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APPENDICES

(a)

APPENDIX A

A.1 Calculation of molar absorptivity of polymer

$$X \text{ ppm} = \underbrace{X \times 10^{-3}}_{\text{Molecular weight}} \text{ moles of polymer/1000 mL}$$

$$= \underbrace{X \times 10^{-3} \times (\text{Repeating unit})}_{\text{Molecular weight}} \text{ moles of monomer/1000 mL}$$

A.2 Calculation of percent penetration of octyl methoxycinnamate

$$A = \varepsilon bc$$

Where A is absorbance

b is the cell path length (1 cm)

c is the concentration of the absorbing species in mol per litre

Molar absorptivity (ϵ) of octyl methoxycinnmate are 23.000 M⁻¹cm⁻¹.

$$c_{OMC} = \underline{A}$$

$$23000$$

$$= \mathbf{X} \quad \text{mole/ litre}$$

Receptor volume is 13 mL, and molecular weight of OMC is 290.4:

Weight of Penetrated OMC = $X \times \frac{13}{1000} \times 290.4$

Weight of initial OMC is 0.005 g:

Percent penetration = $\frac{\text{Weight of penetrated OMC}}{\text{Weight of initial OMC}} \times 100$



Figure B.1 ¹H-NMR (CDCl₃) spectrum of p-(2-hydroxy ethoxy) cinnamic acid (M-2)



Figure B.2 ¹³C-NMR (DMSO- d_6) spectrum of p-(2-hydroxy ethoxy) cinnamic acid (M-2)



Figure B.3 ¹H-NMR (DMSO- d_6) spectrum of p-(2-hydroxy ethoxy) cinnamic acid (M-2)



Figure B.4 IR spectrum of *p*-(2-hydroxy ethoxy) cinnamic acid (M-2)







Figure B.7 ¹³C-NMR (DMSO- d_6) spectrum of p-(3-hydroxy propoxy) cinnamic acid (M-3)



Figure B.8 ¹H-NMR (DMSO- d_6) spectrum of p-(3-hydroxy propoxy) cinnamic acid (M-3)



Figure B.9 IR spectrum of *p*-(3-hydroxy propoxy) cinnamic acid (M-3)







Figure B.12 ¹³C-NMR (DMSO- d_6) spectrum of p-(6-hydroxy hexyloxy) cinnamic acid (M-6)



Figure B.13 ¹H-NMR (DMSO- d_6) spectrum of p-(6-hydroxy hexyloxy) cinnamic acid (M-6)



Figure B.14 IR spectrum of *p*-(6-hydroxy hexyloxy) cinnamic acid (M-6)



Figure B.15 Mass spectrum of *p*-(6-hydroxy hexyloxy) cinnamic acid (M-6)





Figure B.17 ¹³C-NMR (DMSO- d_6) spectrum of p-(11-hydroxy undecyloxy) cinnamic acid (M-11)



Figure B.18 ¹H-NMR (DMSO- d_6) spectrum of p-(11-hydroxy undecyloxy) cinnamic acid (M-11)



Figure B.19 IR spectrum of *p*-(11-hydroxy undecyloxy) cinnamic acid (M-11)



Figure B.20 Mass spectrum of *p*-(11-hydroxy undecyloxy) cinnamic acid (M-11)





Figure B.22 IR spectrum of poly(p-ethoxy cinnamate), P-2





Figure B.24 IR spectrum of poly(*p*-propoxy cinnamate), P-3 (soluble part)



Figure B.25 IR spectrum of poly(p-propoxy cinnamate), P-3 (insoluble part)



Figure B.26¹H-NMR (CDCl₃) spectrum of poly(*p*-propoxy cinnamate), P-3 irradiated





Figure B.28 IR spectrum of poly(p-hexyloxy cinnamate), P-6



Figure B.29 ¹H-NMR (CDCl₃) spectrum of poly(*p*-hexyloxy cinnamate), P-6 irradiated





Figure B.31 IR spectrum of poly(*p*-undecyloxy cinnamate), P-11



Figure B.32 ¹H-NMR (CDCl₃) spectrum of poly(*p*-undecyloxy cinnamate), P-11 irradiated



Figure B.33 ¹H-NMR (CDCl₃) spectrum of poly(*p*-undecyloxy cinnamate), P-11dil



Figure B.34 IR spectrum of poly(*p*-undecyloxy cinnamate), P-11dil





Figure B.36 IR spectrum of poly(p-propoxy cinnamate)-co-(p-undecyloxy cinnamate), P-3/11



Figure B.37¹H-NMR (CDCl₃) spectrum of poly(*p*-propoxy cinnamate)-*co*-(*p*-undecyloxy cinnamate), P-3/11irradiated



Figure B.38 ¹H-NMR (CDCl₃) spectrum of poly(pentaethylene glycol cinnamate), PPGC



Figure B.39¹³C-NMR (CDCl₃) spectrum of poly(pentaethylene glycol cinnamate), PPGC



Figure B.40 IR spectrum of poly(pentaethylene glycol cinnamate), PPGC

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Figure B.41 ¹H-NMR (CDCl₃) spectrum of poly(pentaethylene glycol cinnamate), PPGC irradiated

VITA

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Experience

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