

Electronic Supplementary Information

(9 pages)

A quest for supramolecular gelators: Silver(I) complexes with quinoline-urea derivatives.

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Contents

XRPD for compounds PQ5U and PQ8U	Page 1
¹ H-NMR spectra for compounds PQ6U, PiQ5U, DQ5U and PPT4U	page 2
Solution from powder data and refinement for compound PQ6U	page 3
XRPD for compounds PiQ5U, DQ5U and PPT4U Form I and II	page 3
Map of fluorescence spectra vs excitation wavelength	Page 5
XRPD for the complex [Ag(PQ6U) ₂]NO ₃ ·CH ₃ CN and TGA trace	page 6
XRPD for the complex [Ag(PQ6U) ₂]NO ₃ Form I and Form II	page 7
ORTEP drawings for compounds PiQ5U, DQ5U, PPT4U Form 1 and 2 and for complexes [Ag(PQ6U) ₂]NO ₃ ·CH ₃ CN and [Ag(PQ6U) ₂]NO ₃ Form 1 and 2	page 8
	page 9

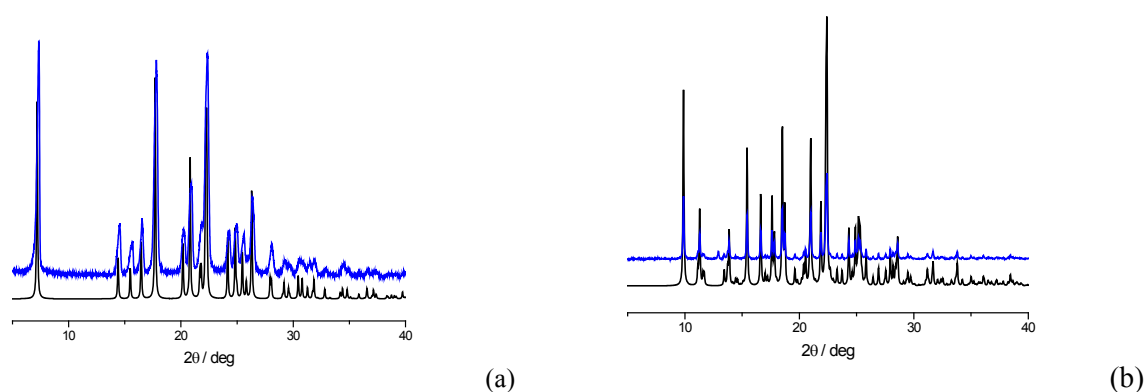


Figure ESI-0. Comparison between diffraction patterns for (a) PQ5U and (b) PQ8U, calculated on the basis of single crystal structure (black line), and experimental recorded at RT (blue line).

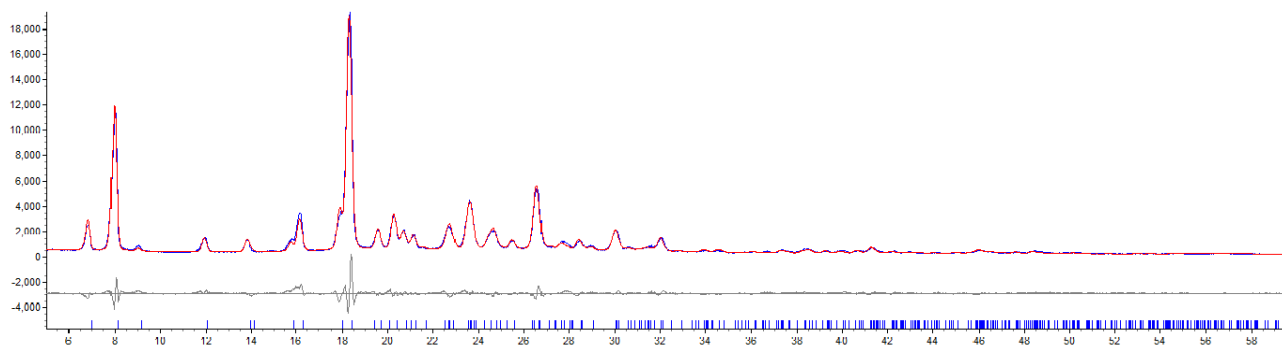


Figure ESI-5. Experimental (blue), calculated (red) and difference (grey) patterns for PQ6U; x axis is in degrees of 2θ .

