

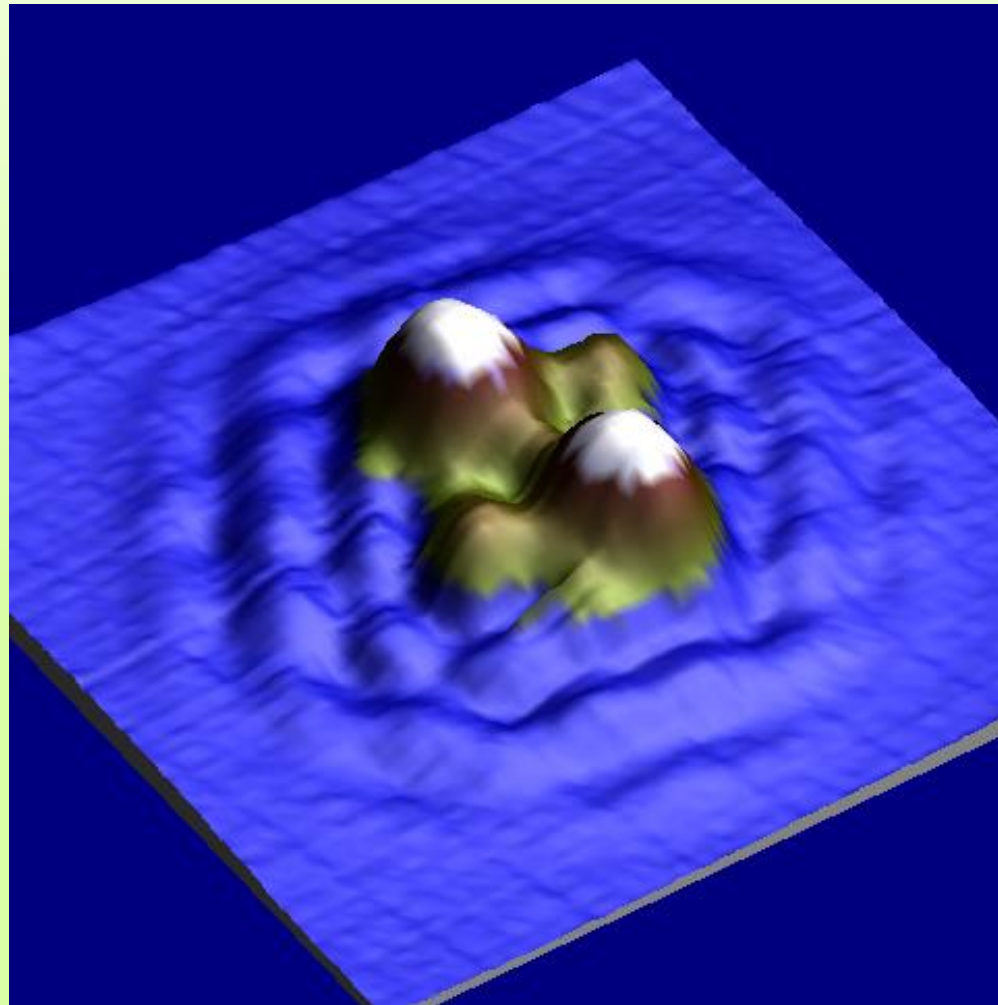
Scanning tunnelling microscopy:

Using a quantum effect to investigate quantum effects

RUB

Karina Morgenstern

Physical Chemistry I, Ruhr-Universität Bochum, Germany

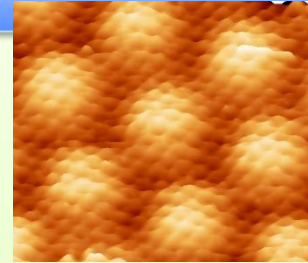


Inelastic Friedel-Like
Surface Oscillations
Nano Lett. 11, 2720
(2011)

Outline

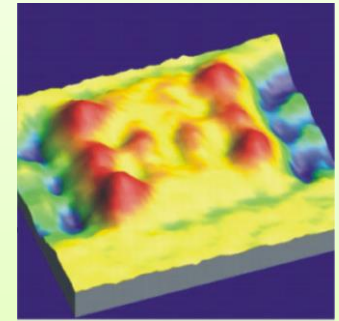
1) STM, STS, IETS, and manipulation

An extended introduction



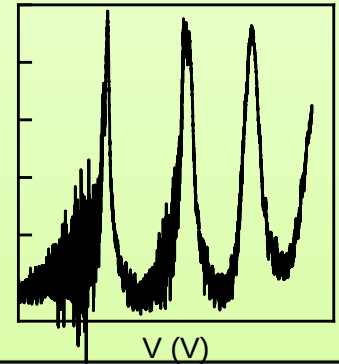
2) Quantum effects on surface imaged by STM

Electron in a box I



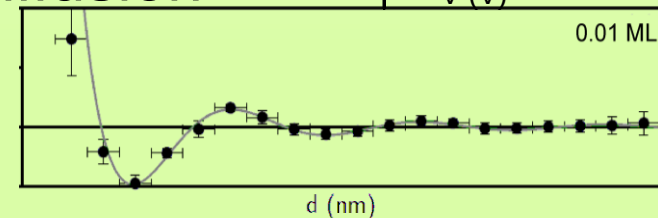
3) Quantum effects on surfaces measured by STS

Electron in a box II



4) Electron interference effects surface diffusion

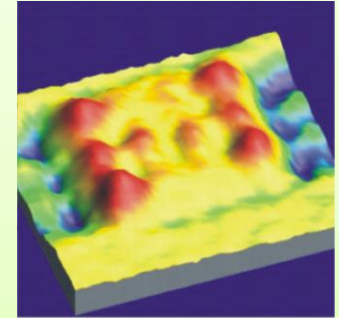
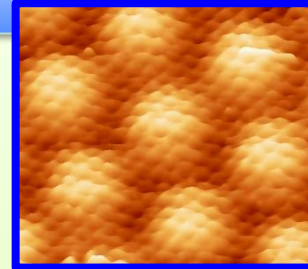
Influence of Friedel oscillations on diffusivity



Outline

1) STM, STS, IETS, and manipulation

An extended introduction

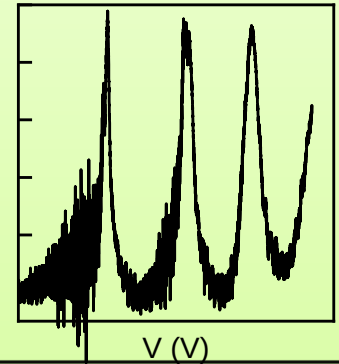


2) Quantum effects on surface imaged by STM

Electron in a box I

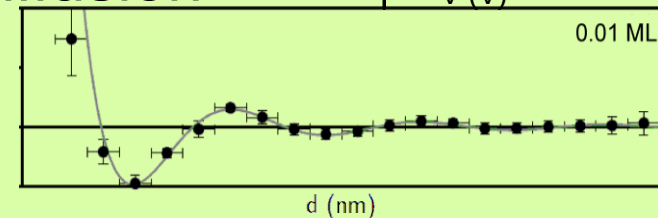
3) Quantum effects on surfaces measured by STS

Electron in a box II

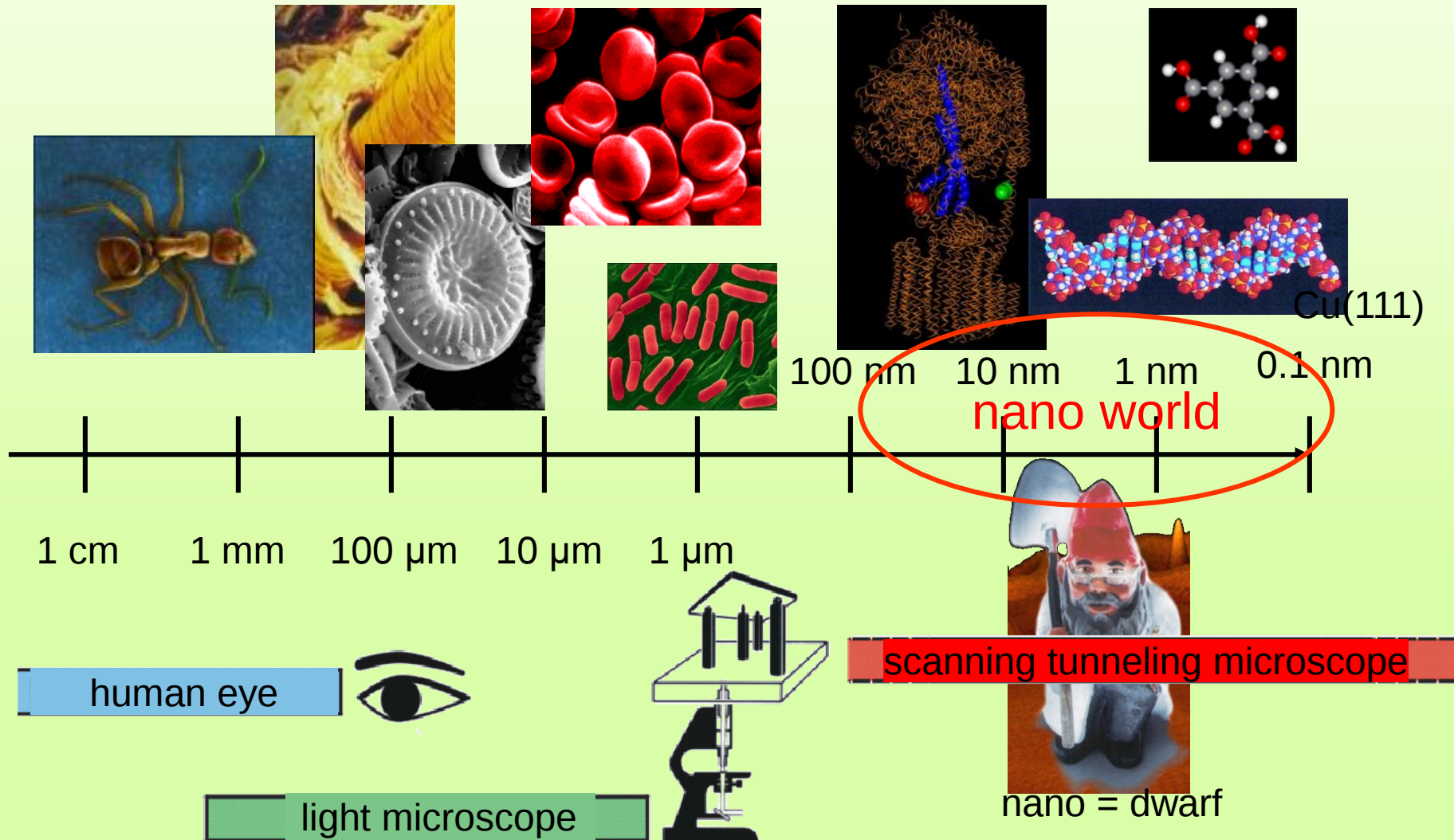


4) Electron interference effects surface diffusion

Influence of Friedel oscillations on diffusivity

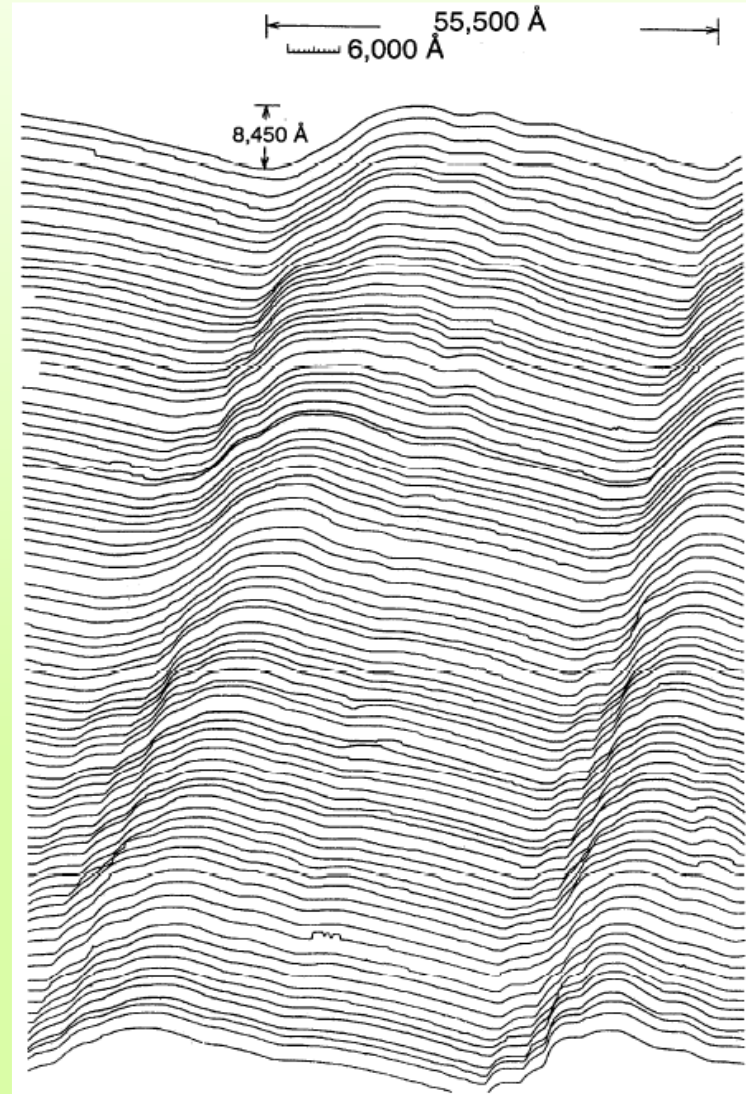
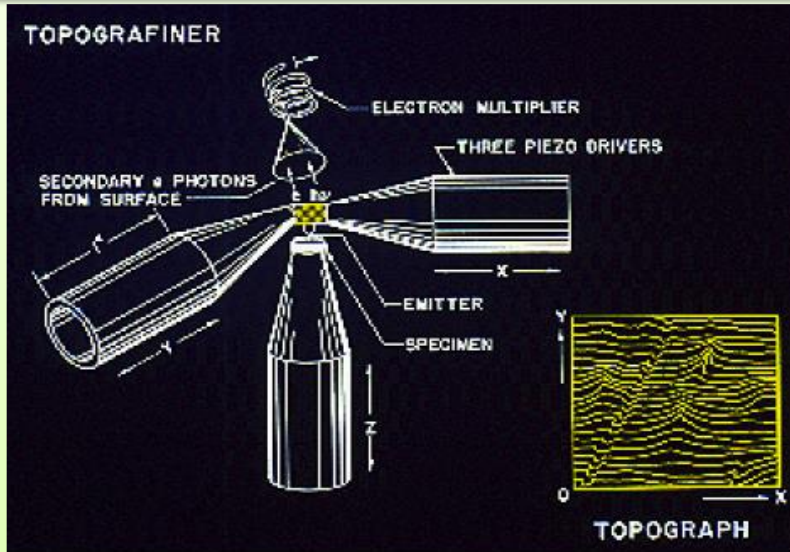


Methods for imaging the nanoworld



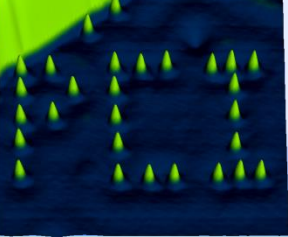
Topographiner

RUB



R. Young, J. Ward, F. Scire, Rev. Sci. Instrum. 43, 999 (1972)

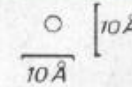
tip-sample distance: 20 nm



Scanning tunnelling microscopy

RUB

planar Au surface



1981



Press Release: The 1986 Nobel Prize in Physics

KUNGL. VETENSKAPSAKADEMIEN
THE ROYAL SWEDISH ACADEMY OF SCIENCES

15 October 1986

The Royal Swedish Academy of Sciences has decided to award the 1986 Nobel Prize in Physics by one half to

Professor, **Ernst Ruska**, Fritz-Haber-Institut der Max-Planck-Gesellschaft, Berlin, Federal Republic of Germany, **for his fundamental work in electron optics, and for the design of the first electron microscope**

and the other half, jointly to

Dr **Gerd Binnig** and Dr **Heinrich Rohrer**, IBM Research Laboratory, Zurich, Switzerland, **for their design of the scanning tunnelling microscope.**



The Nobel Prize in Physics 1986

"for his fundamental work in electron optics, and for the design of the first electron microscope" "for their design of the scanning tunneling microscope"



Ernst Ruska



Federal Republic of Germany

Fritz-Haber-Institut der Max-Planck-Gesellschaft
Berlin, Germany

1906 - 1988



Gerd Binnig



Federal Republic of Germany

IBM Zurich Research Laboratory
Rüschlikon, Switzerland

1947 -



Heinrich Rohrer

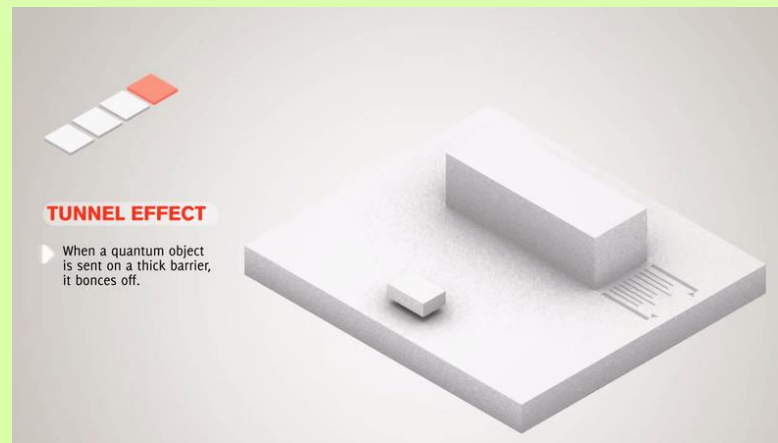
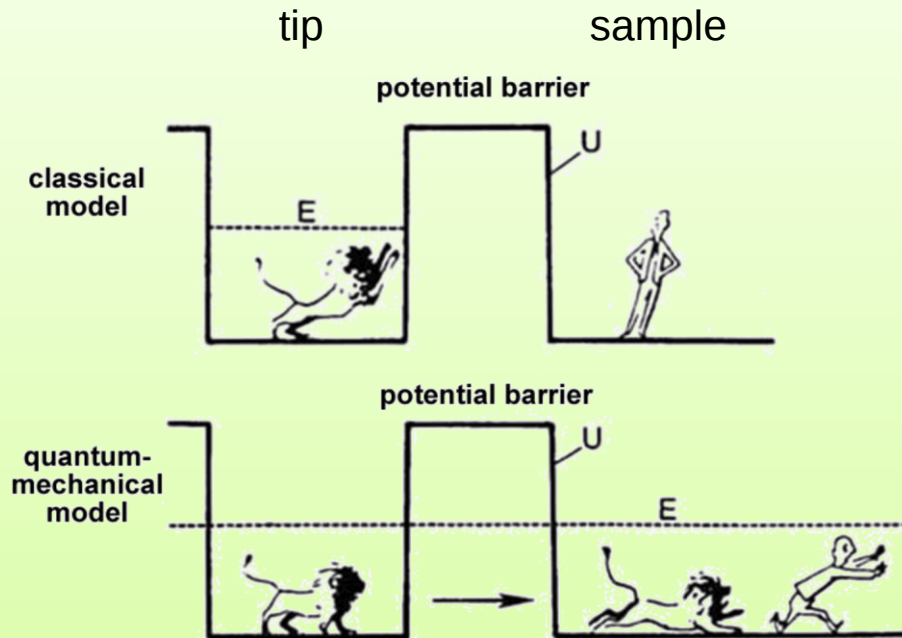


Switzerland

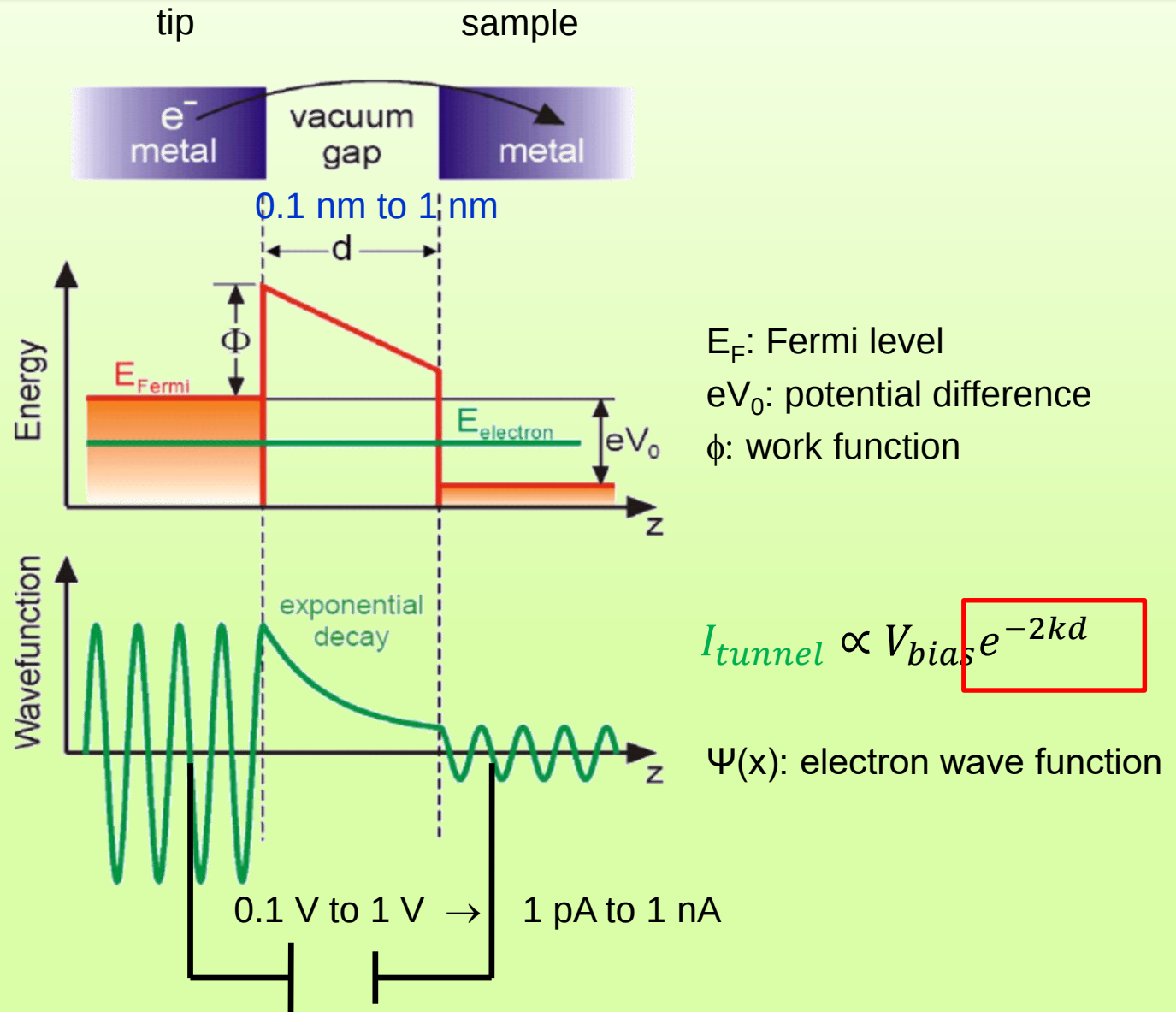
IBM Zurich Research Laboratory
Rüschlikon, Switzerland

1933 -

Scanning tunneling microscopy



Scanning tunneling microscopy



Scanning tunneling microscopy

RUB



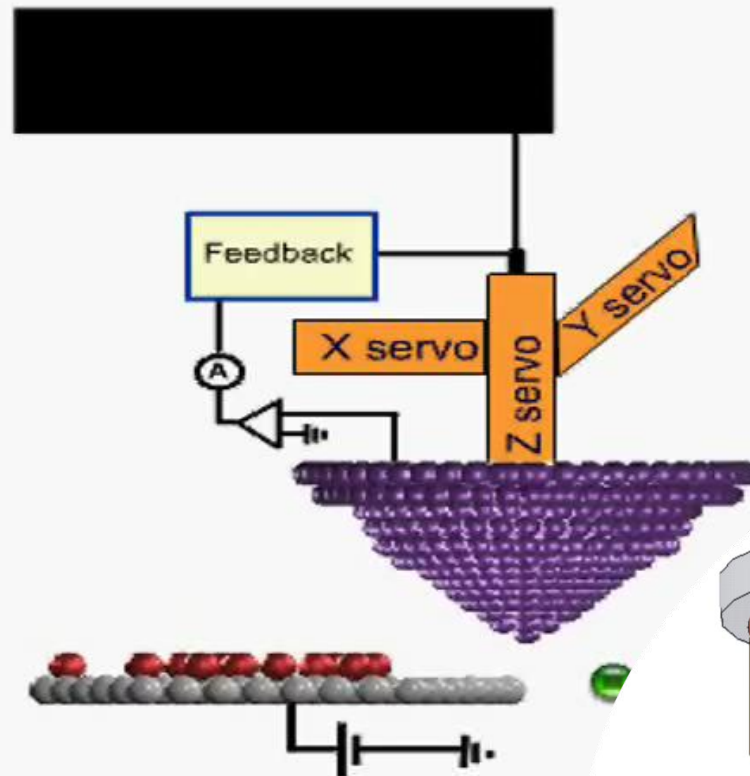
constant
current
mode

Pentacene/Au/Mica

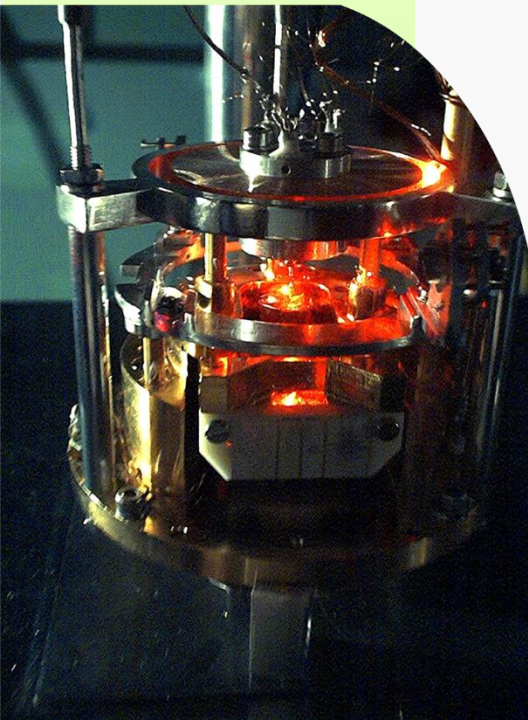
Lars Ruppel, PCI, R.U.B.

