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Supporting information for article:

Synthesis, crystal structure and studies on the interaction with albumin of a new silver(I) complex based on 2-(4-nitrobenzene-sulfonamido)benzoic acid

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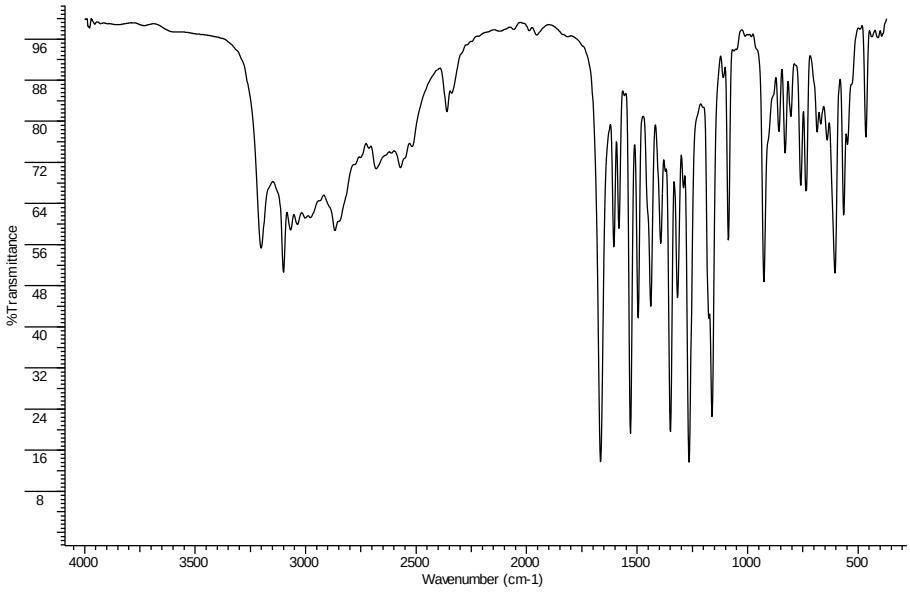
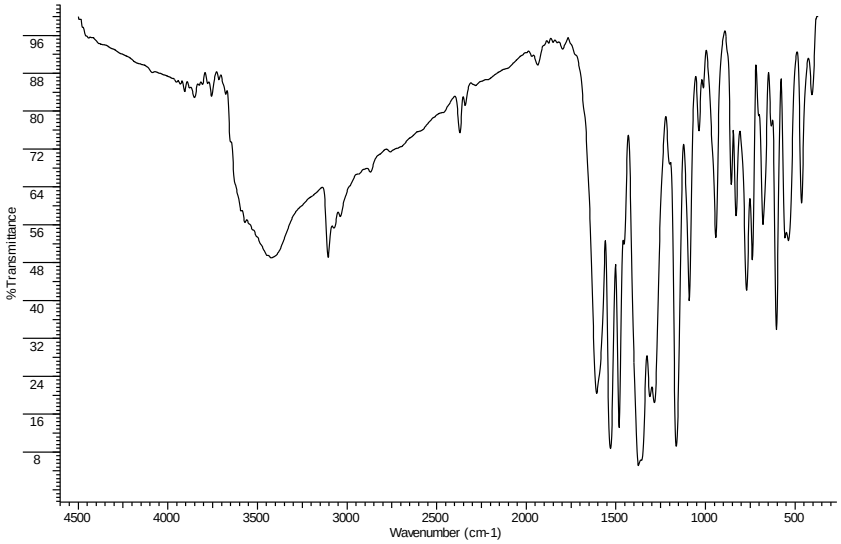
S1. Infrared spectra in the range of 4000 - 400 cm^{-1}.

