

Supplementary Materials

Selective C-N σ Bond Cleavage in Azetidinyl Amides under Transition Metal Free Conditions

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Copies of ^1H -NMR, ^{13}C -NMR 2a–2n and HRMS Spectrum of 2g and 2n

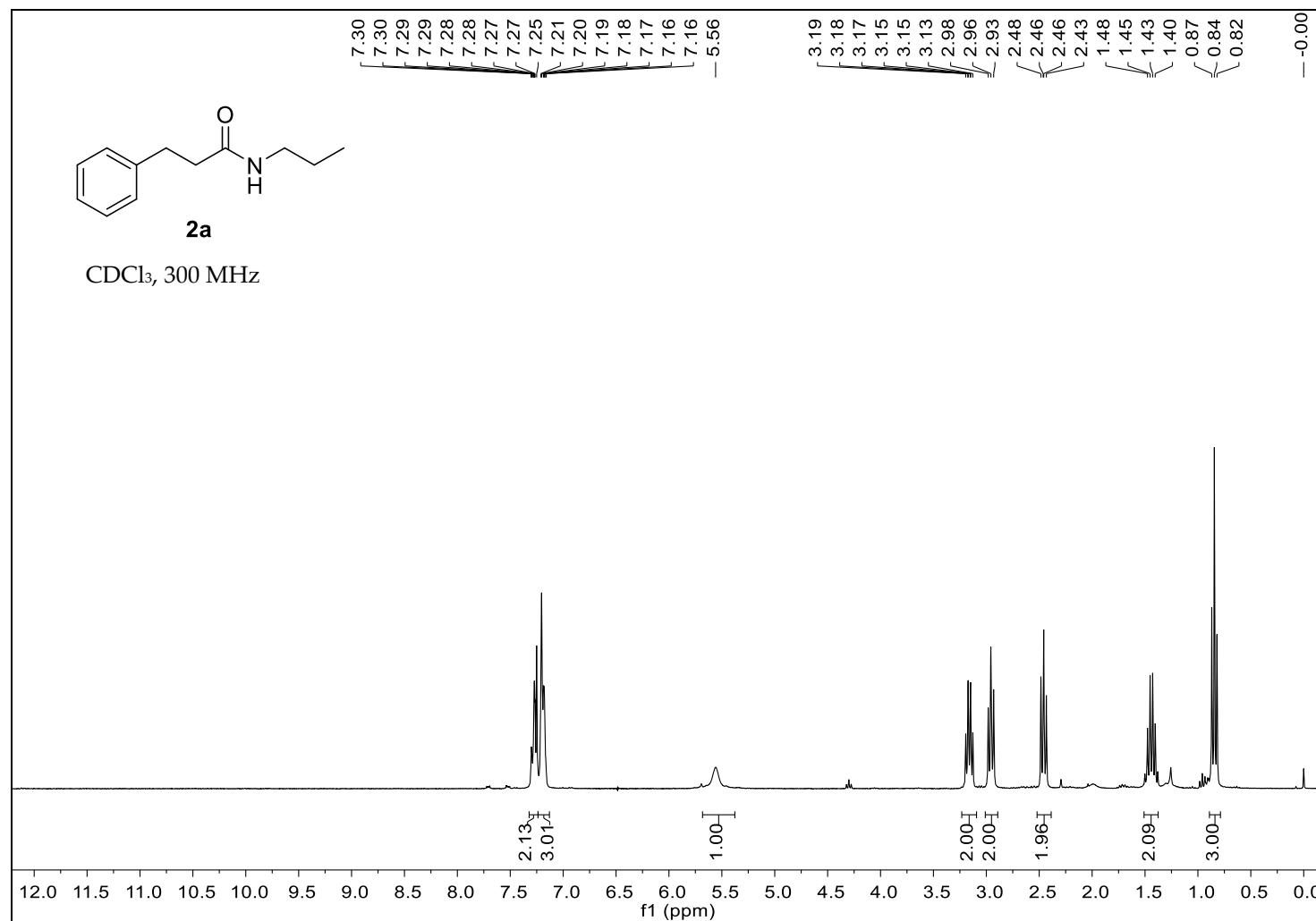


Figure S1. ¹H-NMR of 3-Phenyl-*N*-propylpropanamide (**2a**).

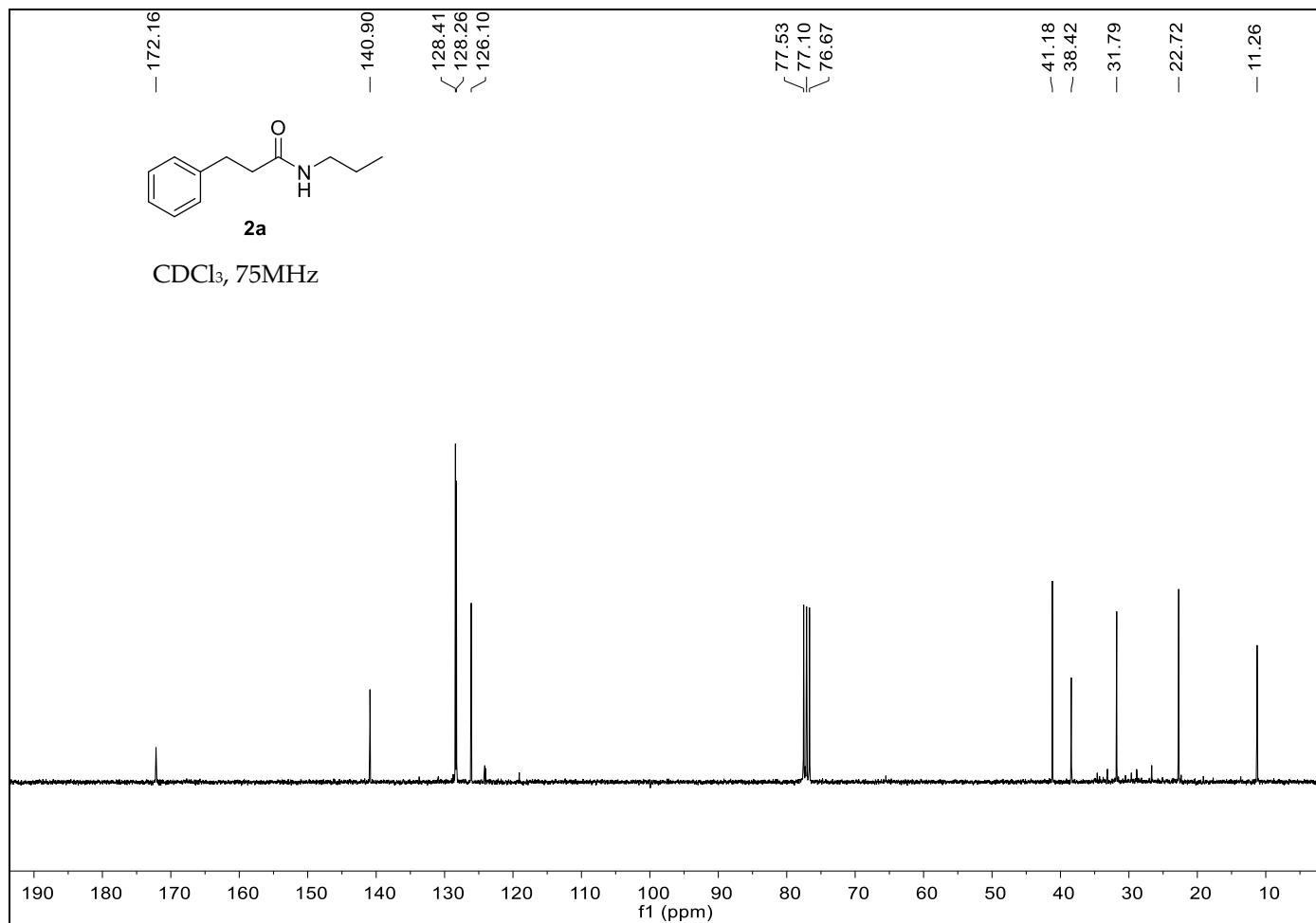


Figure S2. ^{13}C -NMR of 3-Phenyl-N-propylpropanamide (**2a**).

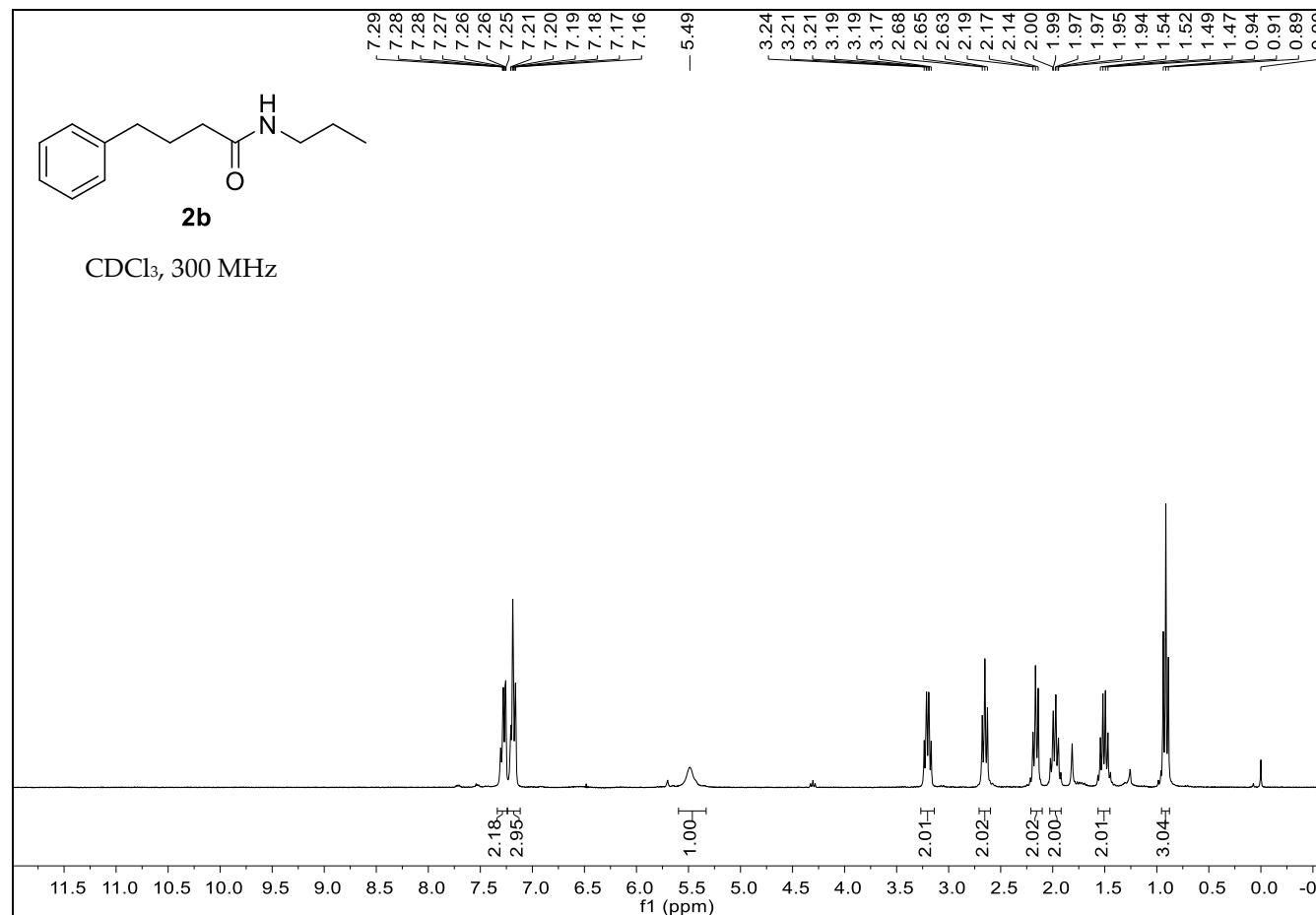


Figure S3. ^1H NMR of 4-Phenyl-N-propylbutanamide (**2b**).

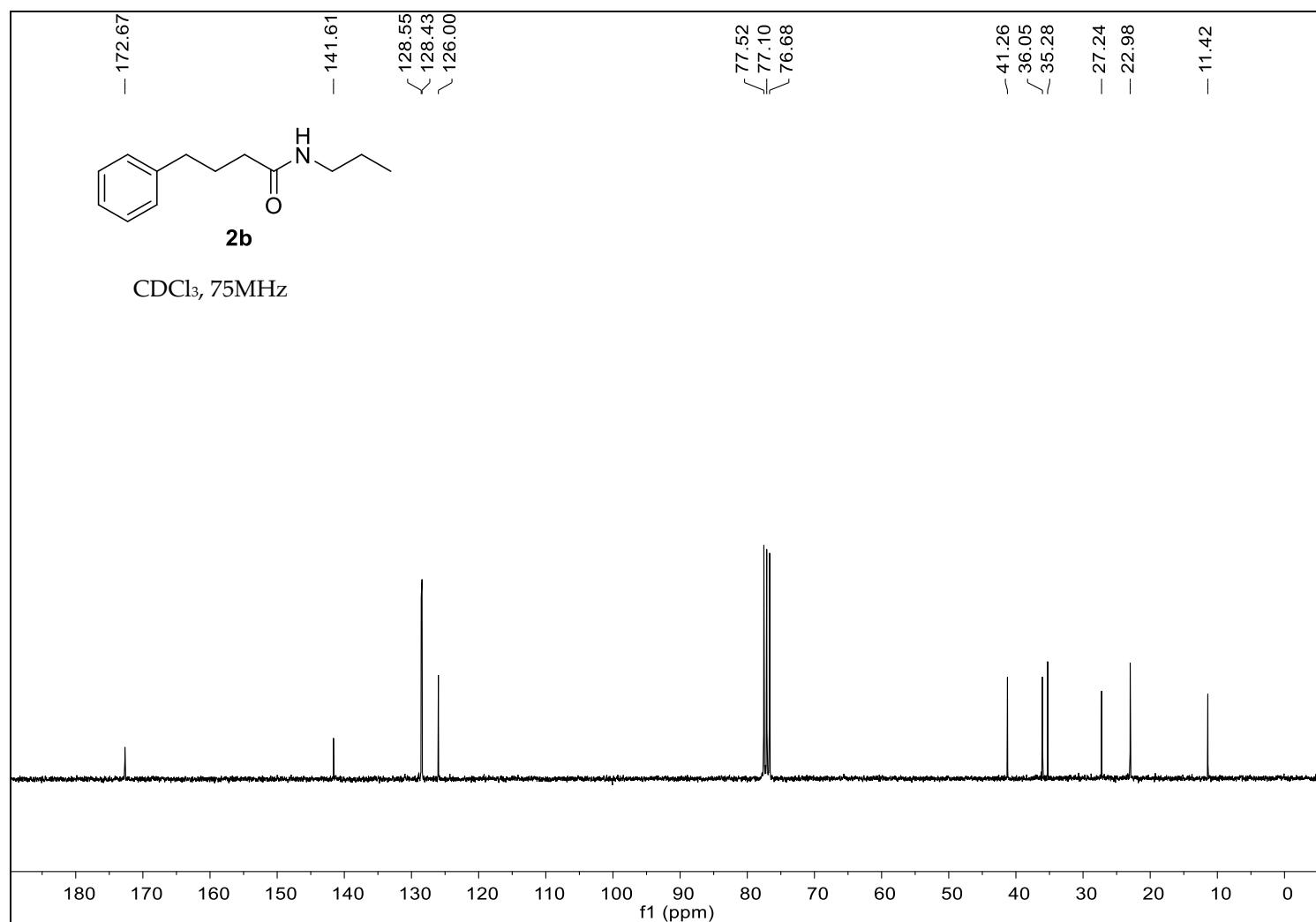


Figure S4. ^{13}C -NMR of 4-Phenyl-N-propylbutanamide (**2b**).

