

SUPPLEMENTARY INFORMATION

Surfactant-free polymeric nanoparticles composed of PEG, cholic acid and a sucrose moiety

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1. MALDI TOF mass spectra of the Suc-PEG-Chol conjugates 10 and 11

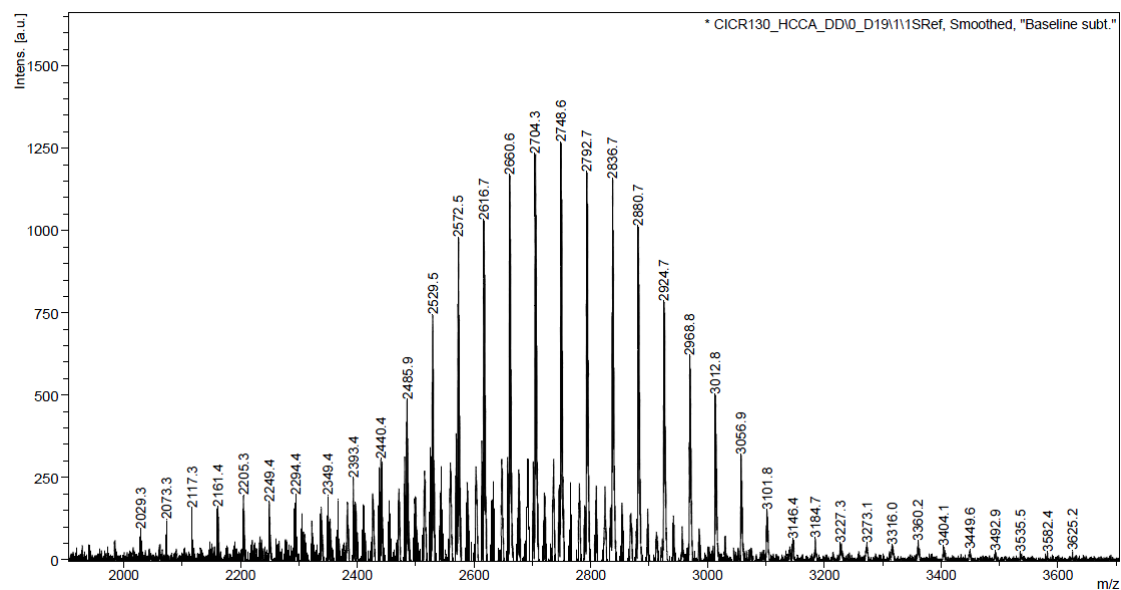


Fig. S 1 - Matrix-assisted laser desorption ionization time-of-flight (MALDI-TOF) of the Suc-PEG₂₀₀₀-Chol (**10**) on a Bruker Autoflex with α -cyano-4-hydroxycinnamic acid (4-HCCA) matrix.

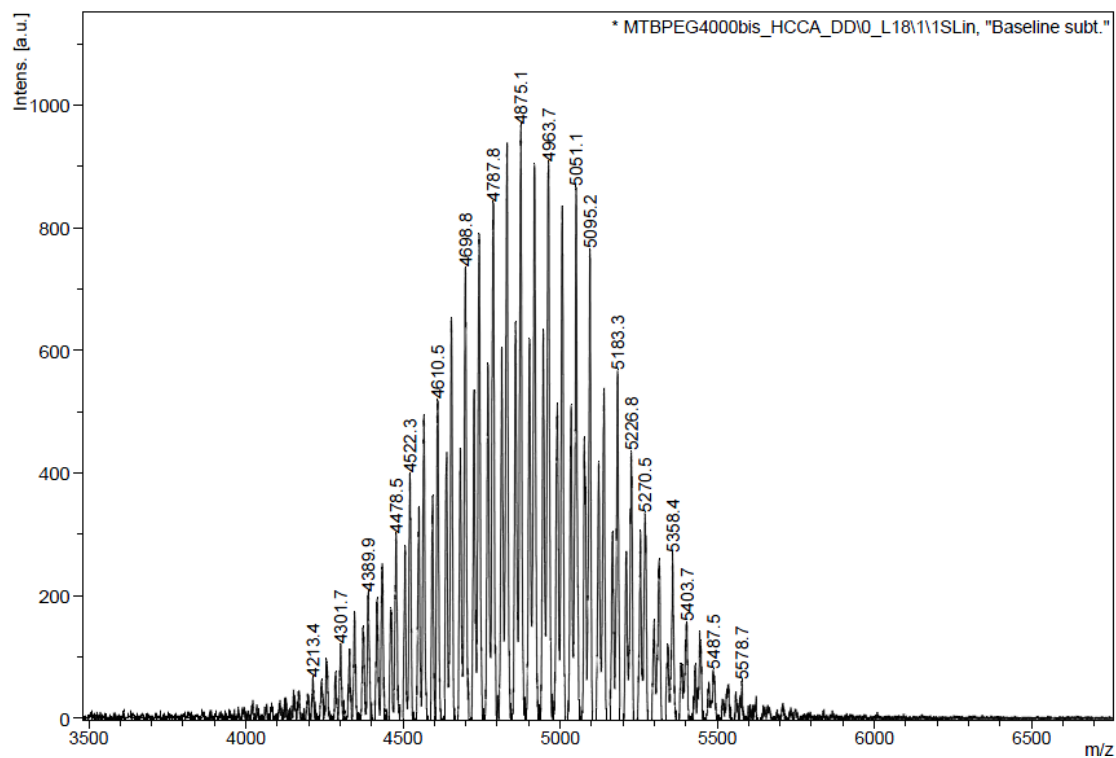


Fig. S 2 - Matrix-assisted laser desorption ionization time-of-flight (MALDI-TOF) of the Suc-PEG₄₀₀₀-Chol (**11**) on a Bruker Autoflex with α -cyano-4-hydroxycinnamic acid (4-HCCA) matrix.

