

## Supporting Information

### The Thermofluoric Behavior of Poly(fluorenetolyldiphenylamine)-Oxadiazole Pair in Polymer Matrix.

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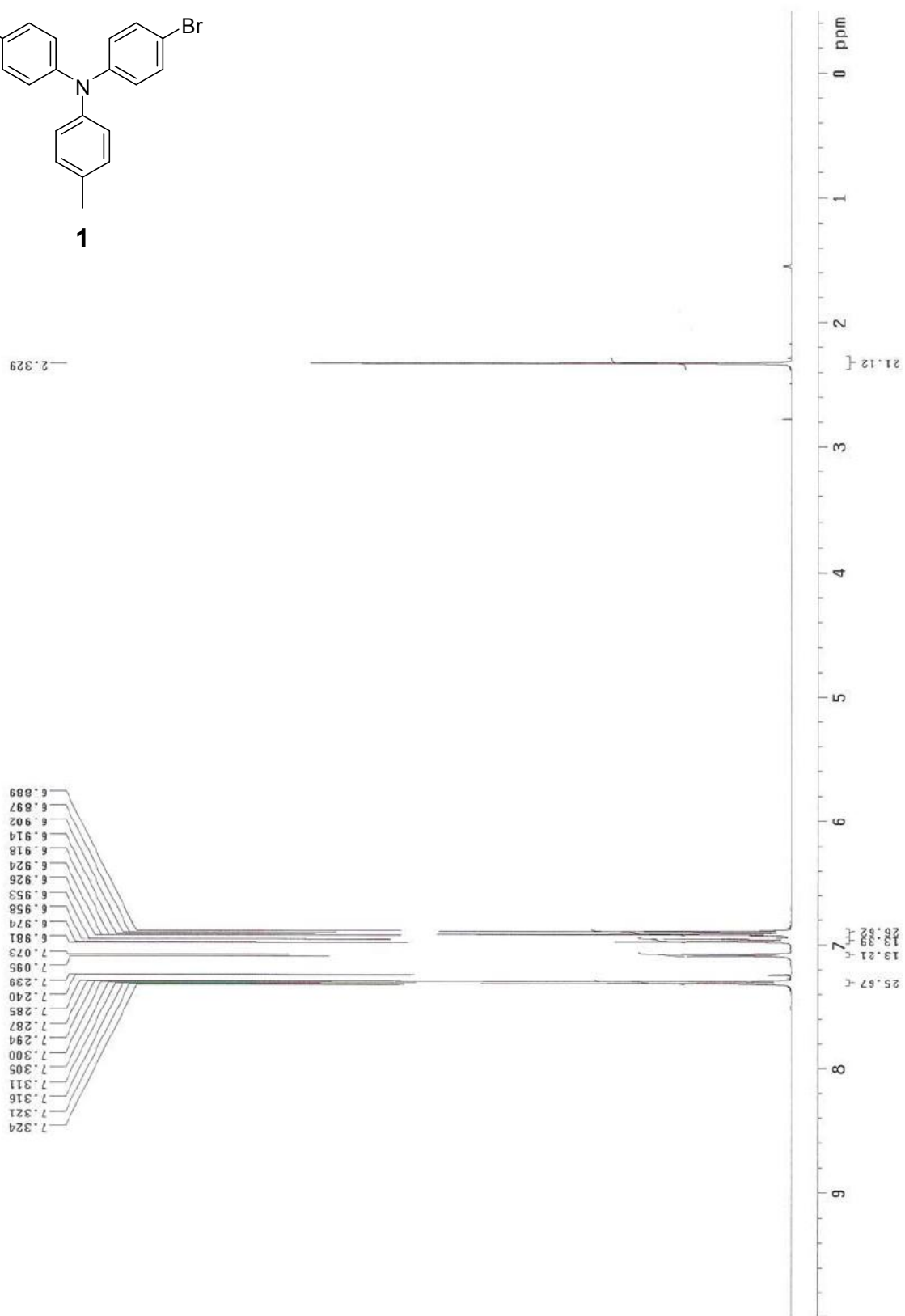
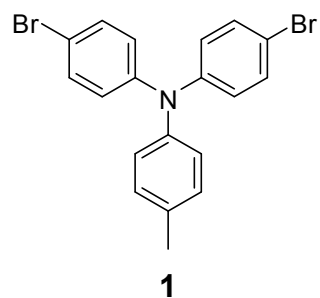
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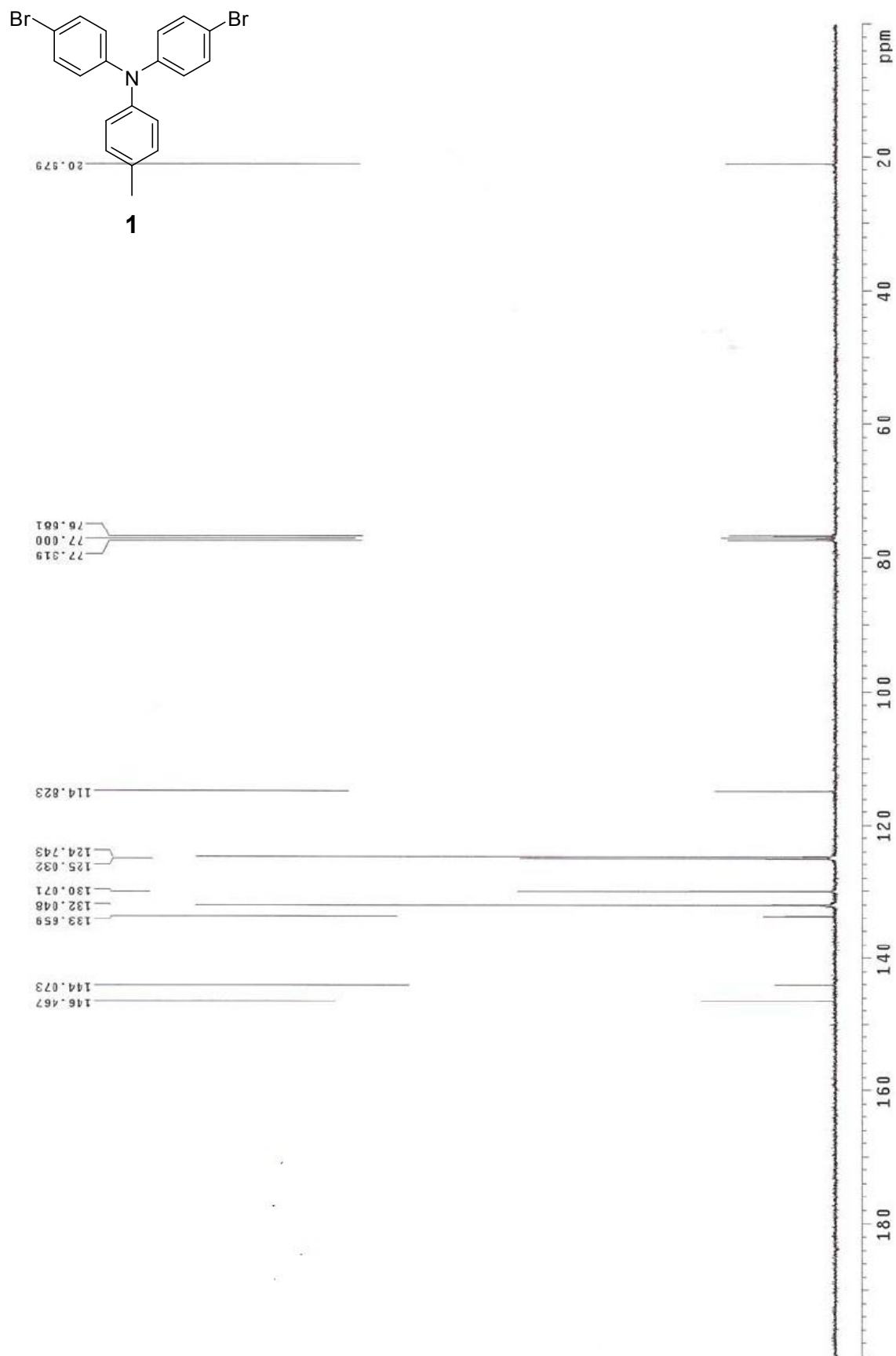
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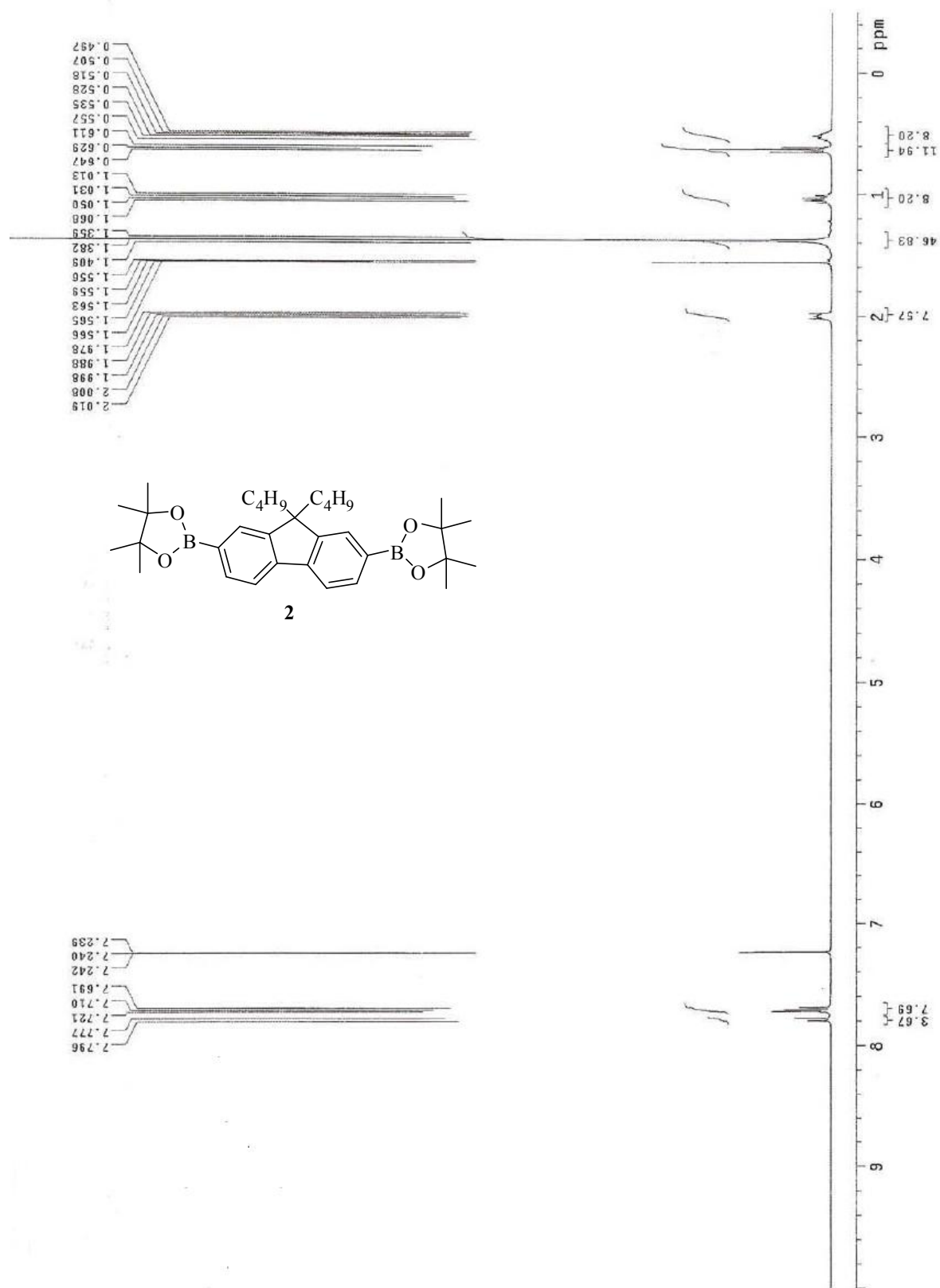
### Supporting Figure S1. $^1\text{H}$ NMR Spectrum of **1**



