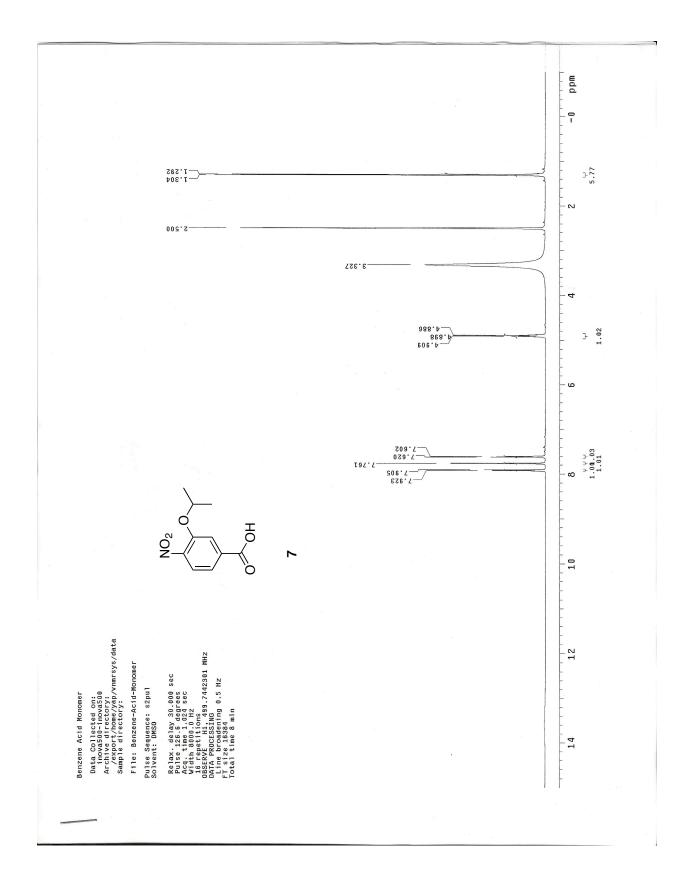
Relaxation of the rigid backbone of an oligoamide-foldamerbased α-helix mimetic: identification of potent Bcl-x_L inhibitors

Jeremy L. Yap,^a Xiaobo Cao,^b Kenno Vanommeslaeghe,^a Kwan-Young Jung,^a Chander S. Peddaboina,^b Paul T. Wilder,^c Anjan Nan,^a Alexander D. MacKerell, Jr.,^{a,d} W. Roy Smythe^b and Steven Fletcher^{*,a,d}

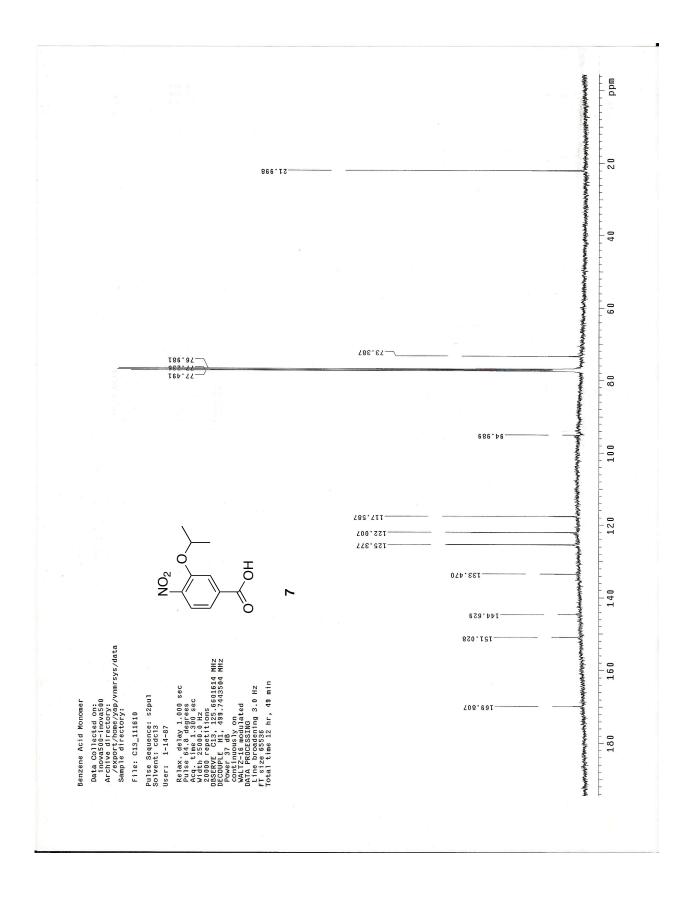
^aDepartment of Pharmaceutical Sciences, University of Maryland School of Pharmacy, 20 N. Pine St., Baltimore, MD 21201, USA; ^bScott & White Memorial Hospital, 702 SW HK Dodgen loop Room 219, Temple, TX 76504, USA; ^cDepartment of Biochemistry and Molecular Biology, University of Maryland School of Medicine, 108 N. Greene Street, Baltimore, MD 21201, USA; ^dUniversity of Maryland Marlene and Stewart Greenebaum Cancer Center, 22 S. Greene St., Baltimore, MD 21201, USA

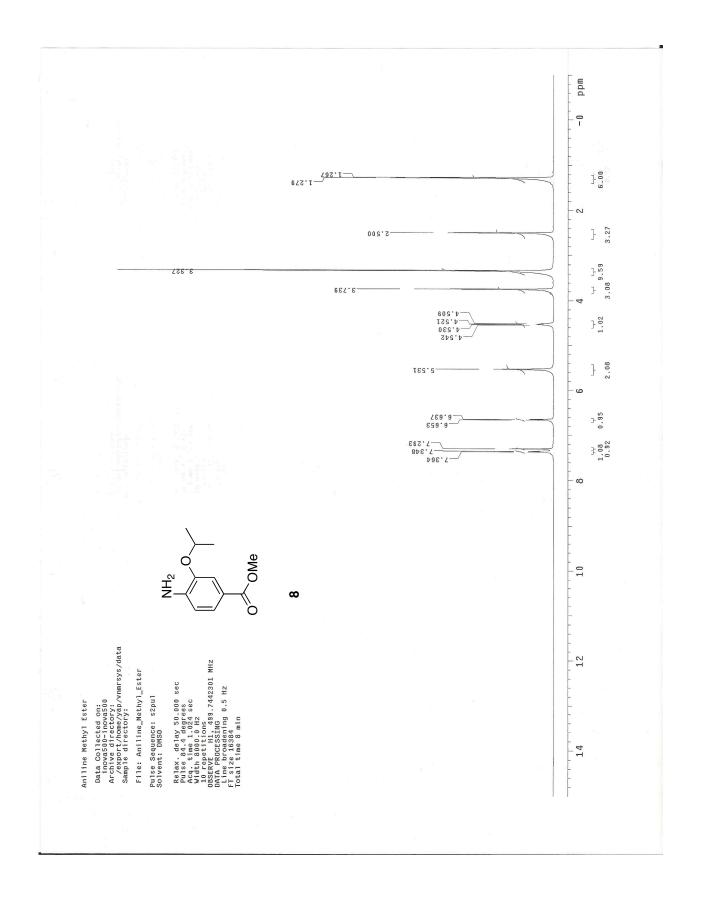
E-mail: sfletche@rx.umaryland.edu

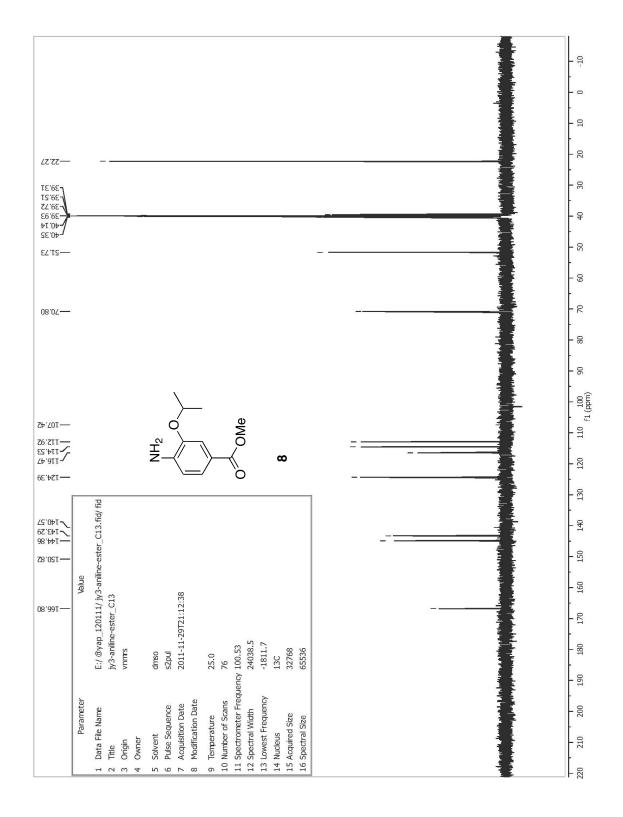
Supplementary Information: ¹H and ¹³C NMR Spectra for Intermediates 7 – 20f and Target Molecules 1 – 6