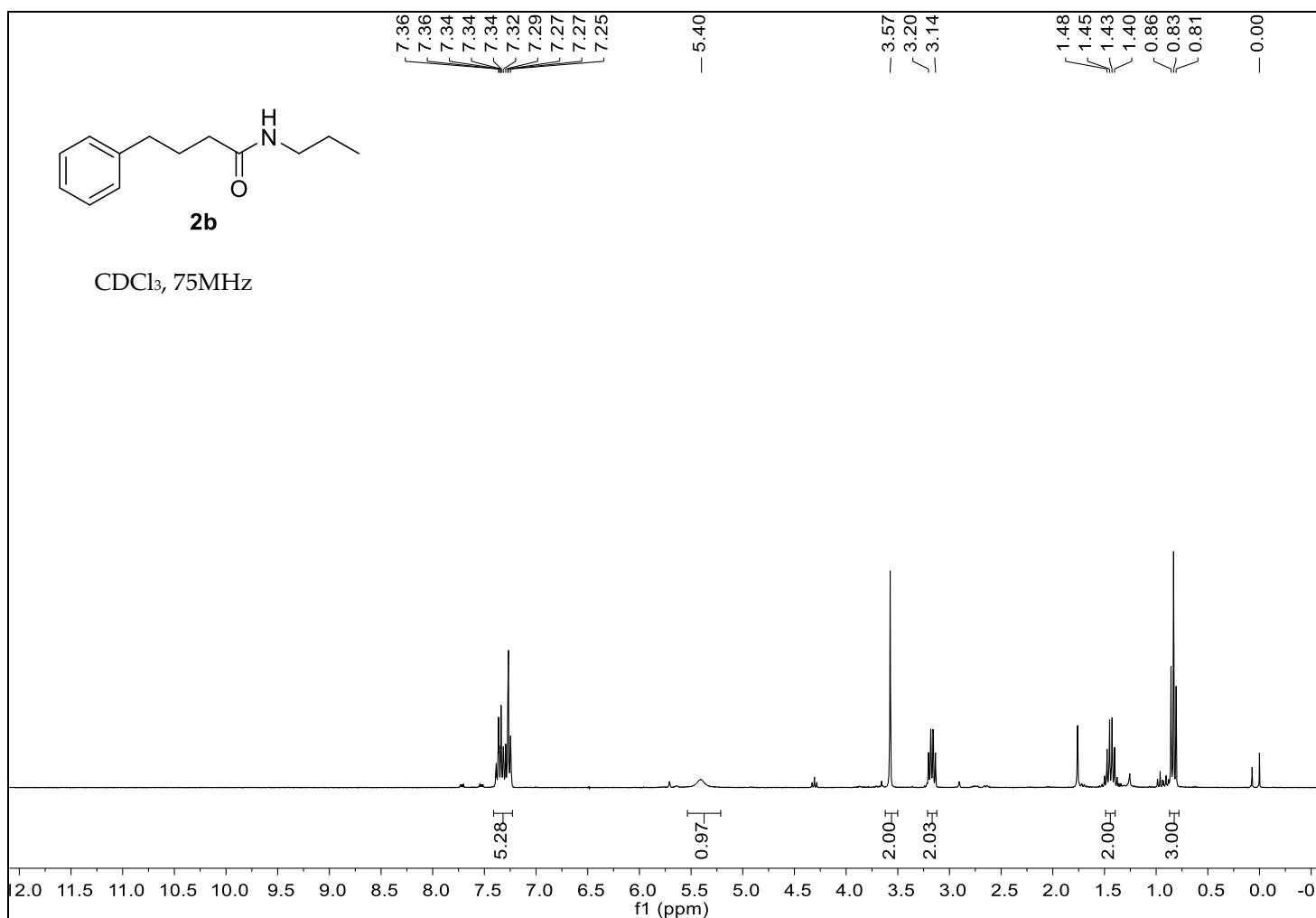


Figure S5. ¹H-NMR of 2-Phenyl-N-propylacetamide (2c).

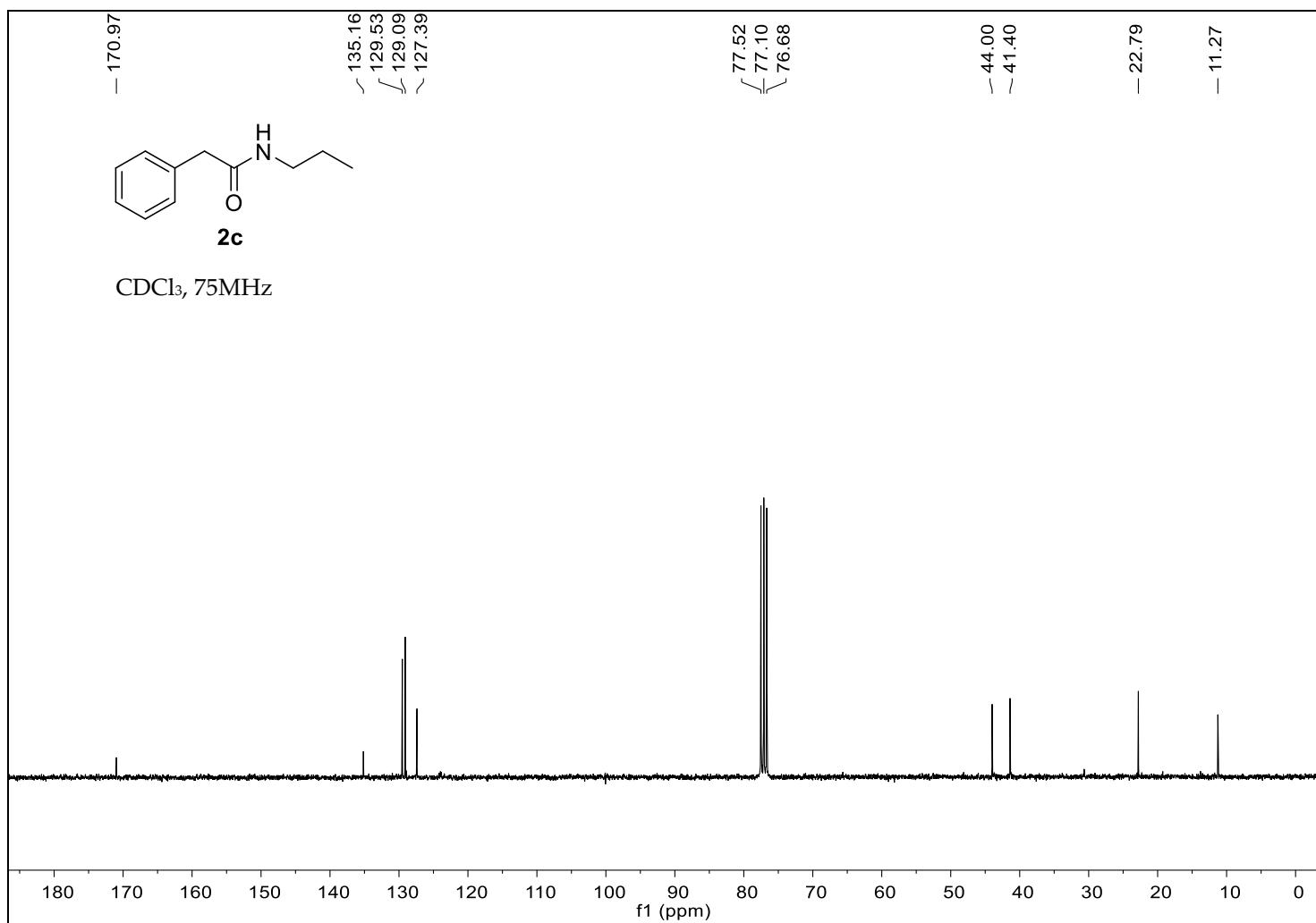


Figure S6. ¹³C-NMR of 2-Phenyl-N-propylacetamide (**2c**).

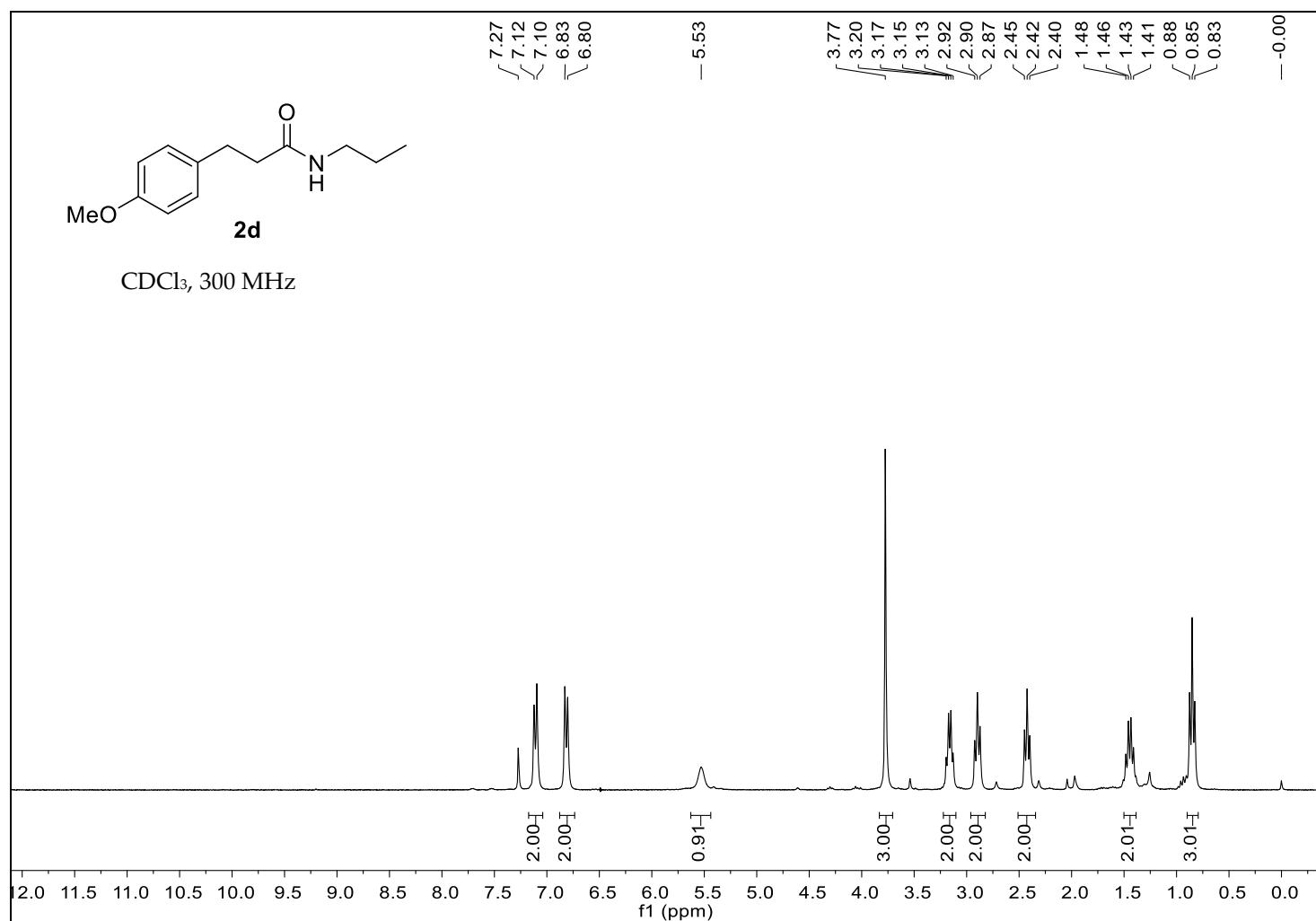


Figure S7. ¹H-NMR of 3-(4-Methoxyphenyl)-N-propylpropanamide (**2d**).