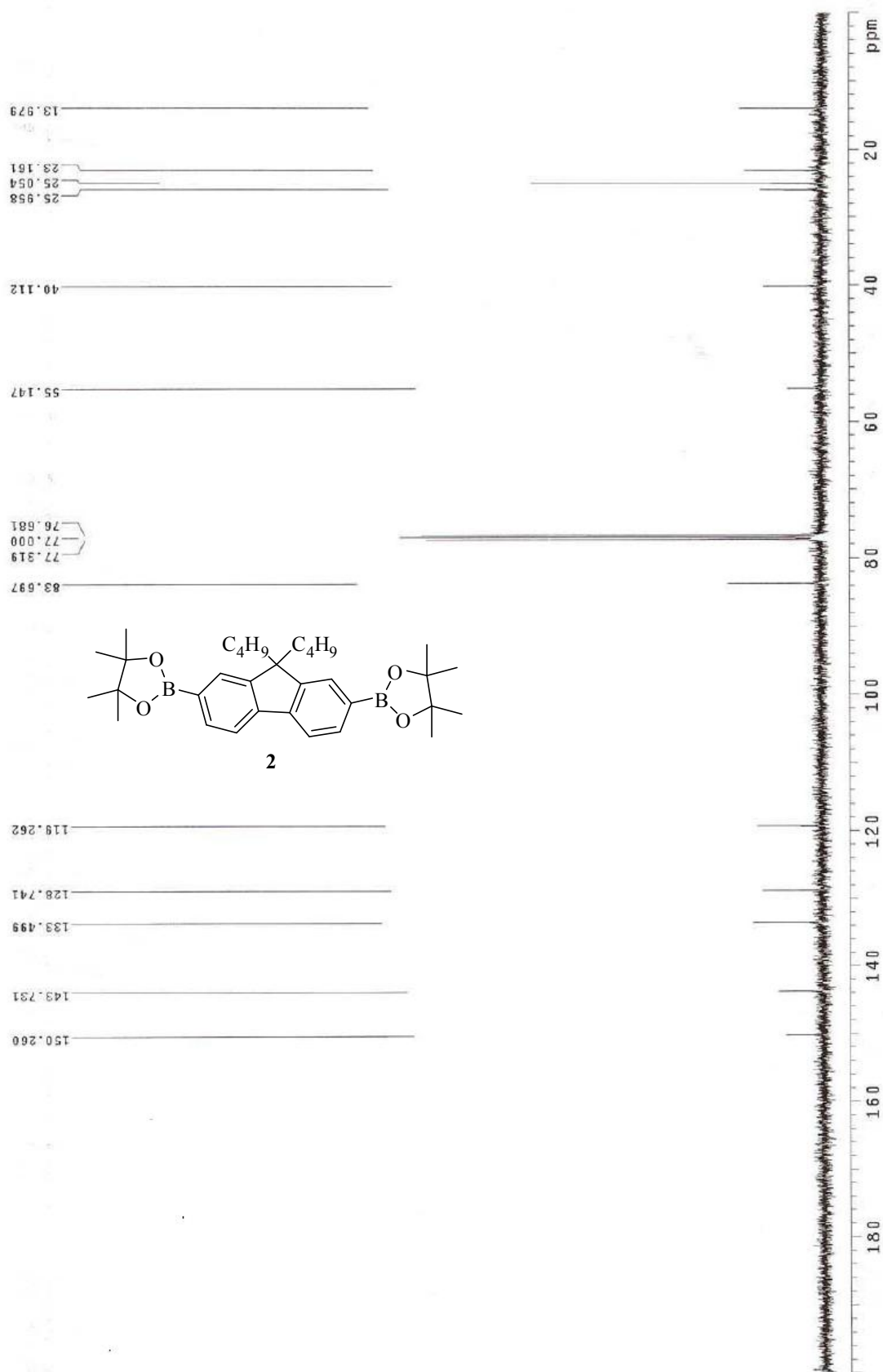
### Supporting Figure S2. $^{13}\text{C}$ NMR Spectrum of **1**



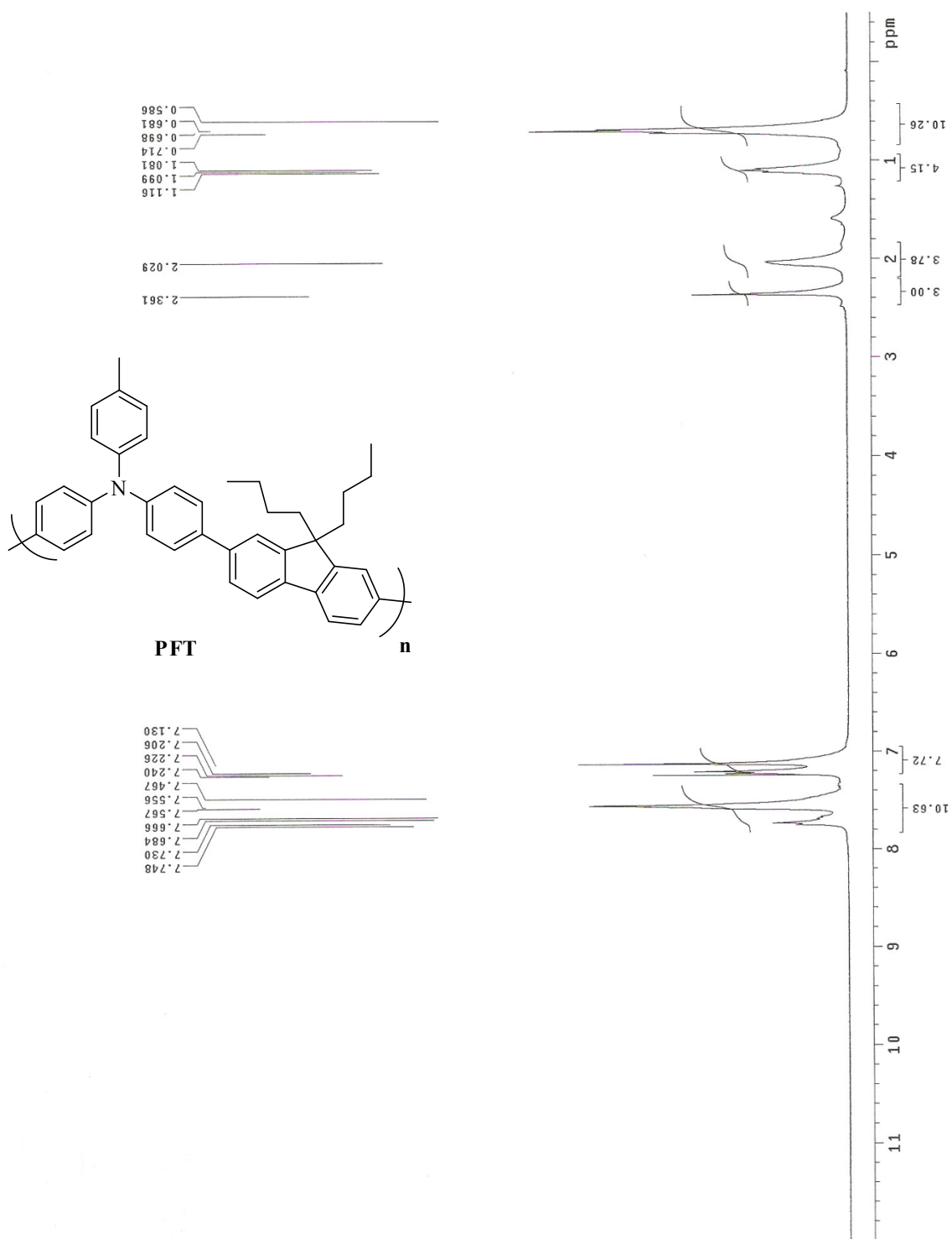
### Supporting Figure S3. $^1\text{H}$ NMR Spectrum of 2



