

Supporting Information

Laterally Stretched Polycyclic Aromatic Hydrocarbons: Synthesis of Dibenzophenanthroheptaphene and Tetrabenzotriphenyleno-pyranthrene Derivatives.

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- II. UV-Emission spectra of 9.
- III. II. Computational Details (3D optimized structures and their HOMO-LUMO configurations).

I. Experimental details MALDI 1a-c, 2a-b, 4a-c, 9 and 10a,b.

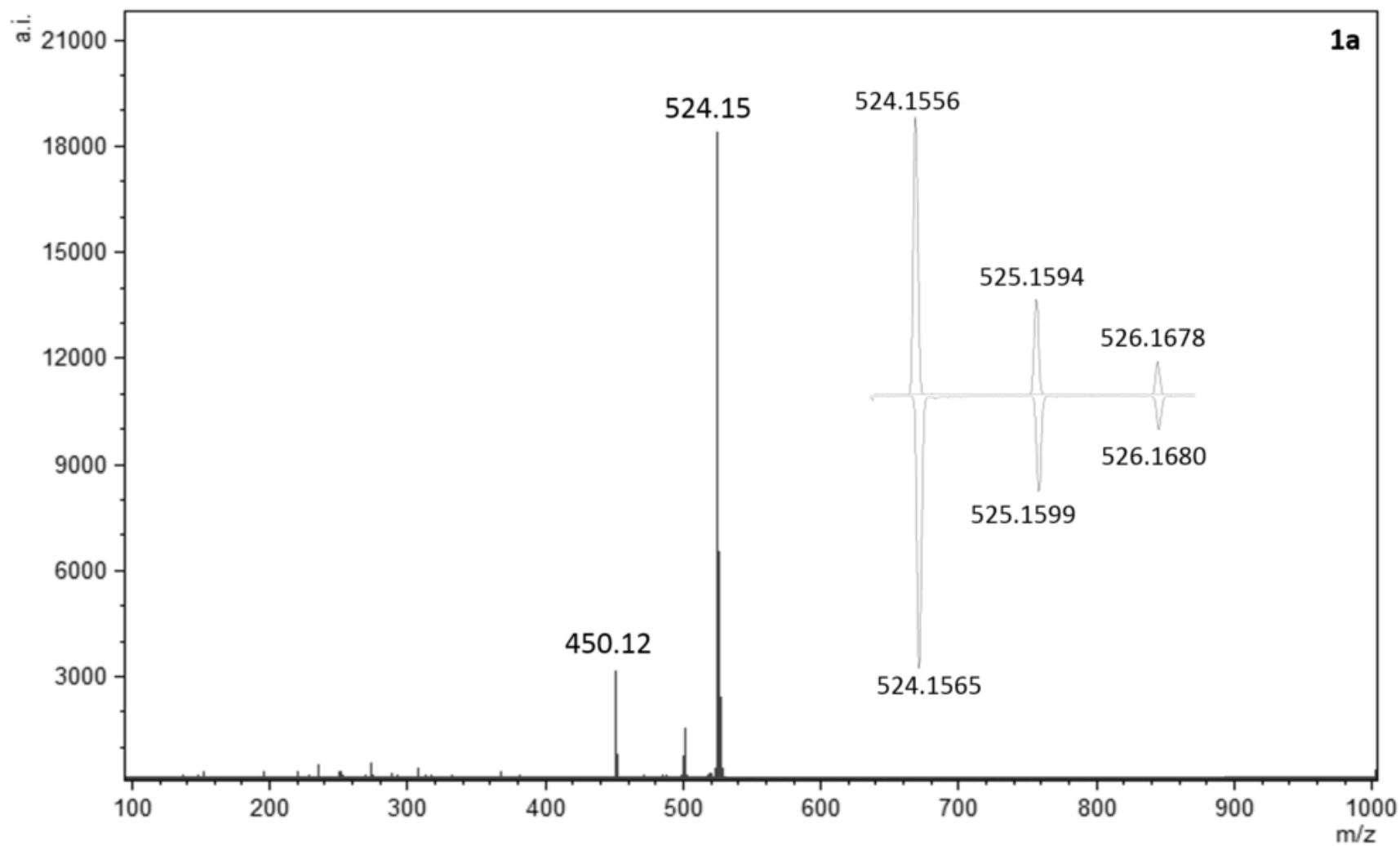


Figure S1. HR-MALDI spectrum of 1a

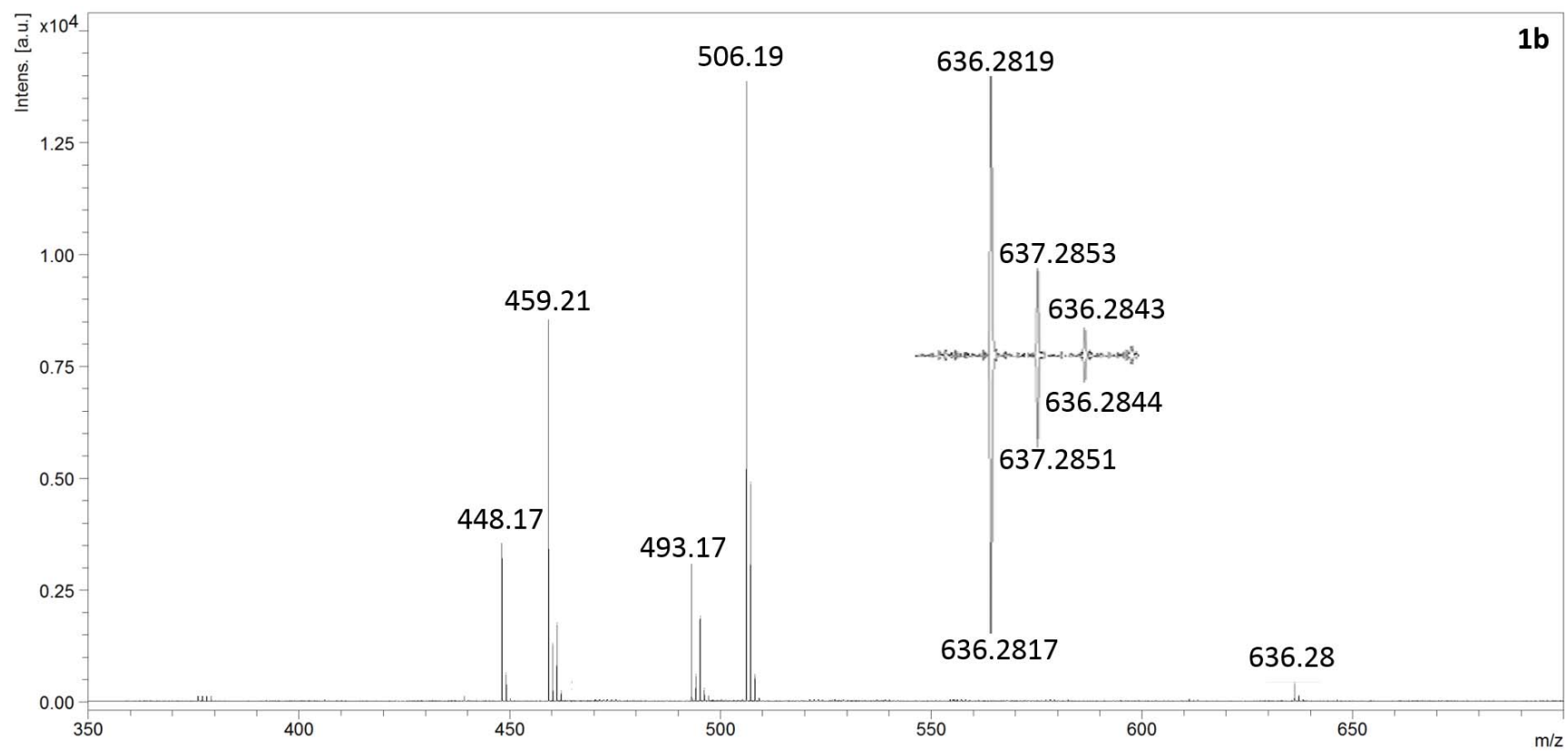


Figure S2. HR-MALDI spectrum of **1b**

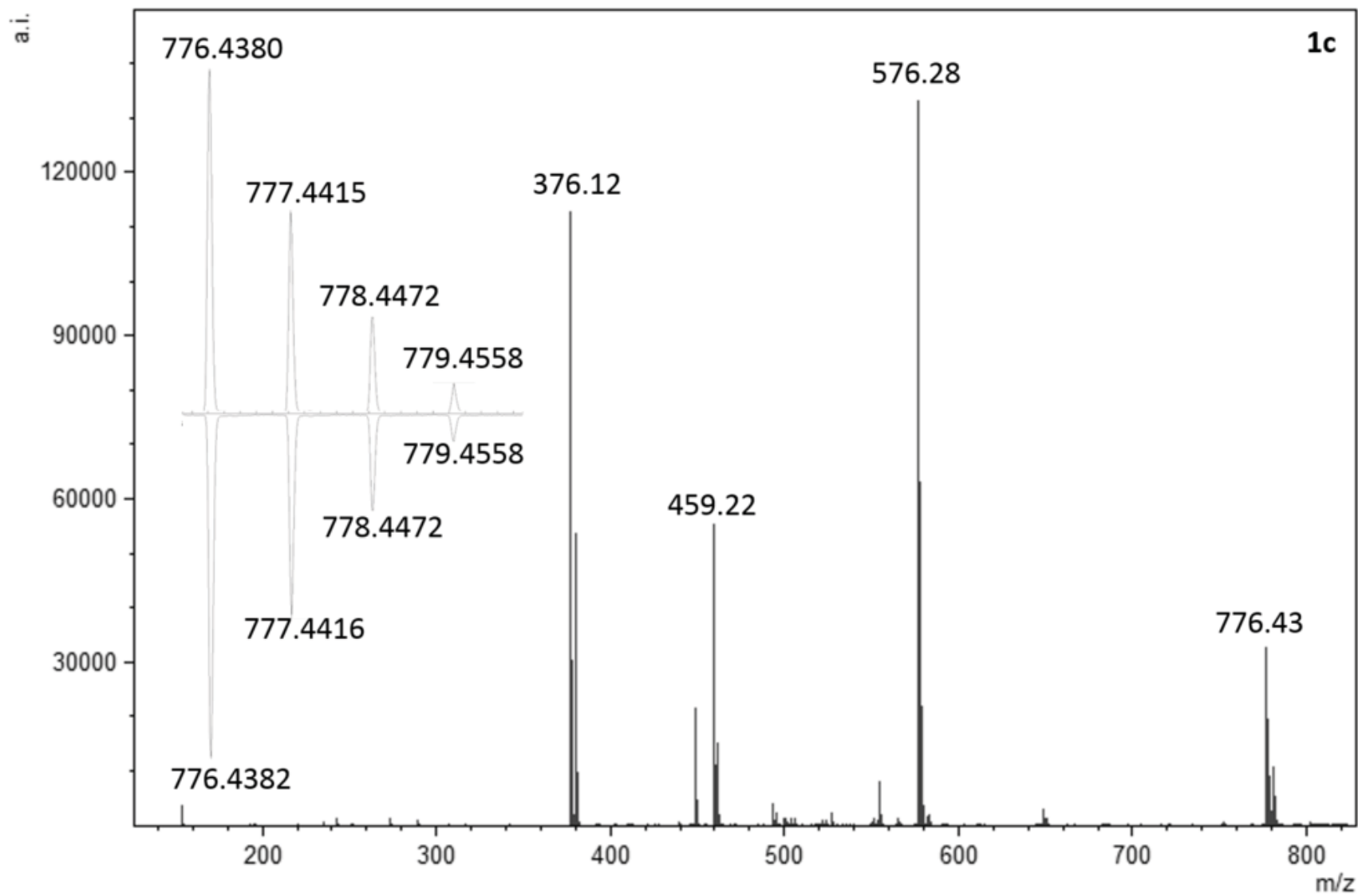


Figure S3. HR-MALDI spectrum of **1c**

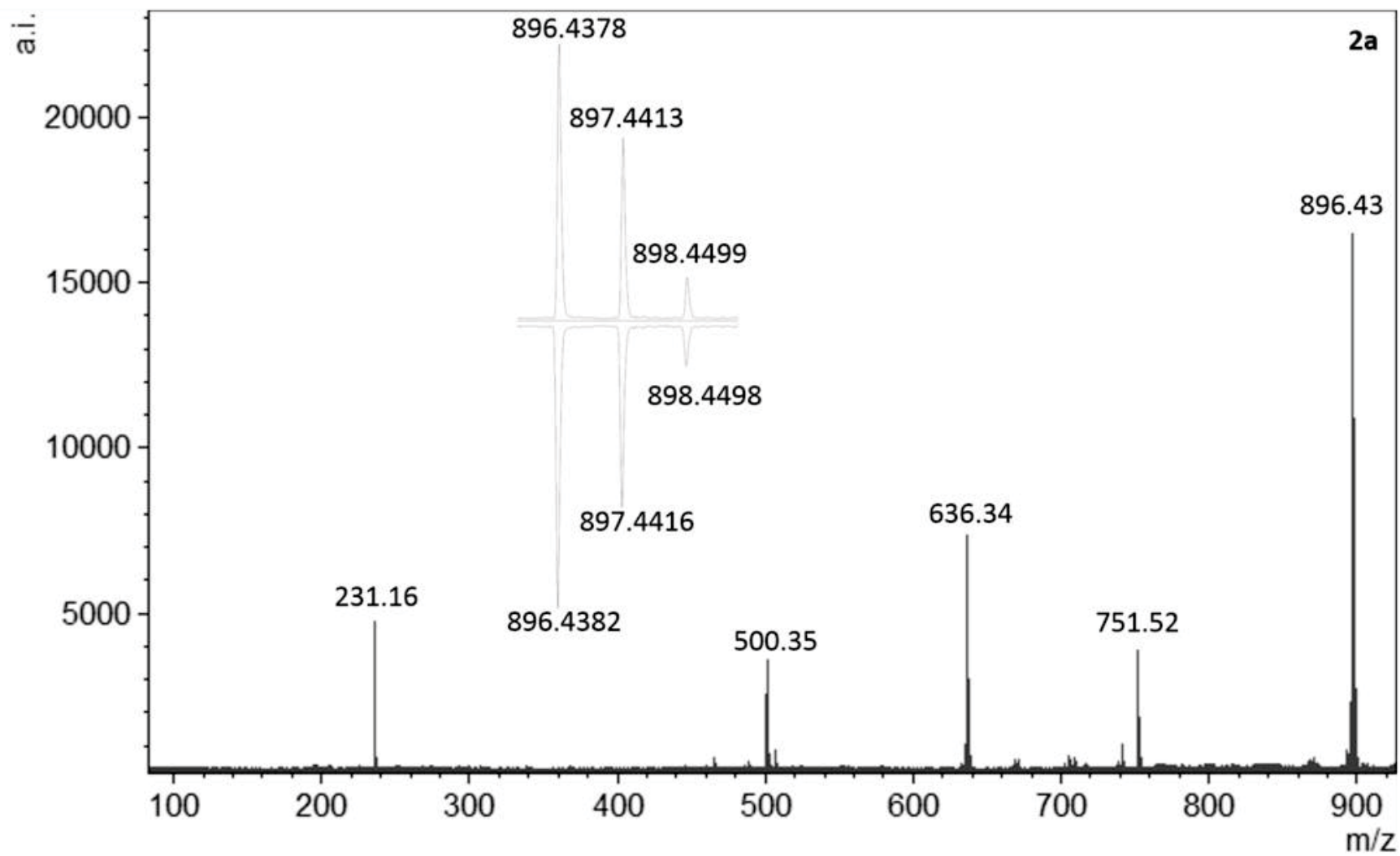


Figure S4. HR-MALDI spectrum of 2a

