

Crosslinking of the Pd(acacCN)₂ building unit with Ag(I) salts: dynamic 1D polymers and an extended 3D network[†]

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Supplementary Information

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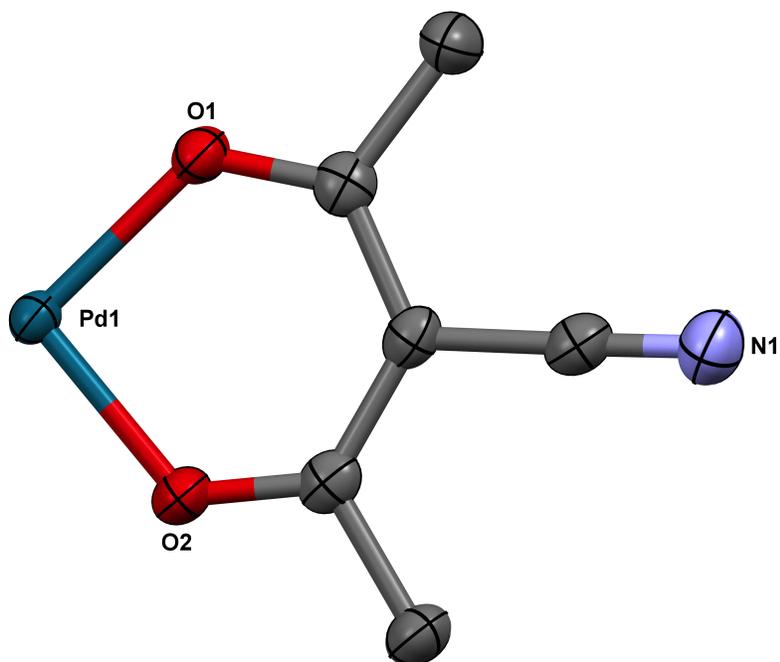


Fig. S1 Displacement ellipsoid plot of the asymmetric unit in **1**. Ellipsoids are drawn at 75 % probability and hydrogen atoms are omitted for clarity.

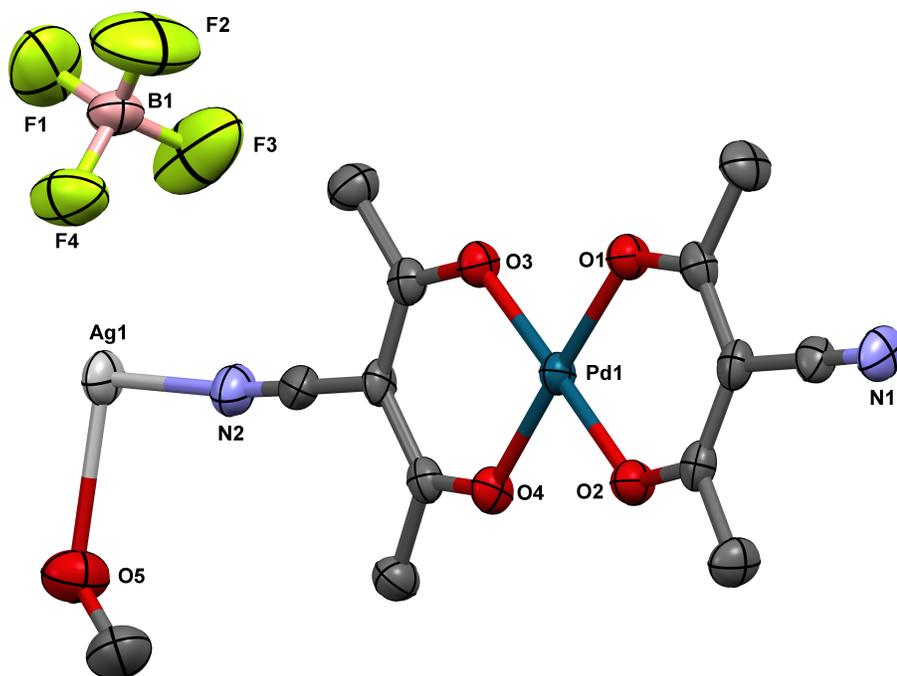


Fig. S2 Displacement ellipsoid plot of the asymmetric unit in **2α**. Ellipsoids are drawn at 75 % probability and hydrogen atoms are omitted for clarity.