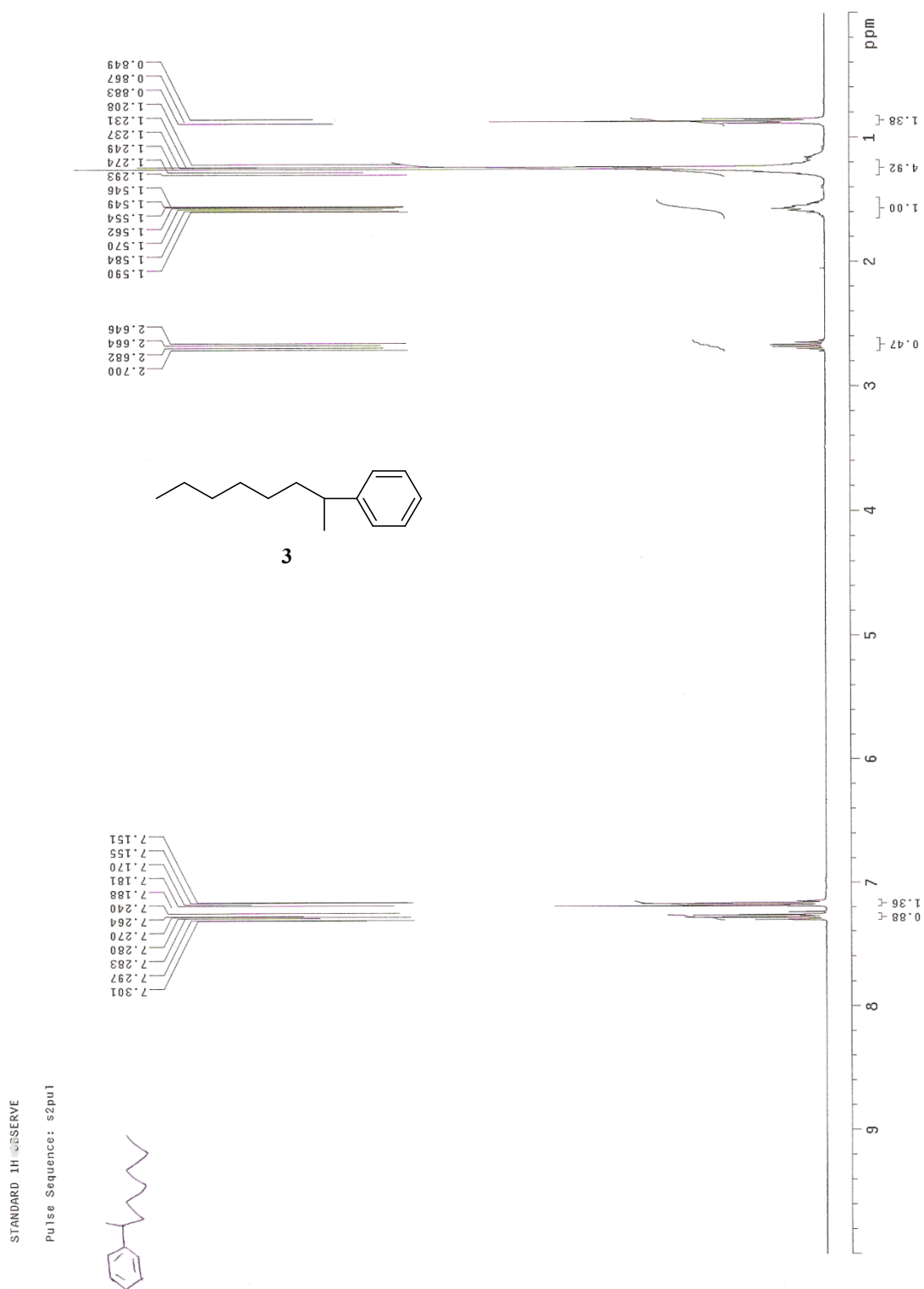
### Supporting Figure S4. $^{13}\text{C}$ NMR Spectrum of **2**