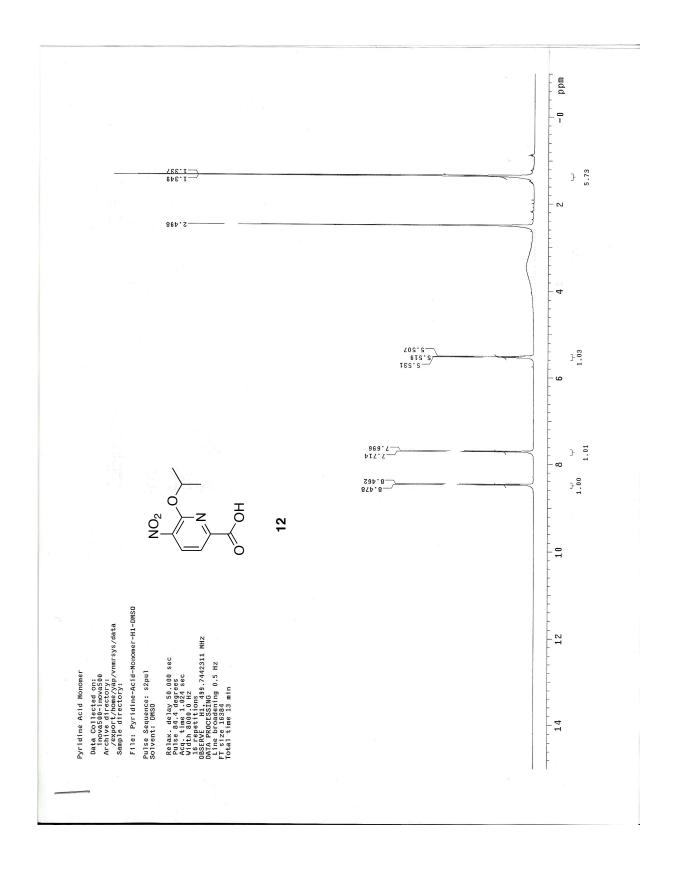


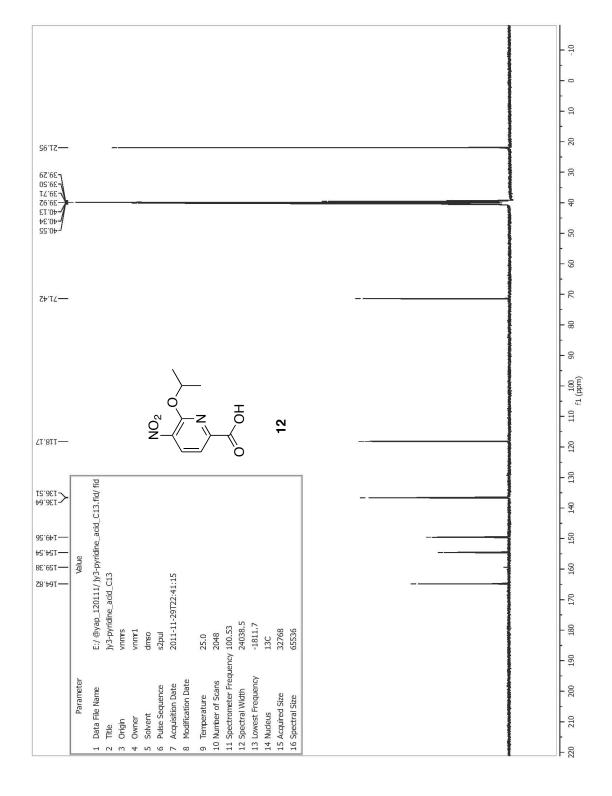
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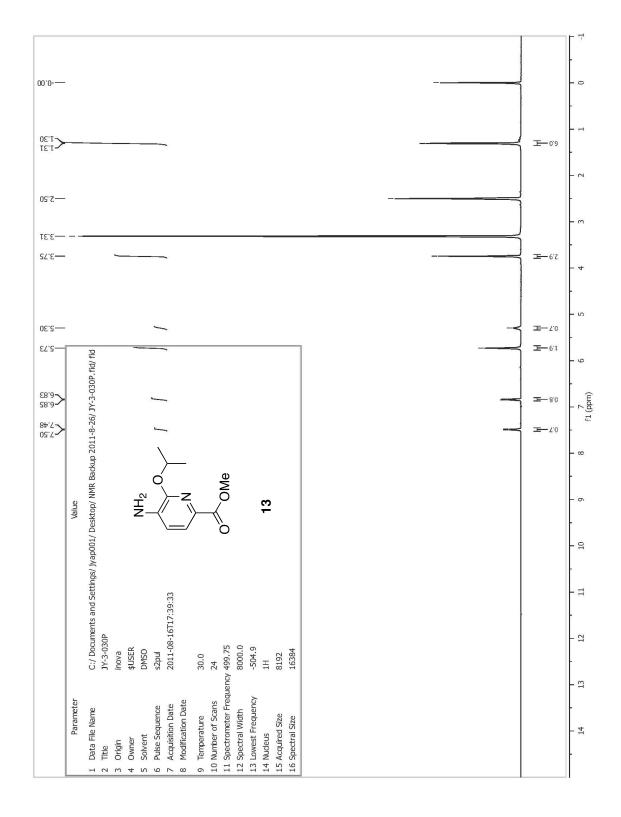