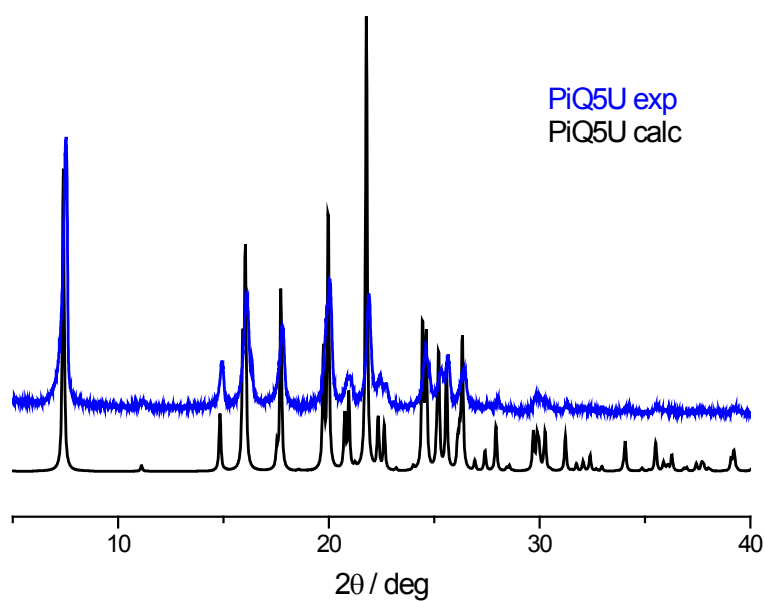


Figure ESI-6. Comparison between diffraction patterns for PiQ5U, calculated on the basis of single crystal structure (black line), and experimental recorded at RT (blue line).

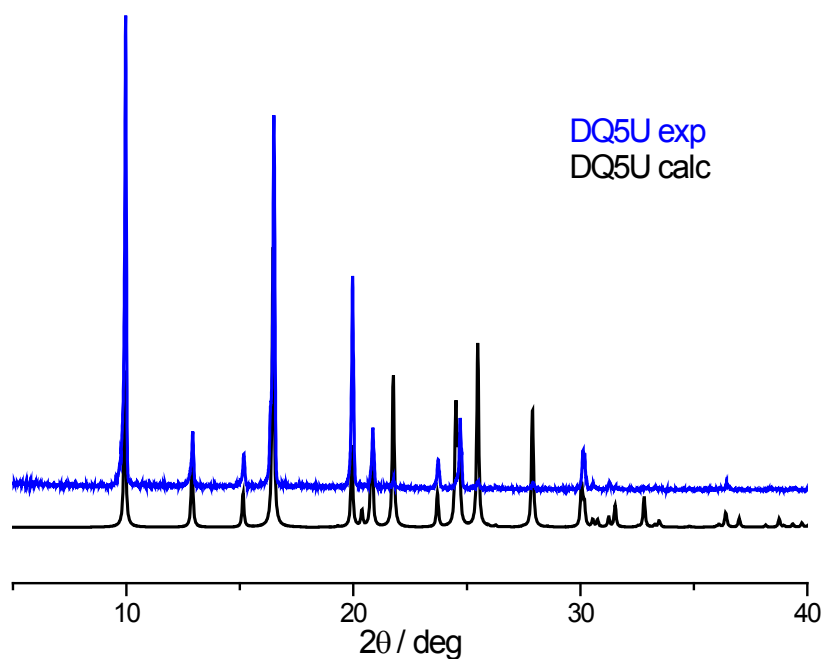


Figure ESI-7. Comparison between diffraction patterns for DQ5U, calculated on the basis of single crystal structure (black line), and experimental recorded at RT (blue line).

