

Supporting Information

Pure Shift Approach for Fast and Accurate Extraction of Heteronuclear Couplings

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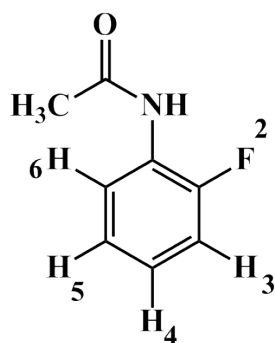
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S1: (a) ^1H NMR of 2-Fluoroacetanilide in CDCl_3 ; (b) Pure shift NMR spectrum of the same molecule depicting only $^nJ_{\text{HF}}$ couplings.

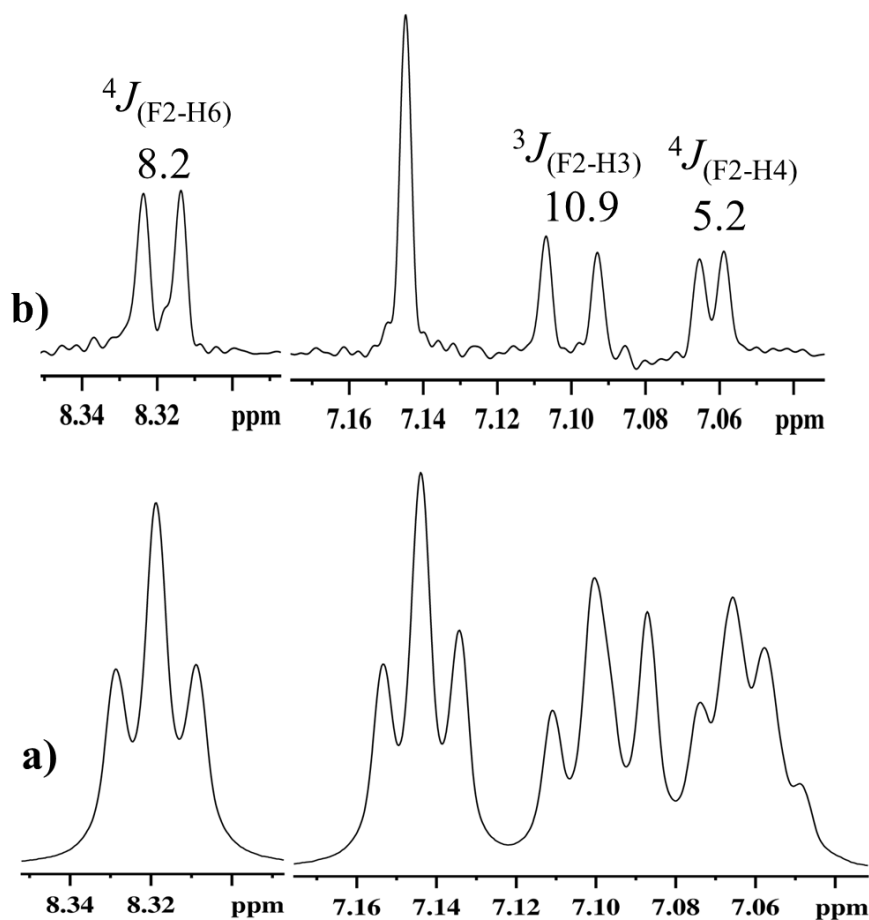
S2: (a) ^1H NMR of 2-fluoropyridine in CDCl_3 ; (b) Pure shift NMR spectrum showing only $^nJ_{\text{HF}}$ couplings.

S3: 2,3-difluoro nitrophenol (e) ^1H NMR data in CDCl_3 . (a) ^1H NMR of 1; (b) Pure shift NMR spectrum showing only $^nJ_{\text{HF}}$ couplings.

(a) The ^1H NMR spectrum of 2-fluoroacetanilide in CDCl_3 . (b) Pure shift spectrum of the same molecule depicting only $^nJ_{\text{HF}}$ couplings. The measured coupling values (doublet separations are reported in Hz) are given at the top of each chemically non-equivalent proton.

