

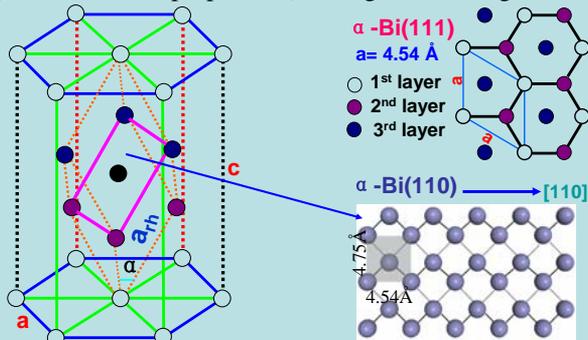
Self-assembly of Bi nanostructure on HOPG, MoS₂ and silicon nitride

Chu, X.J.; Zhang, H.L.; Kushvaha, S.S.; Tang, Z.; Tang P. and Wang, X.S.

Department of Physics, Faculty of Science, National University of Singapore, Singapore 117542
(g0600514@nus.edu.sg)

INTRODUCTION

◆ Bismuth, a group-V semimetal, has unique atomic (*Rhombohedral* lattice) and electronic properties (de Broglie wavelength ~ 40 nm).

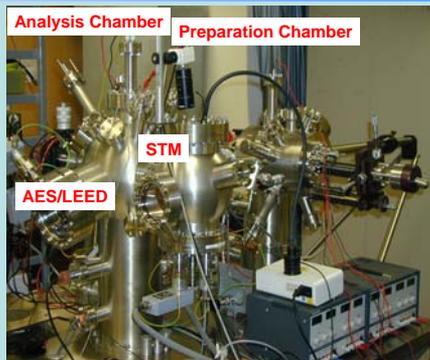


◆ HOPG (Highly Oriented Pyrolytic Graphite) and MoS₂ are easy-to-prepare inert conductive substrate for growing nearly *free-standing* nanostructures, sometimes *1D or quasi 1D* nanostructures.

◆ Si-based inert surfaces: dielectric layers (SiO₂, Si₃N₄, SrTiO₃) on Si, close to real applications.

◆ In this work, Bismuth was deposited on HOPG, MoS₂ and Si₃N₄. *Nanorods, nanoribbons and nanoclusters* were formed.

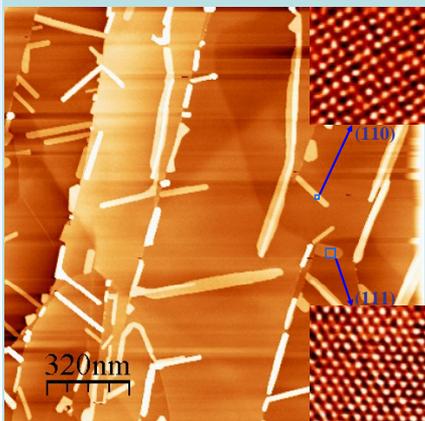
EXPERIMENTAL DETAILS



- ◆ Base Pressure: 1×10^{-10} mbar
- ◆ Bi evaporator boat
- ◆ Flux calibrated with STM & AES
- ◆ HOPG and MoS₂ sample cleaved in air and degassed in vacuum at 700~800K
- ◆ Si₃N₄ prepared by thermal nitridation of Si(111)-7x7
- ◆ STM images at RT

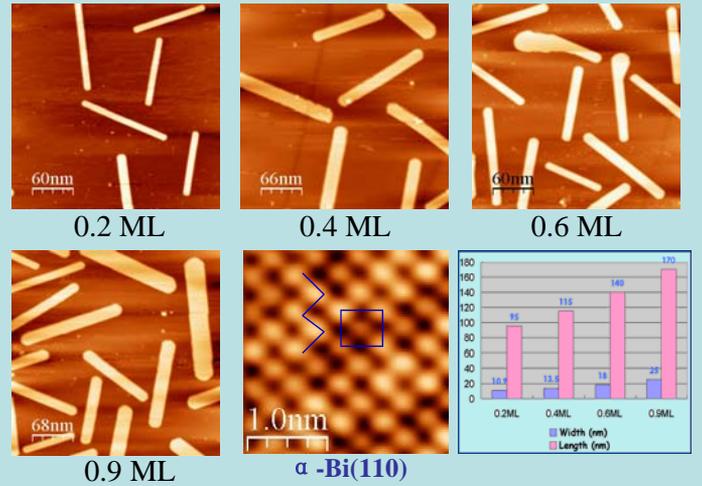
RESULTS AND DISCUSSIONS

I. Bi nanorods on HOPG



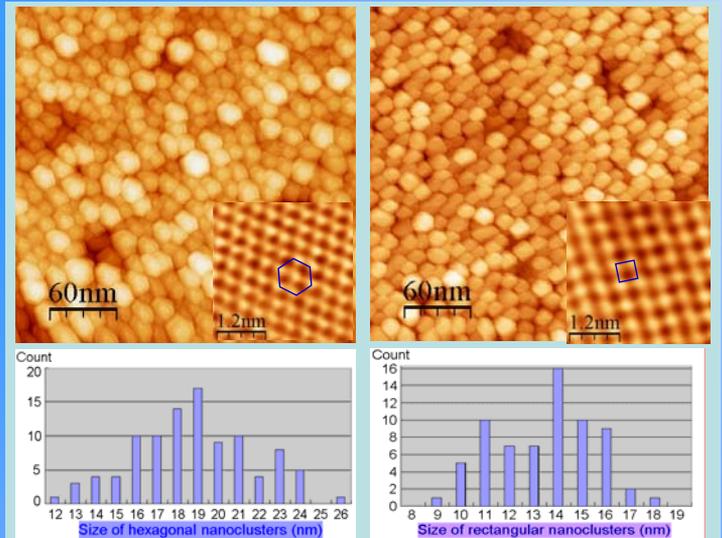
- Low surface coverage
- Edge-decoration with Bi nanorods
- Heights range from 6.6 Å to 33 Å
- Nanorods with height of 8 Å: α -Bi(111) surface
- Nanorods with height of even number of 3.3 Å (one atomic layer spacing): α -Bi(110) surface

II. Bi nanoribbons on MoS₂



- Uniform height of 6.6 Å
- Angles between the nanoribbons are 0°, 60° or 120°, corresponding to the *three-fold symmetry* of the substrate.

III. Bi nanoclusters on Si₃N₄



- Silicon nitride surface was *passivated with Bi*
- Bi forms single crystal faceted clusters at *room temperature*
- Coexistence of *hexagonal* and *rectangular* facet clusters

CONCLUSION

- ◆ The Bi(110) islands with even-number layers is stabilized by forming a puckered-layer structure
- ◆ There is a natural tendency for faster diffusion along $\langle 110 \rangle$ directions in a low flux environment
- ◆ The direction $\langle 110 \rangle$ corresponding to the direction of the long zigzag chains of covalently bonded atoms, likely produces very elongated structures.

Jointly Organized by