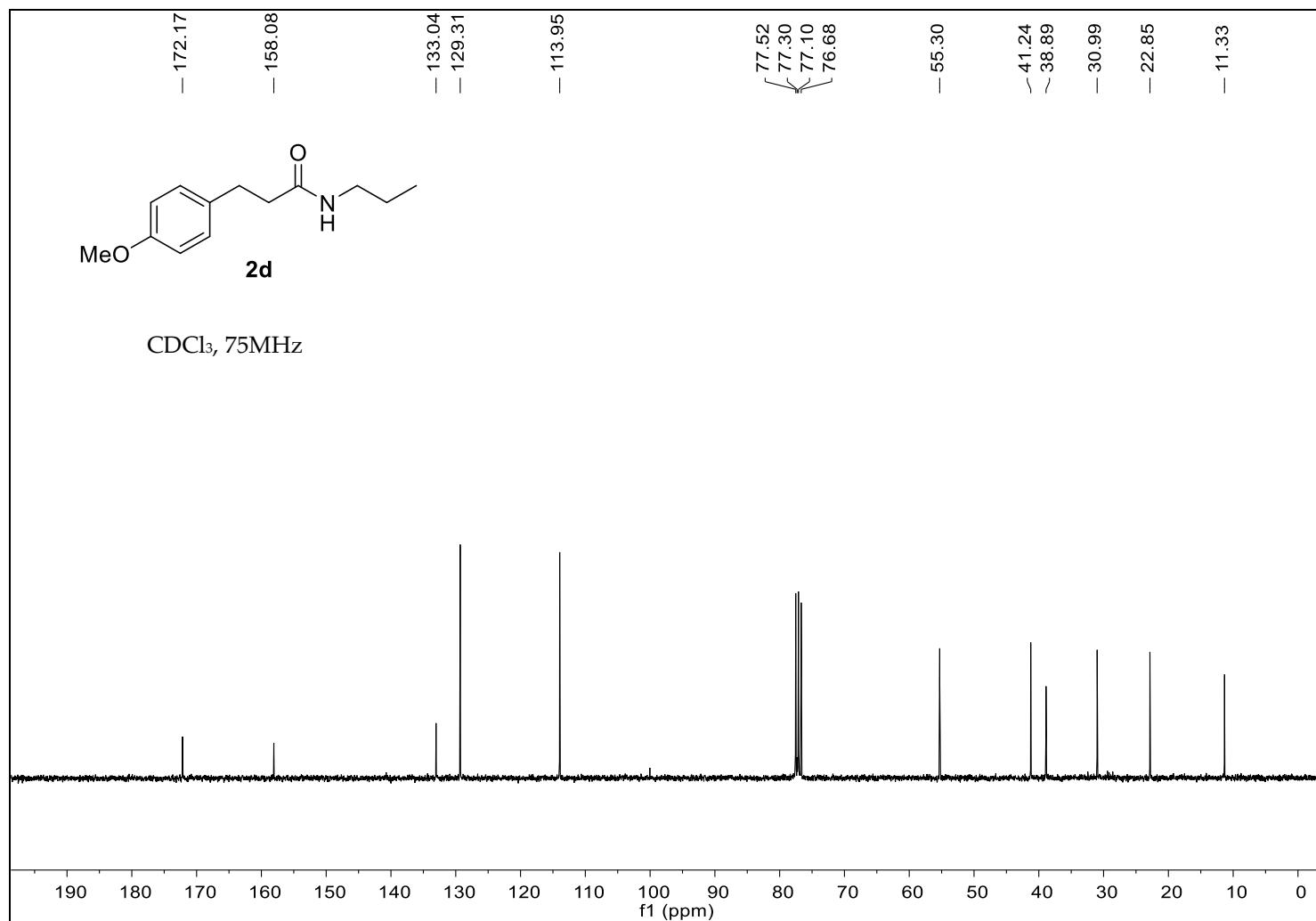


Figure S8. ¹³C-NMR of 3-(4-Methoxyphenyl)-N-propylpropanamide (**2d**).

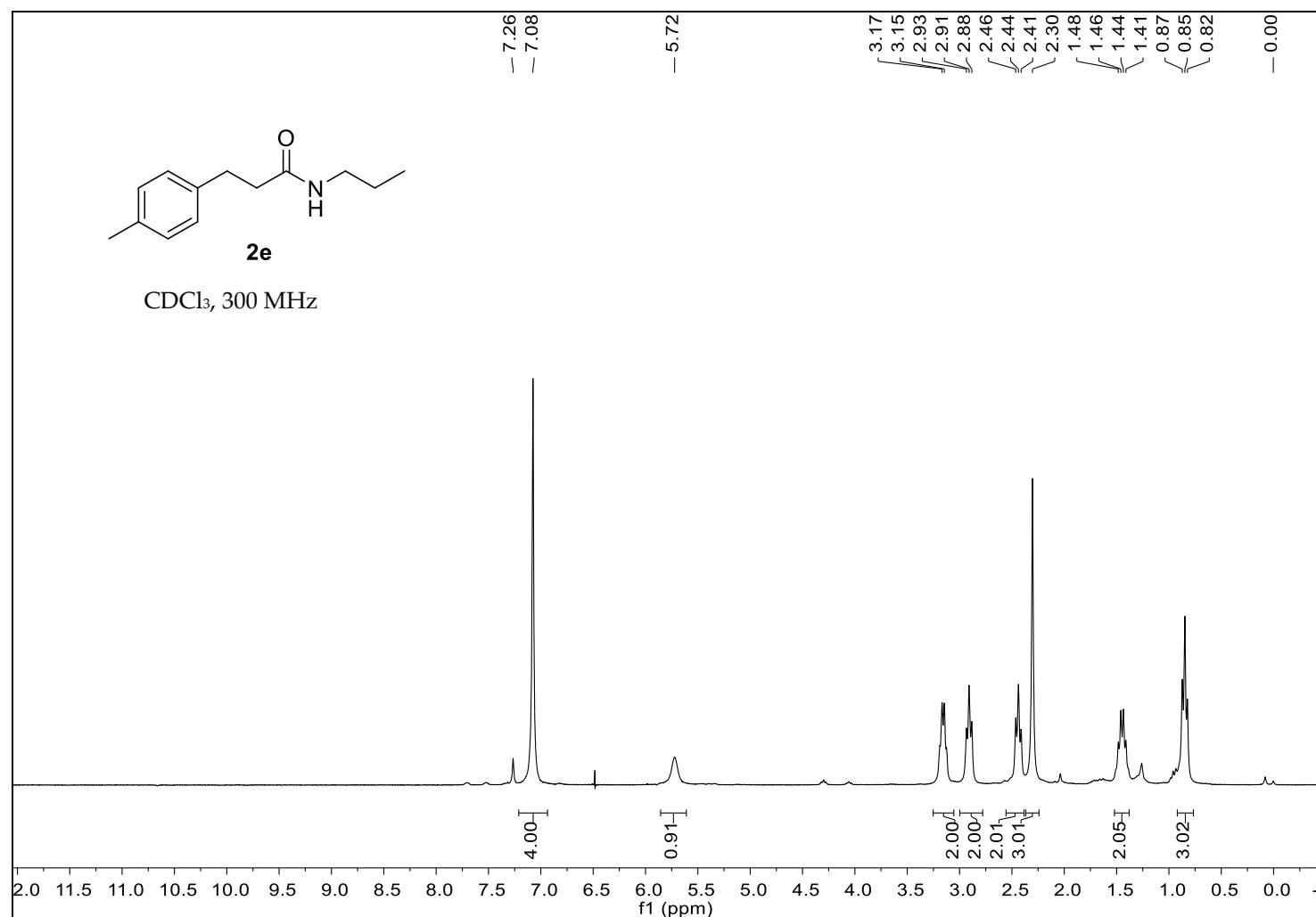


Figure S9. ¹H-NMR of *N*-Propyl-3-(*p*-tolyl)propanamide (**2e**).

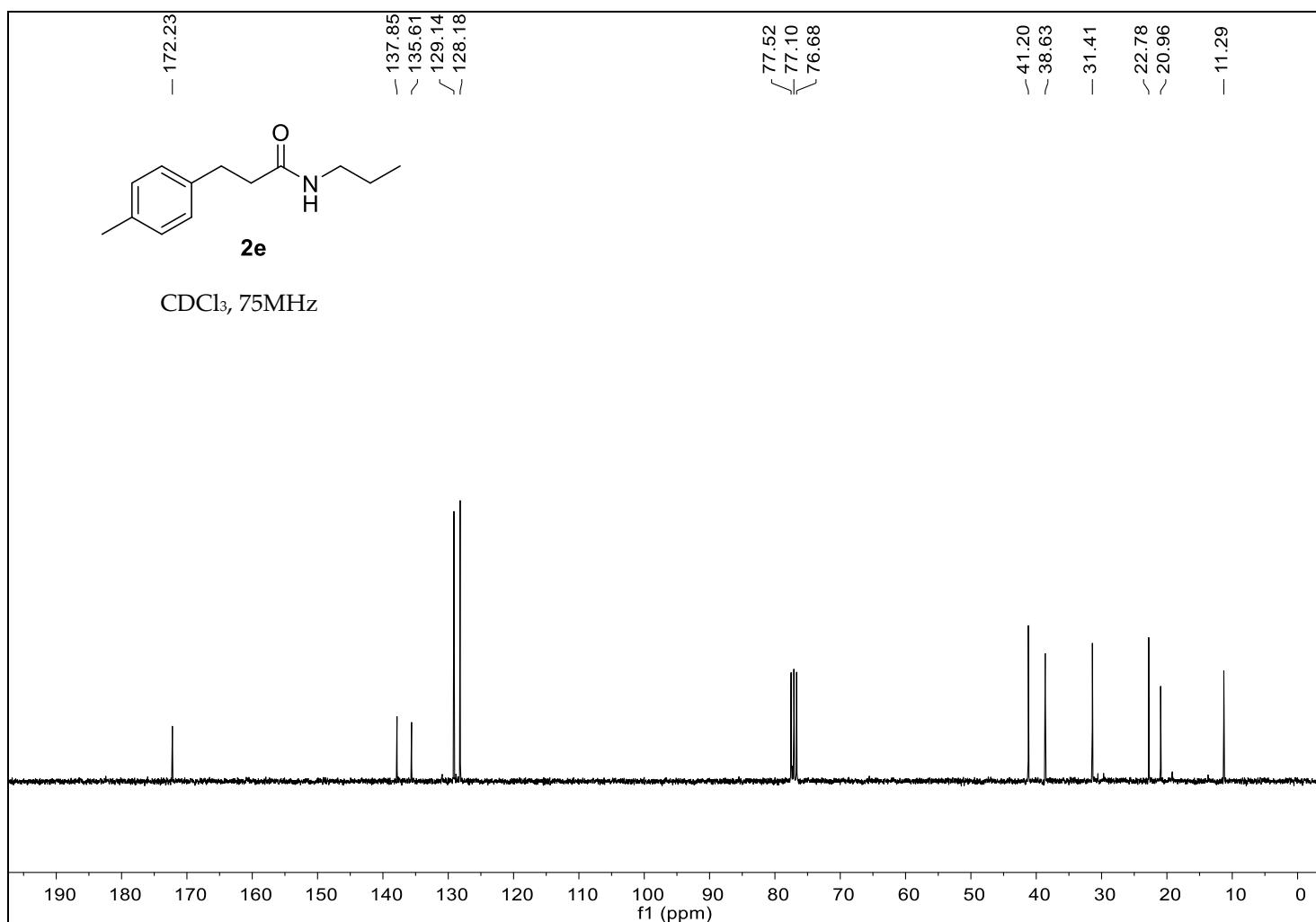


Figure S10. ^{13}C -NMR of *N*-Propyl-3-(*p*-tolyl)propanamide (**2e**).

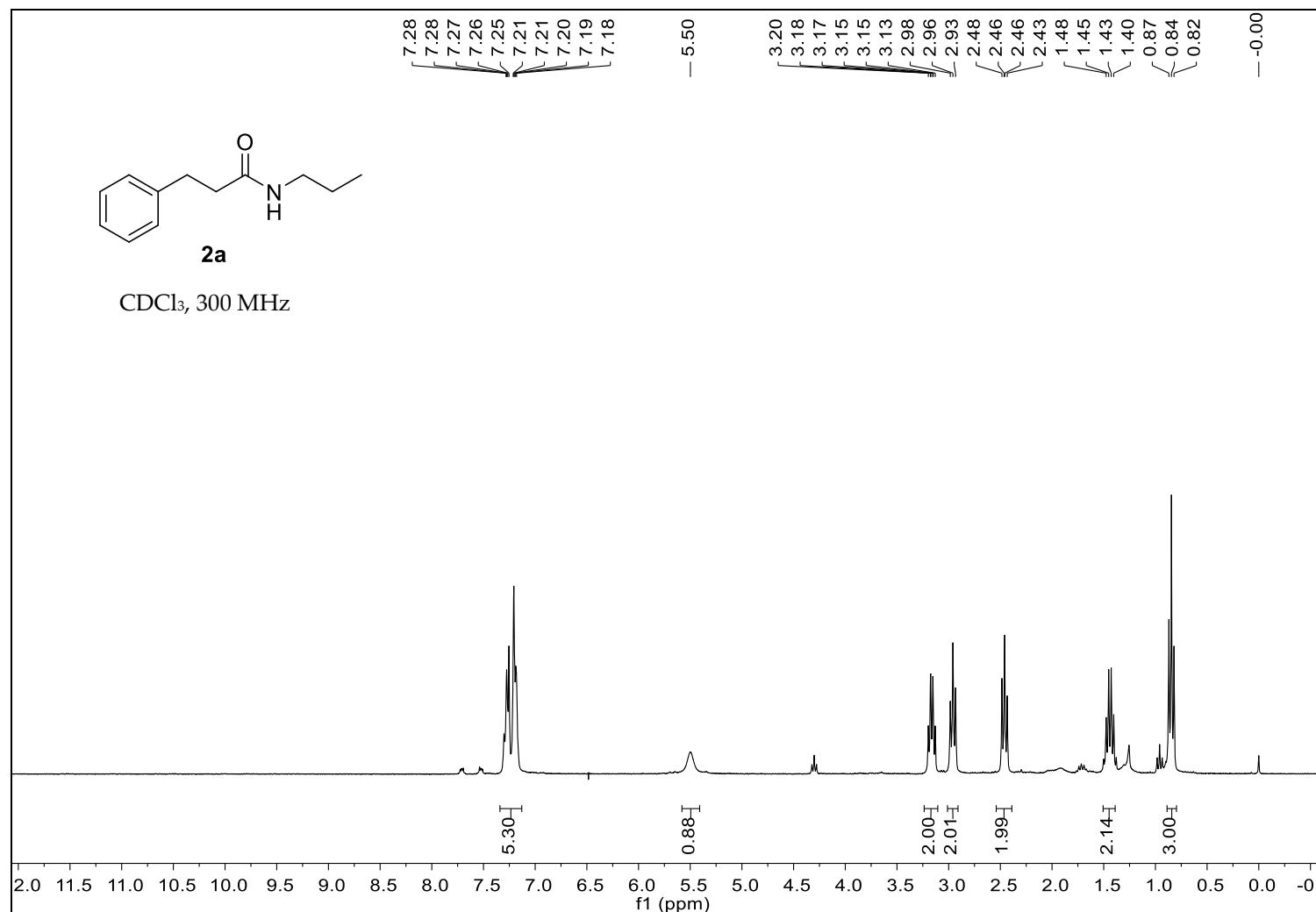


Figure S11. ¹H-NMR of 3-Phenyl-N-propylpropanamide ((2a) (derived from 1f)).