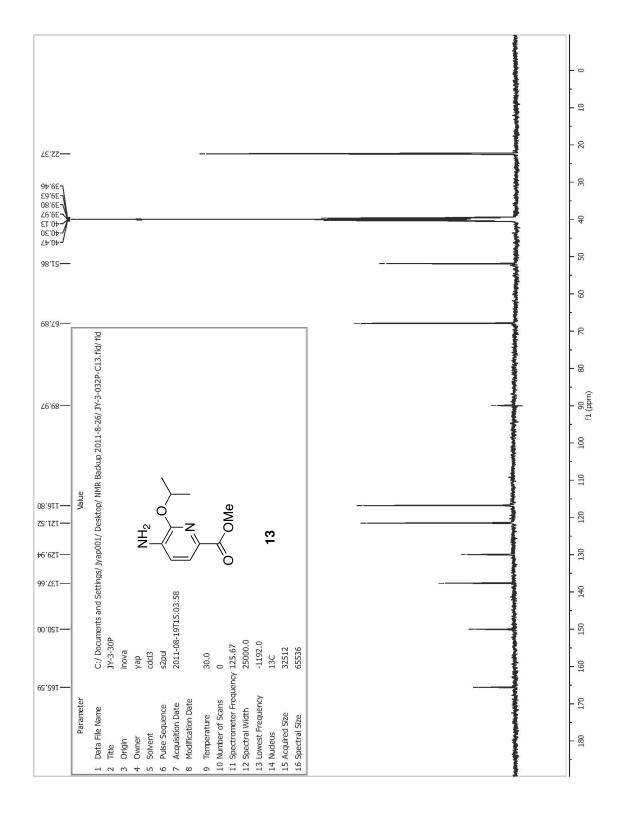


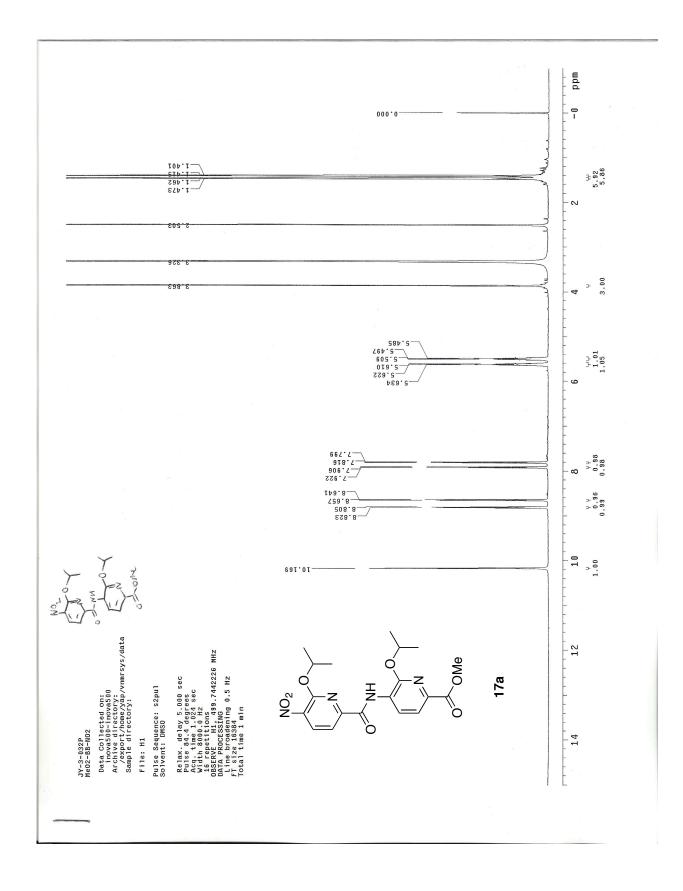


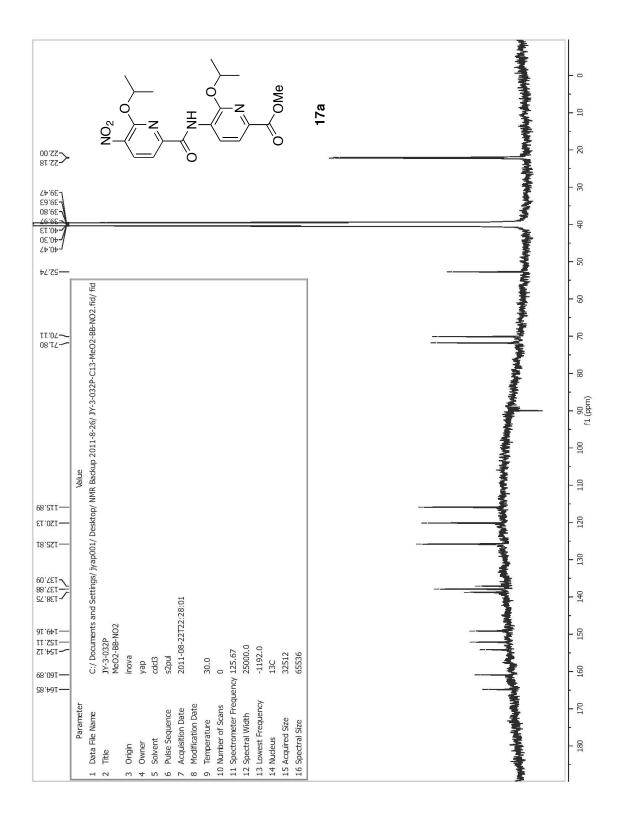


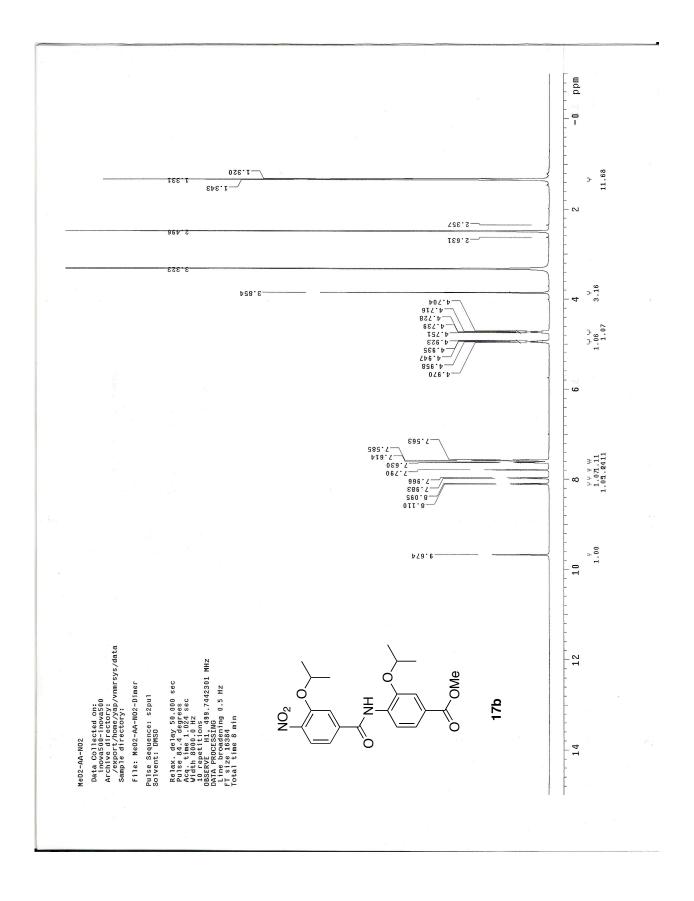


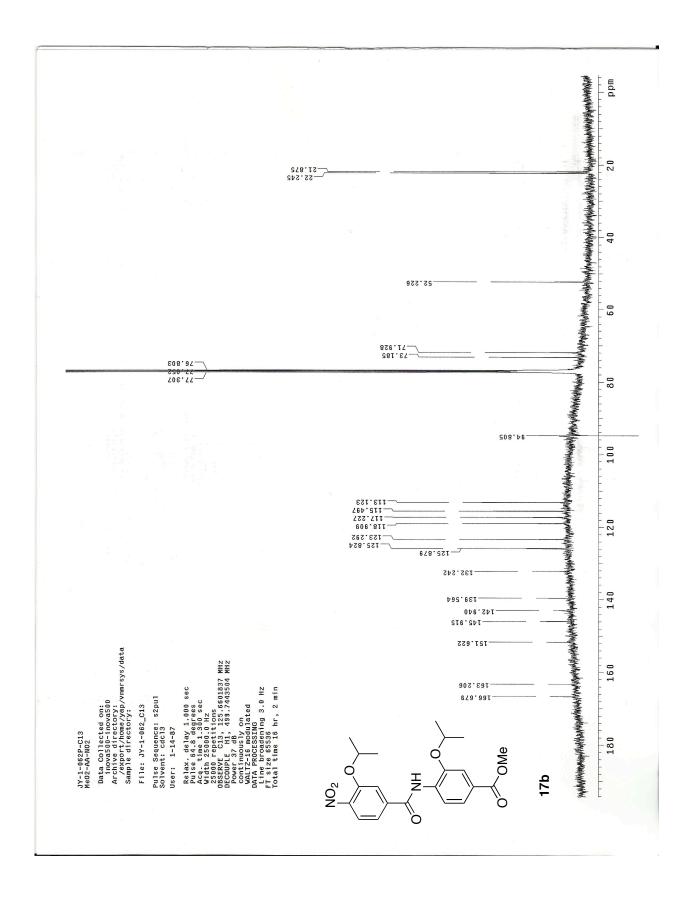


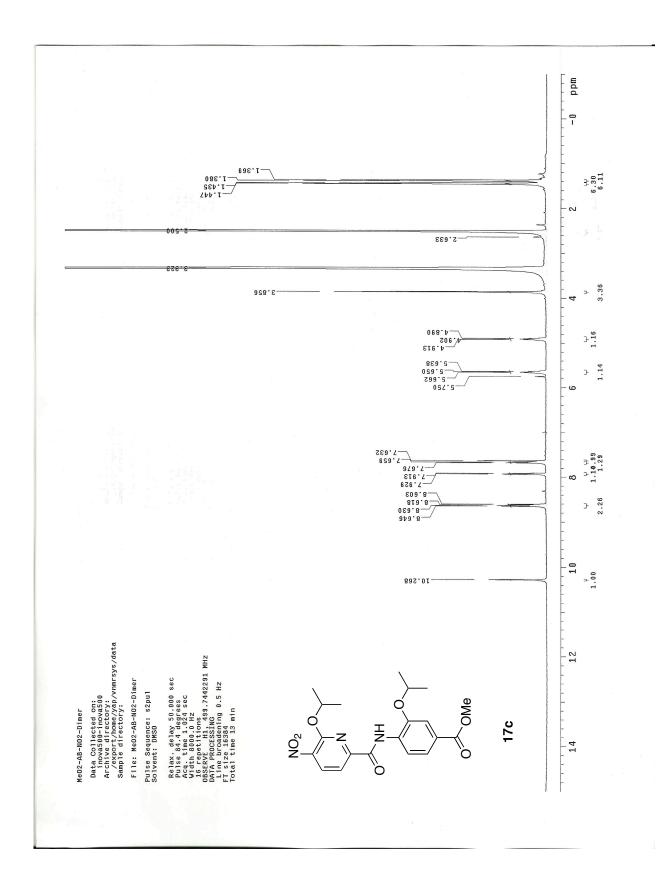


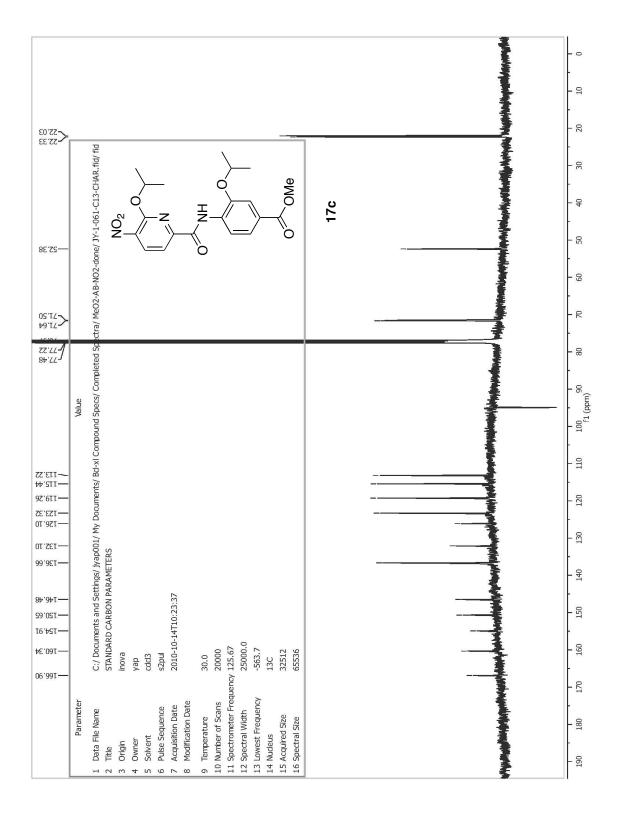


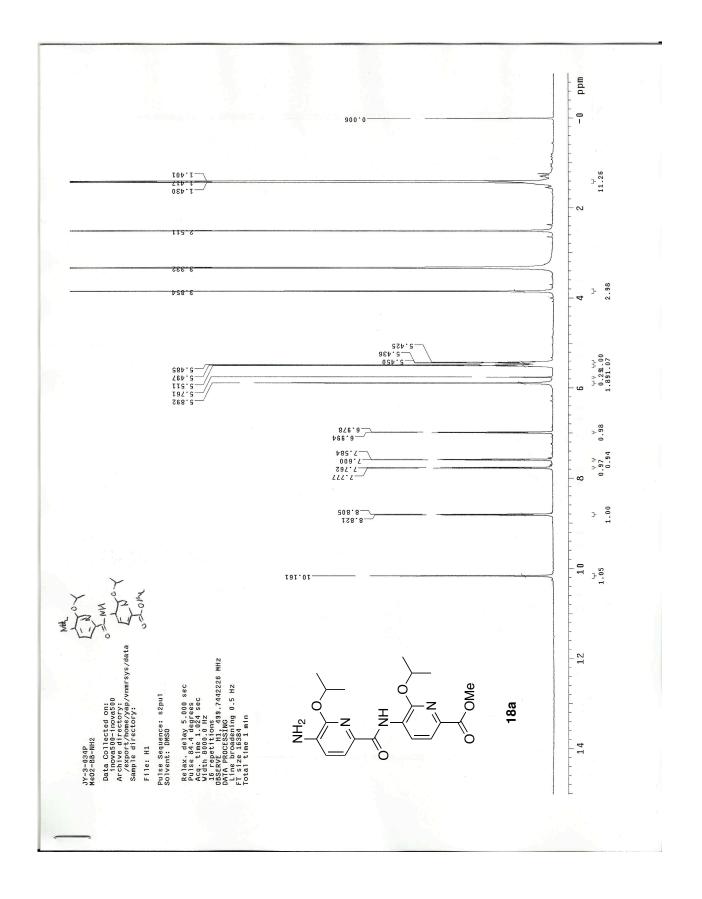