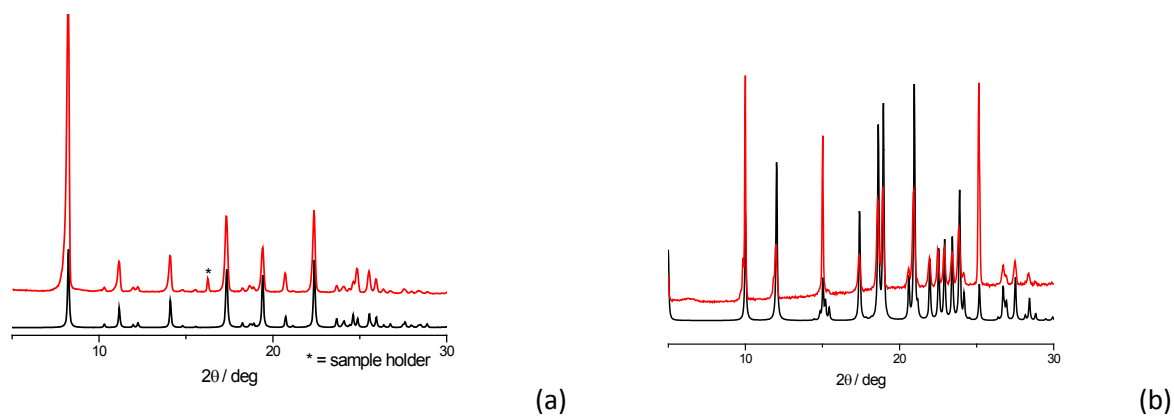


Figure ESI-8. Comparison between diffraction patterns for (a) PPT4U Form I and (b) PPT4U Form II, calculated on the basis of single crystal structure (black line) and experimental recorded at RT (blue line).

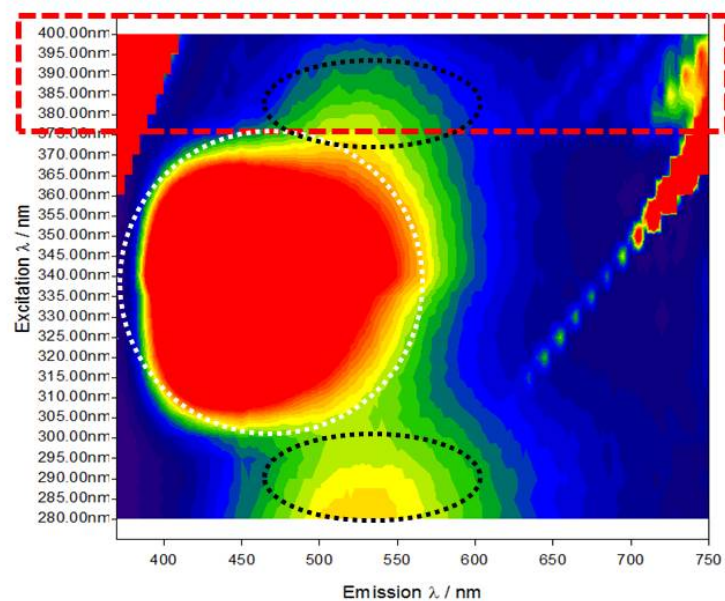


Figure ESI-9. Map of fluorescence spectra versus excitation wavelength for the AgNO_3 :DQ5U gel. White and black dashed circles evidence the emission bands of DQ5U and of the fibrillar aggregates respectively. The red dashed rectangle shows the spectral response of the gel when excited within our microscope setup. As it can be seen, the fibrils are selectively selected and visualized.