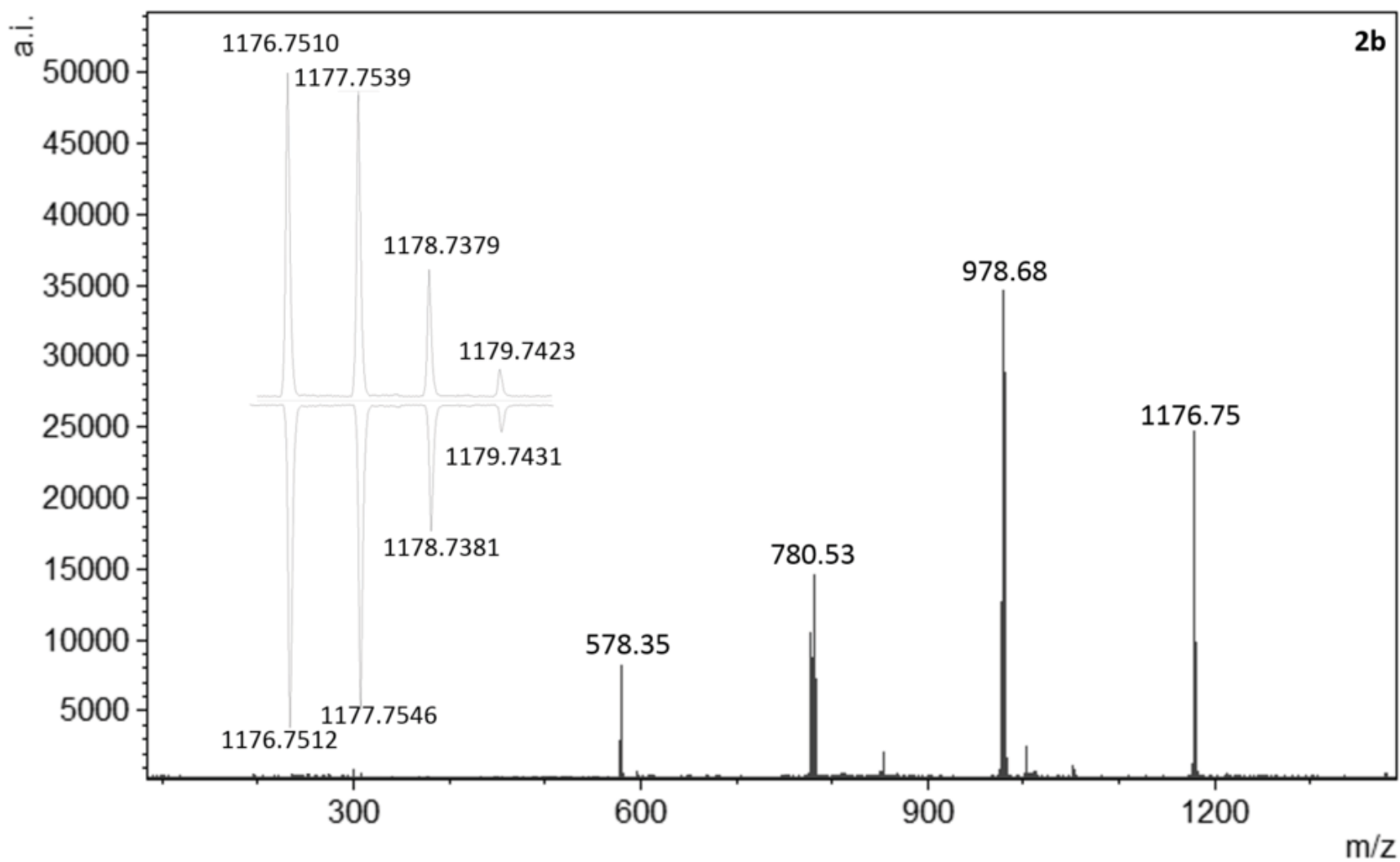


Figure S5. HR-MALDI spectrum of **2b**

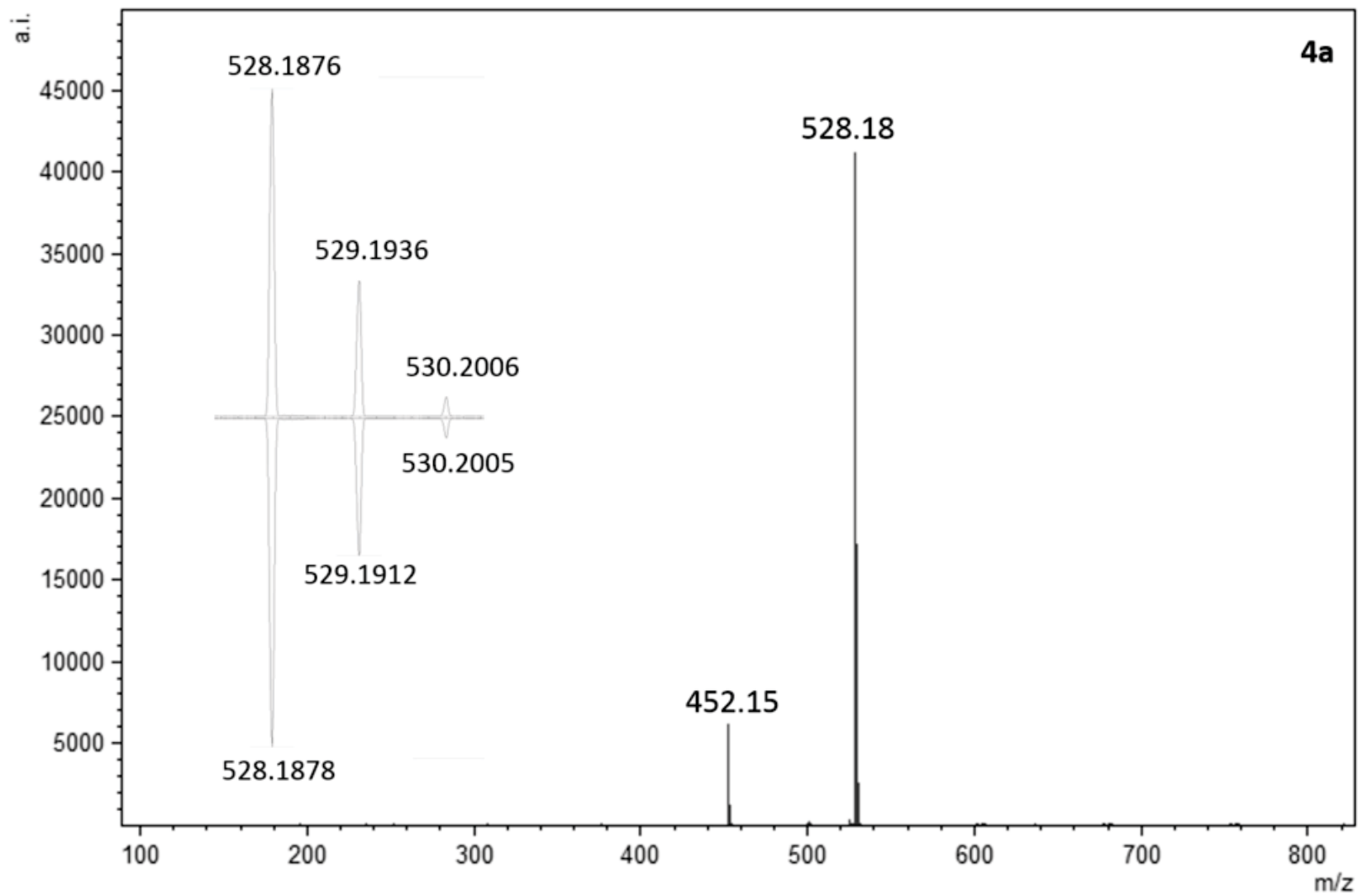


Figure S6. HR-MALDI spectrum of 4a

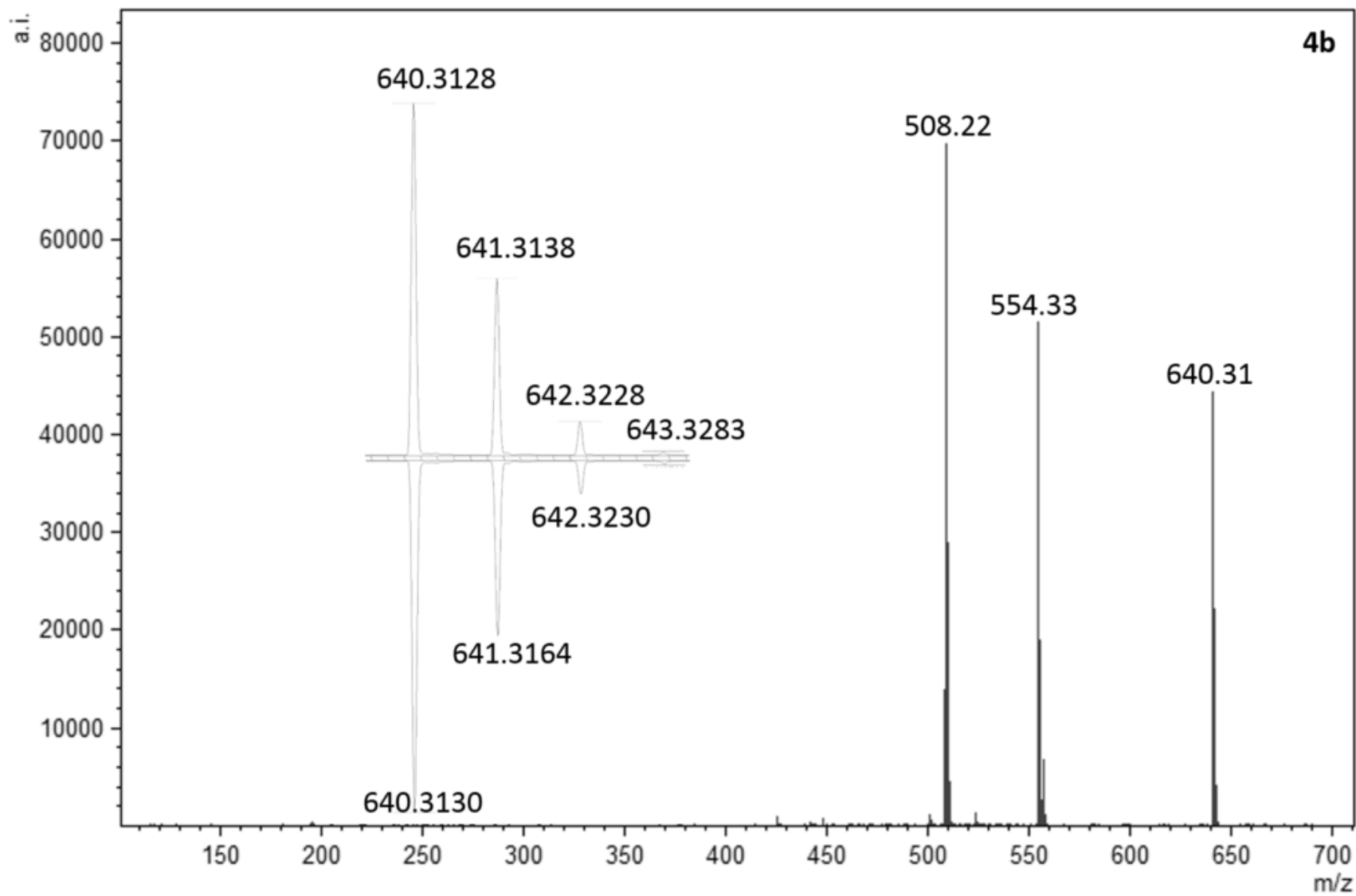


Figure S7. HR-MALDI spectrum of 4b

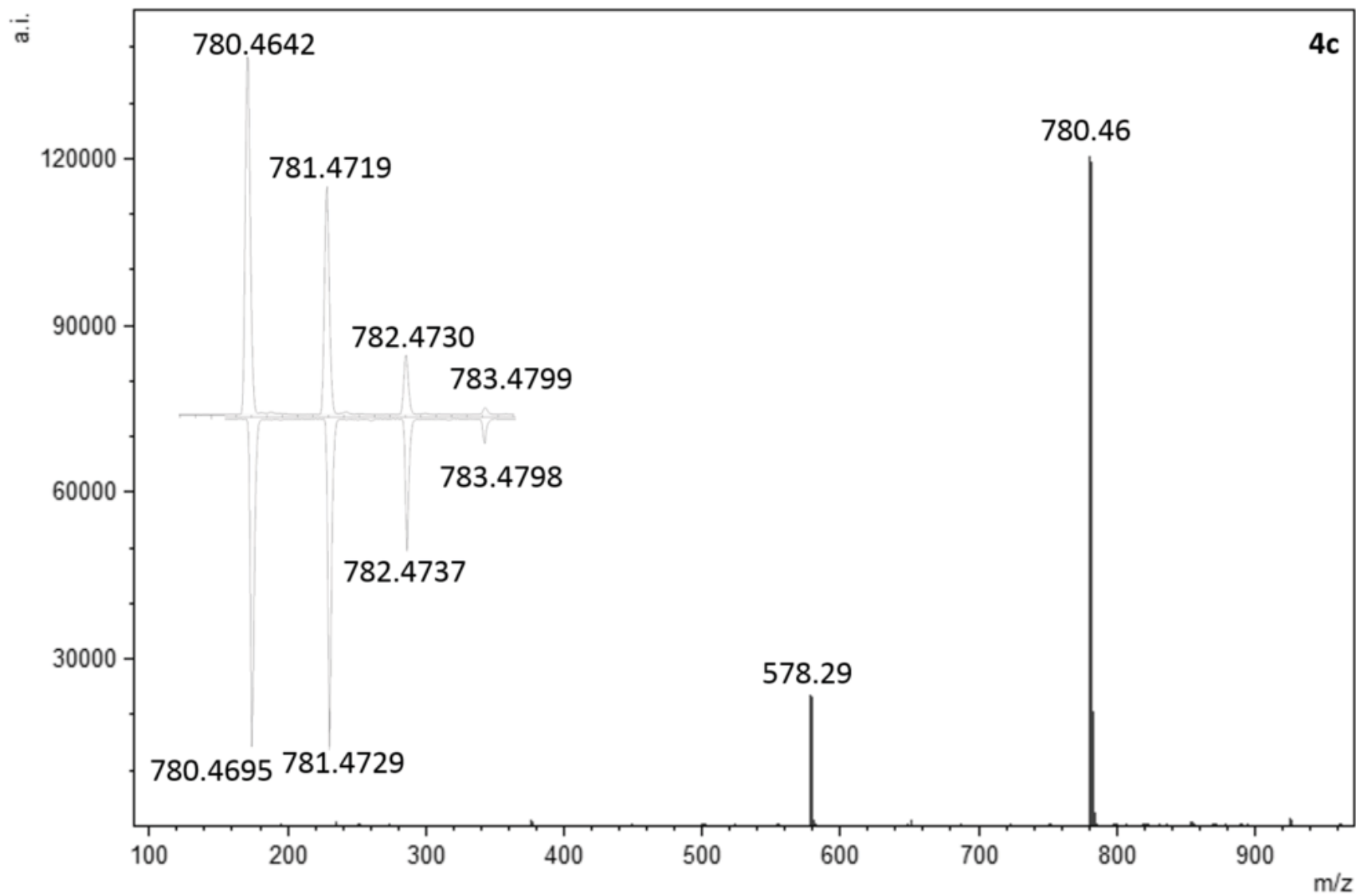


Figure S8. HR-MALDI spectrum of 4c

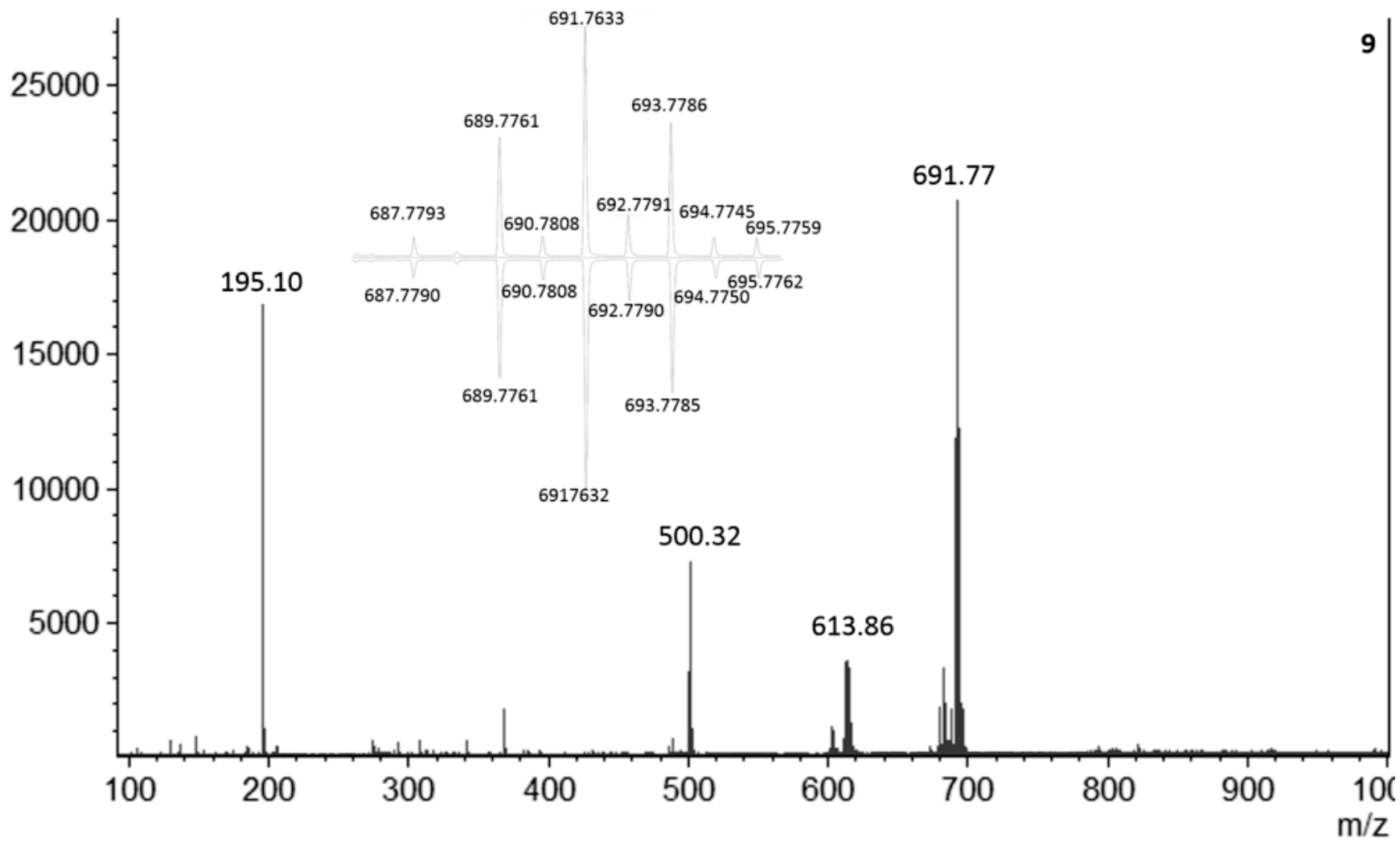


Figure S9. HR-MALDI spectrum of 9