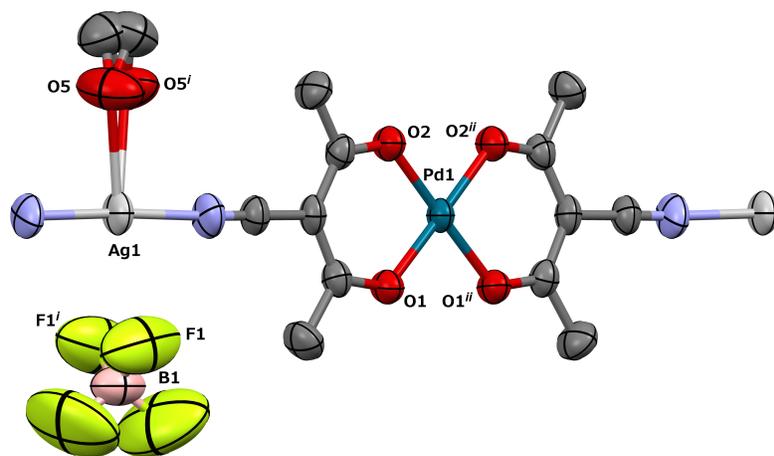


Fig. S3 Displacement ellipsoid plot of 1D chain in 2β . Ellipsoids are drawn at 50 % probability and hydrogen atoms are omitted for clarity. Symmetry operators: $i = 2-x, y, 0.5-z$; $ii = 1-x, y, 1.5-z$.

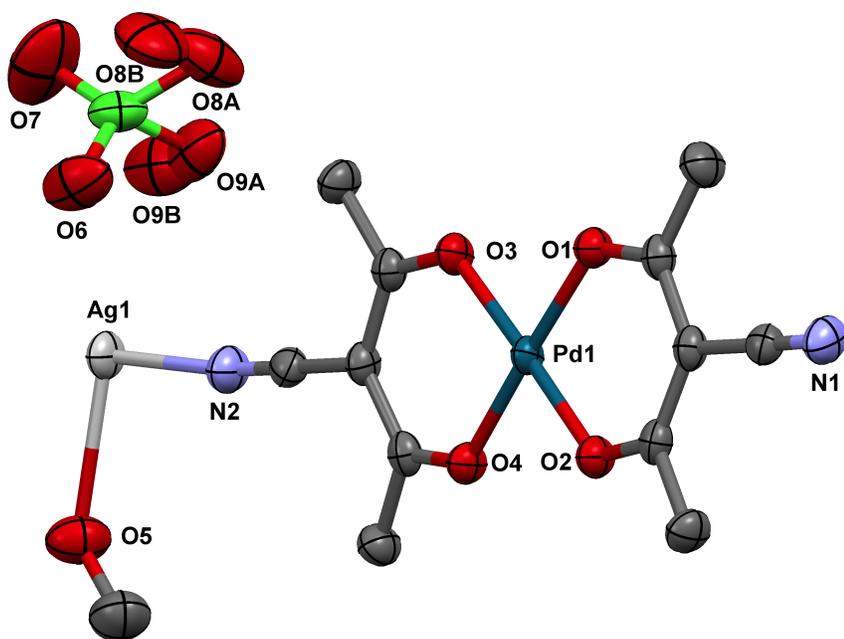


Fig. S4 Displacement ellipsoid plot of the asymmetric unit in 3α . Ellipsoids are drawn at 75 % probability and hydrogen atoms are omitted for clarity.

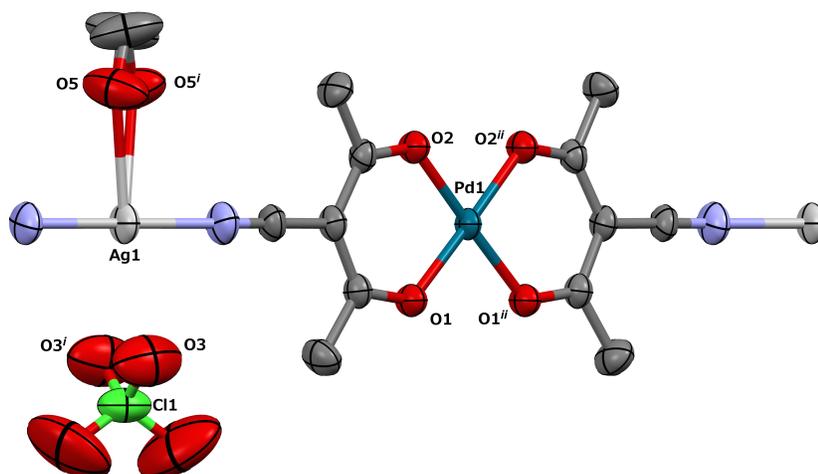


Fig. S5 Displacement ellipsoid plot of the 1D chain in 3β . Ellipsoids are drawn at 50 % probability and hydrogen atoms are omitted for clarity. Symmetry operators: $i = 2-x, y, 0.5-z$; $ii = 1-x, y, 1.5-z$.

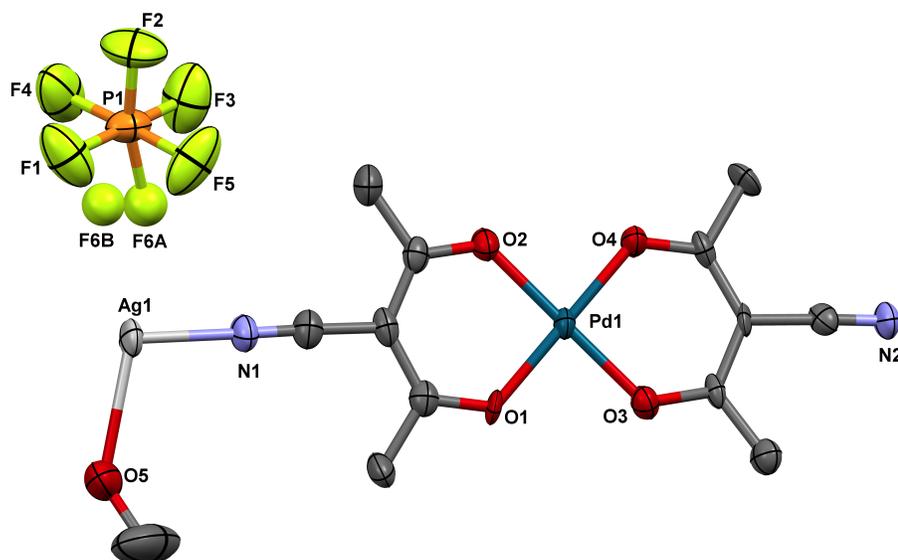


Fig. S6 Displacement ellipsoid plot of the asymmetric unit in 5 . Ellipsoids are drawn at 75 % probability and hydrogen atoms are omitted for clarity.

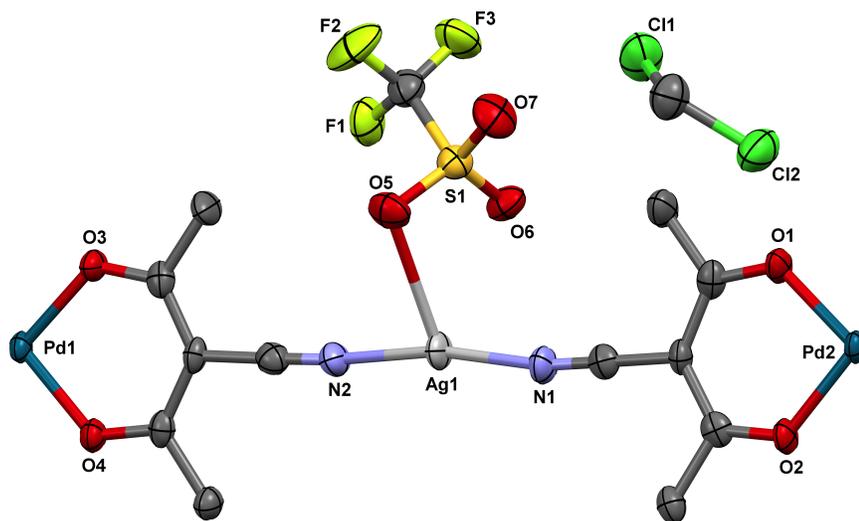


Fig. S7 Displacement ellipsoid plot of the asymmetric unit in **6**. Ellipsoids are drawn at 75 % probability and hydrogen atoms are omitted for clarity.

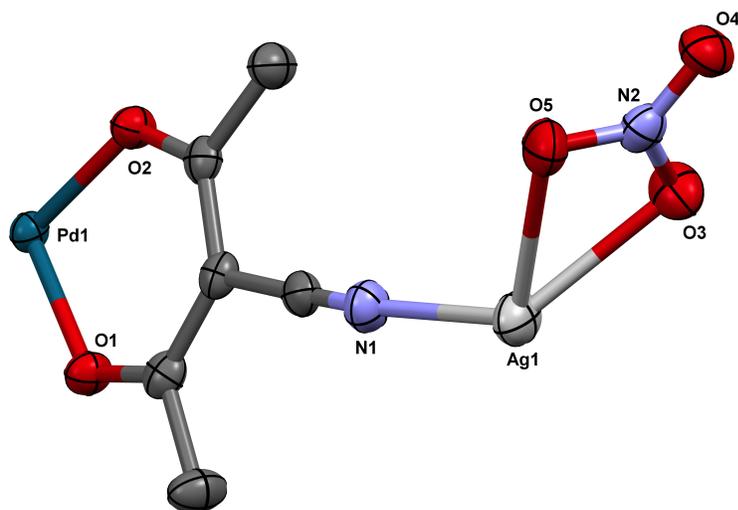


Fig. S8 Displacement ellipsoid plot of the asymmetric unit in **7**. Ellipsoids are drawn at 75 % probability and hydrogen atoms are omitted for clarity.

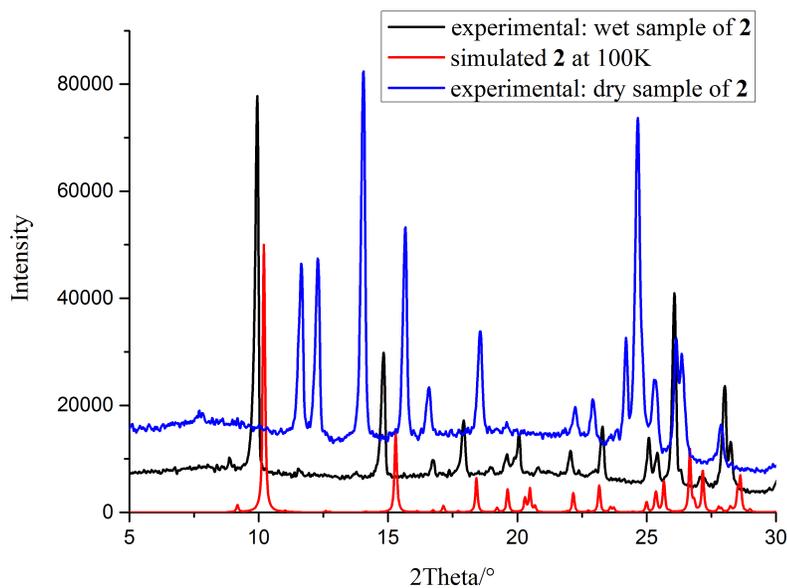


Fig. S9 Simulated and experimental powder patterns for wet and dry samples of **2**.

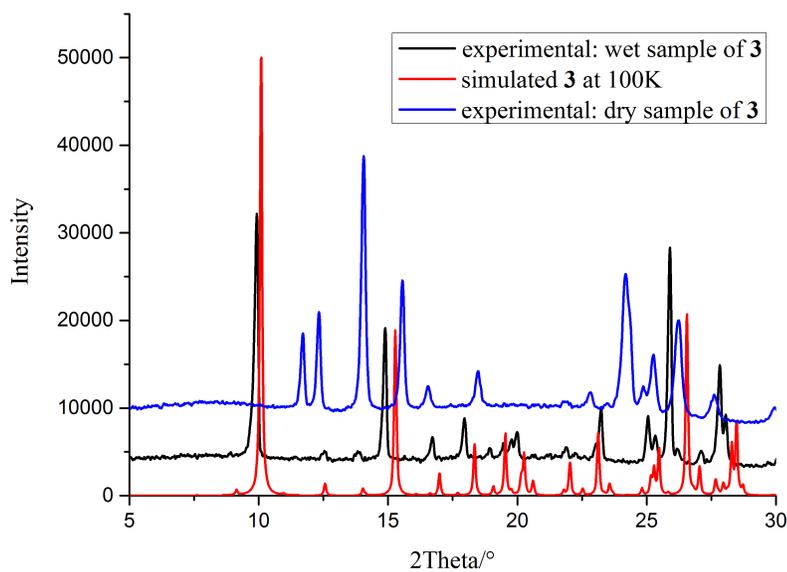


Fig. S10 Simulated and experimental powder patterns for wet and dry samples of **3**.

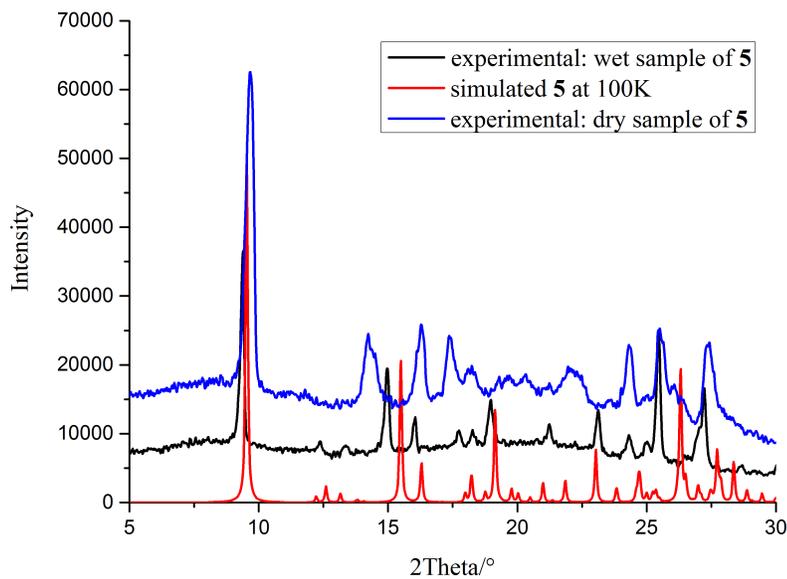


Fig. S11 Simulated and experimental powder patterns for wet and dry samples of **5**.

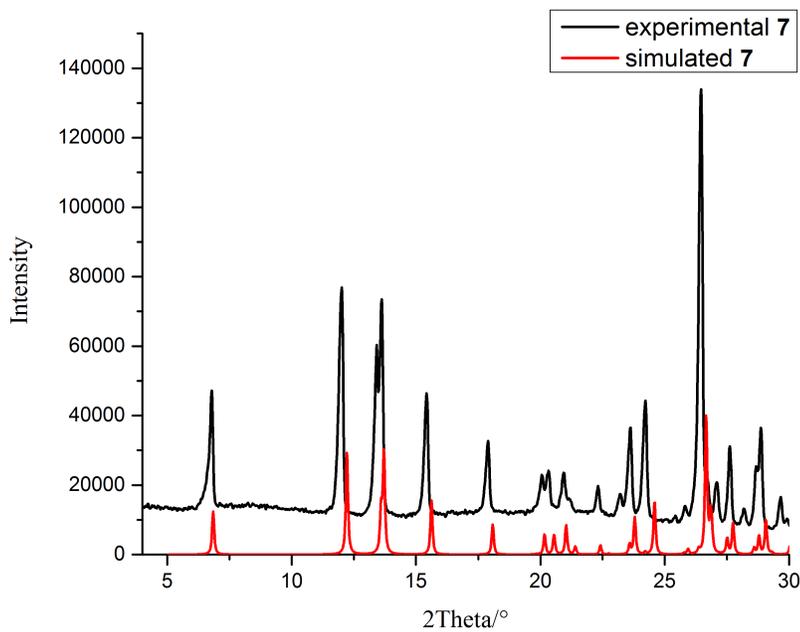


Fig. S12 Simulated and experimental powder patterns of **7**.

