

Supporting Information

RAFT-prepared α -Difunctional Poly(2-vinyl-4,4-dimethylazlactone)s Scaffolds and their Derivatives: Synthesis and Effect of End Groups on Aqueous Inverse Temperature Solubility

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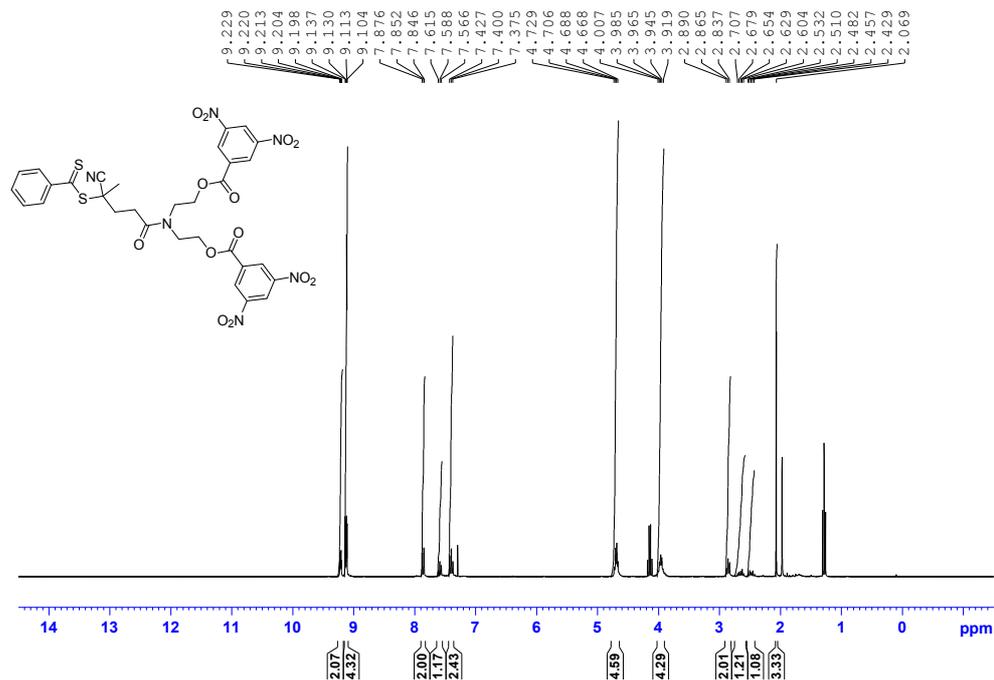
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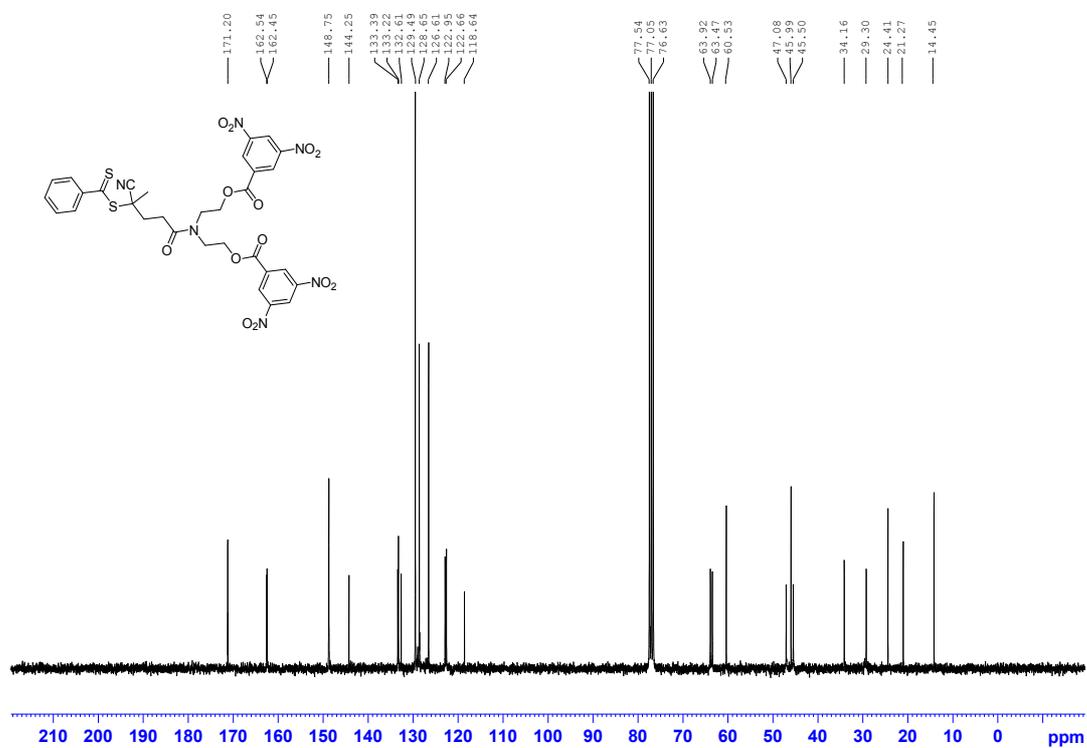
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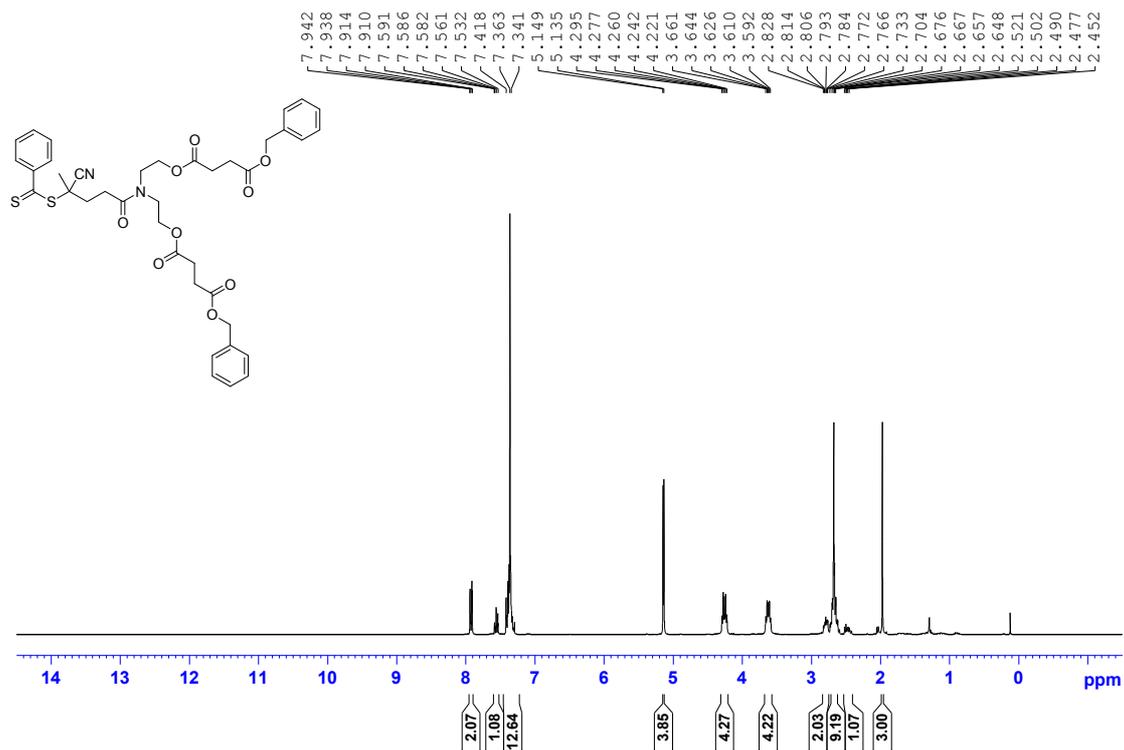
¹H and ¹³C-NMR results of branched RAFT agents