Scanning tunneling microscopy

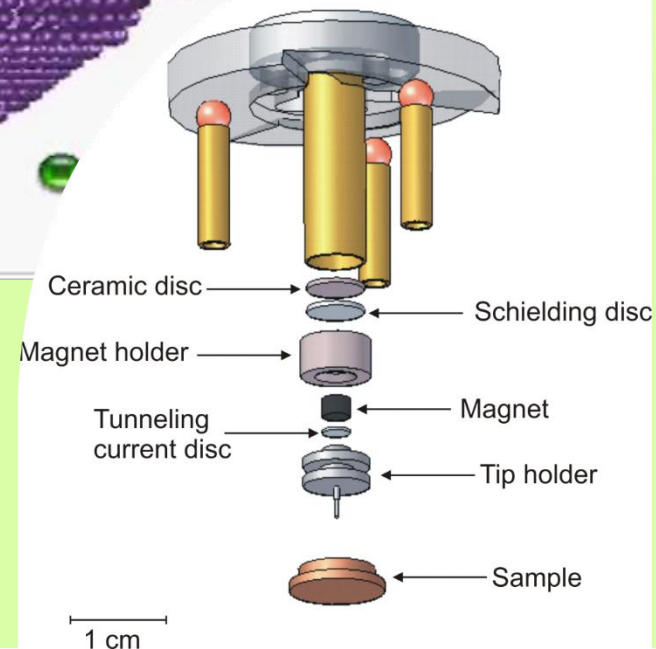
RUB



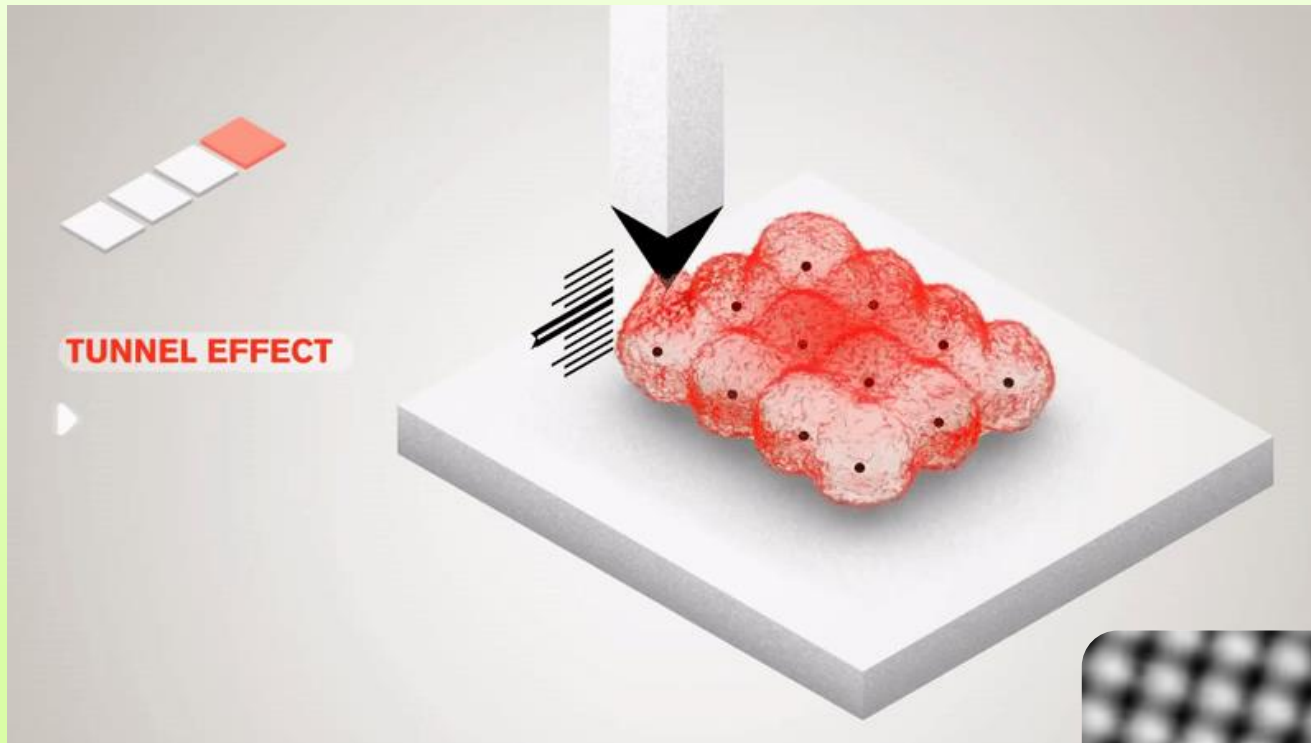
constant
current
mode



imaging at ~ 5 K
in ultra-high vacuum

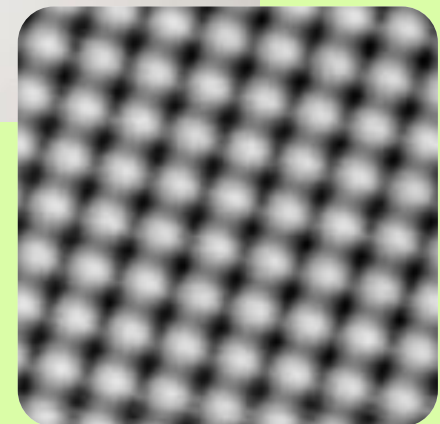


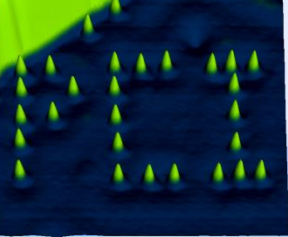
Tunneling effect in STM



constant
height mode

Ag(100)

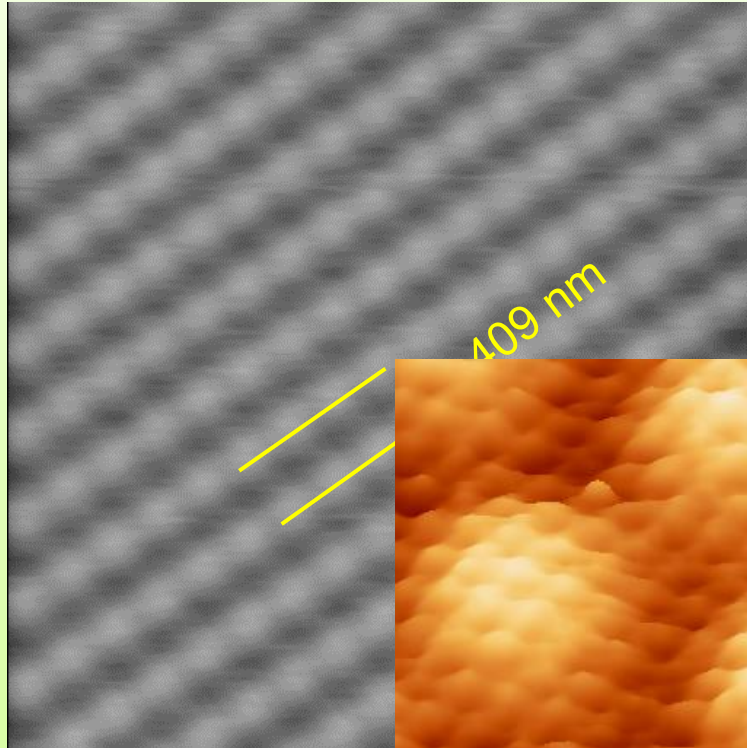




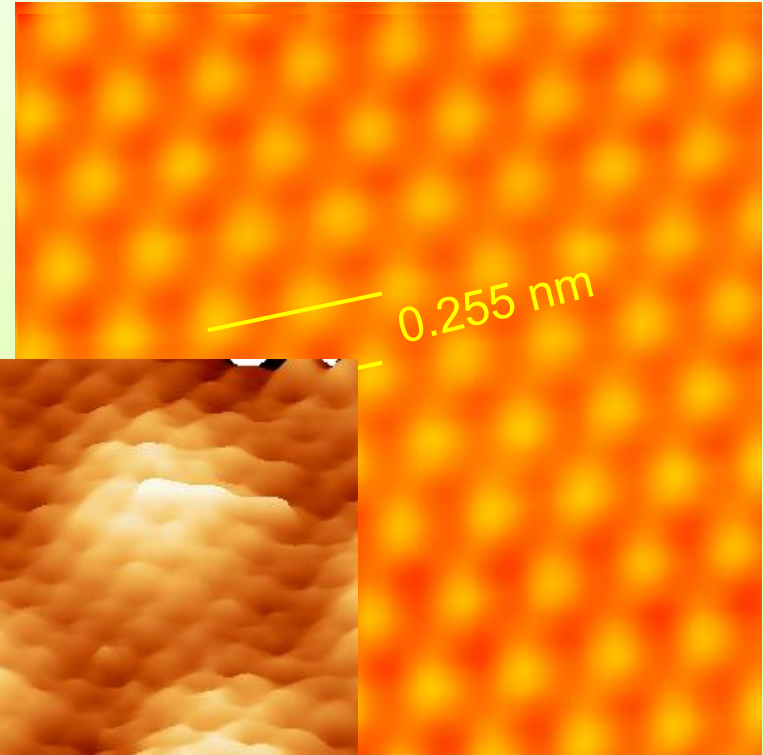
STM: Atomic resolution

RUB

Ag(110)

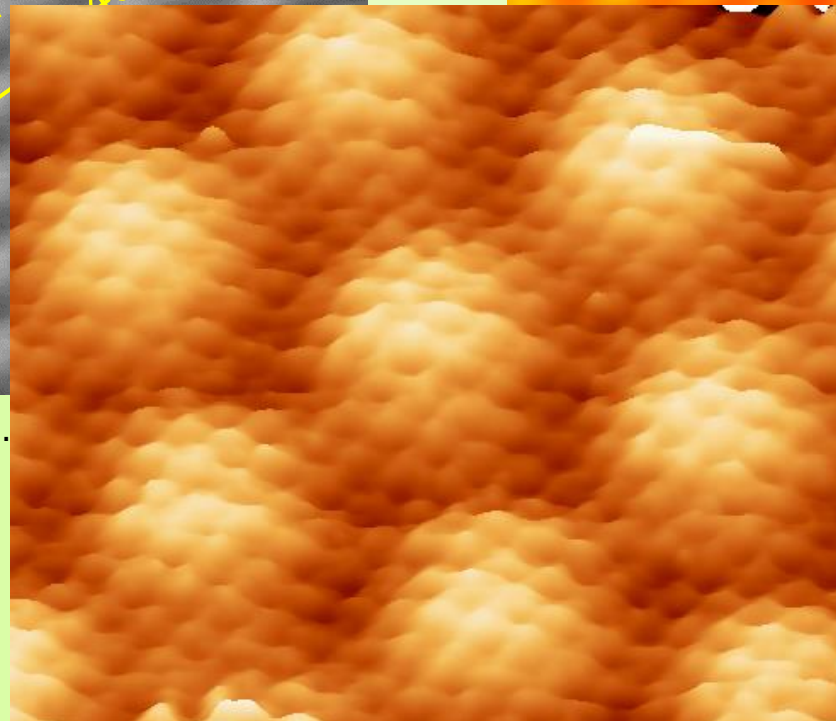


Cu(111)



25 mV, 30 pA, 5 K

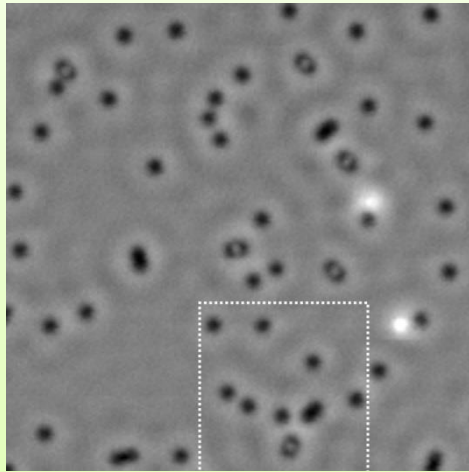
0.



moiré pattern of ZnO/Ag(111)

J. Phys. Chem. C 126, 12500 (2022)

STM: Images of molecules

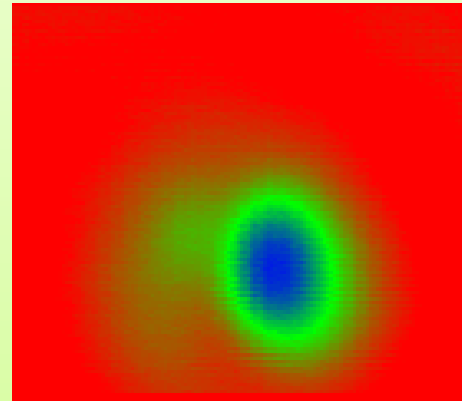


CO / Cu(111)

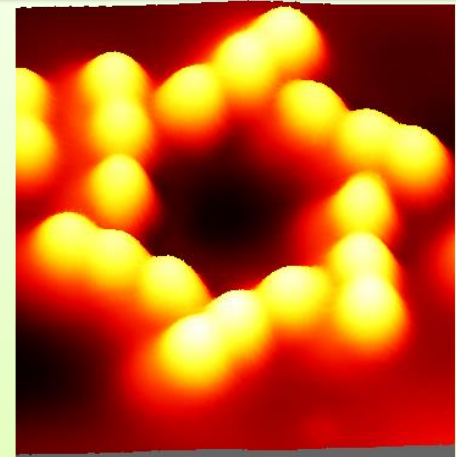
Where are the atoms,
what do we see?



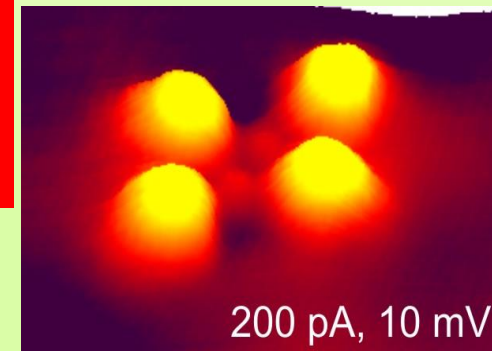
9 water molecules
/ Ag(111)



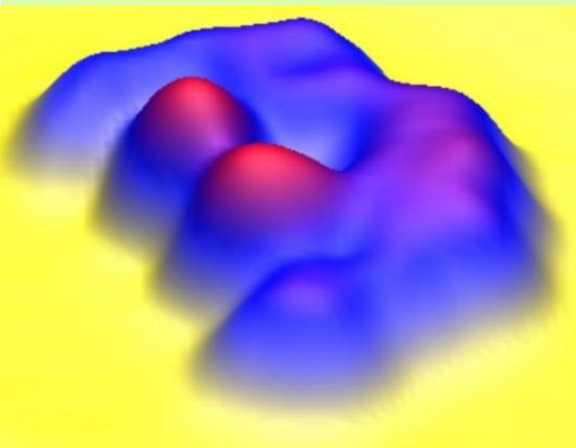
one chloronitro-
benzene / Au(111)



six anilino-amino-
azobenzene / Au(111)

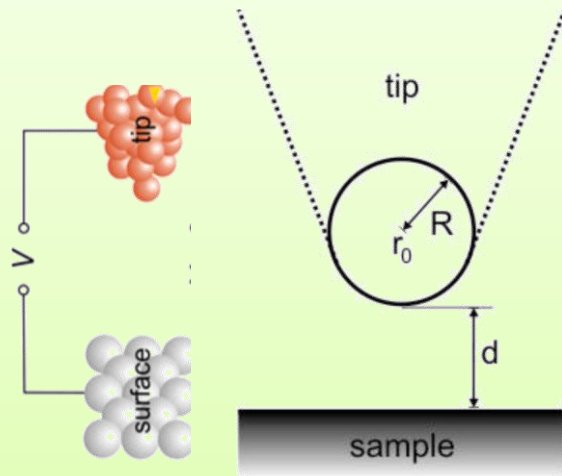


two amino-nitro-
azobenzene / Au(111)



one azobenzene derivative solvated by water / Ag(111)

Tersoff-Hamann theory



Where are the atoms,
what do we see?

$$I(r_0, V) = \frac{2\pi}{\hbar} \int_{E_F}^{E_F + eV} dE \rho_s(r_0, E) T(r_0, E, eV) \rho_t(r_0, E - eV)$$

J. Tersoff, D.R. Hamann

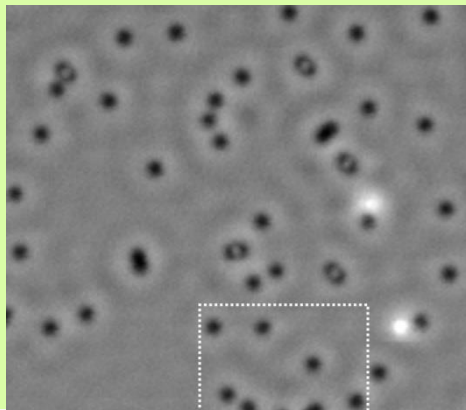
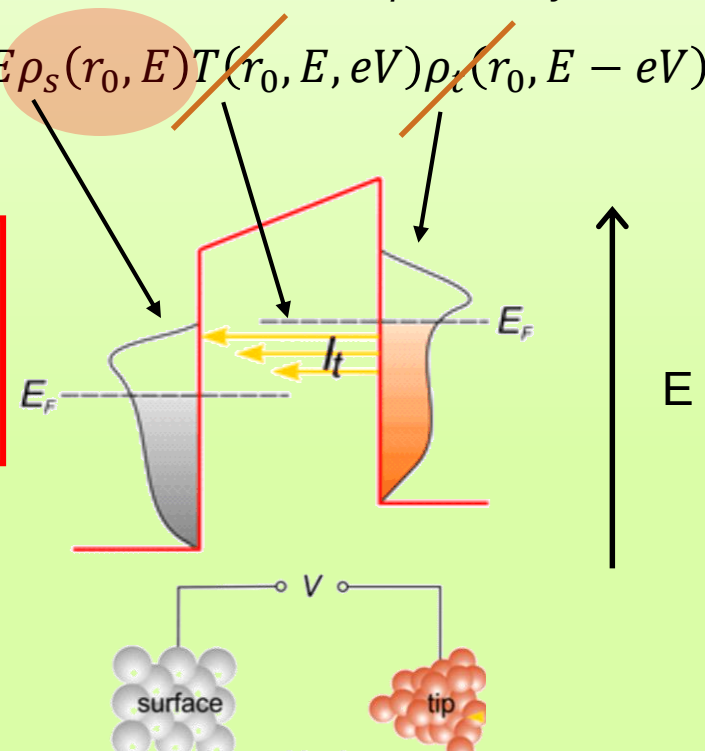
Phys. Rev. Lett. 50,1998 (1983)

based on J. Bardeen

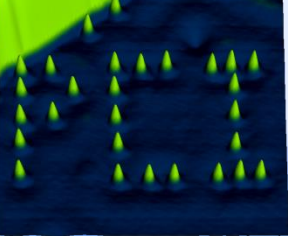
Phys. Rev. Lett. 6, 57 (1961)

ρ : density-of-states

⇒ STM measures the local density of states of the sample from the Fermi level up to the bias voltage energy at the tip position



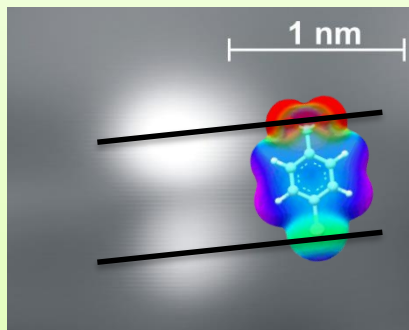
CO / Cu(111)



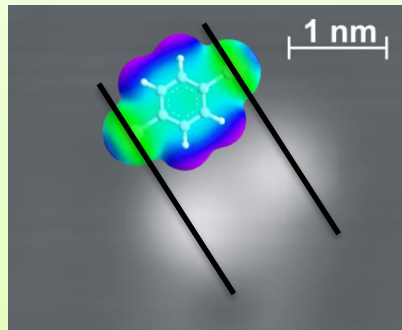
STM: Images of molecules

RUB

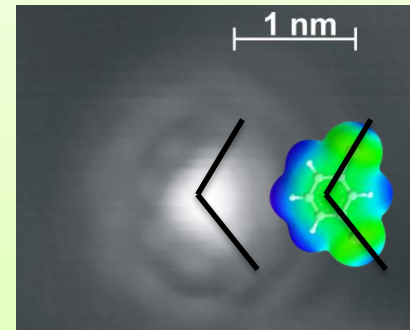
chloronitrobenzene
para



dichlorobenzene
para

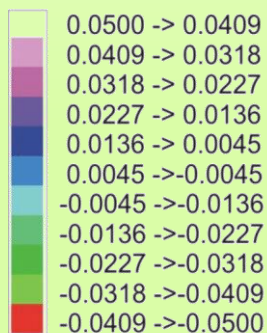


meta

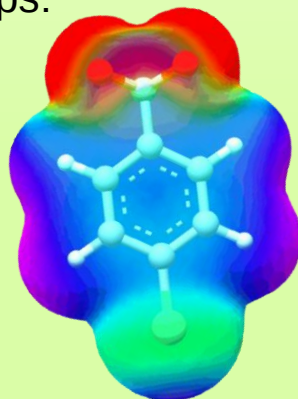


⇒ STM images: Convolution of LDOS with topography

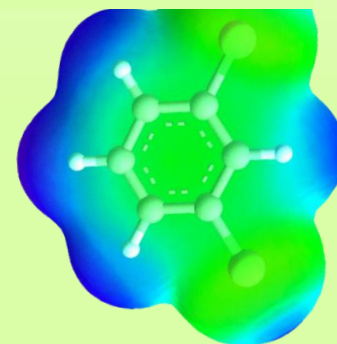
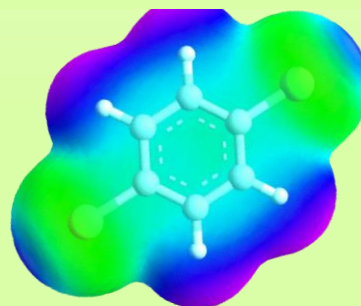
electrostatic potential maps:



Hartree



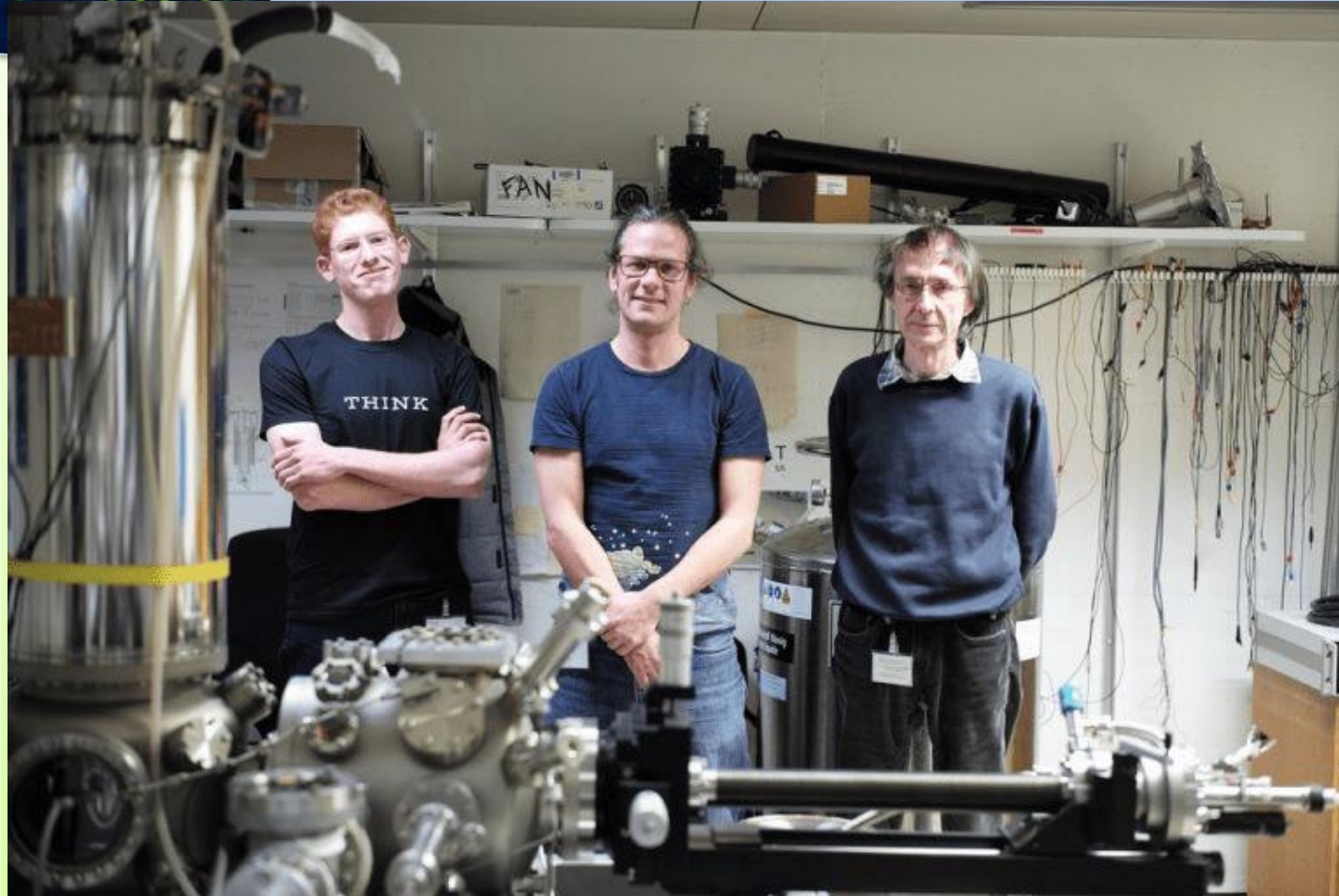
⊗ Pauli repulsion (for molecular tips)



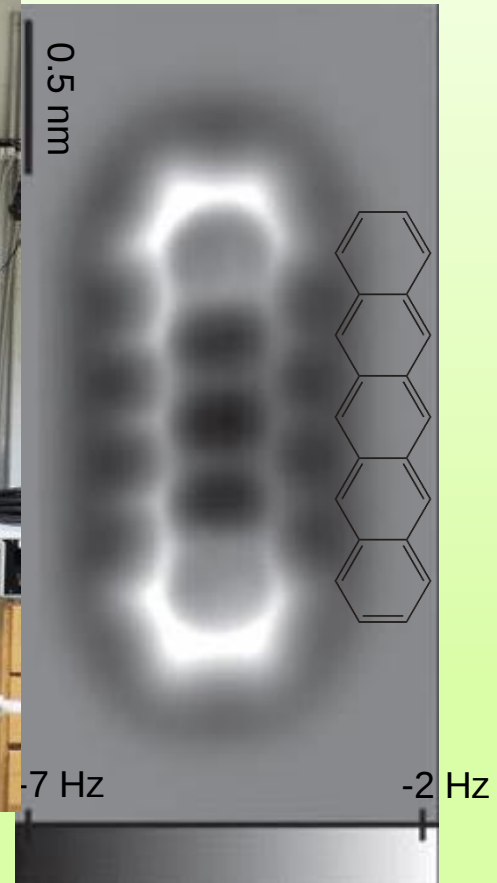
evaporation @ 17 K
measurement @ 5 K

Bond pictures

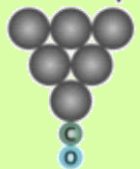
RUB



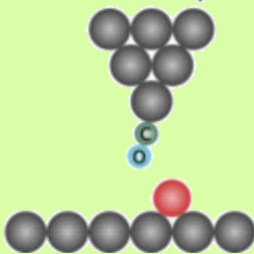
AFM - CO tip



AFM tip



AFM tip



Temirov et al.
New J. Phys. 10, 053012 (2008)

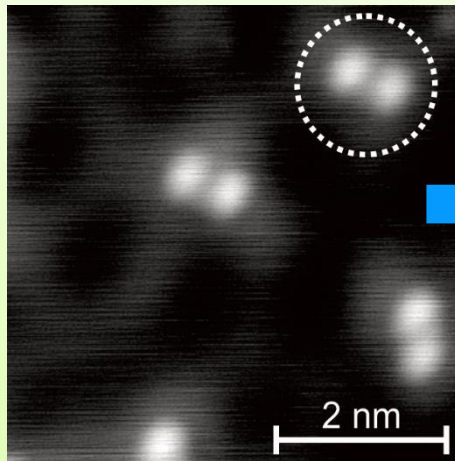
Jelínek
J. Phys.: Condens. Matter
29, 343002 (2017)

CH: 170mV, 110 pA

Pentacene
Gross et al.
Science 325, 5944 (2009)

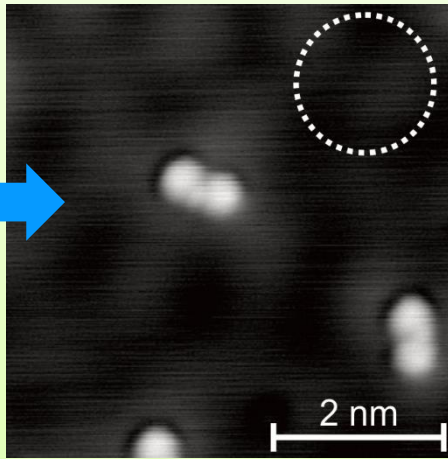
STM: Bond pictures

metallic tip

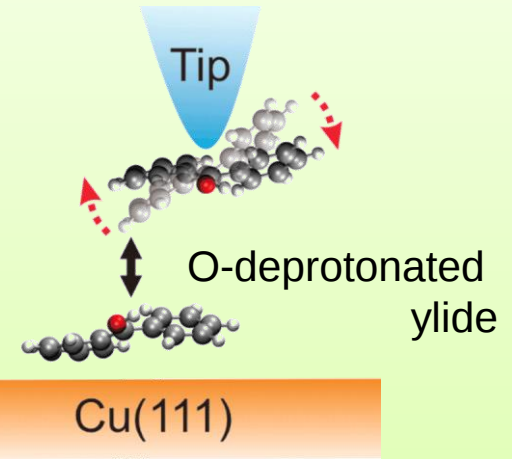


-10 mV, 10 pA

molecular tip

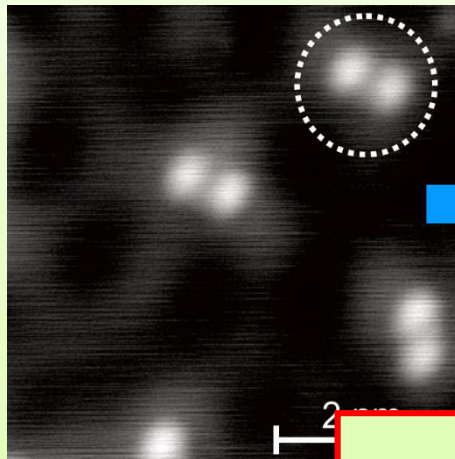


-100 mV, 10 pA



STM: Bond pictures

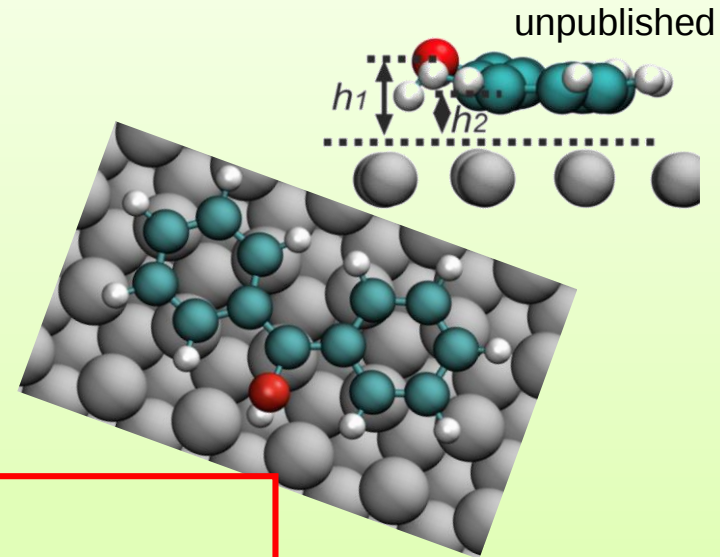
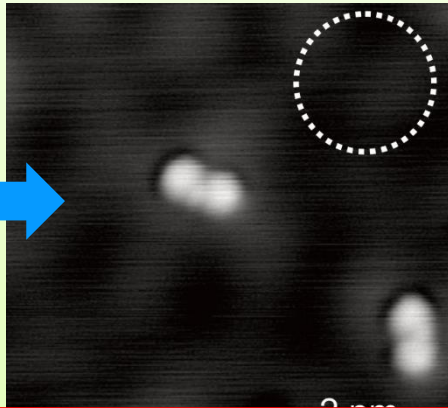
metallic tip



-10 mV, 10 pA

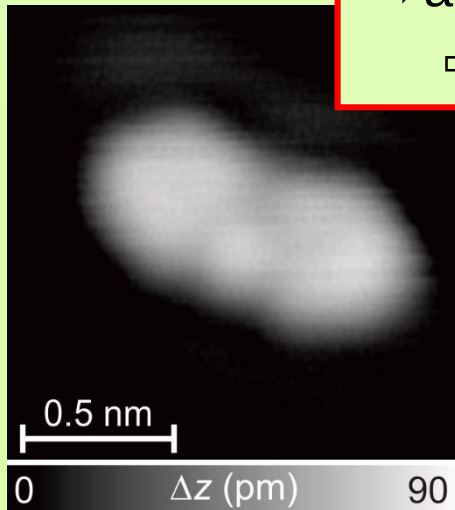
constant current

molecular tip

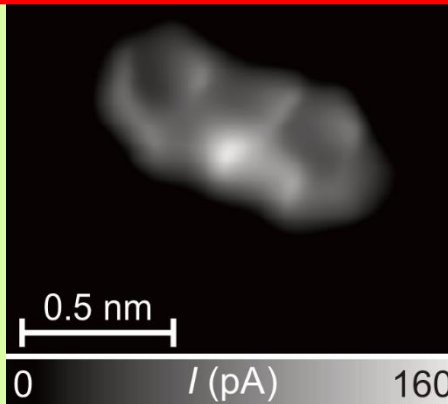


STM

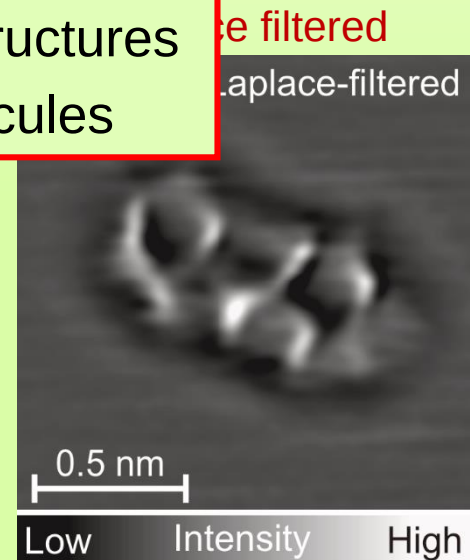
- ⇒ atomic resolution of bulk structures
- ⇒ bond resolution of molecules



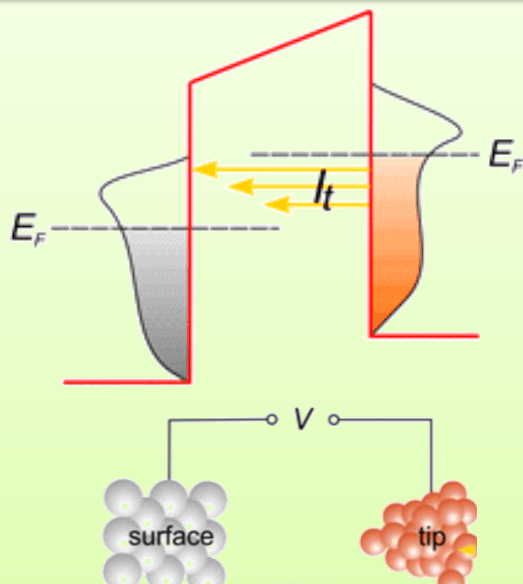
5 mV, 10 pA



5 mV, 10 pA, $\Delta z = -50$ pm



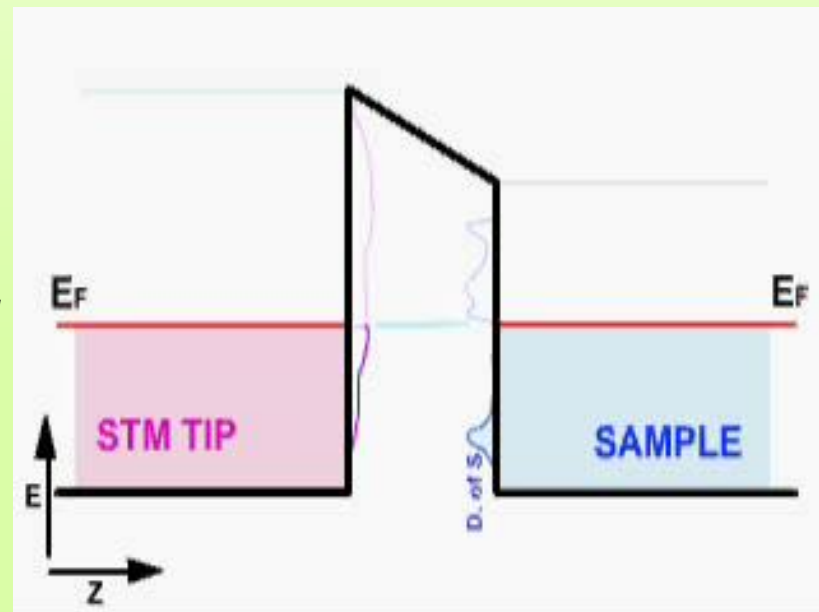
STS: Scanning Tunneling Spectroscopy

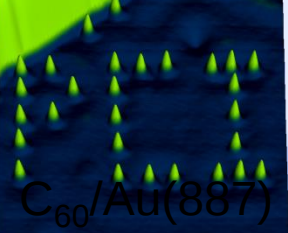


$$I(r_0, V) = \frac{2\pi}{\hbar} \int_{E_F}^{E_F + eV} dE \rho_s(r_0, E) T(r_0, E, eV) \rho_t(r_0, E)$$

$$\rightarrow \frac{dI}{dV} \propto \rho_s(r_0, E)$$

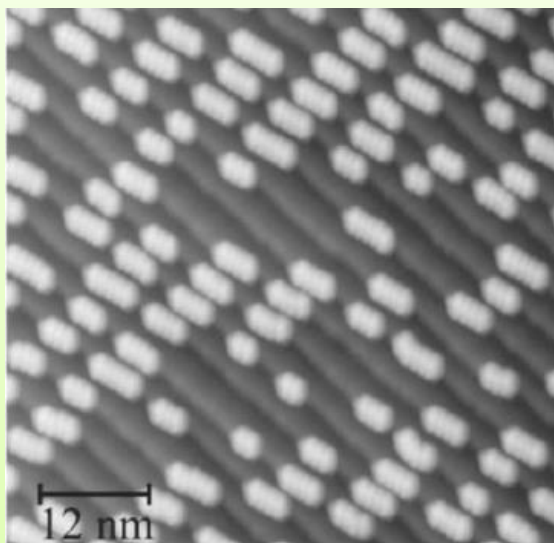
local density of states
of the sample
at the position of the tip



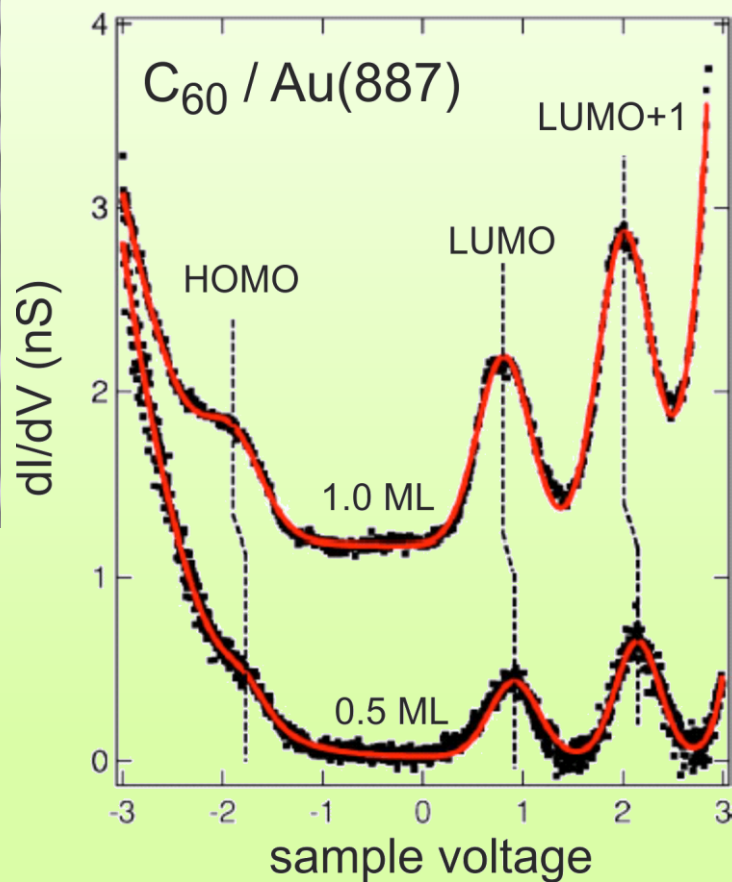


HOMO-LUMO gap

RUB

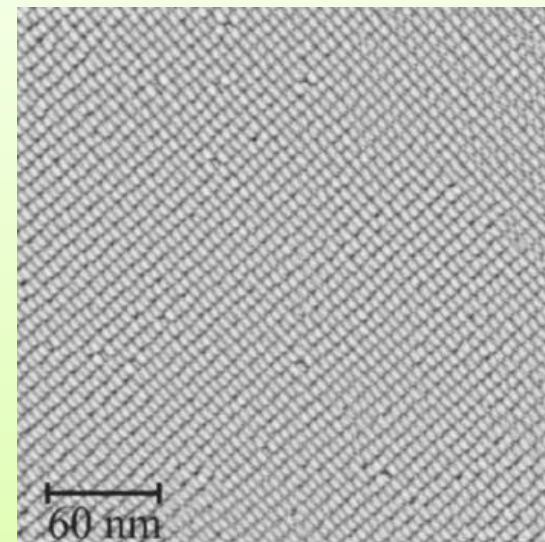


0.5 ML

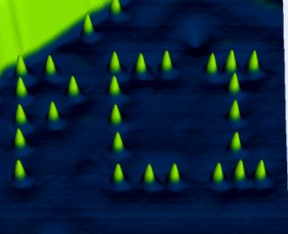


← E_F →

occupied states unoccupied states



1 ML



Identification of molecular orientation

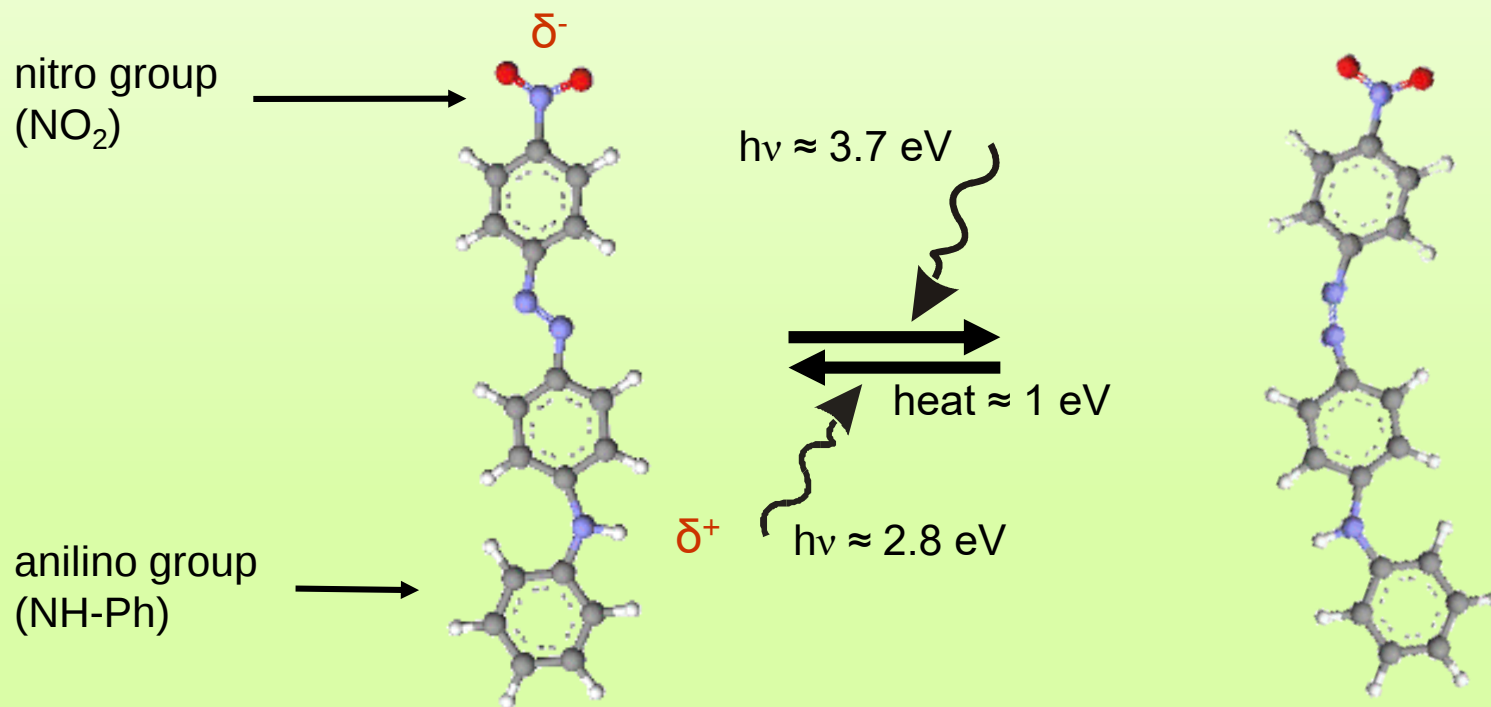
RUB

PCCP 12, 6035 (2010)