### Supporting Figure S5. $^1\text{H}$ NMR Spectrum of PFT

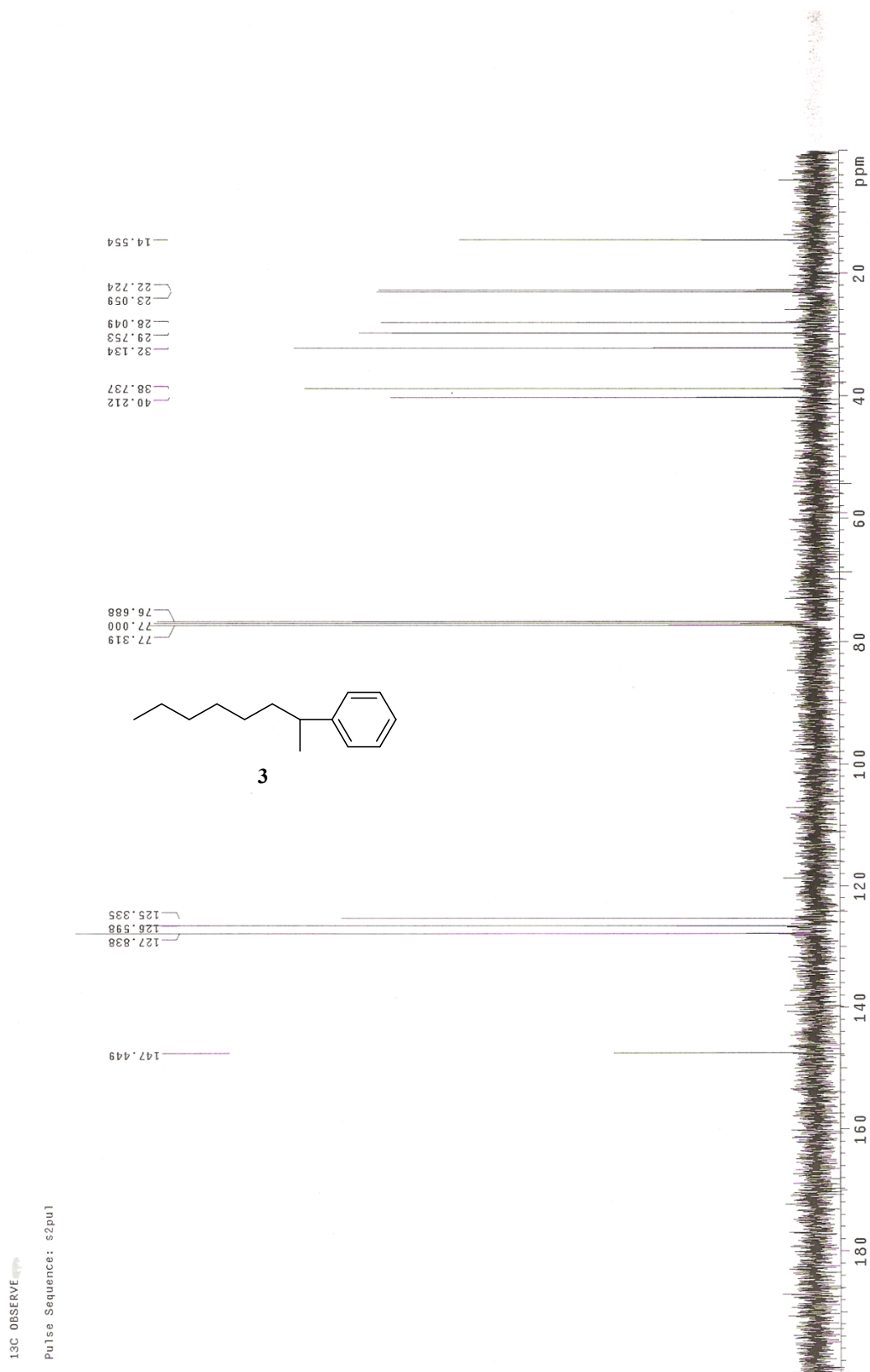


### Supporting Figure S6. <sup>1</sup>H NMR Spectrum of 3

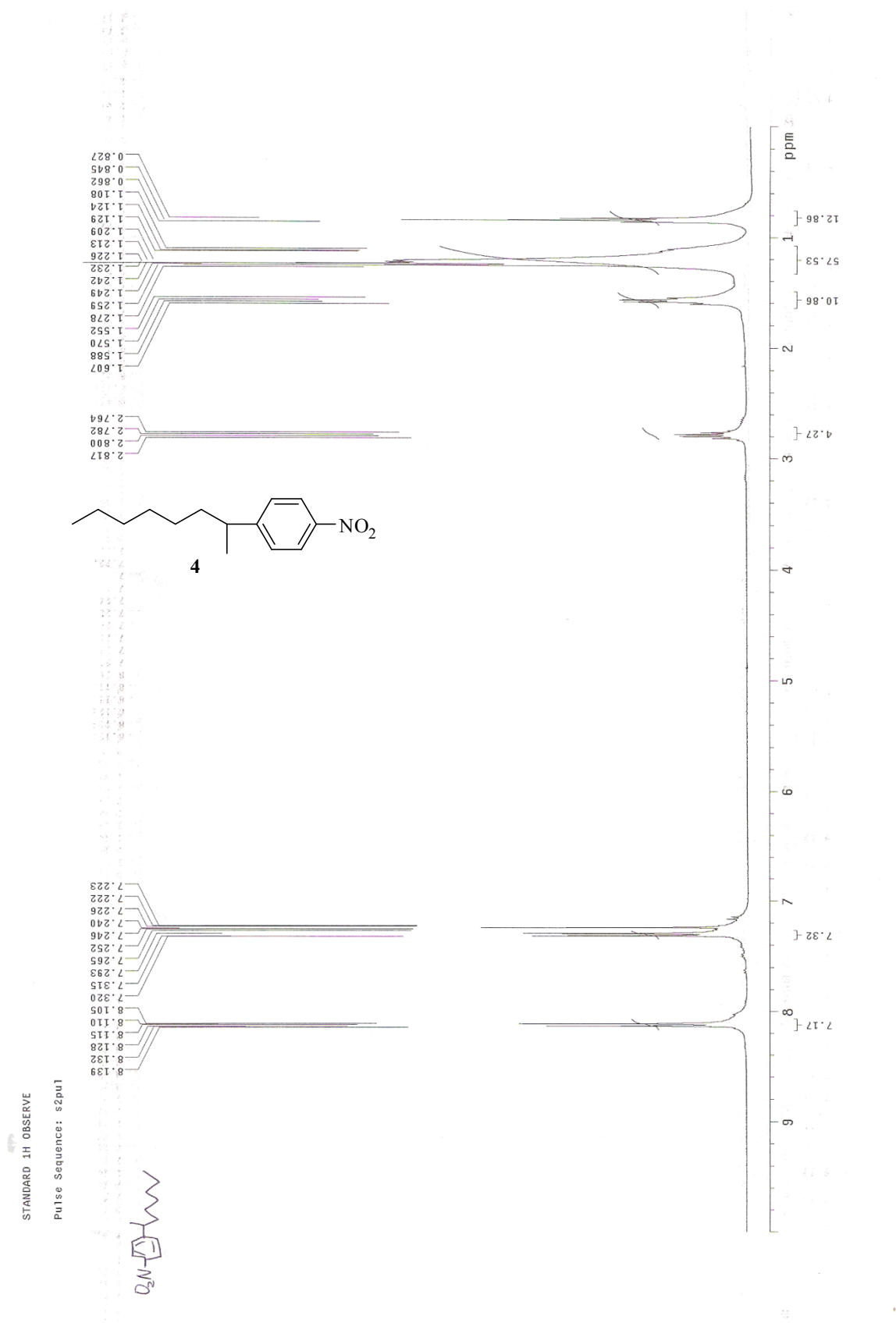




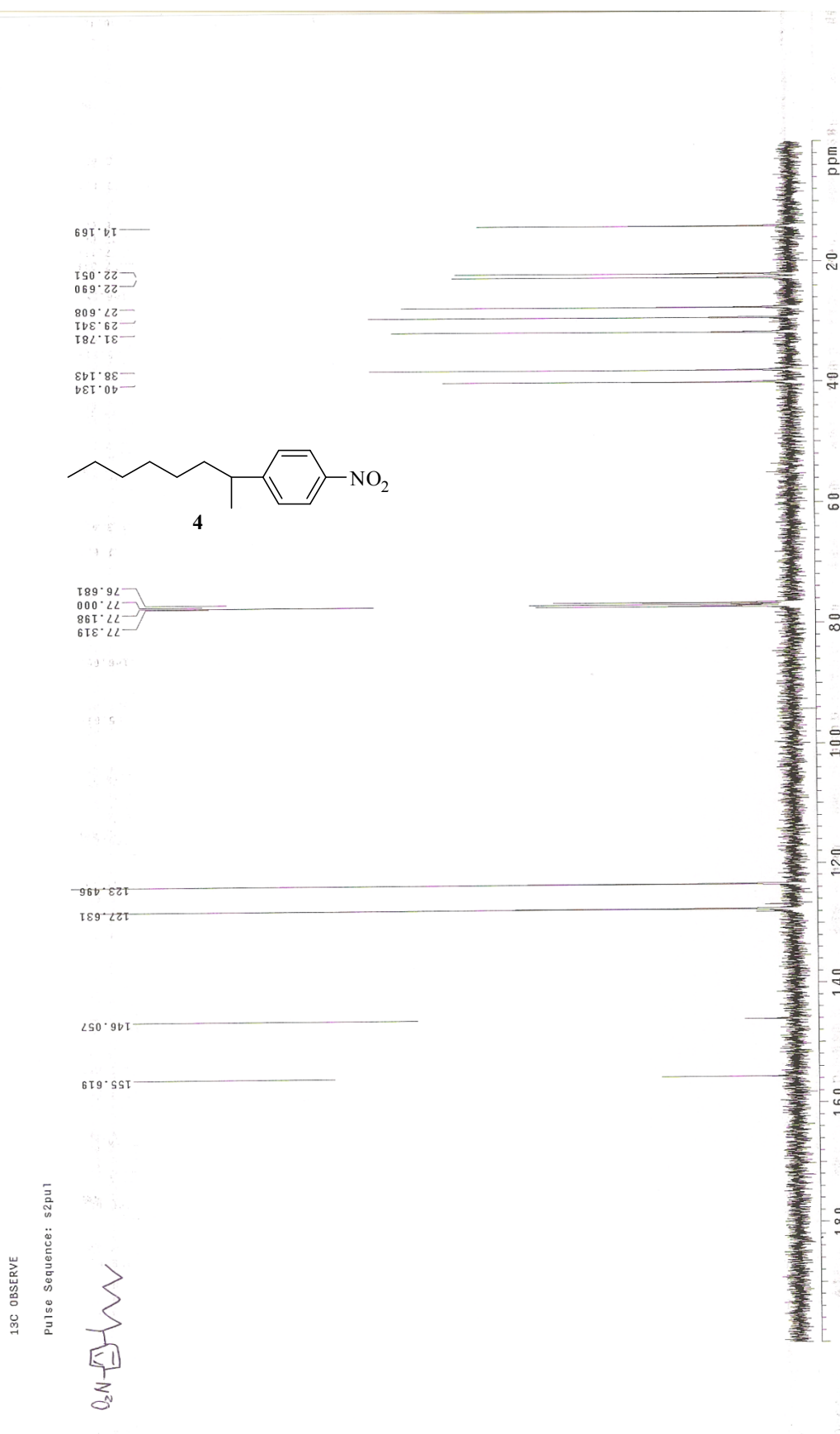
### Supporting Figure S7. <sup>13</sup>C NMR Spectrum of 3



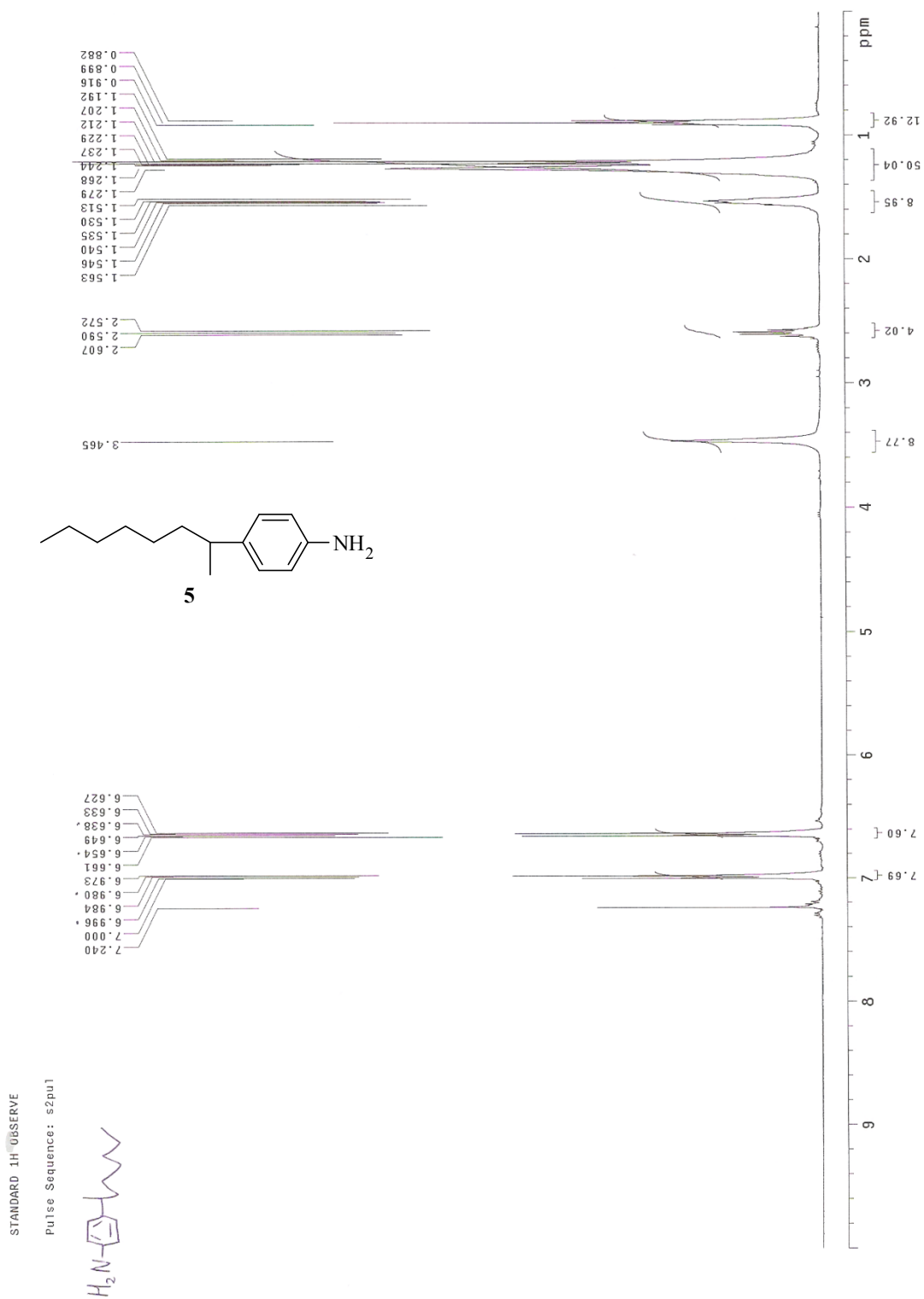
### Supporting Figure S8. <sup>1</sup>H NMR Spectrum of 4



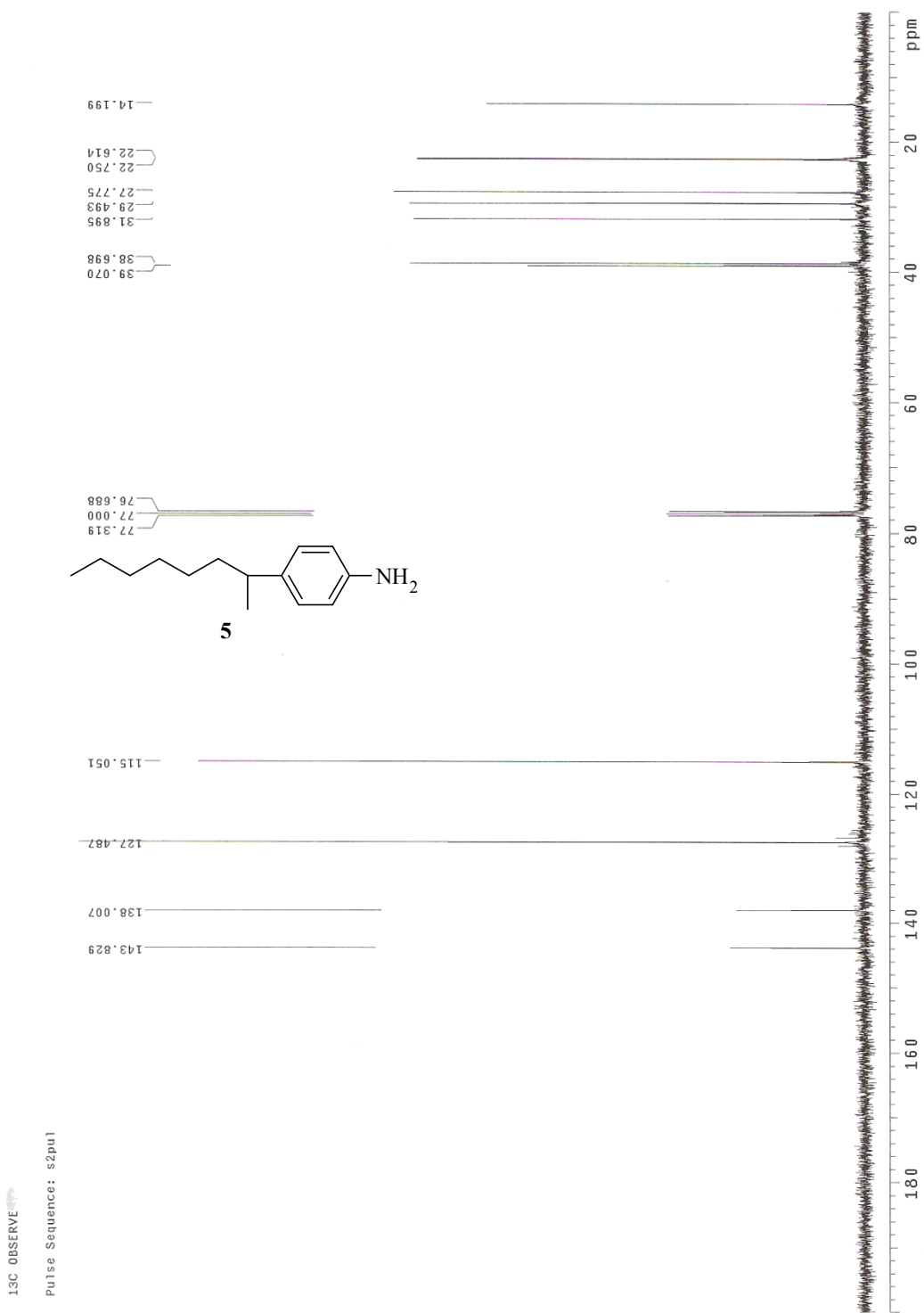
### Supporting Figure S9. $^{13}\text{C}$ NMR Spectrum of **4**