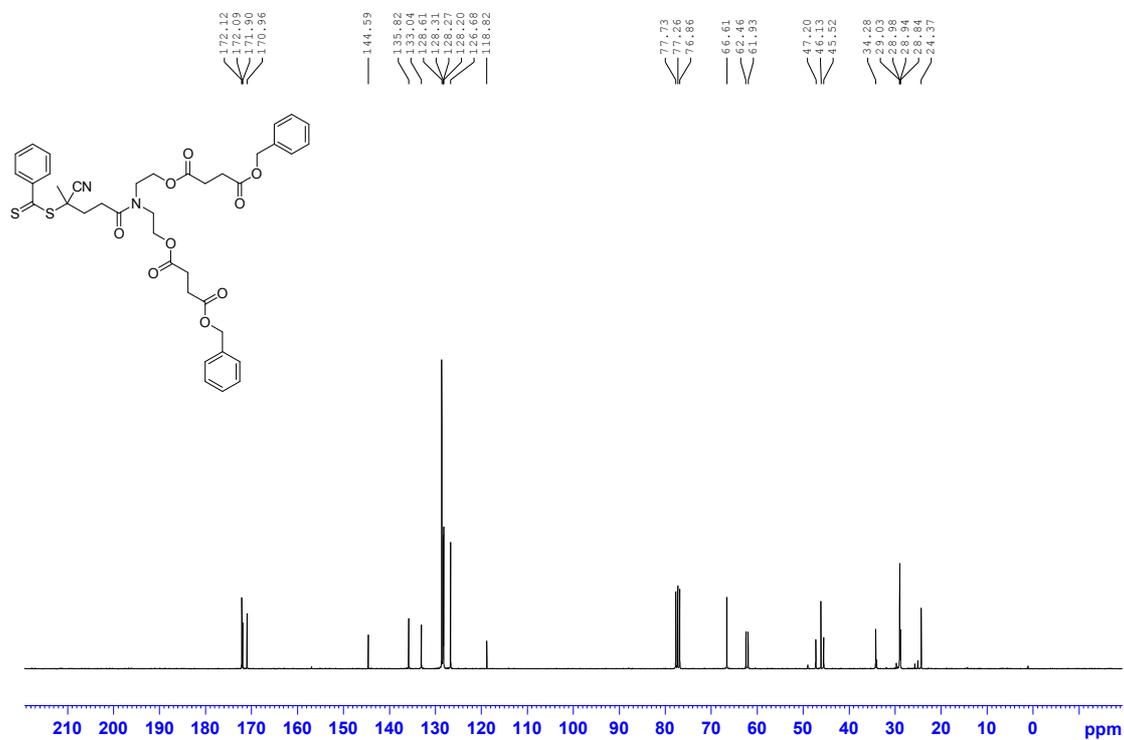
¹H-NMR of ((4-cyano-4-((phenylcarbo thioyl)thio)pentanoyl)azanediy)bis(ethane-2,1-diyl) bis(3,5-dinitrobenzoate) (CTA1)