4-anilino-4'-nitro azobenzene

trans-configuration

cis-configuration

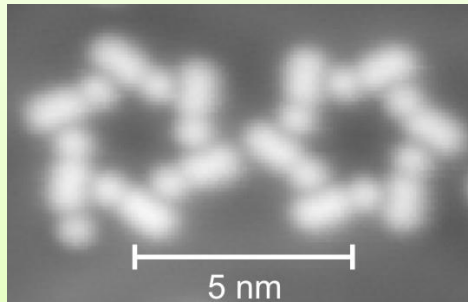


Identification of molecular orientation

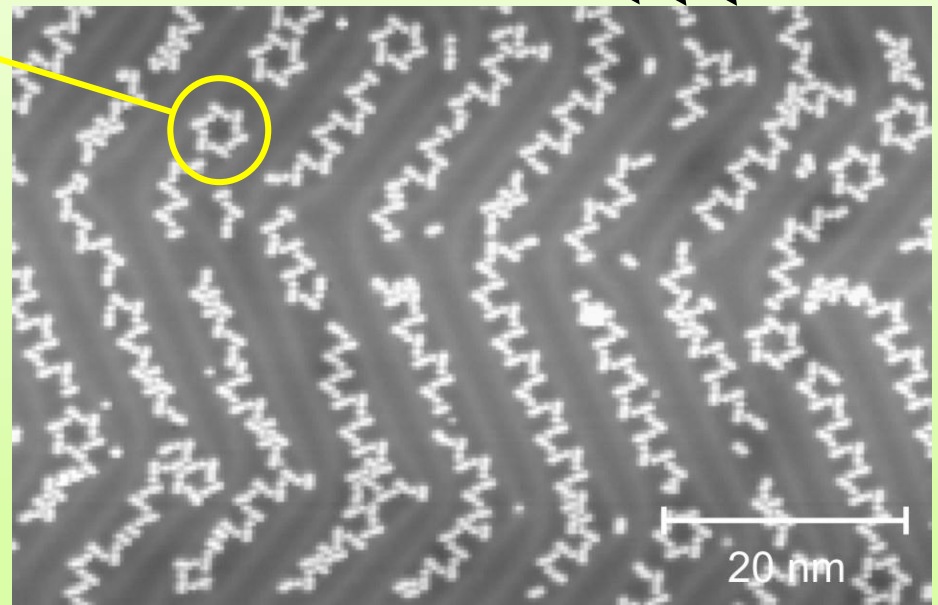
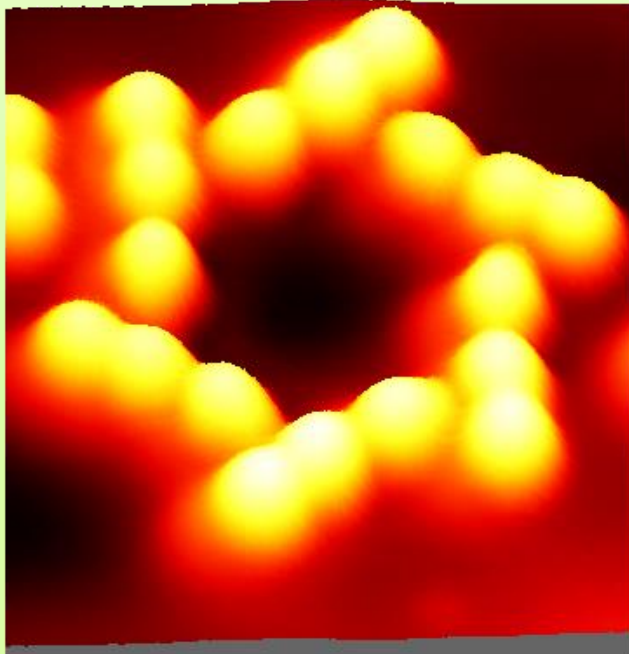
PCCP 12, 6035 (2010)

star structures

4-anilino-4'-nitro azobenzene / Au(111)



310 pA, 125 mV



20 nm

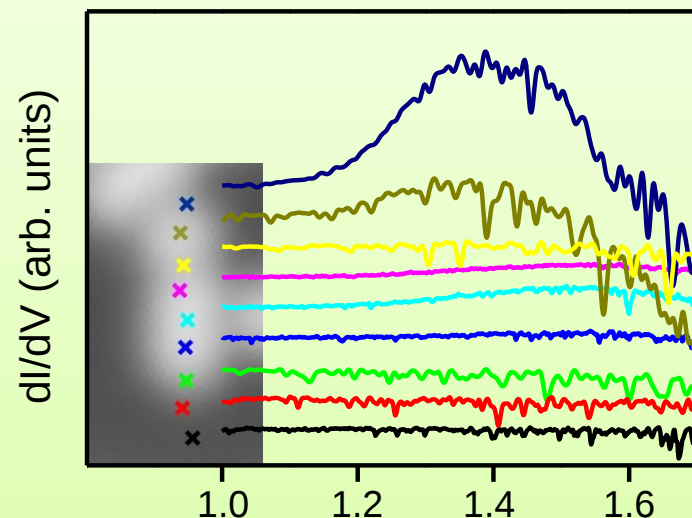
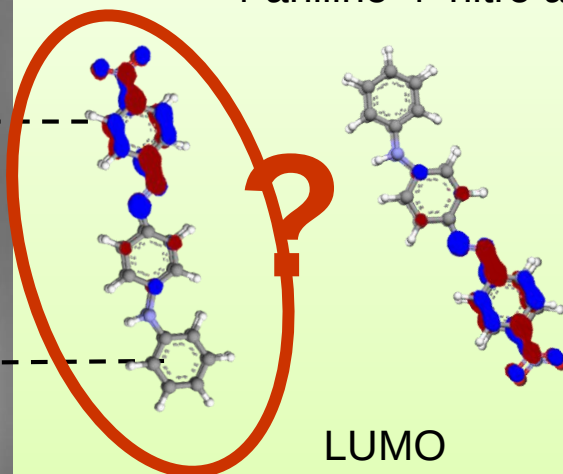
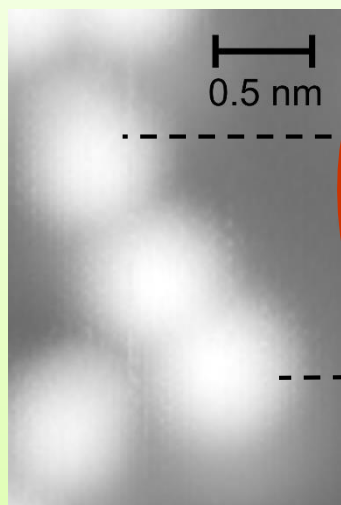
27 pA, 96 mV

hcp fcc hcp

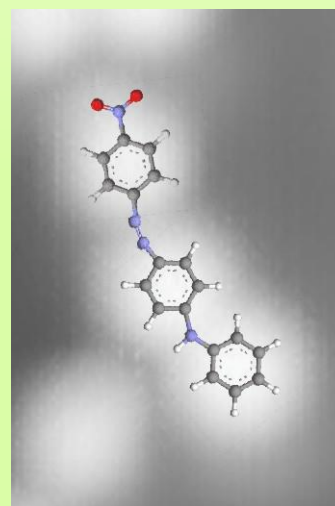
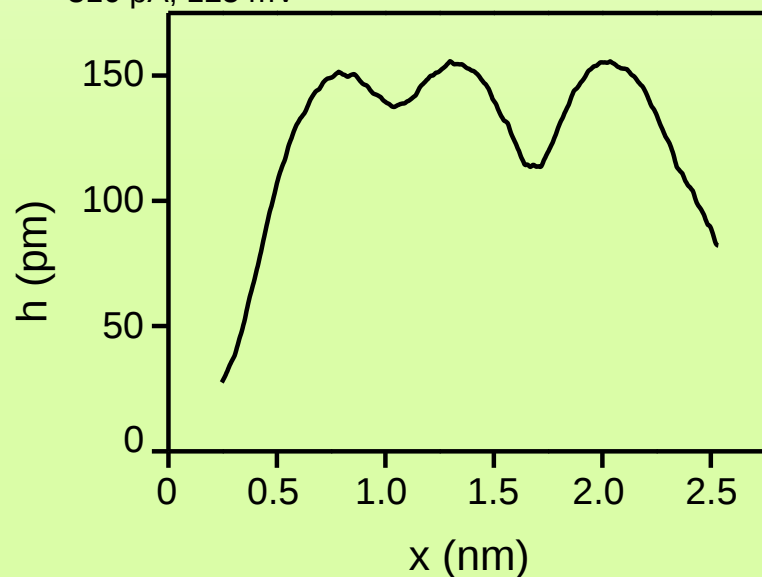
Identification of molecular orientation

4-anilino-4'-nitro azobenzene / Au(111)

PCCP 12, 6035 (2010)

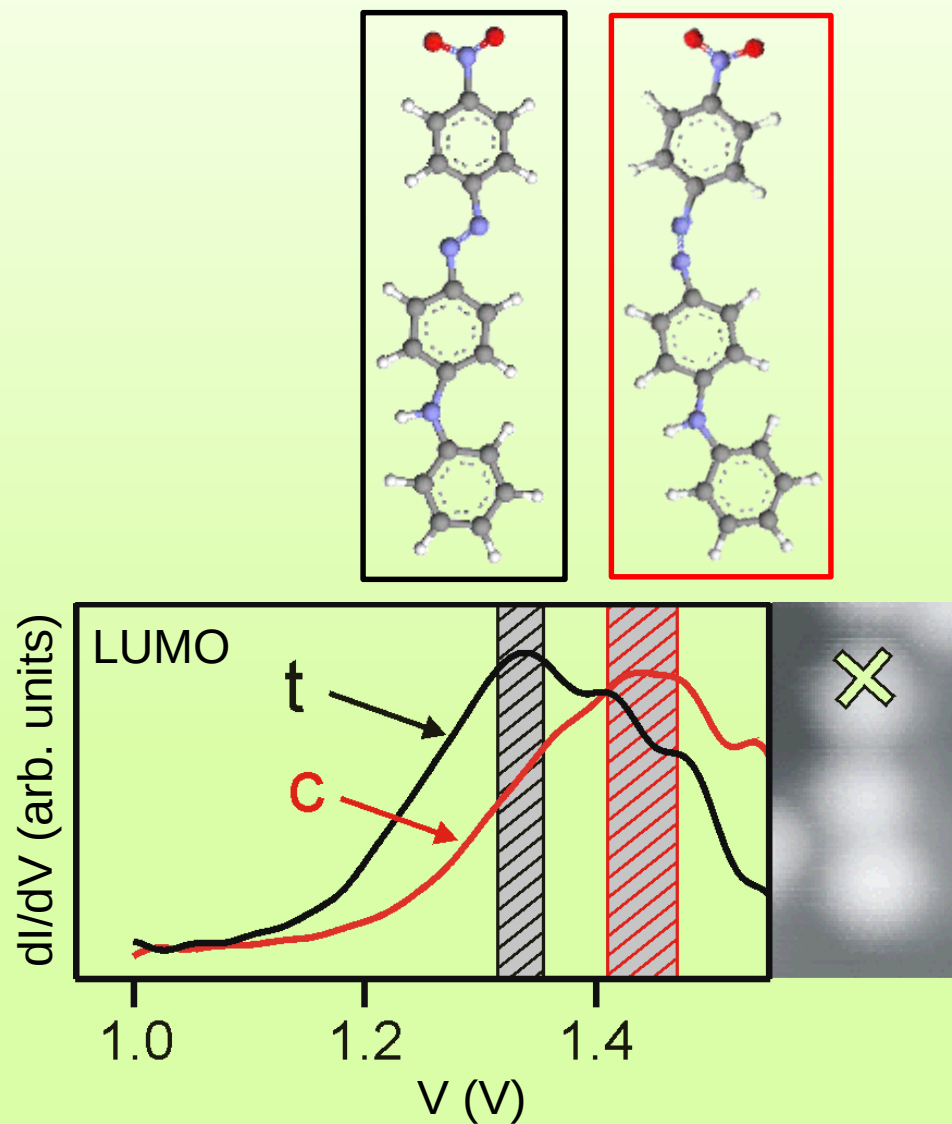


310 pA, 125 mV

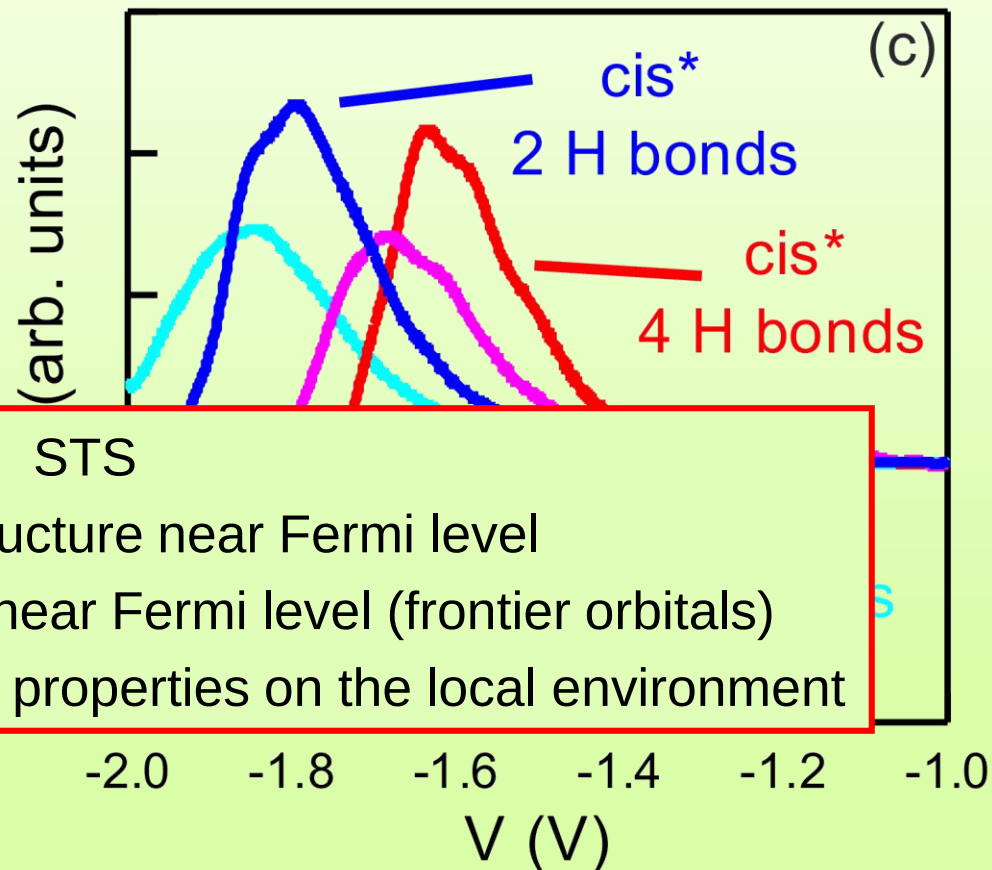
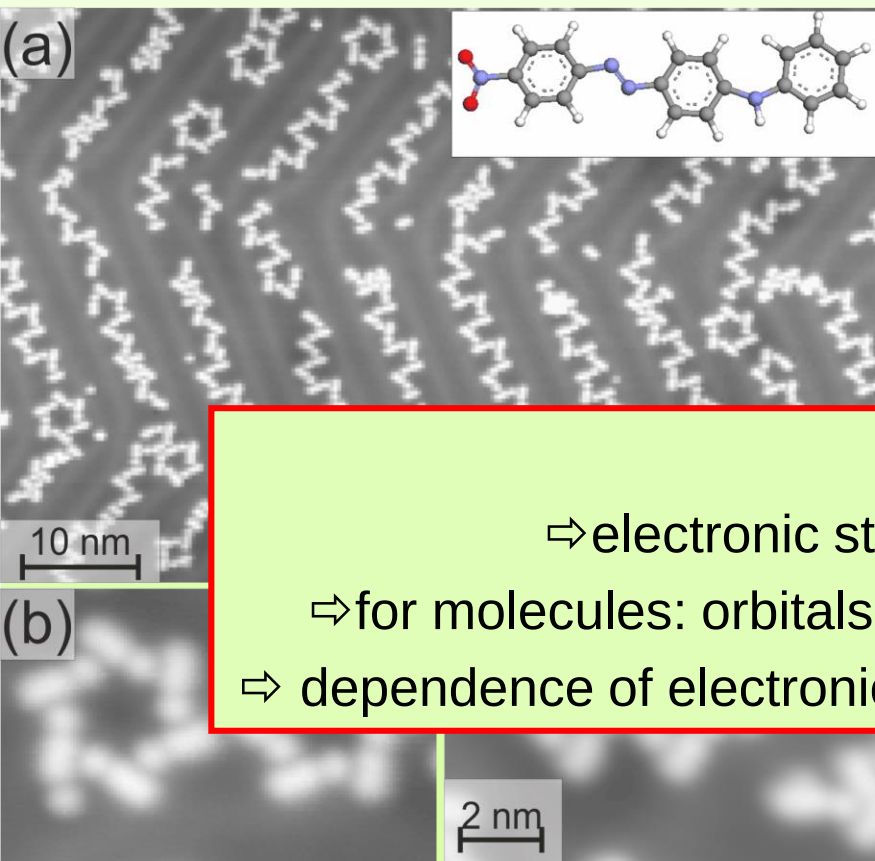


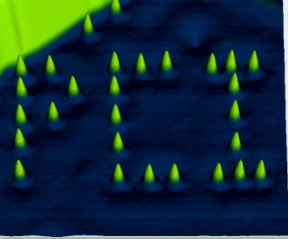
V (V) $V_{\text{mod}} = 26 \text{ mV}, 917 \text{ Hz}$

Electronic structure of conformers



Influence of bonding on HOMO



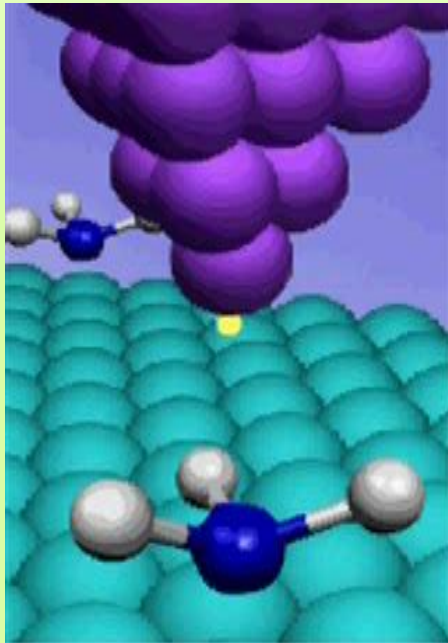
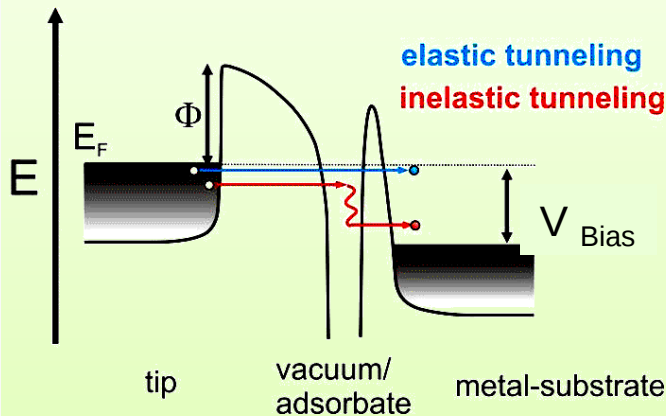


Inelastic electron tunneling spectroscopy (IETS)

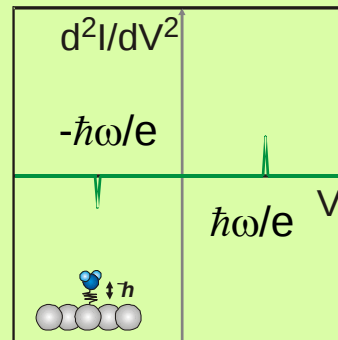
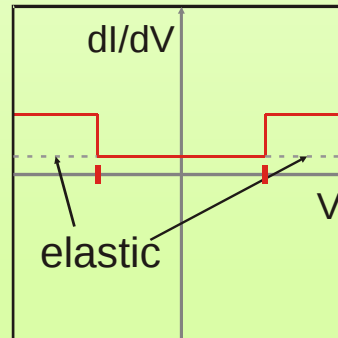
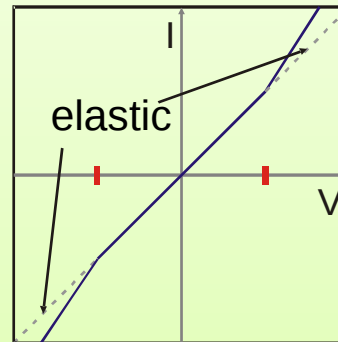
RUB