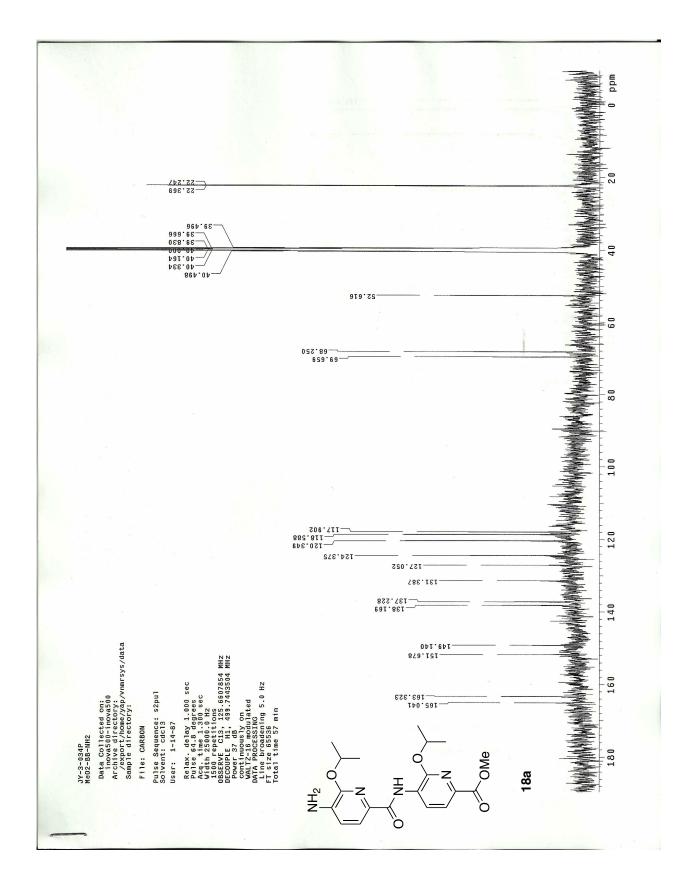


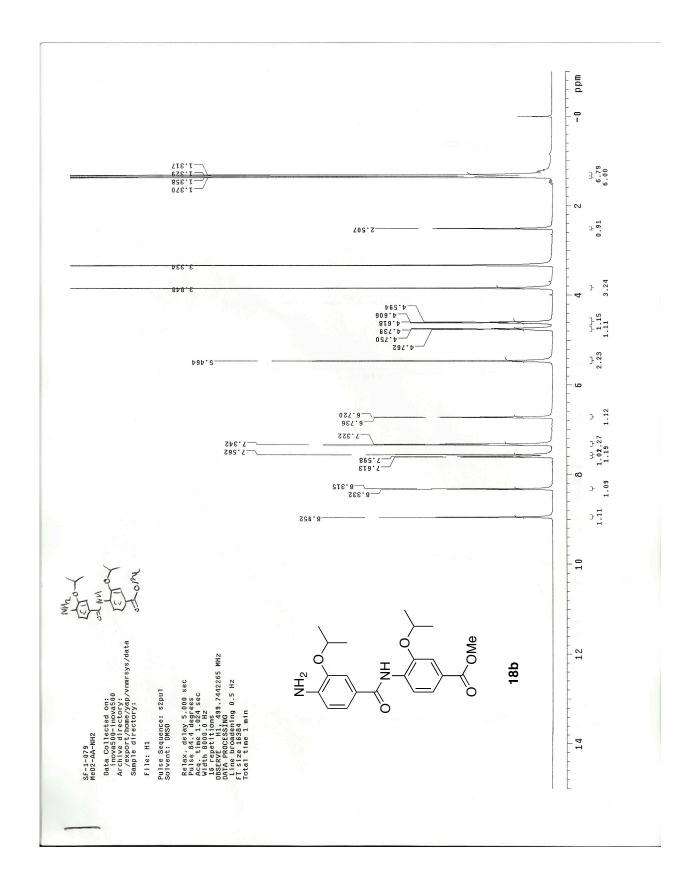


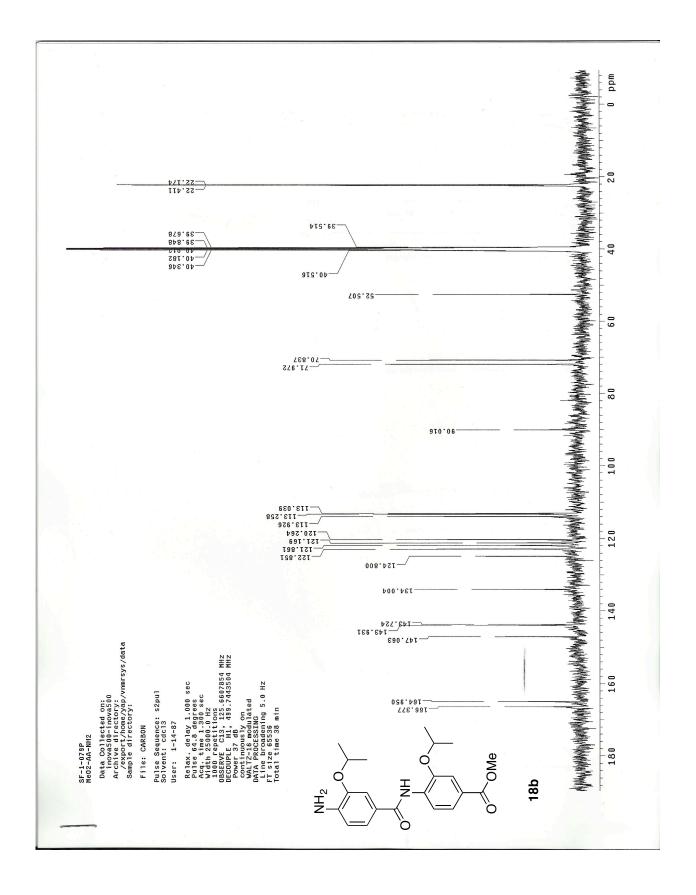


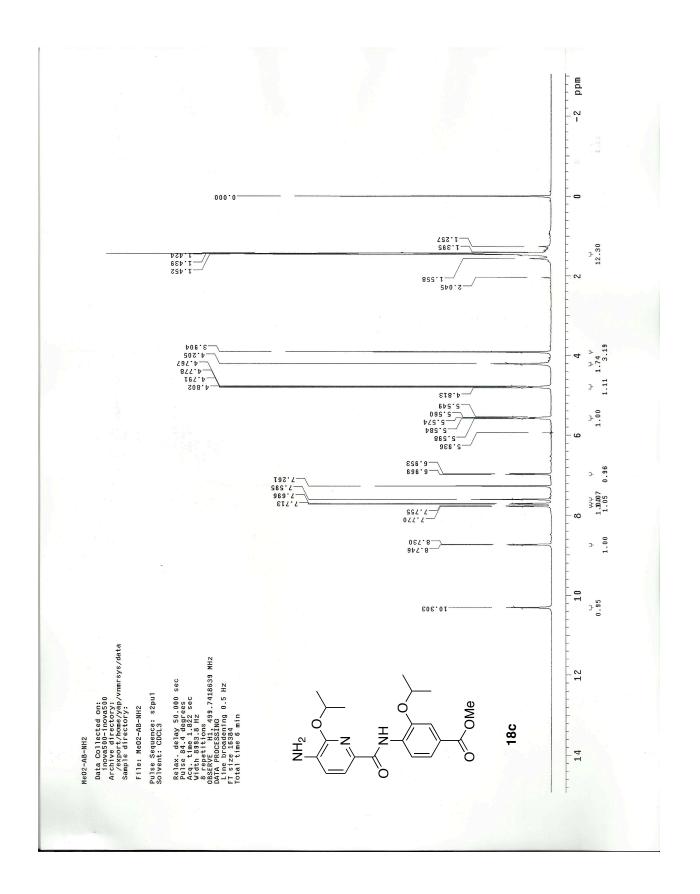


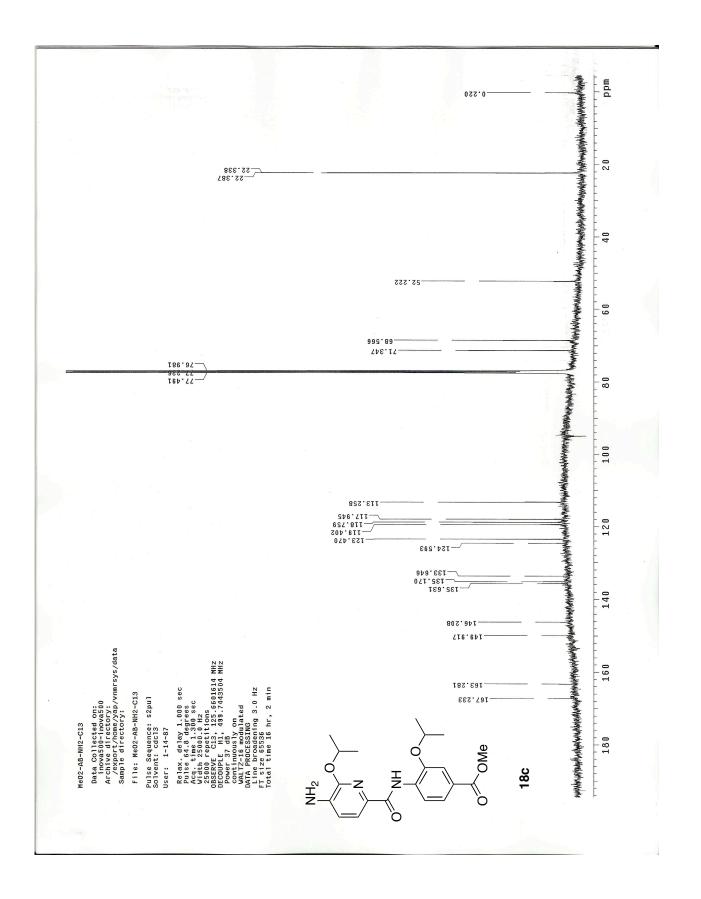


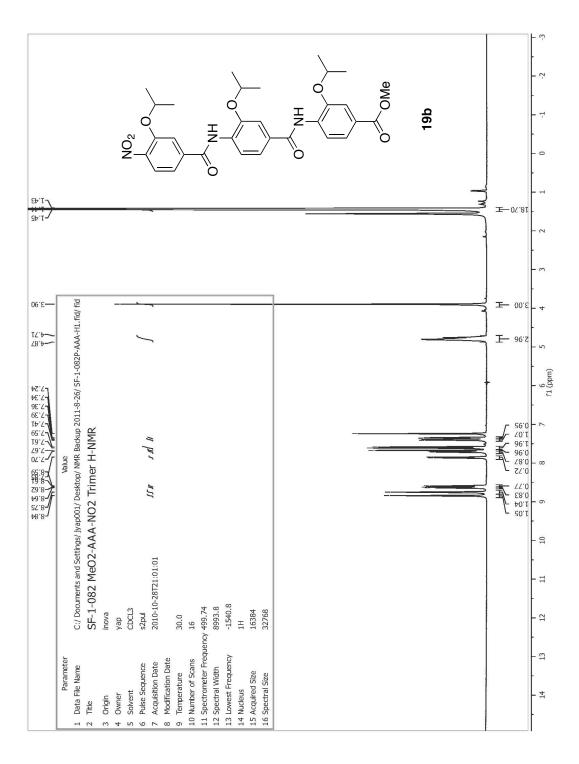


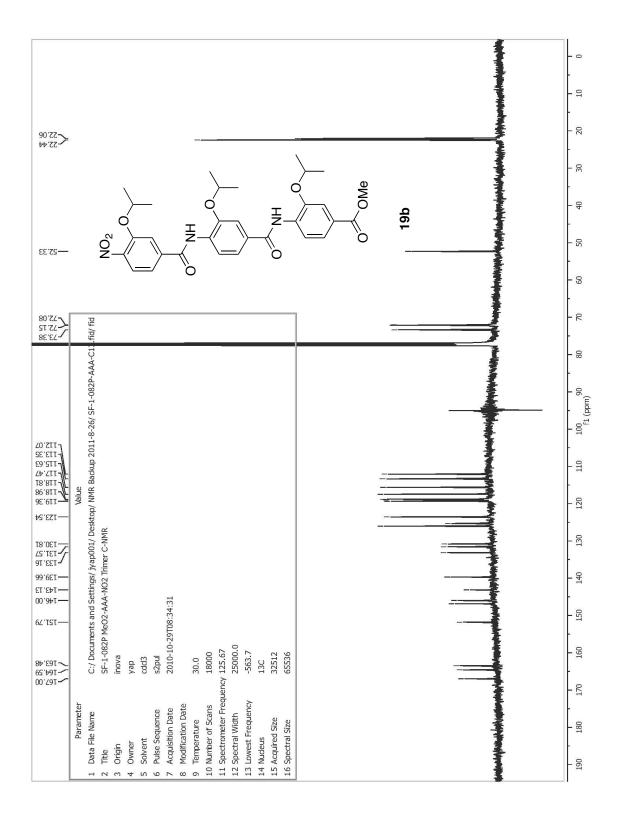


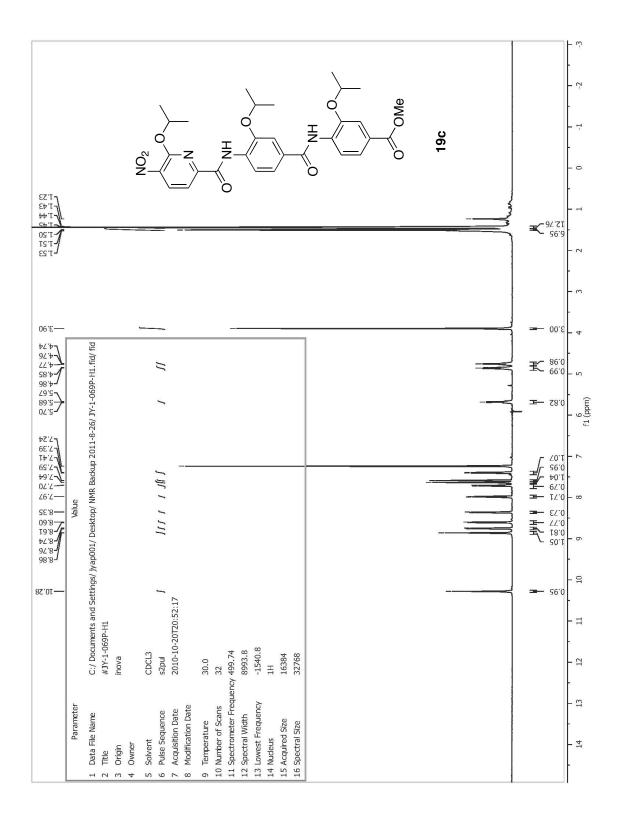


