

Poster Presentation

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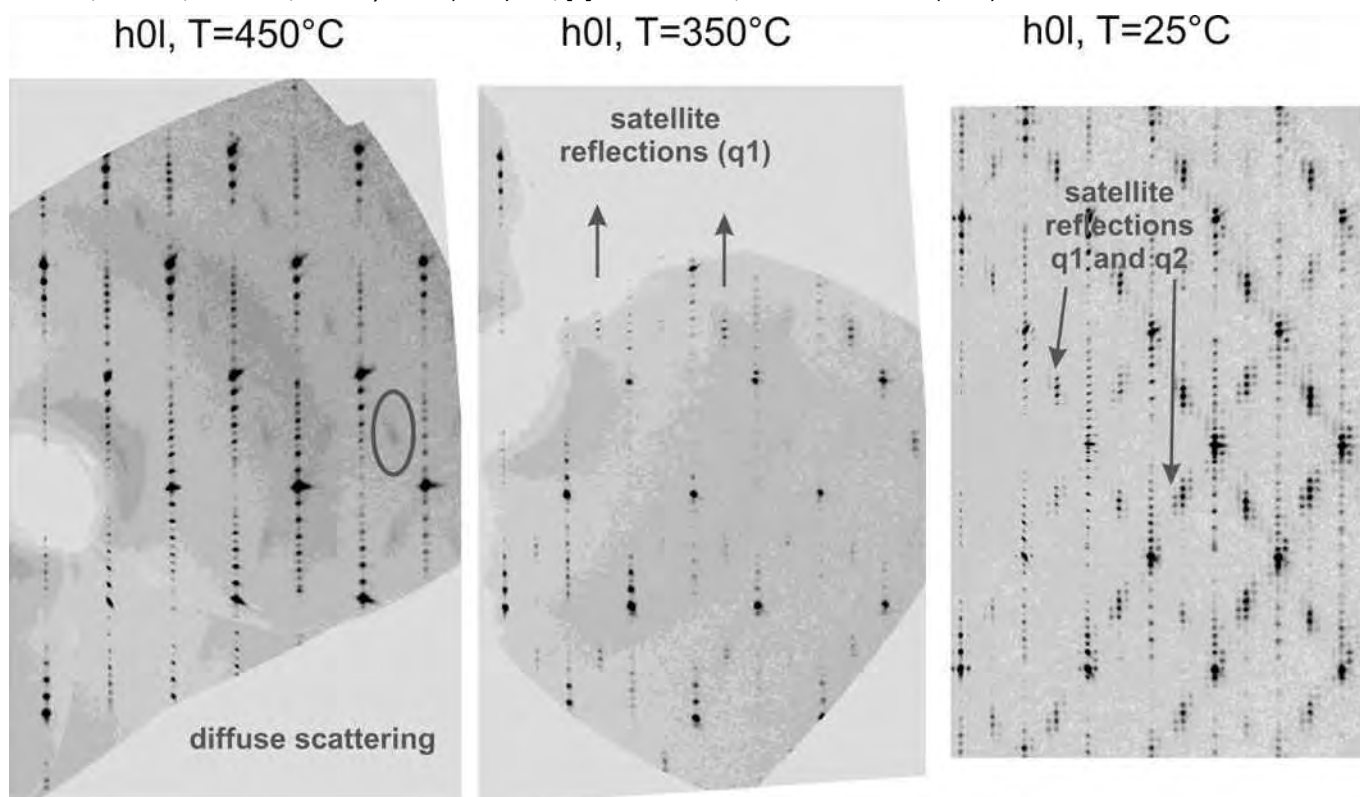
Phase transitions toward charge density wave in the $m=11.5$ Mono phosphate bronze

O. Pérez¹, K. Kolincio¹, A. Pautrat¹

¹CRISMAT, Ensicaen / CNRS, Caen, France

Mono Phosphate Tungsten Bronzes (MPTB), $(\text{PO}_2)_4(\text{WO}_3)_{2m}$ (m ranging from 4 to 14), are a large family of conductors of low dimensionality[1]. Their structure may be described as a regular stacking of WO_3 -type slabs with a thickness function of m , joined by slices of tetrahedral PO_4 phosphate. Successive Peierls transitions towards charge-density wave (CDW) or spin-density wave (SDW) states are observed below different critical temperatures (T_{C1} , T_{C2} ...). Structural transitions toward incommensurate modulated phases are associated with these complex electronic states. Recently an original member of MPTB, the member $m=11.5$ with formula $(\text{PO}_2)_4(\text{WO}_3)_{11}(\text{WO}_3)_{12}$, resulting in a regular intergrowth between $m=11$ and $m=12$ member, has been synthesized. An accurate investigation of the reciprocal space versus the temperature using single crystal X-ray diffraction shows the existence of two phase transitions at $T_{C1}=430^\circ\text{C}$ and $T_{C2}=280^\circ\text{C}$. The ground state structure may be described with the following cell parameters $a=5.3431(7)$ $b=6.5901(9)$ $c=40.884(6)$ $\alpha=93.096(2)$ $\beta=93.734(2)$ $\gamma=90.000(2)$ and the SG P-1. Both transitions associated with the occurrence of CDW are characterized by the appearance of modulation vectors (q_1 and q_2 associated to T_{C1} and T_{C2} respectively) and an increasing of the dimension of the superspace group (see figure 1). A structural study of the $m=11.5$ member is performed above T_{C1} and below T_{C1} and T_{C2} ; an interpretation of the CDW state is then proposed. The evolution of the intensity of the satellite and main reflections around T_{C1} and T_{C2} is analyzed for characterizing the order parameter of the transitions. Finally, resistivity and magnetoresistance measurements of a large single crystal sample of $m=11.5$ MPTB are performed in a temperature interval from 2K to 300K. These measures are revealing at $\sim 50\text{K}$ a possible SDW transition due to electron-electron interactions playing significant role in that material[2]. Further neutron diffraction experiment at very low temperature shall be performed to clarify this point.

[1] P. Roussel, O. Pérez, Ph. Labbé, *Acta Cryst B* 57 (2000) 603, [2] M. Greenblatt, *Acc. Chem. Res.* 29 (1996) 219



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