Wilson Ho

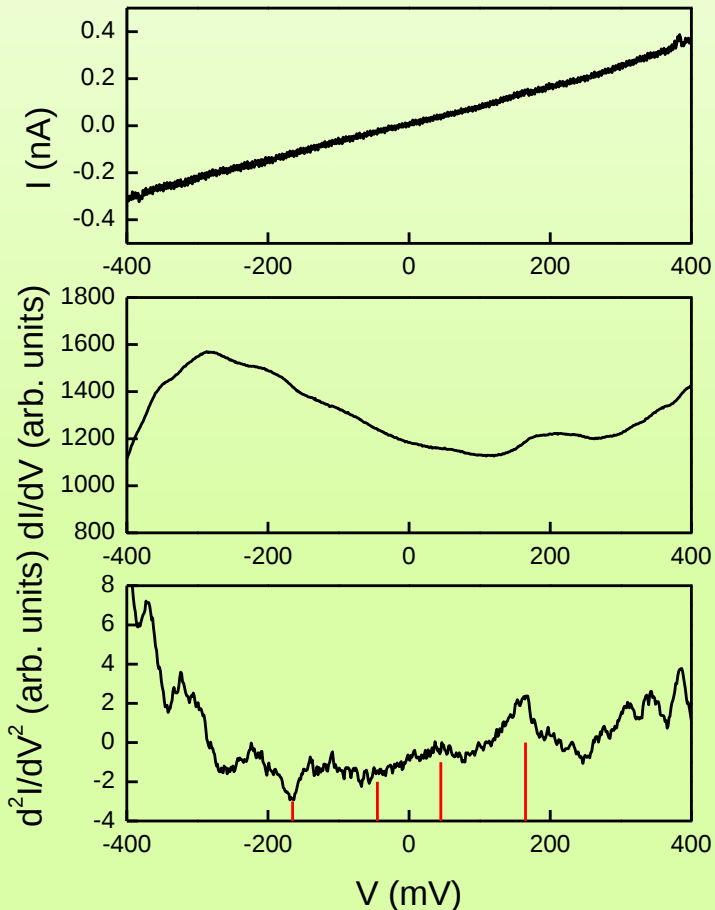
J. Chem. Phys. 117, 11033 (2002)



with courtesy of I. J. Pascual



$H_2O / Cu(111)$

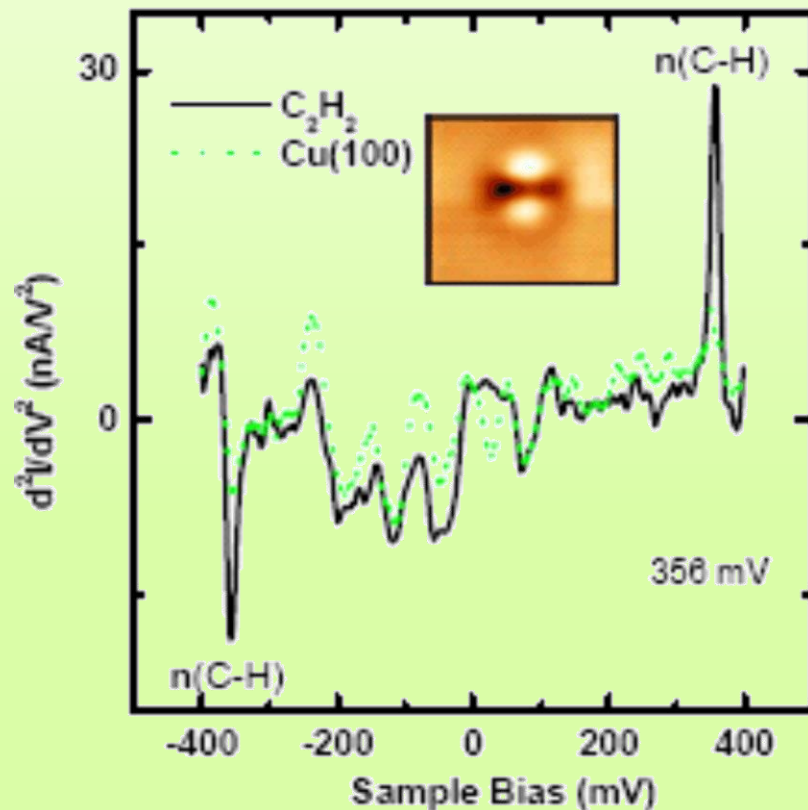


J. Chem. Phys. 116, 5746 (2002)

IETS: Internal and external molecular modes

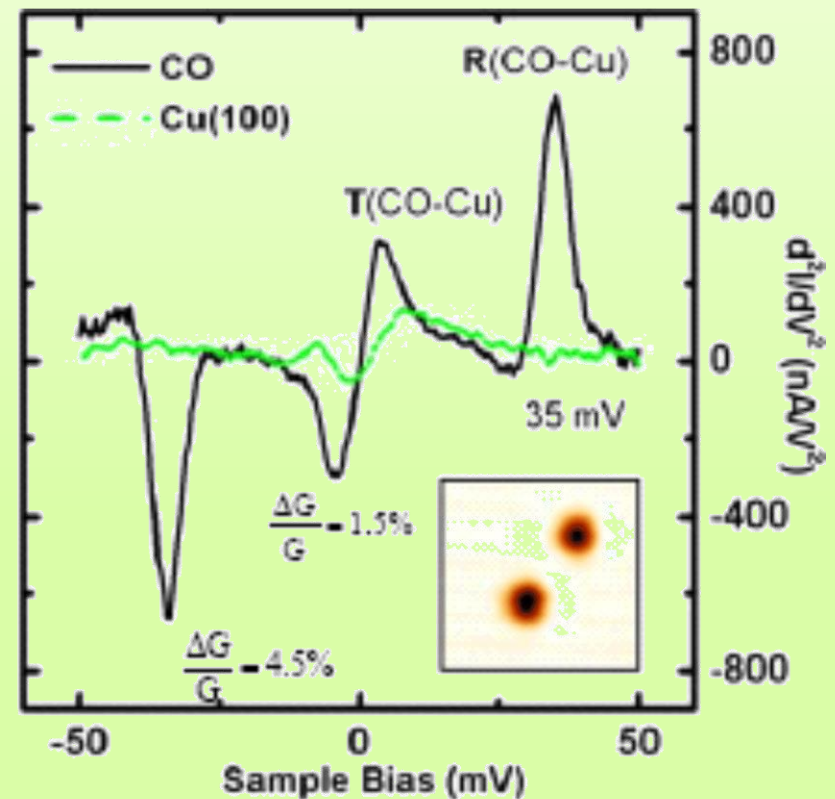
Intramolecular
vibrational modes

C_2H_2 / Cu(100)



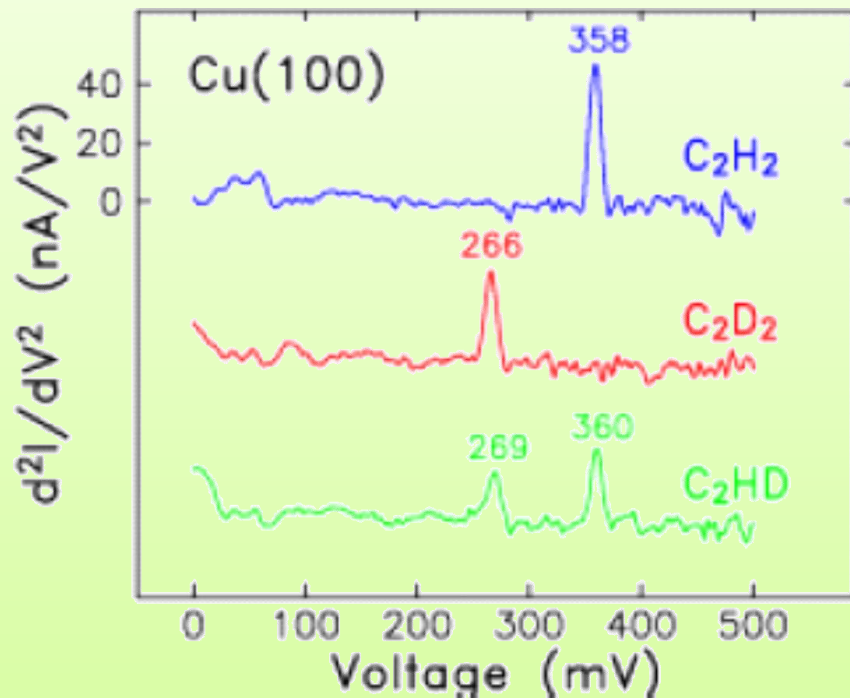
Adsorbate-surface
vibrational modes

CO / Cu(100)



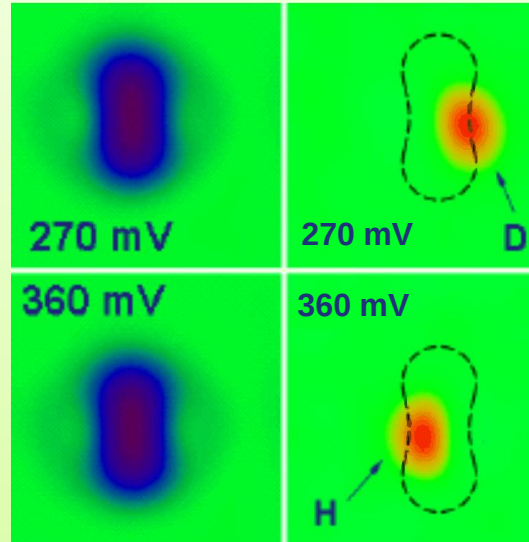
IETS: Isomer identification

Wilson Ho, UC Irvine

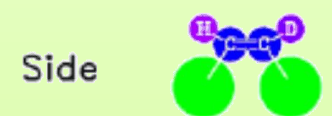
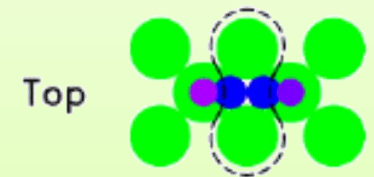


STM

d^2I/dV^2 map



C_2HD / $Cu(100)$



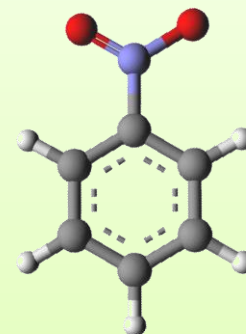
⇒ submolecular vibrational contrast

IETS of organic molecules

nitrobenzene on Cu(111)

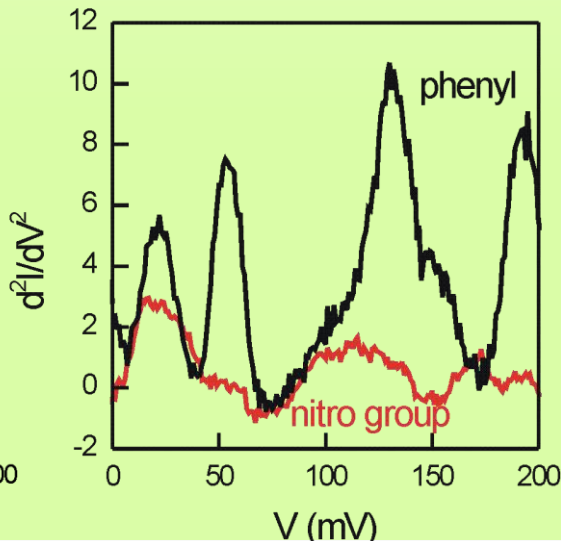
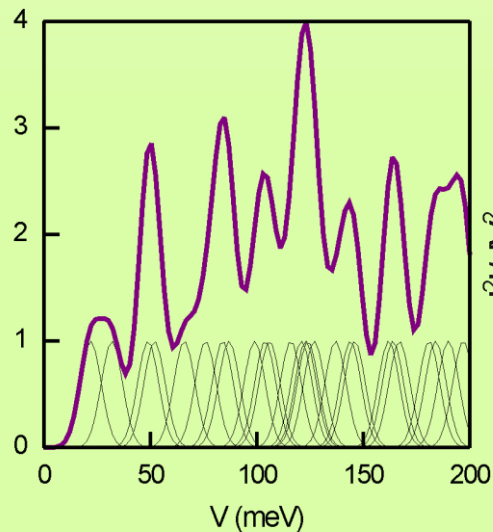
J. Phys.: Condens. Mat. 23, 484007 (2011)

Ring related	symmetry	Energy (meV)	Nitro group related	symmetry	Energy (meV)
C-H stretch	A1	386/382/378			
	B2	384/376			
C-C stretch	A1	197/184			
	B2	197/181/164			
C-H in-plane bent	A1	146/127	NO ₂ stretch	A1	167
	B2	162/144/133		B2	190
breathing	A1	124	C-NO ₂ stretch	A1	137
C-H out-of-plane bent	A2	121/104	NO ₂ bent	A1	106
	B1	123/116/99	NO ₂ Out-of-plane bent	B1	87
Out-of-plane deformation	A2	49	NO ₂ in-plane bent	A2	66
	B1	84/52	C-NO ₂ in-plane bent	B2	32
In-plane deformation	A1	84/49	C-NO ₂ out-of-plane bent	B1	22
	B2	76			



35 modes

IRAS: D. Syomin et al. Surf. Sci. Lett. 495 (2001) L827

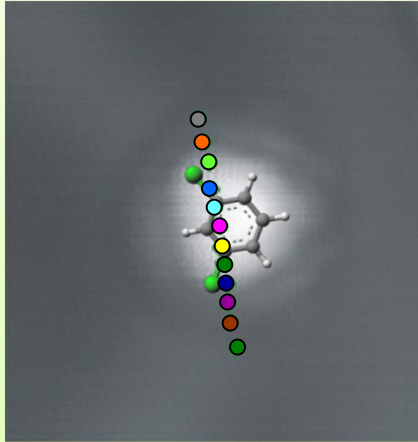


limited energy resolution
due to thermal and
modulation broadening
⇒ vibrational bands

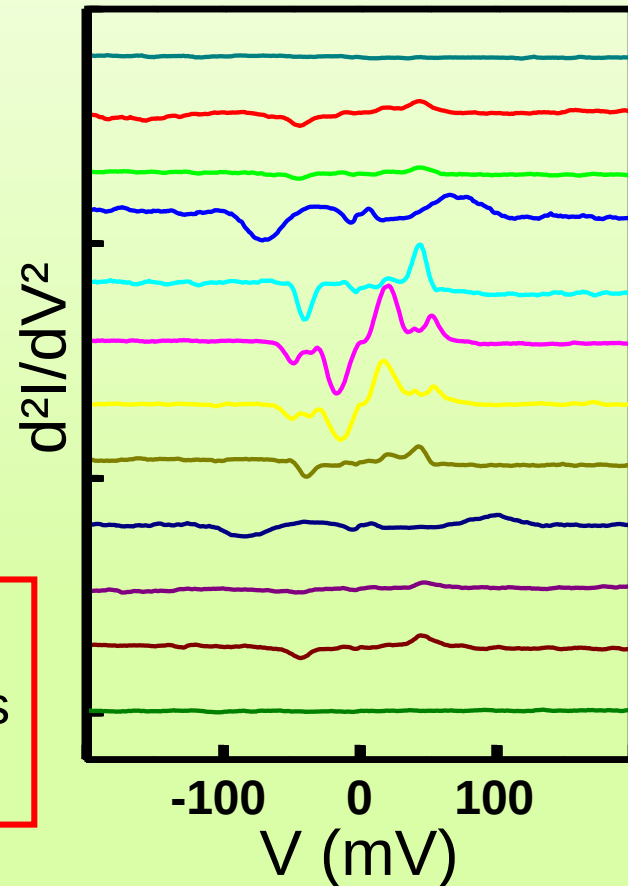
IETS: Spatial resolution

dichlorobenzene / Au(111)

Phys. Rev. B 89, 125420 (2014)



210 pA, -246 mV

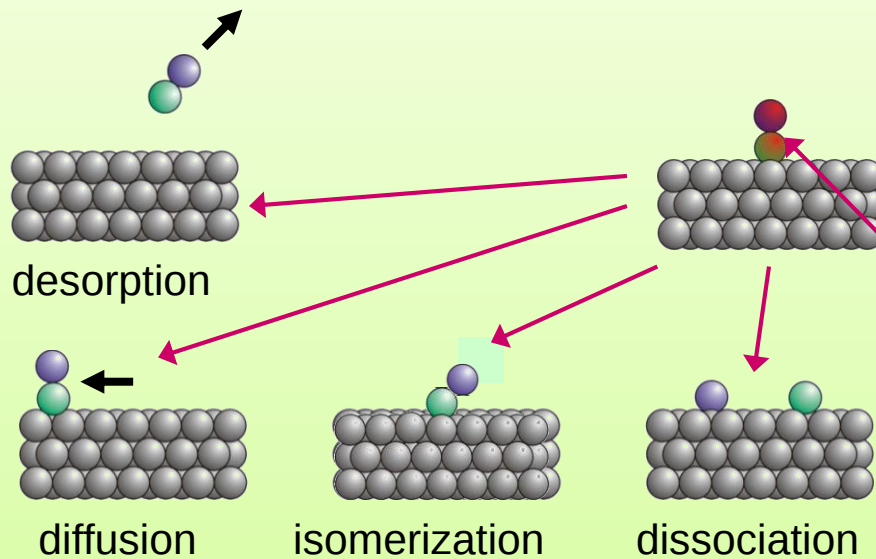


IETS

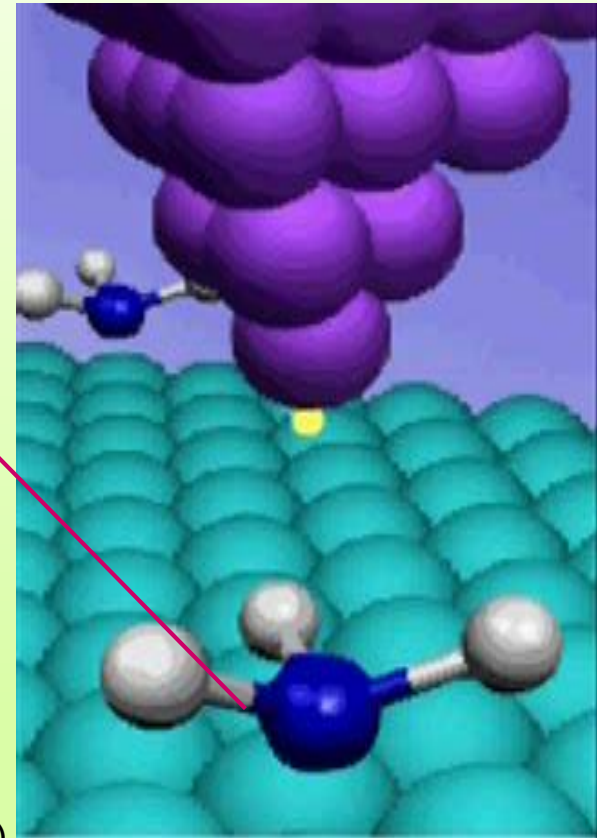
- ⇒ local excitation of molecular vibrations
- ⇒ chemical and isomeric identification

IET manipulation

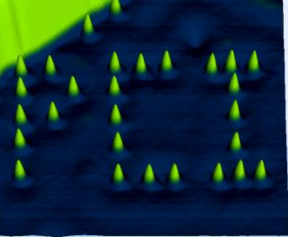
Phys. Status Solidi B, 250, 1671 (2013)



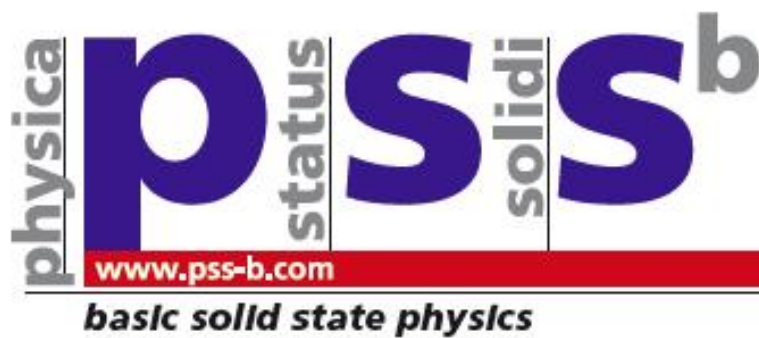
O₂ dissociation on Pt(111)
B.C. Stipe et al.,
Phys. Rev. Lett. 78, 4410 (1997)



with courtesy of
I.J. Pascual



ISSN 0370-1972
Phys Status Solidi B
250 · No. 9 September
1665–1948 (2013)



RUB

9
2013

Review Article

Controlled manipulation of single atoms and small molecules

Karina Morgenstern, Nicolas Lorente, and Karl-Heinz Rieder



Progress in Surface Science 86 (2011) 115–161



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journal homepage: www.elsevier.com/locate/progsurf

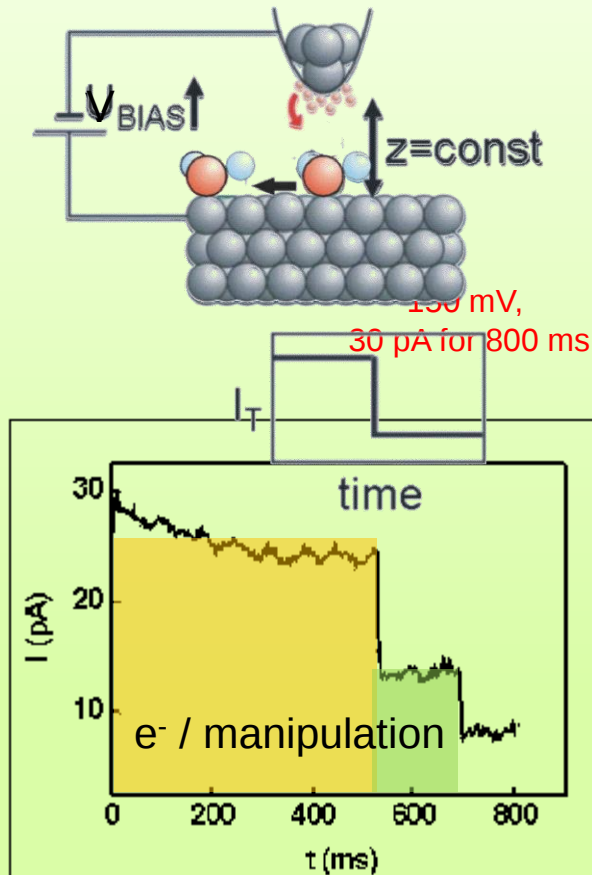


Review

Switching individual molecules by light and electrons: From isomerisation to chirality flip

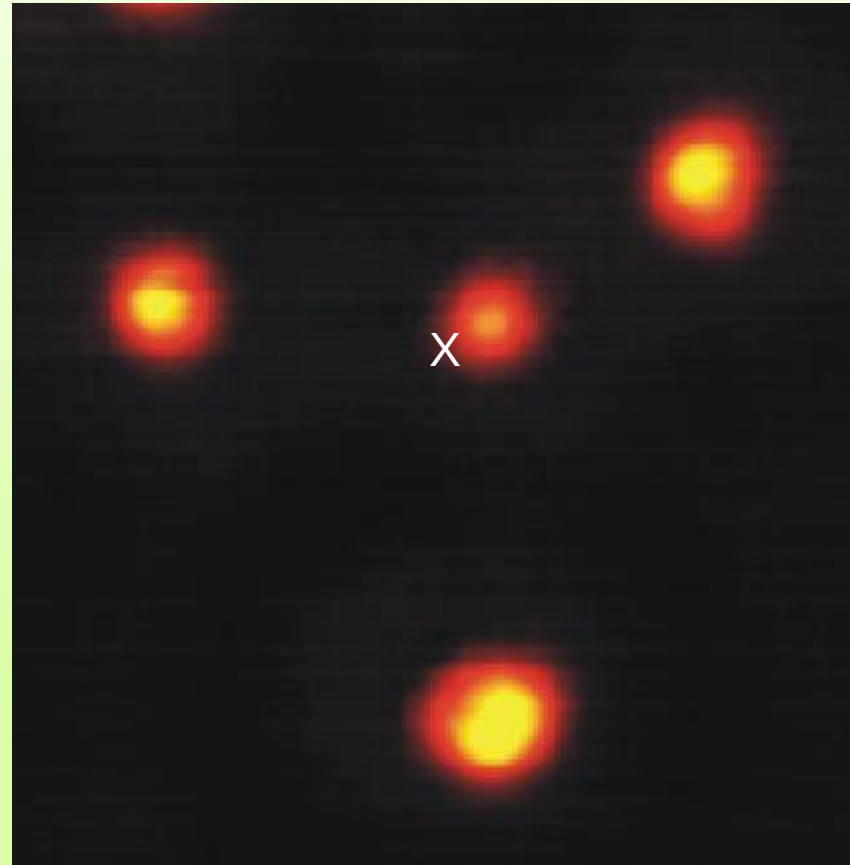
Karina Morgenstern *

IET manipulation



D₂O on Ag(111)

