

Supporting Information

Synthesis, Characterization and Antimicrobial Activities of Water-Soluble Amphiphilic Copolymers Containing Ciprofloxacin and Quaternary Ammonium Salts

Man He,^{a,b} Huining Xiao,^{*a} Yuming Zhou^b and Peng Lu^a

^a Department of Chemical Engineering & Limerick Pulp and Paper
Centre, University of New Brunswick, Fredericton, NB E3B 5A3,
Canada

^b School of Chemistry and Chemical Engineering, Southeast University,
Nanjing 211189, China

Correspondence to: H. Xiao (E-mail: hxiao@unb.ca)

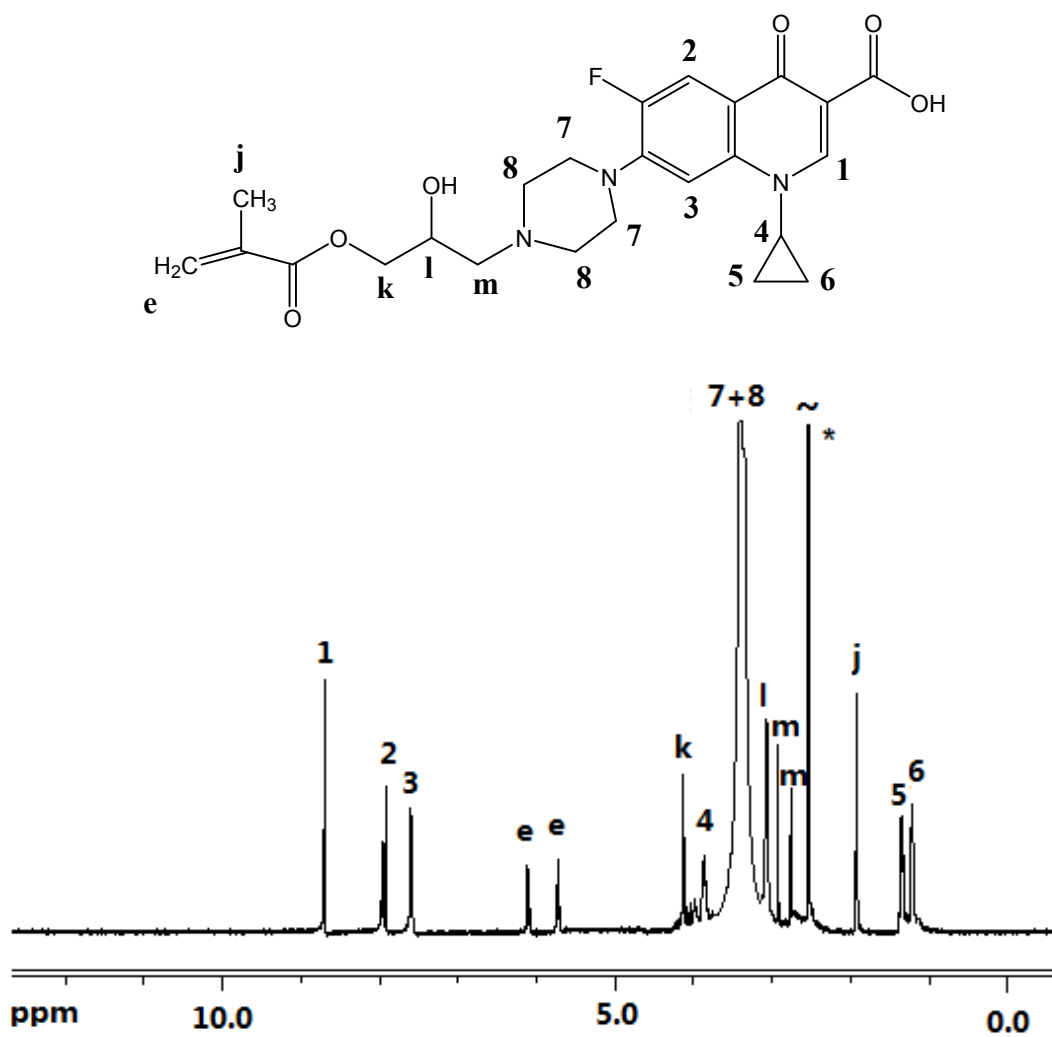


Figure S1. ¹H-NMR spectrum of GMA-CPF

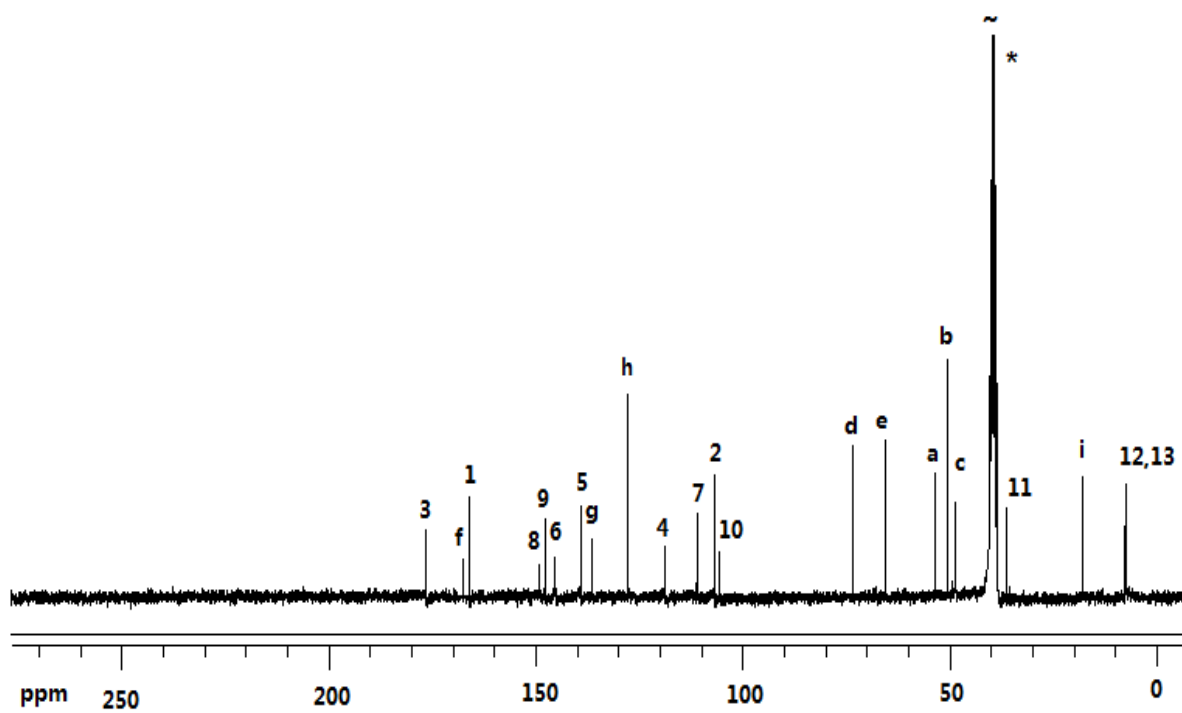
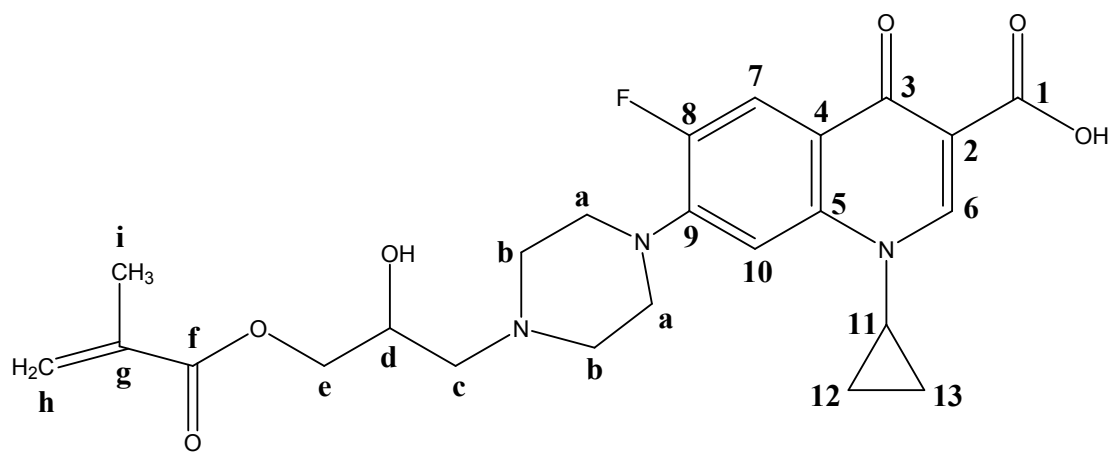


