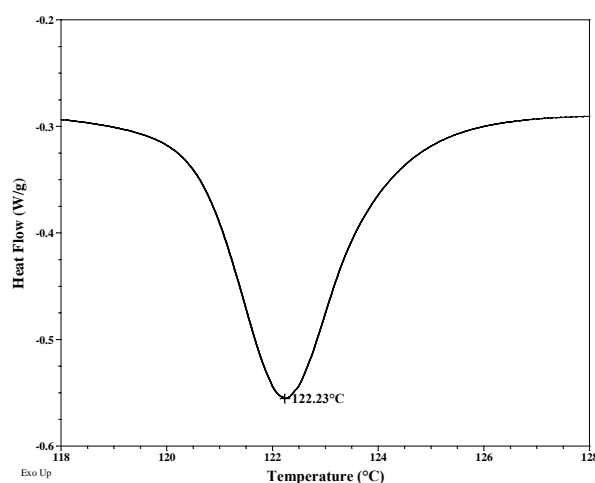
## Thermal analysis results at a glance

Sample	$T_m$ (°C)	$T_c$ (°C)	$T_g$ (°C)
DMS block	Not distinct	Not distinct	-127 (Lit)
<b>4CNBP-11C-MA block</b>	122	118	-

## Melting and crystallization curve for the sample

The melting temperature ( $T_m$ ) was taken as the maximum of the endothermic peak where as the crystallization temperature ( $T_c$ ) was considered as the minimum of the exothermic peak.

## Melting curve for 4CNBP11CMA block:



## Crystallization curve for 4CNBP-11C-MA block:

