

Neutron Scattering Study of the Charge and the Magnetic Ordering in $\text{La}_{2-x}\text{Sr}_x\text{NiO}_4$

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Abstract

We have observed the charge and the magnetic ordering in $\text{La}_{2-x}\text{Sr}_x\text{NiO}_4$ ($0.28 \leq x \leq 0.46$) by neutron scattering. In the *c*-plane, holes and spins order as a two dimensional system, where holes form stripes and they are inserted into the antiferromagnetic ordering. Along the *c*-axis, the two dimensional ordering stacks. However, the coherence along the *c*-axis is much weaker than that in the *c*-plane. We discuss how two dimensional ordering stacks and why the coherence along the *c*-axis becomes weak.

Key words: neutron scattering; nickelate; stripe structure; magnetic ordering

There has been much interest in charge and magnetic ordering of $\text{La}_{2-x}\text{Sr}_x\text{NiO}_4$, a hole-doped two-dimensional antiferromagnet on a square-lattice with a hole concentration $n_{\text{H}} = x$. When n_{H} is small, holes form stripe structures, which are inserted into the antiferromagnetic ordering. [1] Yoshizawa *et al.* found that the optimized hole concentration for the stripe structure is $\frac{1}{3}$. At this concentration, the phase transition temperature is the highest and the coherent lengths in the *c*-plane is the longest for both charge and magnetic orderings. [2] On the other hand, Lee *et al.* reported that the coherent length along the *c*-axis becomes short at $n_{\text{H}} = \frac{1}{3}$. [3] To explore why the coherence along the *c*-axis becomes poor at $n_{\text{H}} = \frac{1}{3}$, We have carried out neutron diffraction measurements on the $\text{La}_{2-x}\text{Sr}_x\text{NiO}_4$ ($0.28 \leq x \leq 0.46$) crystal samples.

Single crystals were grown by the floating zone method. Neutron diffraction measurements were carried out with the ISSP triple-axis spectrometer GPTAS installed at the 4G experimental port of JRR-3M in JAERI (Tokai). We used neutrons with $k_{\text{I}} = 3.83 \text{ \AA}^{-1}$

obtained by the (0 0 2) reflection from pyrolytic graphite and employed the collimation of 40'-40'-40'-40' in the triple-axis mode. To remove higher order neutrons, we used a pyrolytic graphite filter.

Figure 1 shows line profiles along $(4 - 2\epsilon 0 l)$ for charge ordering and $(1 + \epsilon 0 l)$ for magnetic ordering, where $\epsilon = x$. Neither the line width of the charge ordering nor that of the magnetic ordering seem to become wider particular at $n_{\text{H}} = \frac{1}{3}$. However, the line profile shows some sample dependence at $n_{\text{H}} = \frac{1}{3}$.

Figure 2 shows the line profiles along $(1 + \epsilon 0 l)$ for two different samples with $n_{\text{H}} = \frac{1}{3}$, the A sample and the B sample. Both samples ϵ are exactly $\frac{1}{3}$, indicating that the hole concentrations of the both samples agree well. However, the line profiles are significantly different. The B sample does not exhibit even peaks and the line width of the odd peaks of the B sample is much wider than that of the A sample.

Figure 3 shows the schematic picture of the magnetic ordering at $n_{\text{H}} = \frac{1}{3}$ along the *c*-axis if only the Coulomb repulsion force between stripes in the neighbouring *c*-plane is relevant. This force translates the neighbouring stripe structure with either δ or $-\delta$ at

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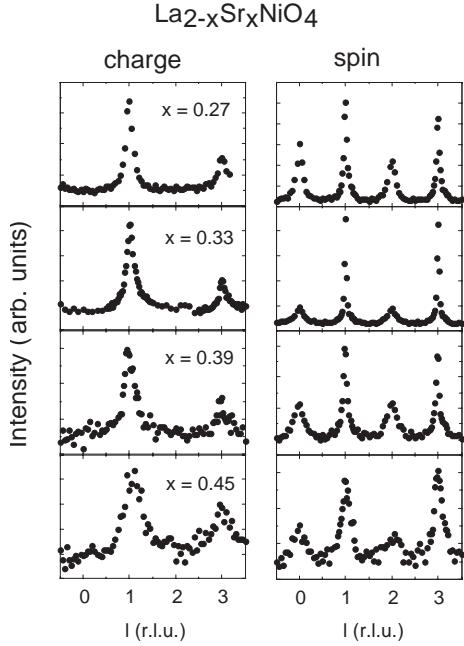


Fig. 1. The hole concentration dependence of line profiles along $(4 - 2\epsilon_0 l)$ for charge ordering and $(1 + \epsilon_0 l)$ for magnetic ordering.