2. DSC of the Suc-PEG-Chol conjugates 10 and 11

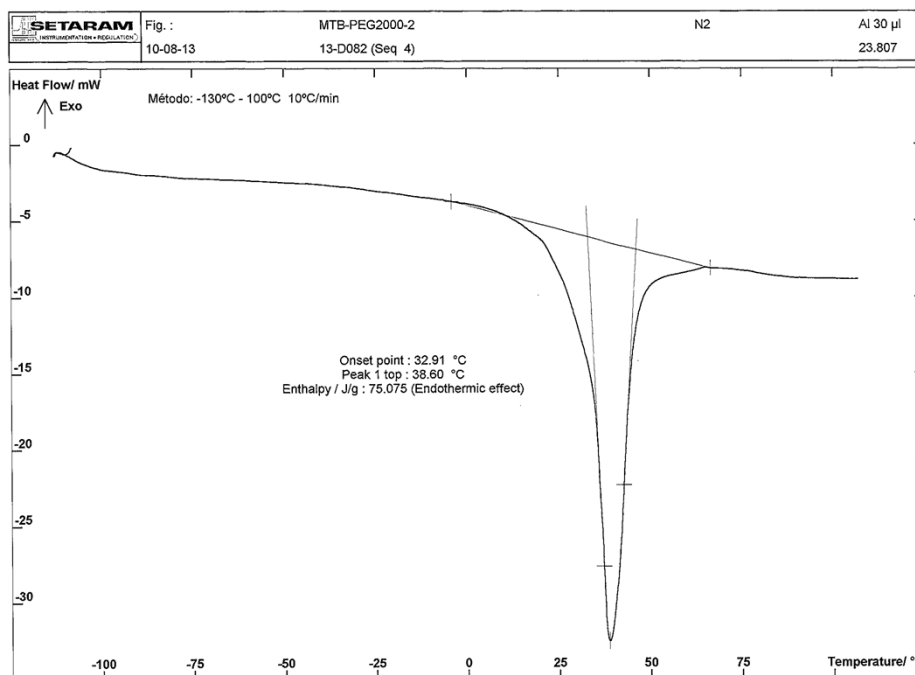


Fig. S 3 – Differential scanning calorimetry (DSC) of Suc-PEG₂₀₀₀-Chol (10).

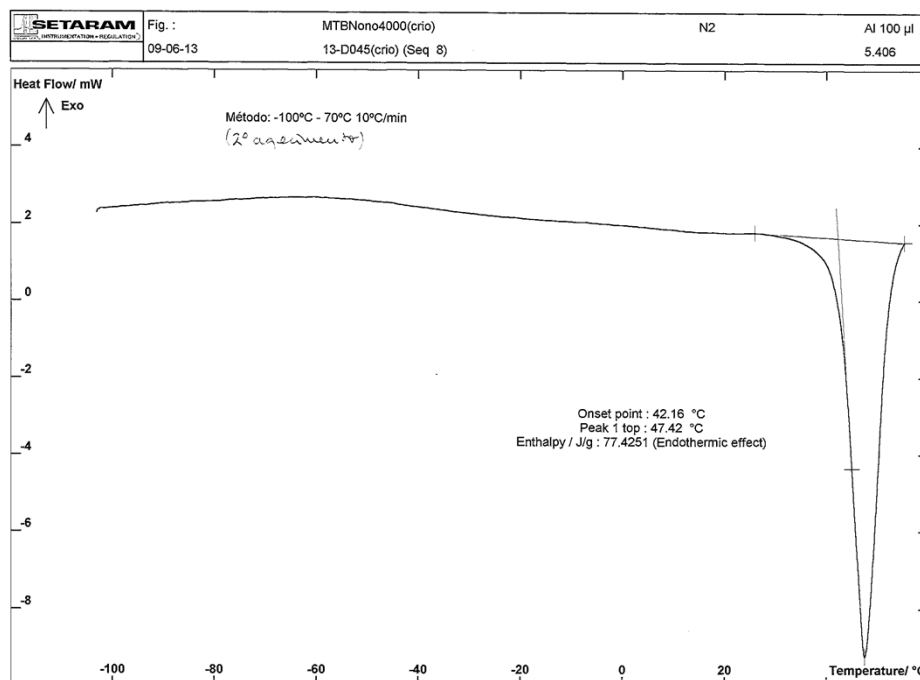


Fig. S 4 – Differential scanning calorimetry (DSC) of Suc-PEG₄₀₀₀-Chol (11).

3. Determination of critical aggregation concentration (CAC) for Suc-PEG₄₀₀₀-Chol conjugate

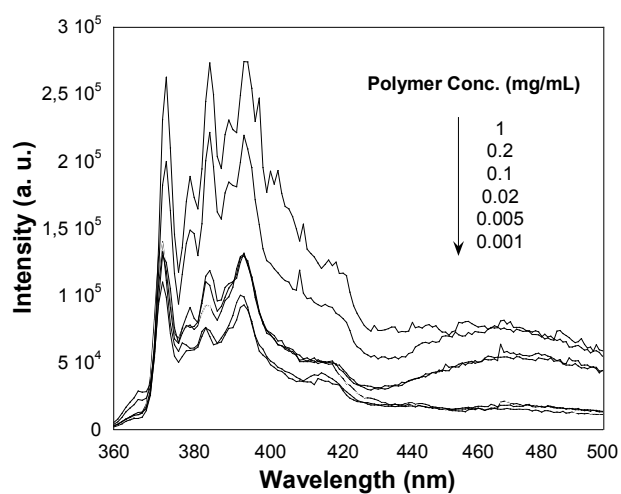


Fig. S 5 - Fluorescence emission spectra of pyrene/Suc-PEG₄₀₀₀-Chol against concentration of Suc-PEG₂₀₀₀-Chol in distilled water.

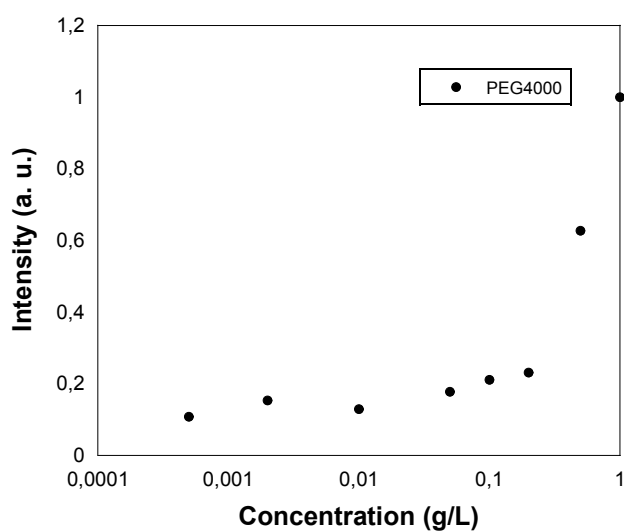


Fig. S 6 - Plots of the area of the pyrene emissions spectra vs. Suc-PEG₄₀₀₀-Chol polymer concentration.