Figure S2. ¹³C-NMR spectrum of GMA-CPF

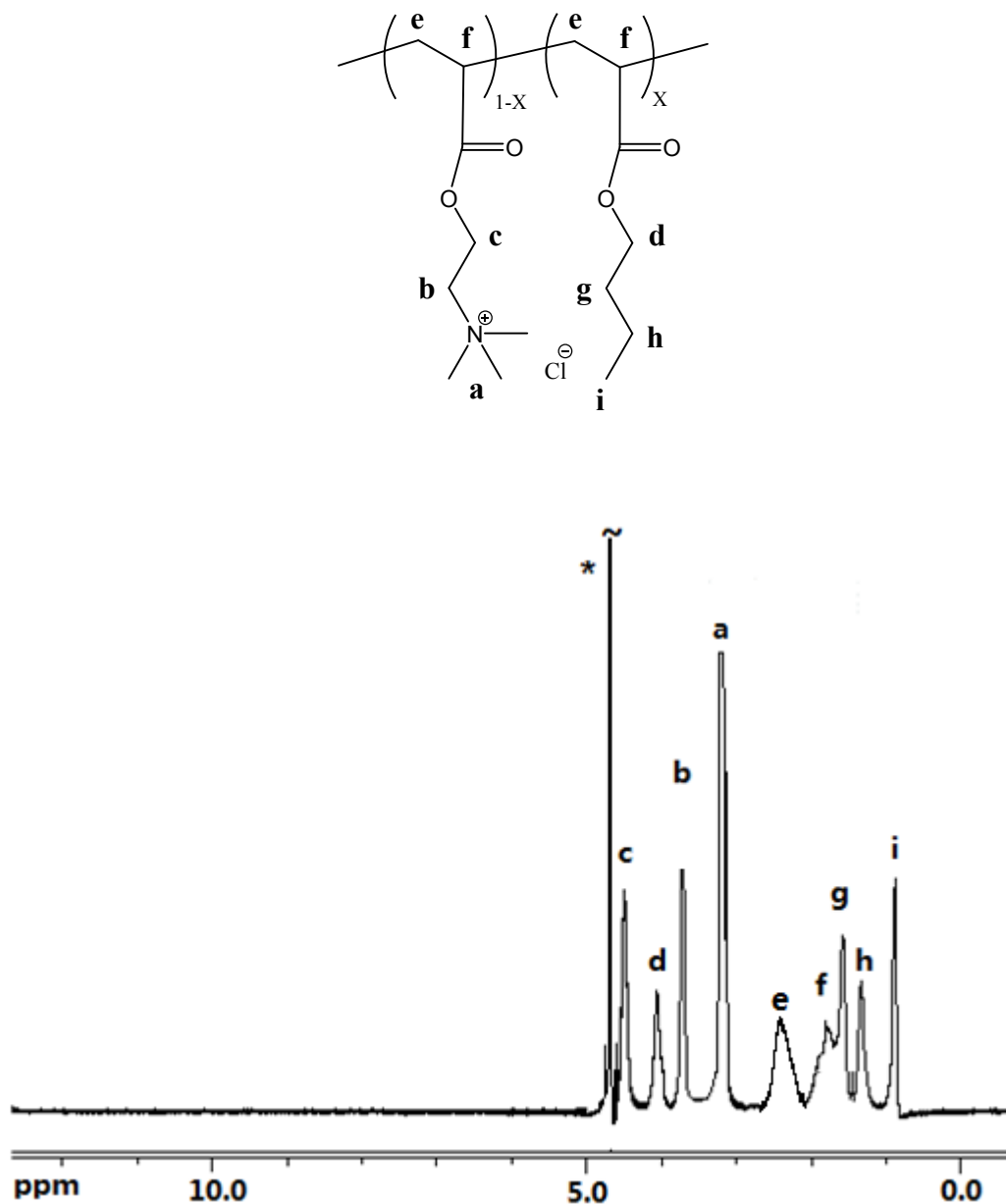


Figure S3. $^1\text{H-NMR}$ spectrum of DMAEAMC/BA copolymer; (1-x): x=59: 41

The copolymer compositions for this system were determined by $^1\text{H-NMR}$ spectroscopy analysis. The $^1\text{H-NMR}$ peak at 3.10 ppm (Figure S3, a) corresponds to the methyl group of DMAEAMC; the peak at 1.32 ppm (Figure S3, h) corresponds to the methylene group, $-\text{O}-(\text{CH}_2)_2\text{CH}_2\text{CH}_3$, in the BA side chain. Let m_1 be the mole fraction of DMAEAMC and $m_2 = (1-m_1)$ that of the BA unit.

$$\frac{\text{Area}_{3.10\text{ppm}}}{\text{Area}_{1.32\text{ppm}}} = \frac{9m_1}{2m_2} = C_1 \quad \text{Equation 1a}$$

Where $\text{Area}_{3.10\text{ppm}}$ =area of the peak at 3.10ppm and $\text{Area}_{1.32\text{ppm}}$ =area of the peak at 1.32ppm.

On simplification,

$$m_1 = \frac{2C_1}{2C_1 + 9} \quad \text{Equation 1b}$$

Where m_1 and m_2 are the copolymer molar compositions. The results are presented in Table 1, which are in agreement with the values calculated by apparent charge density.

