

Low-dimensional compounds containing bioactive ligands. Part XIX: Crystal structures and biological properties of copper complexes with halogen and nitro derivatives of 8-hydroxyquinoline

Martina Kepeňová ¹, Martin Kello ², Romana Smolková ³, Michal Goga ⁴, Richard Frenák ⁴, Ľudmila Tkáčiková ⁵, Miroslava Litecká ⁶, Jan Šubrt ⁶ and Ivan Potočnýák ^{1,*}

¹ Institute of Chemistry, P. J. Šafárik University in Košice, Moyzesova 11, SK-04154 Košice, Slovakia

² Department of Pharmacology, P. J. Šafárik University in Košice, Trieda SNP 1, 040 11 Košice, Slovakia

³ Department of Ecology, Faculty of Humanities and Natural Sciences, University of Prešov, Ulica 17. novembra 1, Prešov, 081 16, Slovakia

⁴ Department of Botany, Institute of Biology and Ecology, Faculty of Science, P. J. Šafárik University in Košice, Mánesova 23, 040 01 Košice, Slovakia

⁵ Department of Chemistry, Biochemistry and Biophysics, University of Veterinary Medicine and Pharmacy, Komenského 73, 041 81 Košice, Slovakia

⁶ Centre of Instrumental Techniques, Institute of Inorganic Chemistry of the CAS, Husinec-Řež č.p. 1001, CZ-25068 Řež, Czech Republic

* Correspondence: ivan.potocnak@upjs.sk; Tel.: +421 55 234 2335; fax: +421 55 62 221 24

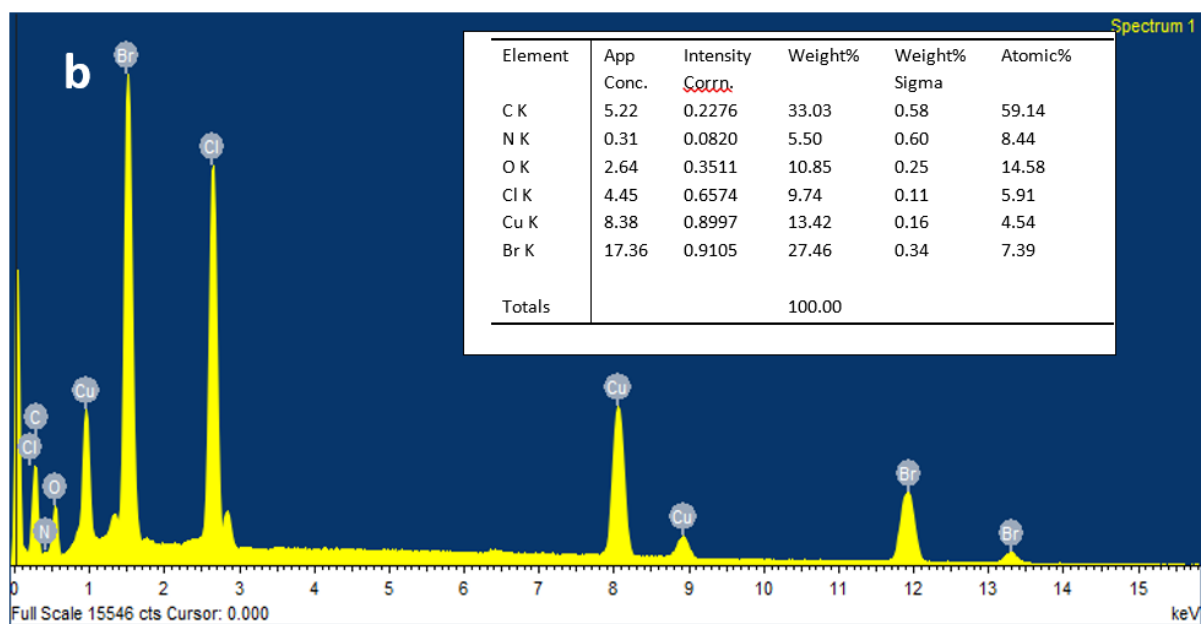
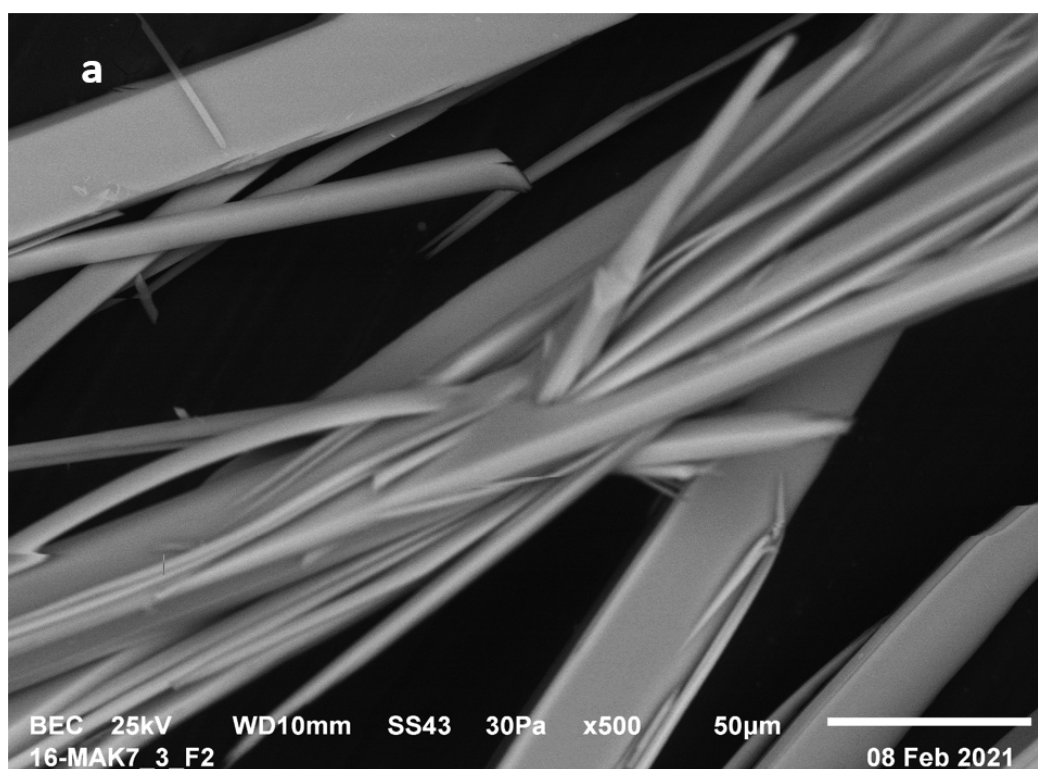