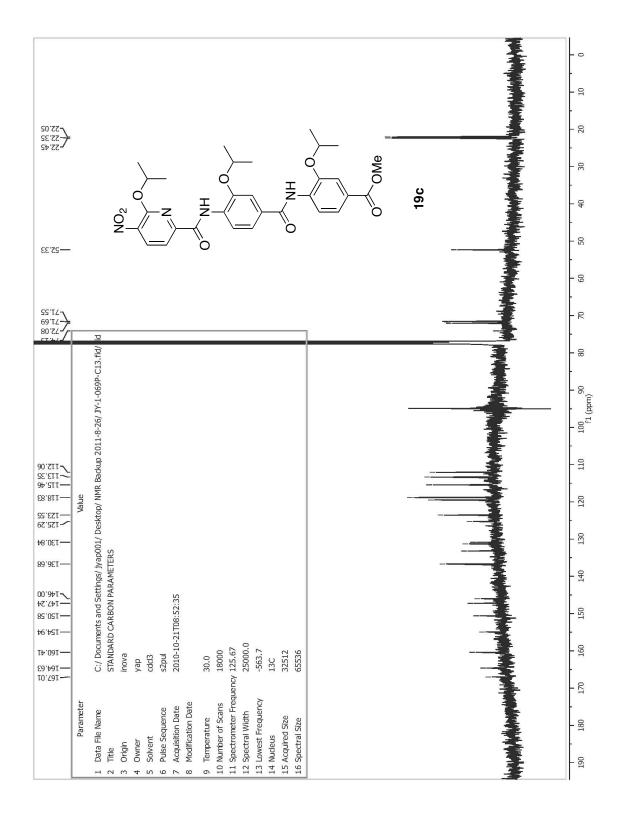


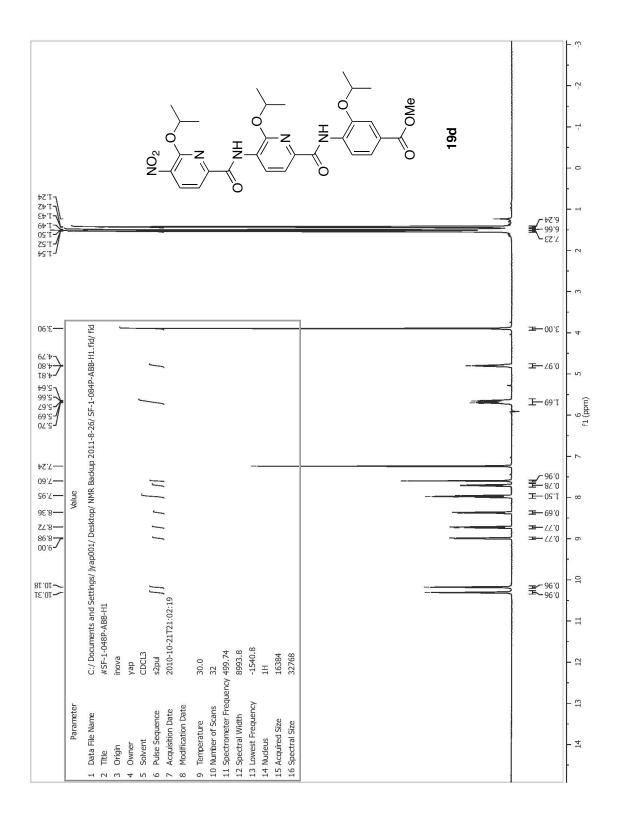


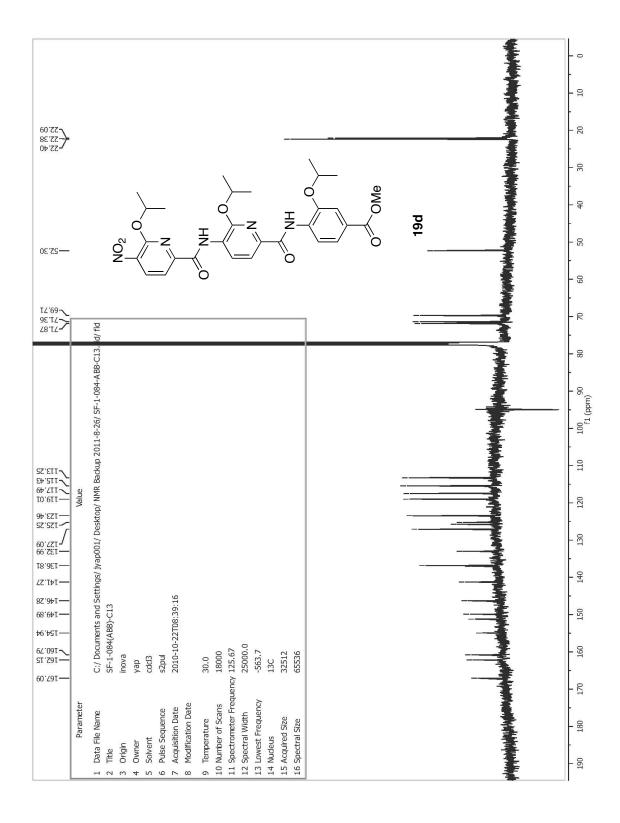


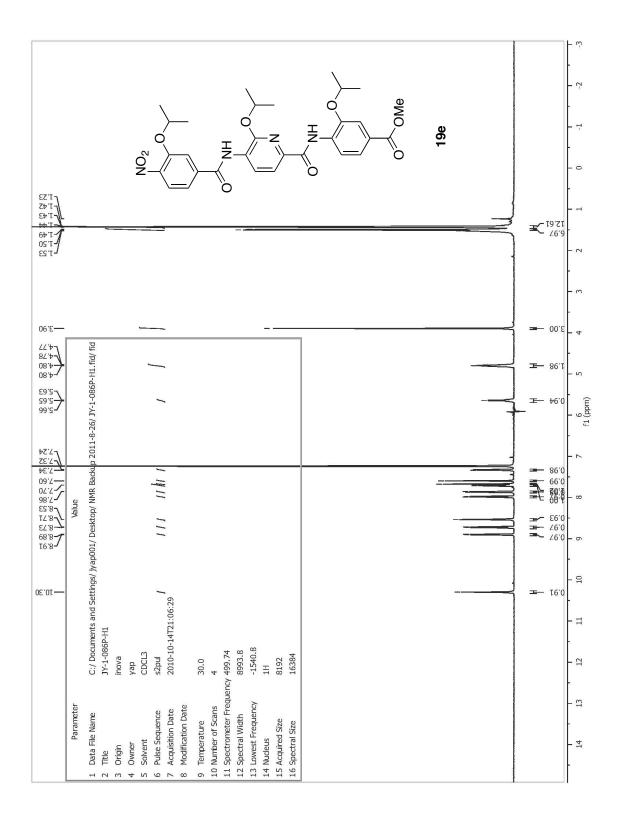


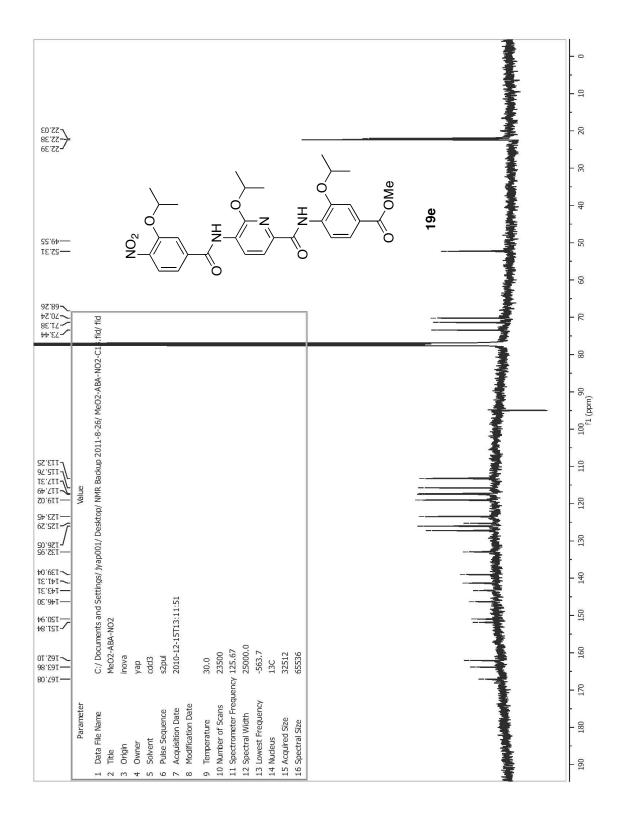


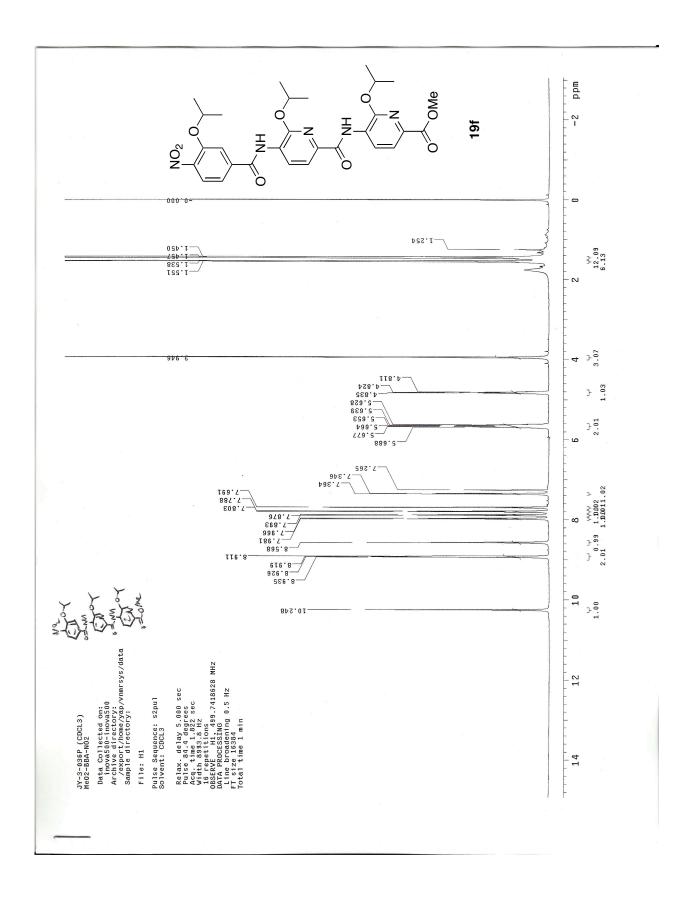


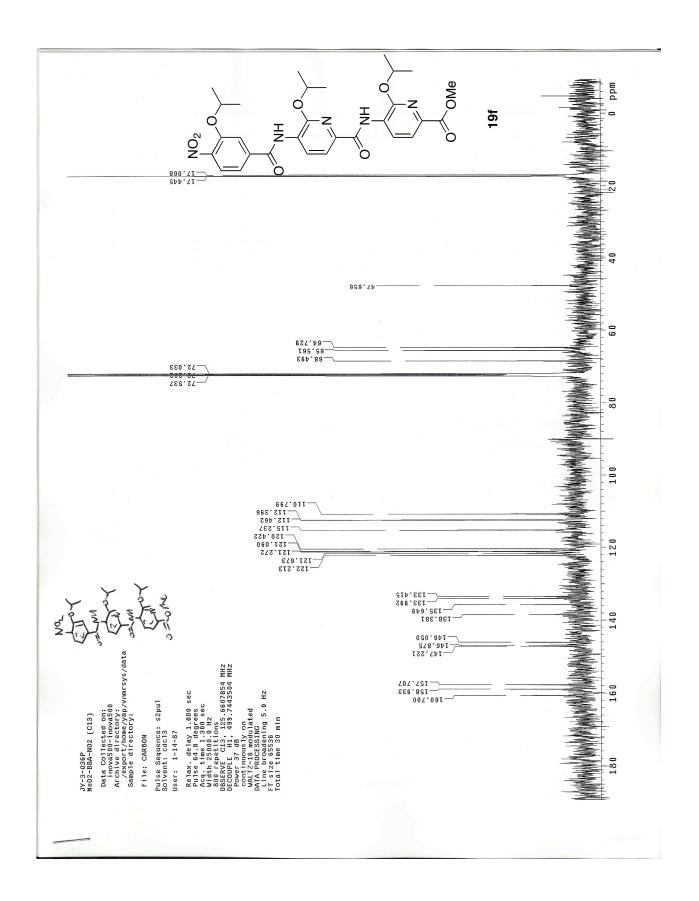


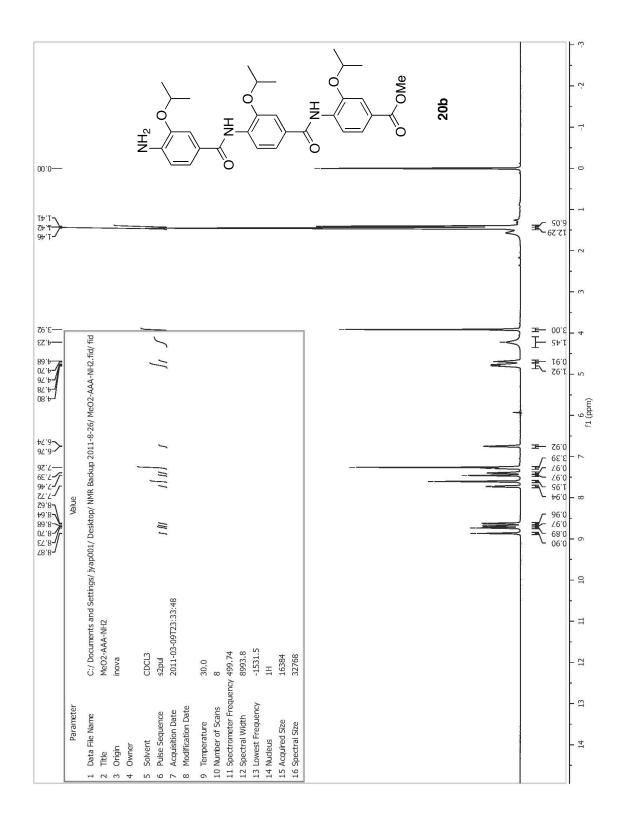


