

Poster Presentations

[MS10-P04] Crystal structure of $\text{Na}_x\text{P}_2\text{W}_{10}\text{O}_{34}$, the $m=10$ member of the series of $\text{Na}_x(\text{PO})_4(\text{WO})_3$.

Amar Benmoussa, Wahiba Bouchelaghem,
Mohamed-Rida Benloucif.

*Laboratoire de Chimie des Matériaux
Inorganiques. Université Badji Mokhtar
Annaba, Algeria.*

E-mail: amar.benmoussa@univ-annaba.org

The phosphates of transition metals M ($M=W, \text{Mo}, \text{Ti}, \text{V}, \dots$) form a huge family of compounds. They are characterized by a framework formed by MO_6 octahedra and $\text{M}'\text{O}_4$ tetrahedra ($\text{M}'=\text{P}, \text{Si}$). A new member, $m=10$ of the series $\text{Na}_x(\text{PO})_4(\text{WO})_3$ has been isolated and studied by single crystal X-ray diffraction. The structure corresponds to that of the Monophosphates Tungsten Bronzes with hexagonal tunnels (MPTBh's) [1-5]. It has been solved and refined to conventional $R=0.0466$, with 665 independent reflections with $I>4\sigma(I)$. The unit cell is monoclinic (space group $P2_1/m$) with $a = 6.651(7)\text{Å}$, $b = 5.293(5)\text{Å}$, $c = 18.125(4)\text{Å}$, $\beta = 94.54(2)^\circ$. The structure is described and compared with those of other members of the series including Na- and K-based analogues.

Références :

- [1]- A. Benmoussa, D. Groult, Ph. Labbé, and B. Raveau, (1984). *Acta Crystallogr. Sect. C*, 40, 573. [2]- J. P. Giroult, M. Goreaud, Ph. Labbé, and B. Raveau, (1982). *J. Solid State Chem.*, 44, 407.
[3]- M. Lamire, M. Goreaud, Ph. Labbé, and B. Raveau, (1987). *J. Solid State Chem.*, 66, 64.
[4]- B. Domengès, M. Goreaud, Ph. Labbé, and B. Raveau, (1983) *J. Solid State Chem.*, 50, 173.
[5]- P. Roussel, A. C. Masset, B. Domengès, A. Maignan, D. Groult and Ph. Labbé, (1998). *J. Solid State Chem.*, 139, 362.

Keywords : Monophosphates, Bronzes, X-Ray Diffraction.