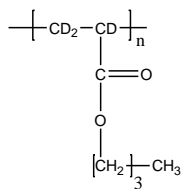


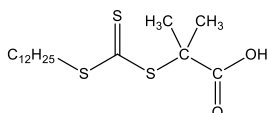
Backbone protons are deuterated

Structure:



Mn x 10 ³	PDI
6.5	1.12

RAFT used in this synthesis :



Chemical Formula: $C_{17}H_{32}O_2S_3$
Molecular Weight: 364.6

Chemical structure of the polymer repeat unit is shown above the spectrum. The structure includes a long aliphatic chain and a side chain containing a carboxylic acid group. The assignments and integrations for the peaks are as follows:

- 0.86: Methyl protons (integration 1.19-1.32)
- 1.19-1.32: Aliphatic chain protons (integration 1.36)
- 1.36: Methyl protons (integration 1.64)
- 1.64: Methyl protons (integration 3.25)
- 3.25: CH₂ protons (integration 1.68)
- 6.80: CH proton (integration 1.68)
- 1.68: Carboxylic acid proton (integration 1.68)

The spectrum shows a broad peak at 11.0 ppm (OH), a sharp peak at 6.80 ppm (CH), and a complex multiplet between 0.8 and 3.3 ppm.

The product was characterized by size exclusion chromatography (SEC) and ¹H NMR.

Agilent GPC/SEC Software