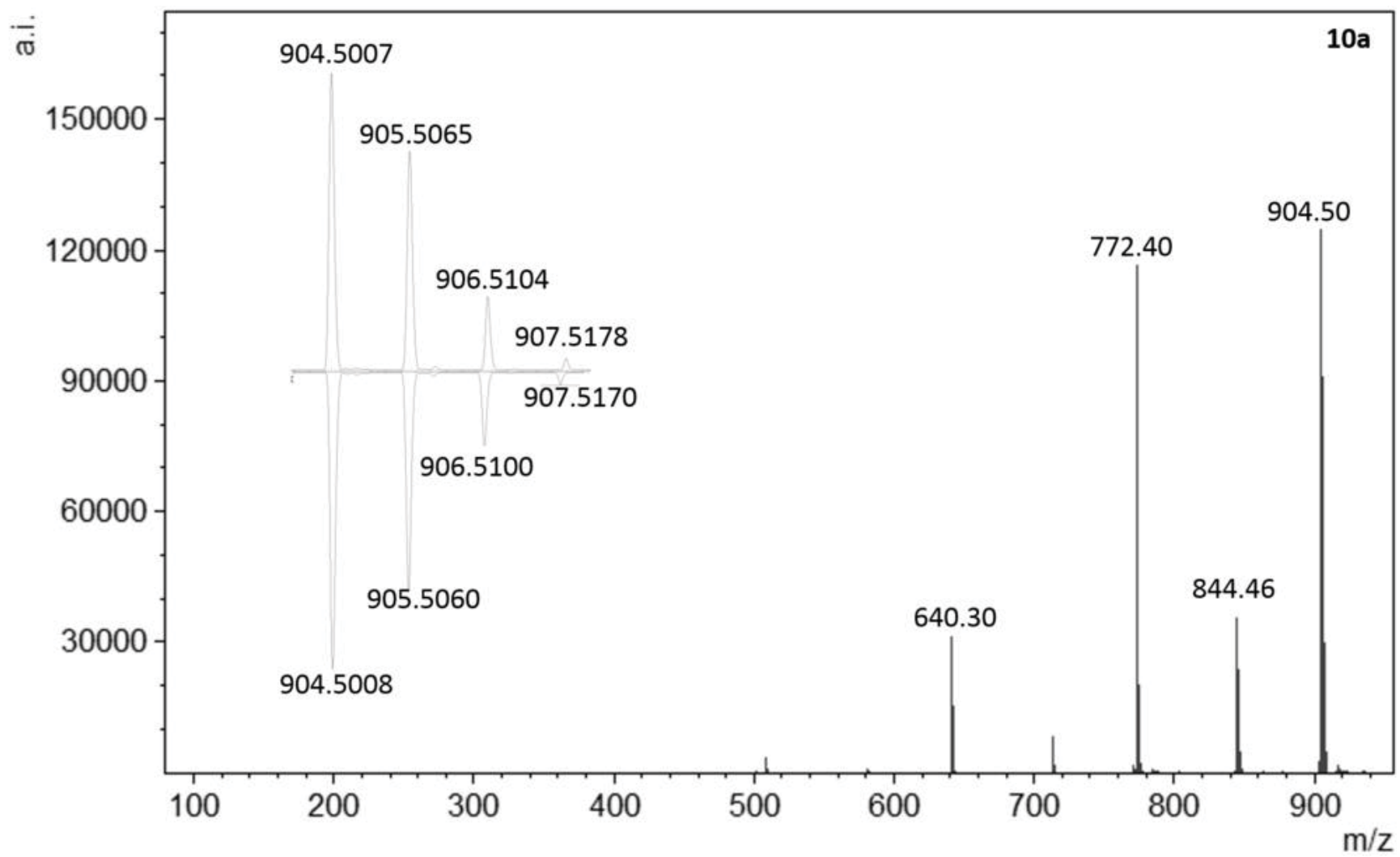


Figure S10 HR-MALDI spectrum of 10a

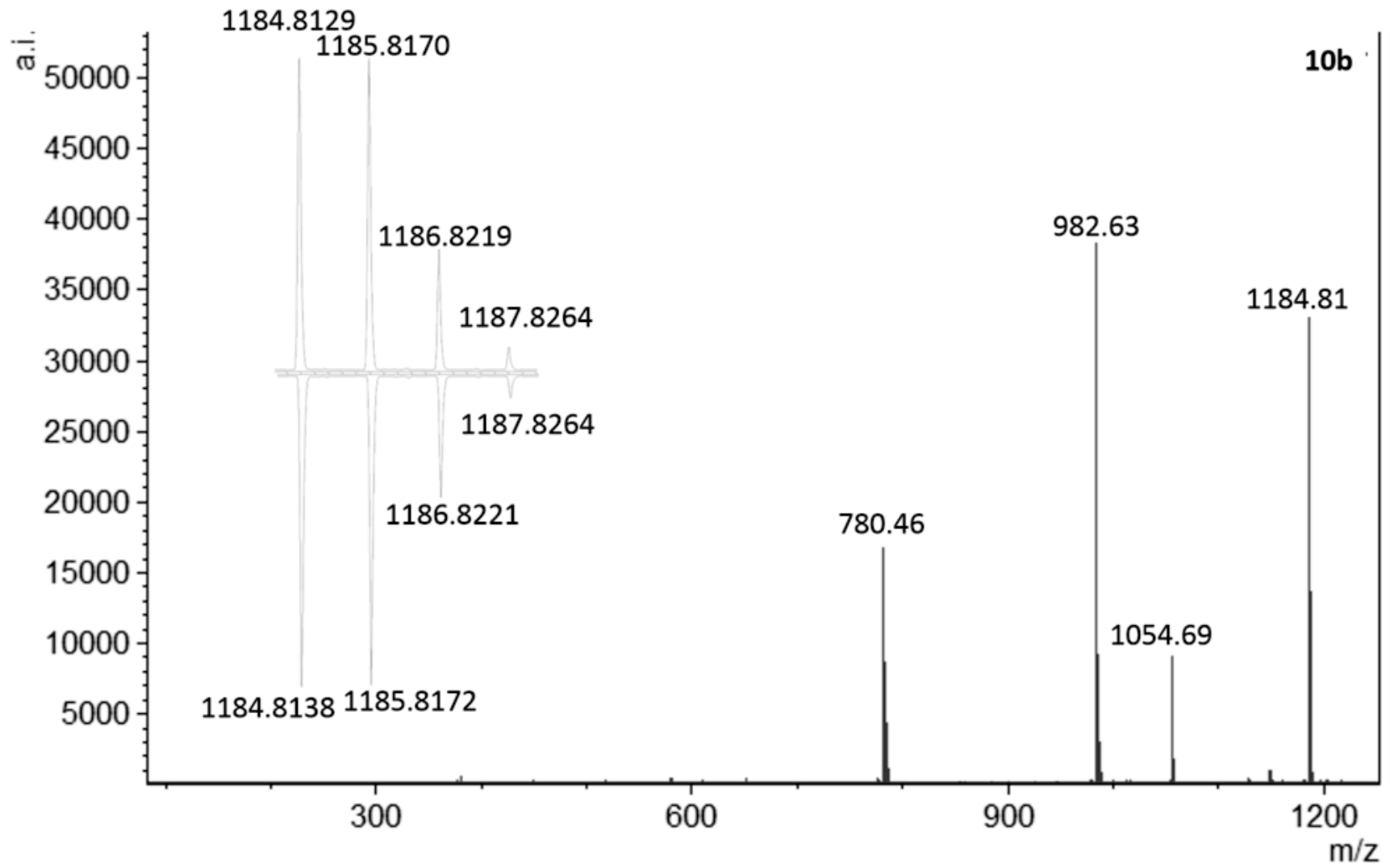


Figure S11. HR-MALDI spectrum of **10b**

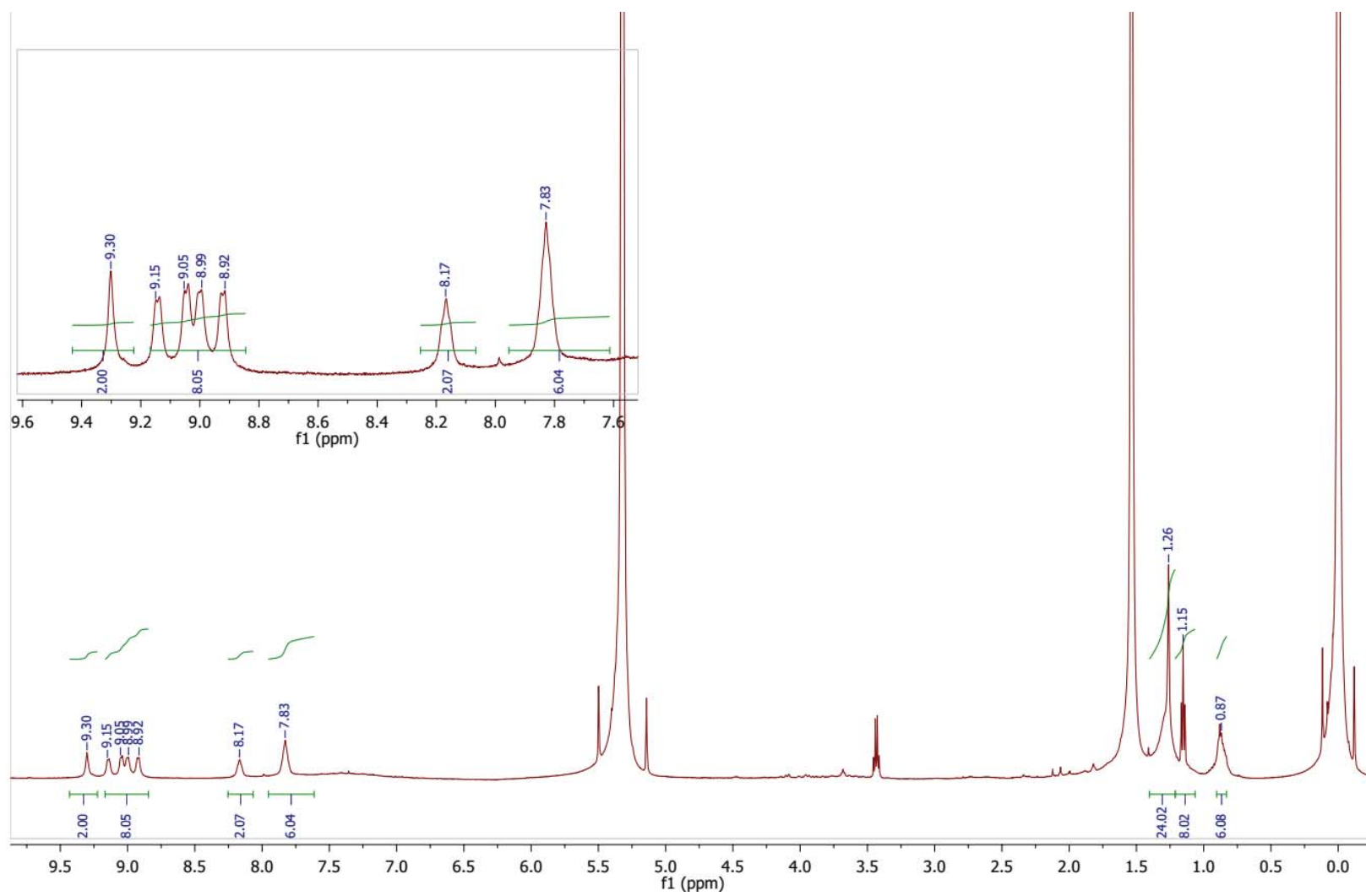


Figure S12. ^1H NMR spectrum of **1c**

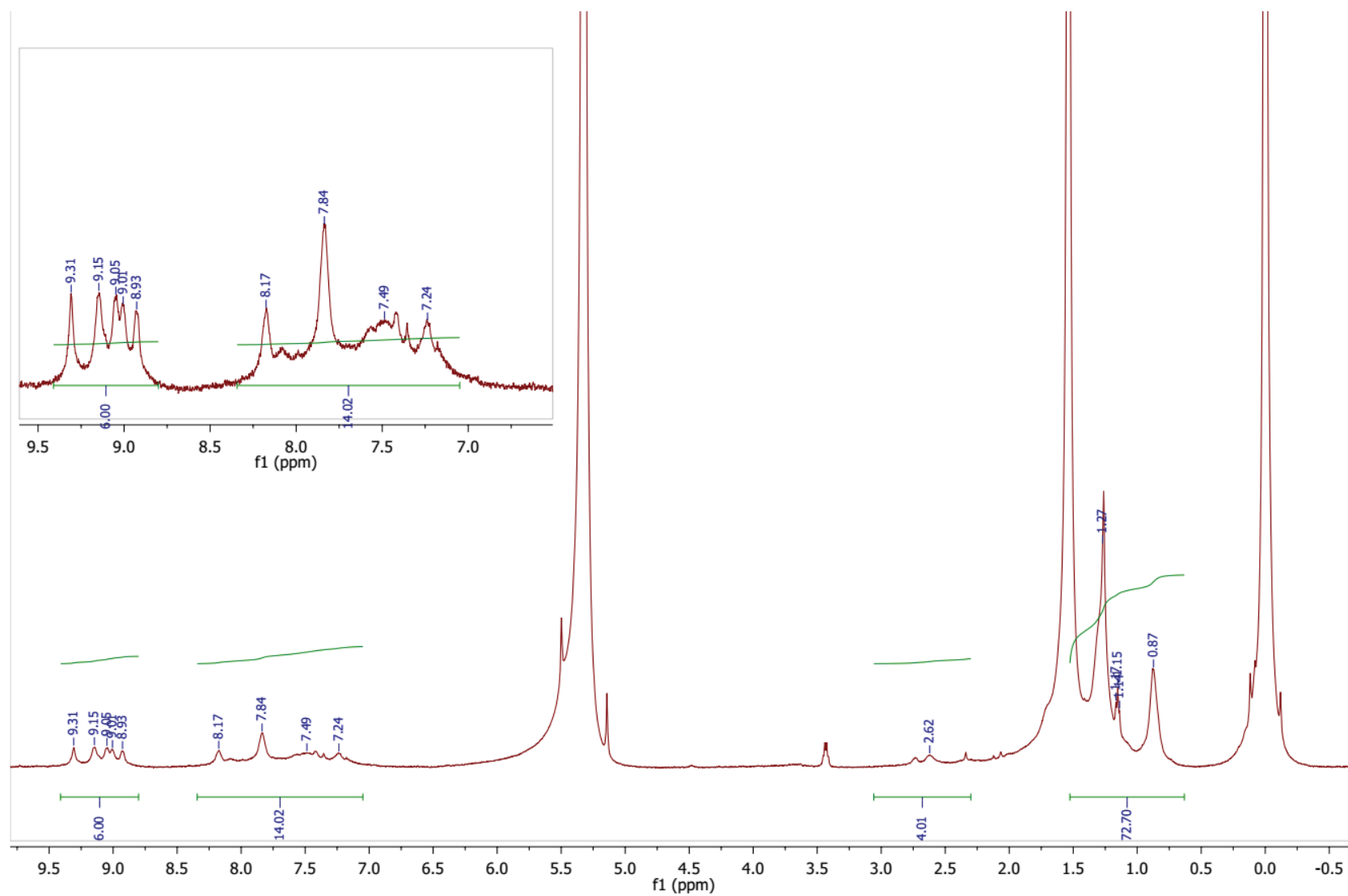


Figure S13. ¹H NMR spectrum of **2b**

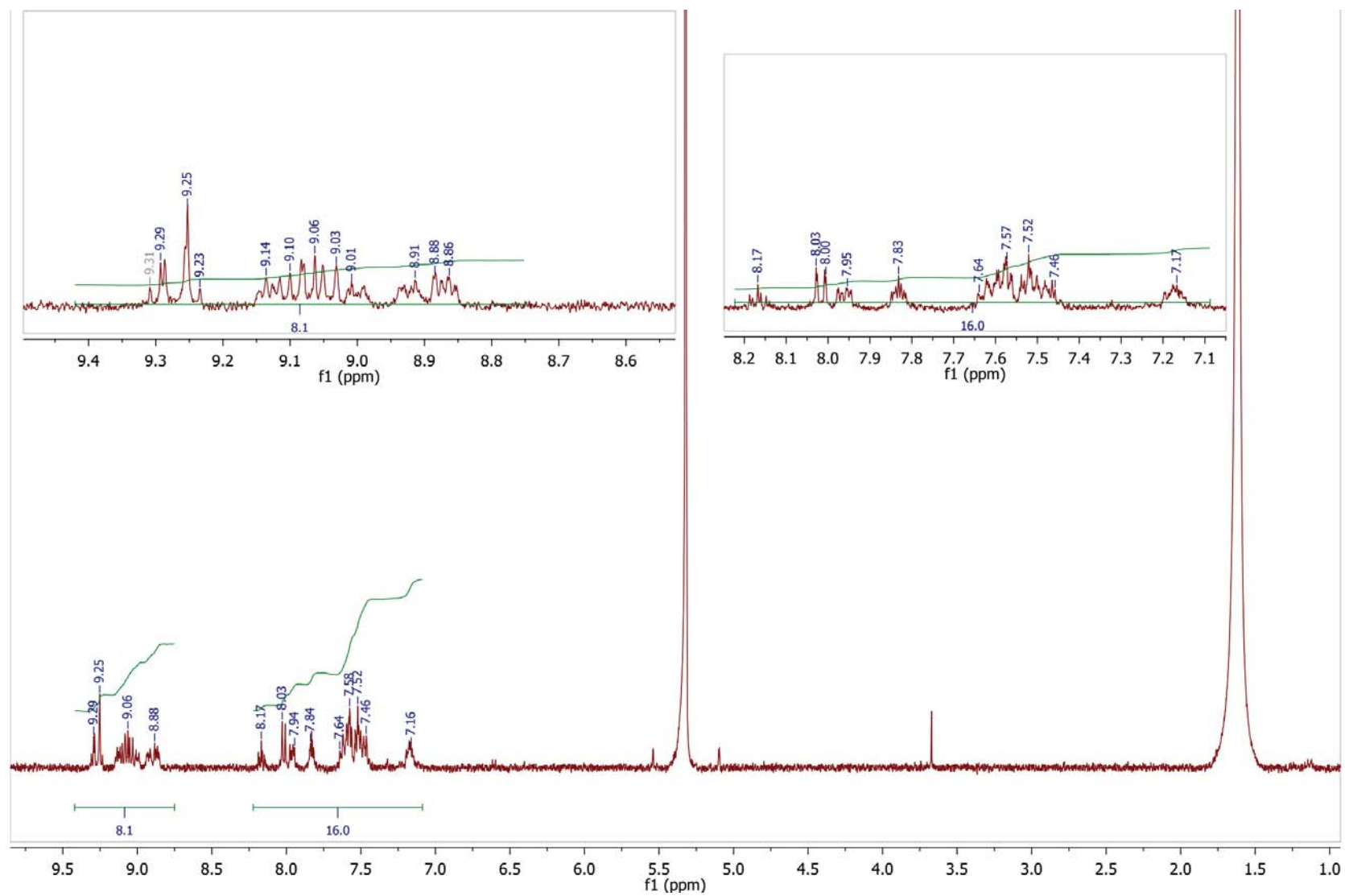