J. Chem. Phys. 116, 5746 (2002)



71 mV, 12 pA

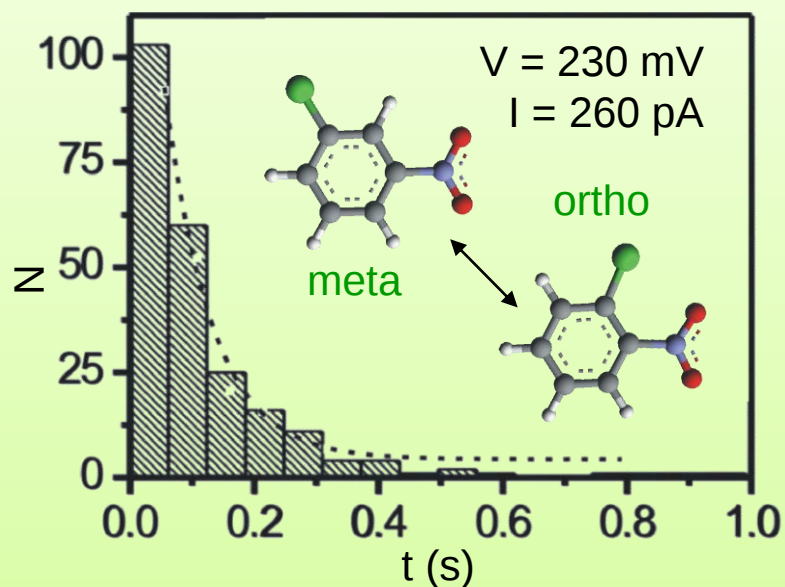
⇒ induced diffusion

IET manipulation: Reaction order

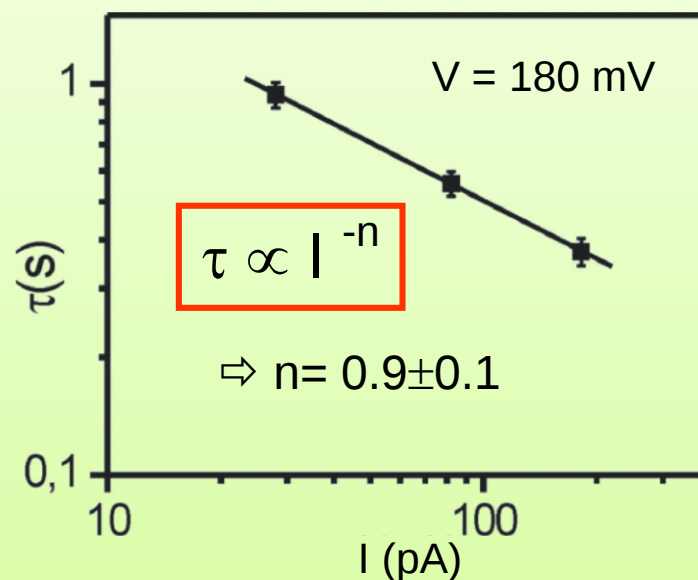
chloronitrobenzene on Cu(111)

Phys. Rev. Lett. 98, 116102 (2007)

Angew. Chem. Int. Ed. 48, 6041 (2009)



residence time
before isomerisation
 $\Rightarrow$ decay constant τ

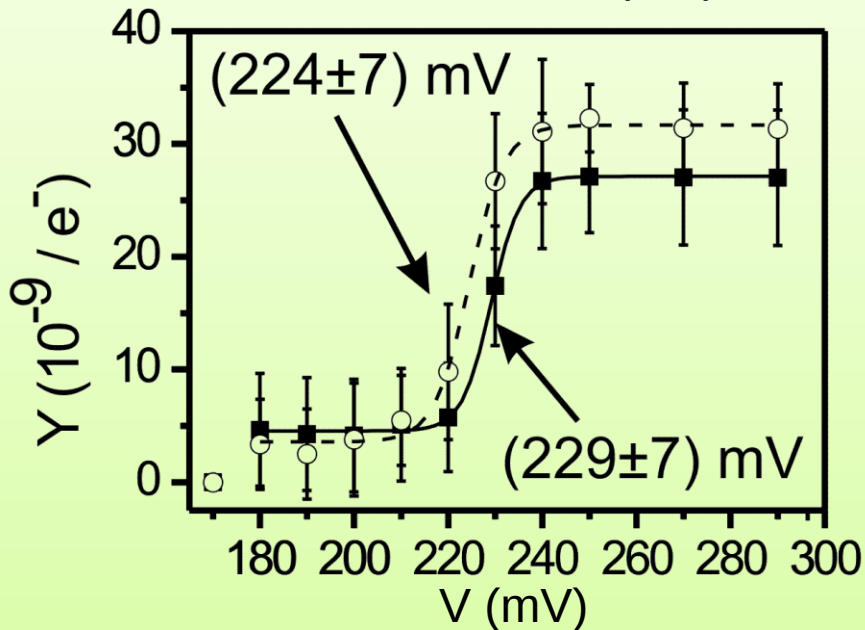


reaction order n
 $\Rightarrow$ one electron process

Action spectroscopy

RUB

chloronitrobenzene on Cu(111)



⇒ energy of molecular vibrations

⇒ action spectroscopy

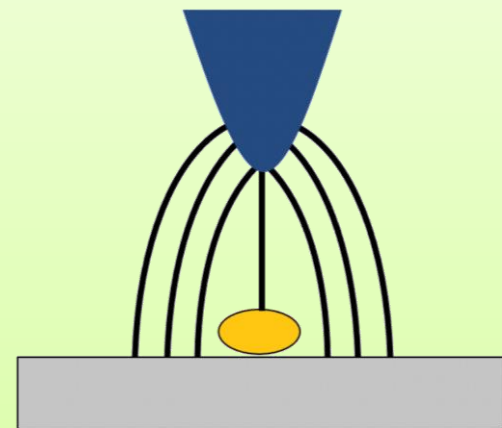
M. Kawai et al.

Phil. Trans. R. Soc. Lond. A 362, 1163 (2004)

Phys. Rev. Lett. 98, 116102 (2007)

Phys. Rev. Lett. 101, 136102 (2008)

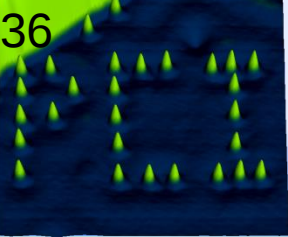
Angew. Chem. Int. Ed. 48, 6041 (2009)



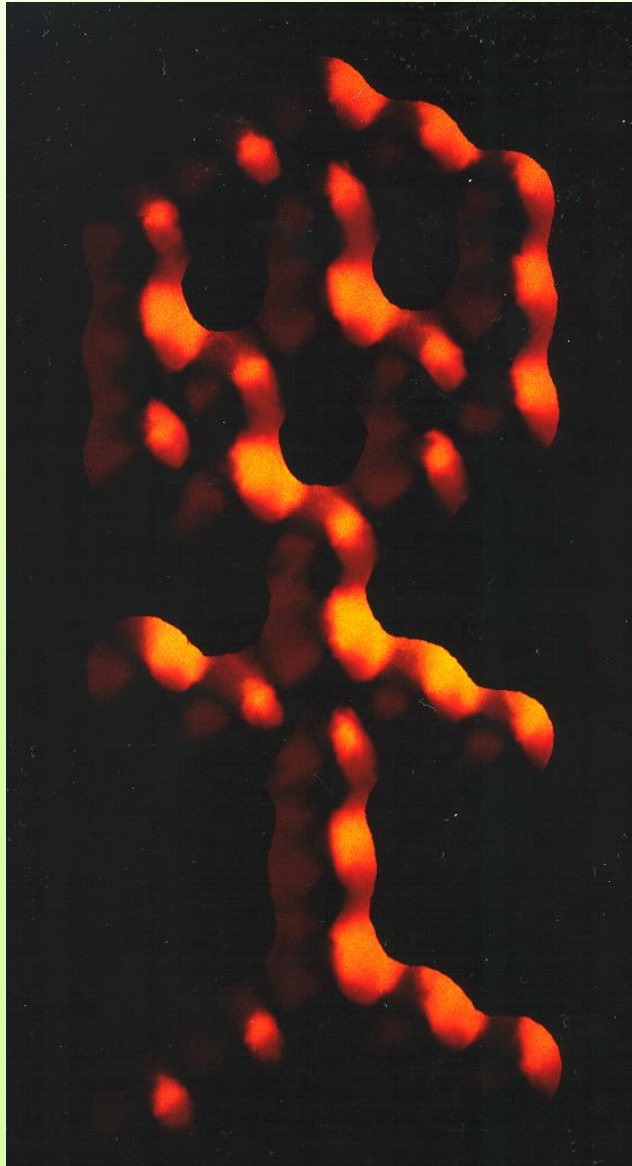
100 mV over 1 nm

⇒ $E = 10^8$ V/m

⇒ influence of the
electric field
on IET manipulation
and spectroscopy

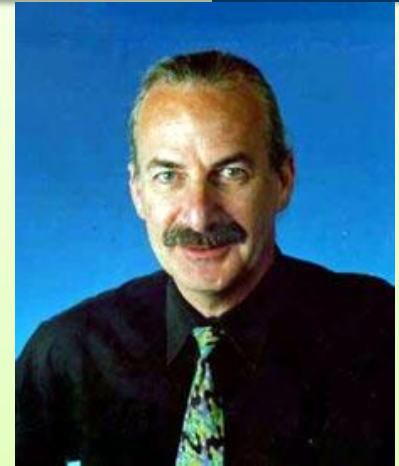


Manipulation by tip interaction



P. Zeppenfeld et al.
Ultramicroscopy 42-44, 128 (1992)

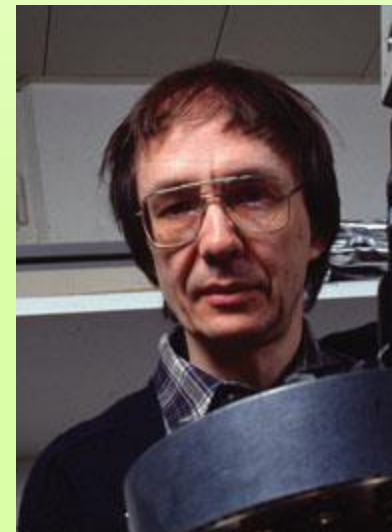
Don Eigler

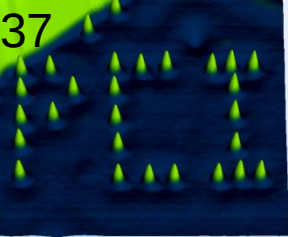


Karl-Heinz Rieder

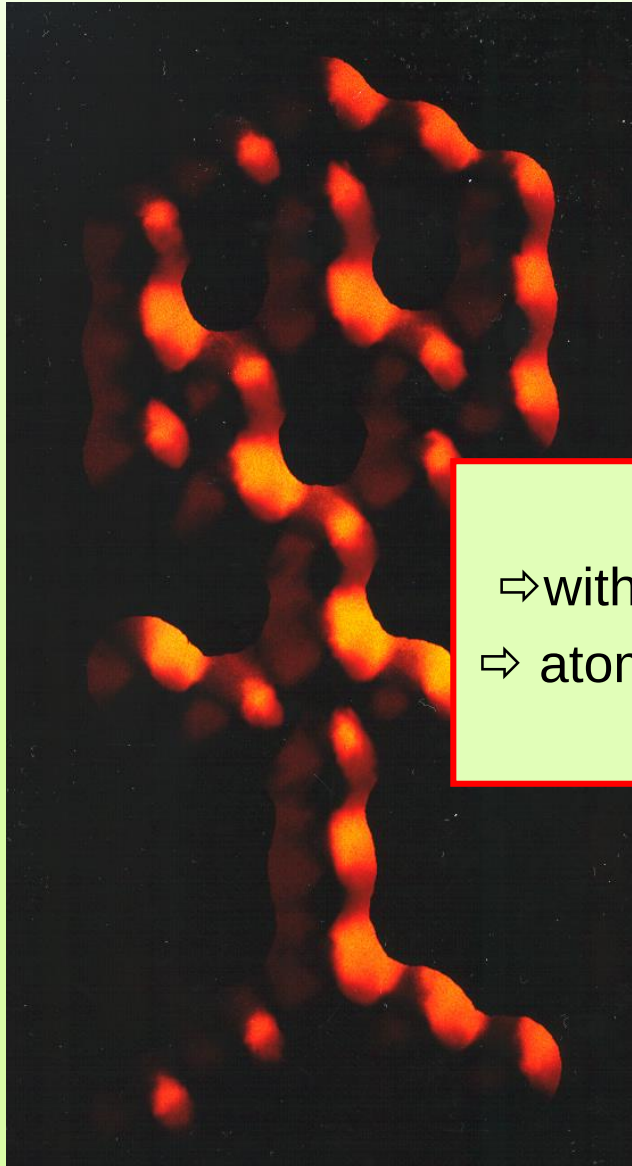


Gerhard Meyer

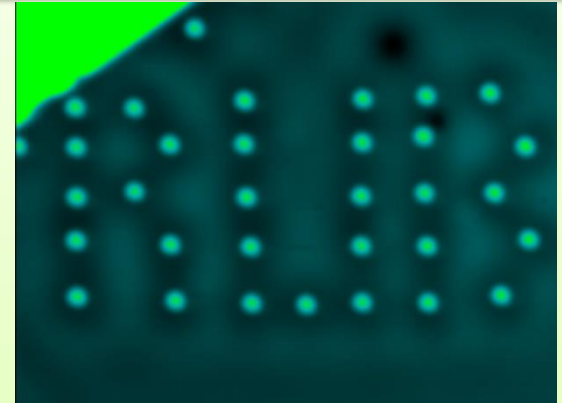




Manipulation by tip interaction



P. Zeppenfeld et al.
Ultramicroscopy 42-44, 128 (1992)



manipulation:

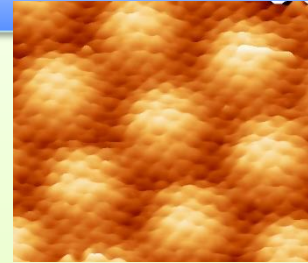
- ⇒ with electrons, electric field, or force
- ⇒ atoms, molecules, or molecular parts
- ⇒ molecular dynamics



Outline

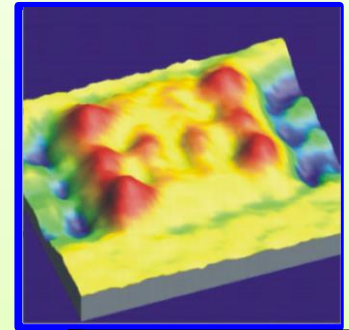
1) STM, STS, IETS, and manipulation

An extended introduction



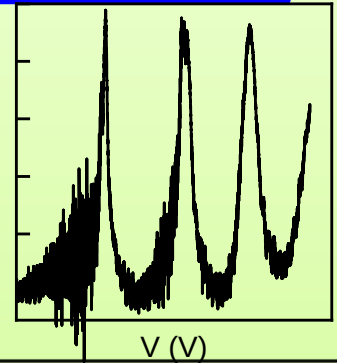
2) Quantum effects on surface imaged by STM

Electron in a box I



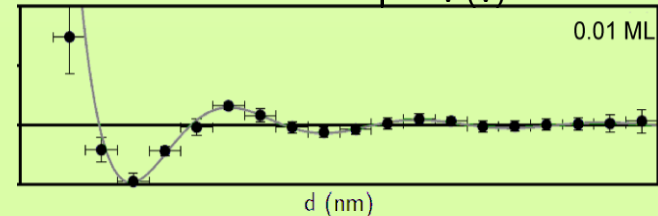
3) Quantum effects on surfaces measured by STS

Electron in a box II

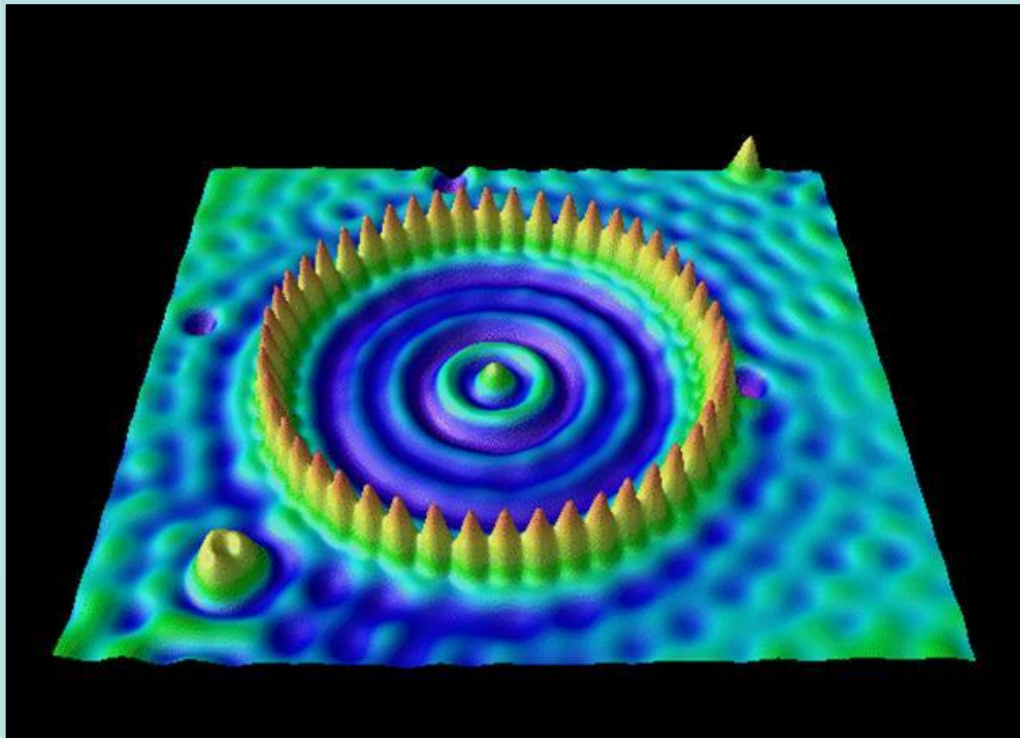


4) Electron interference effects surface diffusion

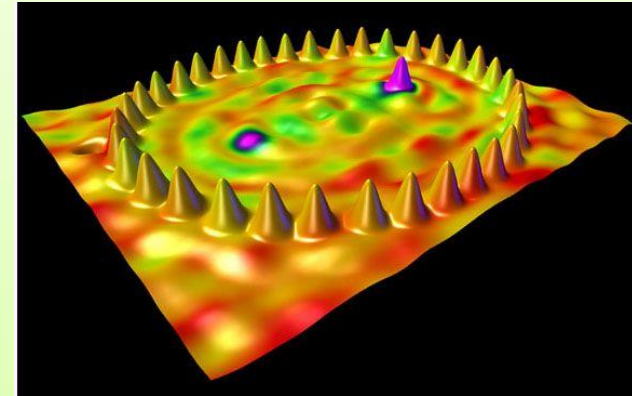
Influence of Friedel oscillations on diffusivity



Quantum corral/quantum mirage

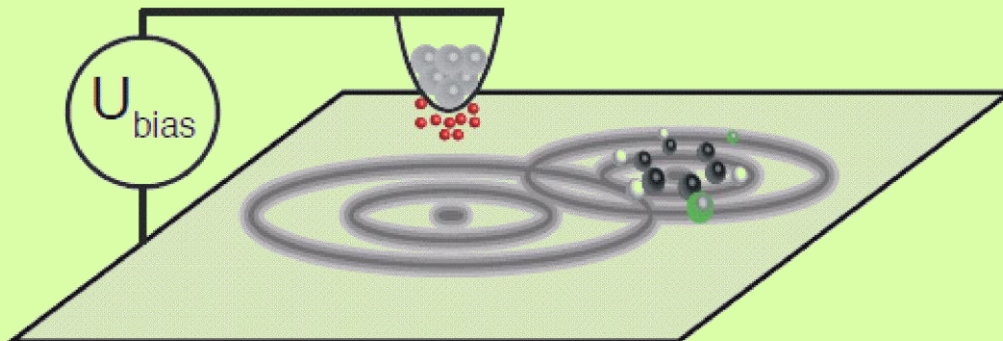


M. F. Crommie et al.,
Science 262, 218 (1993)

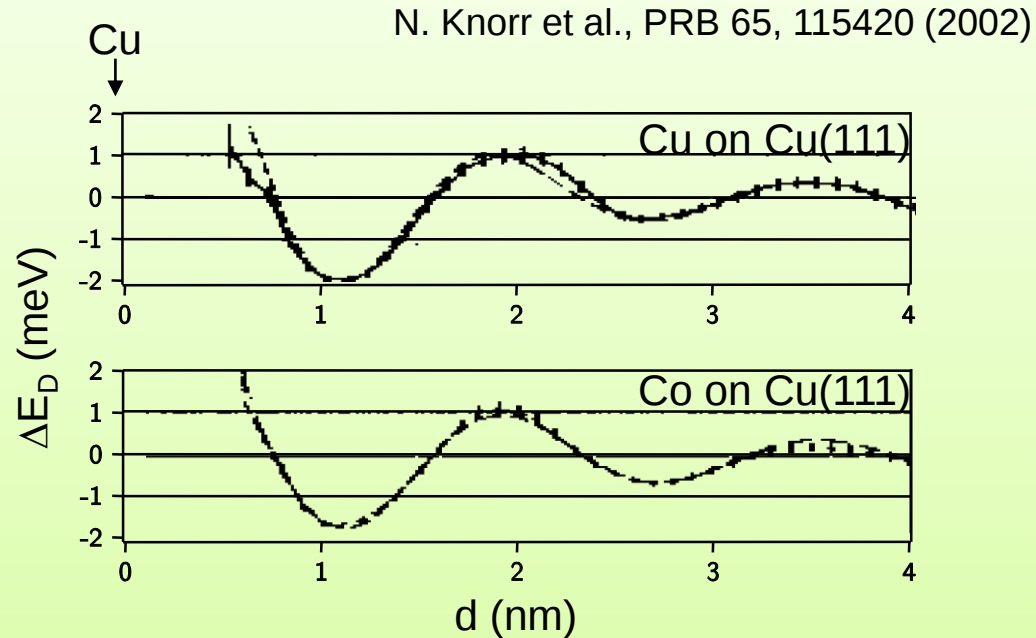
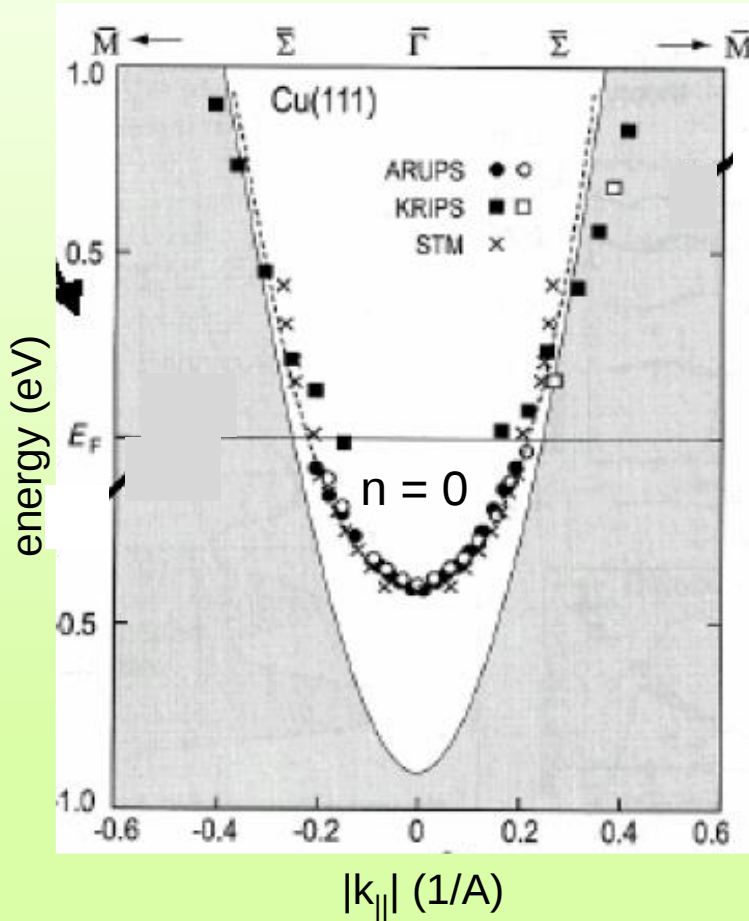