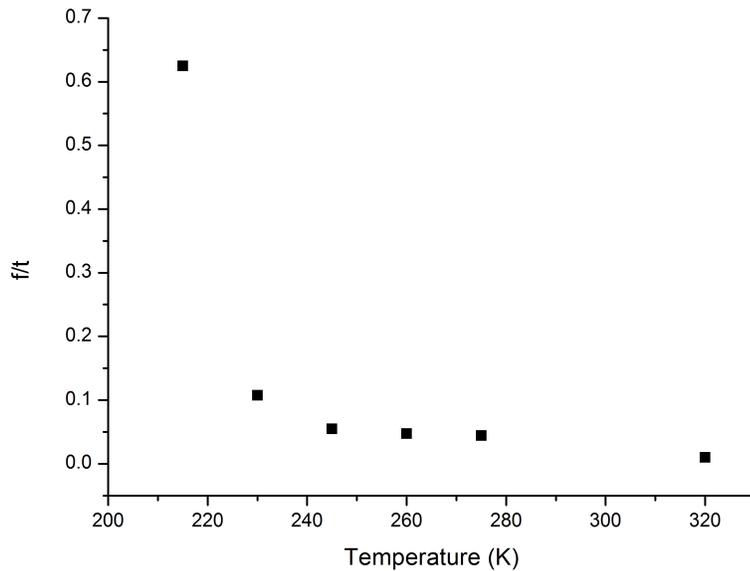


Fig. S13 Ratio $I(f/t)$ as a function of temperature for the sum of all reflection intensities in **2**. In this preliminary experiment, the phase transition temperature was determined approximately.

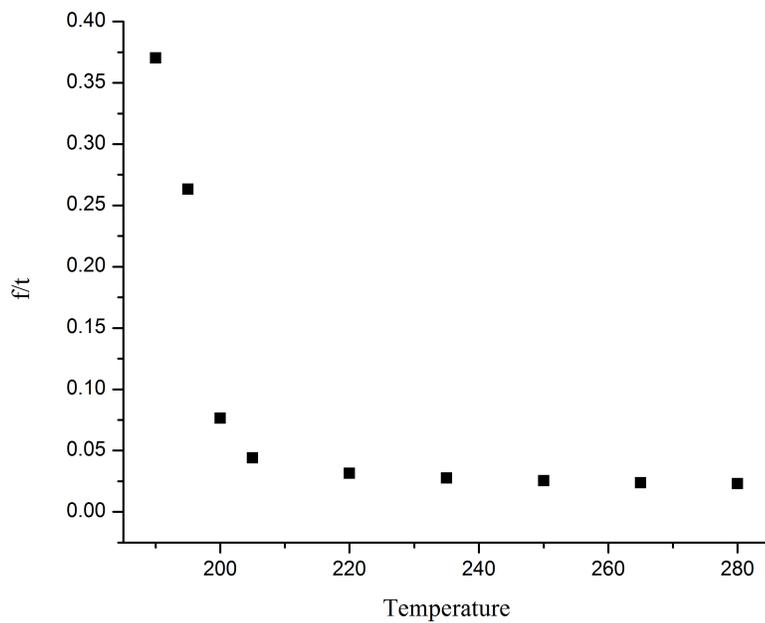


Fig. S14 Ratio $I(f/t)$ as a function of temperature for the sum of all reflection intensities in **3**. In this preliminary experiment, the phase transition temperature was determined approximately.

