

A new chain-like heteropolytungstate formed by Keggin cluster units: Synthesis and structure of $[\text{H}_2\text{bpy}]_3[\text{SiMnW}_{11}\text{O}_{39}]\cdot 1.25\text{H}_2\text{O}$

Jing Xin Meng, Ying Ma, Xin Xin Xu, Yang Guang Li, En Bo Wang*

Key Laboratory of Polyoxometalate Science of Ministry of Education Department of Chemistry,
Northeast Normal University Ren Min Street No.5268, Changchun, Jilin, 130024, PR China

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Abstract

A new heteropolytungstate, $[\text{H}_2\text{bpy}]_3[\text{SiMnW}_{11}\text{O}_{39}]\cdot 1.25\text{H}_2\text{O}$ (**1**), has been prepared under mild hydrothermal conditions and structurally characterized by single crystal X-ray diffraction. It is the first characterized compound containing 1D zigzag chain of transition metal substituted Keggin heteropolytungstate, which is connected through a common oxygen atom. The elemental analysis and IR of it are also described.

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Keywords: Heteropolytungstate; Keggin clusters; Zigzag chain

As a rich and diverse class of organic–inorganic hybrid materials, transition metal substituted polyoxotungstates based on Keggin framework have received considerable attention for their fascinating structural, electrochemical, catalytic, magnetic and photophysical properties [1–4]. Although these transition metal substituted Keggin derivatives are generally reported as discrete entities, several literatures have shown that they are capable of acting as inorganic building blocks to form one-dimensional chain like self-assembly compounds, such as compounds $(\text{ET})_8[\text{PMnW}_{11}\text{O}_{39}]\cdot 2\text{H}_2\text{O}$ (ET = bis (ethylenedithio) tetrathiofulvalene) [5], $[\text{NEt}_3\text{H}][\text{XCoW}_{11}\text{O}_{39}]\cdot 3\text{H}_2\text{O}$ (X = P, As) [6] and $[\text{Co}(\text{dpa})_2(\text{OH}_2)_2]_2[\text{Hdpa}][\text{PCoW}_{11}\text{O}_{39}]$ (dpa = di-2-pyridylamine) [7] they all possess straight chain. In addition, Er^{III} mono-substituted $[\alpha\text{-SiW}_{11}\text{O}_{39}]^{8-}$ polyoxotungstate containing one-dimensional zigzag chainlike structure, which was connected by two common oxygen atoms, has been synthesized by Niu et al. [8].

On the basis of our previous works [9], here we report the synthesis and structure of the first characterized compound containing 1D zigzag chains of transition metal substituted Keggin heteropolytungstate **1**, which is connected through a common oxygen atom. It shows a new zigzag chain structure, which is different to the straight chain structure in the reported 1D transition metal substituted Keggin heteropolytungstate [5–7]. Furthermore, the zigzag chains of **1** form layers through short inter-species contact and these layers are stacked parallel and form a three-dimensional structure with dumbbell-like 1D channel.

* Corresponding author.

E-mail address: wangenbo@public.cc.jl.cn (E.B. Wang).

1. Experimental

All organic solvents and materials used for synthesis were reagent grade and used without further purification. $K_8[\beta_2\text{-SiW}_{11}\text{O}_{39}]\cdot 14\text{H}_2\text{O}$ was prepared according to the literature method [10]. The identity was confirmed by IR spectrum.