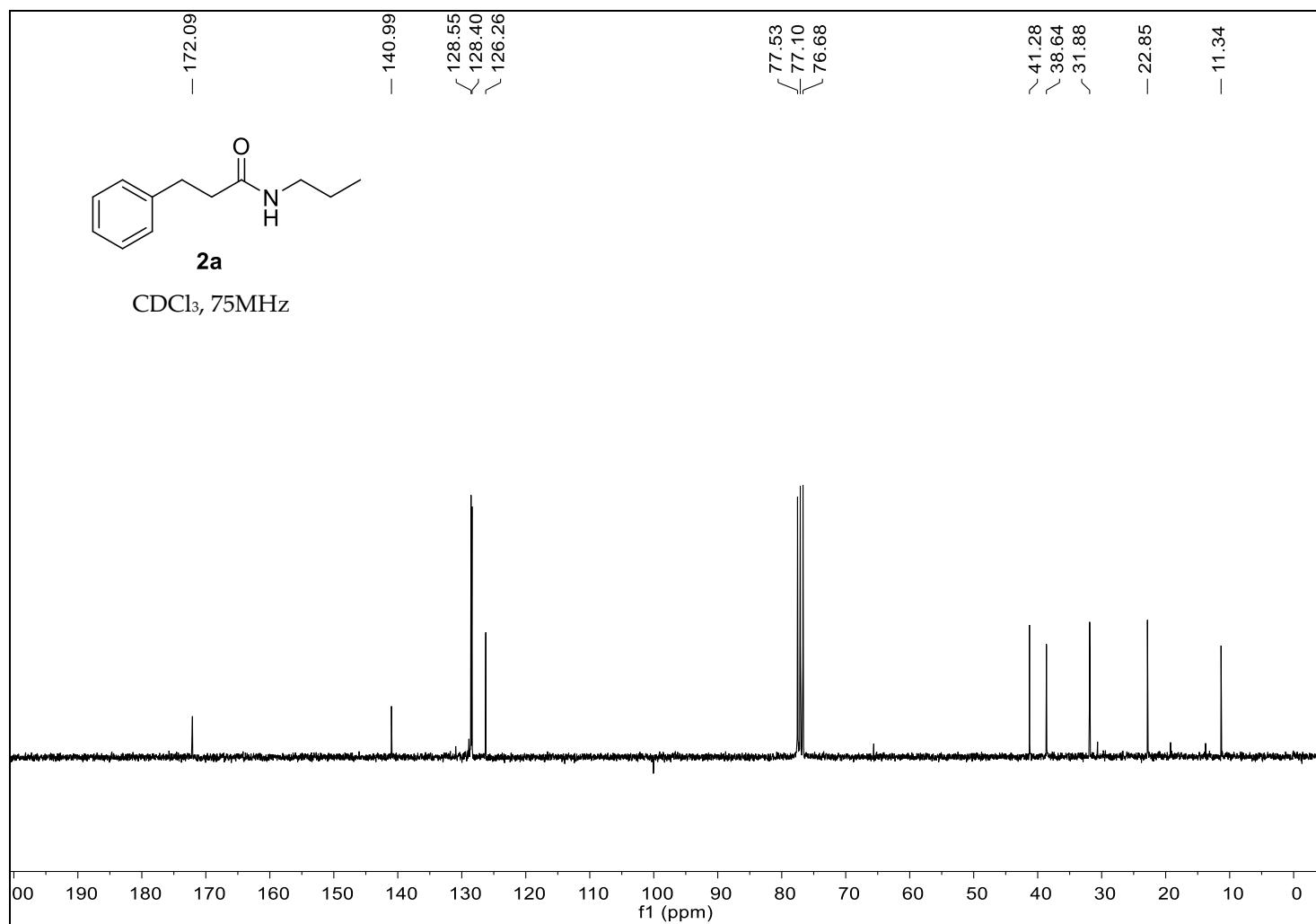


Figure S12. ^{13}C -NMR of 3-Phenyl-*N*-propylpropanamide ((**2a**) (derived from **1f**).

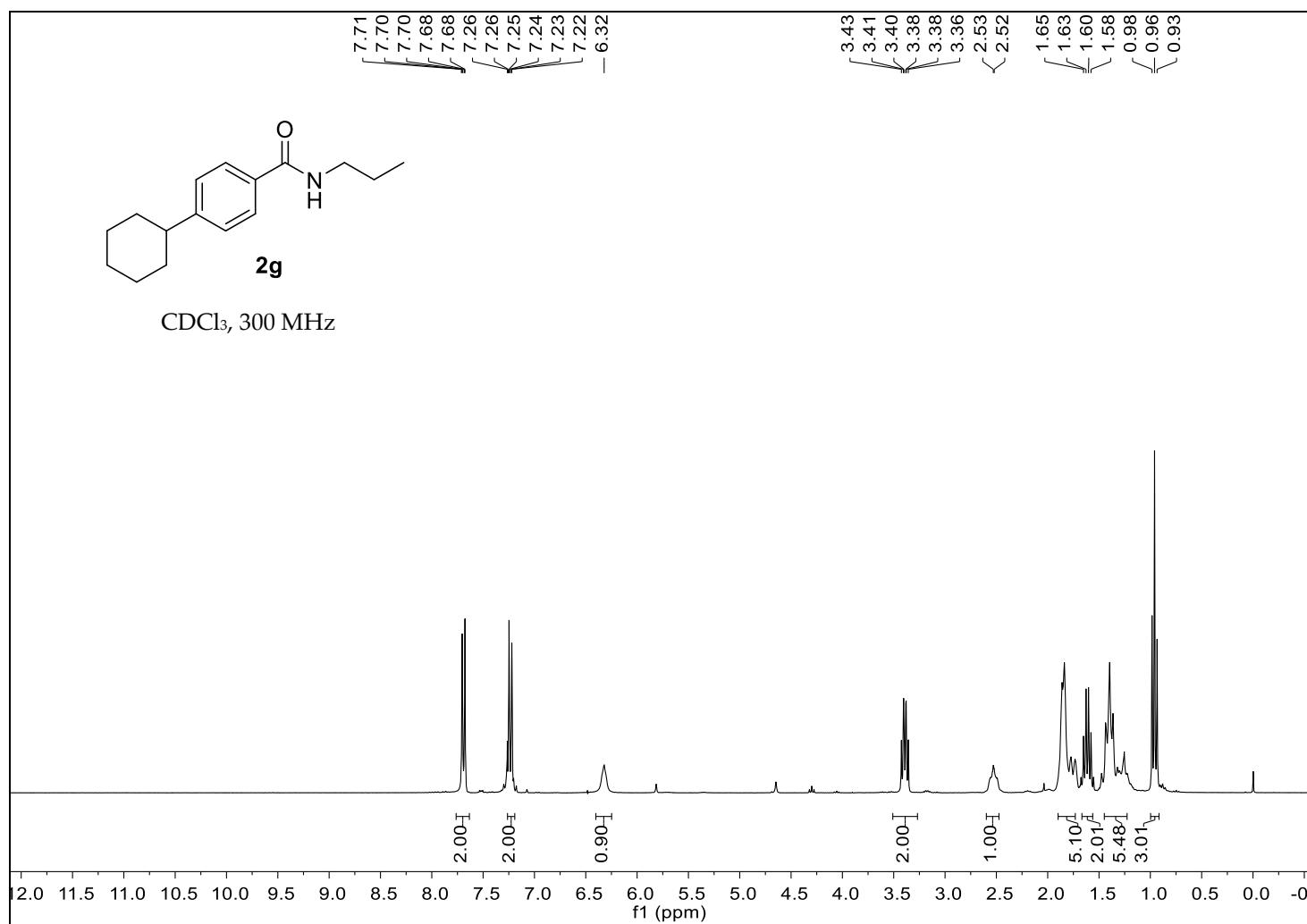


Figure S13. ¹H-NMR of 4-Cyclohexyl-N-propylbenzamide (2g).

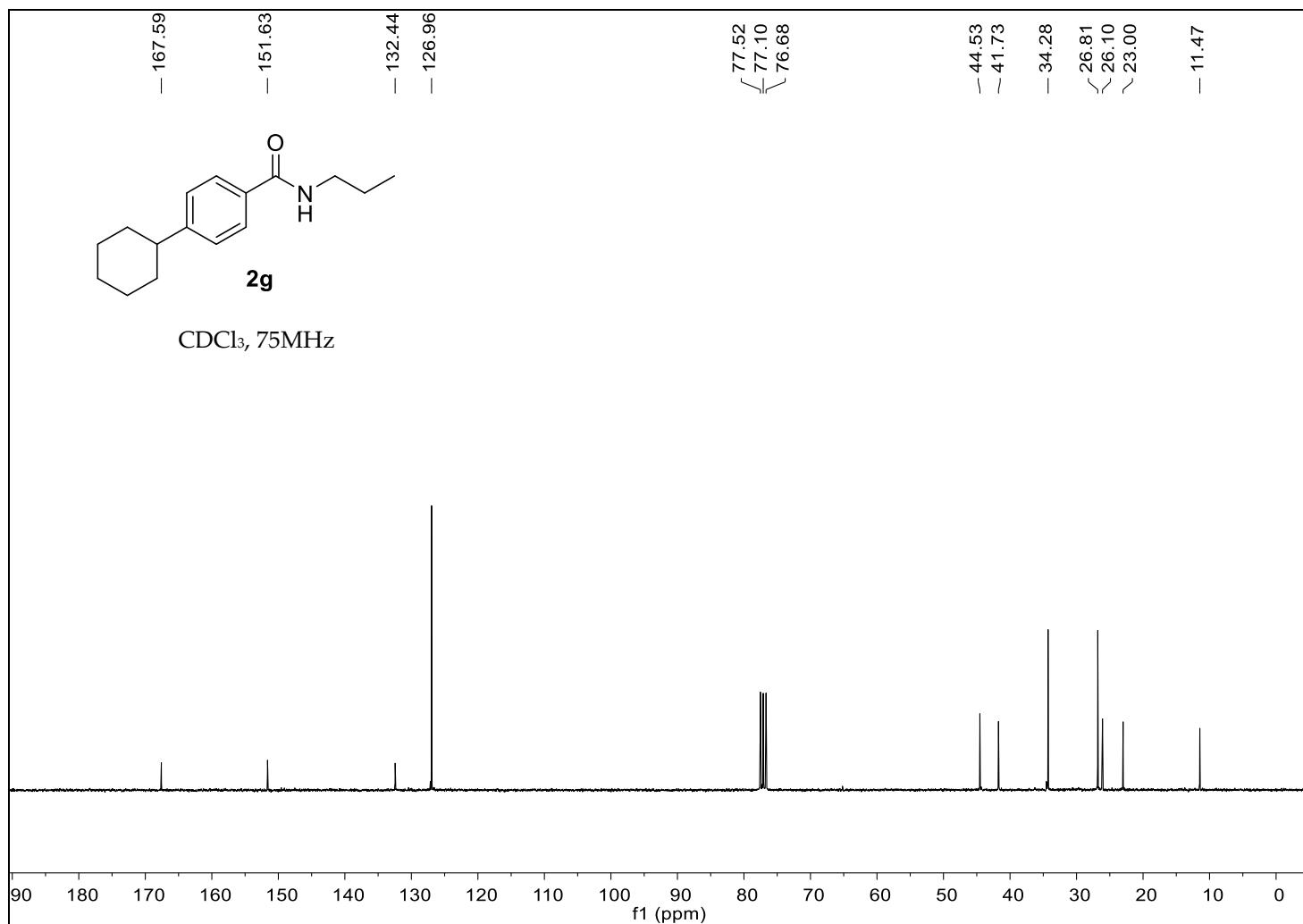


Figure S14. ^{13}C -NMR of 4-Cyclohexyl-N-propylbenzamide (**2g**).

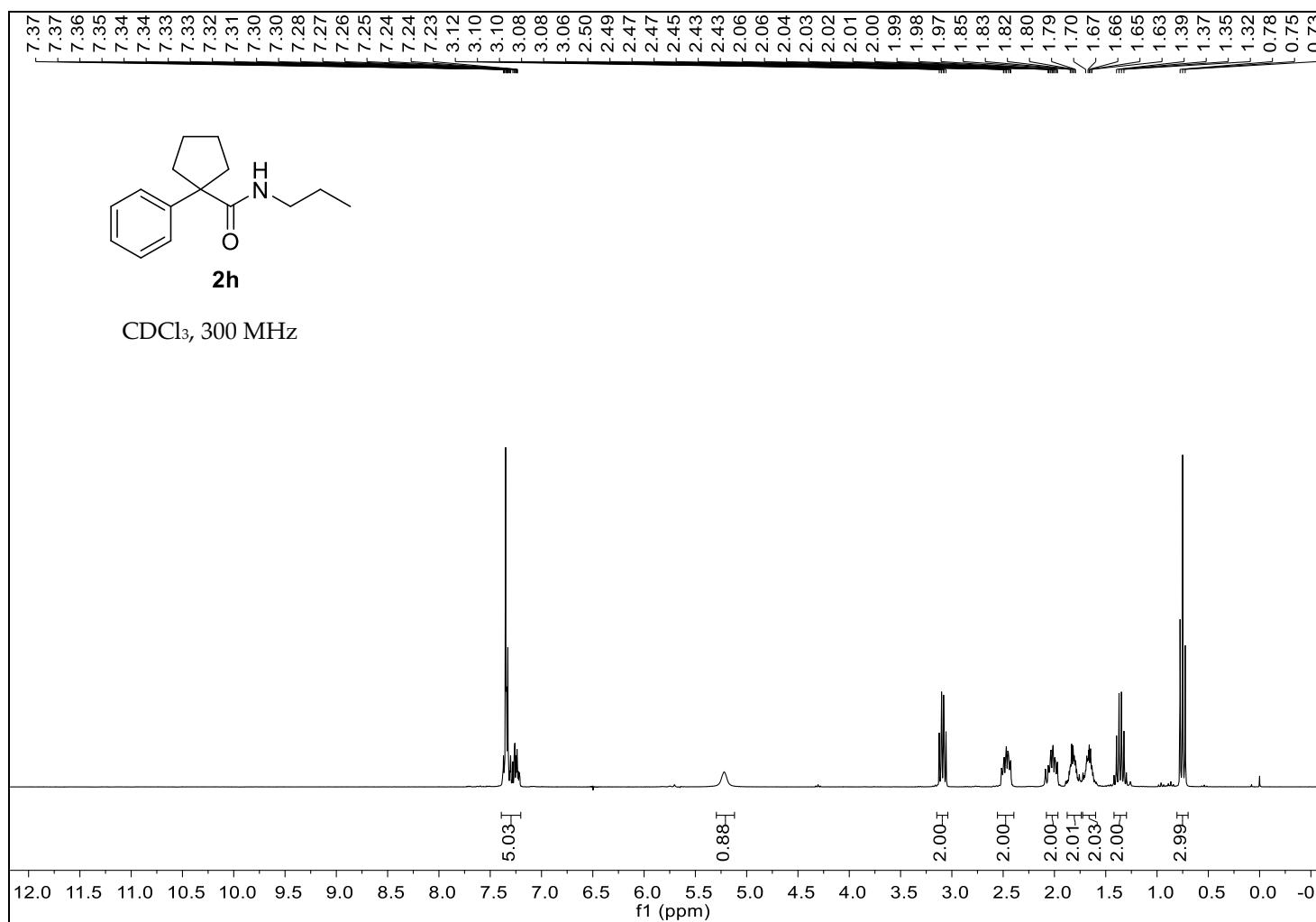


Figure S15. ¹H-NMR of 1-Phenyl-N-propylcyclopentane-1-carboxamide (**2h**).