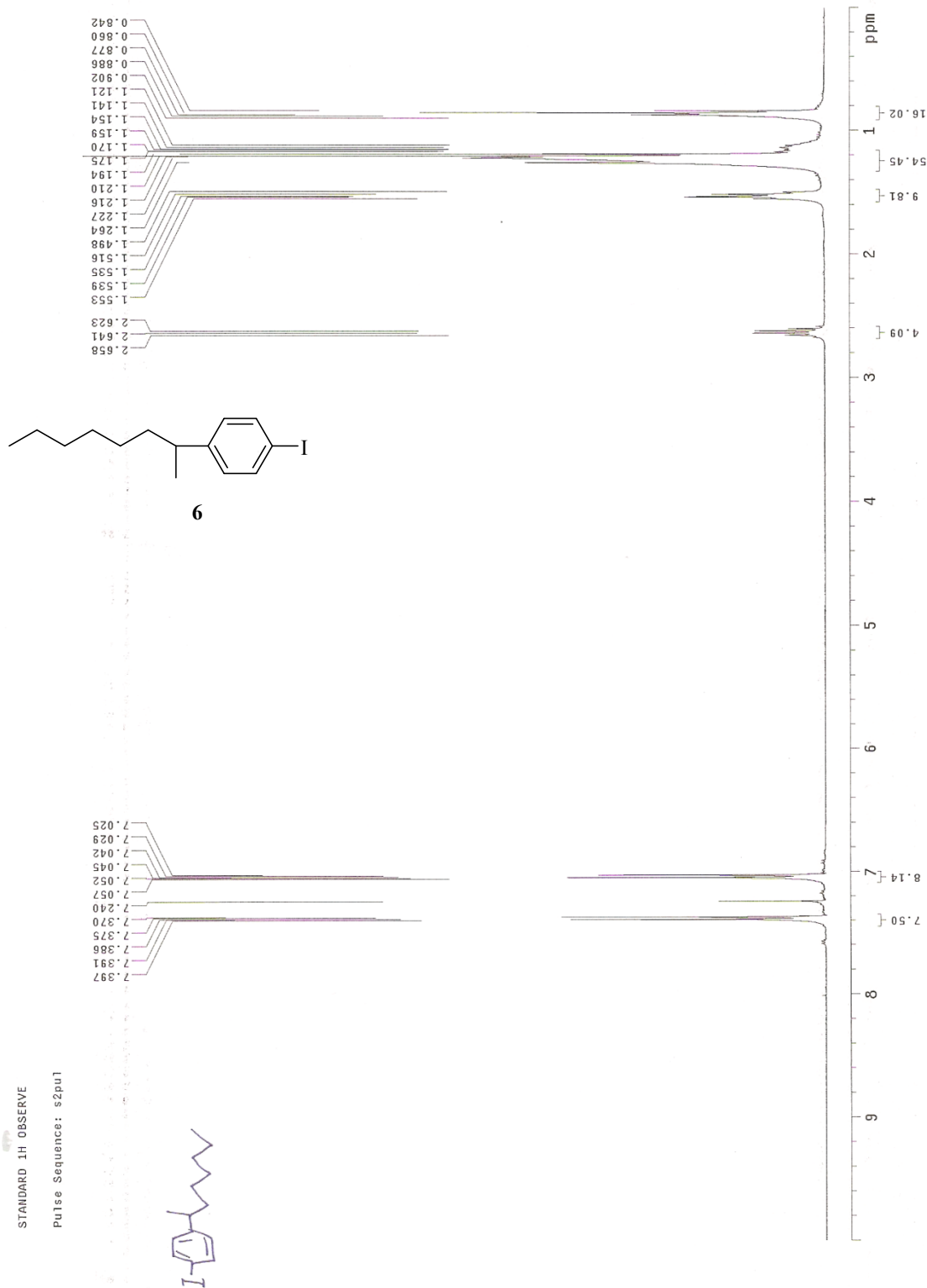
### Supporting Figure S10. <sup>1</sup>H NMR Spectrum of 5



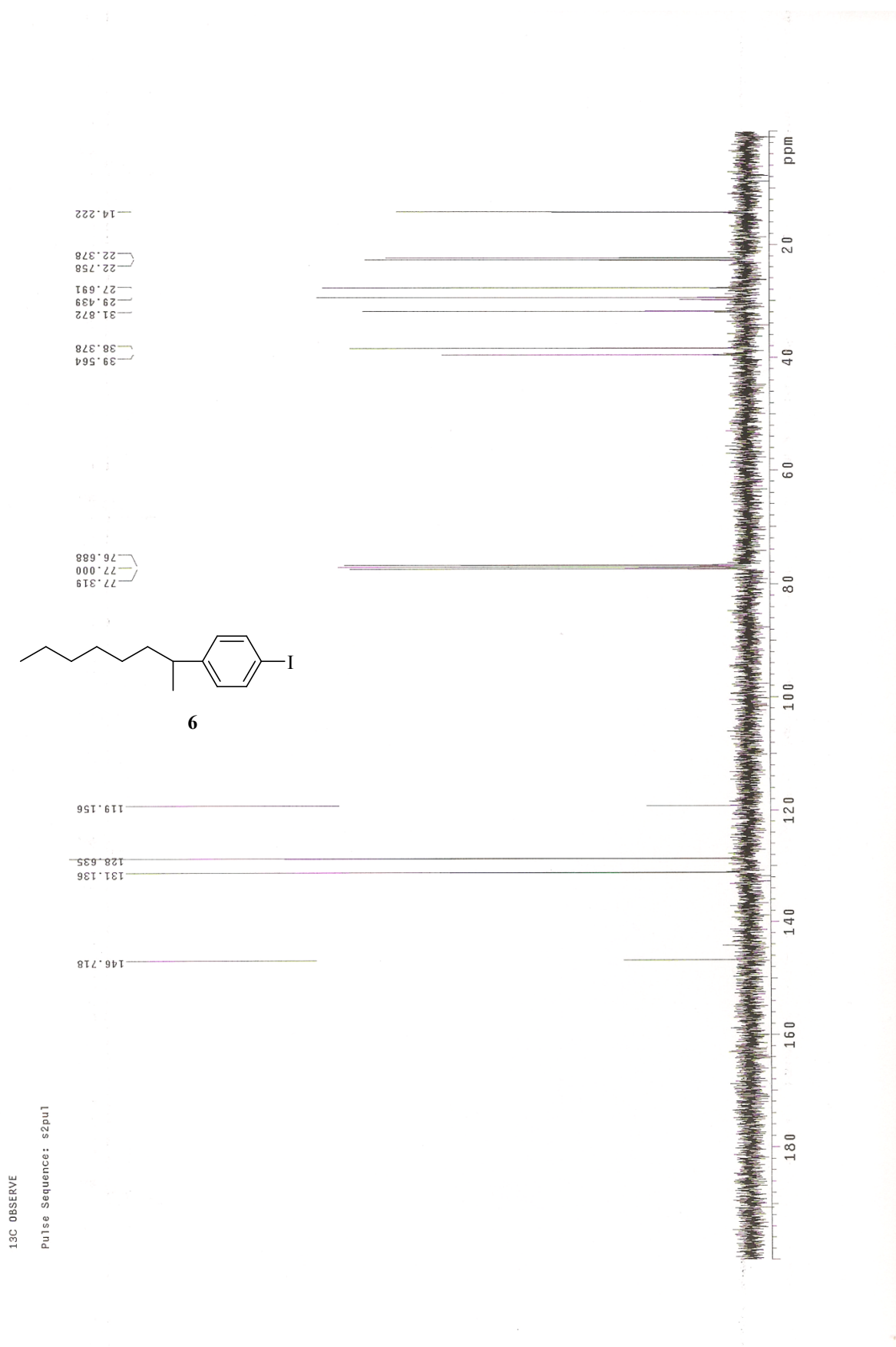
### Supporting Figure S11. $^{13}\text{C}$ NMR Spectrum of 5