Synthesis of the title compound: a mixture of $\text{MnCl}_2\cdot 4\text{H}_2\text{O}$ (0.25 g), 4,4'-bpy (0.032 g), $K_8[\beta_2\text{-SiW}_{11}\text{O}_{39}]\cdot 14\text{H}_2\text{O}$ (0.15 g) and H_2O (4 mL) was adjusted to pH (5–6) by adding HCl (1 mol/L), then the mixture was stirred for 20 min in air. The mixture was then transferred to a teflon-lined autoclave (20 mL) and kept at 130 °C for 5 days. After slow cooling to room temperature, red crystals were filtered, washed with distilled water and dried in a desiccator at room temperature to give a yield of 39% based on W. The elemental analysis (%) found: C, 11.20; H, 1.10; N, 2.55, calcd: C, 11.17; H, 1.02; N, 2.60. The ICP analysis (%) showed that **1** contained W, 62.50; Mn, 1.82; Si, 0.88, calcd: W, 62.68; Mn, 1.70; Si, 0.87. IR spectra: 3430, 1592, 1521, 1492, 1476, 1414, 1393, 1360, 1249, 1232, 1206, 1101, 1064, 1007, 961 and 914 cm^{-1} .

2. X-ray crystallography

A red single crystal of **1** was carefully selected under a polarizing microscope and glued at the tip of a thin glass fiber with cyanoacrylate (super glue) adhesive. Single crystal structure determination by X-ray diffraction was performed on a *R*-axis RAPID IP diffractometer equipped with a normal focus, 18 kW sealed tube X-ray source (Mo $K\alpha$ radiation, $\lambda = 0.71073 \text{ \AA}$) operating at 50 kV and 200 mA. A total of 40,638 reflections were collected with 9320

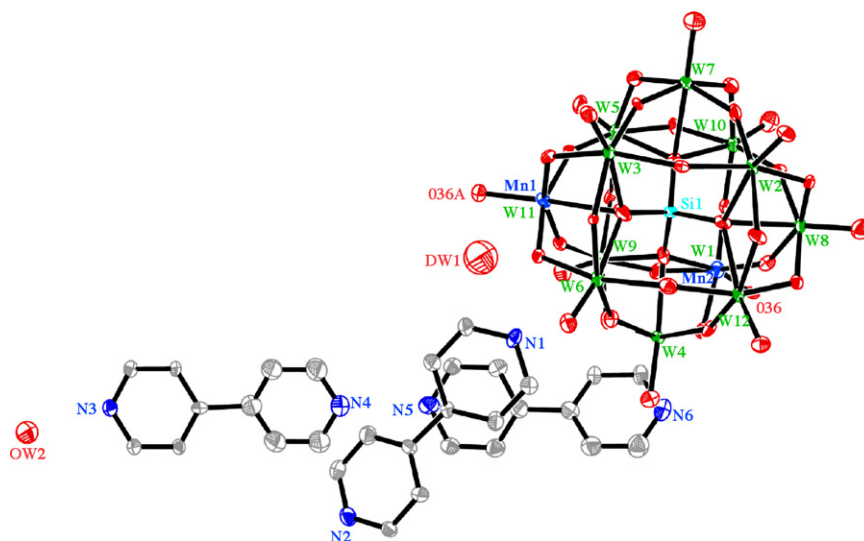


Fig. 1. ORTEP drawing of **1** showing the labeling of atoms with thermal ellipsoids at 50% probability. Hydrogen atoms are omitted for clarity.

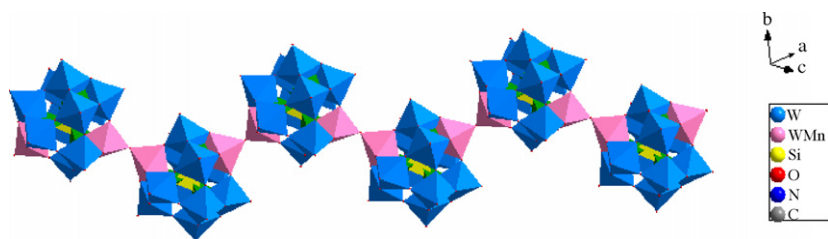


Fig. 2. A polyhedral view of the 1D chain in **1**. The pink octahedra are CuO_6 , blue octahedra are WO_6 and yellow tetrahedra are SiO_4 . (For interpretation of the references to color in this figure legend, the reader is referred to the web version of the article.)

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