Synthesis, structure and cytotoxicity of a series of Dioxidomolybdenum(VI) complexes featuring Salan ligands.

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Seven hexacoordinated cis-dioxidomolybdenum(VI) complexes [MoO₂L¹⁻⁷] (1-7) derived from various tetradentate diamino bis(phenolato) "salan" ligands, N,N'-dimethyl-N,N'-bis-(2hydroxy-3-X-5-Y-6-Z-benzyl)-1,2- diaminoethane $\{(X = Br, Y = Me, Z = H (H_2L^1); X = Me, L^2L^2\}$ Y|Cl, Z =H (H_2L^2); X= i Pr, Y= Cl, Z= Me (H_2L^3)} and N,N'-bis-(2-hydroxy-3-X-5-Y-6-Zbenzyl)-1,2-diaminopropane $\{(X = Y = {}^{t}Bu, Z = H (H_2L^4); X = Y = Me, Z = H (H_2L^5); X = {}^{i}Pr, \}$ Y|Cl, Z =Me (H_2L^6) ; X =Y =Br, Z =H (H_2L^7) } containing O-N donor atoms, have been isolated and structurally characterized. The formation of cis-dioxidomolybdenum(VI) complexes was confirmed by elemental analysis, IR, UV-vis and NMR spectroscopy, ESI-MS and cyclic voltammetry. X-ray crystallography showed the O₂N₂ donor set to define an octahedral geometry in each case. The complexes (1-7) were tested for their in vitro antiproliferative activity against HT-29 and HeLa cancer cell line. IC50 values of the complexes in HT-29 follow the order 6 < 7 < < 1 < 2 < 5 < < 3 < 4 while the order was 6 < 7 <5 < 1 < < 3 < 4 < 2 in HeLa cells. Some of the complexes proved to be as active as the clinical referred drugs, and the greater potency of 6 and 7 (IC₅₀ values of 6 are 2.62 and 10.74 µM and that of 7 is 11.79 and 30.48 µM in HT-29 and HeLa cells, respectively) may be dependent on the substituents in the salan ligand environment coordinated to the metal. More in Journal of Inorganic Biochemistry, 2017, 172, 110, https://doi.org/10.1016/j.jinorgbio.2017.04.015

