Dalton Transactions



View Article Online

CORRECTION

Check for updates

Cite this: Dalton Trans., 2024, 53, 3414

Correction: Antiproliferative activity of Ni(μ), Cu(μ), and Zn(μ) complexes of dithiocarbamate: synthesis, structural characterization, and thermal studies

Anupam Singh, 跑 ^a Kunal Shiv, 跑 ^a Ranjeet Singh, 跑 ^b M. K. Bharty, 跑 ^a Partha Pratim Manna 跑 *^b and Lal Bahadur Prasad ២ *^a

DOI: 10.1039/d4dt90018a

Correction for 'Antiproliferative activity of Ni(II), Cu(II), and Zn(II) complexes of dithiocarbamate: synthesis, structural characterization, and thermal studies' by Anupam Singh *et al.*, *Dalton Trans.*, 2024, **53**, 1196–1208, https://doi.org/10.1039/D3DT03724J.

The authors regret the errors in the assignment of the oxidation state and geometry of the iron and cobalt (dithiocarbamate) complexes (1 and 2). In the original paper, the iron and cobalt (dithiocarbamate) complexes were reported in +2 oxidation state with tetrahedral geometry. In fact, these complexes are air-sensitive and tend to oxidize readily into Fe(m) and Co(m) tris (dithiocarbamate) complexes with octahedral geometry.^{1,2}

In view of these inaccuracies, the $Fe(\pi)$ and $Co(\pi)$ dithiocarbamate complexes and their properties throughout the paper including the title (as reflected in the title shown here) should be omitted.

This correction will not affect the conclusions or the main finding of the manuscript (kindly refer to the slightly modified version of the conclusion as given below).

Conclusion: In summary, three new complexes have been prepared and characterized by various techniques. Further $Cu(\pi)$ and $Zn(\pi)$ complexes were characterized by single crystal-X-ray data. The structures of $Cu(\pi)$ and $Zn(\pi)$ complexes are stabilized by intermolecular hydrogen bonding. Single-crystal X-ray data reveals that the $Cu(\pi)$ complex has distorted square planar geometry whereas the $Zn(\pi)$ complex has distorted tetrahedral geometry around the central metal ions. In the $Zn(\pi)$ complex, the bonded dithiocarbamate moiety of the ligand uses the second sulfur to bridge the two $Zn(\pi)$ centers resulting in a binuclear dimeric $Zn(\pi)$ complex. The thermal studies of the metal complexes as analyzed by TG-DTA data, showed strong evidence for the formation of their respective metal sulfide at higher temperatures. The cytotoxic efficacy of the ligands and their metal complexes were examined on DL cells. Cytotoxicity assay results indicate that complexes have significant cytotoxic potential as compared to the free ligand. Complexes have significant antitumor activity against malignant lymphoma cells *in vitro* and *in vivo* in an animal model of lymphoma. In comparison to the standard drug cisplatin, the coordinate complexes are superior with reference to the IC₅₀ value, indicating better applicability against malignant lymphoma. The $Zn(\pi)$ complex was significantly tumoricidal both *in vitro* and *in vivo*. The cytotoxicity study suggests that the $Zn(\pi)$ complex was significantly tumoricidal both *in vitro* and *in vivo*. The cytotoxicity study suggests that the $Zn(\pi)$ complex was significantly to be used as an antitumor agent. These applications of complexes will give new insights into the further development of anticancer drugs based on transition metal ions.

The Royal Society of Chemistry apologises for these errors and any consequent inconvenience to authors and readers.

References

- 1 G. Hogarth, Prog. Inorg. Chem., 2005, 53, 71-561.
- 2 G. Hogarth, Mini-Rev. Med. Chem., 2012, 12, 1202-1215.

^aDepartment of Chemistry, Institute of Science, Banaras Hindu University, Varanasi 221005, India. E-mail: lbprasad@bhu.ac.in

^bImmunobiology Laboratory, Department of Zoology, Institute of Science, Banaras Hindu University, Varanasi 221005, India. E-mail: pp_manna@yahoo.com