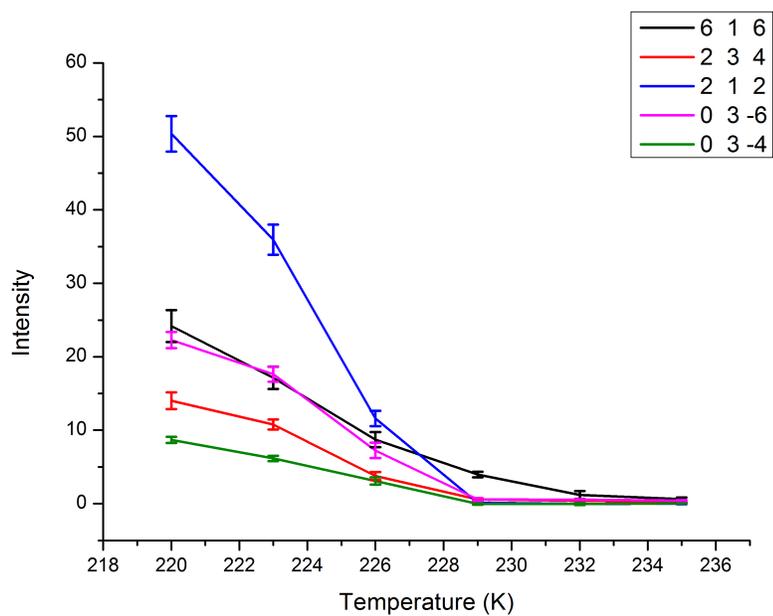


Fig. S15 Intensity of individual reflections hkl , $h+k = 2n+1$ in **2** as a function of temperature.

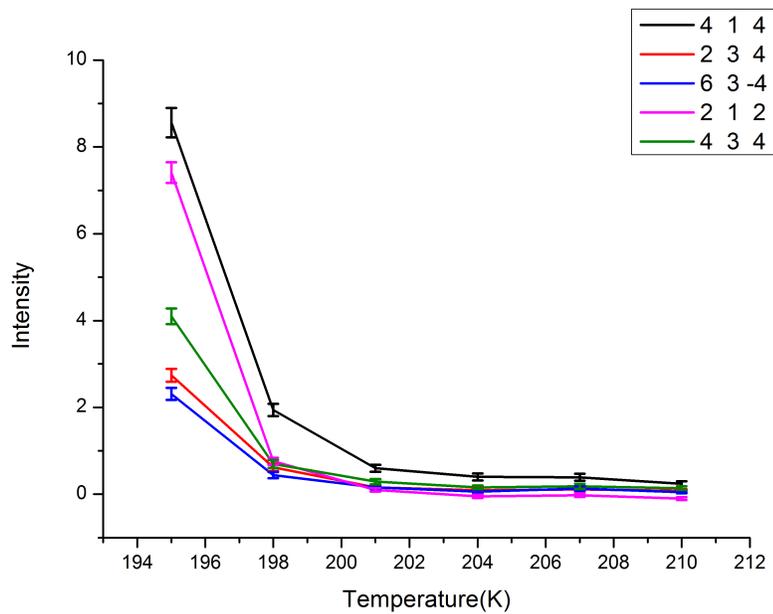


Fig. S16 Intensity of individual reflections hkl , $h+k = 2n+1$ in **3** as a function of temperature.

