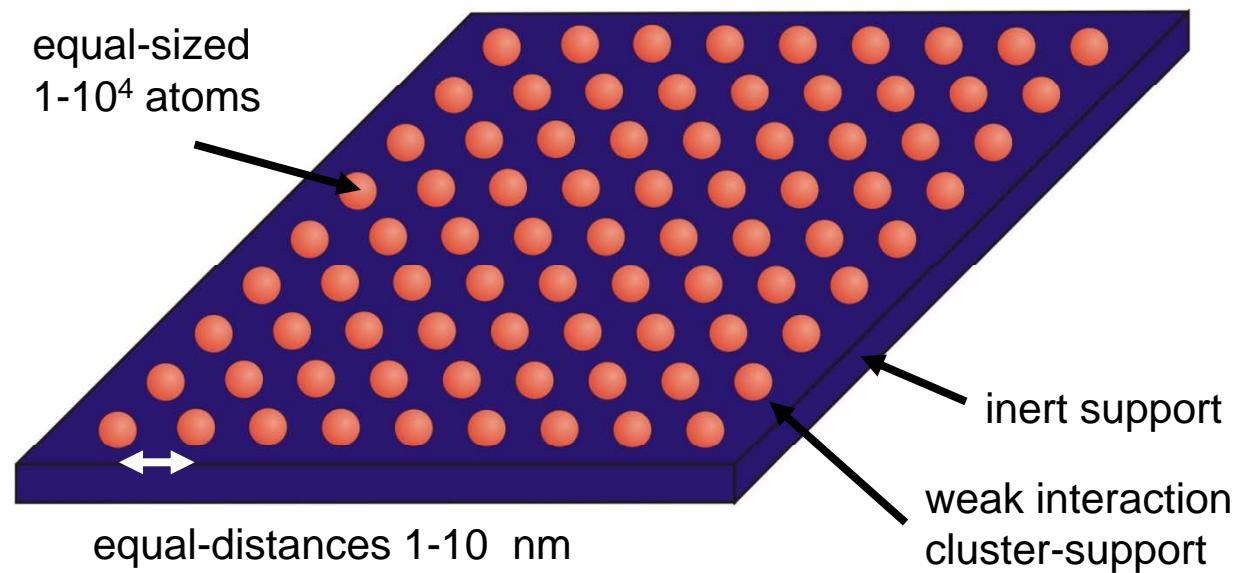


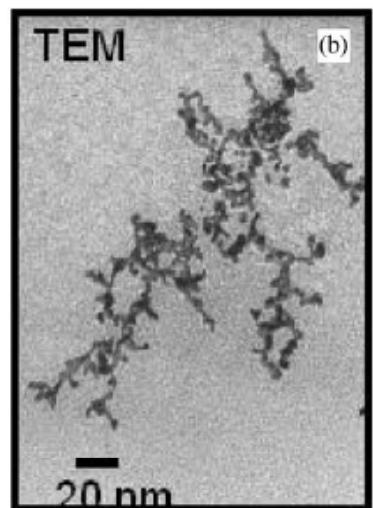
Ideal Cluster Research Material



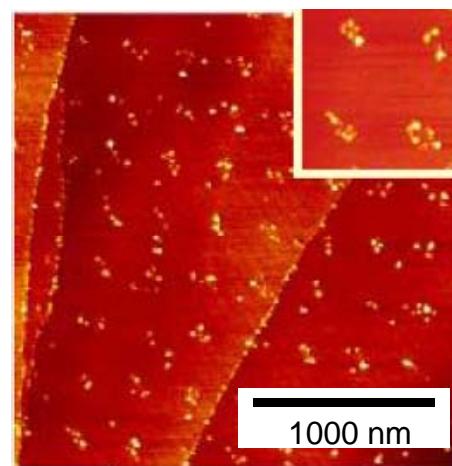
- address single clusters
- collective cluster response
- cluster coupling

Fig. 5.16

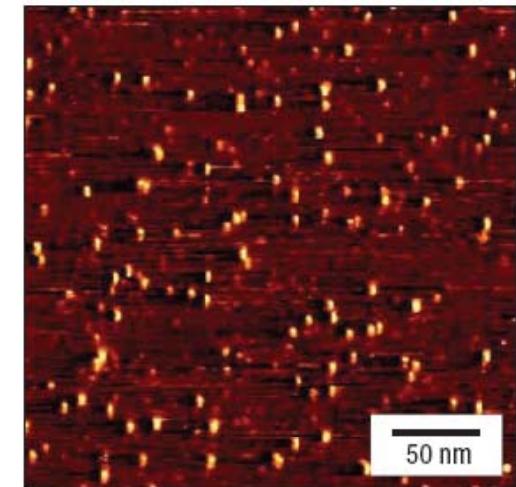
Deposition of Mass-Selected Clusters



Pt-clusters, 0.25 eV/atom
→ aggregation [1]



FIB-prepatterned,
then Au₇₅₀
0.25 eV/atom [1]



Ag₁₄₇, 14 eV/atom [2]

Fig. 5.17

- [1] P. Mélinon et al., J. Crys. Growth 275 (2005) 317,
- [2] R.E. Palmer, S. Pratontep, H.-G. Boyen, Nature Materials 2 (2003) 443

Ir-Clusters on Graphene on Ir(111)

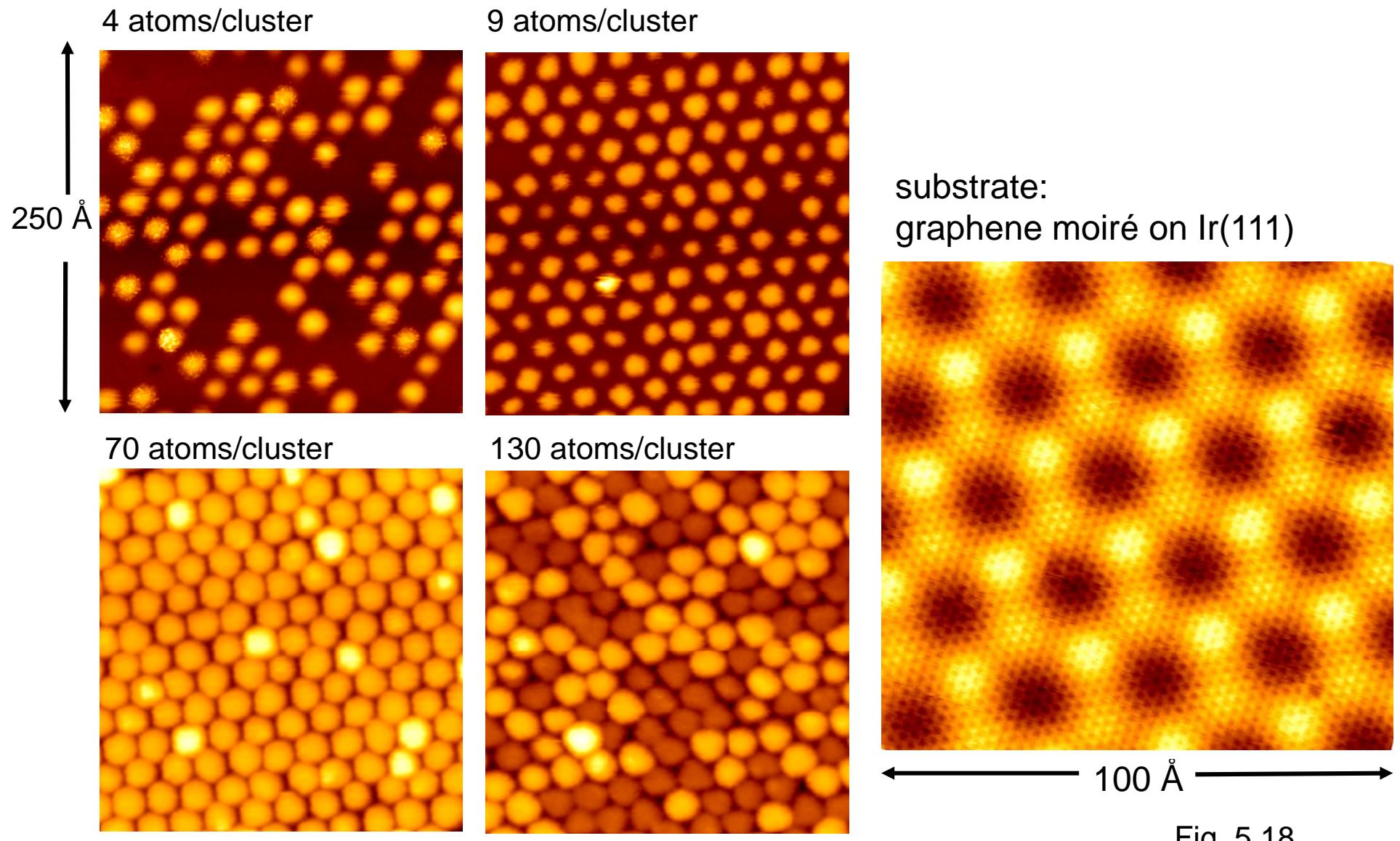
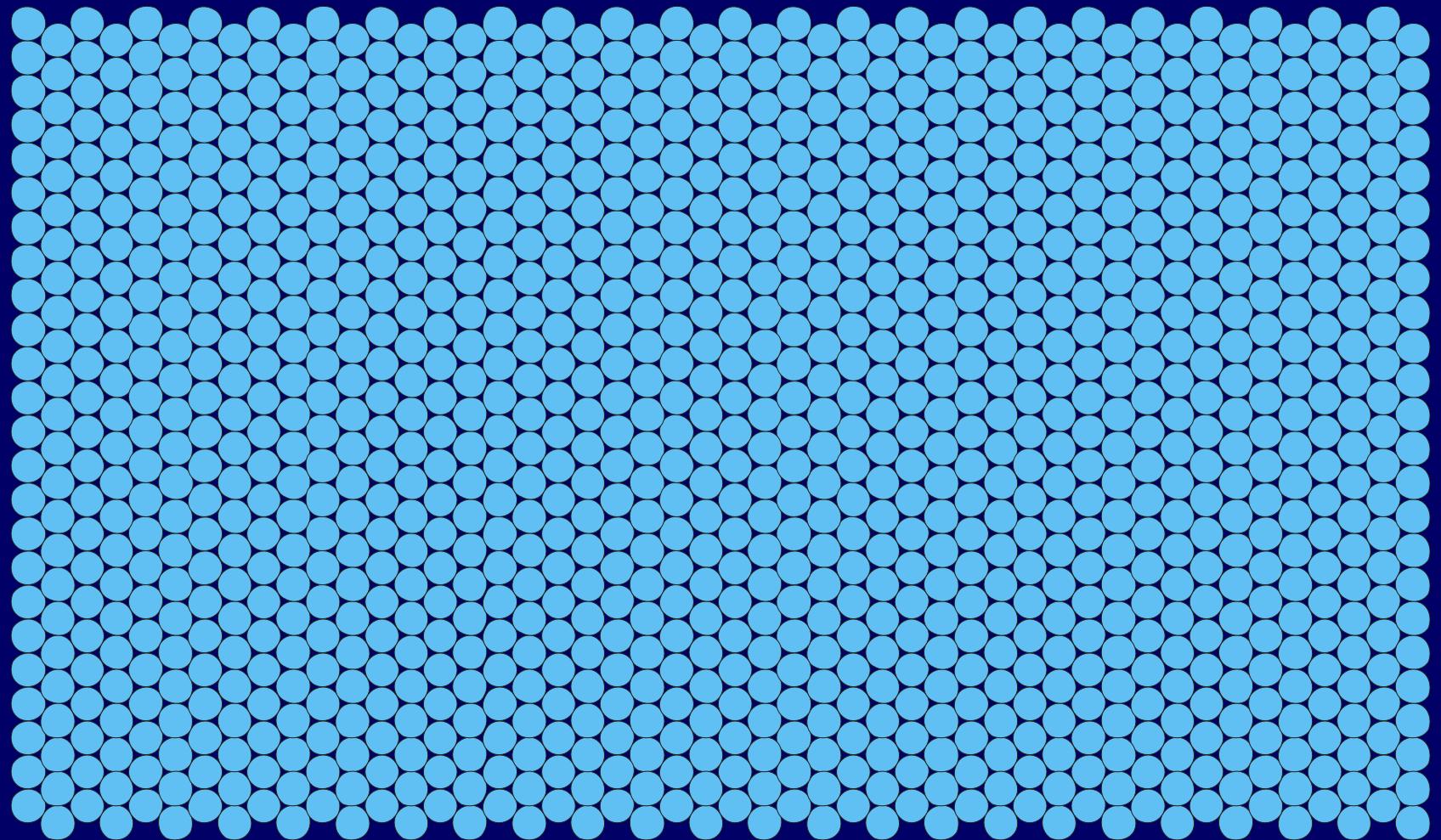