HL

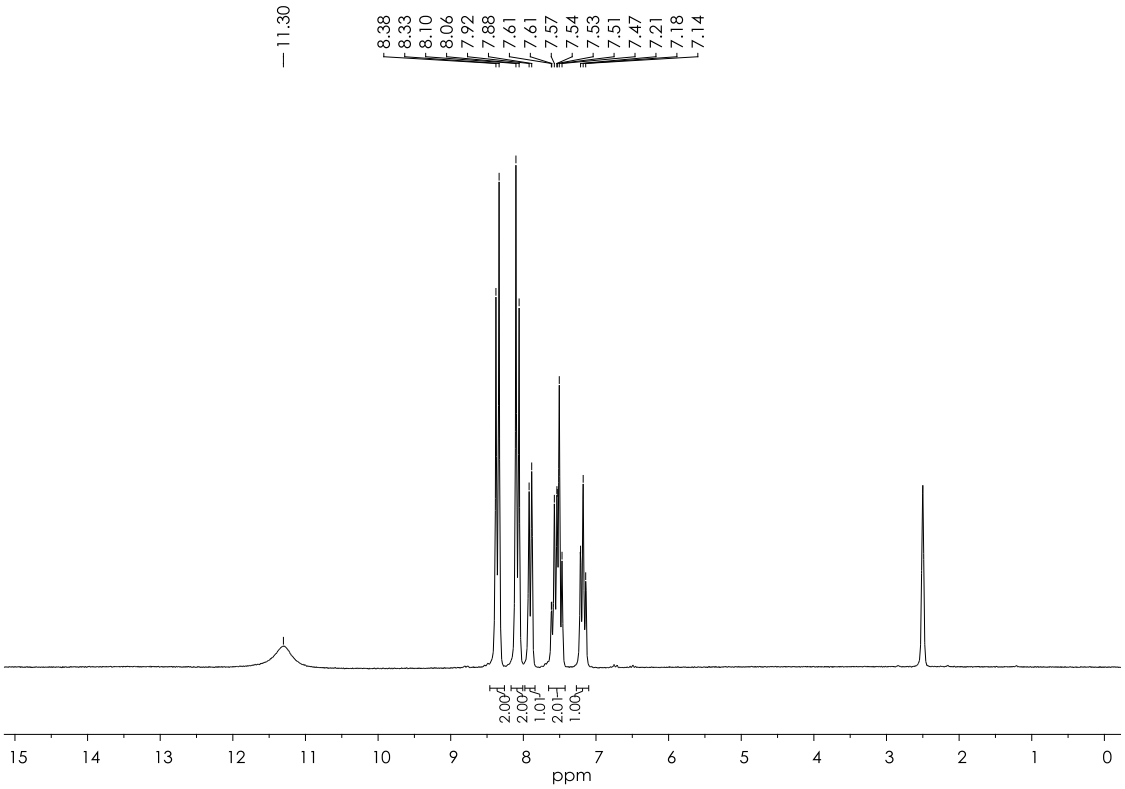


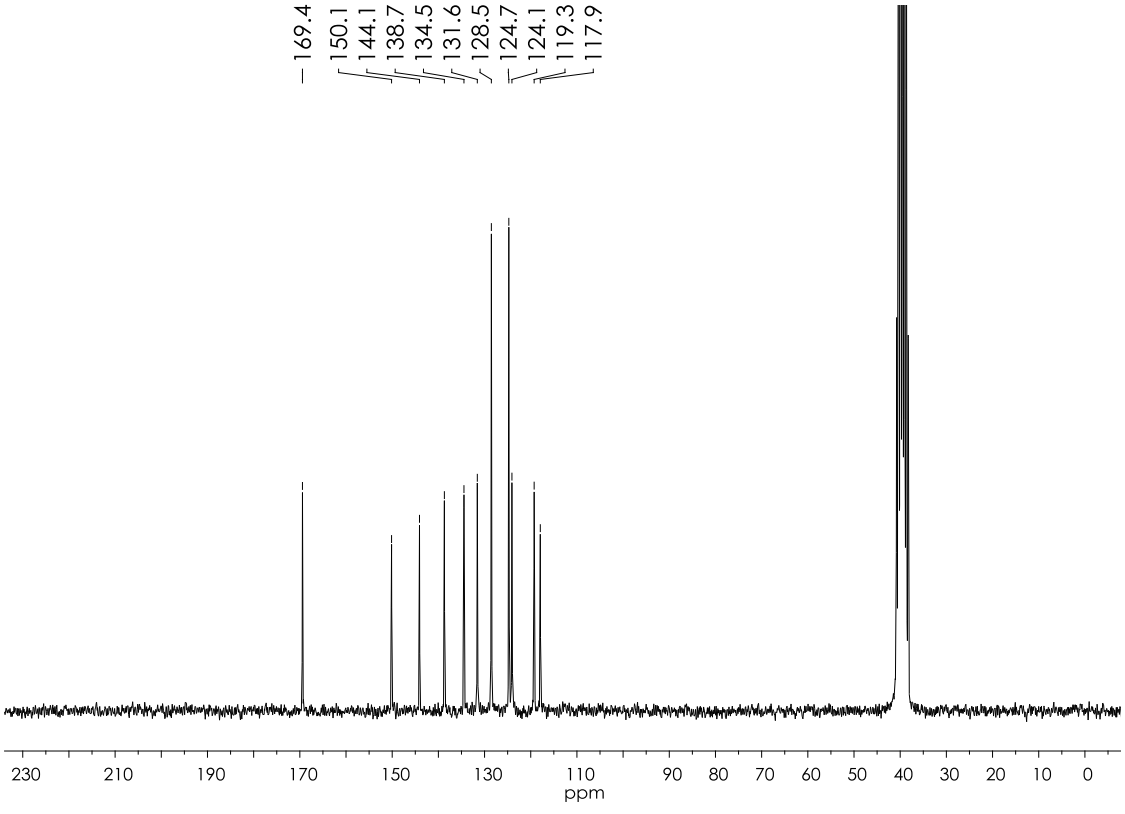
AgL