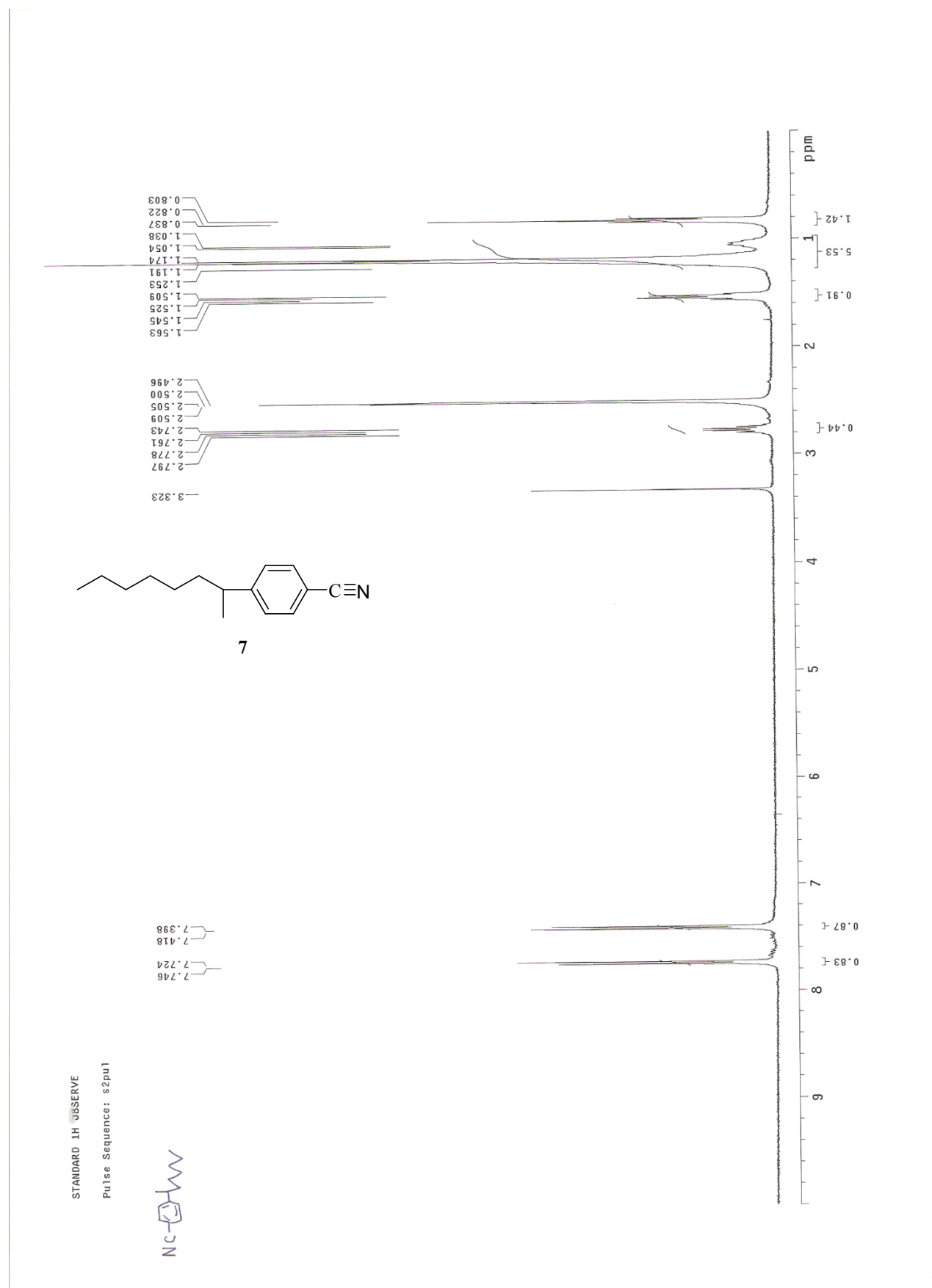
### Supporting Figure S12. <sup>1</sup>H NMR Spectrum of 6