4. Particle size distributions of Suc-PEG-Chol nanoparticles

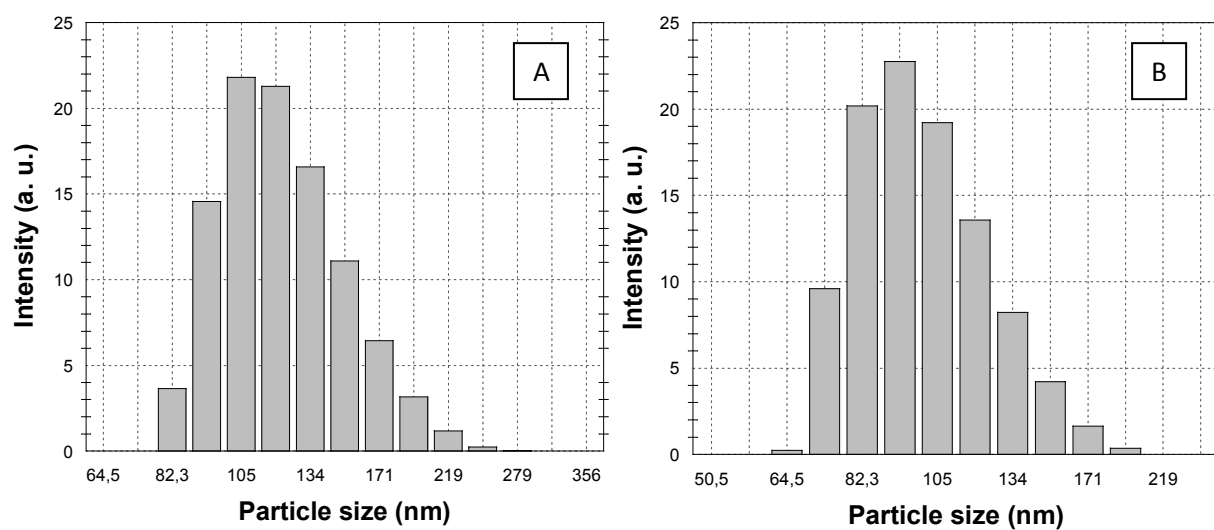


Fig. S 7 – Particle size distributions of Suc-PEG-Chol nanoparticles. (A) Suc-PEG₂₀₀₀-Chol, (B) Suc-PEG₄₀₀₀-Chol.

5. AFM images of Suc-PEG-Chol nanoparticles

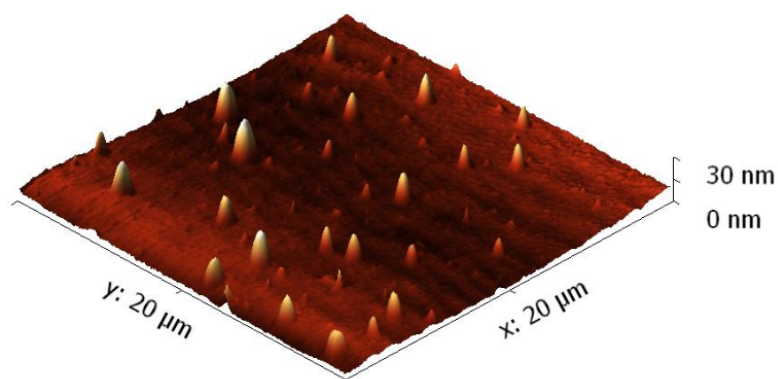


Fig. S 8 - AFM height image of Suc-PEG₄₀₀₀-Chol nanoparticles from a PNPs solution of 0.1mg/mL.

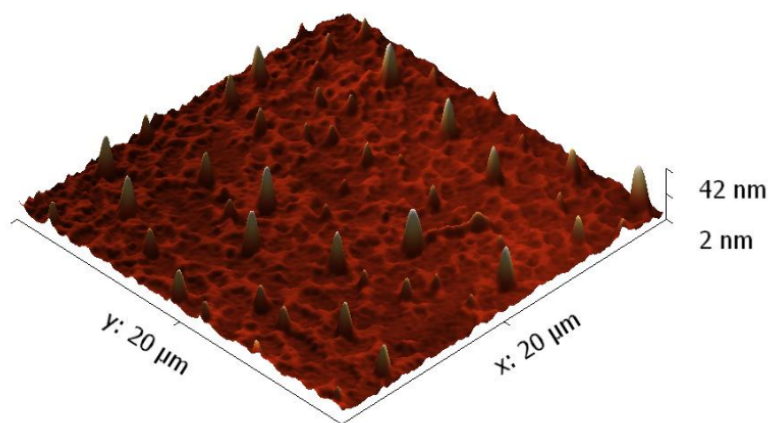


Fig. S 9 - AFM height image of Suc-PEG₂₀₀₀-Chol nanoparticles from a PNPs solution of 0.1mg/mL.