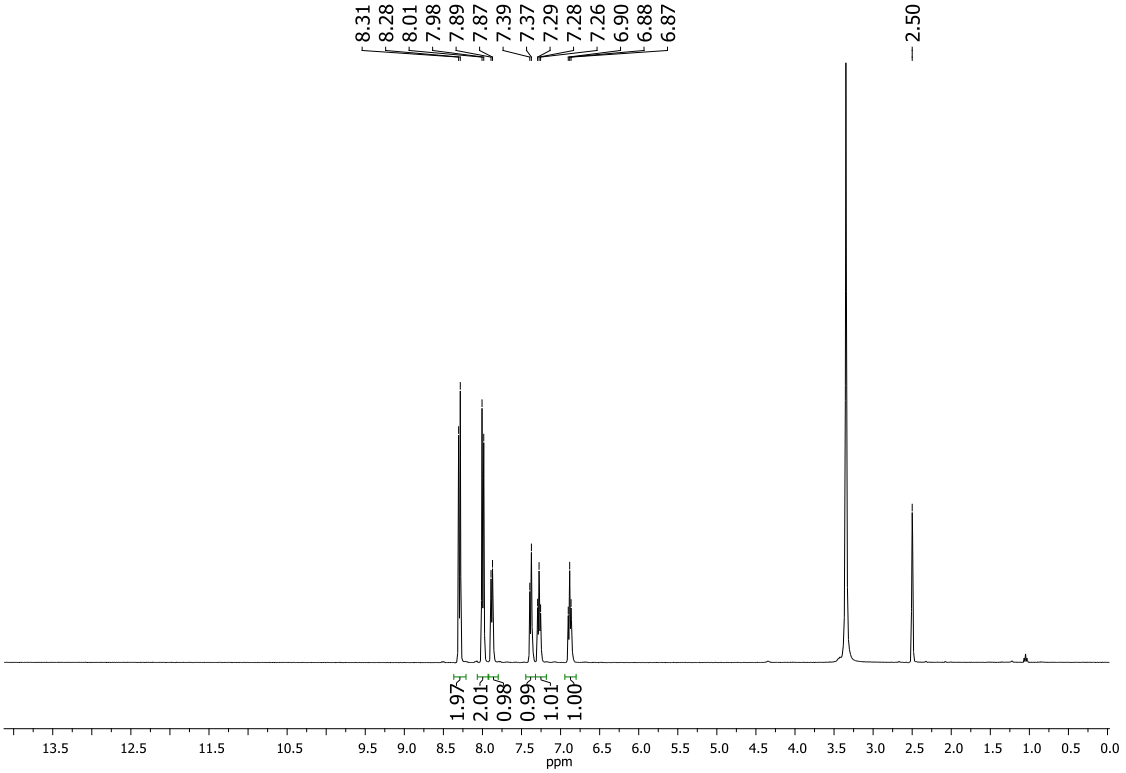
S2. NMR ^{1}H and $^{13}C\{1H\}$ spectra in 400 MHz (DMSO-d_6).