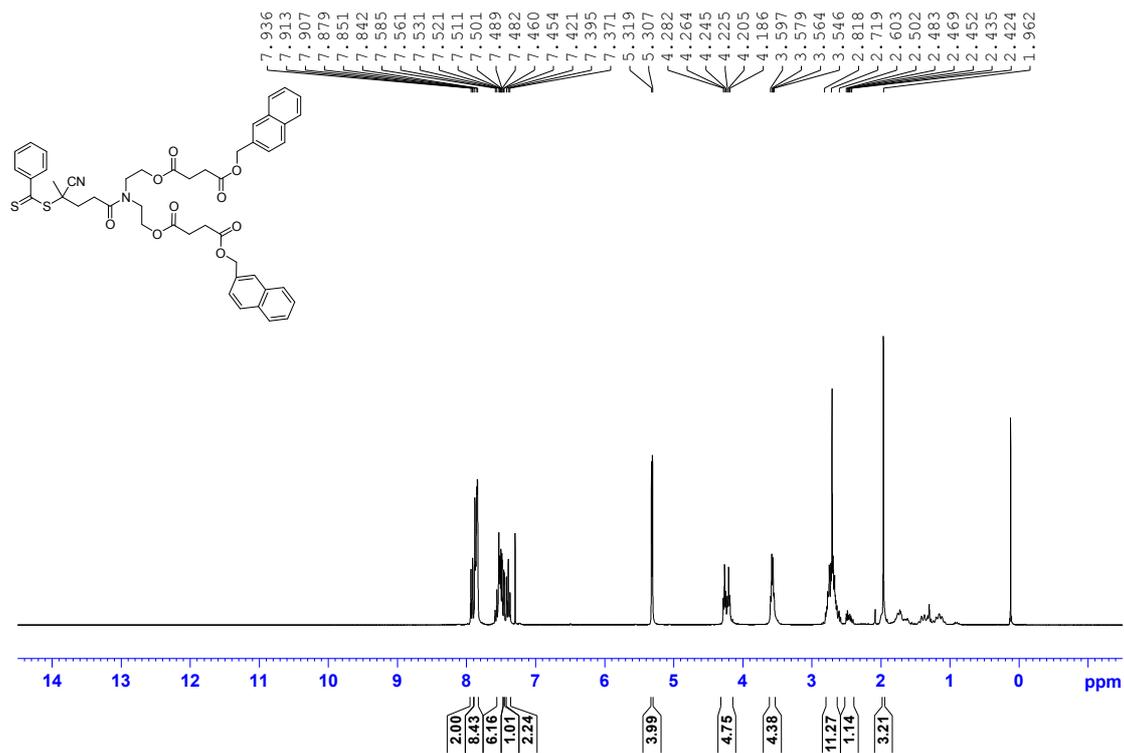
¹³C-NMR of CTA1



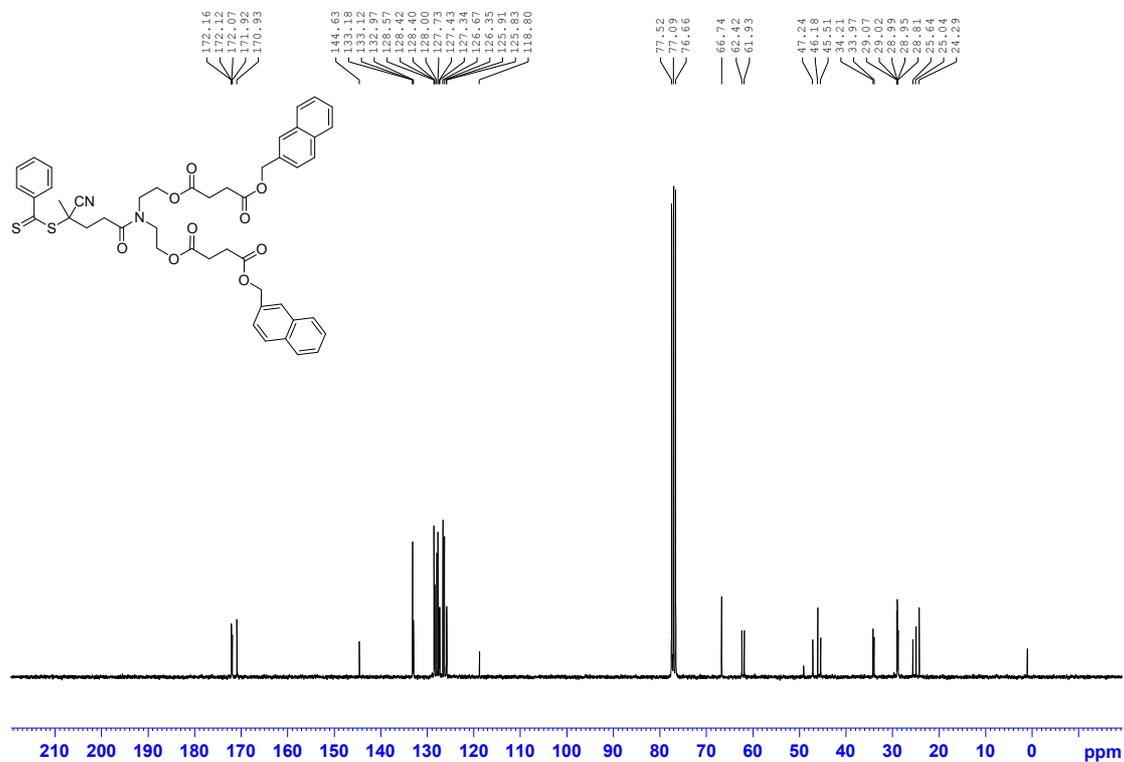
¹H-NMR of dibenzyl O,O'-(((4-cyano-4-(phenylcarbonothioyl)thio)pentanoyl)azanediy) bis(ethane-2,1-diyl) disuccinate (CTA2)