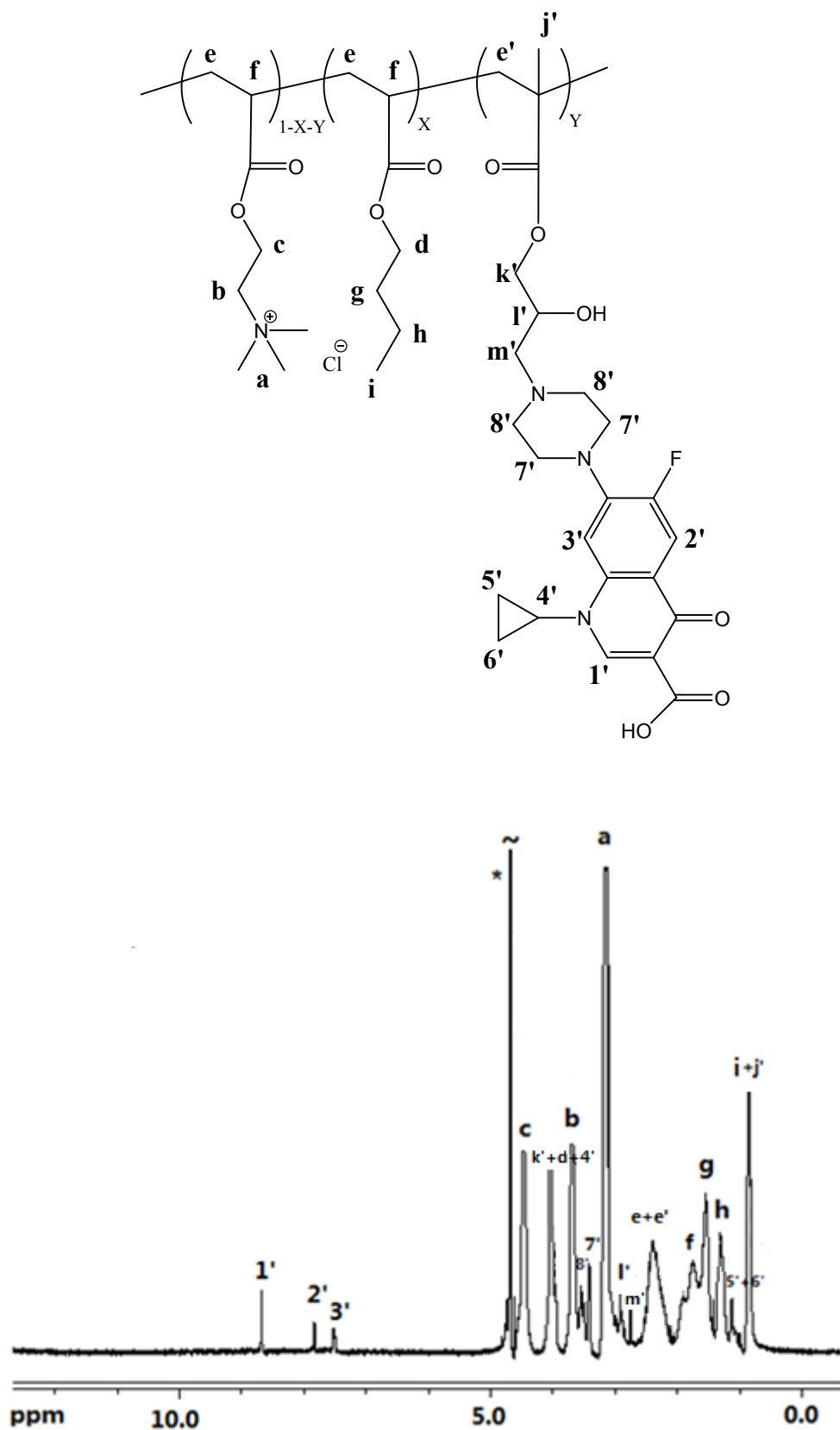


Figure S4. ¹H-NMR spectrum of DMAEAMC/BA/GMA-CPF copolymer; (1-x-y): x:
y=52.8: 42.9: 4.3
The copolymer compositions for this system were determined by ¹H-NMR

spectroscopy analysis. The $^1\text{H-NMR}$ peak at 3.10 ppm (Figure S4, a) corresponds to the methyl group of DMAEAMC; the peak at 1.32 ppm (Figure S4, h) corresponds to the methylene group, $-\text{O}-(\text{CH}_2)_2\text{CH}_2\text{CH}_3$, in the BA side chain; the peaks at 0.85 ppm (Figure S4, i+j') are the methyl group for BA and GMA-CPF. Considering the mole fraction of DMAEAMC as m_1 , the mole fraction of BA as m_2 , and $m_3 = (1-m_1-m_2)$ that of the GMA-CPF unit.

$$\frac{\text{Area}_{3.10\text{ppm}}}{\text{Area}_{1.32\text{ppm}}} = \frac{9m_1}{2m_2} = C_1 \quad \text{Equation 2a}$$

$$\frac{\text{Area}_{1.32\text{ppm}}}{\text{Area}_{0.85\text{ppm}}} = \frac{2m_2}{3m_2 + 3m_3} = C_2 \quad \text{Equation 2b}$$

Where $\text{Area}_{3.10\text{ppm}}$ =area of the peak at 3.10ppm, $\text{Area}_{1.32\text{ppm}}$ =area of the peak at 1.32ppm, and $\text{Area}_{0.85\text{ppm}}$ =area of the peak at 0.85ppm.

On simplification,

$$m_1 = \frac{C_1 C_2}{C_1 C_2 + 3} \quad \text{Equation 2c}$$

$$m_2 = \frac{9C_1}{2C_1 C_2 + 6} \quad \text{Equation 2d}$$

$$m_3 = \frac{6 - 9C_1}{2C_1 C_2 + 6} \quad \text{Equation 2e}$$

Where m_1 , m_2 , and m_3 are the copolymer molar compositions. The results are presented in Table 1, which are in agreement with the values calculated by apparent charge density and UV measurements.