6. ^1H -NMR spectrum of Suc-PEG₄₀₀₀-Chol (11) in D_2O

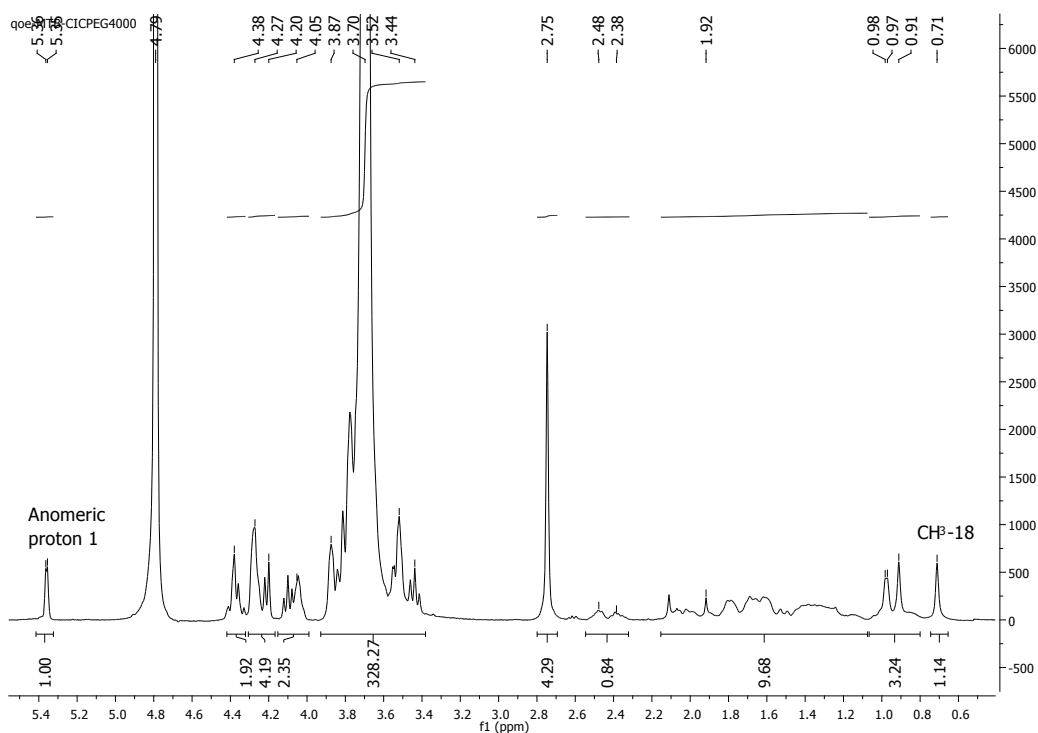


Fig. S 10 - ^1H -NMR spectrum of Suc-PEG₄₀₀₀-Chol (11) in D_2O .

7. ^1H -NMR spectrum of 1',2,3,3',4,4',6-Hepta-O-benzyl-6'-O-succinyl-sucrose (2) in CDCl_3

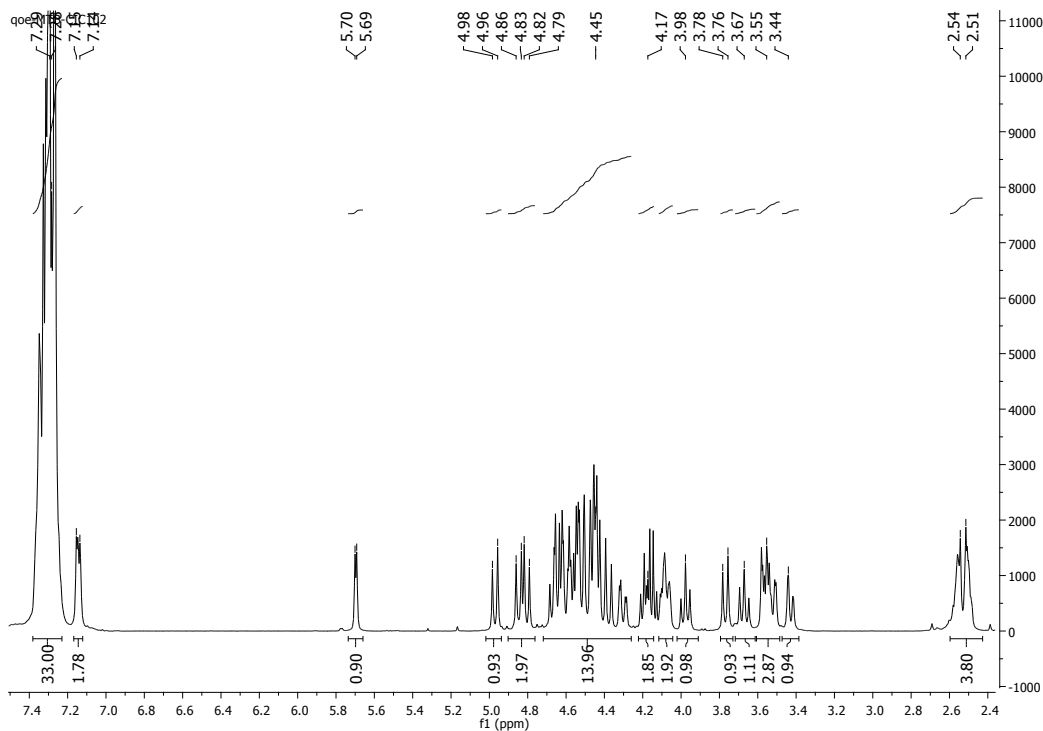


Fig. S 11 - ^1H -NMR spectrum of 1',2,3,3',4,4',6-Hepta-O-benzyl-6'-O-succinyl-sucrose (2) in CDCl_3 .

8. ^1H -NMR spectrum of Methyl 3 α ,7 α ,12 α -trihydroxy-5 β -cholan-24-ate (3) in CDCl_3

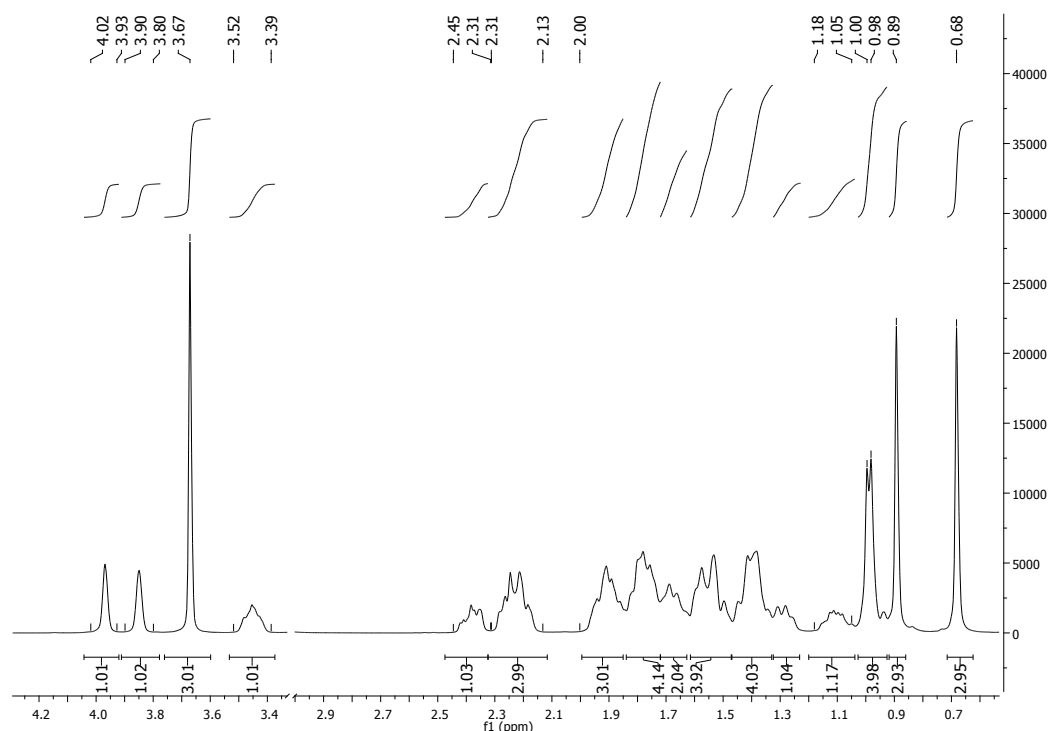


Fig. S 12 - ^1H -NMR spectrum of Methyl 3 α ,7 α ,12 α -trihydroxy-5 β -cholan-24-ate (3) in CDCl_3 .