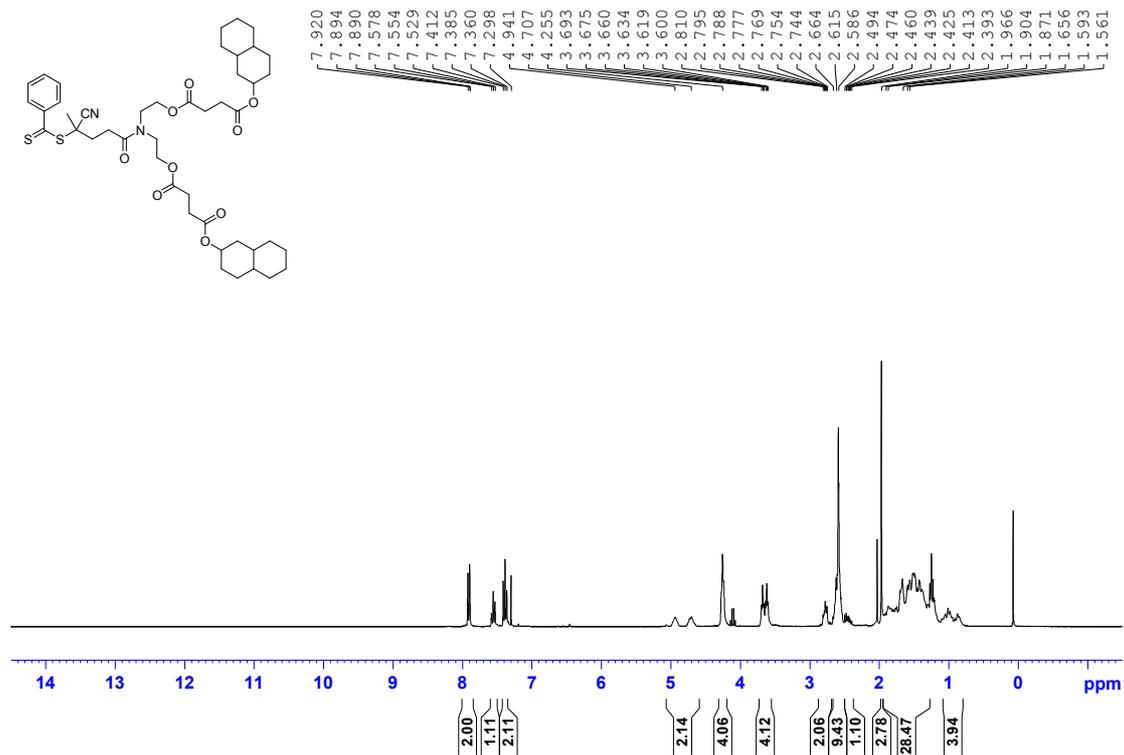
¹³C-NMR of CTA2



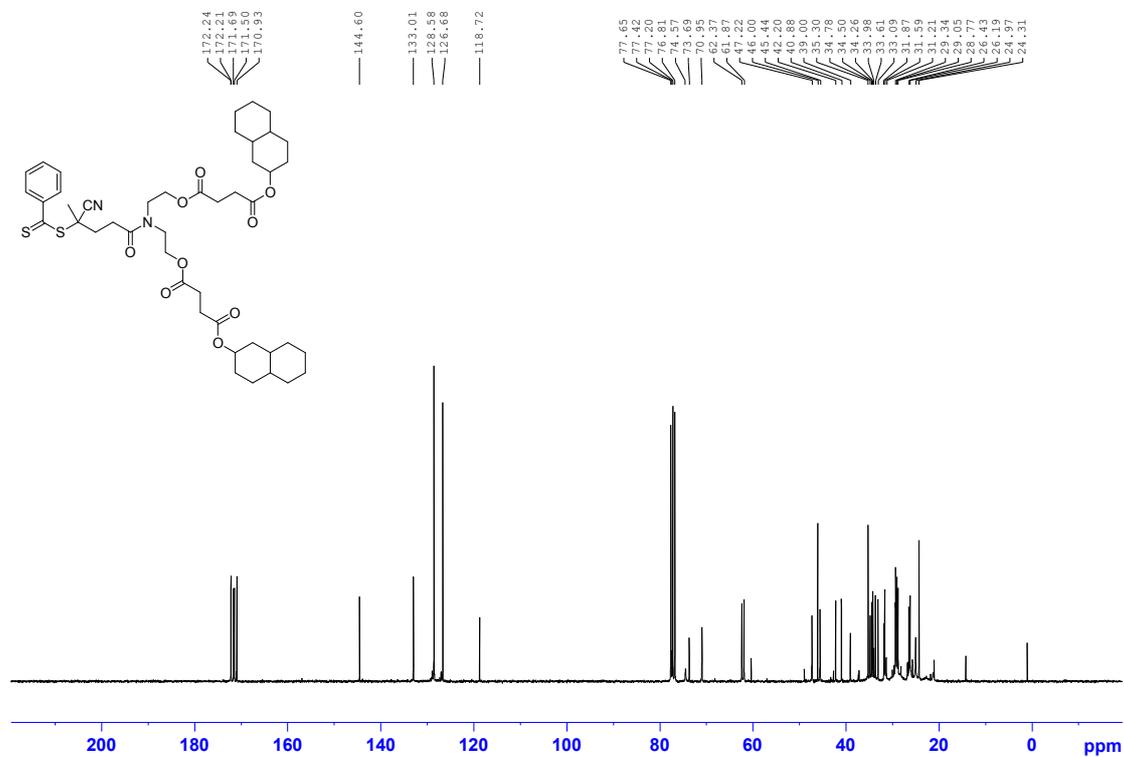
¹H-NMR of O,O'-(((4-cyano-4-((phenyl carbonothioyl)thio)pentanoyl)azanediyl)bis(ethane-2,1-diyl)bis(naphthalen-2-ylmethyl) disuccinate (CTA3)