### Supporting Figure S13. <sup>13</sup>C NMR Spectrum of 6

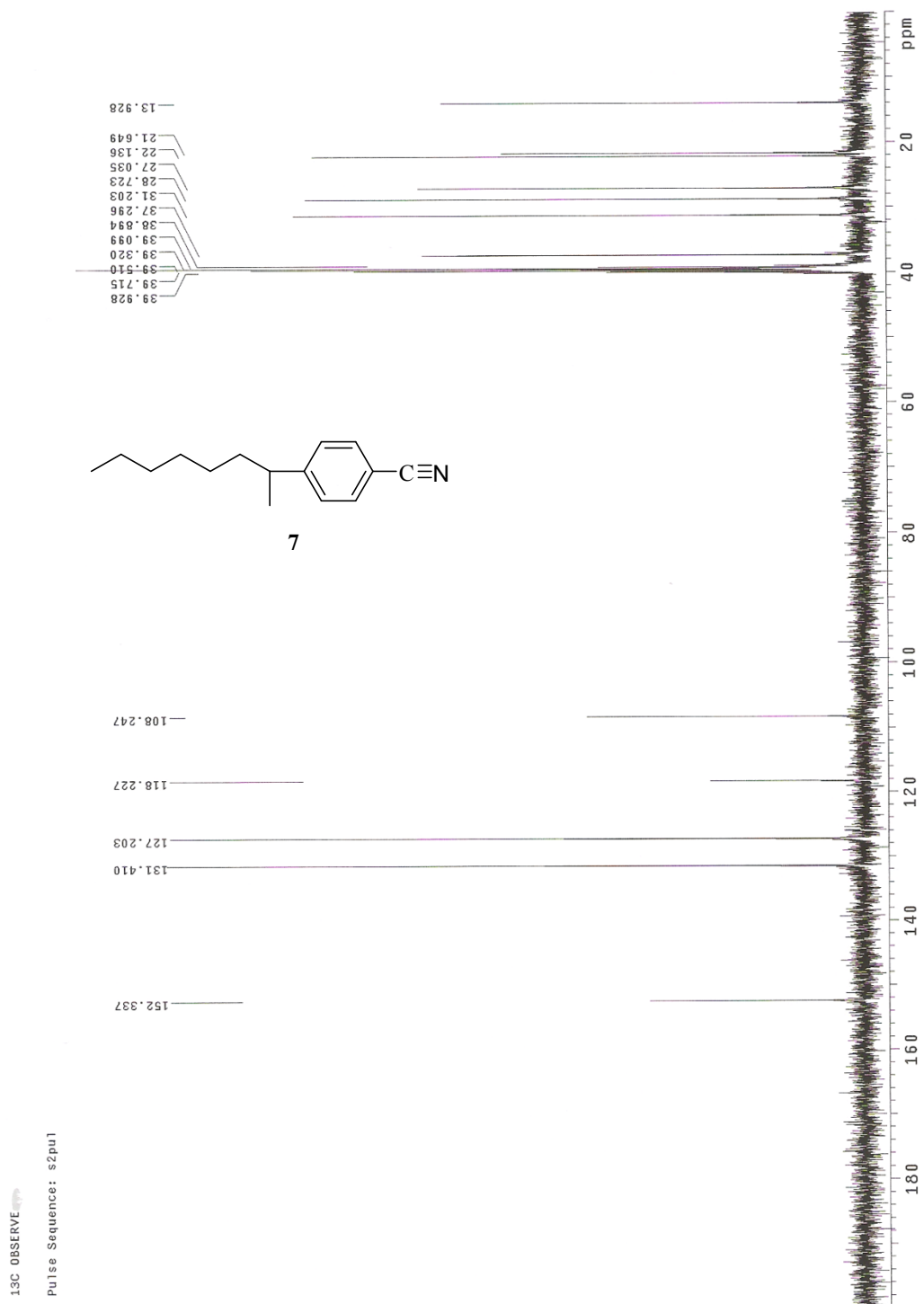


### Supporting Figure S14. <sup>1</sup>H NMR Spectrum of 7

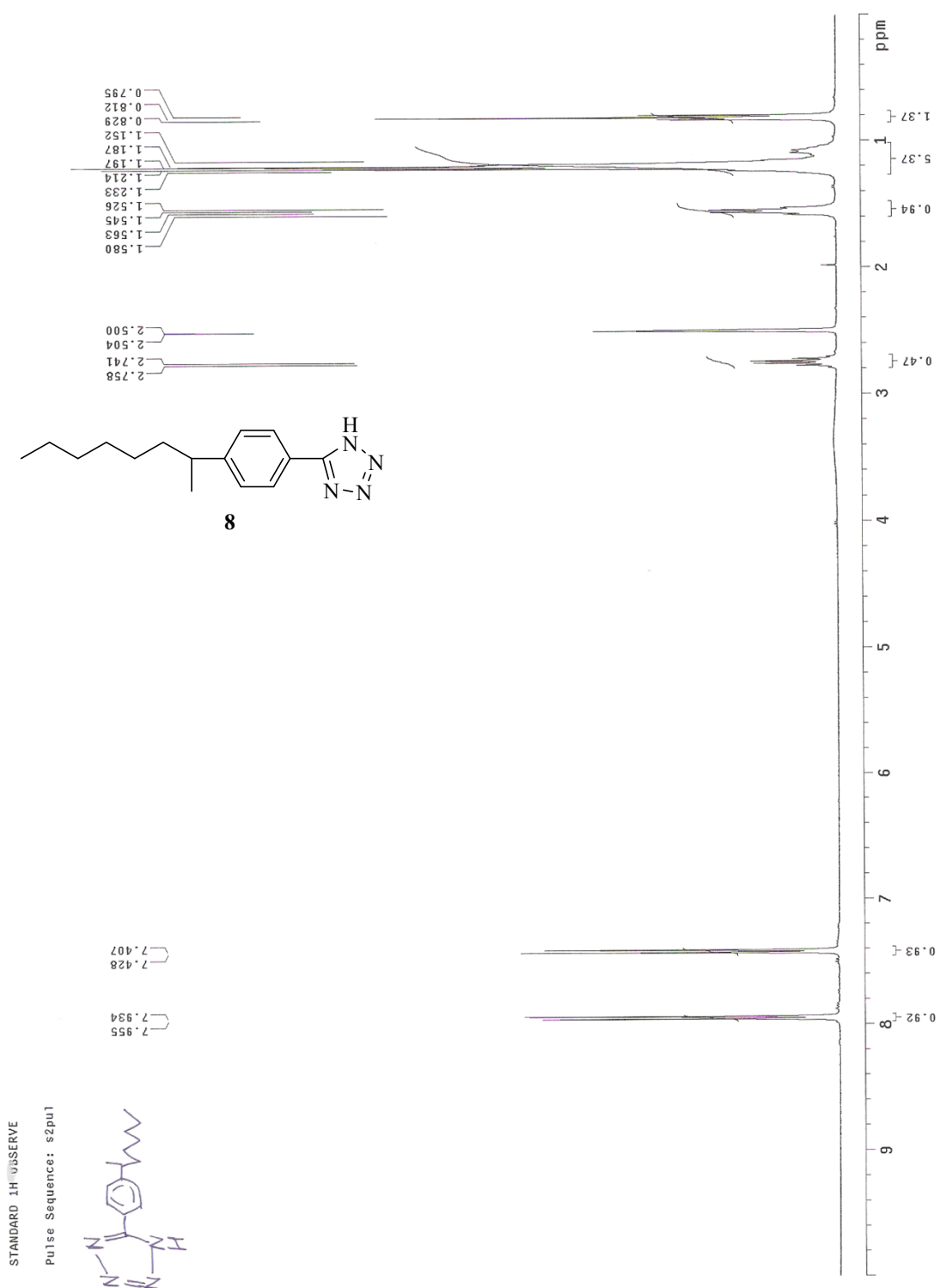




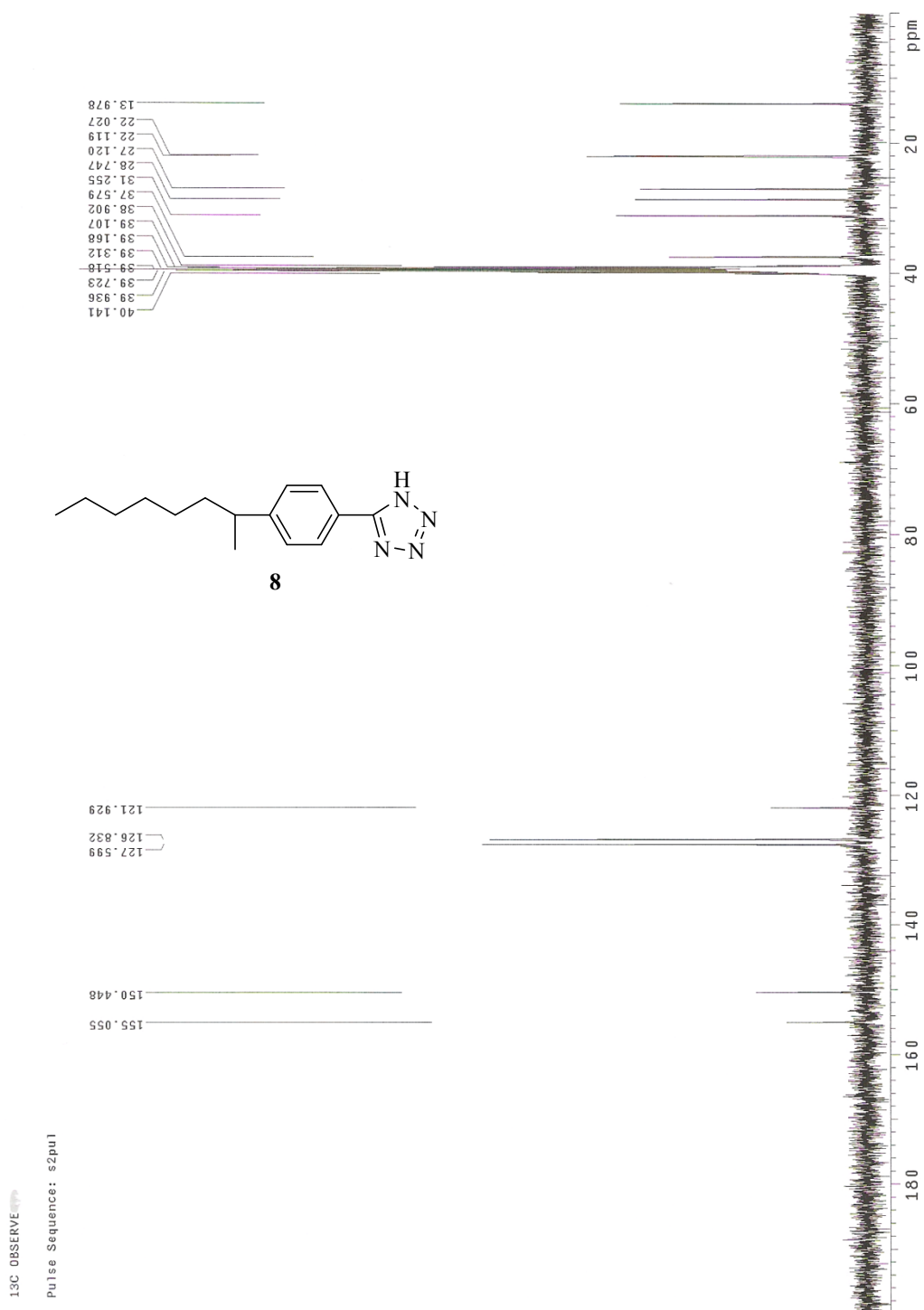
### Supporting Figure S15. <sup>13</sup>C NMR Spectrum of 7



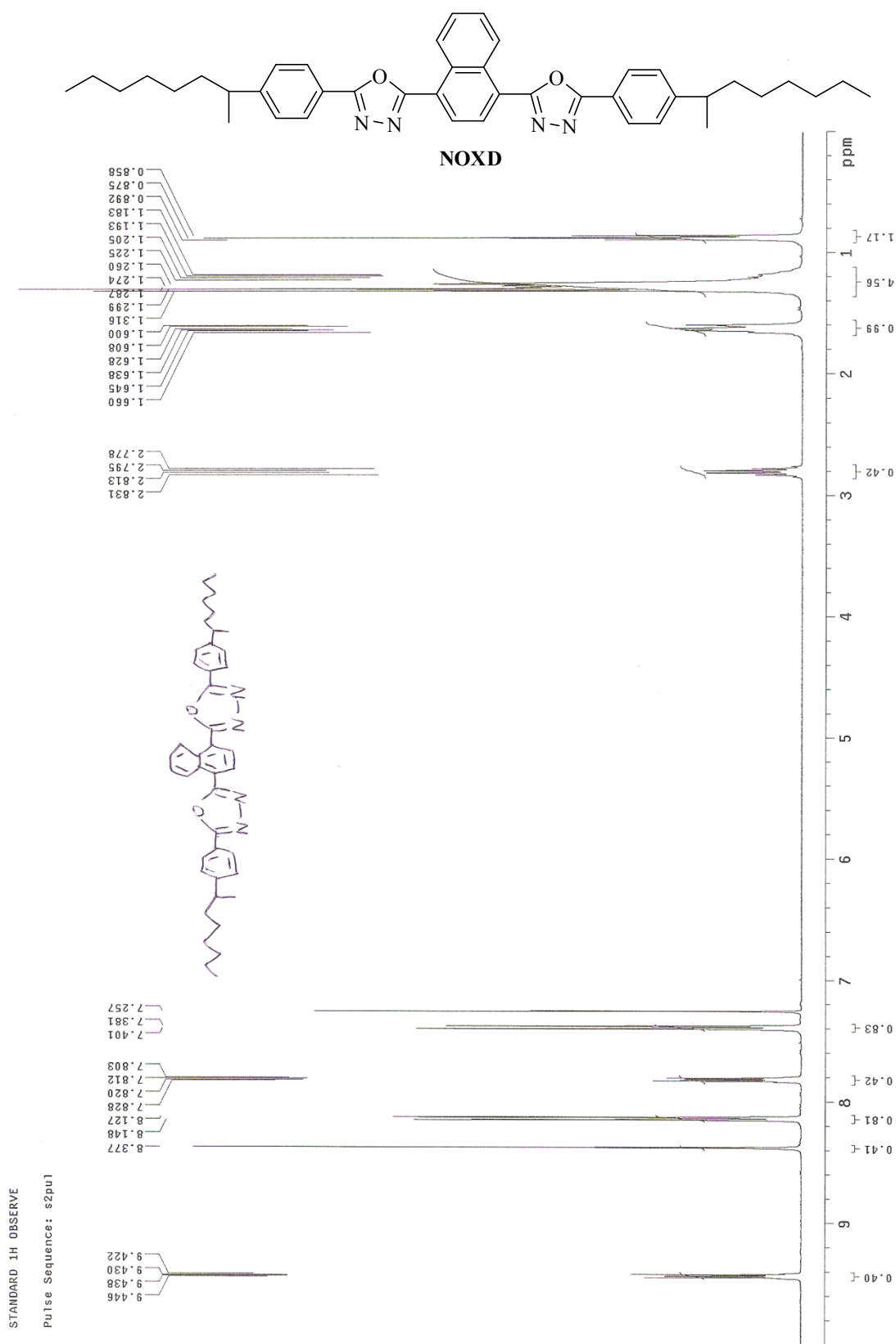
### Supporting Figure S16. <sup>1</sup>H NMR Spectrum of 8



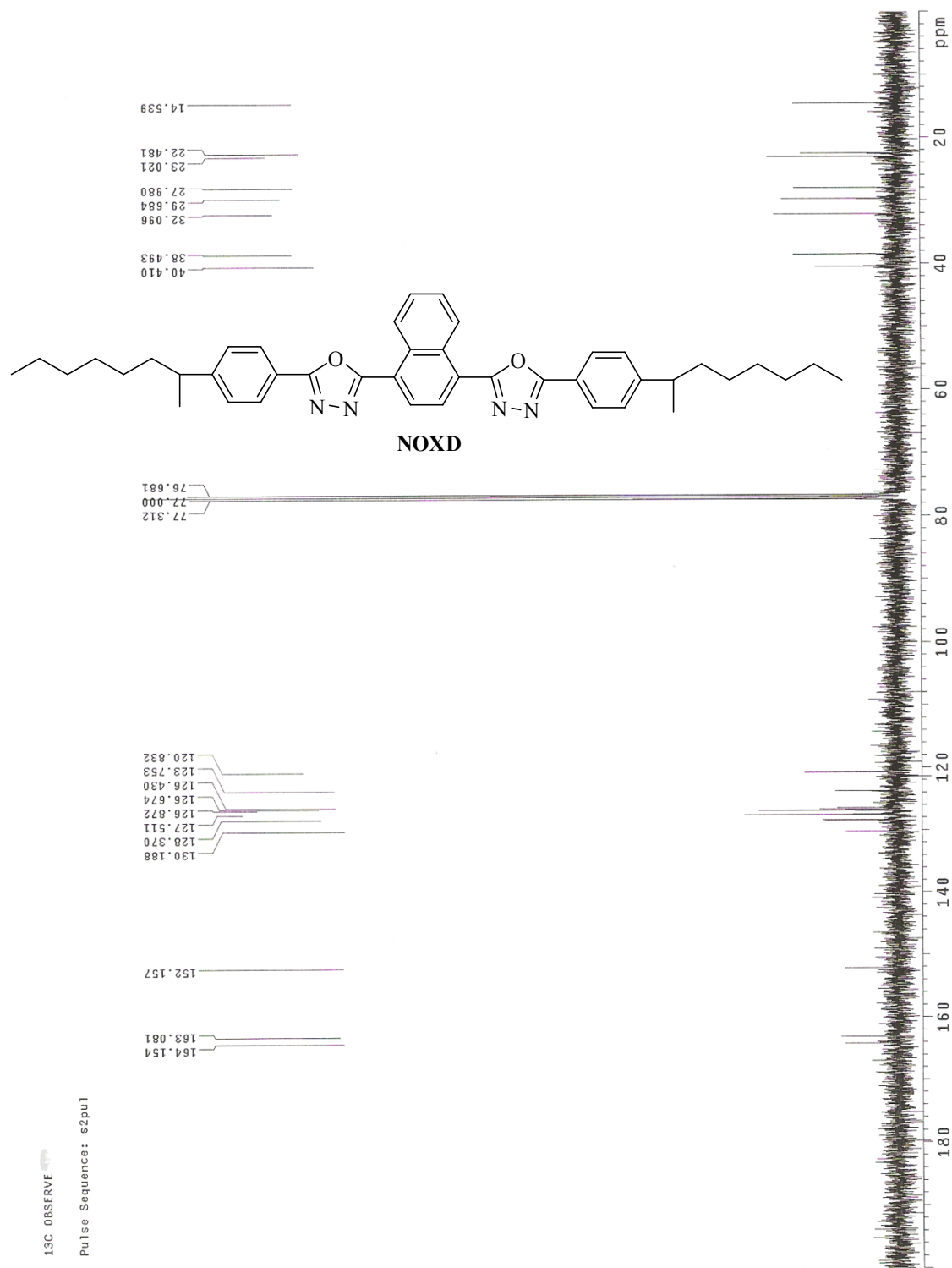
### Supporting Figure S17. $^{13}\text{C}$ NMR Spectrum of **8**