9. ^1H -NMR spectrum of Methyl 3 α -O-benzyl, 7 α , 12 α -dihydroxy-5 β -cholan-24-oate (4) in CDCl_3

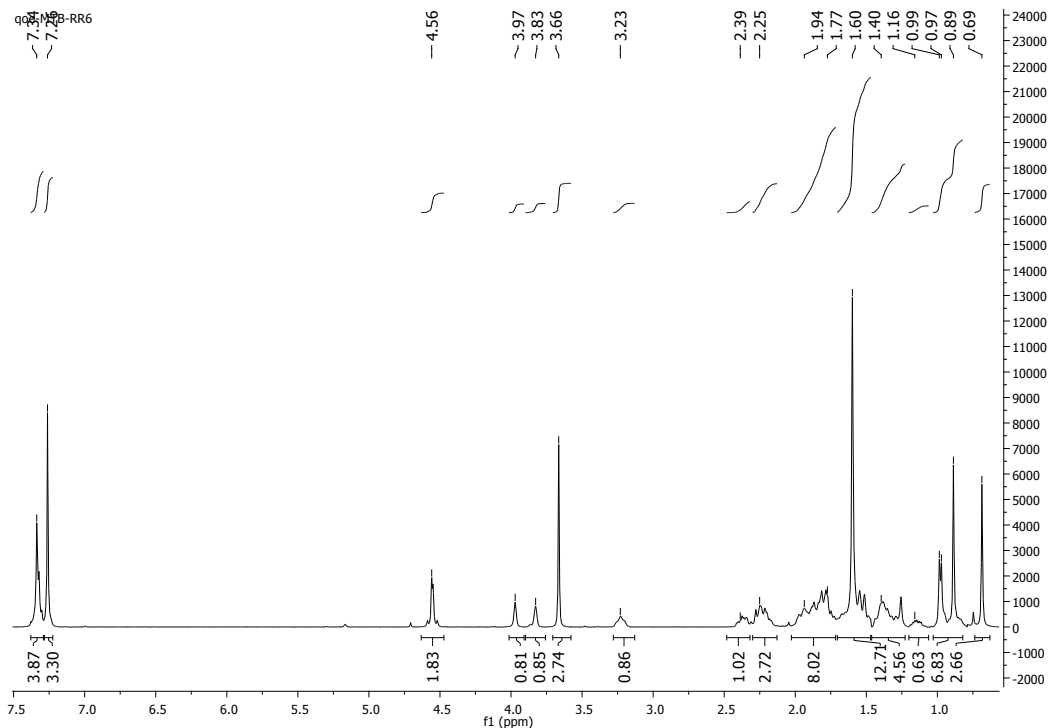


Fig. S 13 - ^1H -NMR spectrum of Methyl 3 α -O-benzyl, 7 α , 12 α -dihydroxy-5 β -cholan-24-oate (4) in CDCl_3 .

10. ^1H -NMR spectrum of 3 α -O-benzyl, 7 α , 12 α -dihydroxy-5 β -cholic acid (5) in CDCl_3

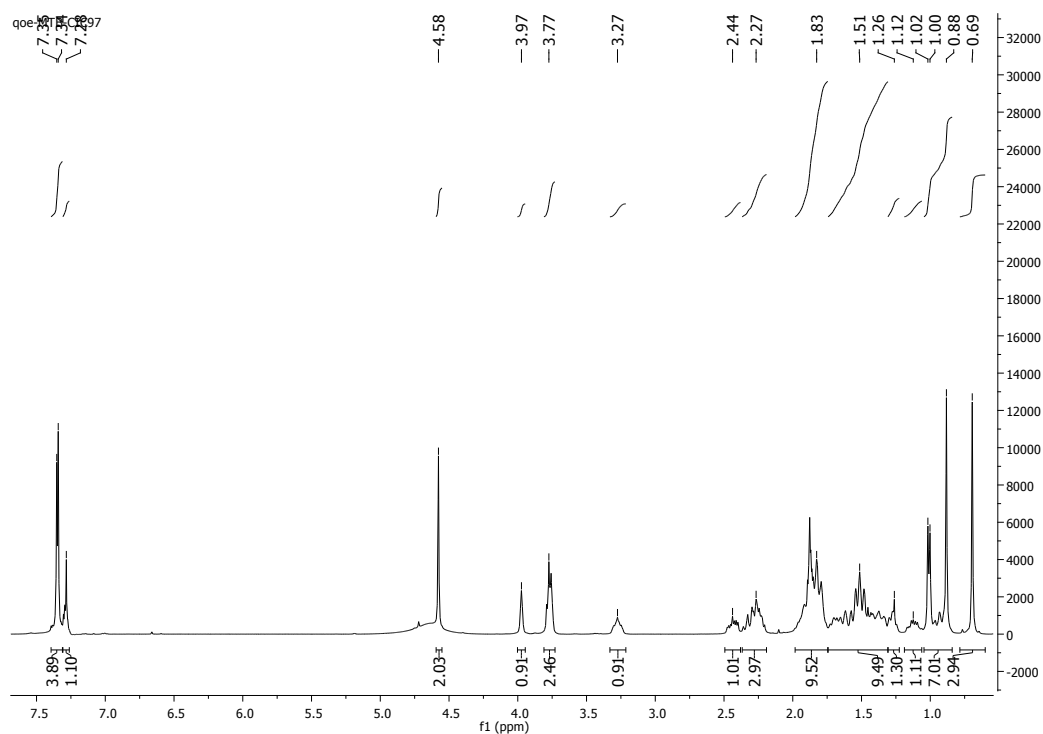


Fig. S 14 - ^1H -NMR spectrum of 3 α -O-benzyl, 7 α , 12 α -dihydroxy-5 β -cholic acid (5) in CDCl_3 .