¹³C-NMR of CTA3

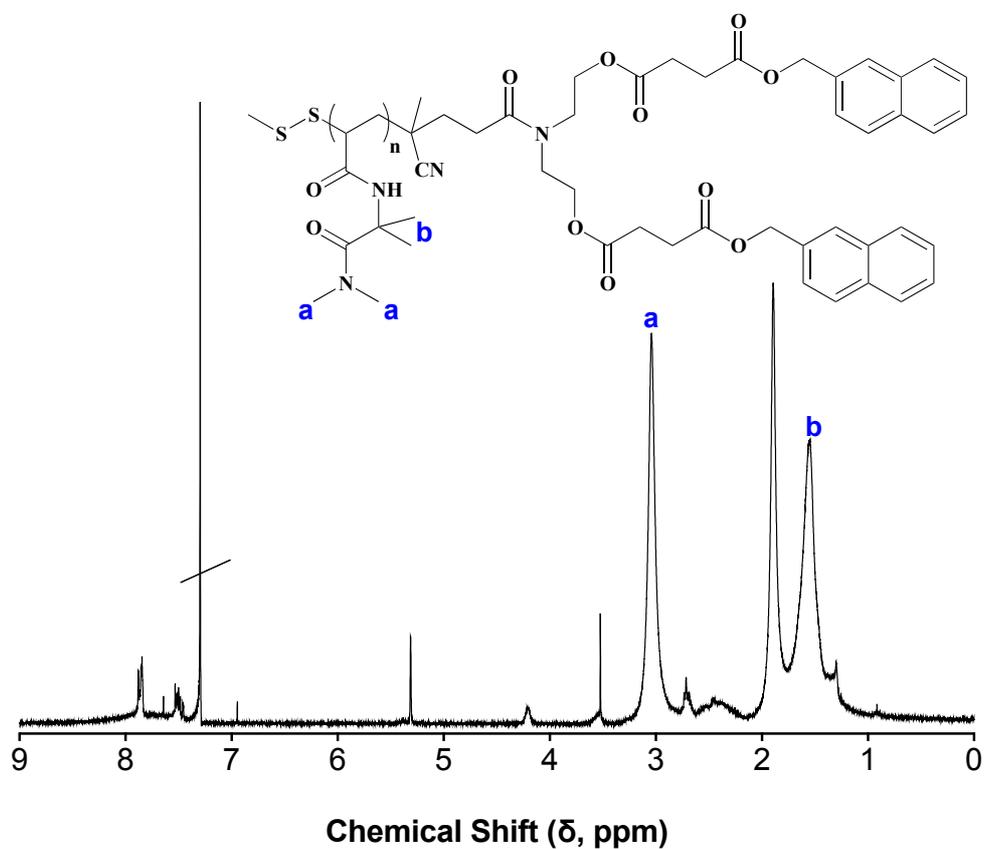


¹H-NMR of O,O'-(((4-cyano-4-((phenyl carbonothioyl)thio)pentanoyl)azanediy)bis(ethane-2,1-diyl)) bis(decahydro naphthalen-2-yl) disuccinate (**CTA4**)

