

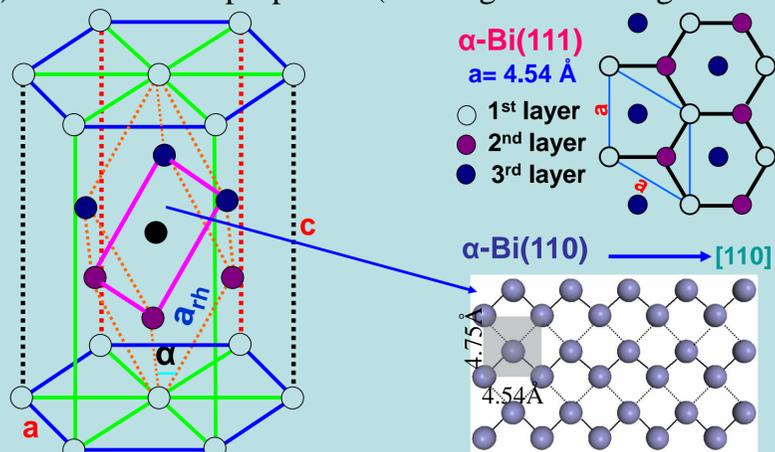
Self-assembly of Bi nanostructure on HOPG, MoS₂ and silicon nitride

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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.

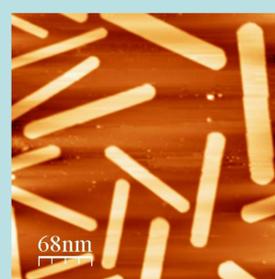
II. Bi nanoribbons on MoS₂



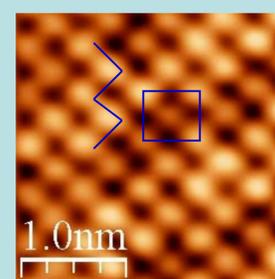
0.2 ML

0.4 ML

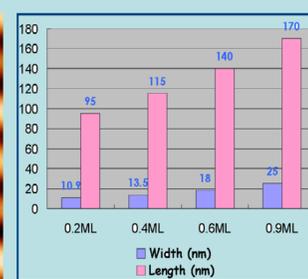
0.6 ML



0.9 ML

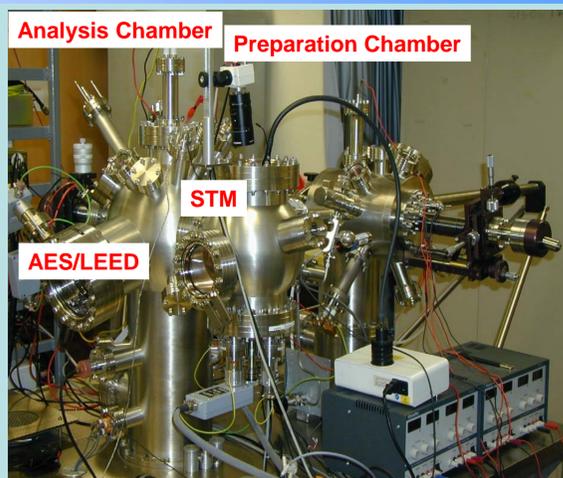


α -Bi(110)



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

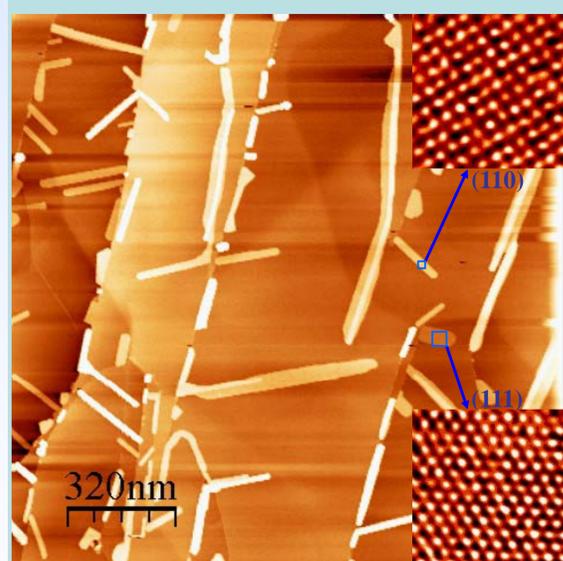
EXPERIMENTAL DETAILS



- ◆ 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*

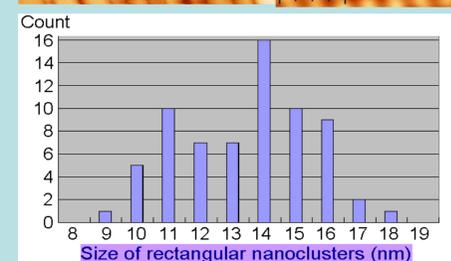
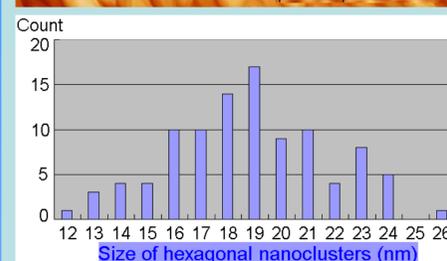
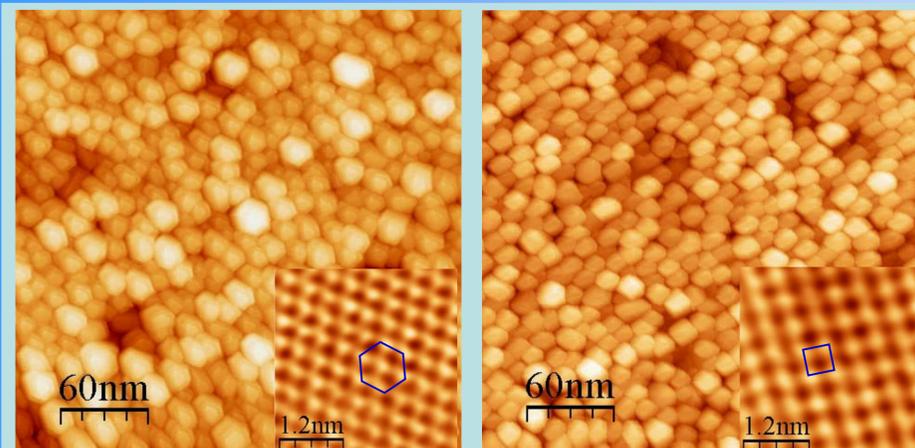
RESULTS AND DISCUSSIONS

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*

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