11. HMQC spectrum of 3 α -O-benzyl, 7 α , 12 α -dihydroxy-5 β -cholic acid (5) in CDCl_3

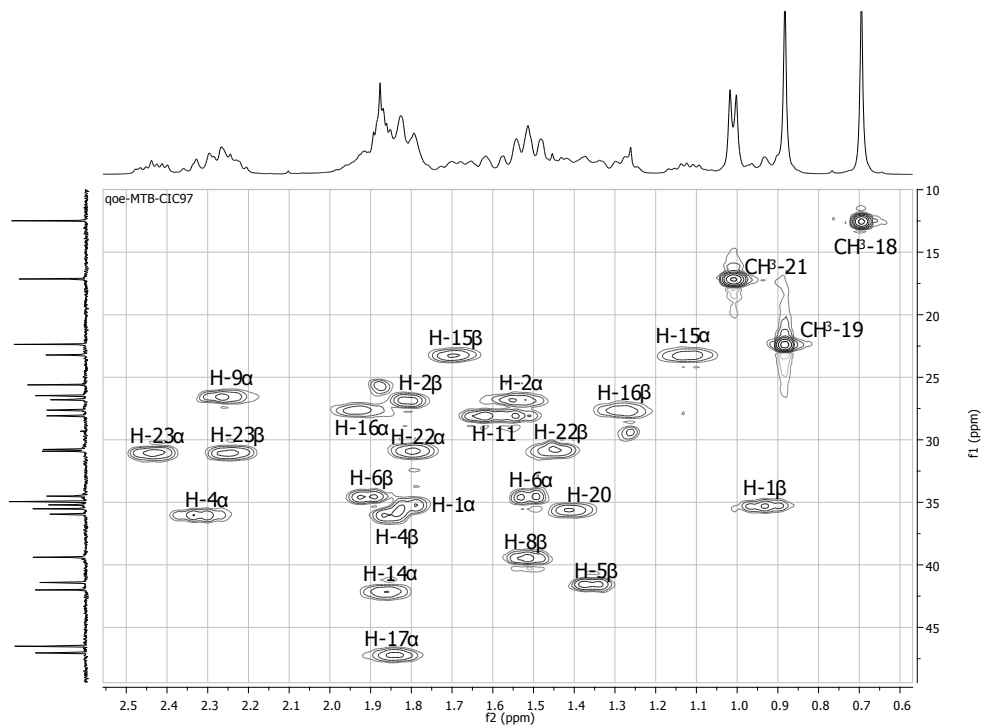


Fig. S 15 - HMQC spectrum of 3 α -O-benzyl, 7 α , 12 α -dihydroxy-5 β -cholic acid (5) in CDCl_3 .

12. ^1H -NMR spectrum of Benzylated Suc-PEG₂₀₀₀-OH (6) in CDCl_3

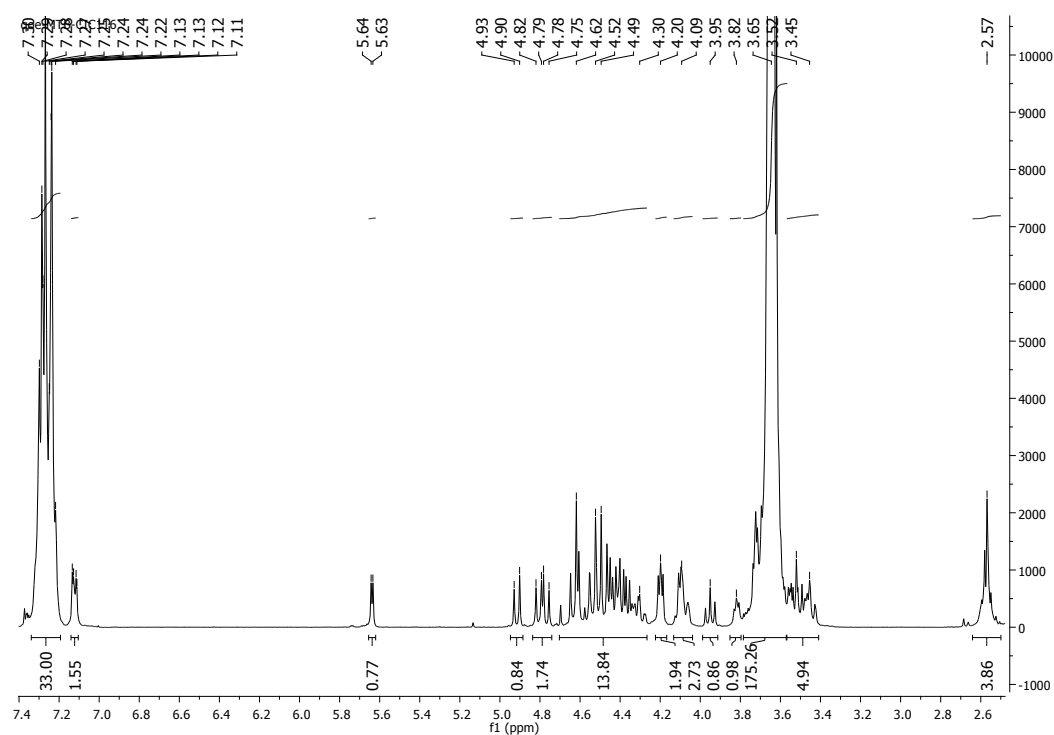


Fig. S 16 - ^1H -NMR spectrum of Benzylated Suc-PEG₂₀₀₀-OH (6) in CDCl_3 .