Figure S1. SEM micrograph (a) of **1a** along with EDS elemental analysis (b).

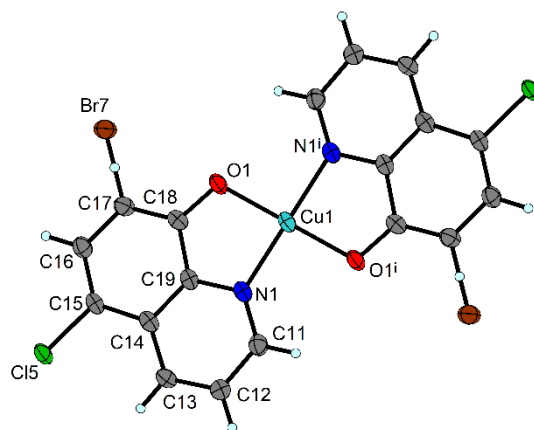


Figure S2. Molecular structure of **1a**. Displacement ellipsoids are drawn at the 50 % probability levels. Symmetry code: i = -x, -y+1, -z+1.

Table S1. Hydrogen bonds [\AA and $^\circ$] for **5** and **6**.

	D–H \cdots A	<i>d</i> (D–H)	<i>d</i> (H \cdots A)	<i>d</i> (D \cdots A)	\angle (DHA)
5	C12–H12 \cdots O2 ⁱⁱ	0.93	2.46	3.154(3)	132.0
	C3A–H3A3 \cdots O52 ⁱⁱⁱ	0.96	2.47	3.348(6)	152.6
6	C12–H12 \cdots O72 ⁱⁱ	0.95	2.58	3.456(3)	154.2

Symmetry codes:

5: ii = $x-1, y, z$; iii = $-x+2, -y+1, z$;

6: ii = $x-1, -y+3/2, z-1/2$.