Fig. 5.18

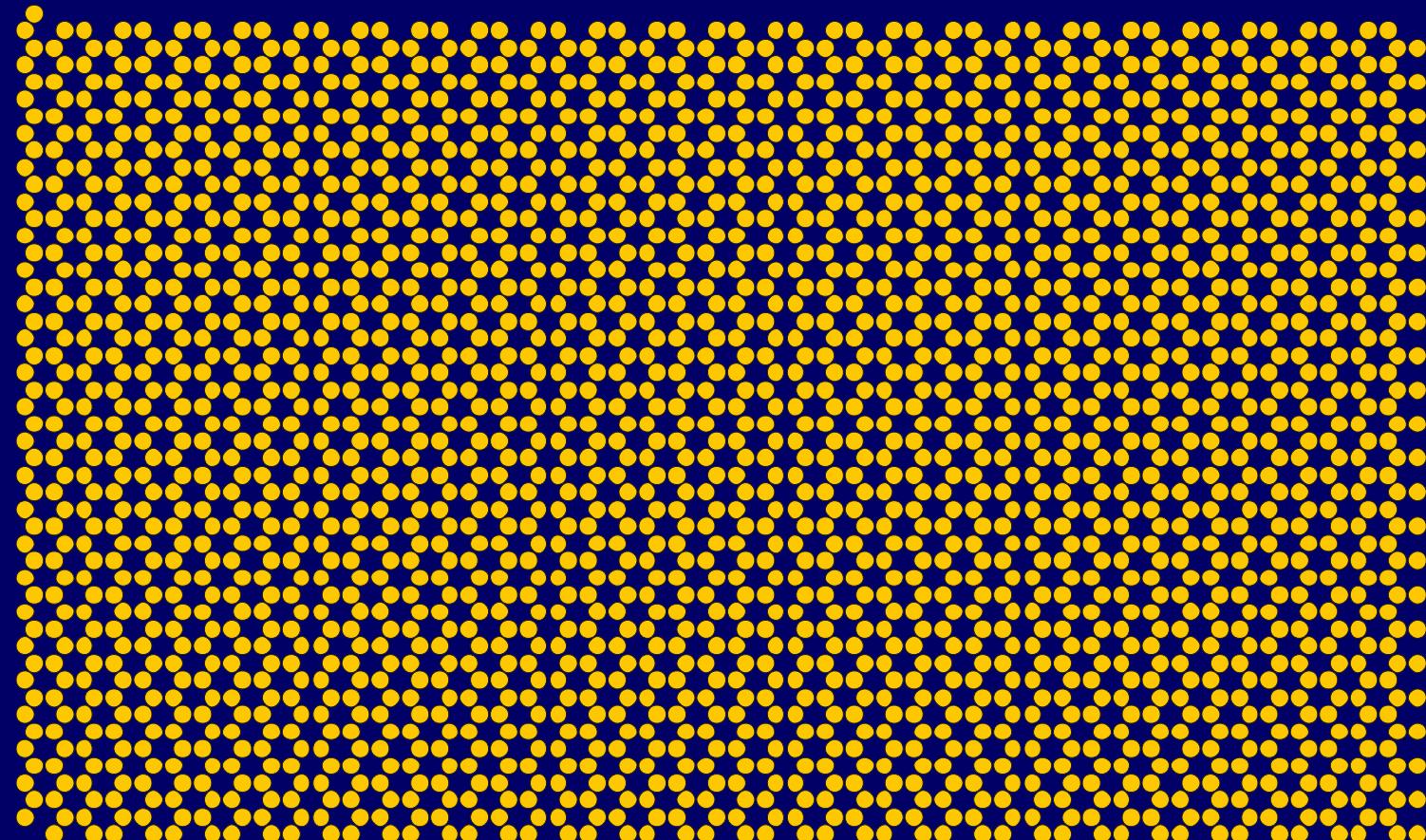
Origin of Moiré



Ir(111)

Fig. 5.19

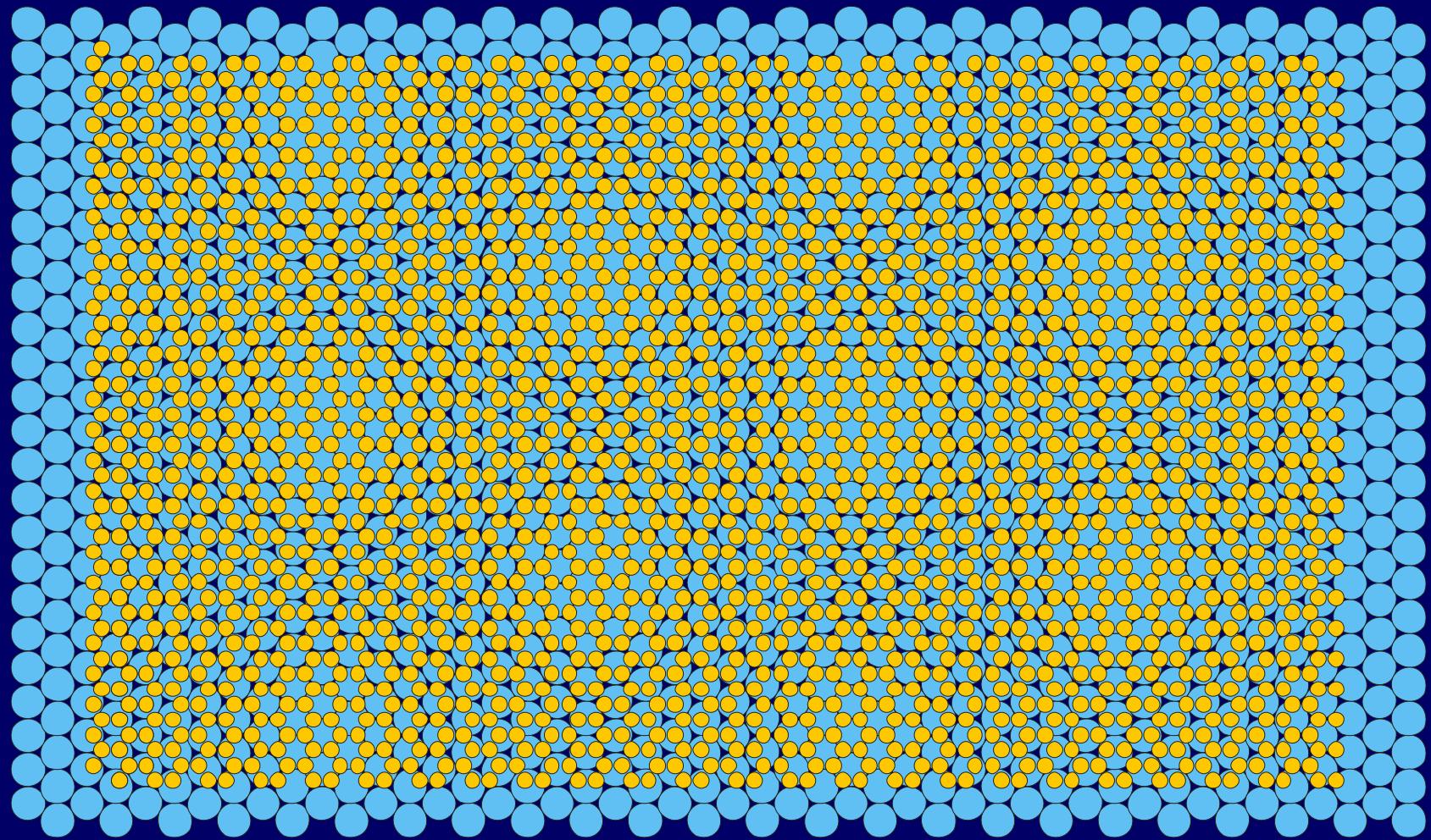
Origin of Moiré



graphene

Fig. 5.20

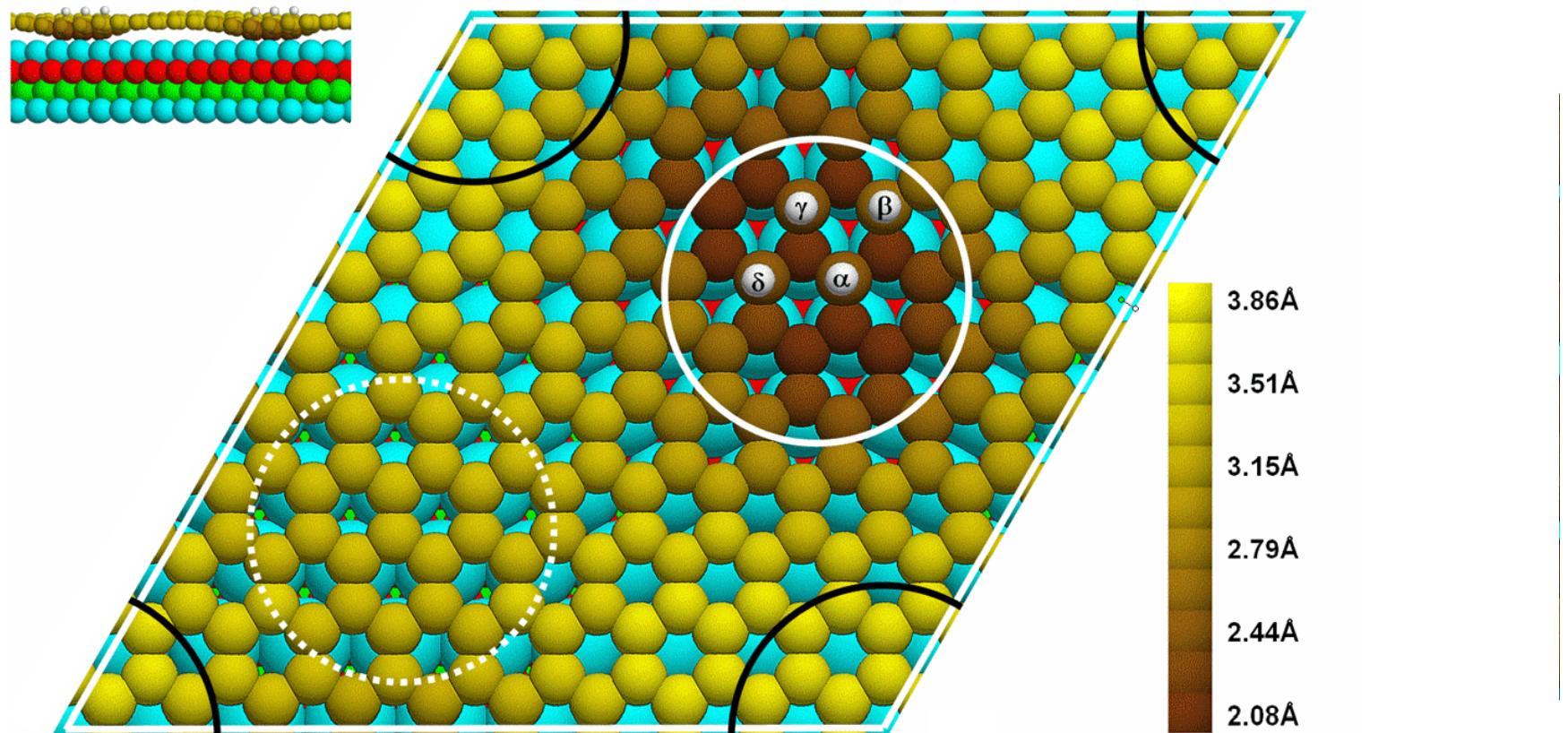
Origin of Moiré



Ir(111) and graphene

Fig. 5.21

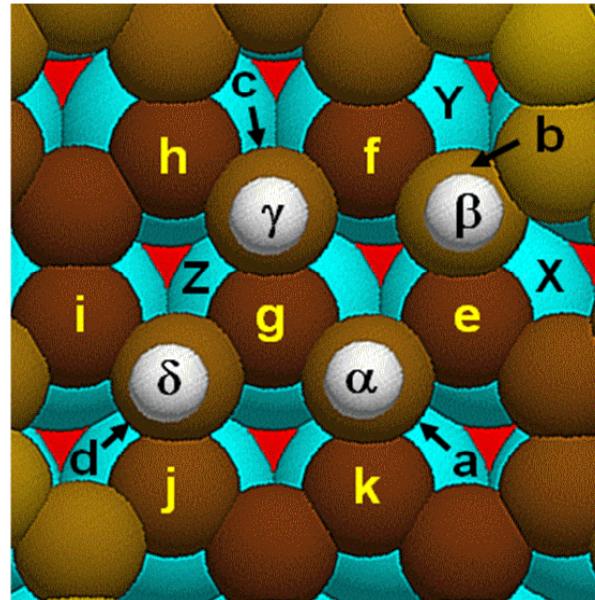
DFT-calculations: Pinning Effect



- short bonds $\approx 2.1 \text{ \AA} \approx$ sum of Ir and C atomic radii
- more than twice C-Ir chemical bonds than adatoms !
- hcp-type preferred by 3.7 eV

Fig. 22

Nano-Diamond Formation Underneath Cluster



bond angles:

$$\begin{aligned} Xea, Xeb, Yfb, Yfc, Zgc, Zgd, Zga &= 111^\circ, 107^\circ, 107^\circ, 110^\circ, 107^\circ, 109^\circ, 108^\circ \\ ea\alpha, eb\beta, fb\beta, fc\gamma, gc\gamma, gd\gamma, gd\delta, gd\alpha &= 106^\circ, 108^\circ, 108^\circ, 105^\circ, 106^\circ, 111^\circ, 106^\circ \end{aligned}$$