13. ^1H -NMR spectrum of Benzylated Suc-PEG₄₀₀₀-OH (7) in CDCl_3

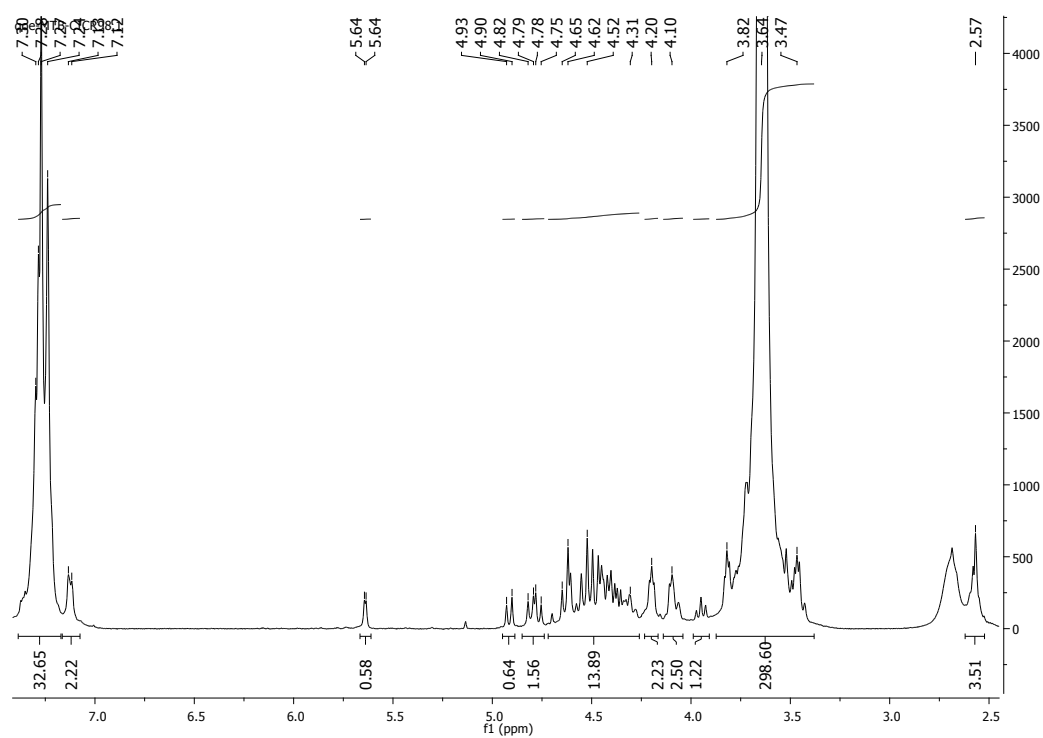


Fig. S 17 - ^1H -NMR spectrum of Benzylated Suc-PEG₄₀₀₀-OH (7) in CDCl_3 .

14. ^1H -NMR spectrum of Benzylated Suc-PEG₂₀₀₀-Chol (8) in CDCl_3

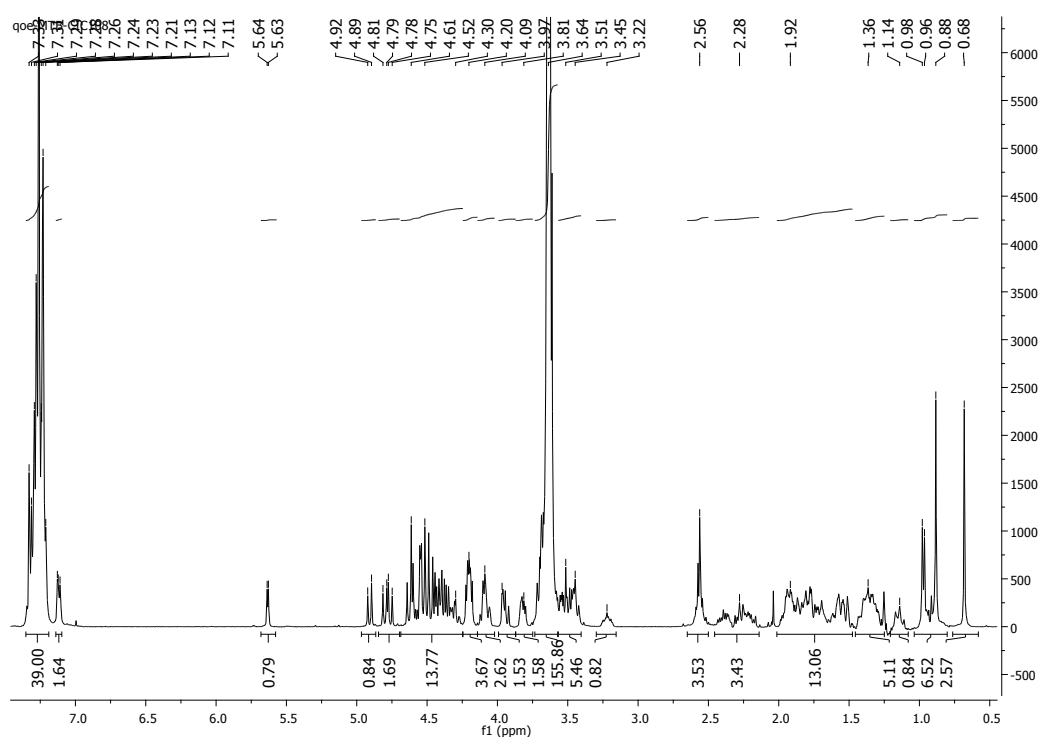


Fig. S 18 - ^1H -NMR spectrum of Benzylated Suc-PEG₂₀₀₀-Chol (8) in CDCl_3 .

15. ^1H -NMR spectrum of Benzylated Suc-PEG₄₀₀₀-Chol (9) in CDCl_3

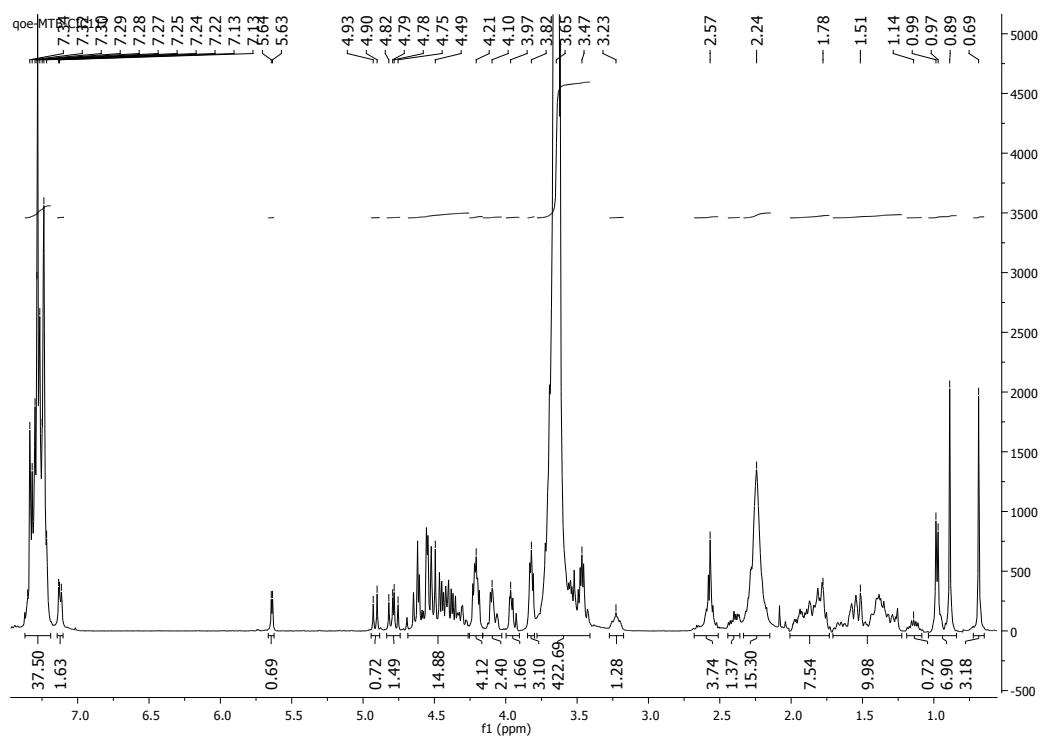


Fig. S 19 - ^1H -NMR spectrum of Benzylated Suc-PEG₄₀₀₀-Chol (9) in CDCl_3 .