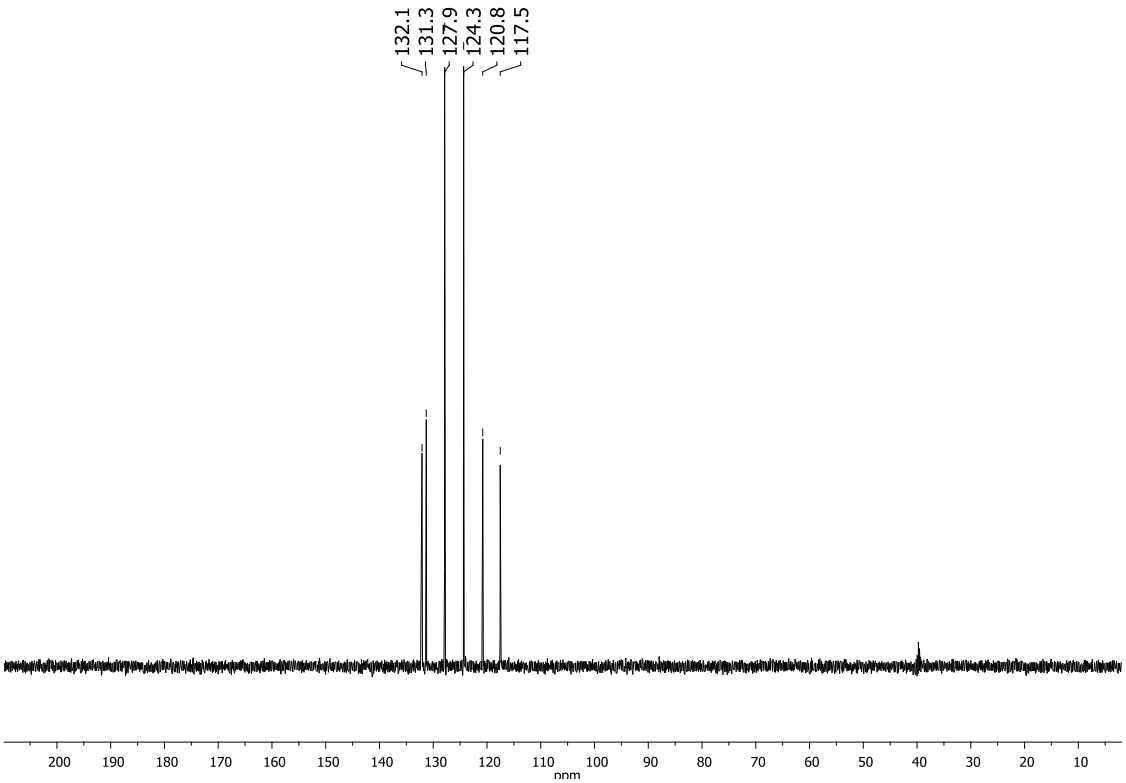
HL


11.30
8.38
8.33
8.10
8.06
7.92
7.88
7.61
7.61
7.57
7.54
7.53
7.51
7.47
7.21
7.18
7.14
2.00
2.00
1.01
2.01
1.00
15 14 13 12 11 10 9 8 7 6 5 4 3 2 1 0
ppm


169.4
150.1
144.1
138.7
134.5
131.6
128.5
124.7
124.1
119.3
117.9
230 210 190 170 150 130 110 90 80 70 60 50 40 30 20 10 0
ppm

AgL





S3. Hirshfeld surface analysis

Hirshfeld surfaces were obtained following F. L. Hirshfeld (Theor. Chim. Acta 44, 129 – 138 (1977)) using $w_i(\mathbf{r}) = 0.5$. d_{norm} is a function of d_i (distance of a point in the surface to the nearest atom inside the surface), d_e (distance of a point in the surface to the nearest atom outside the surface) and the van der Waals radii of these atoms:

$$d_{norm} = \frac{d_i - r_i^{vdW}}{r_i^{vdW}} + \frac{d_e - r_e^{vdW}}{r_e^{vdW}}$$

The CrystalExplorer program following the approach described McKinnon, Spackman and Mitchell (Acta Cryst., B60, 627 -668), was used for mapping the surfaces of the metal and of the ligand as well to draw the fingerprint plots of these surfaces.