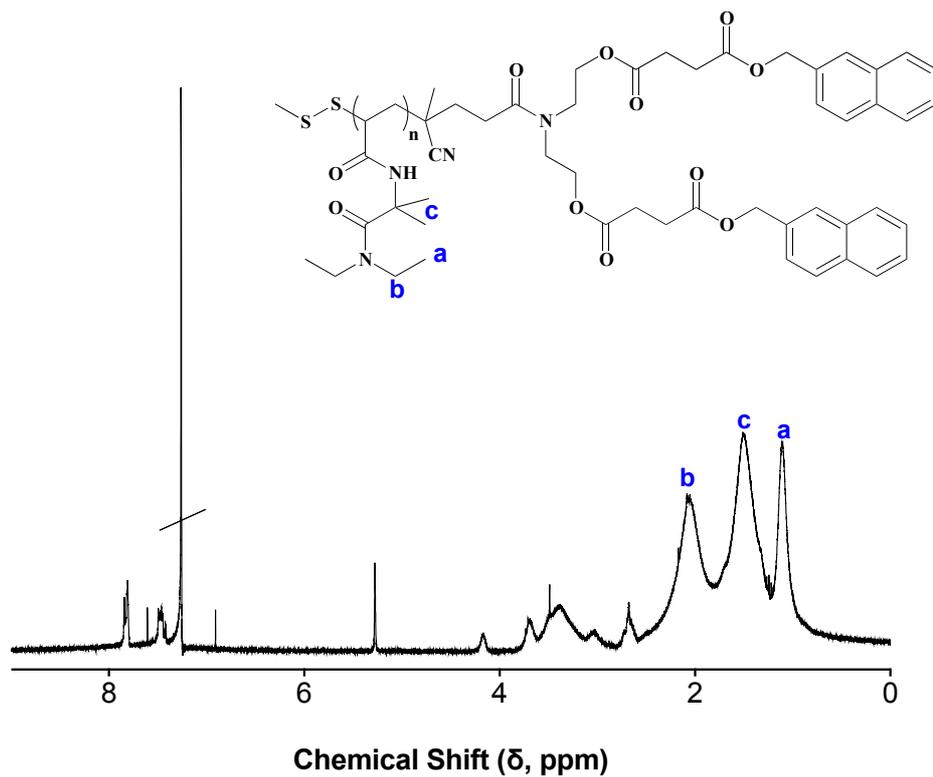


¹³C-NMR of CTA4

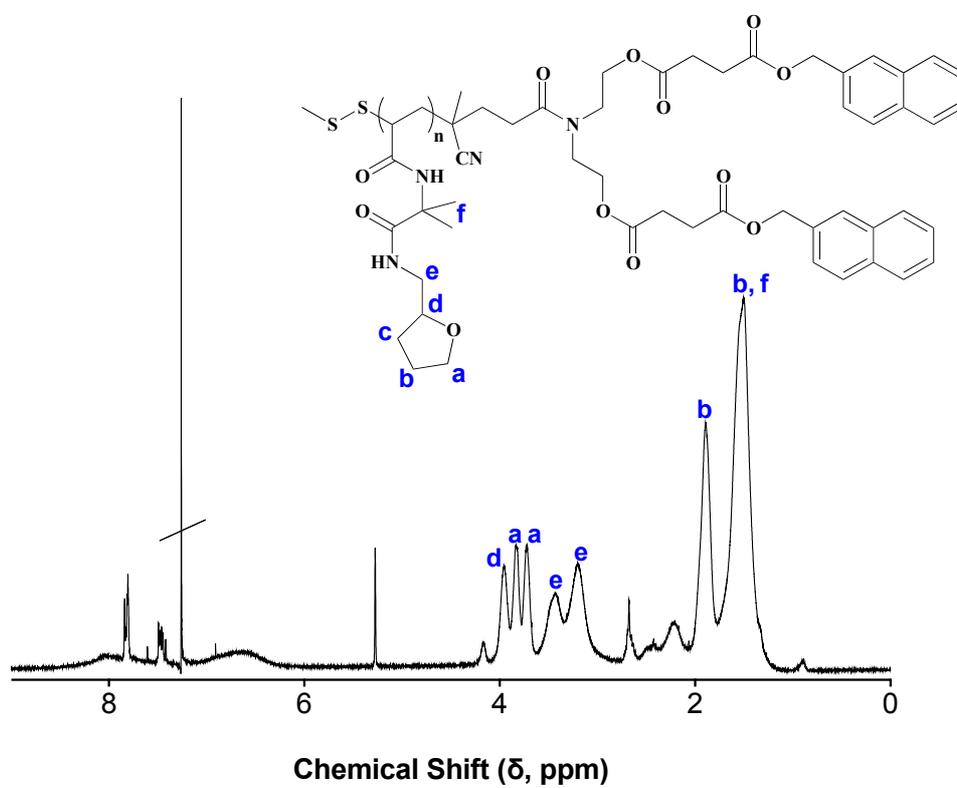
¹H-NMR of post-modified bis- α -naphthalenemethyl end-functional species