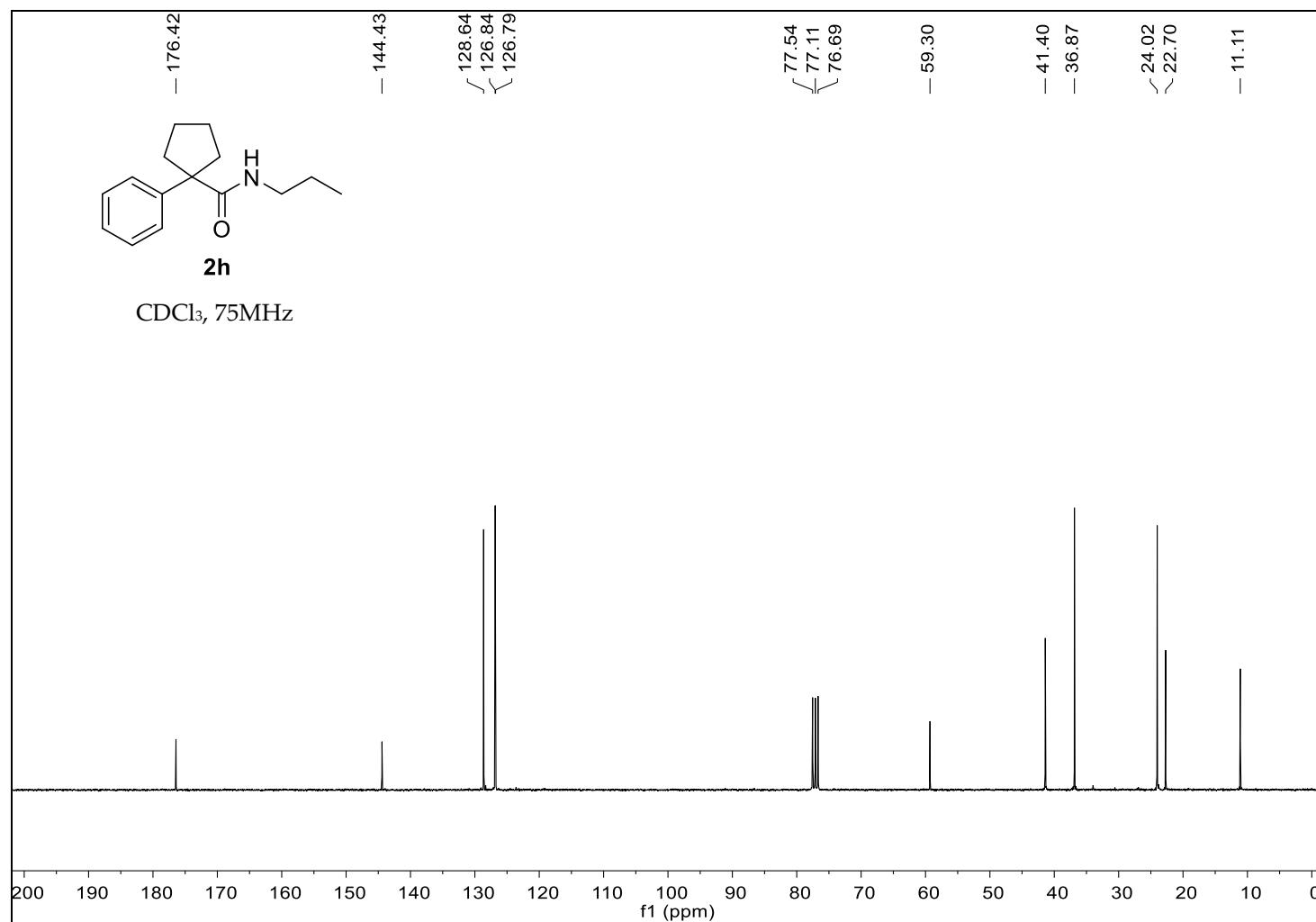


Figure S16. ¹³C-NMR of 1-Phenyl-N-propylcyclopentane-1-carboxamide (**2h**).

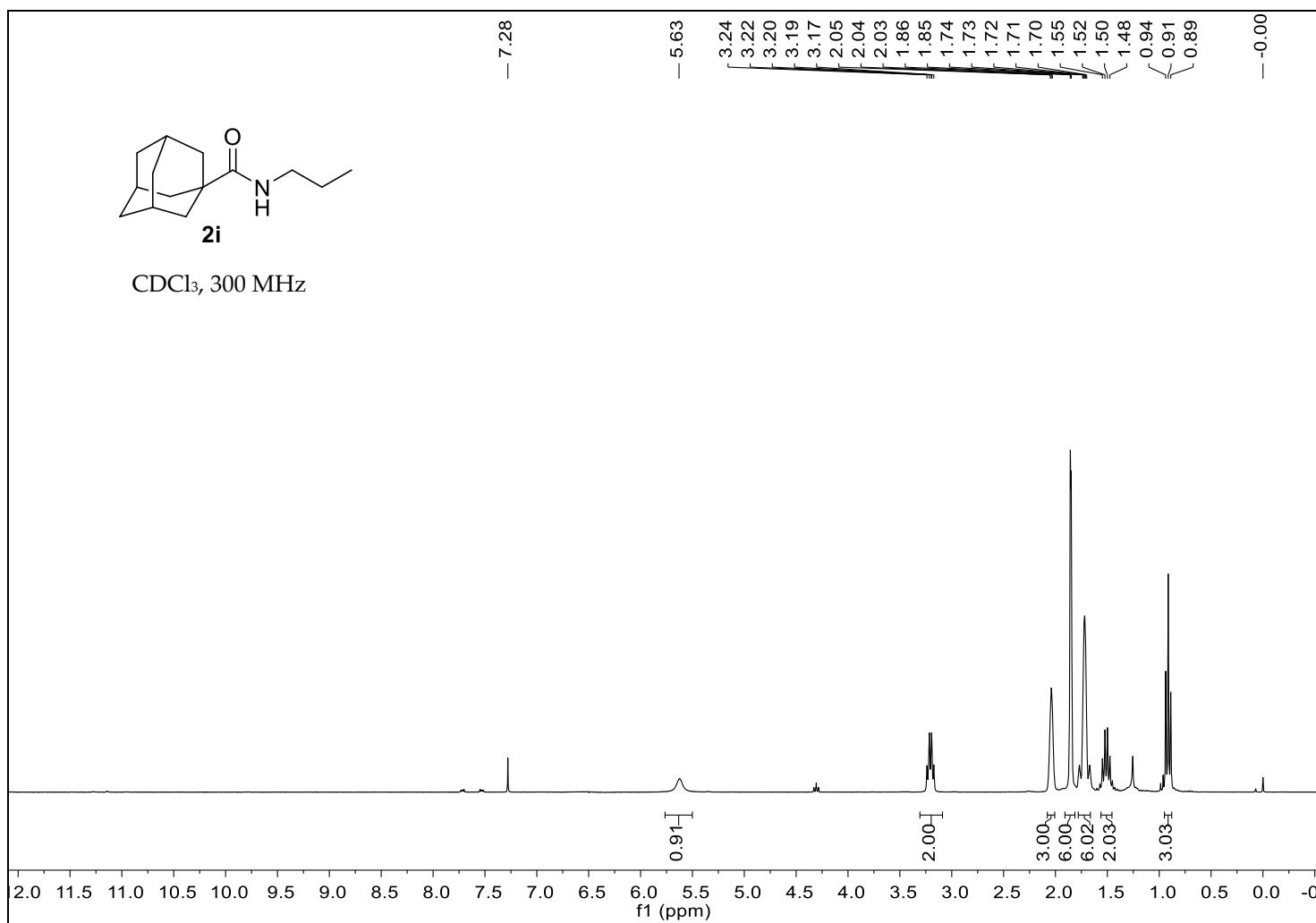


Figure S17. ¹H-NMR of (3*r*,5*r*,7*r*)-*N*-Propyladamantane-1-carboxamide (**2i**).

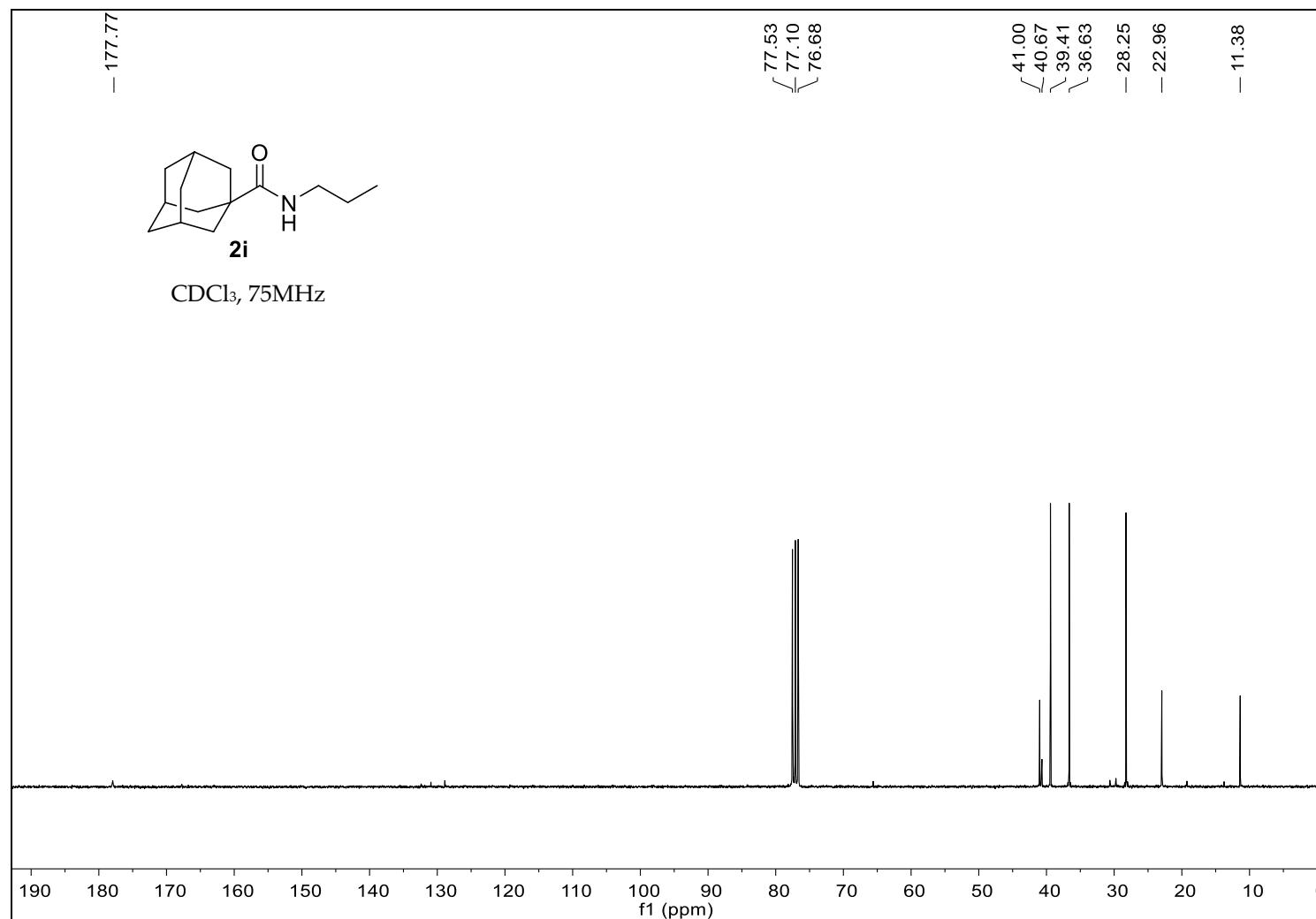


Figure S18. ^{13}C -NMR of (3r,5r,7r)-N-Propyladamantane-1-carboxamide (**2i**).

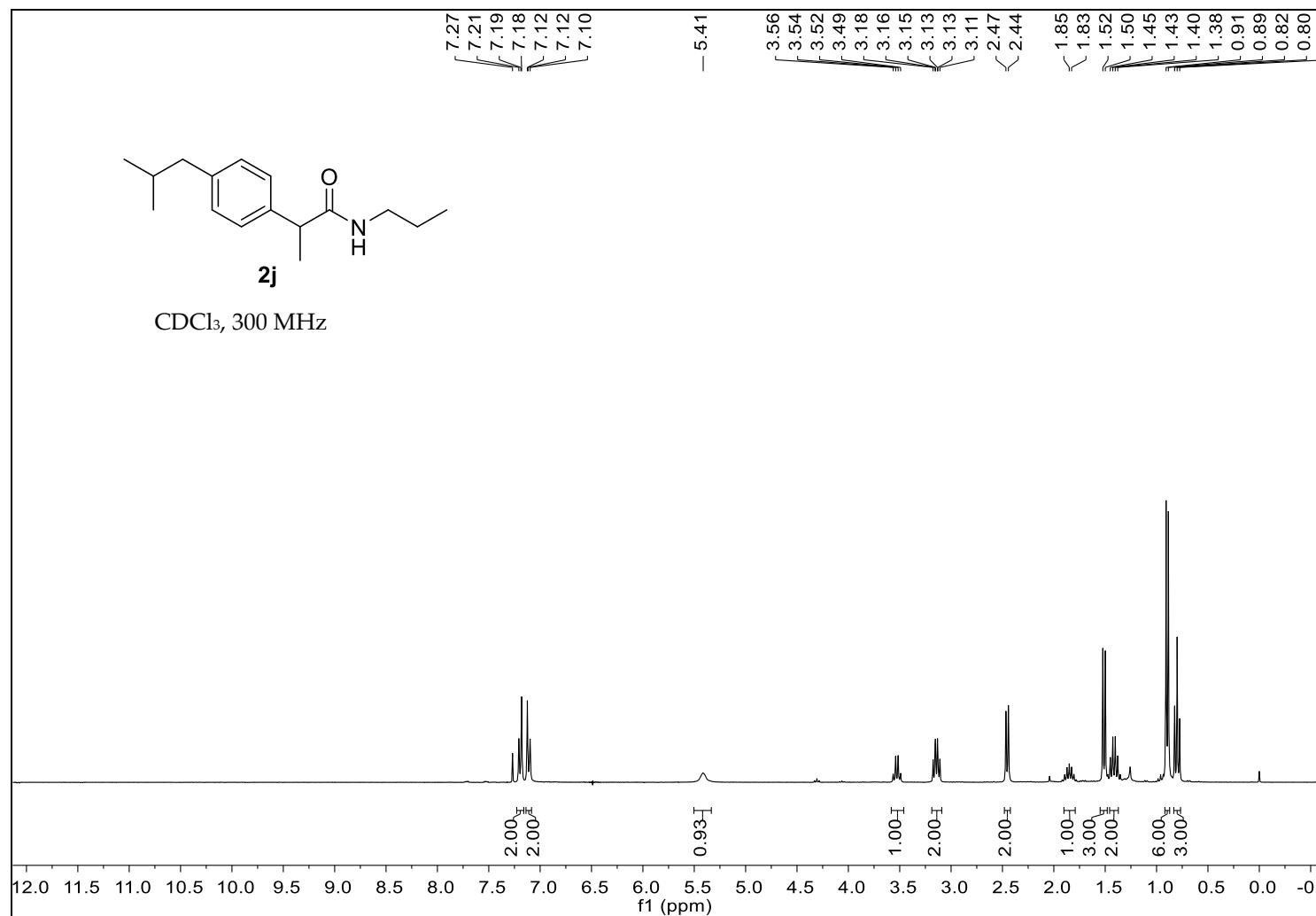


Figure S19. ^1H -NMR of 2-(4-Isobutylphenyl)-*N*-propylpropanamide (**2j**).