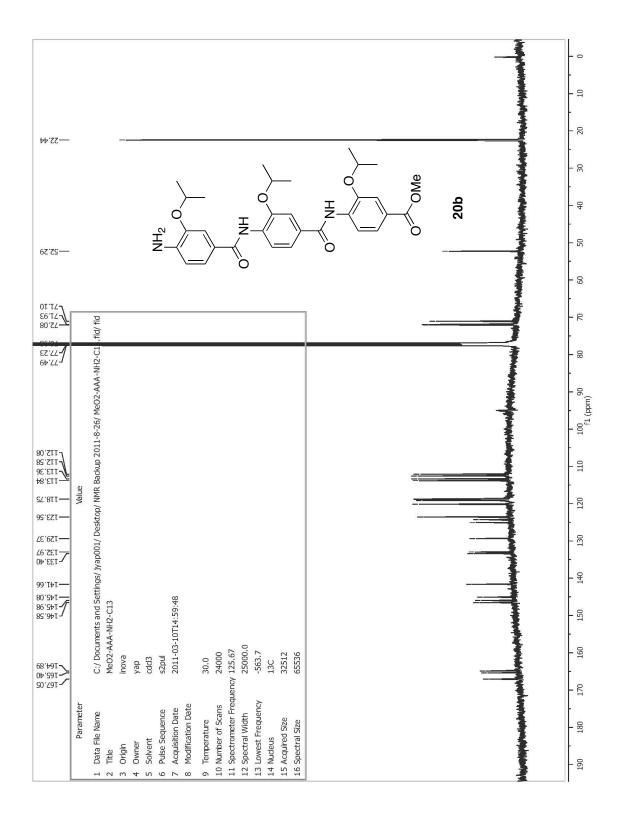


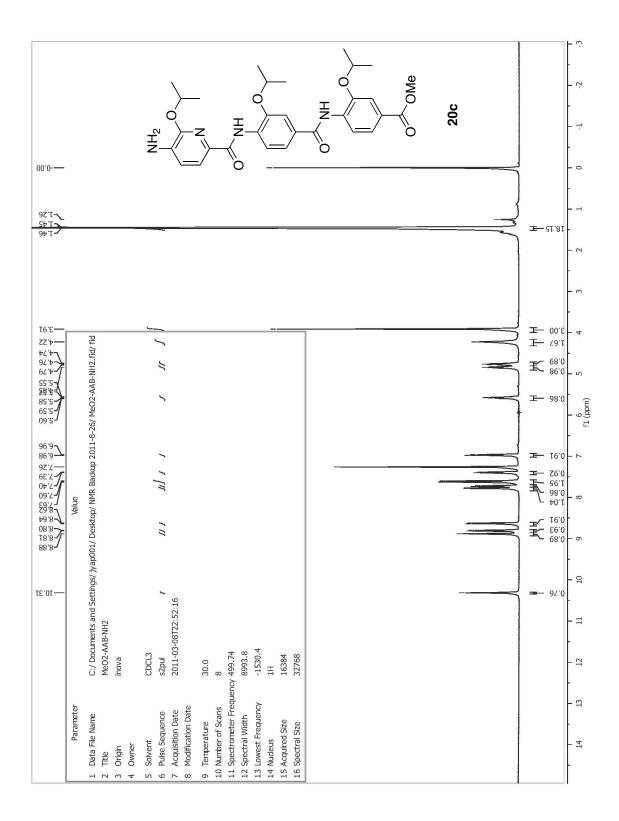


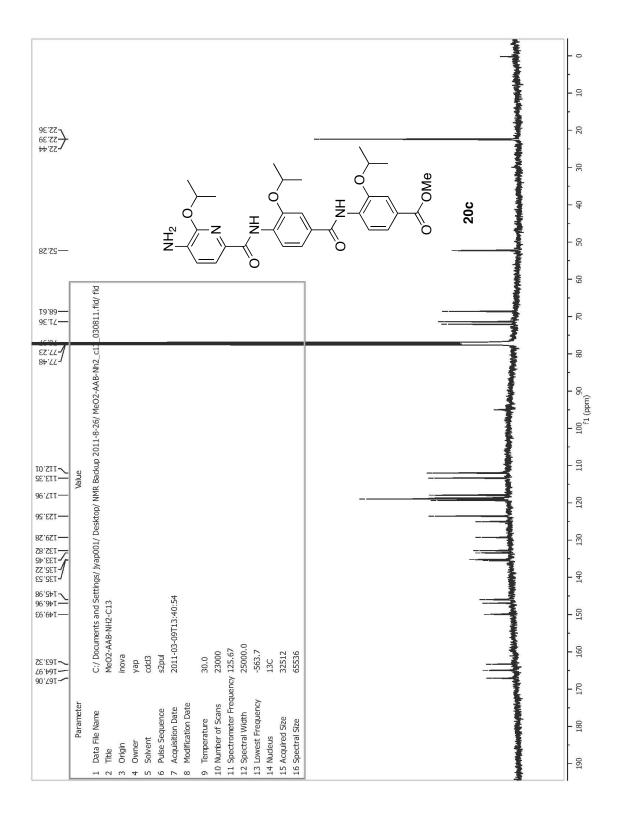


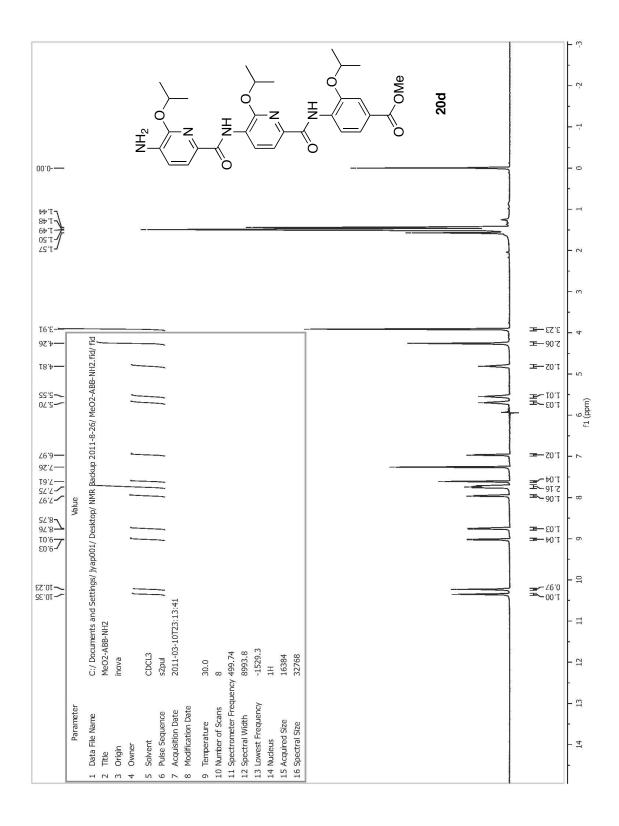


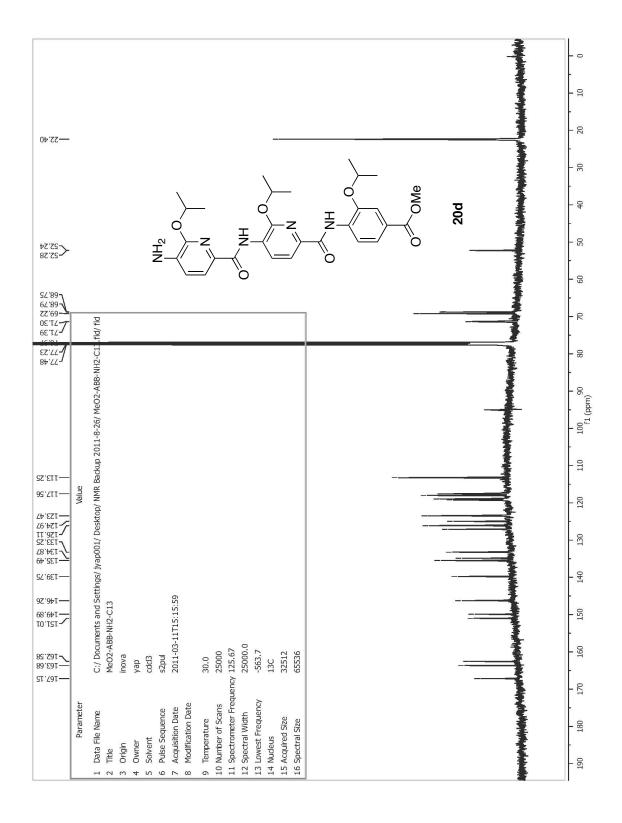


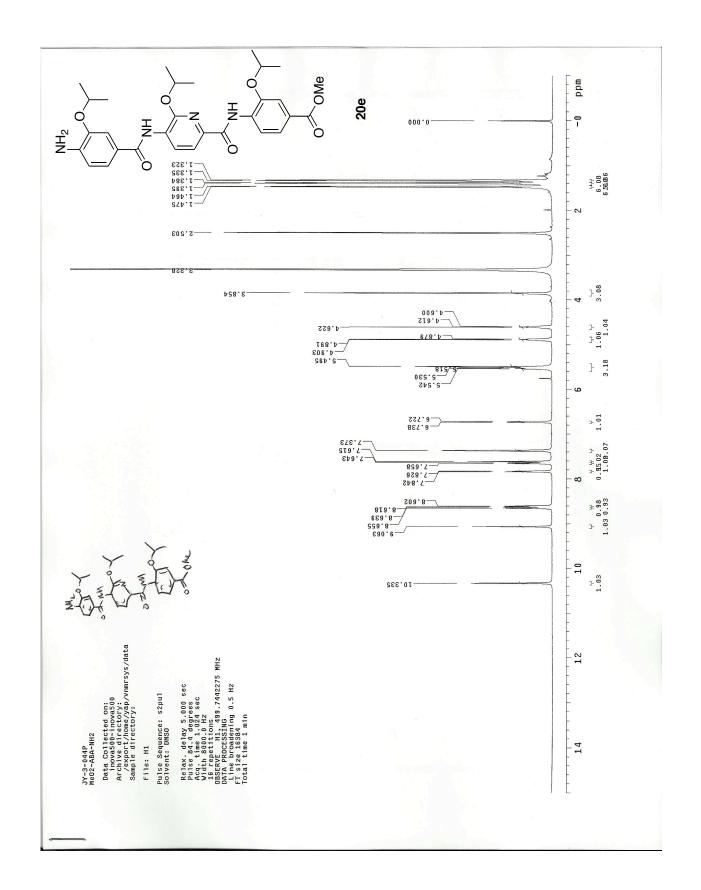


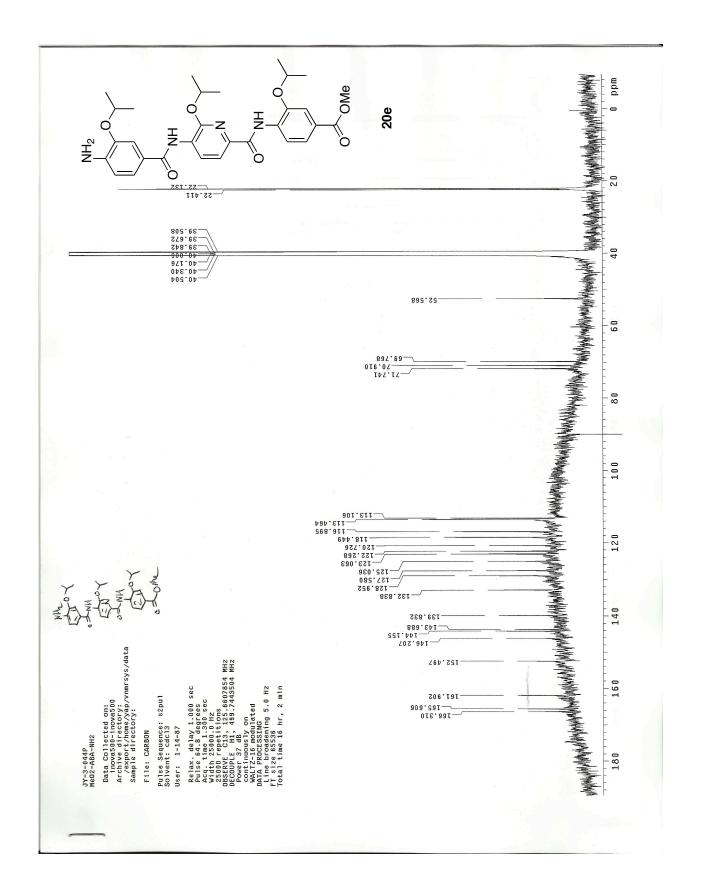


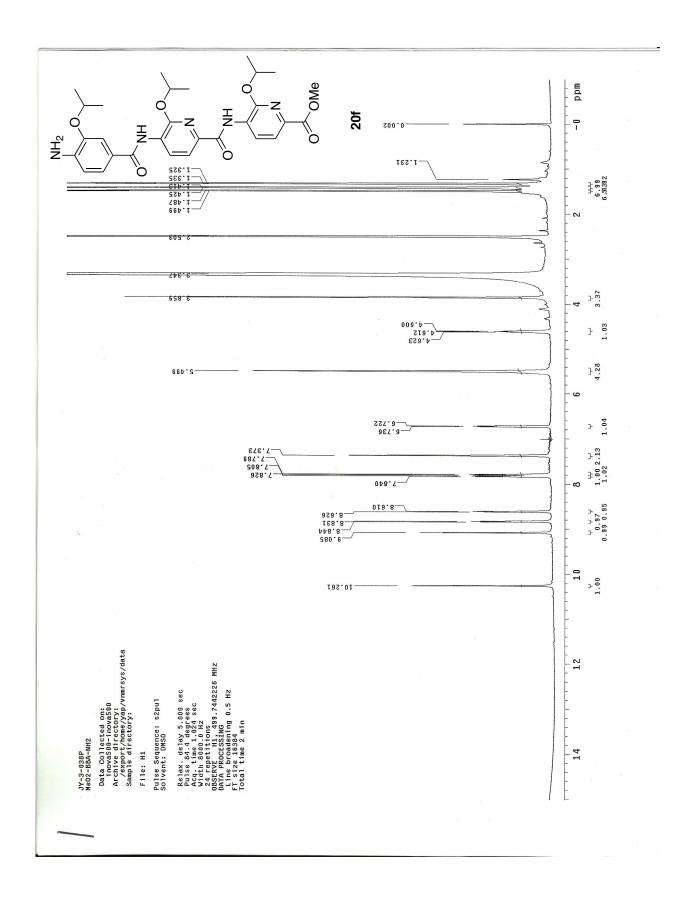


