

# Polarization Dependent XAFS Studies on Low Coverage of Ni at the Al<sub>2</sub>O<sub>3</sub> Surface

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## Introduction

Ni supported on Al<sub>2</sub>O<sub>3</sub> is an important catalyst for CO and alkene hydrogenation reaction. However, the chemical interaction between Ni and Al<sub>2</sub>O<sub>3</sub> is still unknown. In this paper we report our recent results about polarization dependent XAFS investigation on Ni at the Al<sub>2</sub>O<sub>3</sub> (0001) single crystal surface.

## Experimental

### Sample preparation

$\alpha$ -Al<sub>2</sub>O<sub>3</sub>(0001) (20x20x1mm<sup>3</sup>) single crystal was purchased from Earth Jewelry Co. The sample was first treated at 1173 K. Then the sample was put in a UHV chamber and annealed at 873 K. We obtained the (1x1) LEED patterns by this treatment. XPS suggests that little carbon impurity is present. Ni was deposited on  $\alpha$ -Al<sub>2</sub>O<sub>3</sub> by resistive heating of W wire on which Ni wire is wrapped. The Ni coverage was monitored by XPS using Ni2p peak and O KLL Auger peak. The coverage was estimated to be 0.04 ML.

### XAFS measurement

Polarization-dependent XAFS measurements were carried out in a newly-built UHV XAFS chamber for total reflection fluorescence mode. The base pressure of the chamber is  $5 \times 10^{-10}$  Torr. XAFS measurement was carried out in BL9A of Photon Factory in Institute for Material Structure Science, (KEK-IMSS-PF). Synchrotron radiation emitted from 2.5 GeV 400 mA electron storage ring was monochromatized by double crystal monochromator together with focusing mirrors. The X-ray is collimated by a pin-hole slit to avoid needless radiation other than the sample. Fluorescence signal was detected by 19 elements solid state detector (Cambera).

## Results

Figure 1 shows polarization dependent XAFS oscillations for the Ni on Al<sub>2</sub>O<sub>3</sub> at a coverage of 0.04 ML. We carried out curve fitting analysis and found that the Ni-O distances for both directions were 0.203 nm a little shorter than that predicted by the theoretical calculation.<sup>1)</sup> We proceed detailed analysis to provide the model structure of Ni on Al<sub>2</sub>O<sub>3</sub>(0001).

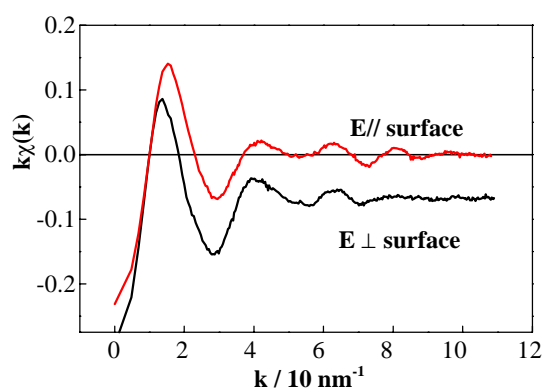


Fig.1 XAFS oscillations of 0.04 ML Ni on Al<sub>2</sub>O<sub>3</sub>(0001) surface.

### References

- 1) Q. Ma, K. Klier, H. Cheng, J. W. Mitchell and K. S. Hayes, *J. Phys. Chem.*, **105**, 2212 (2001).

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