Fig. 5.23

tetrahedral bond angle 109,5°

rehybridization from sp^2 to sp^3

Cluster Lattices of Pt, Re, W, Ir

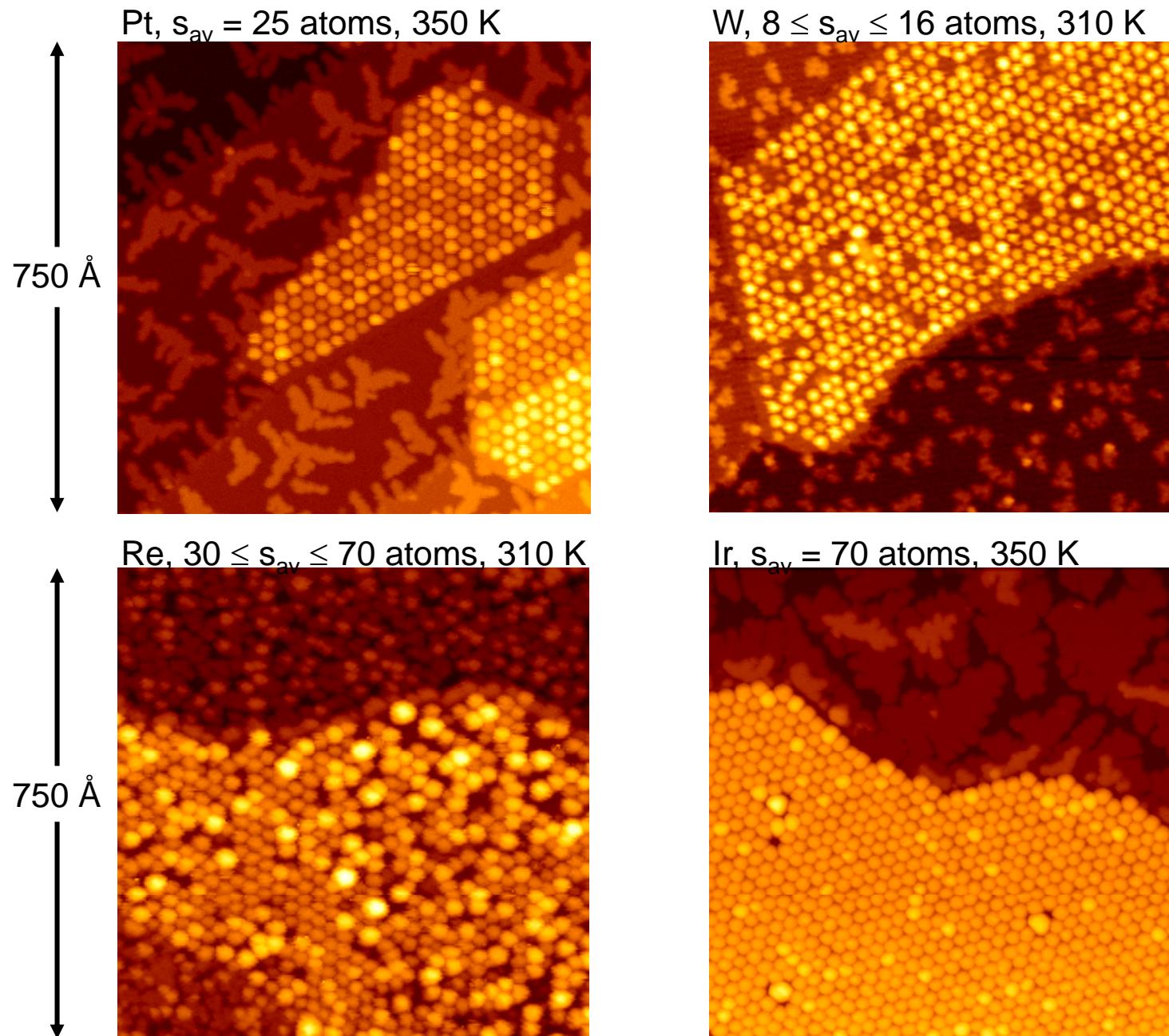
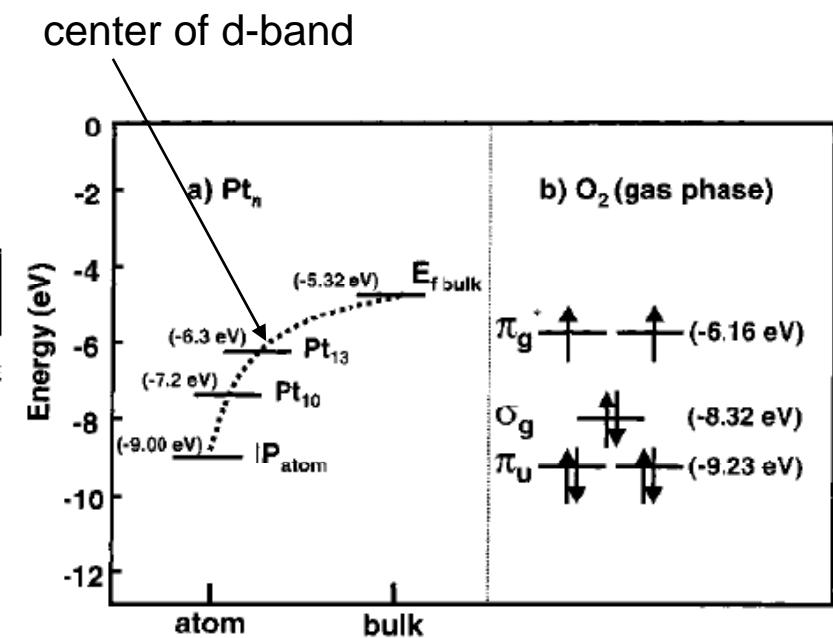
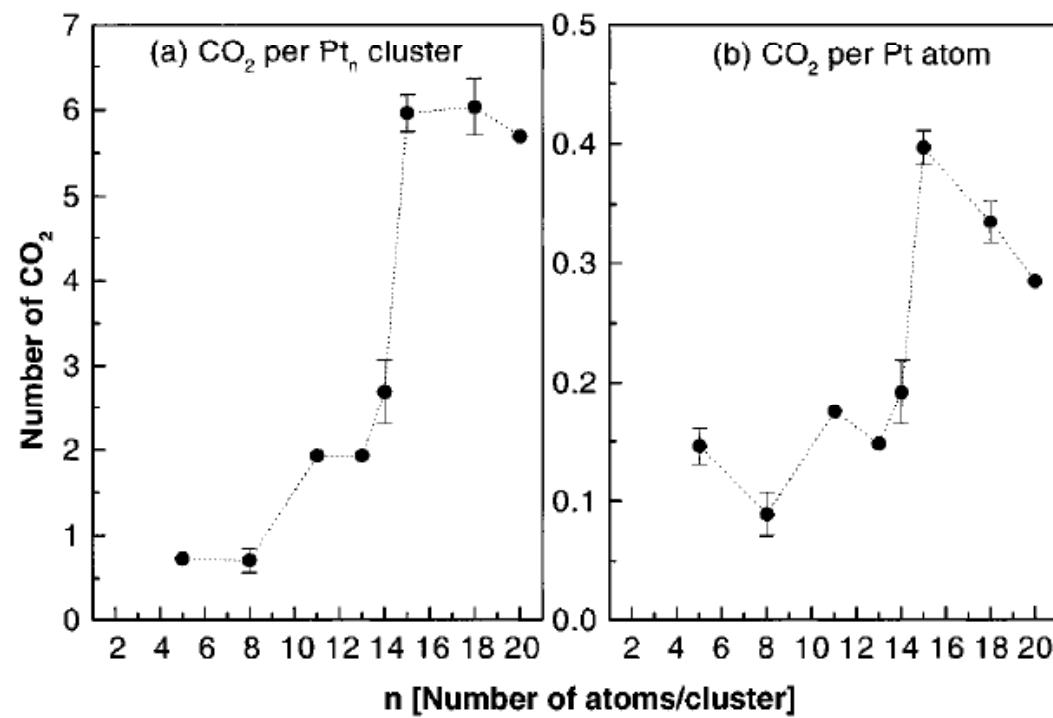


Fig. 5.24

Size Dependence of CO Oxidation by Pt-Clusters on MgO



Size Dependence of CO Oxidation by Au-Clusters on TiO₂

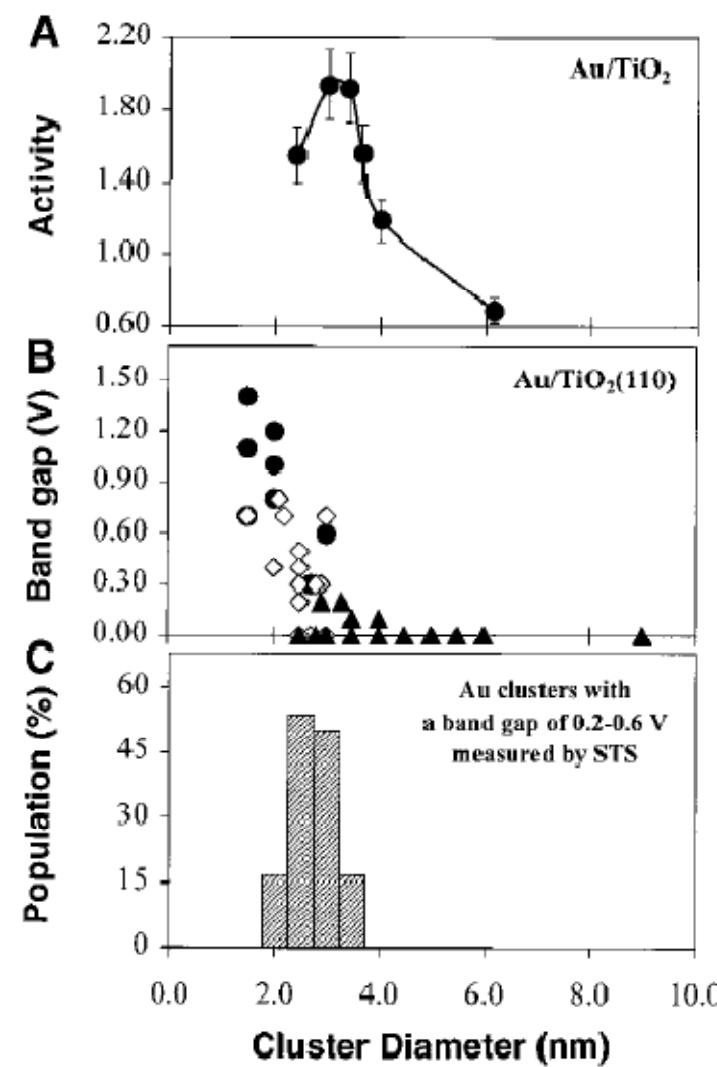
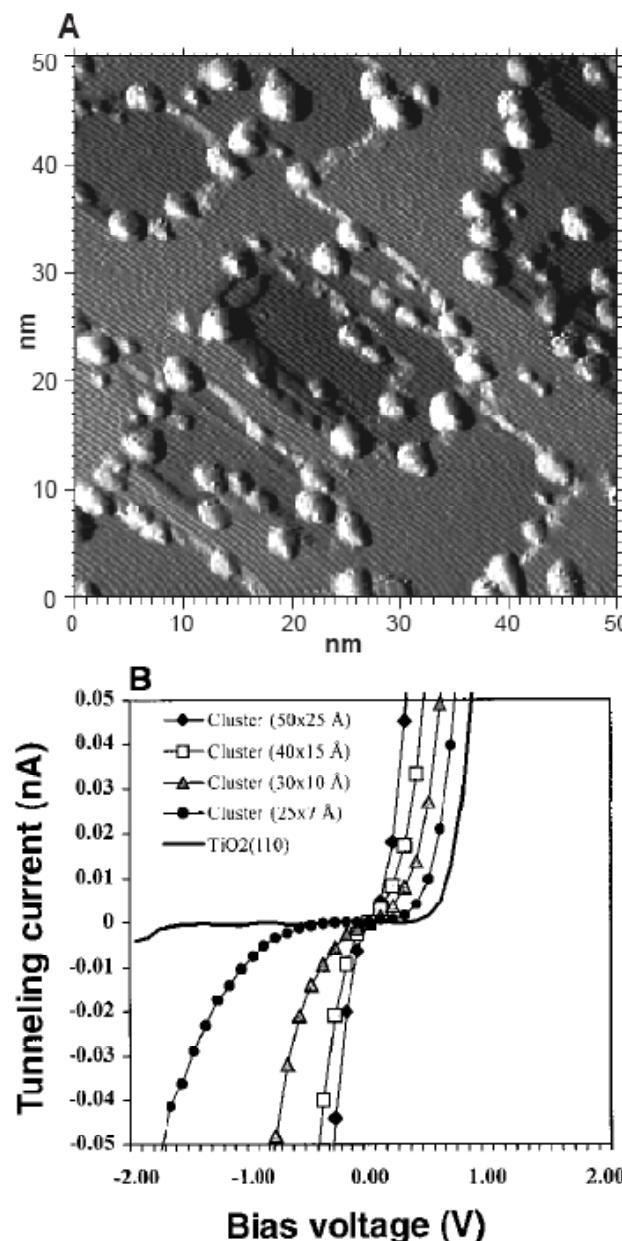
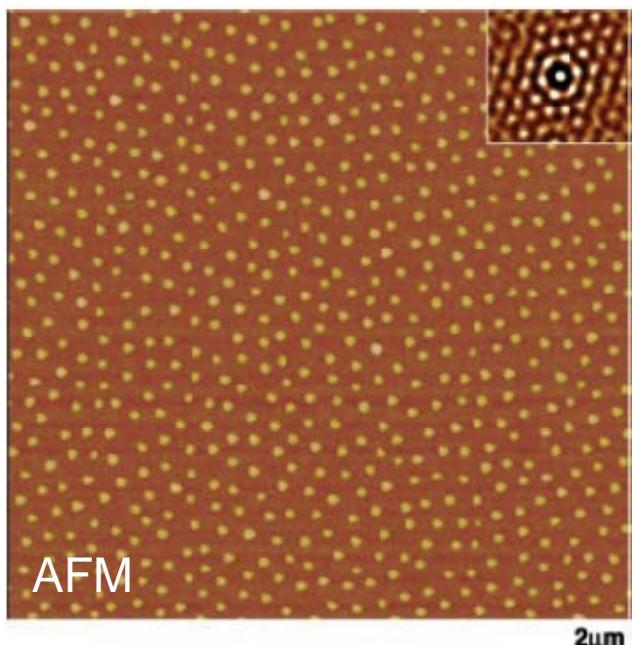