Figure S14. ¹H NMR spectrum of 4a

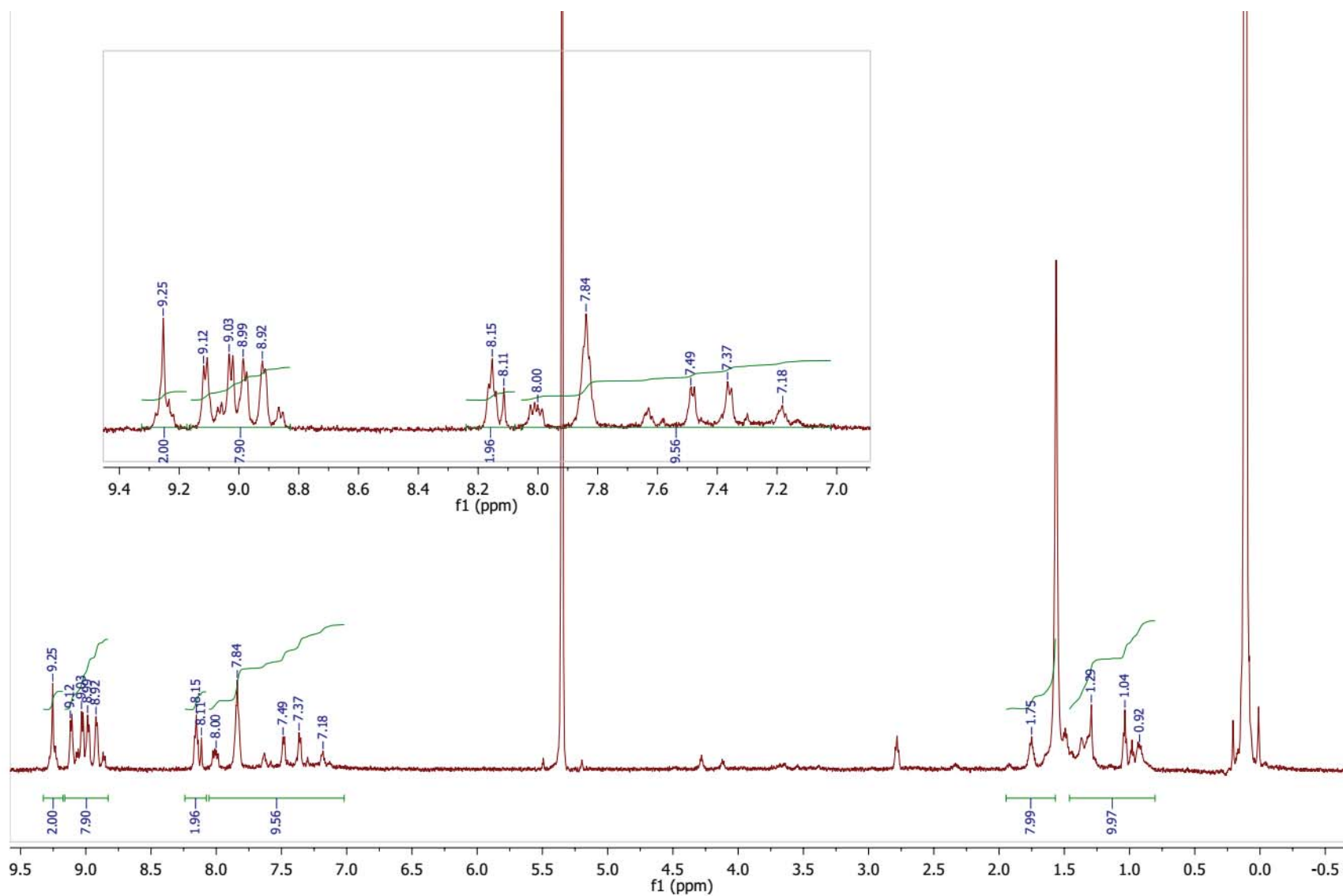


Figure S15. ¹H NMR spectrum of 4b

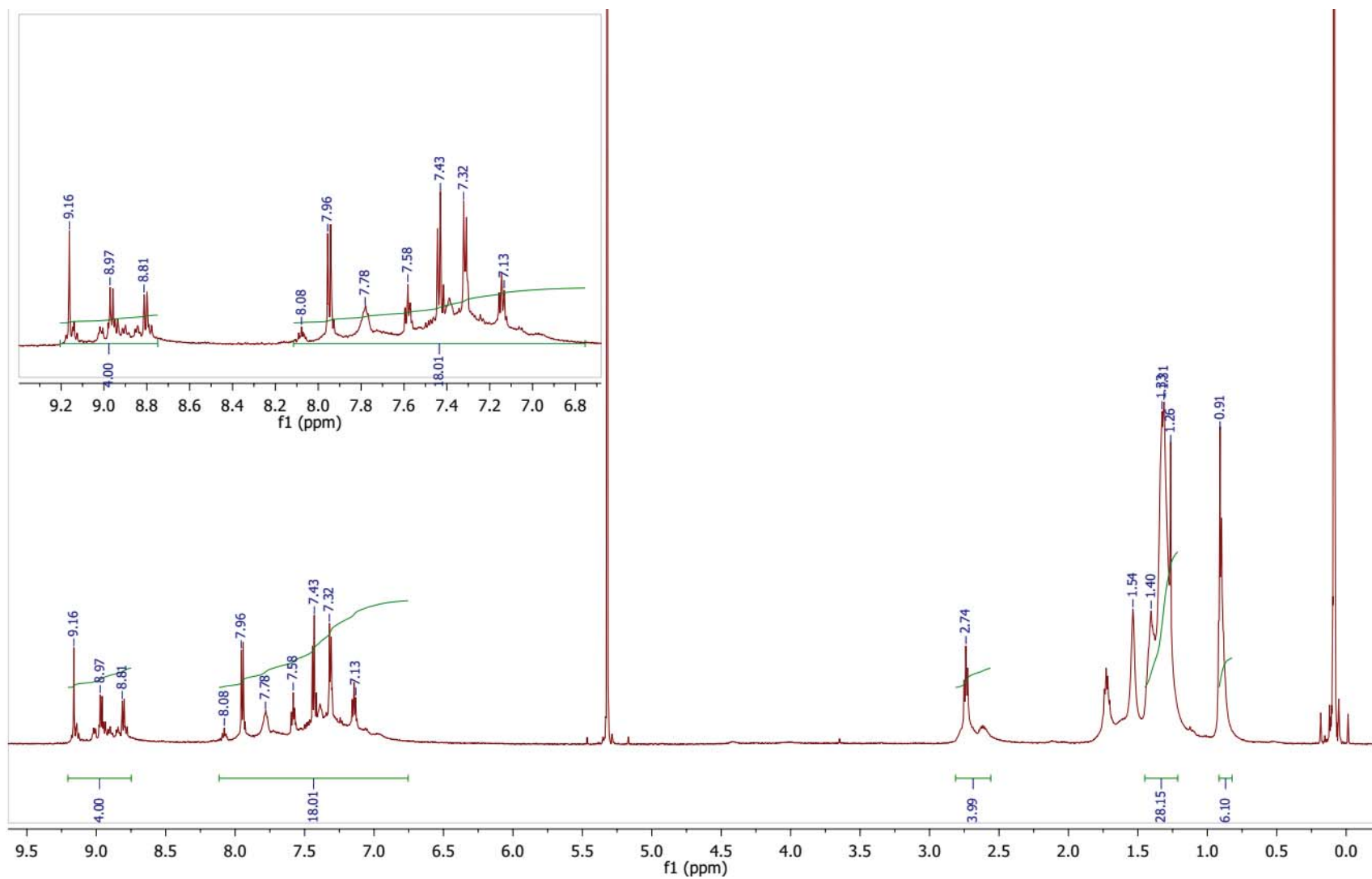


Figure S16. ¹H NMR spectrum of 4c

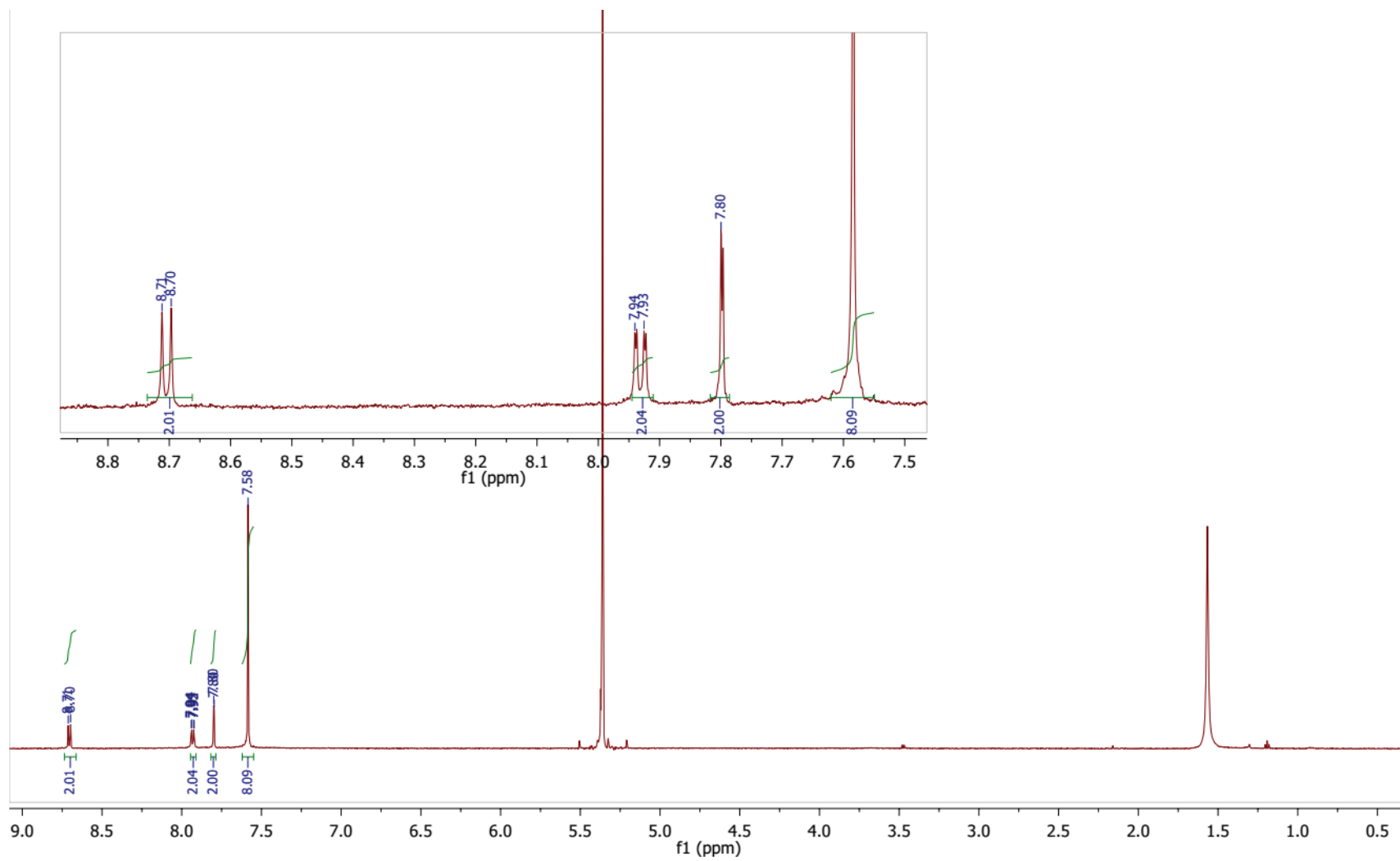


Figure S17. ^1H NMR spectrum of 9

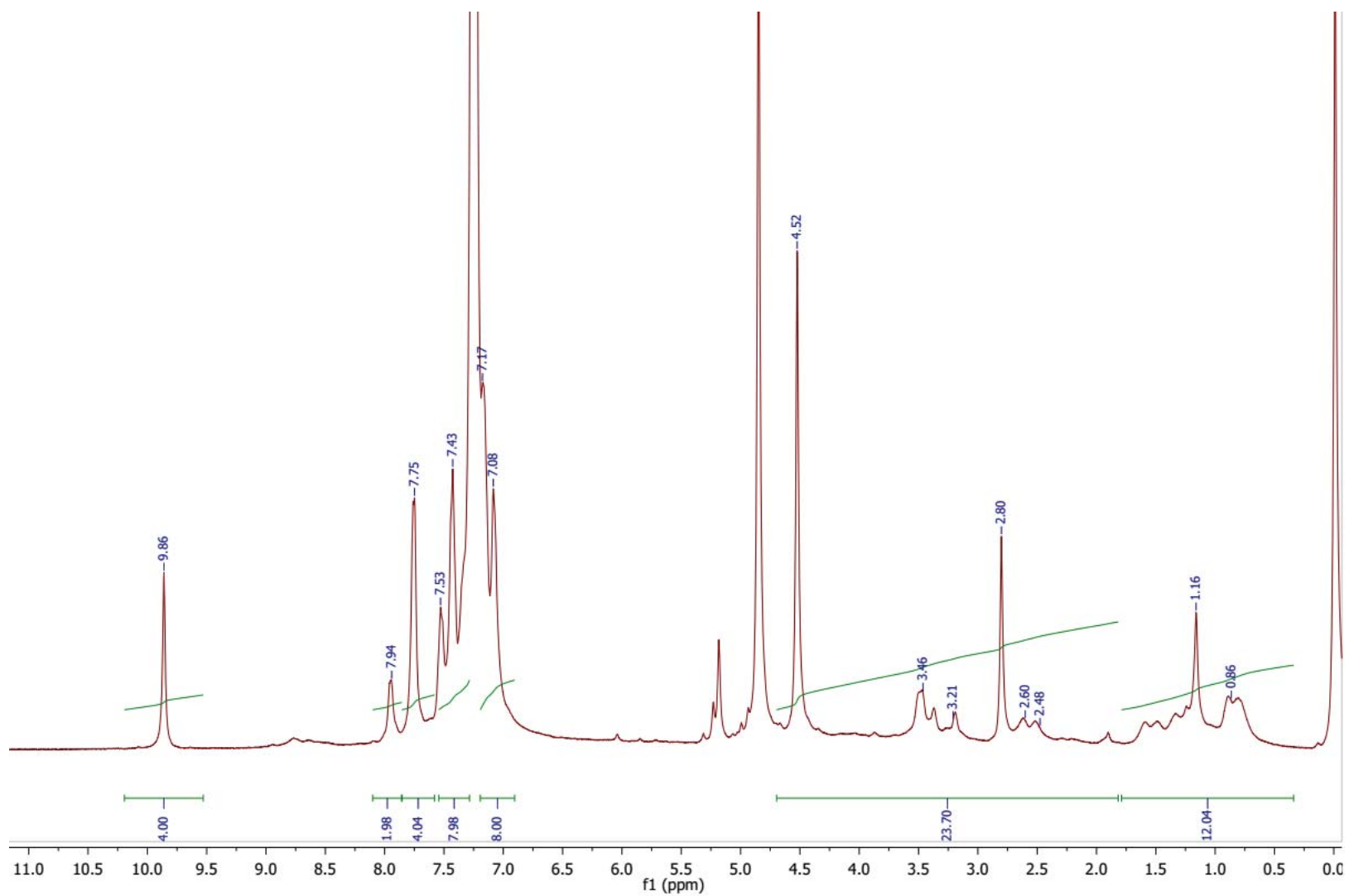


Figure S18. ¹H NMR spectrum of 10a