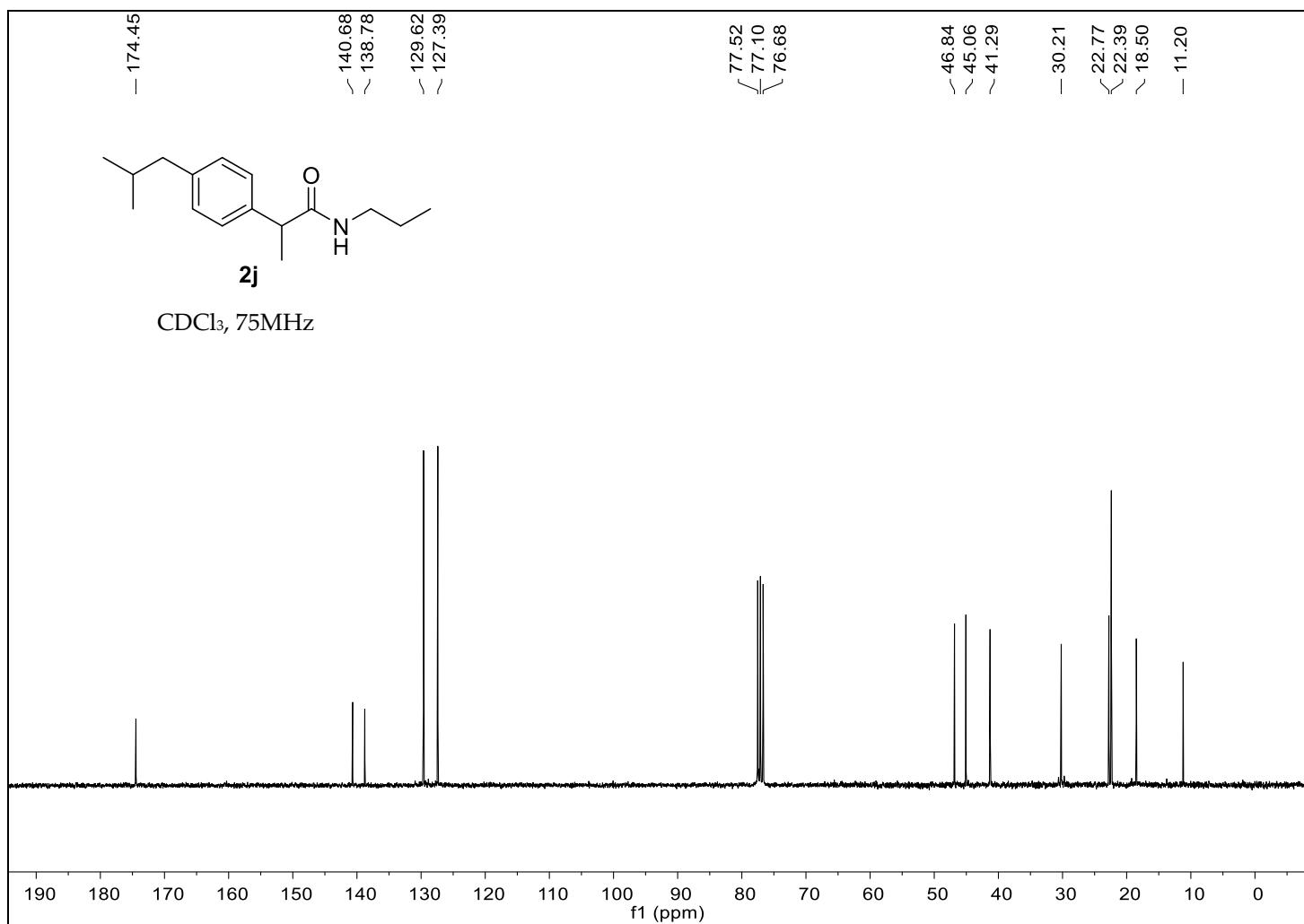


Figure S20. ^{13}C -NMR of 2-(4-Isobutylphenyl)-N-propylpropanamide (**2j**).

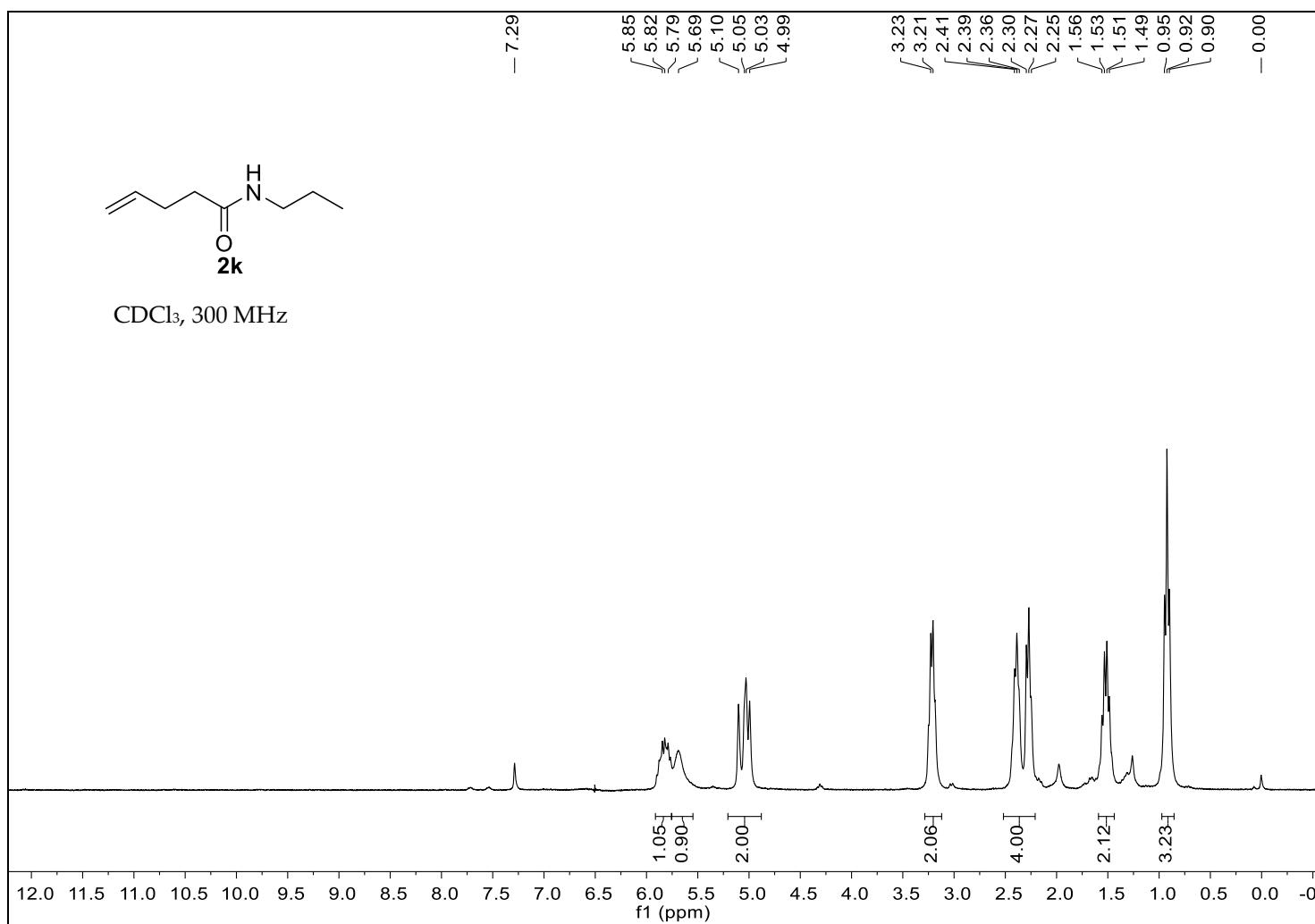


Figure S21. ^1H -NMR of *N*-Propylpent-4-enamide (**2k**).

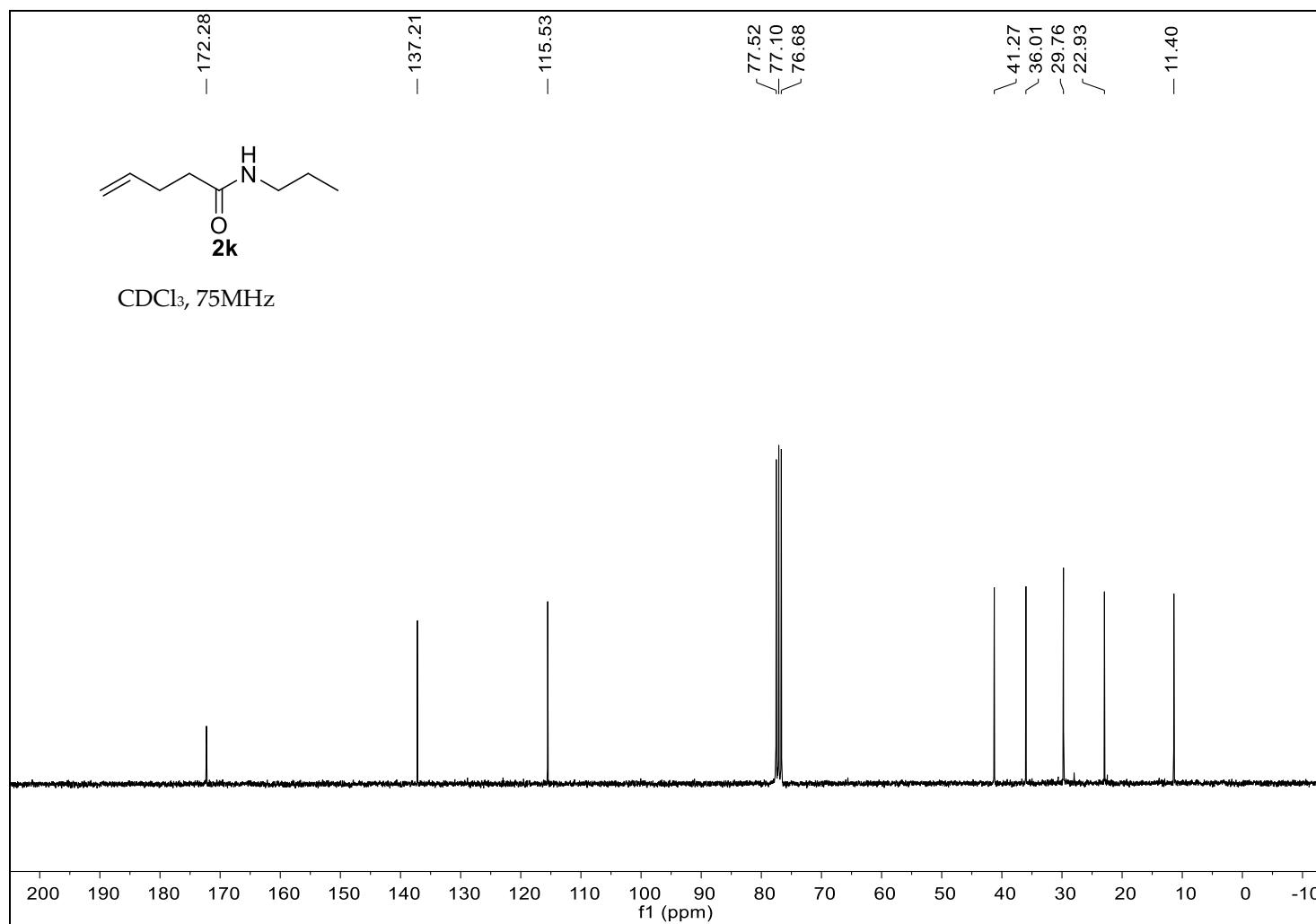


Figure S22. ^{13}C -NMR of *N*-Propylpent-4-enamide (**2k**).