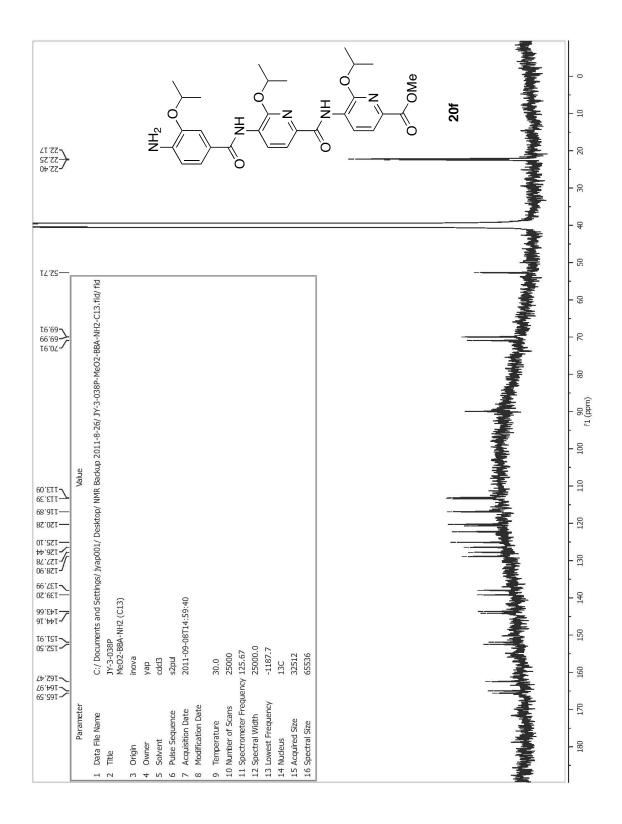


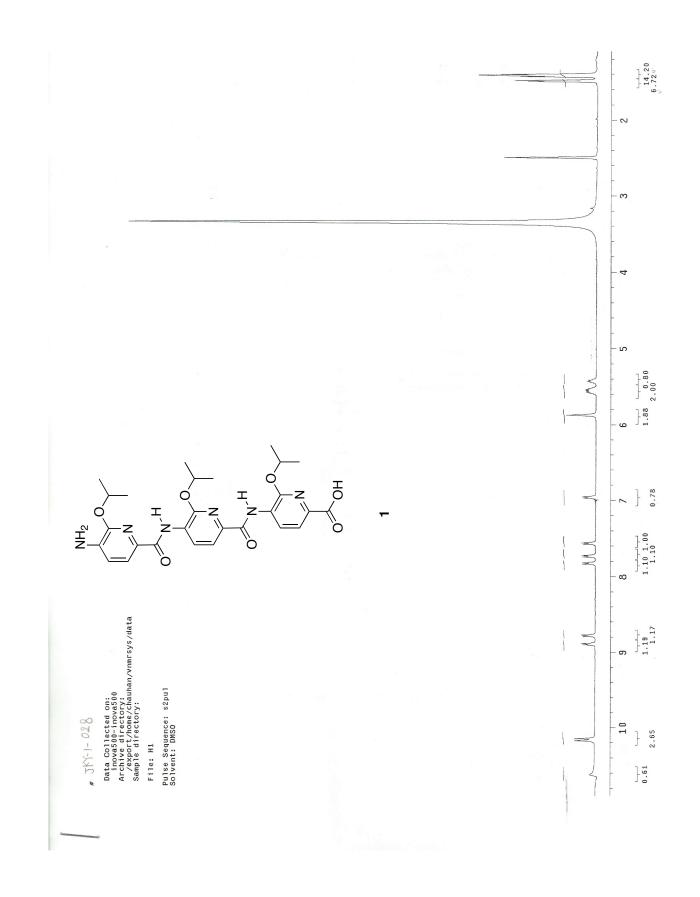


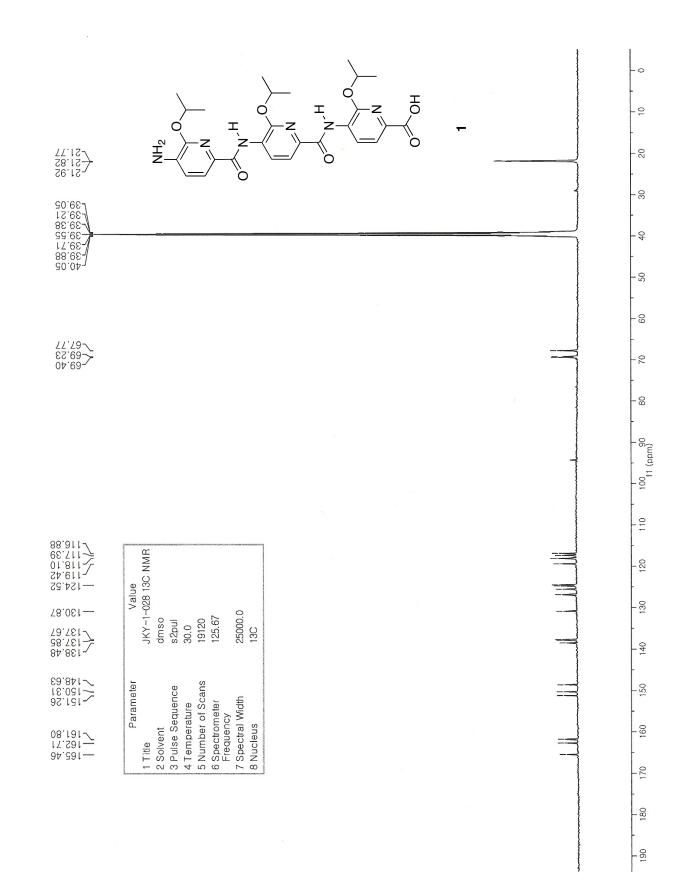




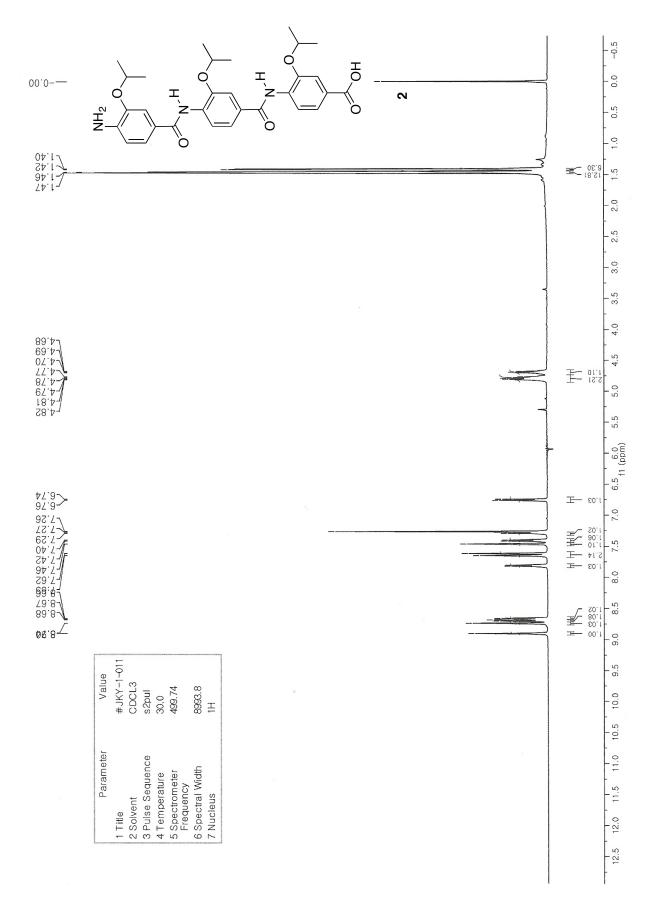


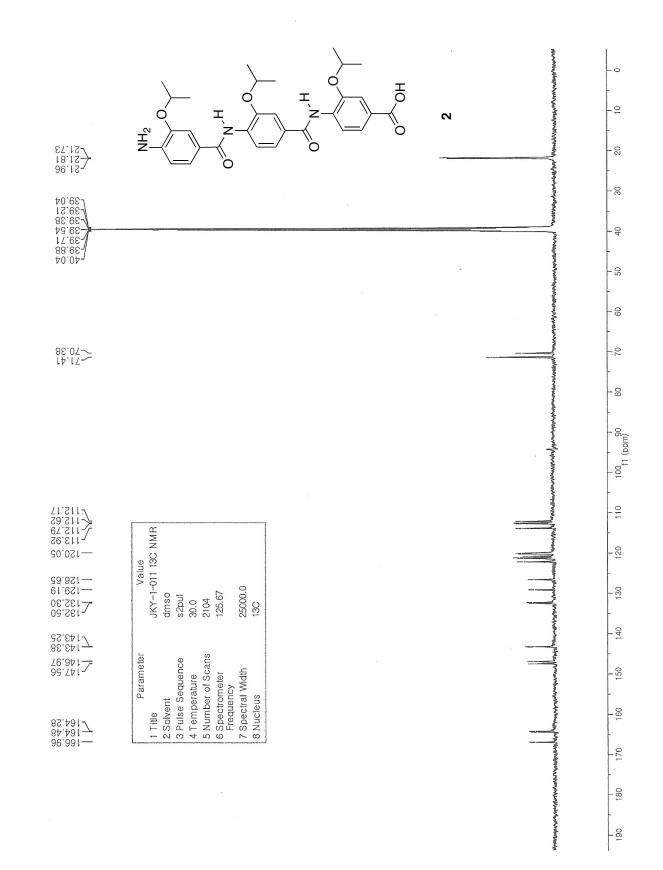


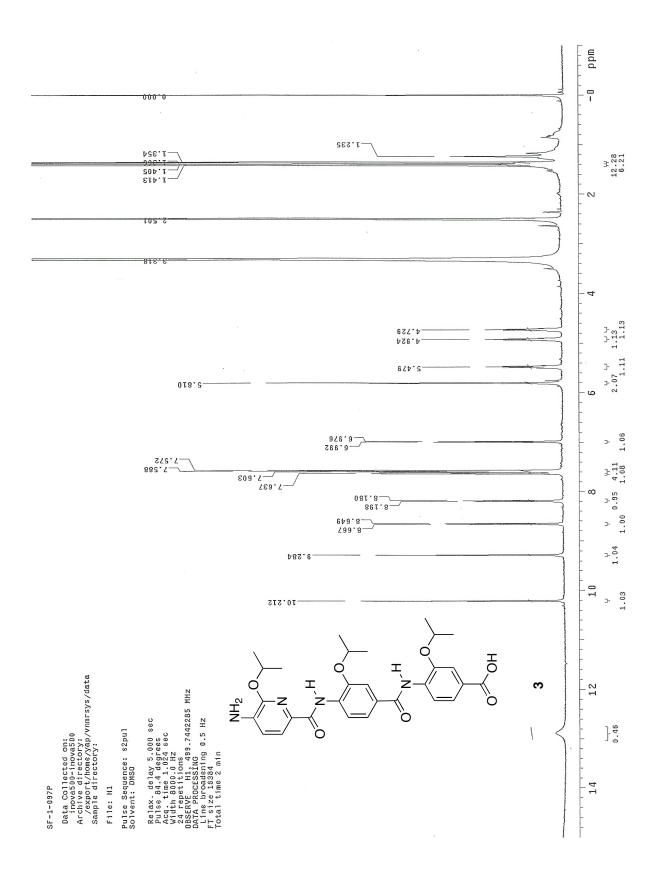


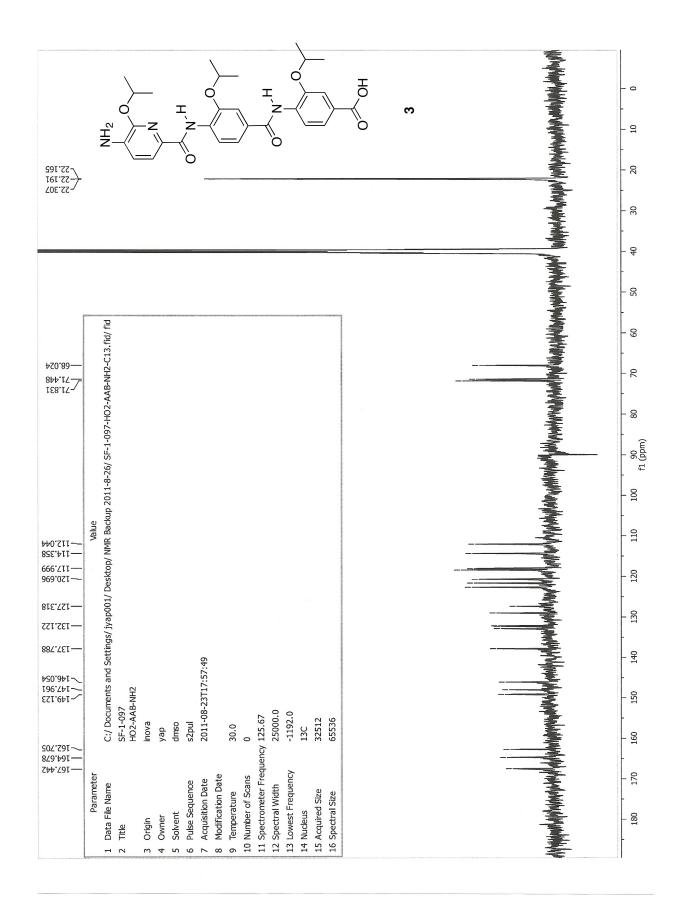


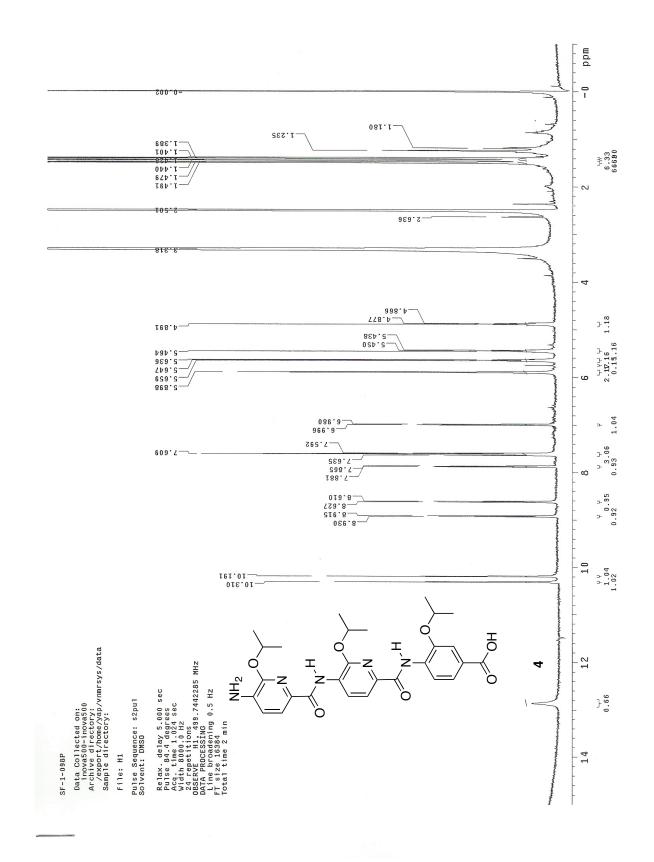