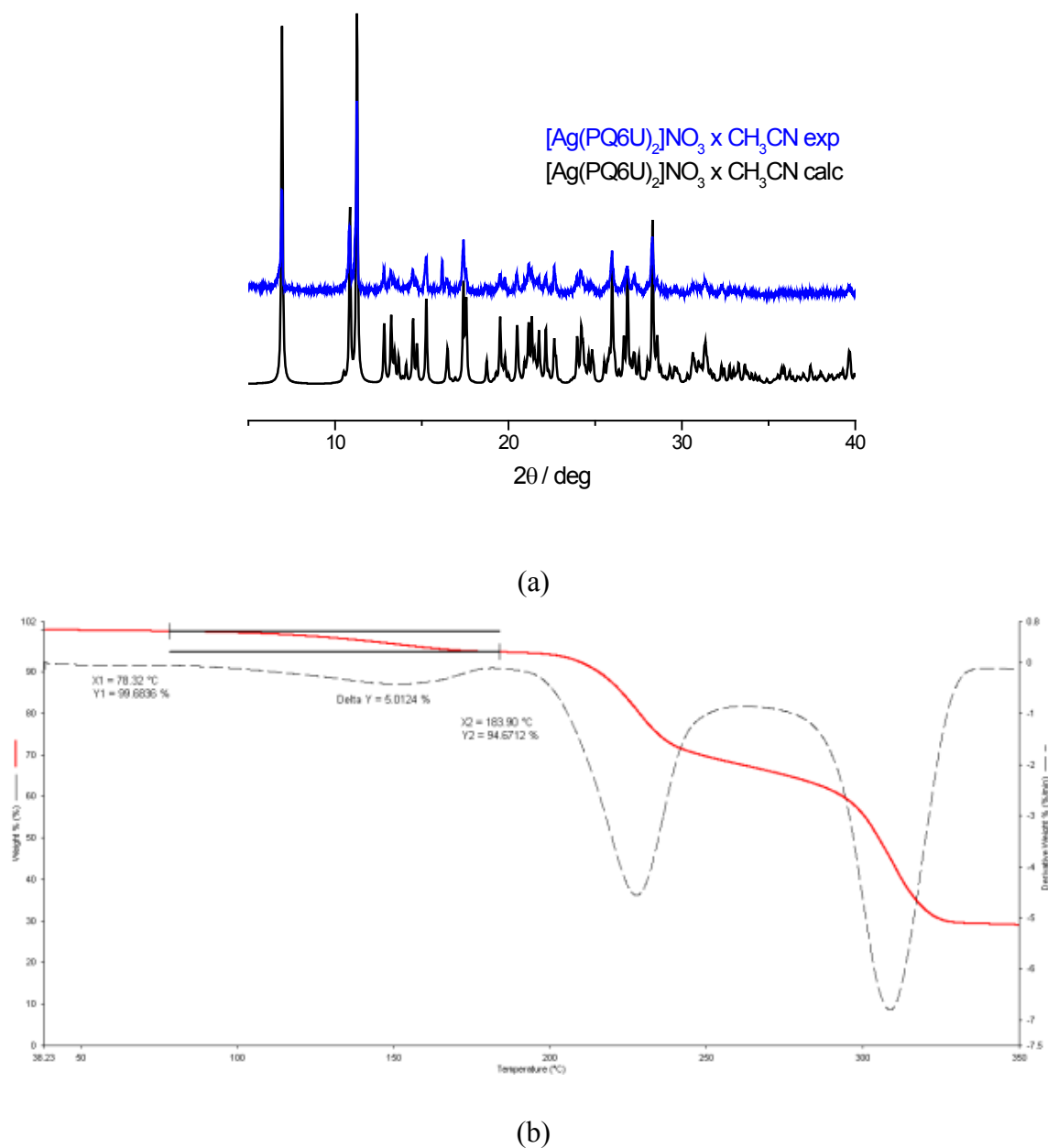
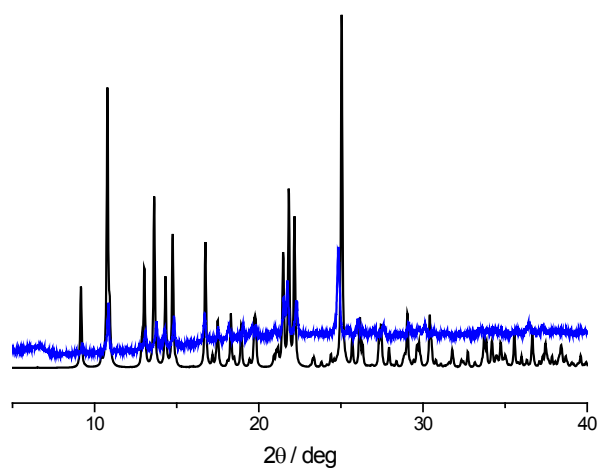
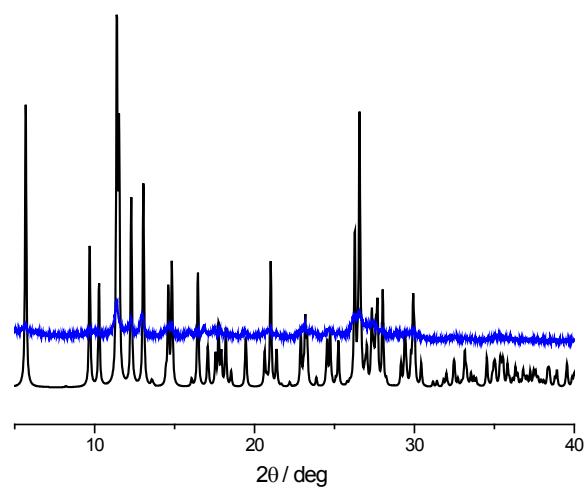


Figure ESI-10. Comparison between diffraction patterns for $[\text{Ag}(\text{PQ6U})_2]\text{NO}_3 \cdot \text{CH}_3\text{CN}$, calculated on the basis of single crystal structure (black line), and experimental recorded at RT (blue line); (b) the TGA trace for complex $[\text{Ag}(\text{PQ6U})_2]\text{NO}_3 \cdot \text{CH}_3\text{CN}$.



