



# Synthesis, Spectral Characterization and Crystal Structures of Dioxidomolybdenum(VI) Complexes Derived from Nicotinoylhydrazones

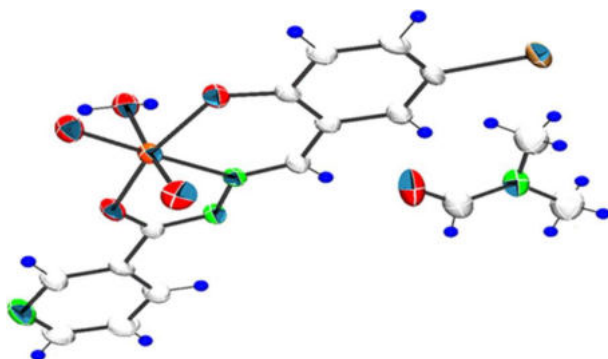
T. M. Asha<sup>1</sup> · M. R. P. Kurup<sup>1,2</sup>

Received: 12 July 2018 / Accepted: 29 November 2018  
© Springer Science+Business Media, LLC, part of Springer Nature 2018

## Abstract

*Cis*-dioxidomolybdenum(VI) complexes [MoO<sub>2</sub>L<sup>1</sup>(DMSO)] (**1**), [MoO<sub>2</sub>L<sup>2</sup>(H<sub>2</sub>O)]·DMF (**2**) and [MoO<sub>2</sub>L<sup>3</sup>(DMF)] (**3**) were synthesized by solution based reactions of aroylhydrazones (H<sub>2</sub>L<sup>1</sup> = 5-chloro-2-hydroxybenzaldehyde nicotinoylhydrazone, H<sub>2</sub>L<sup>2</sup> = 2-hydroxy-5-iodobenzaldehyde nicotinoylhydrazone and H<sub>2</sub>L<sup>3</sup> = 3,5-dichloro-2-hydroxybenzaldehyde nicotinoylhydrazone) with bis(acetylacetonato)dioxidomolybdenum(VI) complex, [MoO<sub>2</sub>(acac)<sub>2</sub>] in presence of DMSO or DMF. The synthesized aroylhydrazones and their molybdenum complexes were characterized by elemental analysis, spectroscopic techniques (FT-IR, UV-Vis, <sup>1</sup>H NMR) and conductivity measurements. Finally the three dimensional structures of the complexes were confirmed by single crystal X-ray diffraction studies. Crystal structures of **1**, **2** and **3** were solved by direct methods and refined with full-matrix least-squares calculations using the SHELXS97 and SHELXL2014 software programs respectively. Complex **1** got crystallized in monoclinic space group, *P*2<sub>1</sub>/*c* with *Z* = 4, whereas complexes **2** and **3** in triclinic space group, *P* $\bar{1}$  with *Z* = 2. The ligands, H<sub>2</sub>L<sup>1-3</sup> showed coordination to the metal ion in a dibasic tridentate manner through deprotonated phenolate oxygen, azomethine nitrogen and enolate oxygen.

## Graphical Abstract



All the three dioxidomolybdenum(VI) complexes exhibit a distorted octahedral geometry around molybdenum atom. ORTEP plot of [MoO<sub>2</sub>L<sup>2</sup>(H<sub>2</sub>O)]·DMF. (Displacement ellipsoids are drawn at 50% probability)

**Keywords** Aroylhydrazones · Molybdenum complexes · X-ray diffraction studies · Crystal structure · Azomethinenitrogen

## Introduction

Aroylhydrazones are significant class of ONO donor ligands which have played an important role in the development of coordination chemistry of molybdenum. Furthermore the study of oxidomolybdenum complexes coordinated with

✉ M. R. P. Kurup  
mrpcusat@gmail.com; mrp@cukerala.ac.in

Extended author information available on the last page of the article