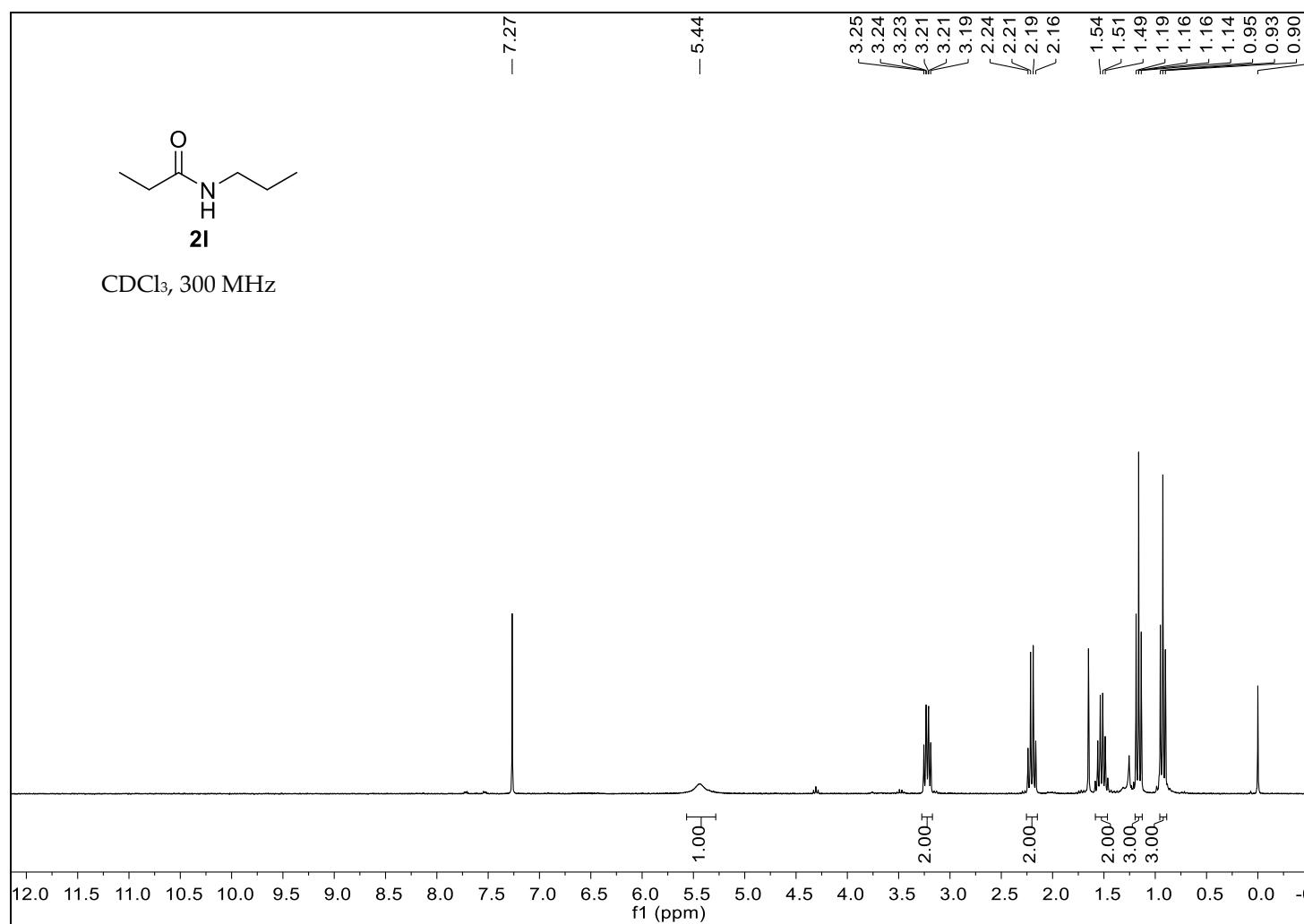


Figure S23. ^1H -NMR of *N*-Propylpropionamide (**2l**).

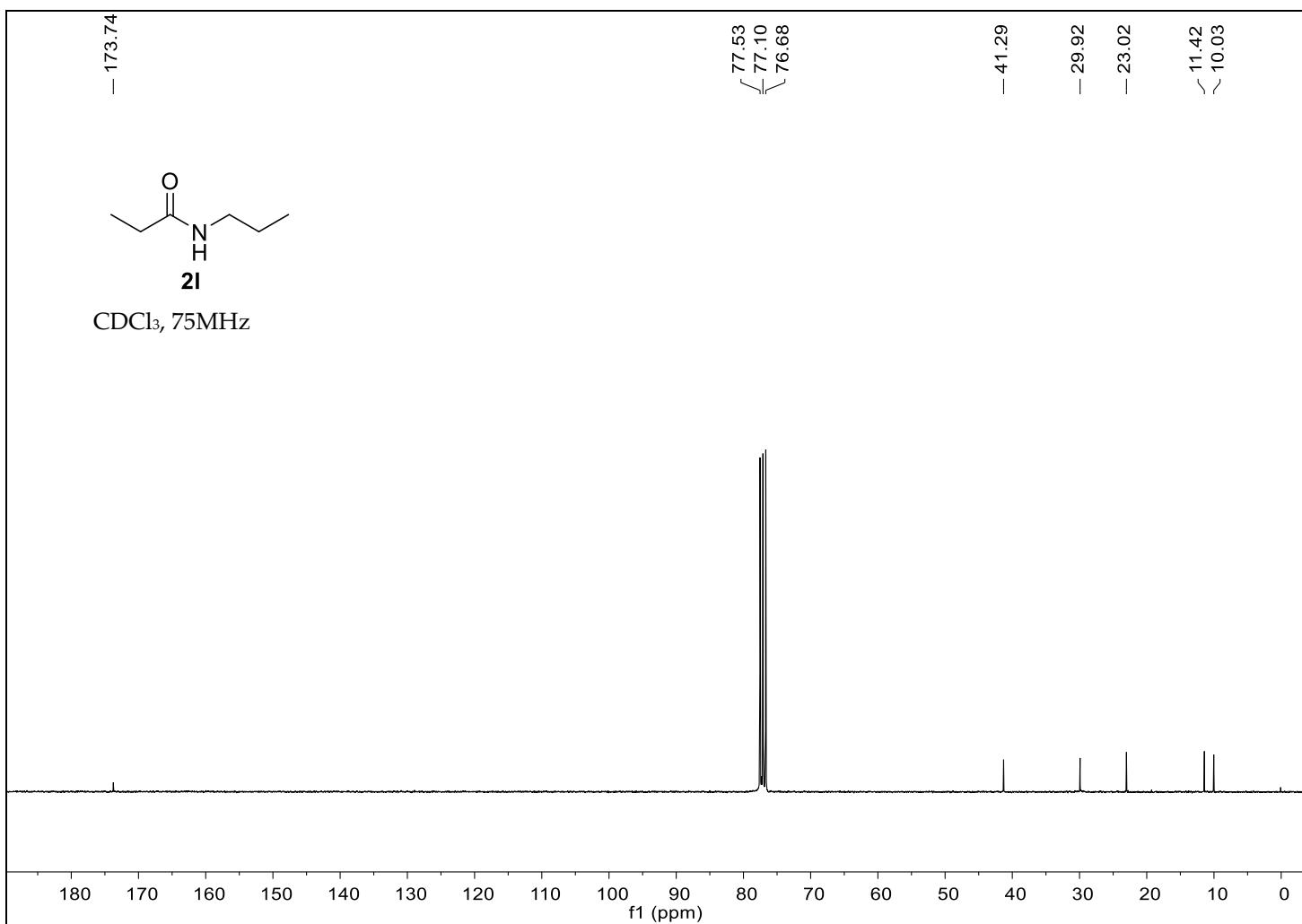


Figure S24. ^{13}C -NMR of *N*-Propylpropionamide (**2l**).

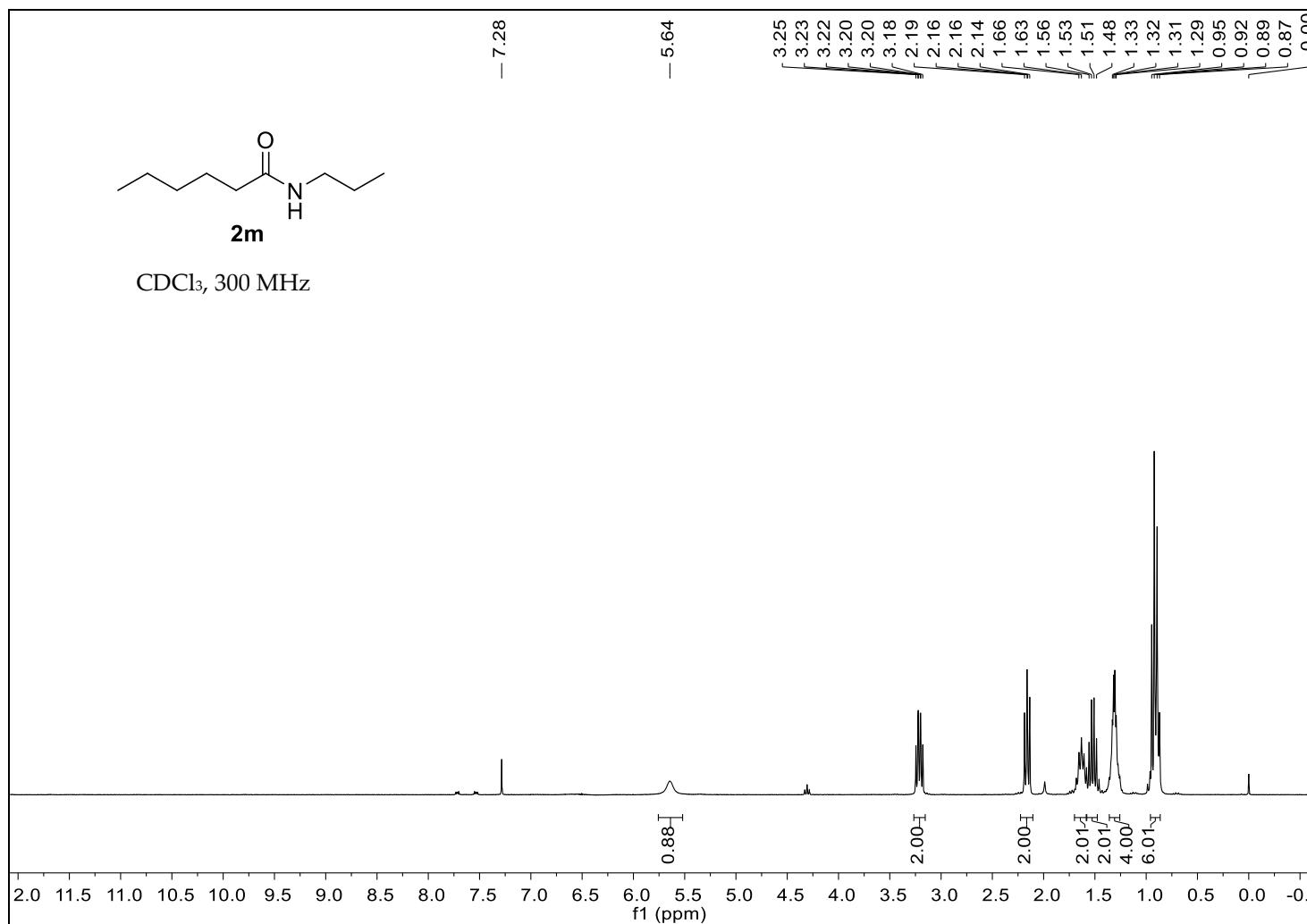


Figure S25. ¹H-NMR of *N*-Propylhexanamide (**2m**).

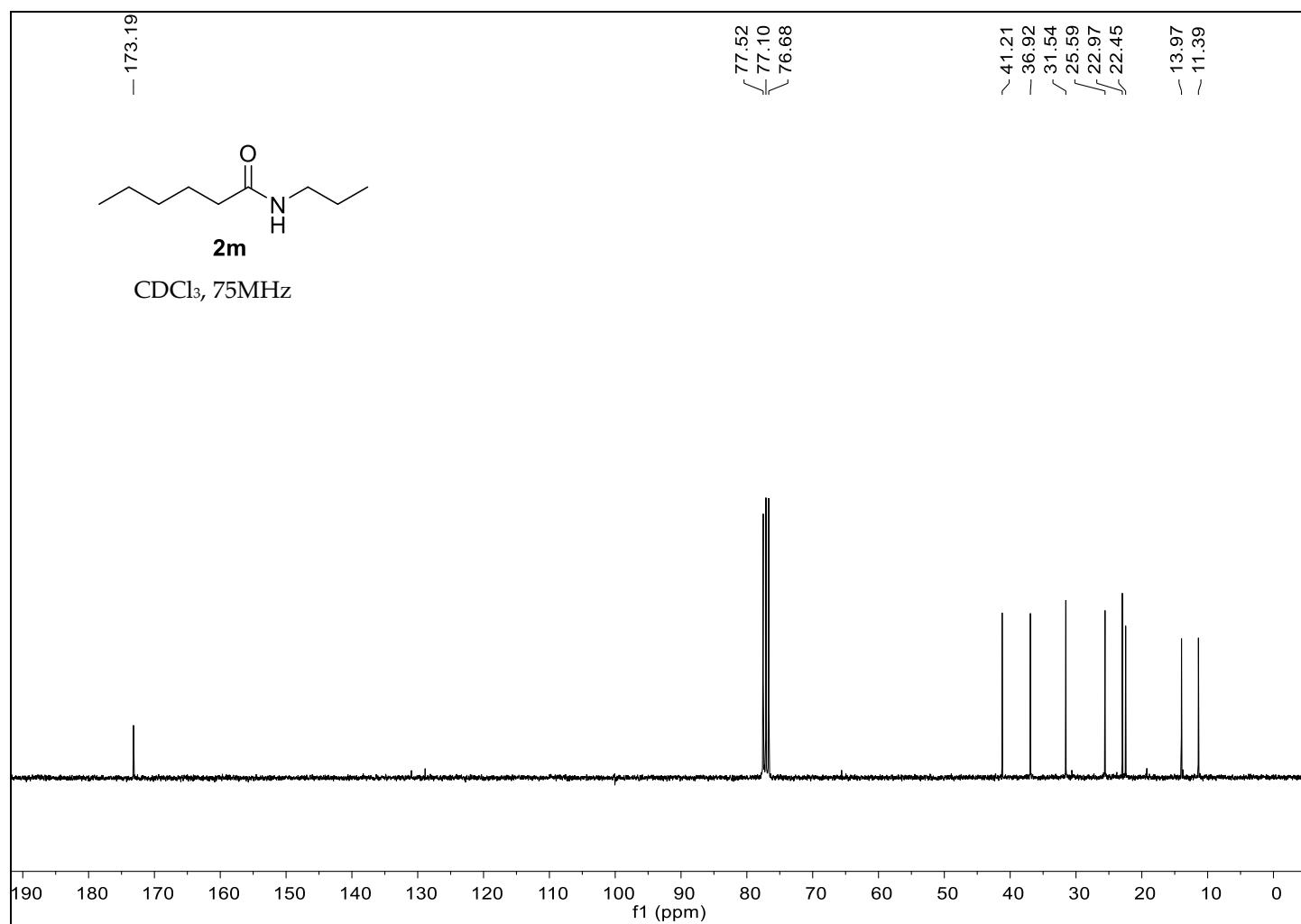


Figure S26. ^{13}C -NMR of *N*-Propylhexanamide (**2m**).