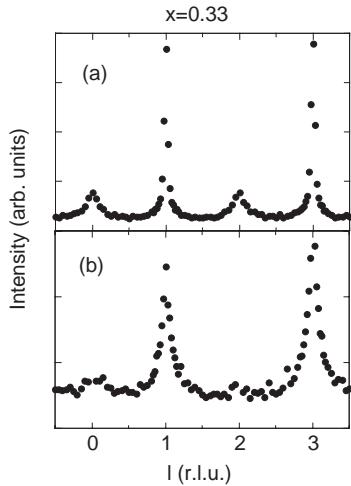


Fig. 2. The sample dependence of line profiles for magnetic ordering at $n_H = \frac{1}{3}$; (a) the A sample and (b) the B sample.

$n_H = \frac{1}{3}$, where δ denotes the spacing between the neighbouring Ni site along the a -axis. Consequently, the long range order cannot be established. A next-nearest "ferro" charge interaction, J_{NN} , is necessary for establishing the long-range order.

Figure 4 shows the calculated line profiles of the $\frac{1}{3}$ sample for strong and weak J_{NN} . If J_{NN} is weak, the coherent length along the c -axis becomes short and depresses the even peak intensity. The line profile of the B sample is similar to that with $J_{NN} = 0$. Thus,

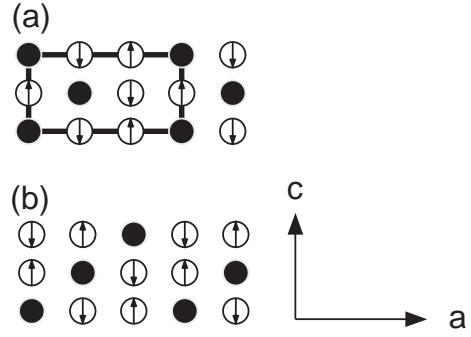


Fig. 3. The schematic picture of two possible positions for the second-nearest stripe structure of the $\frac{1}{3}$ sample. A closed circle denote a hole stripe.

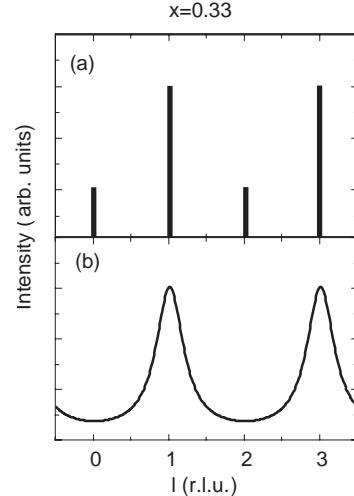


Fig. 4. Calculated magnetic profiles for the $\frac{1}{3}$ sample; (a) the Bragg reflections corresponding to the (a) structure in Fig. 3 built by $J_{NN} >> 0$, and (b) the diffuse signal due to the equivalent mixture of the (a) structure and the (b) structure in Fig. 3 built by $J_{NN} = 0$.

we believe that J_{NN} of the $\frac{1}{3}$ sample is so weak that some subtle perturbed forces (e.g. crystal defects) can disturb the magnetic ordering.

References

- [1] J. M. Tranquada, J. E. Lorenzo, D. J. Buttrey, V. Sachan Phys. Rev. B **52** (1995) 3581
- [2] H. Yoshizawa, T. Kakeshita, R. Kajimoto, T. Katsufuji, T. Tanabe, Y. Tokura, Phys. Rev. B **61** (2000) R854.
- [3] S.-H. Lee, S.-W. Cheong, K. Yamada, C. F. Majkrzak, Phys. Rev. B **63** (2001) 060405 .