

STM study of the TiO₂(110)-(1x2) surface reconstruction

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The study of the metal oxide surface is of a great scientific interest due to their huge technological applications. In particular, the rutile TiO₂ has been the subject of numerous investigations since it is relevant for a wide variety of fields, ranging from catalysis to fundamental chemistry. Amongst the different surfaces of the single-crystalline TiO₂ system, the (110) face is the most stable one being its 1x1 structure extensively studied in the last decades. This well-known (1x1) structure transforms into a long ordered reconstruction with a (1x2) symmetry after reducing the TiO₂(110) surface by annealing up to about 800°C. The understanding of the basic properties of the non-stoichiometric (1x2) TiO₂(110) reconstruction is still nowadays unclear. Recently, the atomic configuration of the 1x2 TiO₂(110) surface has been elucidated by combining scanning tunneling microscopy (STM), quantitative LEED and density functional theory (DFT) [1]. The resulting model consists of added Ti₂O₃ rows along the [001] direction in agreement with a previous proposal by Onishi et al [2].

In the present work, the non-stoichiometric (1x2) TiO₂(110) reconstruction was studied by STM and scanning tunneling spectroscopy (STS). The recorded images show monoatomic steps, wide terraces and no cross-link features as it can be observed in Figure 1. STS current-voltage curves were recorded to obtain information on the local density of states (DOS).

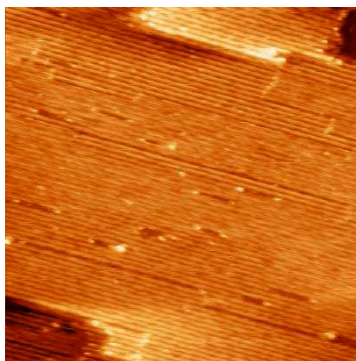


Figure 1: STM image of the 1x2 TiO₂ (110) surface reconstruction

Keywords: Scanning Tunnelling microscopy, titanium oxide

References:

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Topic: Surface Science