¹H-NMR bis- α -naphthalenemethyl dimethylamine polymer

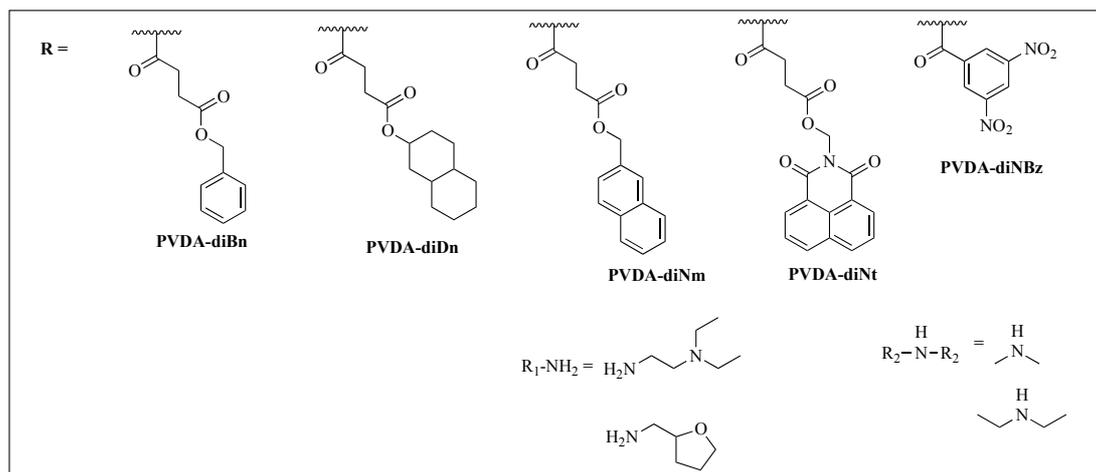
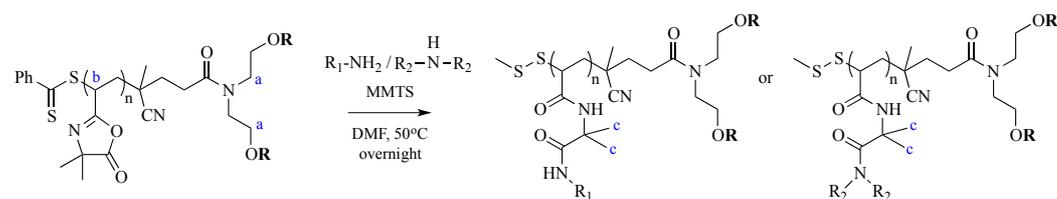


$^1\text{H-NMR}$ bis- α -naphthalenemethyl diethylamine polymer



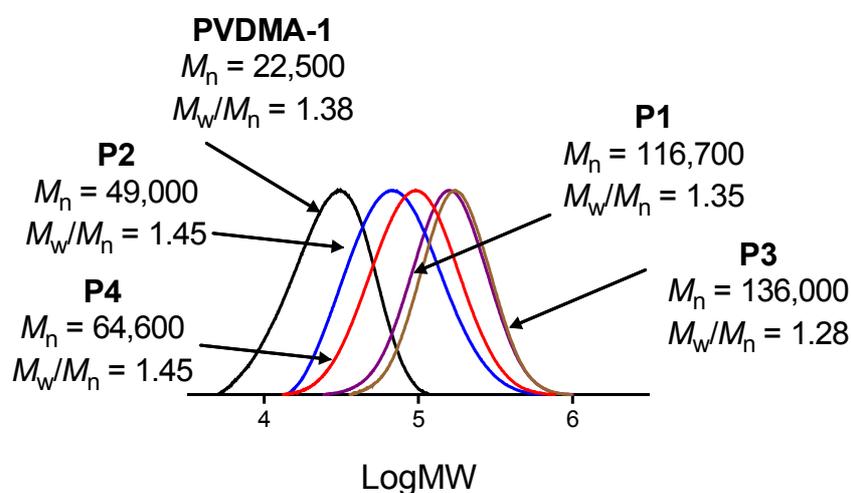
$^1\text{H-NMR}$ bis- α -naphthalenemethyl tetrahydrofurfylamine polymer