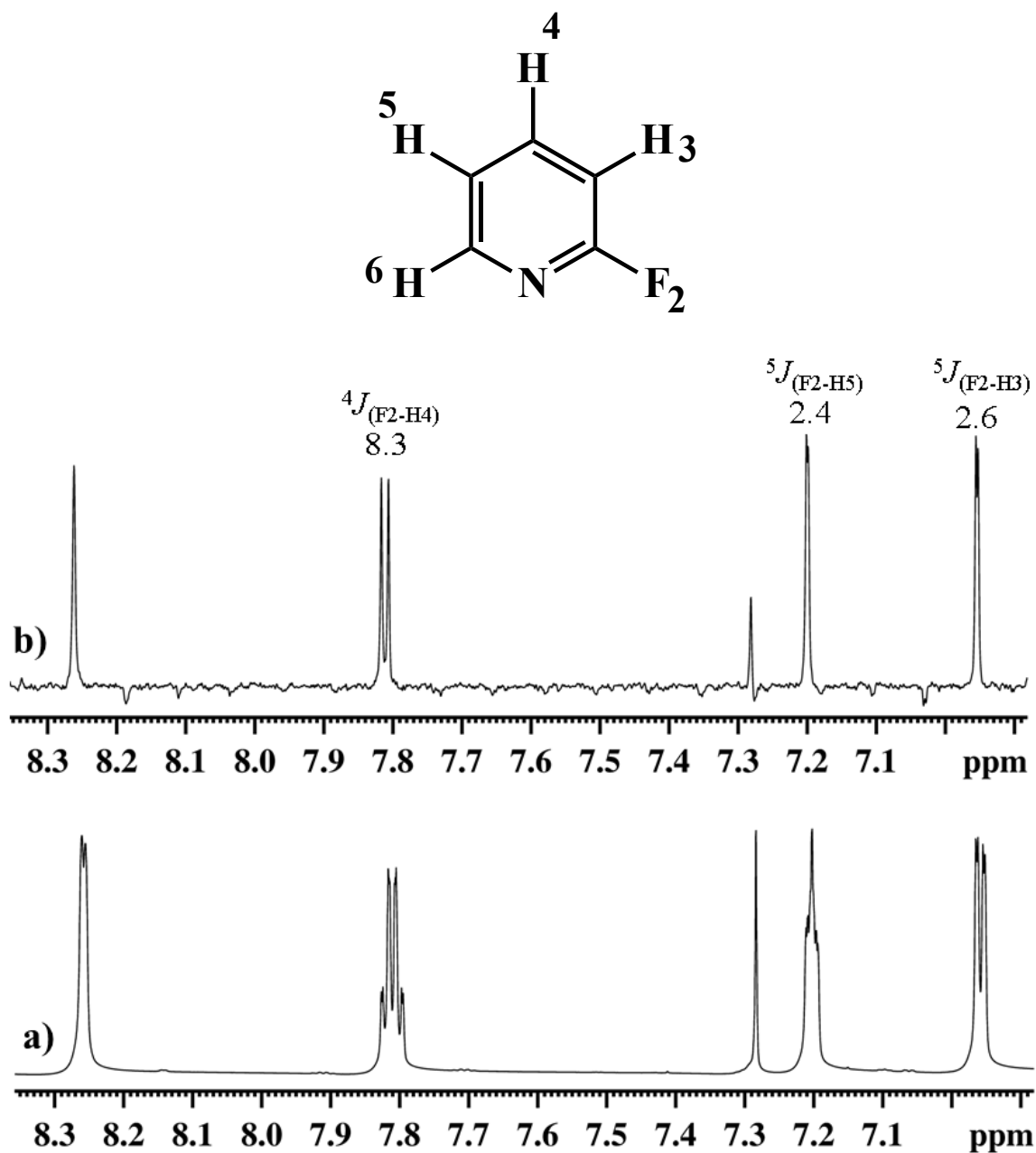


b



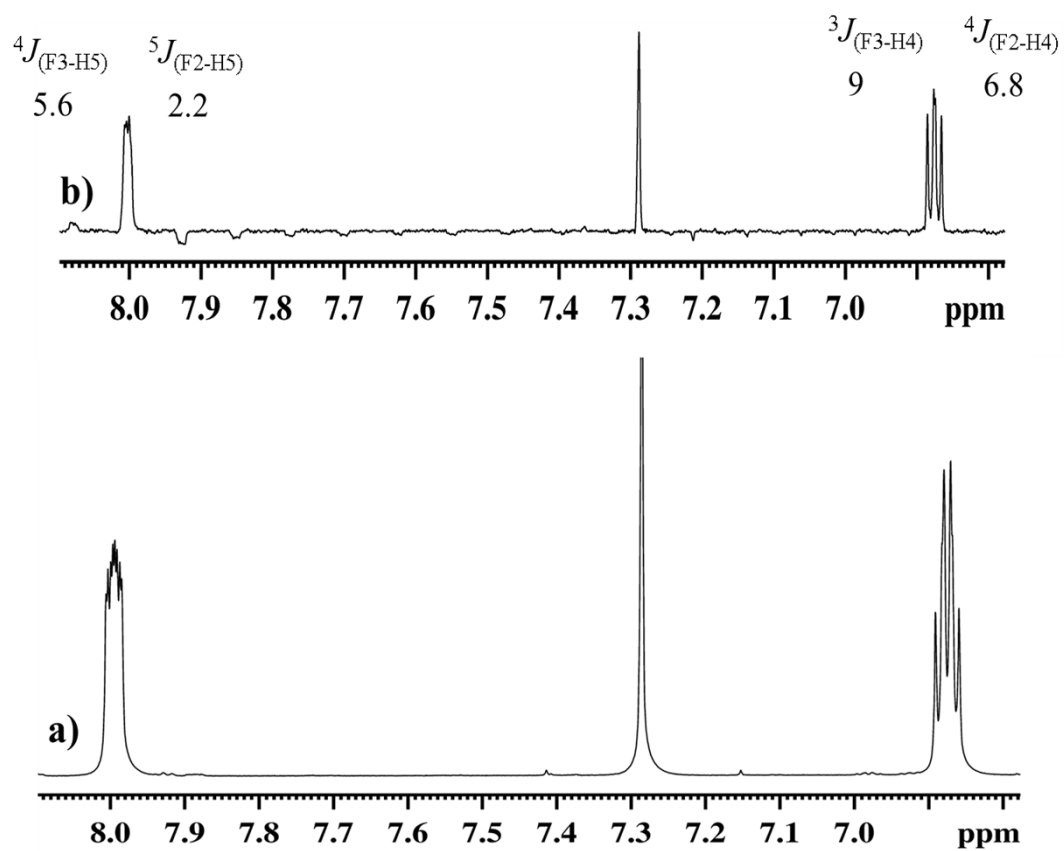
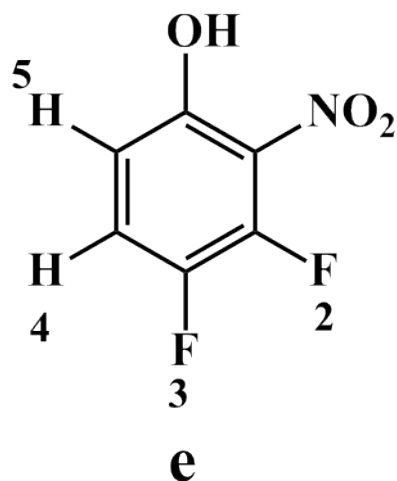
S1

(a) The ^1H NMR spectrum of 2-fluoropyridine in CDCl_3 ; (b) Pure shift spectrum of the same molecule depicting only $^nJ_{\text{HF}}$ couplings. The measured coupling values (doublet separations in Hz) are given at the top of each chemically non-equivalent proton.



S2

(a) The ^1H NMR spectrum of 2,3-difluoro-nitrophenol in CDCl_3 ; (b) Pure shift spectrum of the same molecule depicting only $^nJ_{\text{HF}}$ couplings. The measured coupling values (doublet separations in Hz) are given at the top of each chemically non-equivalent proton.



S3