(a)



(b)

Figure ESI-11. Comparison between diffraction patterns for $[\text{Ag}(\text{PQ6U})_2]\text{NO}_3$ Form I (a) and Form II (b), calculated on the basis of single crystal structure (black line), and experimental recorded at RT (blue line).

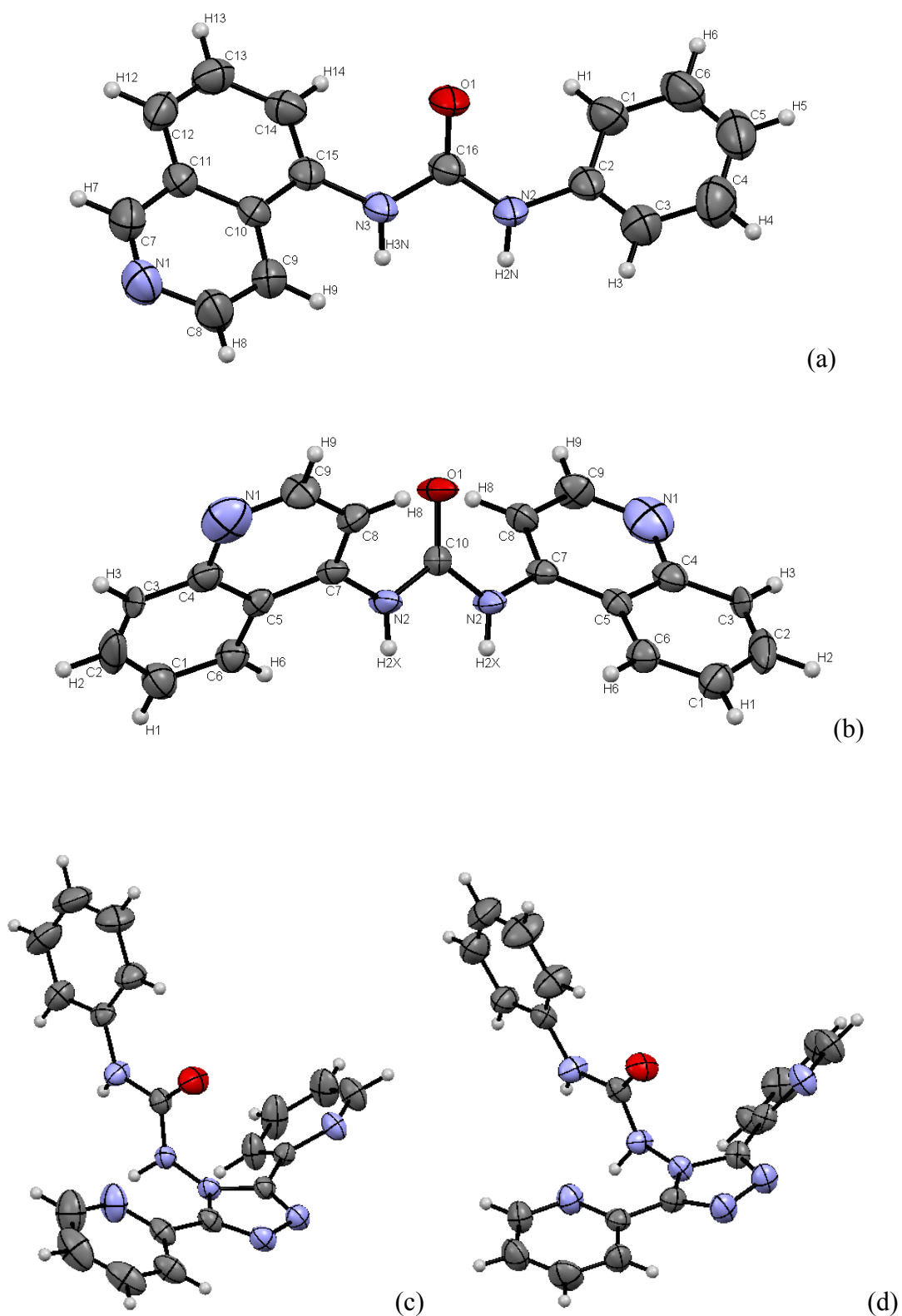


Figure ESI-12. ORTEP drawings for compounds (a) PiQ5U, (b) DQ5U, (c) PPT4U Form 1, and (d) PTT4U Form 2.