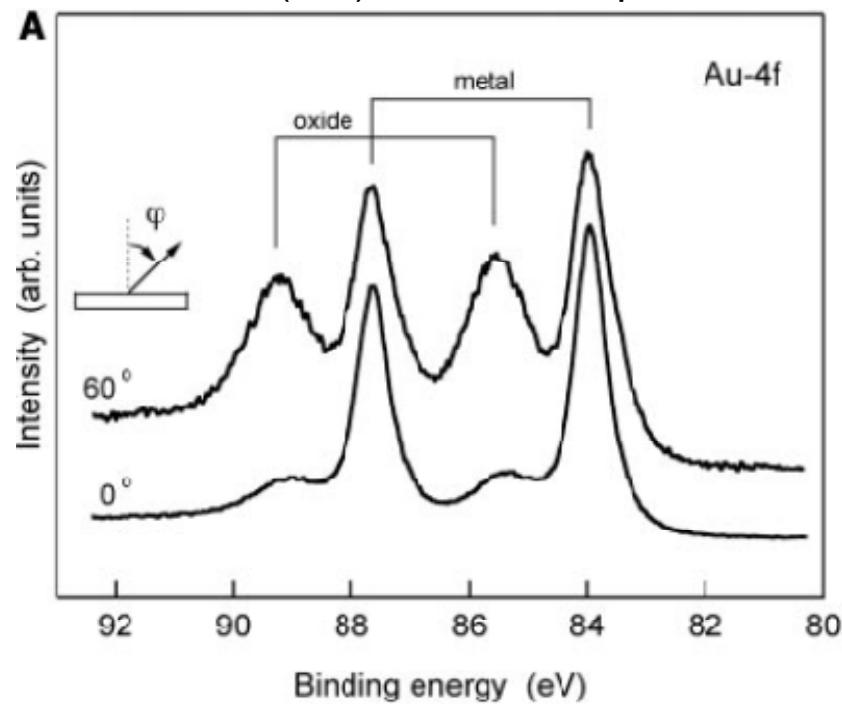
Fig. 3. (A) The activity for CO oxidation at 350 K as a function of the Au cluster size supported on TiO₂(110)-(1×1) assuming total dispersion of the Au. The CO:O₂ mixture was 1:5 at a total pressure of 40 Torr. Activity is expressed as $(\text{product molecules}) \times (\text{total Au atoms})^{-1} \text{ s}^{-1}$. (B) Cluster band gap measured by STS as a function of the Au cluster size supported on TiO₂(110)-(1×1). The band gaps were obtained while the corresponding topographic scan was acquired on various Au coverages ranging from 0.2 to 4.0 ML. (●) Two-dimensional (2D) clusters; (□) 3D clusters, two atom layers in height; (▲) 3D clusters with three atom layers or greater in height. (C) Relative population of the Au clusters (two atom layers in height) that exhibited a band gap of 0.2 to 0.6 V as measured by STS from Au/TiO₂(110).