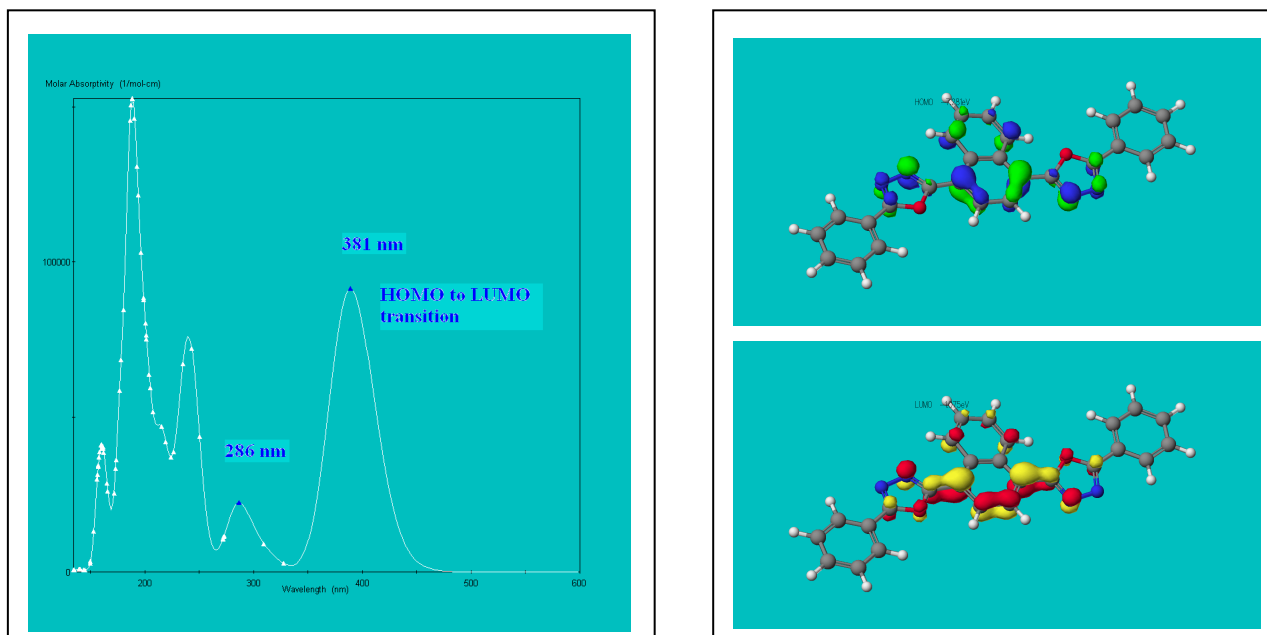
### Supporting Figure S18. <sup>1</sup>H NMR Spectrum of NOXD



### Supporting Figure S19. <sup>13</sup>C NMR Spectrum of NOXD

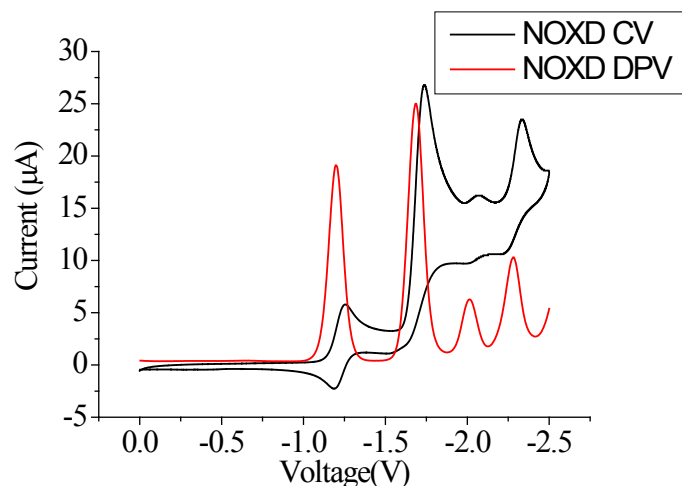


## Supporting Figure S20. Theoretical UV-Vis spectral analysis of NOXD

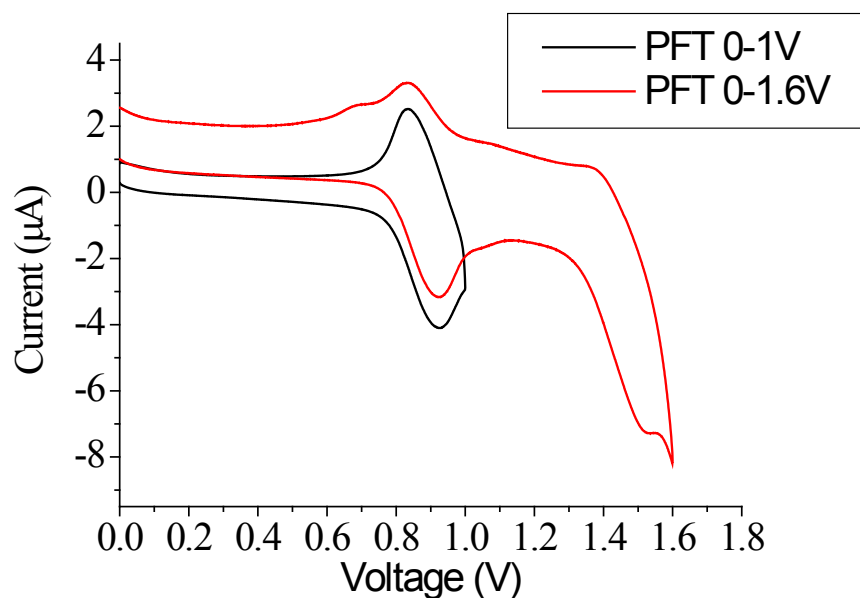


Theoretical UV-Vis spectral assignments are based on the molecular mechanics geometrical optimization and PM5/ZINDO calculations. The theoretical predictions are highly consistent with the experimental spectral data we obtained: The first transition at 360 nm is arising from HOMO/LUMO transition. The second band contains contributions from other 5 transitions.

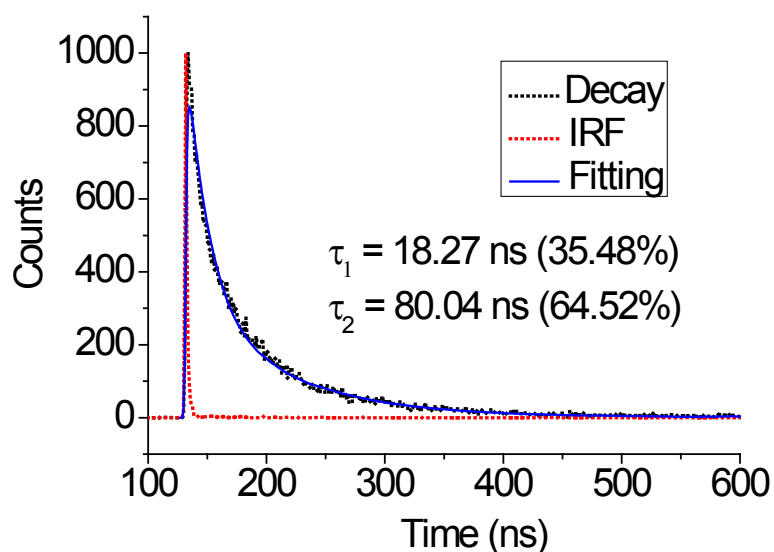
**Supporting Figure S21.** Cyclic voltammograms of NOXD in TBAP/DMF (0.1M) under N<sub>2</sub> with a scan rate of 100 mV/s. A Glassy Carbon electrode was used as the working electrode versus a AgCl/Ag electrode as the reference electrode, and a Pt wire, as the counter electrode. The voltage was calibrated by using ferrocene as the internal standard.



**Supporting Figure S22.** Cyclic voltammograms of PFT in TBAP/CH<sub>2</sub>Cl<sub>2</sub> (0.1M) with a scan rate of 100 mV/s.



**Supporting Figure S23.** Time-resolved photoluminescence measurements of the **NOXD:PFT**.



**EXPERIMENTAL PROCEDURES**

The fluorescence lifetime was detected using a time-correlated single photon counting (TCSPC) unit of an Edinburgh FLSP920 instrument. Samples were excited at 337 nm from a nitrogen pulsed flashlamp with 1 ns fwhm pulse duration transmitted through a Czerny-Turner design monochromator. Photomultiplier tube was employed as the detector. The instrument response function was profiled using a scatter solution and subsequently deconvoluted from the emission data to yield an undisturbed decay. Nonlinear least-squares fitting of the decay curves were performed with the Levenburg-Marquardt algorithm and implemented with the Edinburgh Instruments F900 software. The quality of the fits was judged by the  $\chi^2$  values and distribution of the residuals. The fitting curve based on two single-exponential decay components ( $\tau_1 \sim 18.27 \text{ ns}$ ,  $\tau_2 \sim 80.04 \text{ ns}$ ).