	<b>Experiment title:</b> Charge and structural order parameters and correlated polarons at the metal-insulator transition in NdNiO <sub>3</sub> epitaxial films.	<b>Experiment number:</b> HE-1391
<b>Beamline:</b> ID 20	<b>Date of experiment:</b> from: 04/09/2002                      to: 10/09/2002	<b>Date of report:</b> 20/02/2004
<b>Shifts:</b> 18	<b>Local contact(s):</b> Luigi Paolasini	<i>Received at ESRF:</i>
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Electronic charge localization in 3d transition-metal oxides has attracted considerable interest in the past. These localization phenomena, which lead to metal-insulator transitions, play a crucial role in such materials as high- $T_c$  cuprates and giant magneto-resistive manganites. Recently we have found a charge disproportionation phenomena in NdNiO<sub>3</sub>, which occurs below the metal-insulator [1]. The simultaneous presence of orbital-ordering with this phenomena is under debate.

In order to investigate the presence of orbital-ordering in the insulating phase, we have collected data on the azimuthal behavior (rotation about the wave vector) of particular reflections on a NdNiO<sub>3</sub> [011] thin film.

On and off resonance spectra were collected for the high and low temperature phases in both  $Q_{\parallel}$  and  $Q_{\perp}$  polarization channels. The experimental behavior has been compared with the theoretical one, calculated with the use of the tensor  $\langle T_Q^K \rangle$  which takes into account the contribution of the anisotropy of the atomic scattering factor close to the absorption edges.  $K$  is the rank of the tensor ( $K=0$  is for charge,  $K=1$  is for magnetic and  $K=2$  is for orbital contributions). In the case of NdNiO<sub>3</sub> ( $Pbnm$  space group for the high temperature phase), for forbidden reflection like  $(0 \ 2n+1 \ 1)$  i.e.  $(0 \ 3 \ 3)$ , only the tensor of rank 2 is contributing with  $Q=0, \pm 1$  (i.e.  $T_1^2$ ).

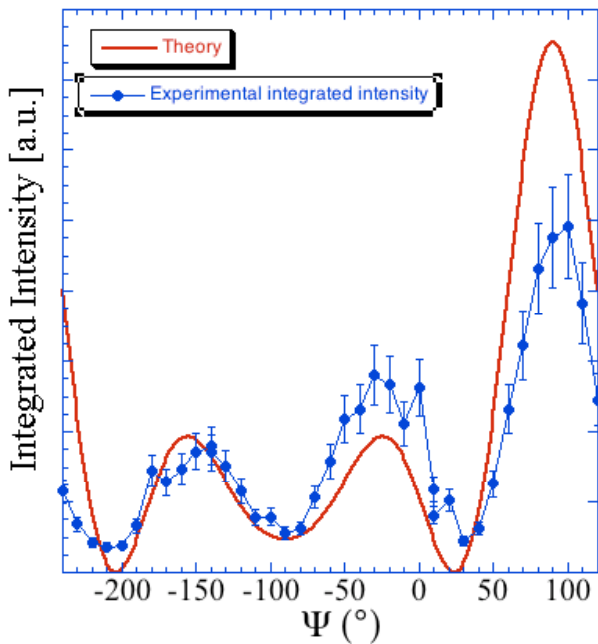
A more detailed explanation can be found in reference [2, 3].

No azimuthal dependence was founded, as expected, in the  $\pi\pi\pi$  channel.

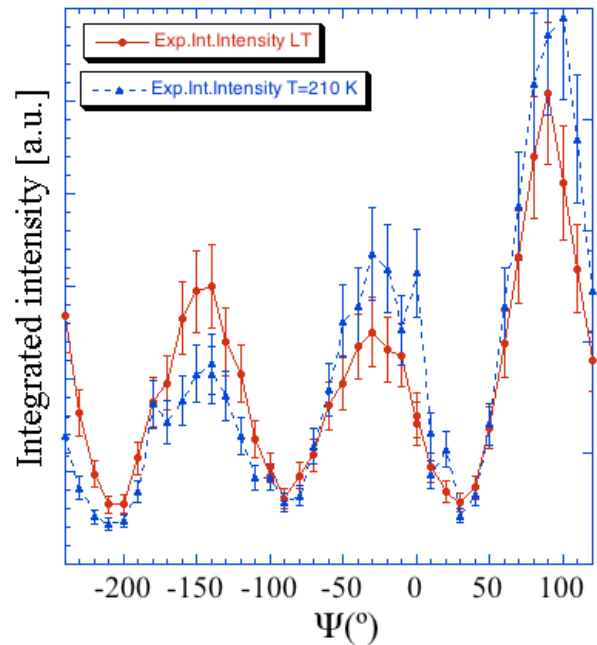
In the case of the  $\pi\pi\pi$  channel there is good agreement between theory and measured integrated intensity (see fig.1).

A similar azimuthal behavior was found in the low temperature phase ( $T < T_{MI}$ ), see fig.2.

Those results suggest that the orbital-ordering is not present or doesn't play any role in determining the charge ordering at metal-insulator transition.



**Fig.1** Experimental behavior of the (033) reflection, in the  $\pi\pi\pi$  channel as a function of the azimuthal angle  $\Psi$  (dotted line) compared with theoretical behavior (continuous line).



**Fig.2** Azimuthal angle dependence of the (033) reflection in the  $\pi\pi\pi$  channel for the low temperature phase (dotted line) and the high temperature one (continuous line).

[1] U. Staub *et al.* Phys. Rev. Lett. 88, 126402 (2002)

[2] U. Staub *et al.* Physica B 345,23 (2004)

[3] V. Scagnoli *et al.* Jour of Magn. and Mag. Mat., in press