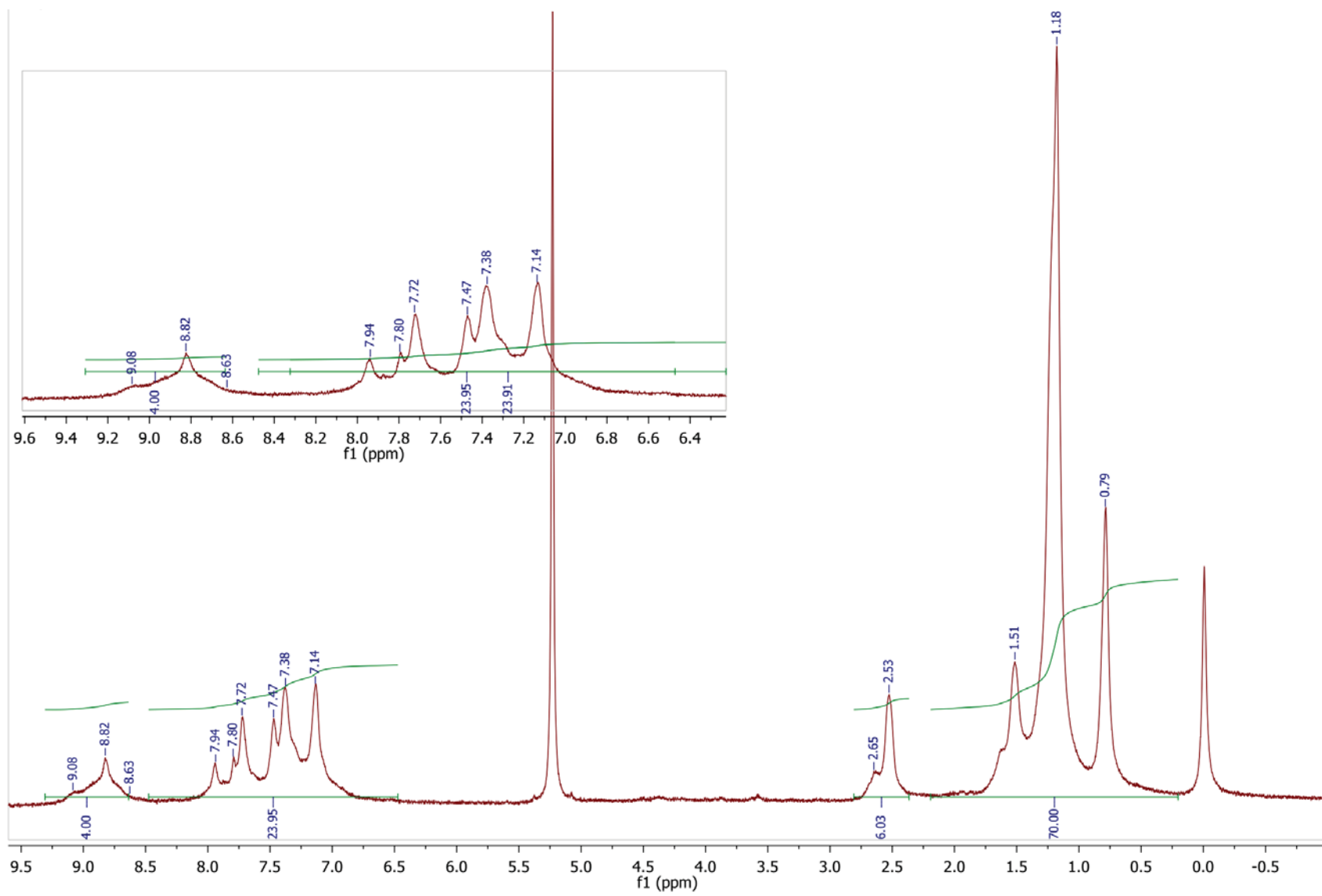


Figure S19. ¹H NMR spectrum of 10b

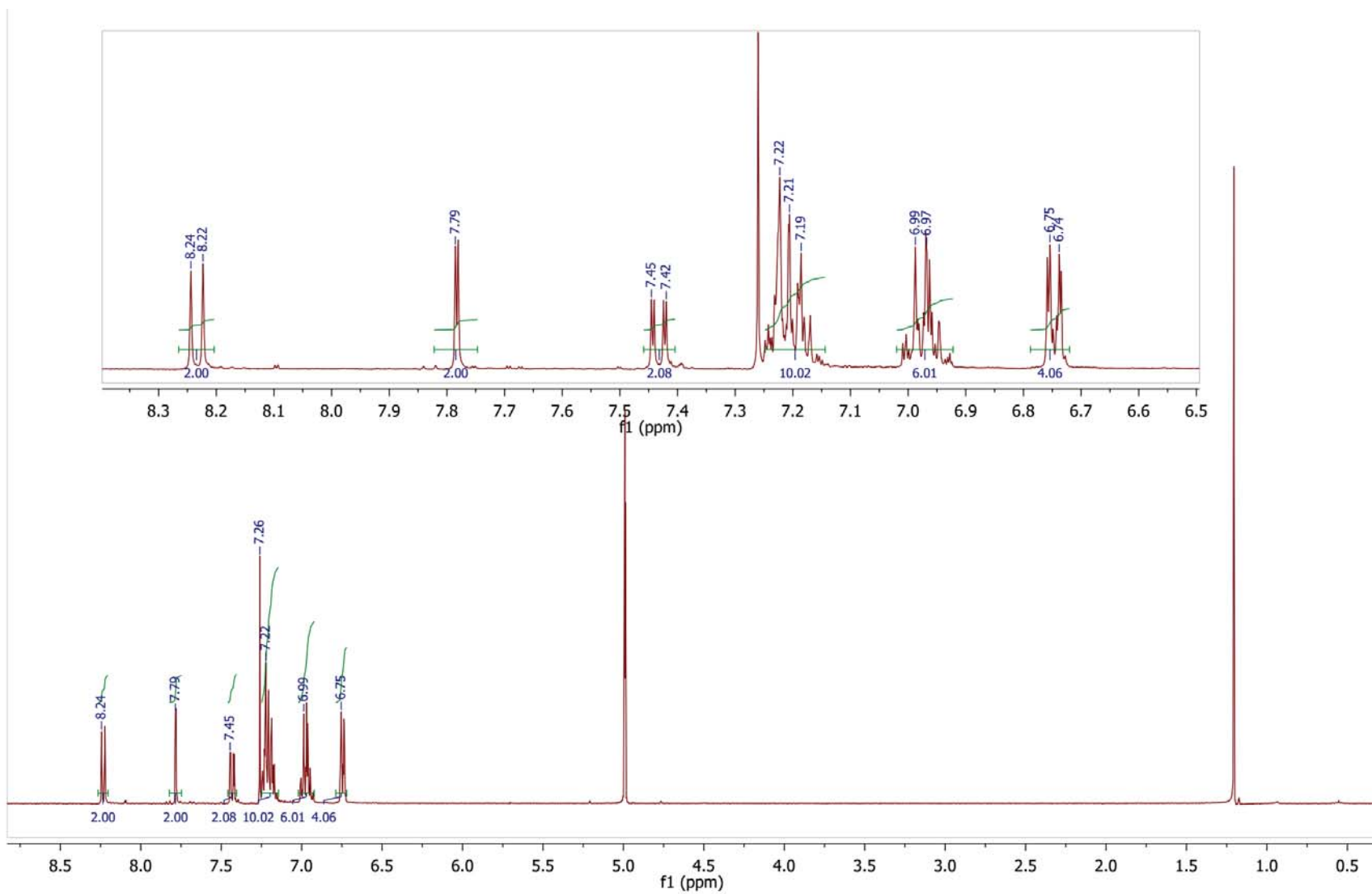


Figure S20. ^1H NMR spectrum of 13a

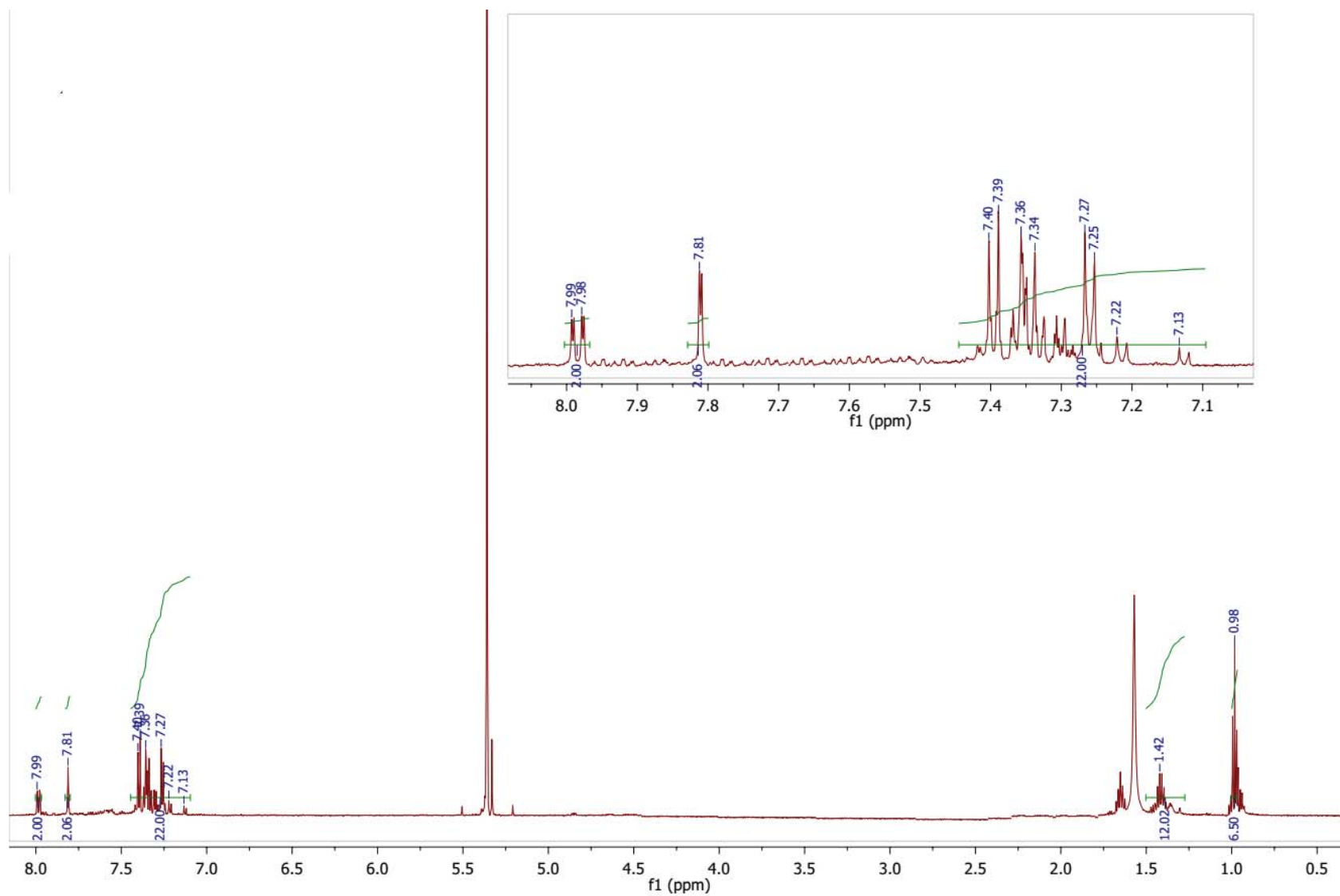


Figure S21. ¹H NMR spectrum of 13b

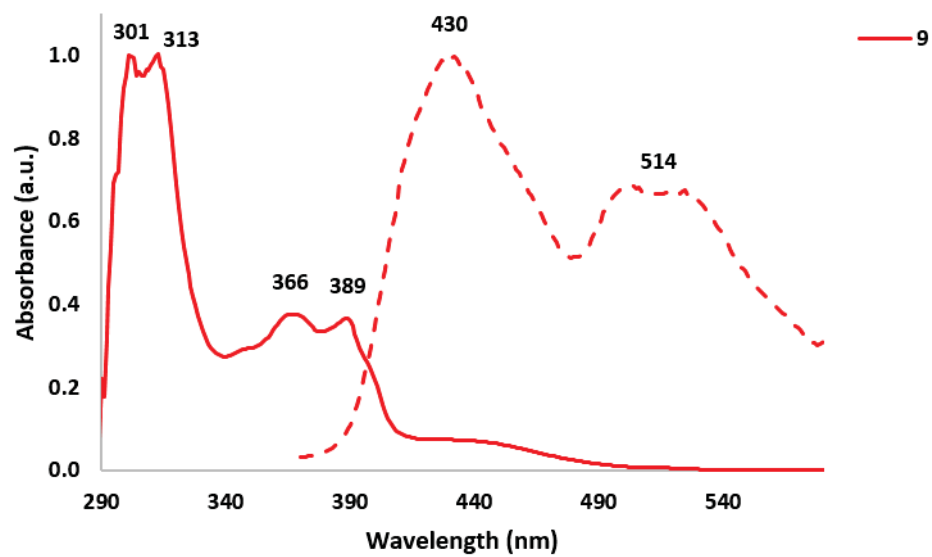
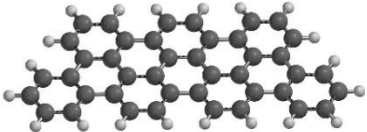
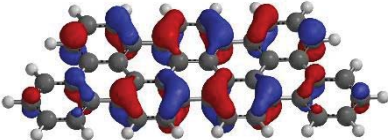
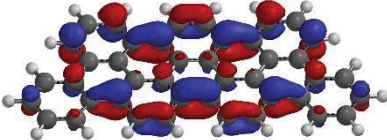