16. ^1H -NMR spectrum of Suc-PEG₂₀₀₀-Chol (10) in DMSO-*d*₆

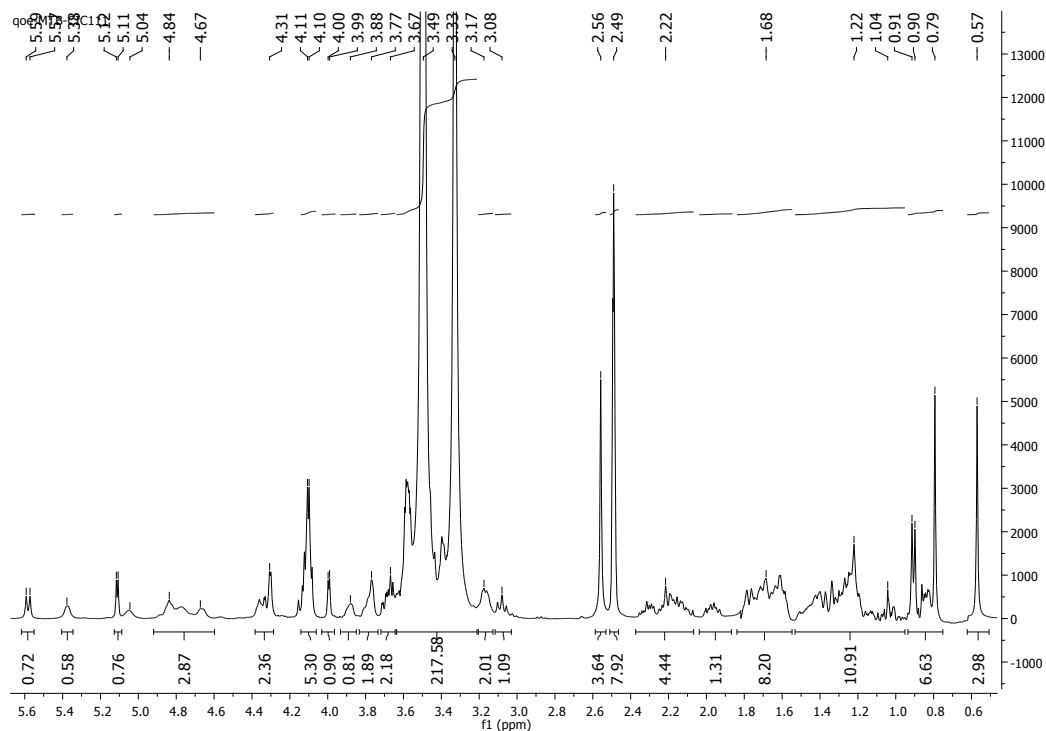


Fig. S 20 - ^1H -NMR spectrum of Suc-PEG₂₀₀₀-Chol (10) in DMSO-*d*₆.

17. ^1H -NMR spectrum of Suc-PEG₄₀₀₀-Chol (11) in DMSO-*d*₆

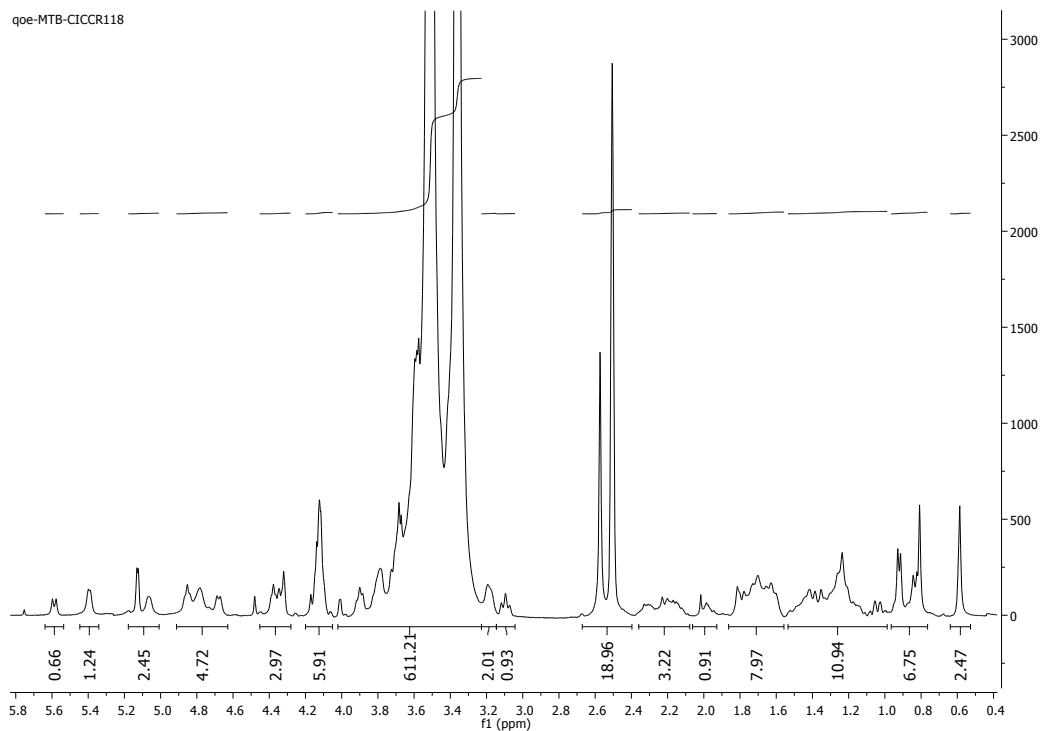


Fig. S 21 - ^1H -NMR spectrum of Suc-PEG₄₀₀₀-Chol (11) in DMSO-*d*₆.