M. Valden, X. Lai, D.W. Goodman, Science 281 (1998) 1647

Fig. 5.26



XPS of Au(111) treated in O plasma



Resistance of Au-Clusters to Oxidation by an O-Plasma

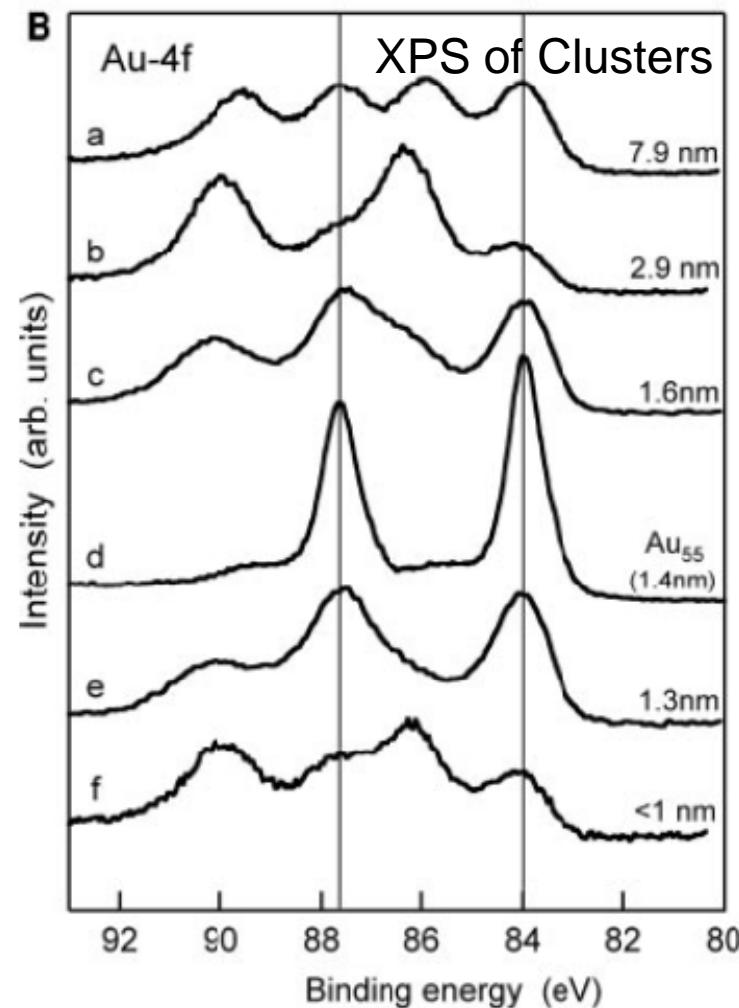


Fig. 5.27

H.-G. Boyen et al., Science 297 (2002) 1533

The Future of Magnetic Data Storage

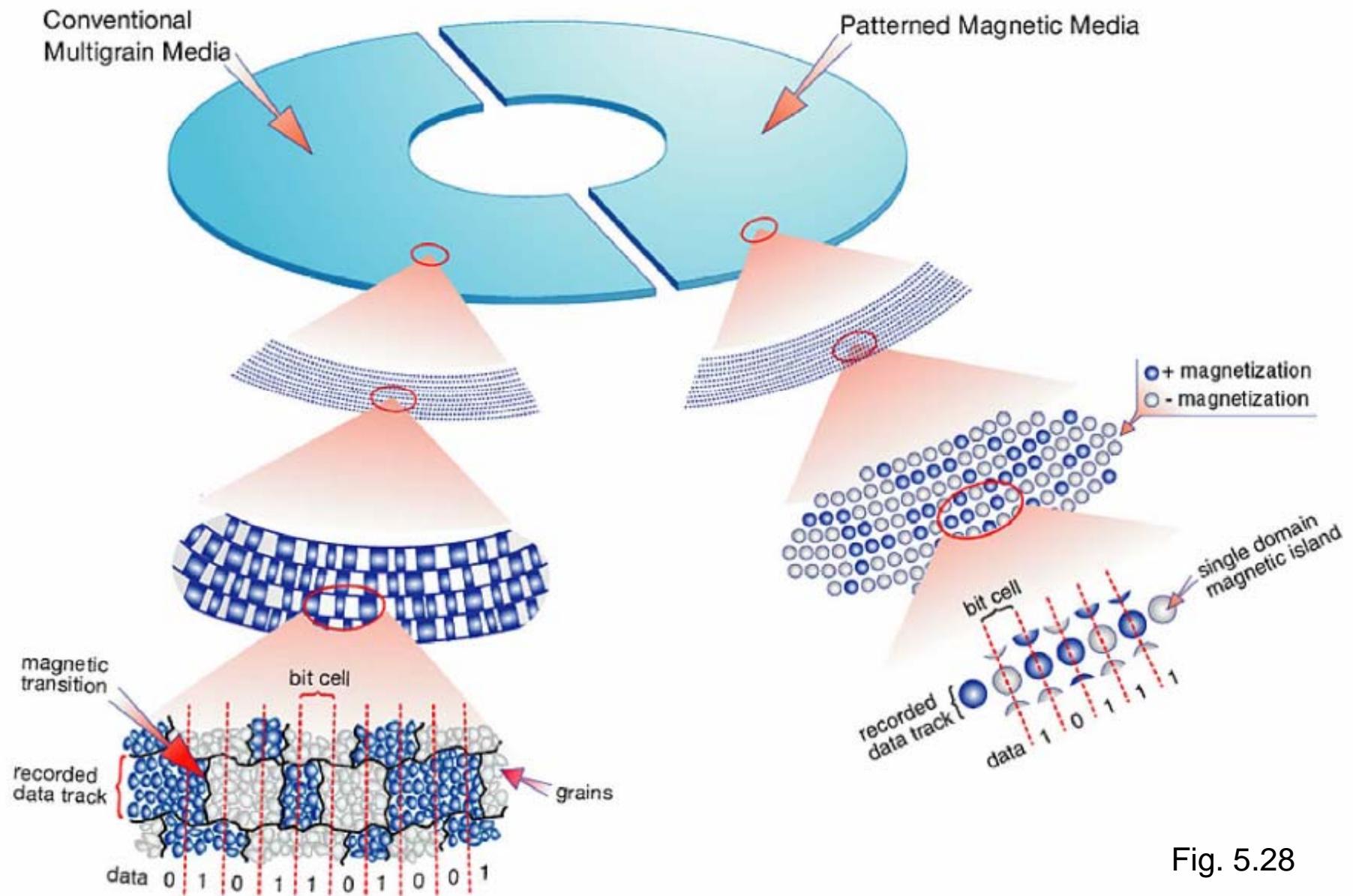
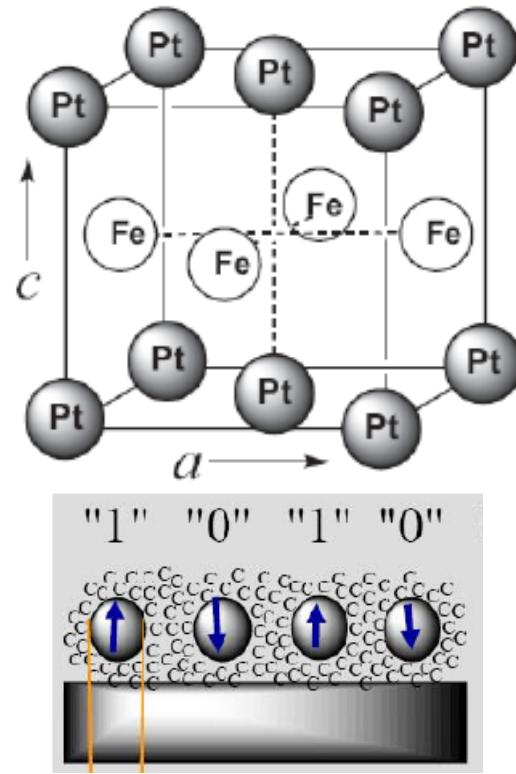


Fig. 5.28



The FePt $\text{L}1_0$ -Phase

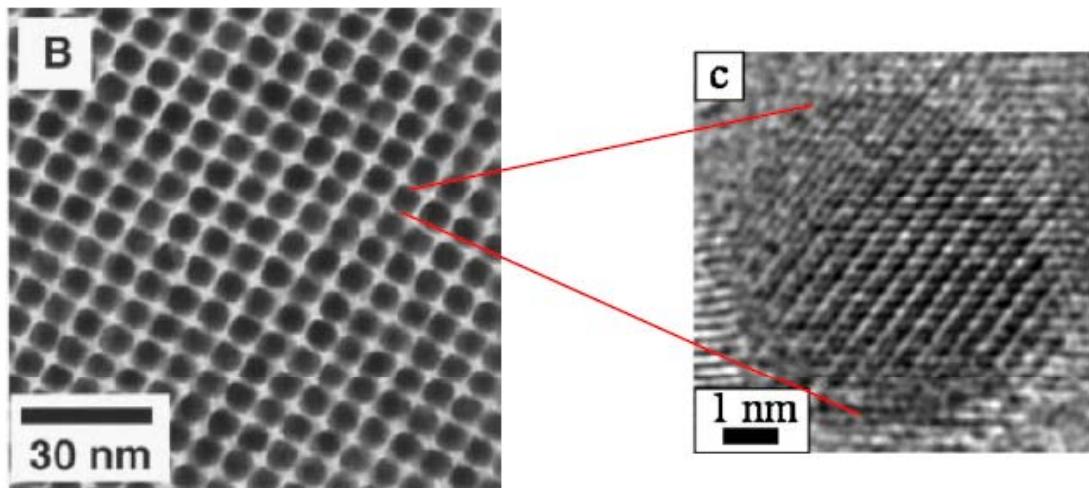
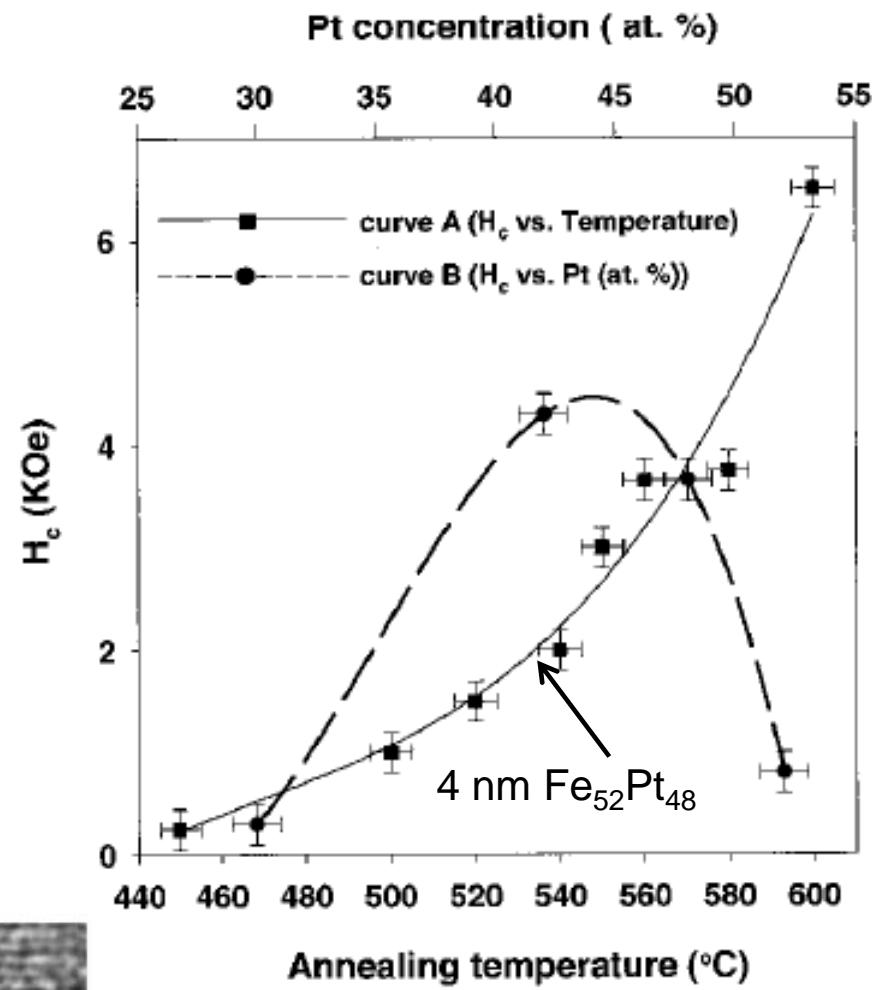
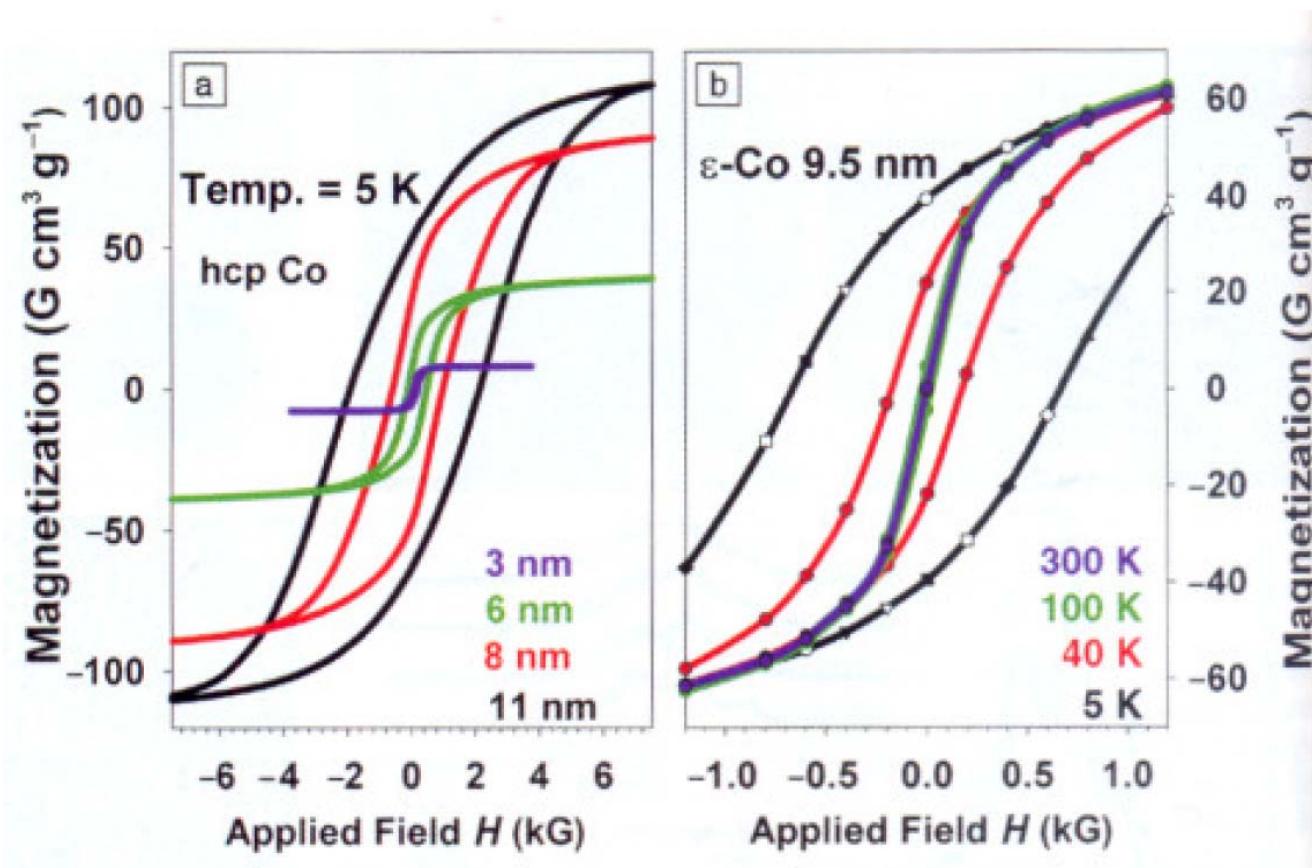