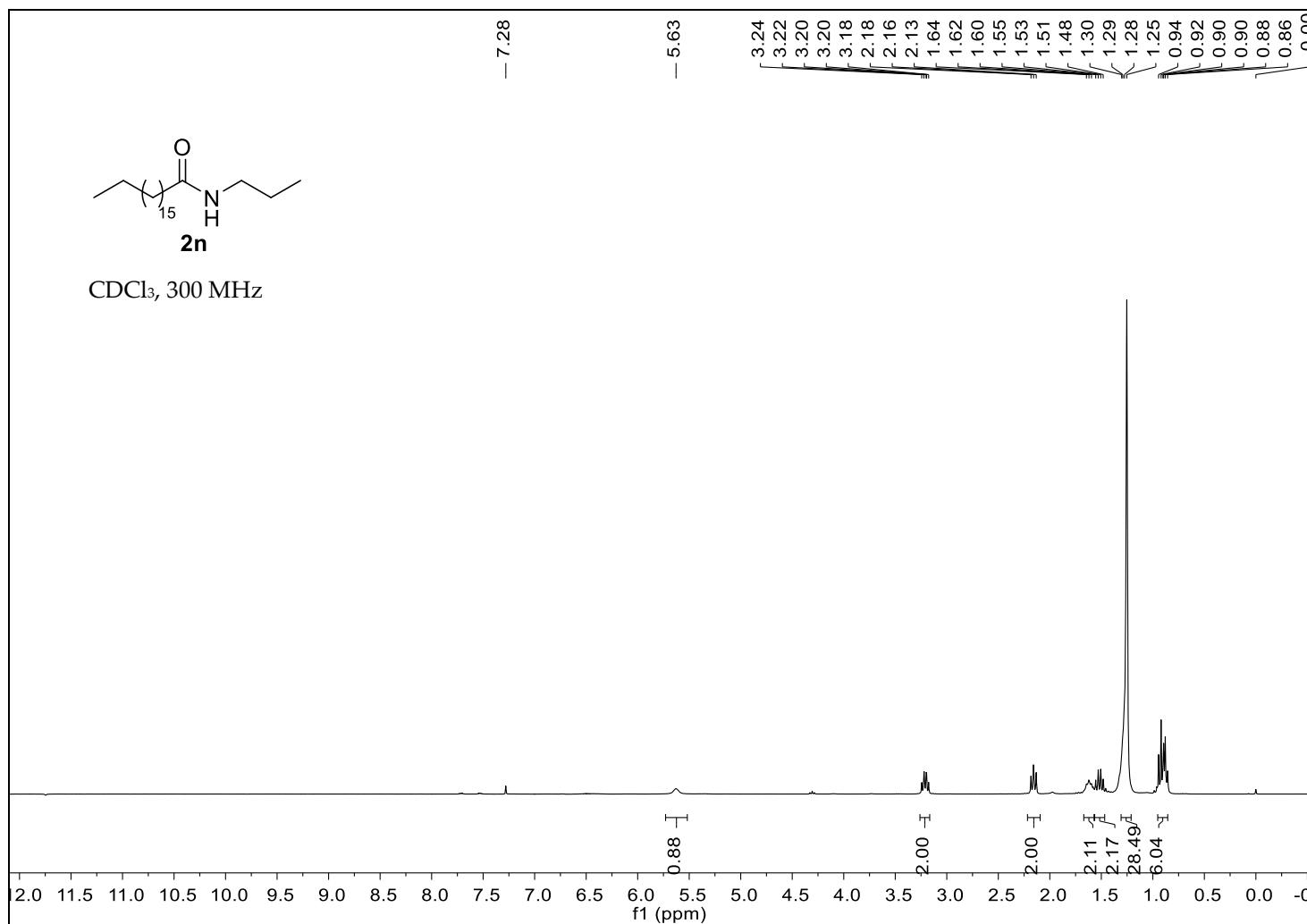


Figure S27. ^1H -NMR of *N*-Propylstearamide (**2n**).

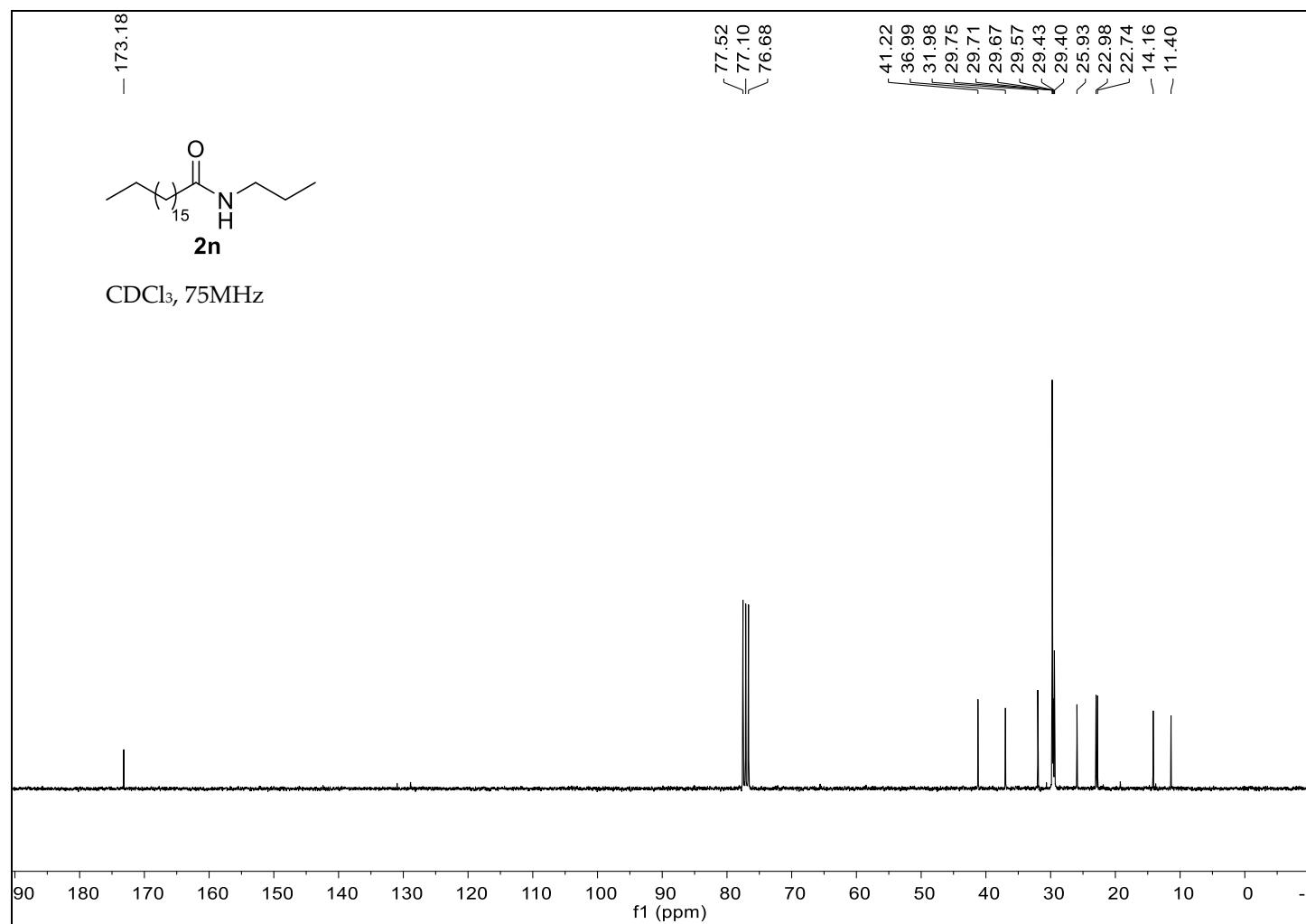


Figure S28. ^{13}C -NMR of *N*-Propylstearamide (**2n**).

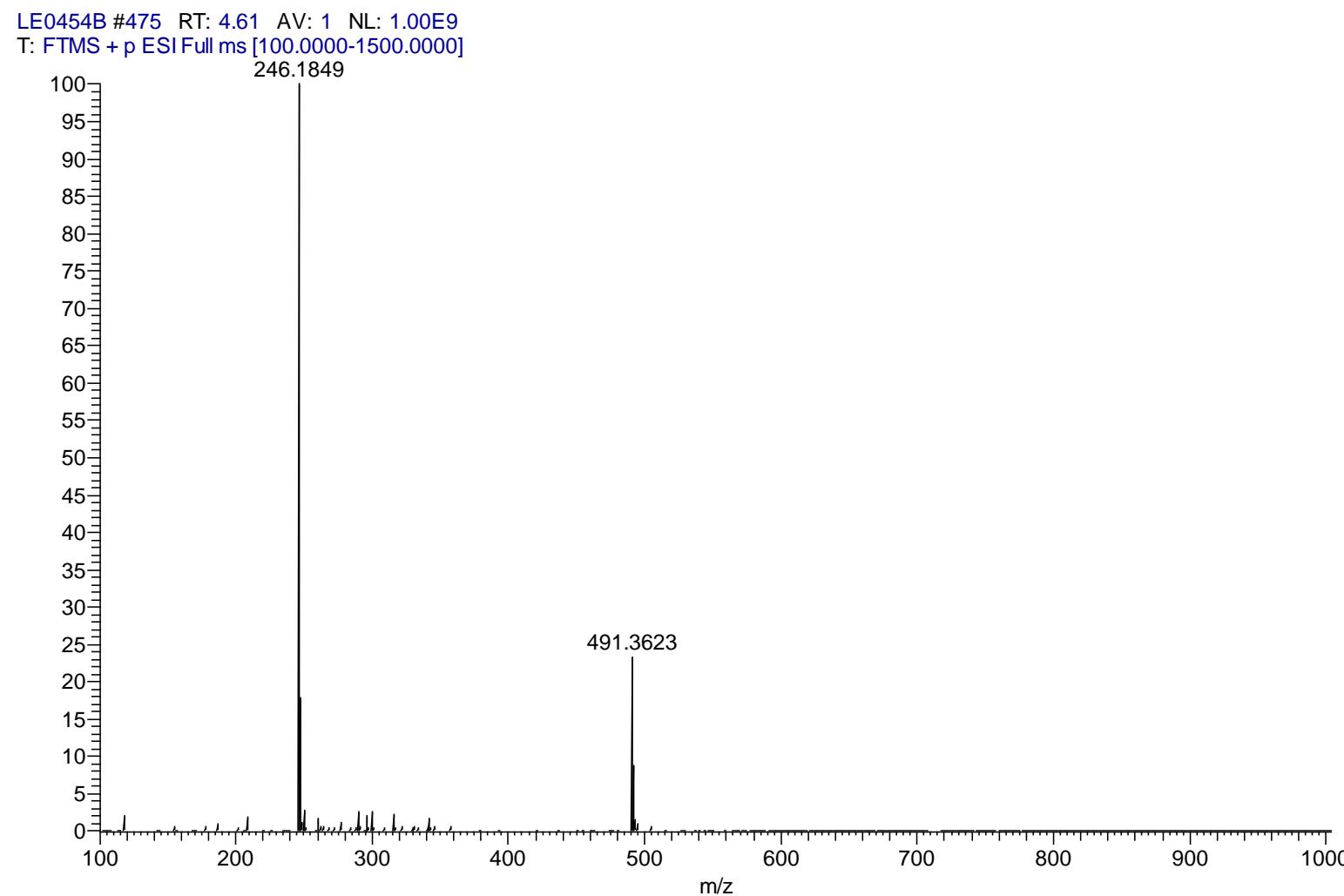


Figure S29. HRMS of 4-Cyclohexyl-N-propylbenzamide (**2g**).

Le0459C #44-81 RT: 0.20-0.36 AV: 38 NL: 5.29E6
T: FTMS + p ESI Full ms [100.0000-500.0000]

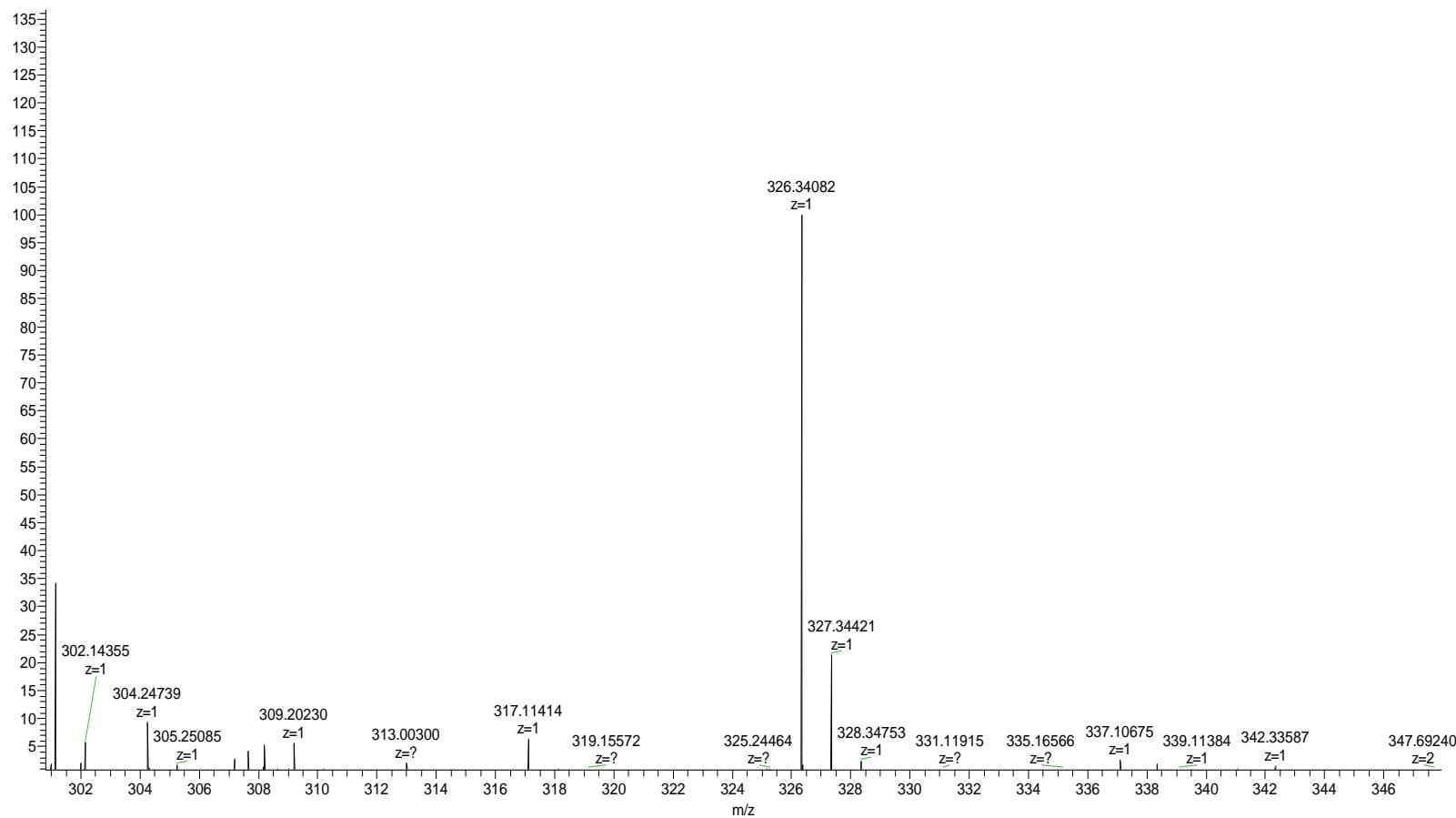


Figure S30. HRMS of *N*-Propylstearamide (**2n**).