Friedel-like oscillations

⇒ interference pattern
of electrons



Surface state on fcc(111) surfaces

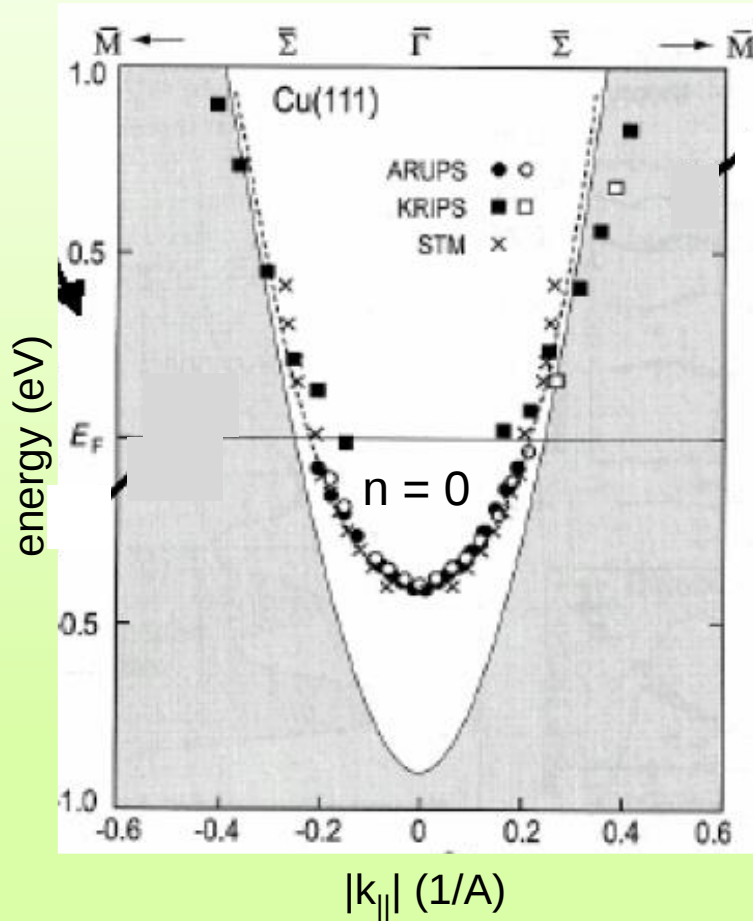


$$k_F = (2.0 \pm 0.1) \text{ nm}^{-1} \quad \phi_F = (0.49 \pm 0.03) \pi$$

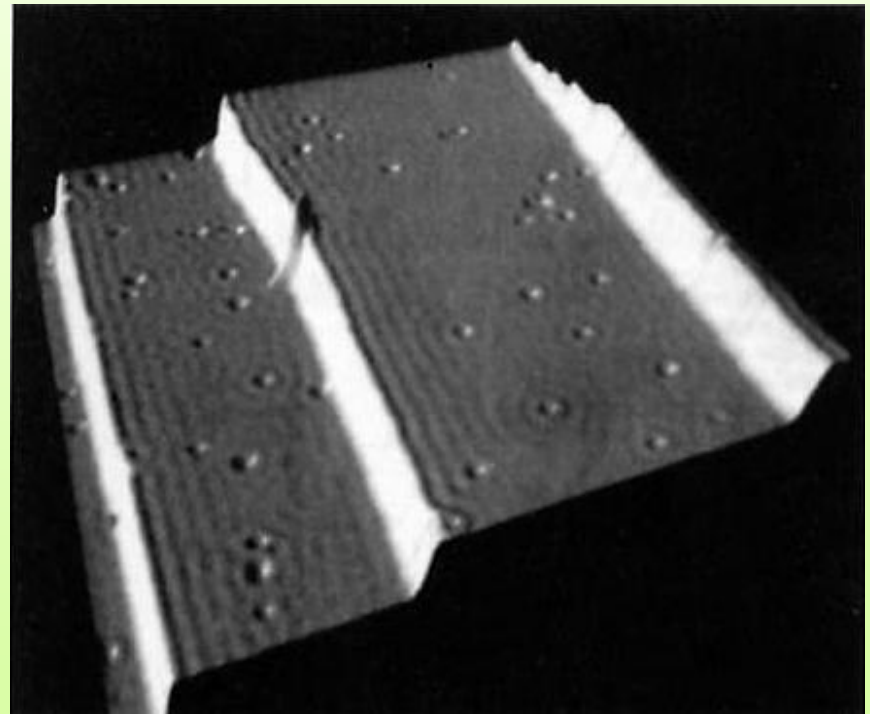
$$\Delta E(R) = -\alpha \cdot \frac{\sin(2k_F R + 2\phi)}{(k_F R)^2}$$

P. Hyldgaard, M. Persson,
J. Phys. Condens. Matter 12, L13 (2000)

Surface state on fcc(111) surfaces



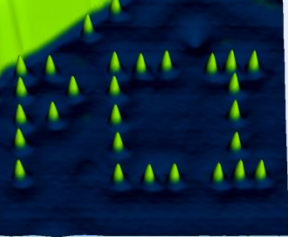
Crommie et al. Nature 363, 524 (1993)



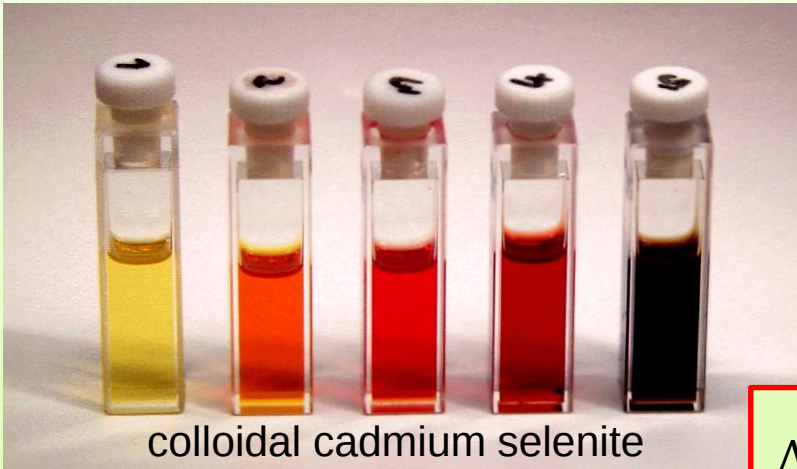
0.1 V; 50 x 50 nm²

$$\Delta E(R) = -\alpha \cdot \frac{\sin(2k_F R + 2\varphi)}{(k_F R)^2}$$

P. Hyldgaard, M. Persson,
J. Phys. Condens. Matter 12, L13 (2000)



Quantum size effects

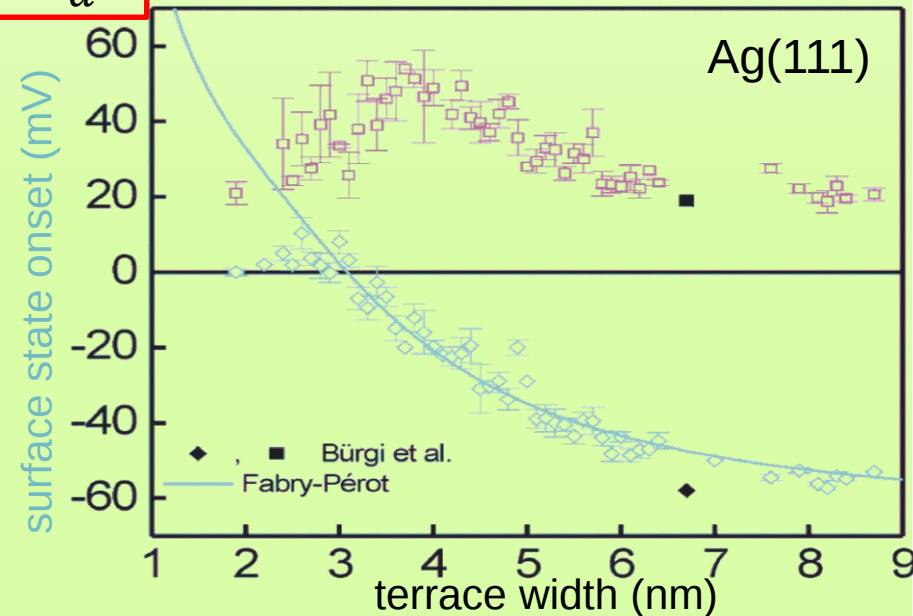
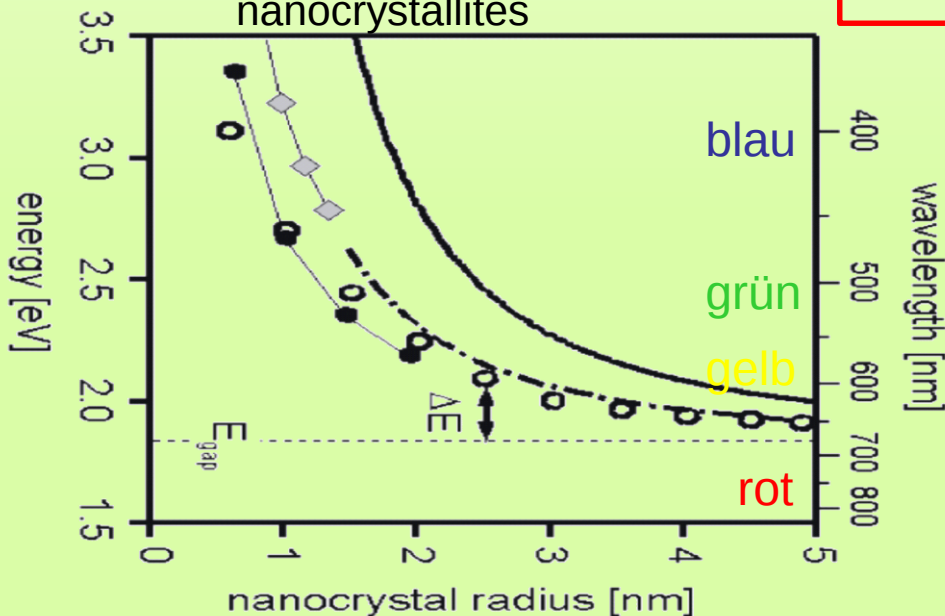


colloidal cadmium selenite
nanocrystallites

$$\Delta E(d) \propto \frac{1}{d^2}$$

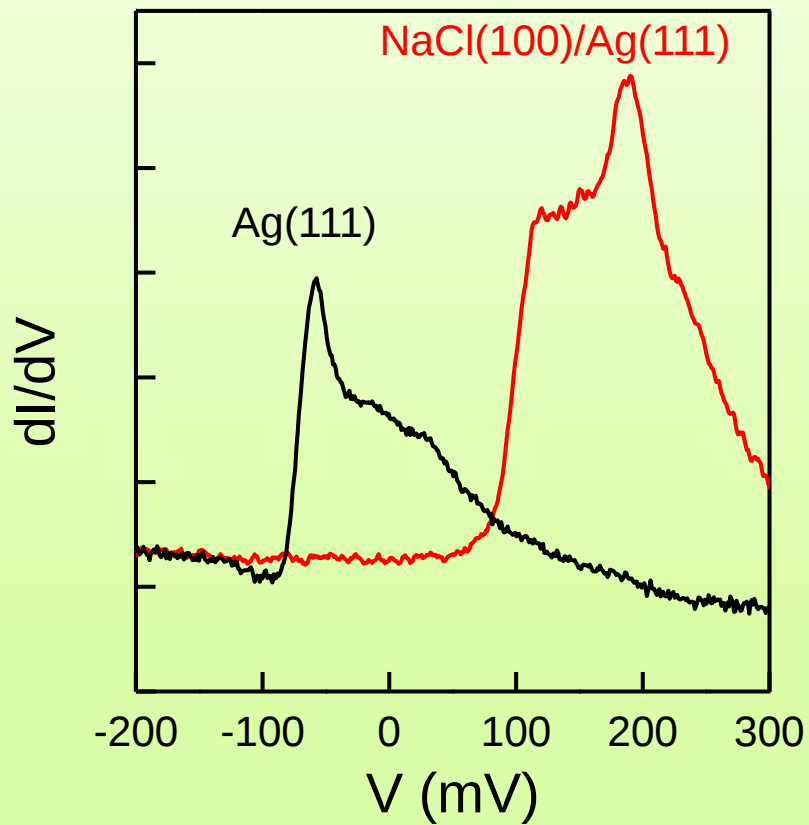


K. Morgenstern et al.
Phys. Rev. Lett. 89, 226801 (2002)

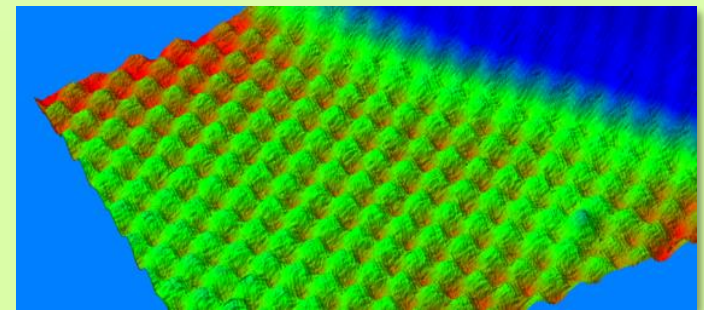
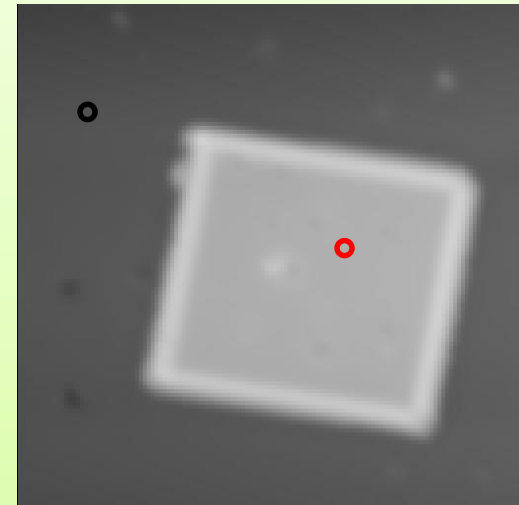


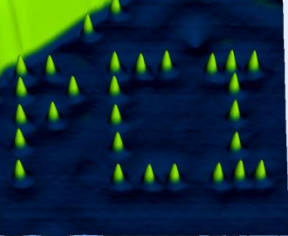
STS of insulator on fcc(111) metal

J Phys. Chem. C. 117, 16094 (2013)



NaCl(100) on Ag(111)

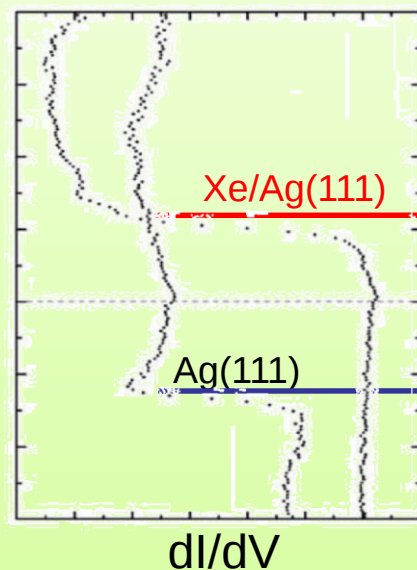
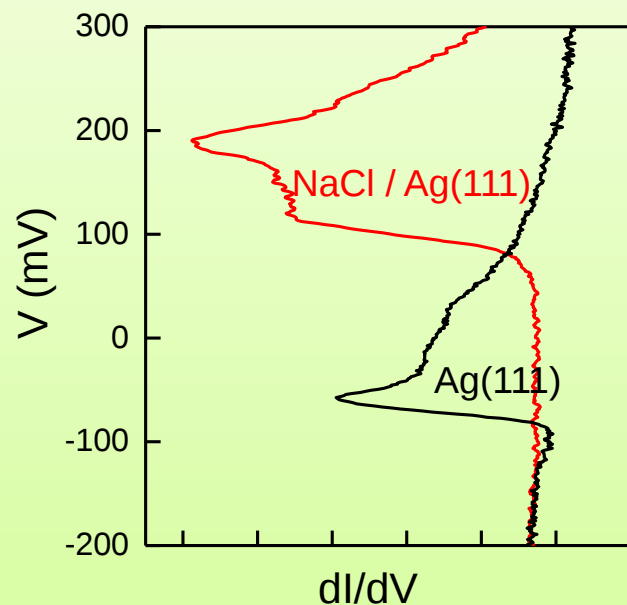




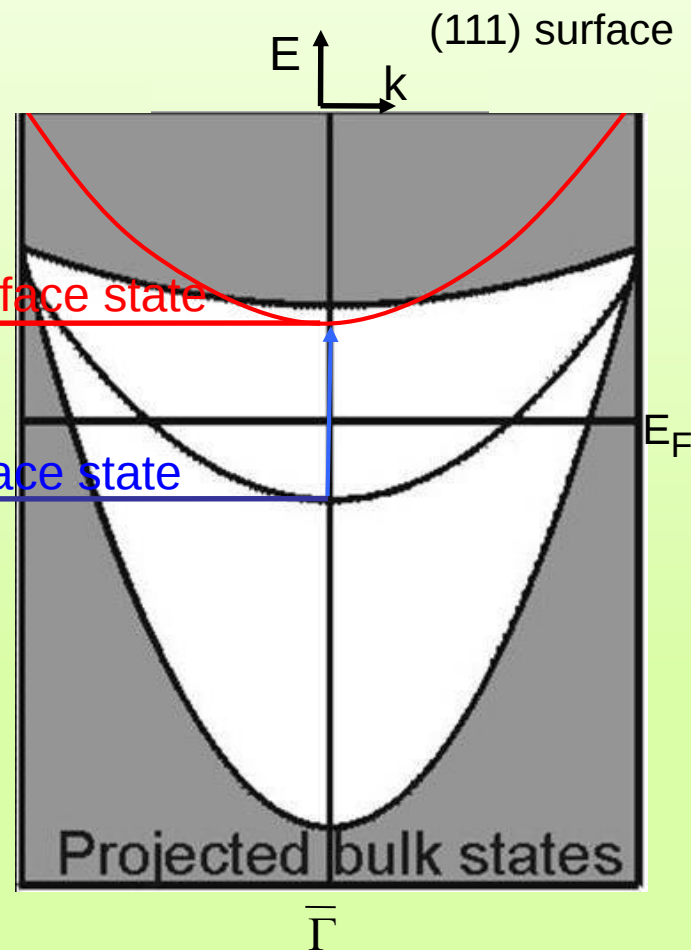
Surface vs. interface state

RUB

J Phys. Chem. C. 117, 16094 (2013)

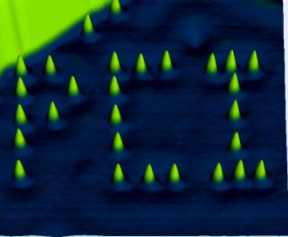


H. Hövel et al.
 Surf. Sci. 477, 43, 2001
 Ag(111) c.f. Bürgi et al.
 Phys. Rev. Lett. 81 (1998) 5370



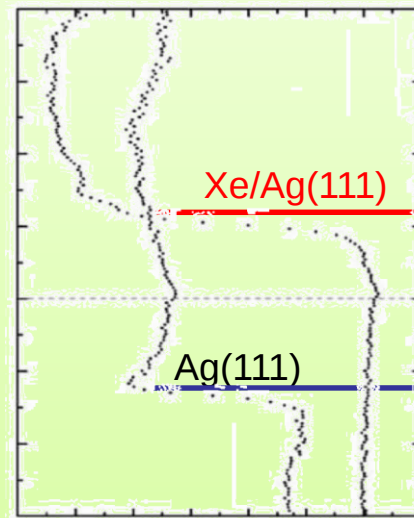
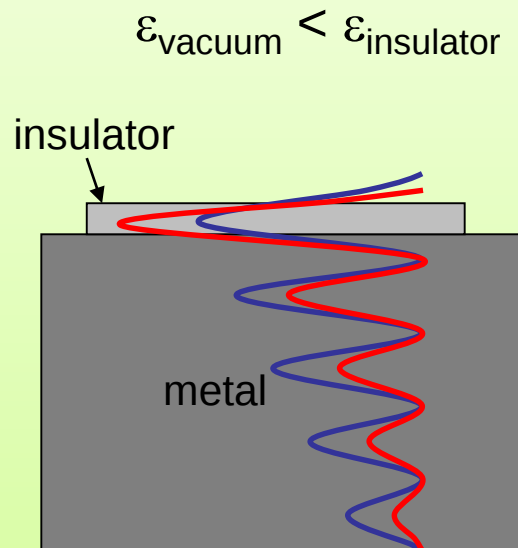
J. Li et al., Phys. Rev. Lett., 80(15), 3332 (1998)

H. Jensen et al., Phys. Rev. B, 71(15), 155417 (2005)

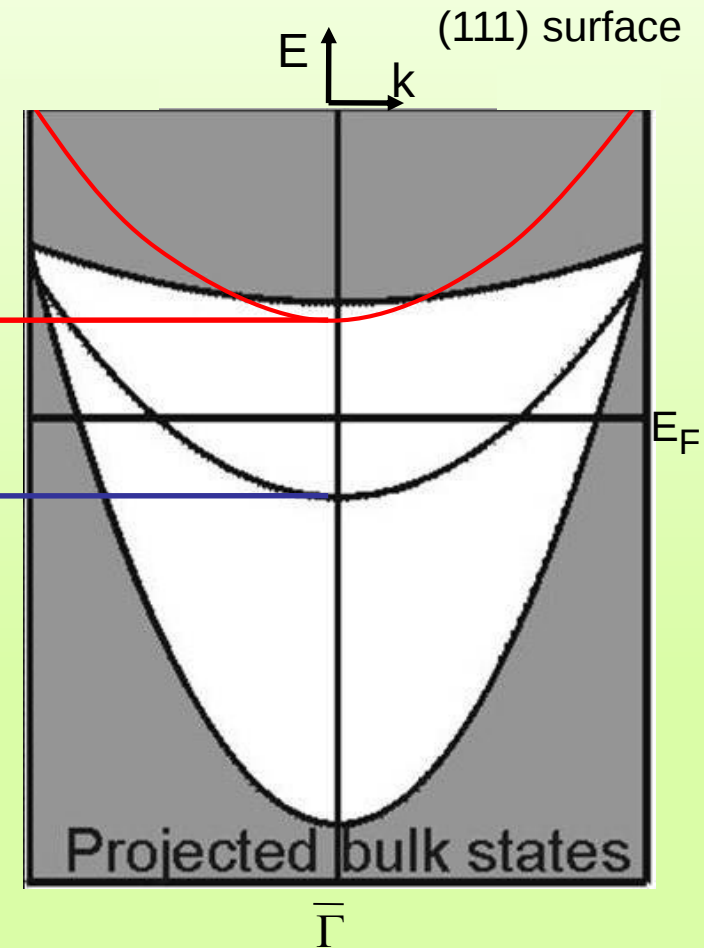


Surface vs. interface state

J Phys. Chem. C. 117, 16094 (2013)

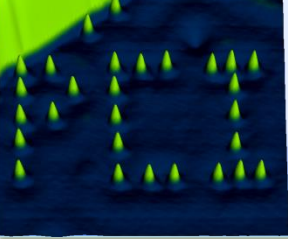


H. Hövel et al.
Surf. Sci. 477, 43, 2001



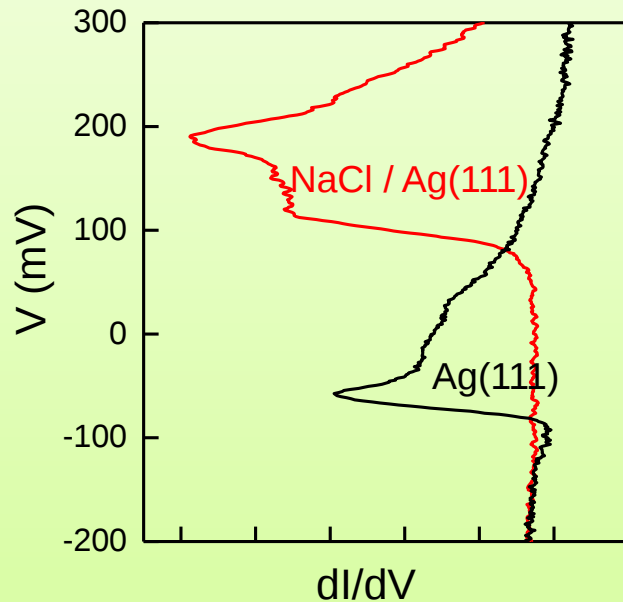
J. Li et al., Phys. Rev. Lett., 80(15), 3332 (1998)