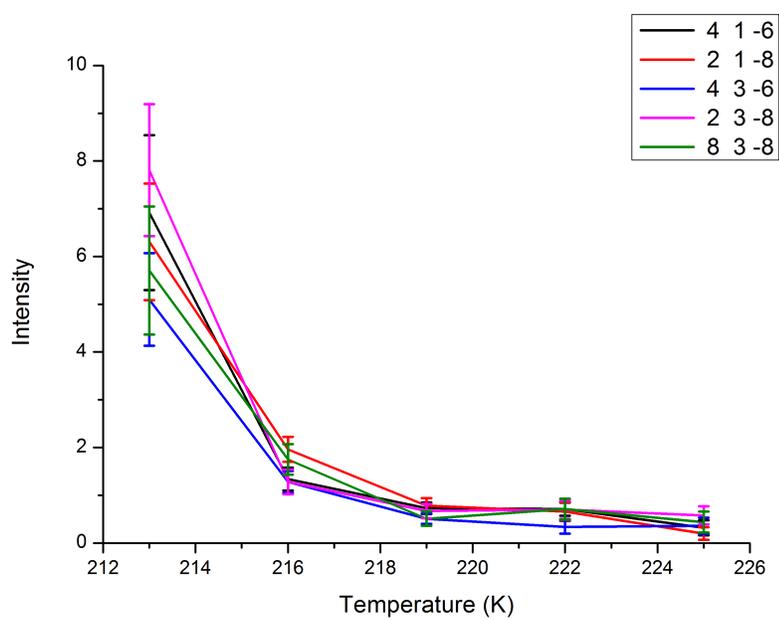


Fig. S17 Intensity of individual reflections hkl , $h+k = 2n+1$ in **4** as a function of temperature.

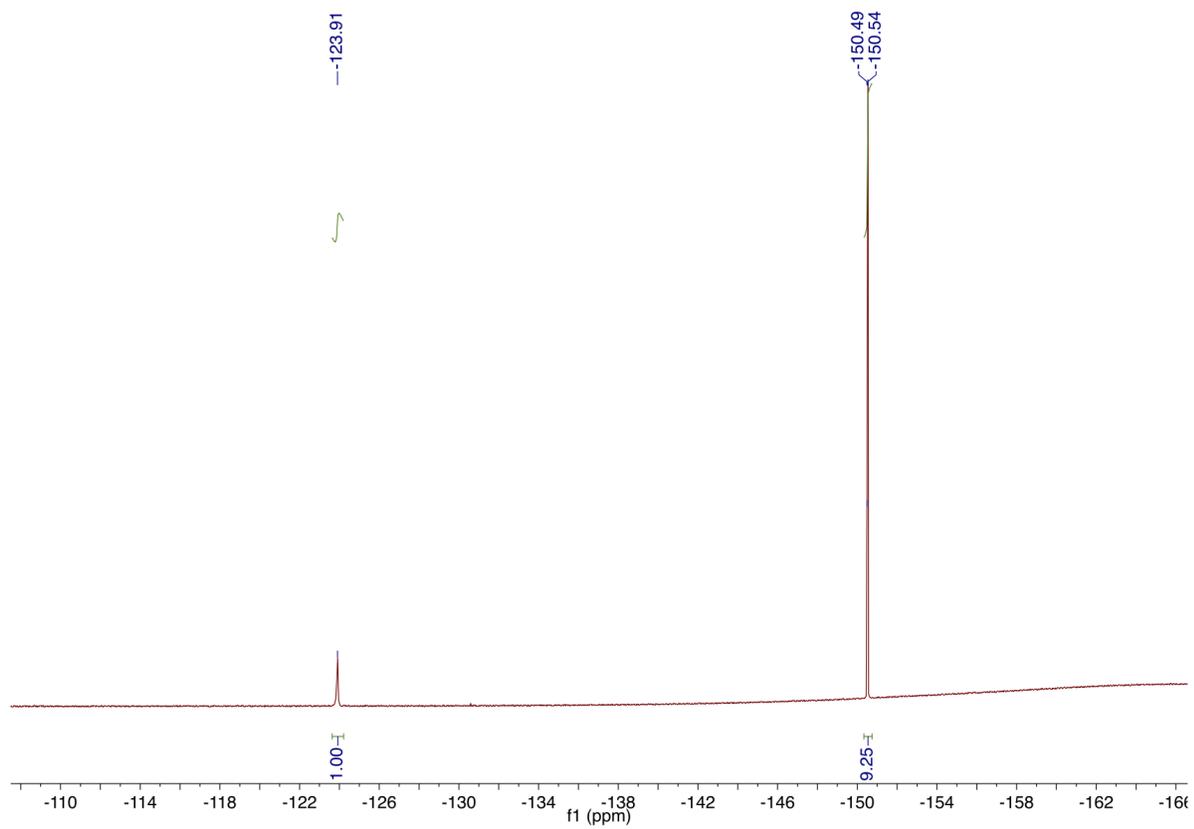


Fig. S18 ^{19}F NMR spectrum of compound **4** in D_2O .