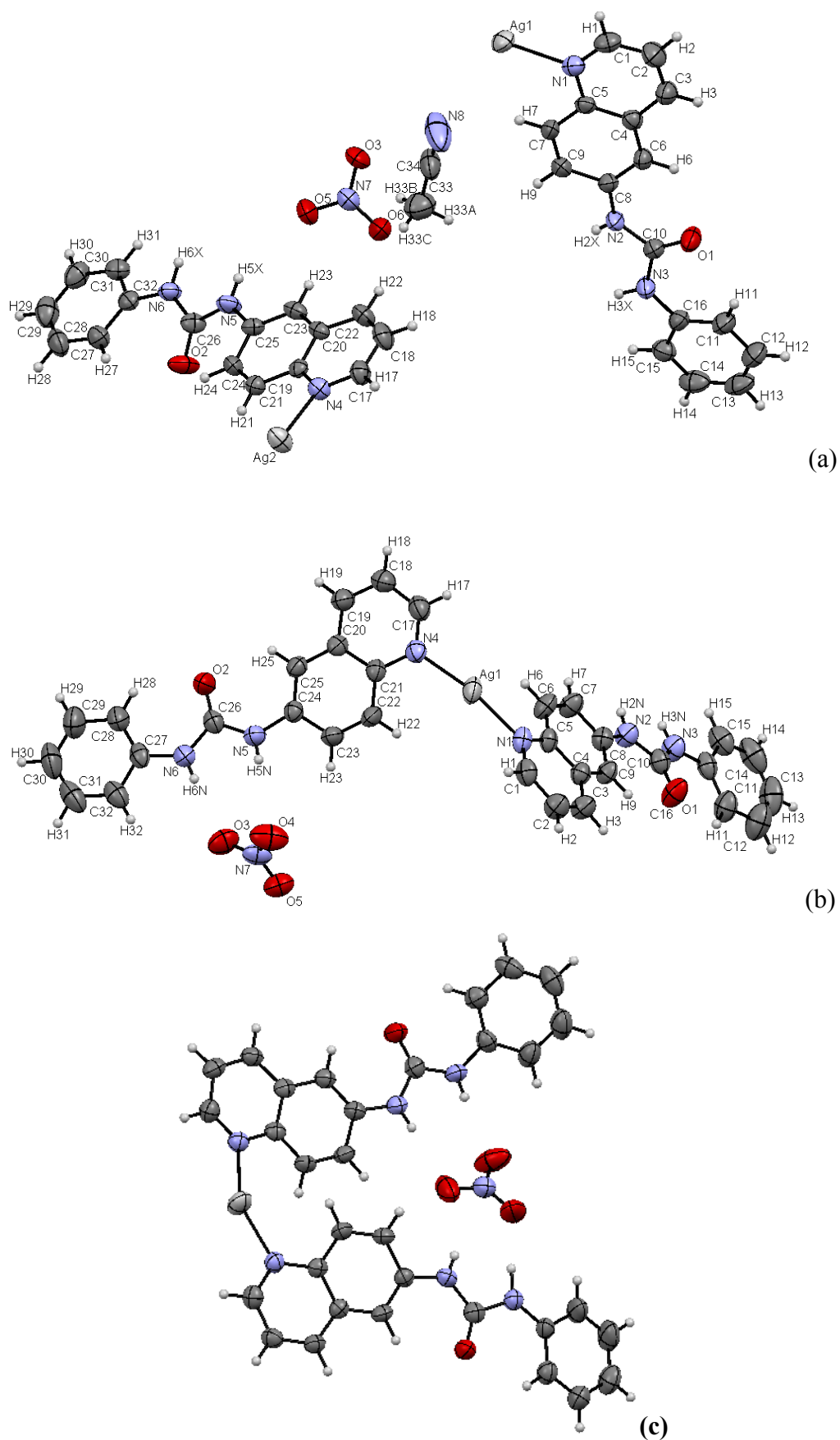


Figure ESI-13. ORTEP drawings for complexes: (a) $[\text{Ag}(\text{PQ6U})_2]\text{NO}_3 \cdot \text{CH}_3\text{CN}$, (b) $[\text{Ag}(\text{PQ6U})_2]\text{NO}_3$ Form 1 and (c) $[\text{Ag}(\text{PQ6U})_2]\text{NO}_3$ Form 2.