Determination of absolute molecular weight by $^1\text{H-NMR}$ results

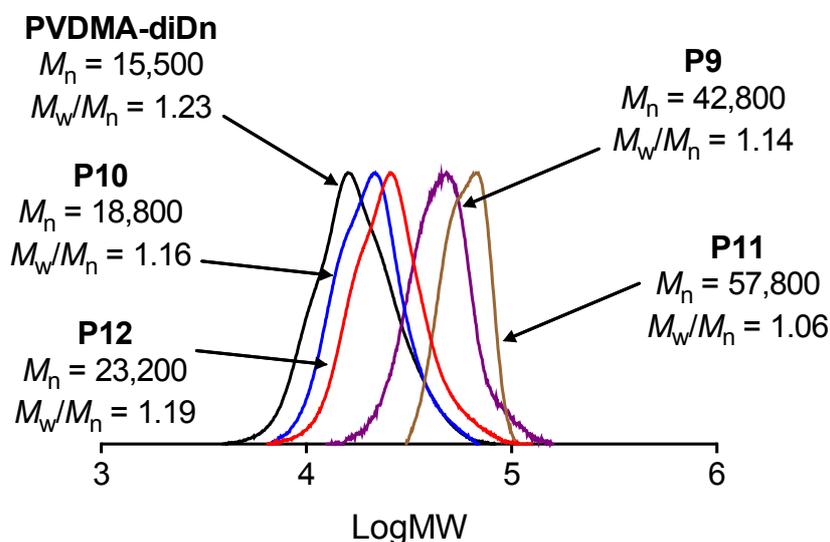


General chemical structure of VDMA homopolymer with protons used for absolute molecular weight determination labelled where (a) is the protons present in the alkyl chains of the RAFT agent ($\sim 4.2\text{ppm}$) and (b) is the proton present in the polymer backbone ($\sim 2.8\text{ppm}$). In the case of the post-modified derivatives, the integration signals for the labelled protons (c) present in the 2-methylalanyl spacer ($\sim 1.5\text{ppm}$) were compared with signals assigned to various introduced side groups suggest the successful and quantitative conversion of VDMA homopolymers to the respective post-modified bisamide derivatives.

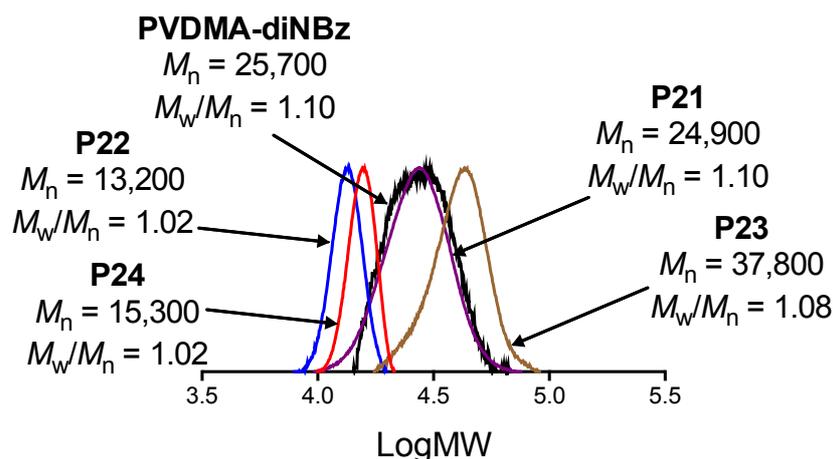
SEC results for various VDMA polymer and its post-modified derivatives



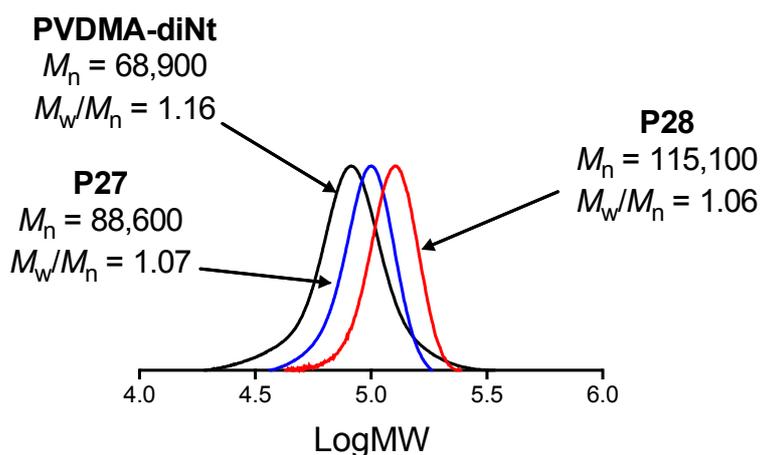
Molecular weight distribution curves for PVDMA-1 (black line), P1 DMA-modified homopolymer (purple line), P2 DEEDA-modified homopolymer (blue line), P3 DEA-modified homopolymer (green line) and P4 THFA-modified homopolymer (red line). SEC determined number average molecular weights (\bar{M}_n)



Molecular weight distribution curves for PVDMA-diDn (black line), P9 DMA-modified homopolymer (purple line), P10 DEEDA-modified homopolymer (blue line), P11 DEA-modified homopolymer (green line) and P12 THFA-modified homopolymer (red line). SEC determined number average molecular weights (\bar{M}_n) and dispersities ($D_M = \bar{M}_w / \bar{M}_n$) are also given.



Molecular weight distribution curves for PVDMA-diNBz (black line), P21 DMA-modified homopolymer (purple line), P22 DEEDA-modified homopolymer (blue line), P23 DEA-modified homopolymer (green line) and P24 THFA-modified homopolymer (red line). SEC determined number average molecular weights (\bar{M}_n) and dispersities ($D_M = \bar{M}_w / \bar{M}_n$) are also given.



Molecular weight distribution curves for PVDMA-diNt (black line), P27 DEEDA-modified homopolymer (blue line) and P28 THFA-modified homopolymer (red line). SEC determined number average molecular weights (\bar{M}_n) and dispersities ($D_M = \bar{M}_w / \bar{M}_n$) are also given.