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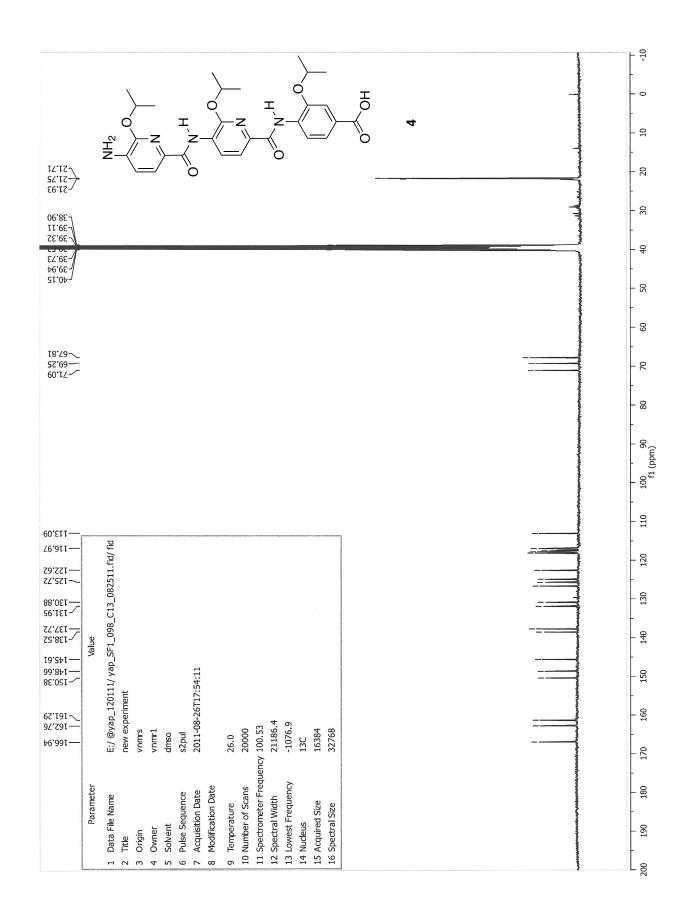




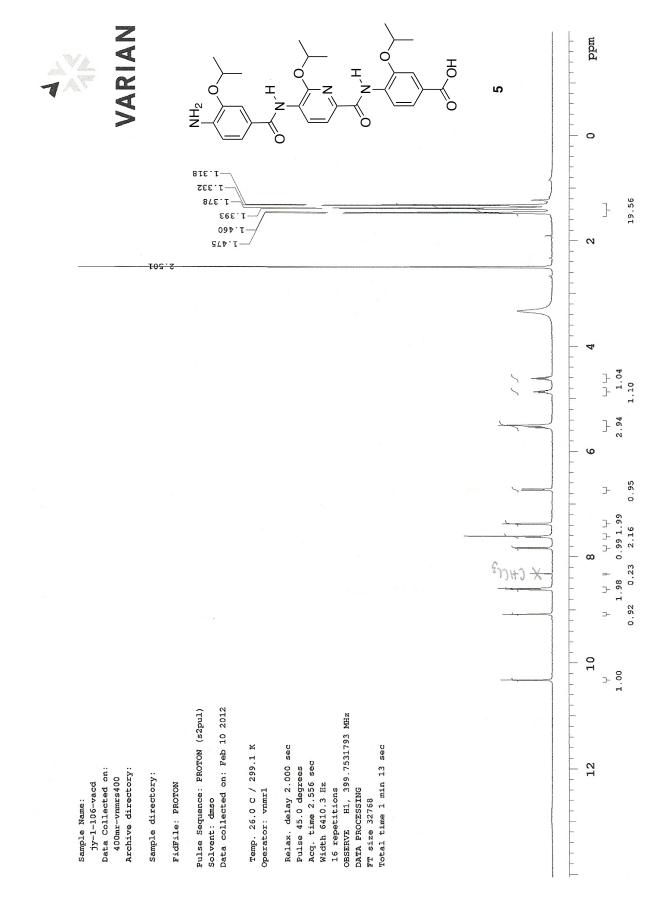




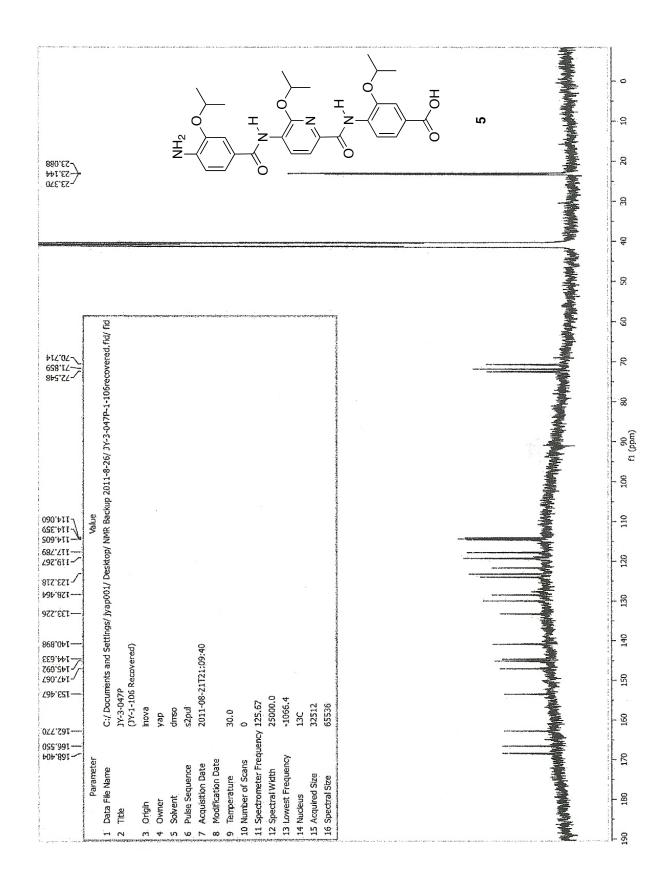


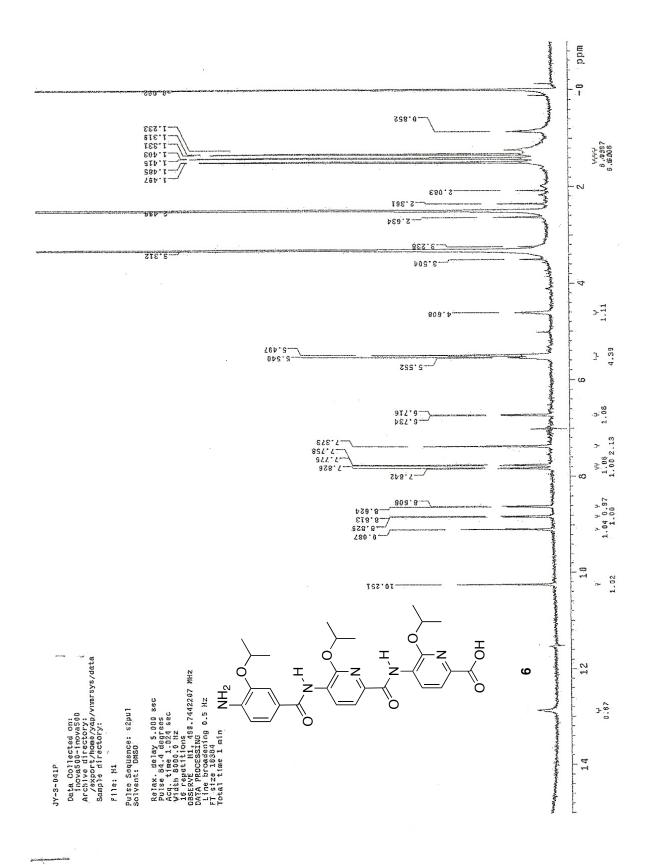


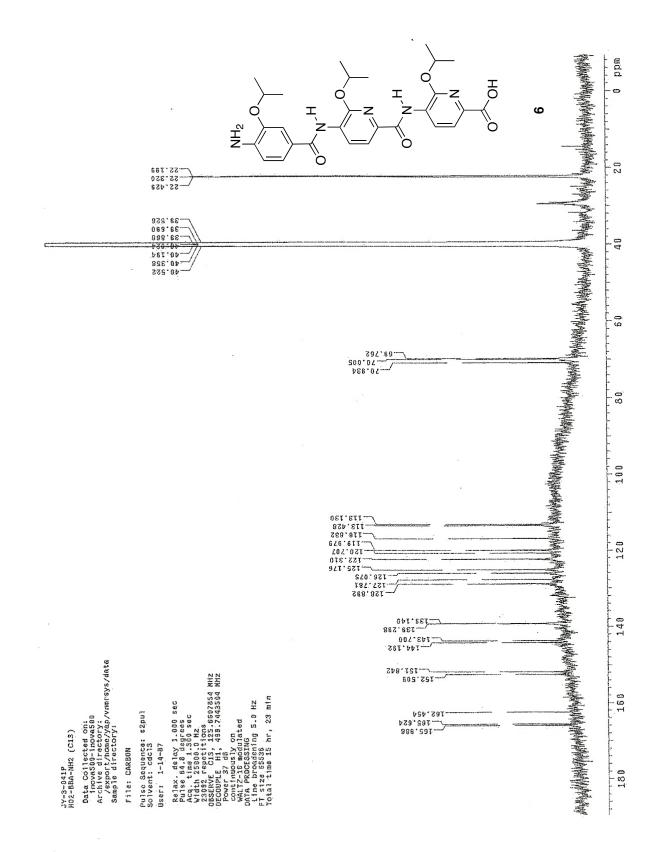
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