

Library Indian Institute of Science Education and Research Mohali



DSpace@IISERMohali (/jspui/)

- / Publications of IISER Mohali (/jspui/handle/123456789/4)
- / Research Articles (/jspui/handle/123456789/9)

Please use this identifier to cite or link to this item: http://hdl.handle.net/123456789/3025

 $\label{thm:continuous} \mbox{Title:} \qquad \mbox{Synthesis, structural characterization, and solution properties of a 1-D Pb(II)-bipyridine}$

coordination polymer

Authors: Dey, D. (/jspui/browse?type=author&value=Dey%2C+D.)

Yadav, H.R. (/jspui/browse?type=author&value=Yadav%2C+H.R.)

 $Choudhury, A.R. \ (/jspui/browse?type=author\&value=Choudhury\%2C+A.R.)$

Keywords: 1-D polymer

Pb(II)

Spectroscopic investigation Supramolecular interactions

Issue Date:

2015

Publisher: Taylor and Francis Ltd.

Citation: Journal of Coordination Chemistry, 68(1) pp. 169-180

 $Abstract: \qquad A \text{ new one-dimensional (1-D) Pb(II) coordination polymer, } [Pb(2,2'-bpy)(NO3)2(H2O)]n \text{ (1) } (2,2'-by)(NO3)2(H2O)]n \text{ (2) } (2,2'-by)(NO3)2(H2O)[n] n \text{ (2) } (2,2'-by)(NO3)2$

bpy = 2,2'-bipyridine), has been synthesized and characterized by different spectroscopic techniques and X-ray single-crystal analysis. From the X-ray crystal structure of 1, the Pb2+ can be best described as a highly distorted pentagonal bipyramid with O4 (water) and O6 (nitrate) at apical positions (O4-Pb-O6 of 143.7(1)°). Variability in bond distances reveals that Pb2+ is unsymmetrically surrounded by two nitrates, one 2,2'-bpy and one water. Nitrates bridge between monomers. The molecule crystallizes in the monoclinic P21/n (14) space group. This is the first example of a 1-D Pb(II) polymer in which nitrates show three different coordination motifs (terminal, chelating, and bridging). Solid state as well as solution phase UV-vis spectral analysis and mass spectrometric studies clearly reveal instability with breakdown of Pb(II) polymer in aqueous solution. The arrangement of the 2,2'-bpy, water, and nitrates leaves a coordination gap at the

Pb(II) occupied probably by a stereo-active lone pair of electrons.

URI: https://www.tandfonline.com/doi/full/10.1080/00958972.2014.985215

(https://www.tandfonline.com/doi/full/10.1080/00958972.2014.985215) http://hdl.handle.net/123456789/3025 (http://hdl.handle.net/123456789/3025)

Appears in Research Articles (/jspui/handle/123456789/9)

Collections:

Files in This Item:

File	Description	Size	Format	
Need to add pdf.odt (/jspui/bitstream/123456789/3025/1/Need%20to%20add%20pdf.odt)		8.63 kB	OpenDocument Text	View/Open (/jspui/bitstream/12345

Show full item record (/jspui/handle/123456789/3025?mode=full)

(/jspui/handle/123456789/3025/statistics)

Items in DSpace are protected by copyright, with all rights reserved, unless otherwise indicated.