Fig. 5.29

Magnetization and Hysteresis in Dependence of Particle Size and Temperature for Co Clusters

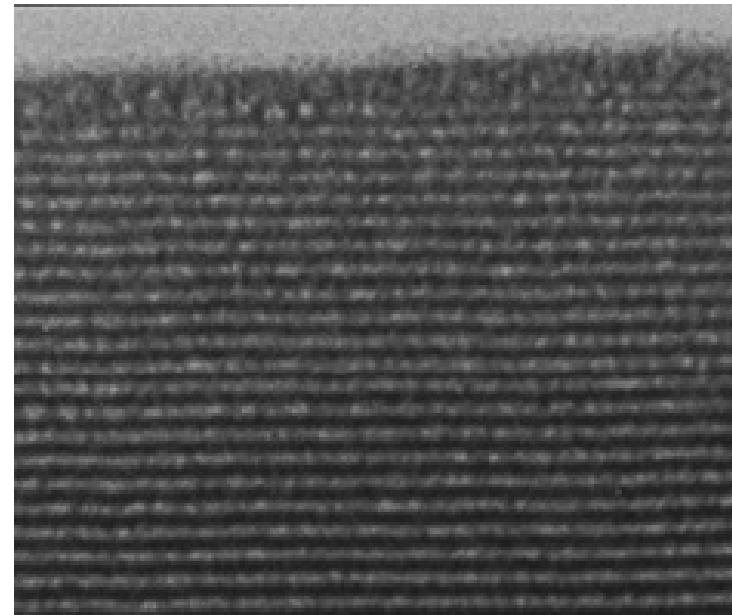


Figures of Chapter 6: Epitaxy and Thin Films

Examples of Thin Film Applications



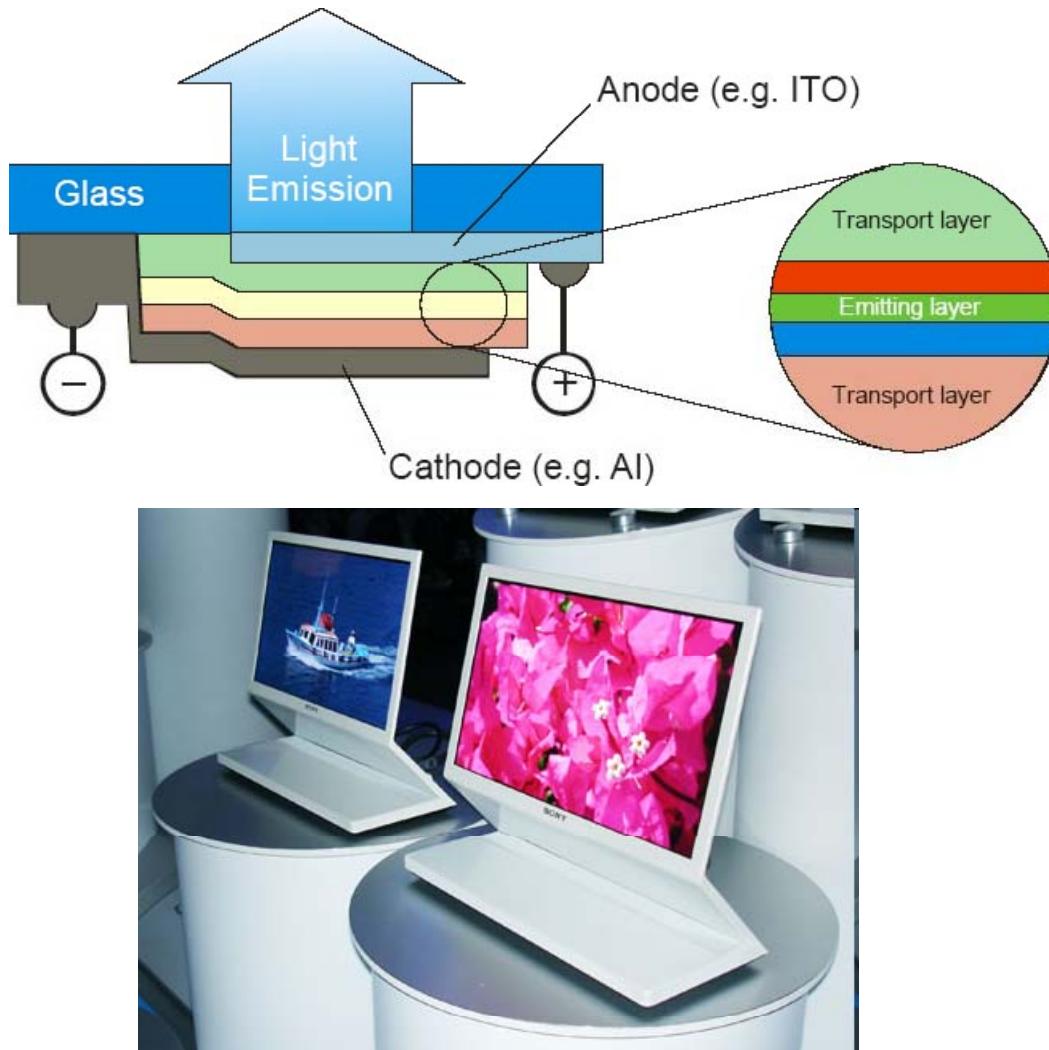
TiN hard-coating for steel cutting tools



TEM of a W-Si X-ray mirror,
periodicity 2.5 nm

Fig. 6.1

Examples of Thin Film Applications



organic light emitting diode OLED

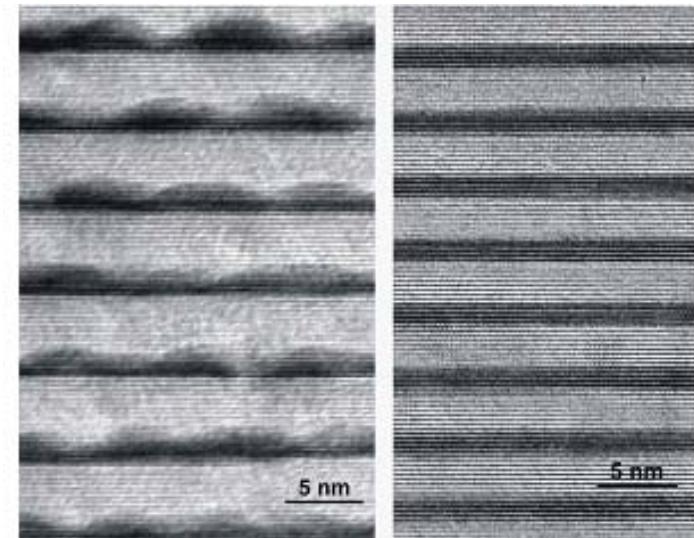


Figure 1: High-resolution transmission electron microscopy images of QD (left) and QW (right) superlattices for ISB absorption at 1.55 μm .