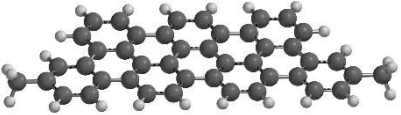
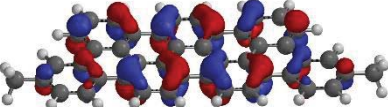
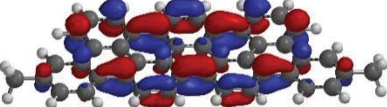
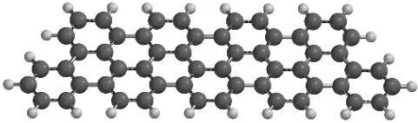
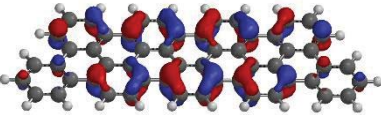
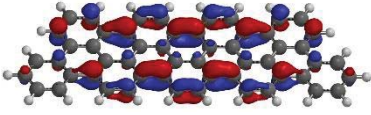


Figure S22. Normalized absorption and emission spectra for the pentaphene derivatives **9** in THF. The solid lines represent absorption spectra and the dotted lines correspond to emission spectra.

II. Computational details

Table S1. Optimized structures and HOMO-LUMO frontier orbitals of DBPH and TBTP at the level of the B₃LYP/6-31 G* basis set with R = H (1a and 2-H) or Me (1-Me and 2-Me) instead of the alkyl groups for simplicity.

No	Optimized structure	HOMO	LUMO
1a			
1Me			
2H			
2Me	