Fig. 6.2

Energy Minimizing Processes – Limited Mobility

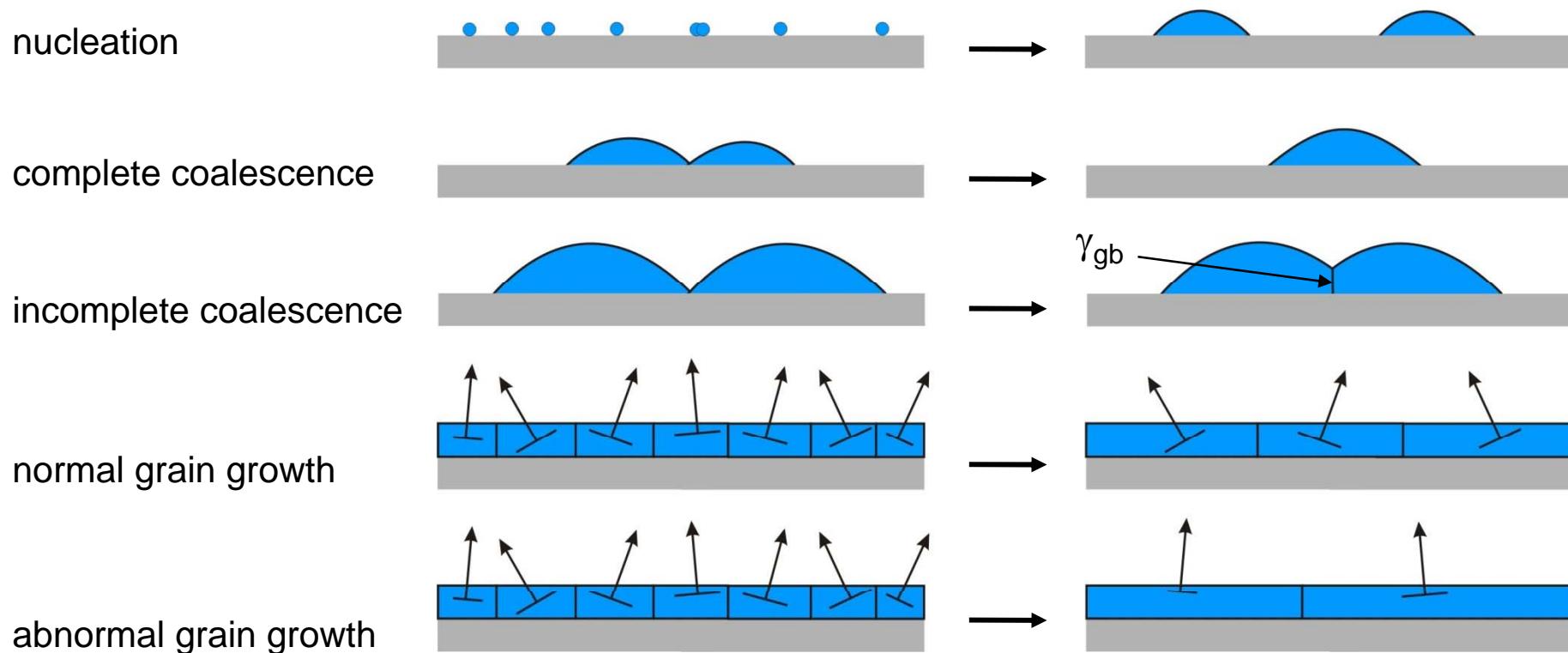
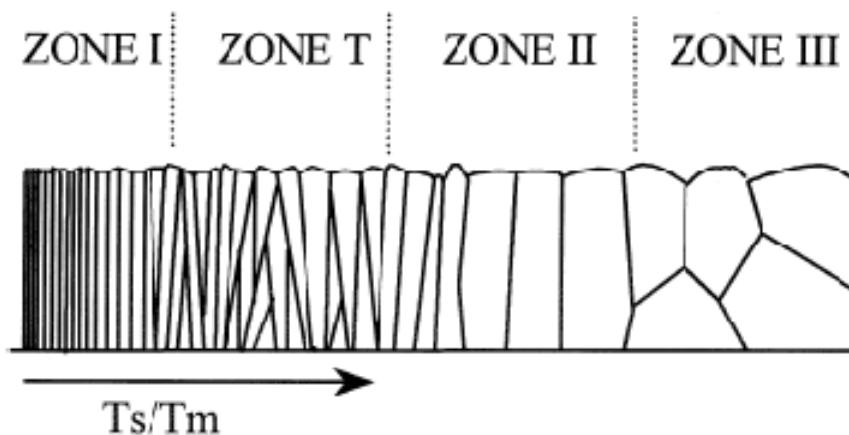
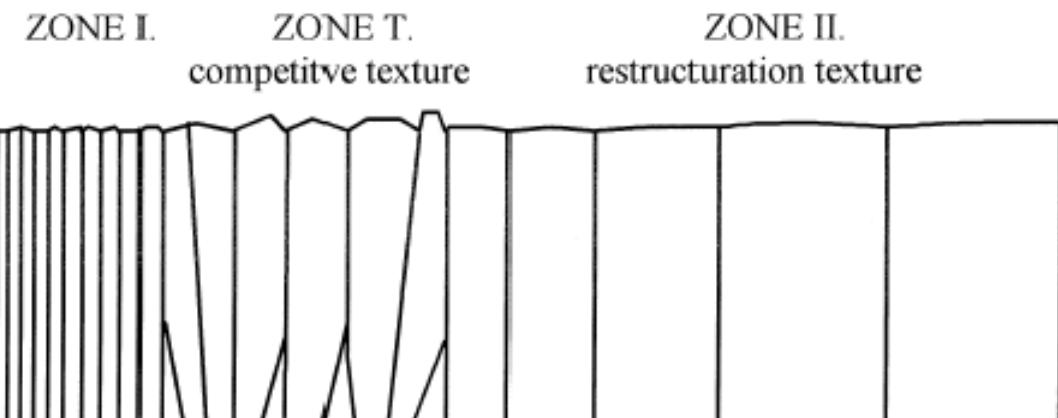


Fig. 6.3

Temperature Dependence of Thin Film Structure



thin film growth in poor vacuum (Thornton model)



thin film growth in good vacuum (Barna model)

Fig. 6.4

Epitaxy of Rutile on Hematite



Fig. 6.5

Epitaxial Relation for Rutile on Hematite

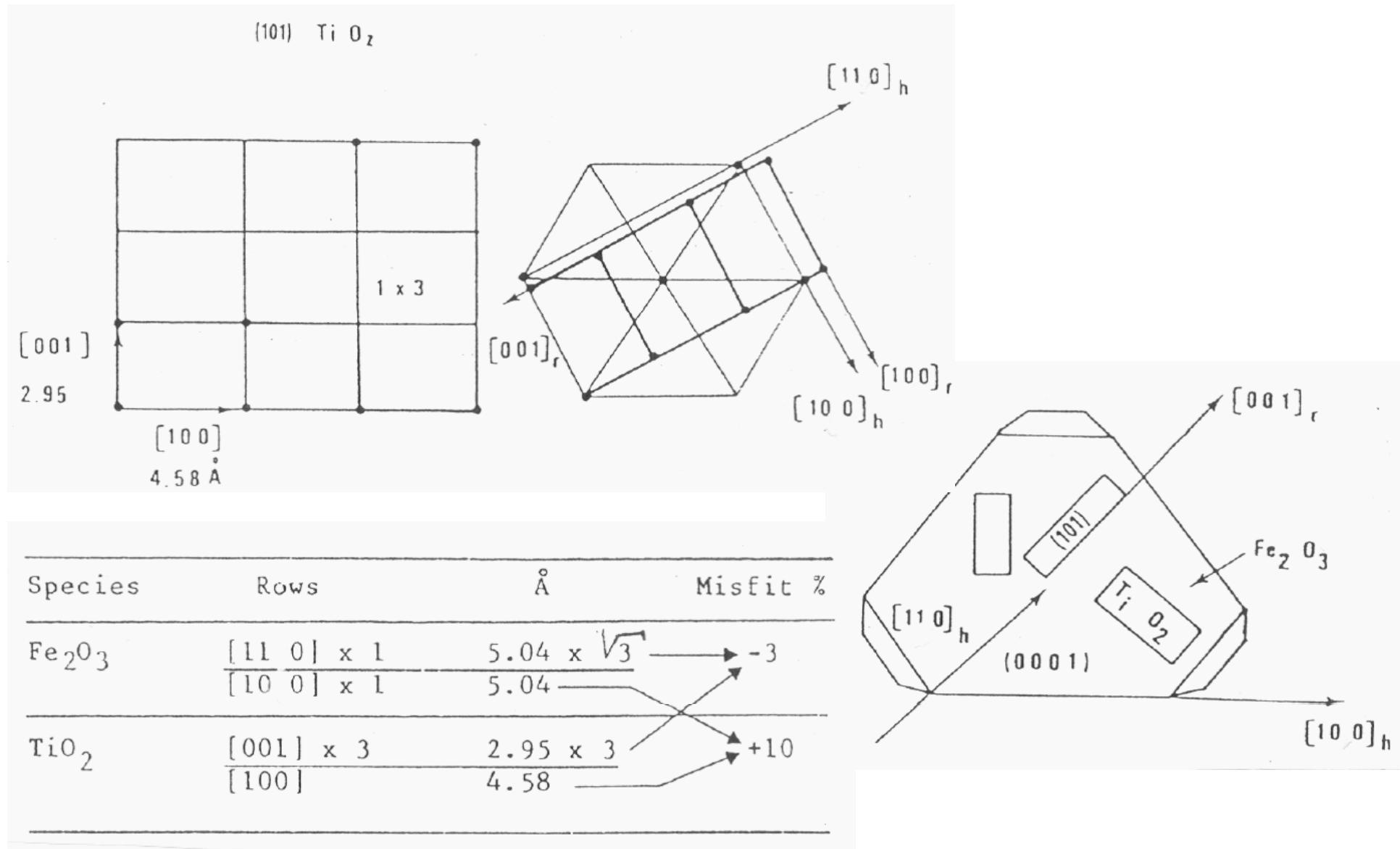


Fig. 6.6