H. Jensen et al., Phys. Rev. B, 71(15), 155417 (2005)



Surface vs. interface state

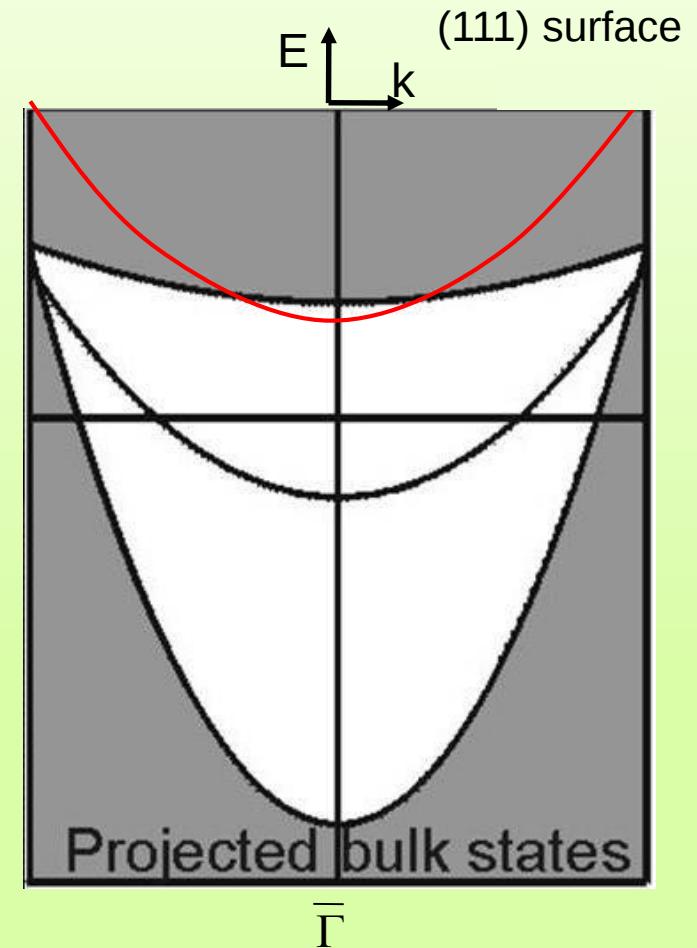
J Phys. Chem. C. 117, 16094 (2013)



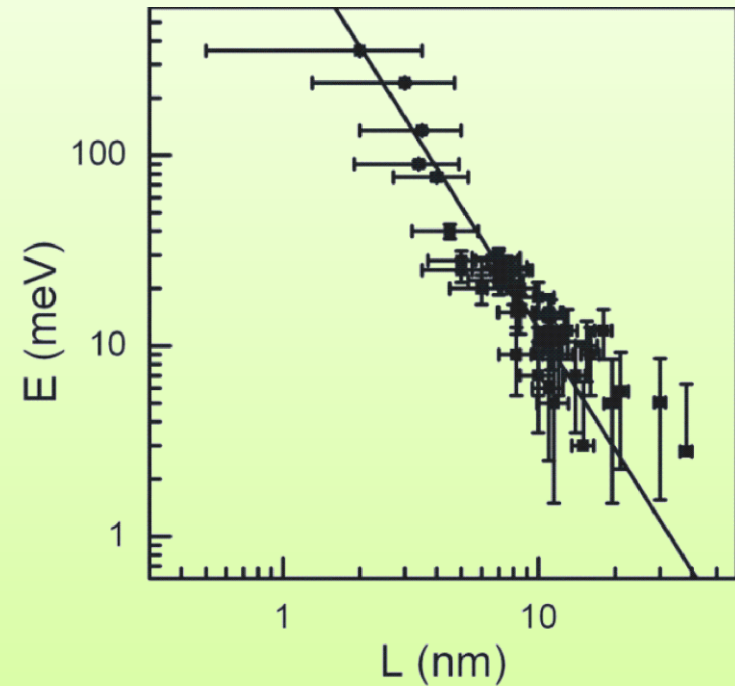
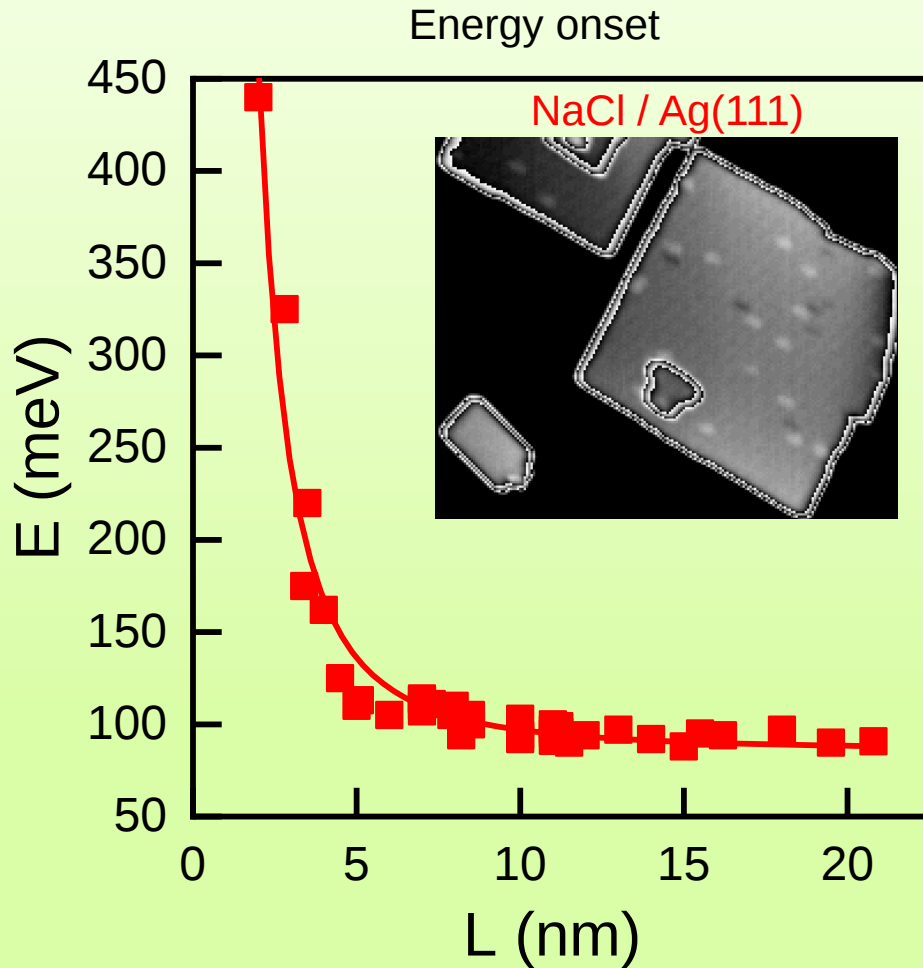
⇒ interface state
@ $+(90 \pm 4)$ meV

⇒ surface state
@ $-(65 \pm 3)$ meV

⇒ ~150 meV shift for NaCl layers



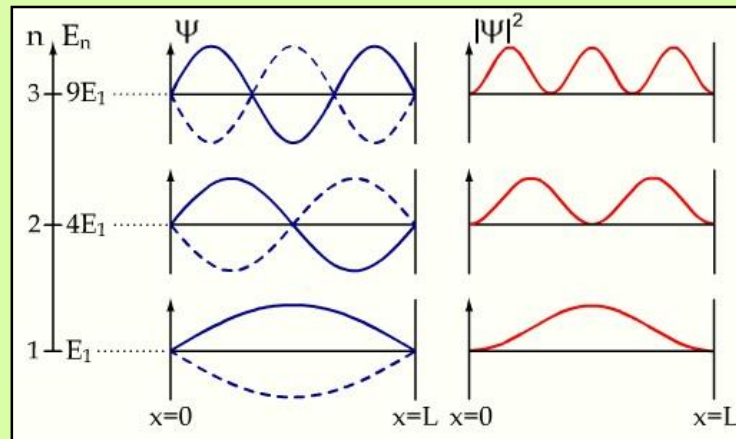
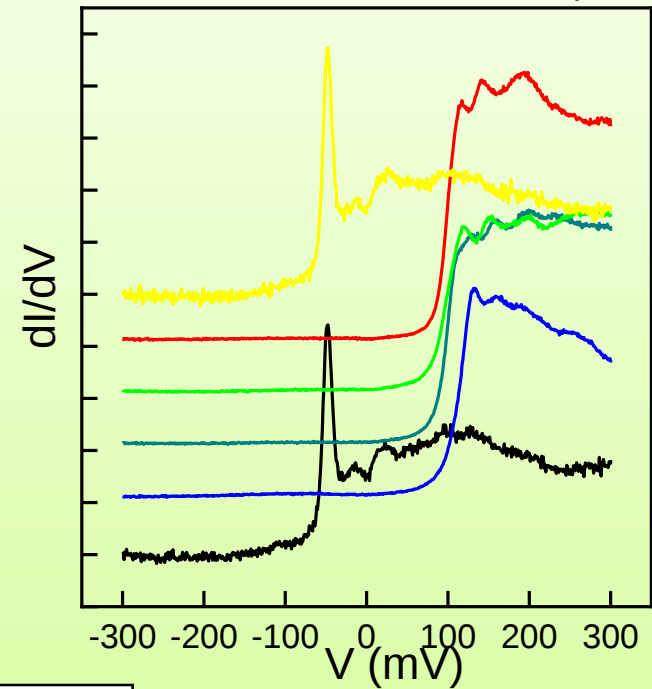
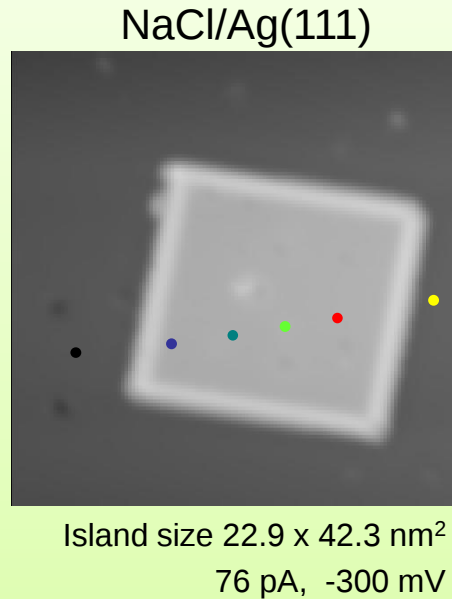
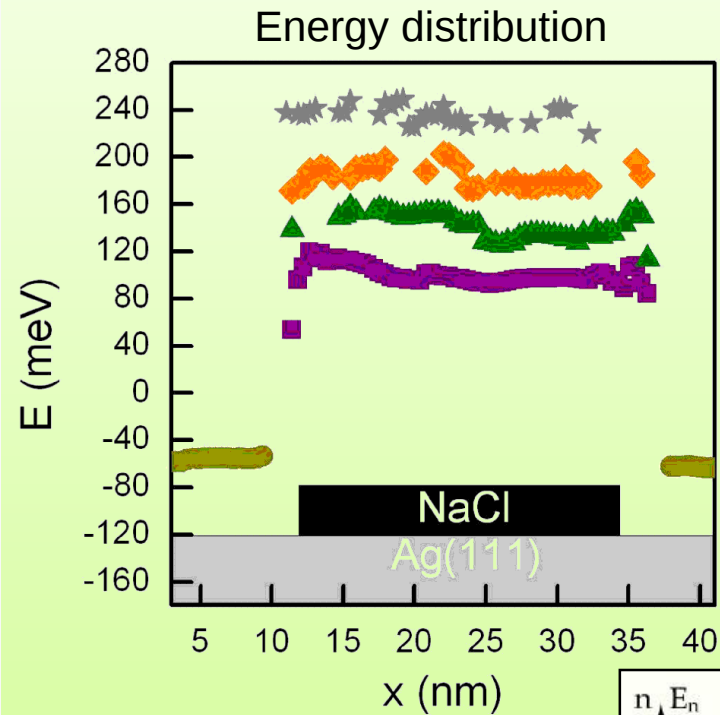
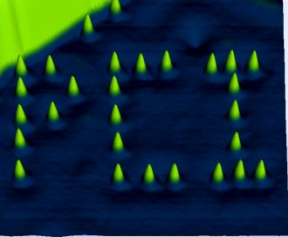
Quantum confinement



$$E \propto 1/L^2$$

$\Rightarrow$ quantum confinement < 8 nm

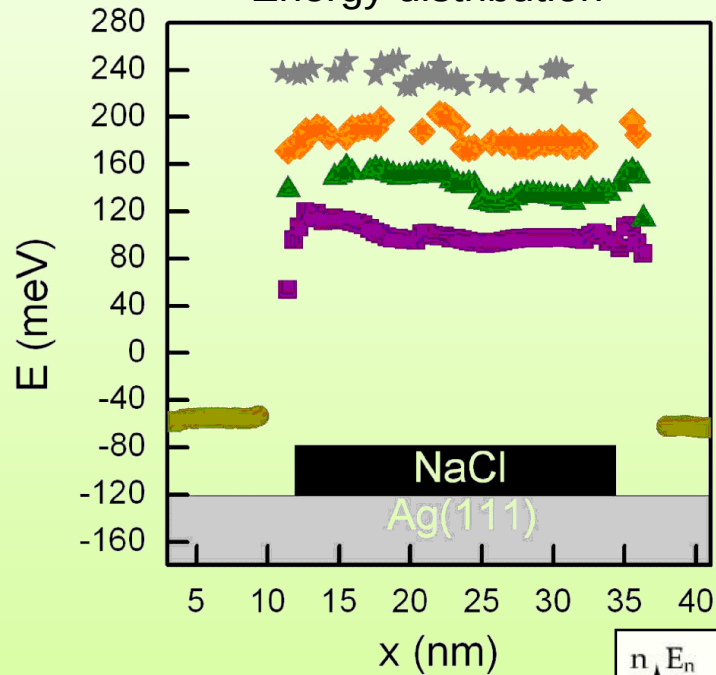
Particle in a box



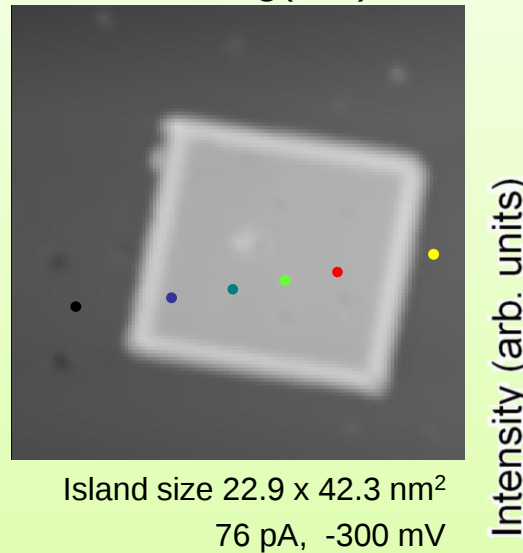
Particle in a box

Nano Lett. 14, 13 (2014)

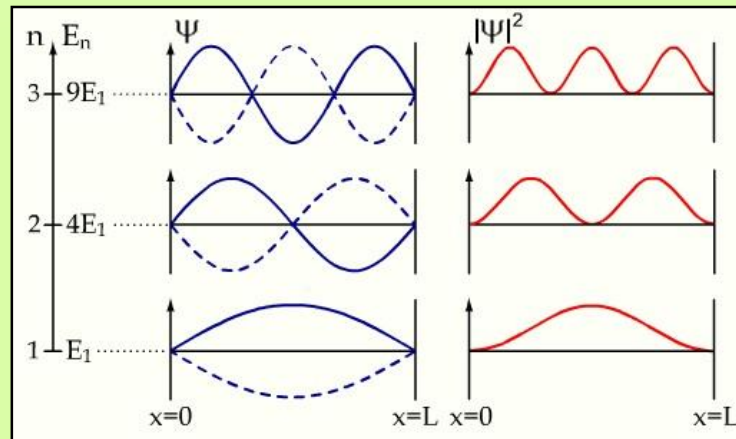
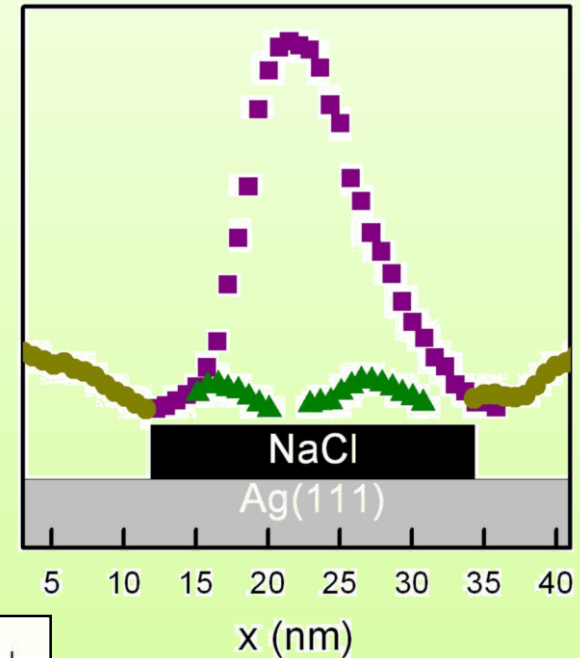
Energy distribution



NaCl/Ag(111)



Intensity distribution

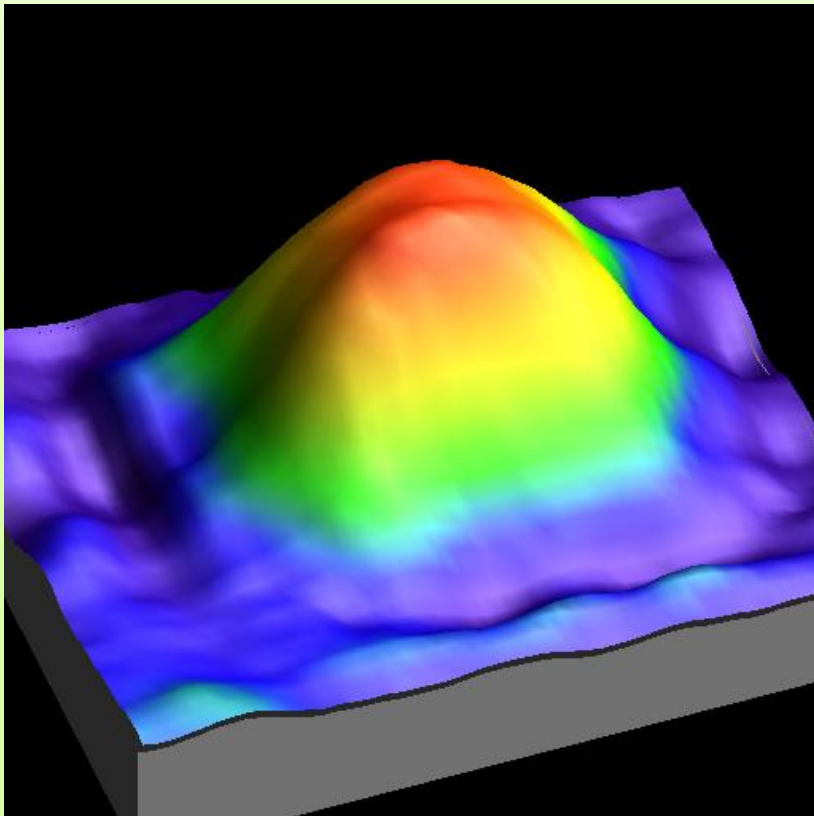


Particle in a box

STS maps

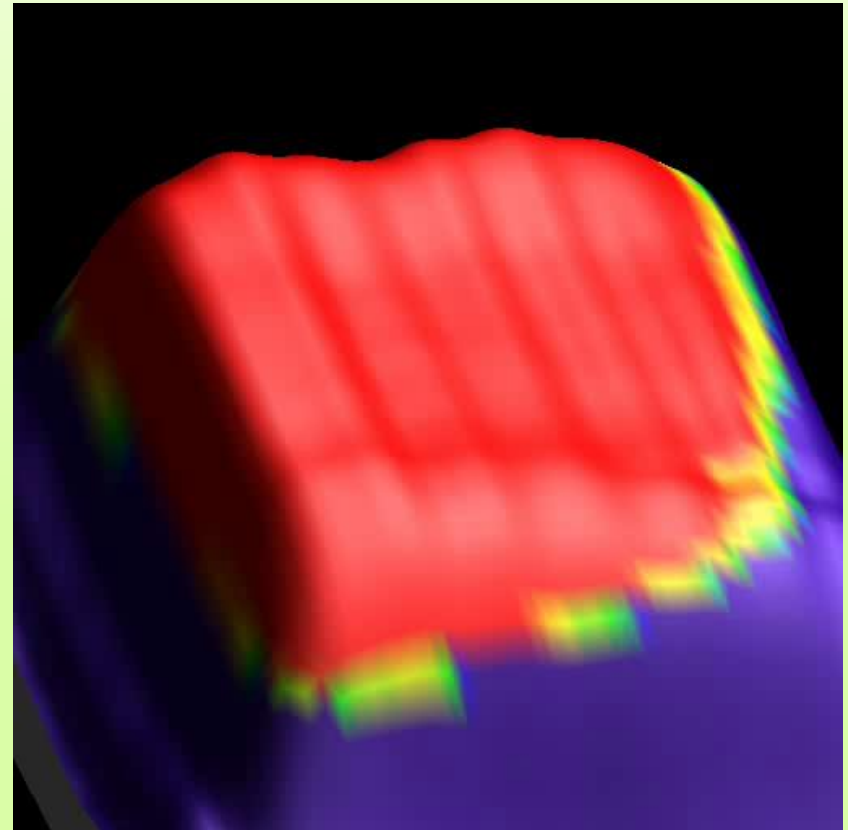
Nano Lett. 14, 13 (2014)

from 70 to 300 meV
Island: $12 \times 10 \text{ nm}^2$
voltage steps: 6 meV



0.1 nA
modulation: 5 meV

from 83 to 300 meV
Island: $21 \times 22 \text{ nm}^2$
voltage steps: 7 meV

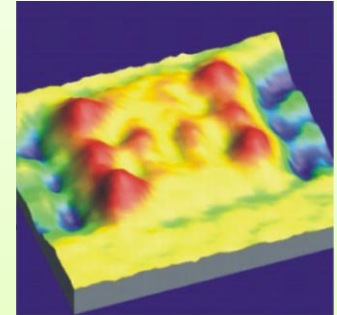
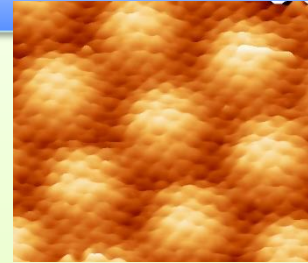


0.1 nA
modulation: 7 meV

Outline

1) STM, STS, IETS, and manipulation

An extended introduction

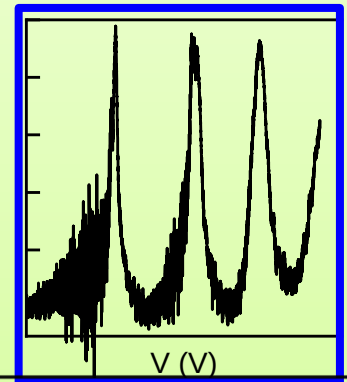


2) Quantum effects on surface imaged by STM

Electron in a box I

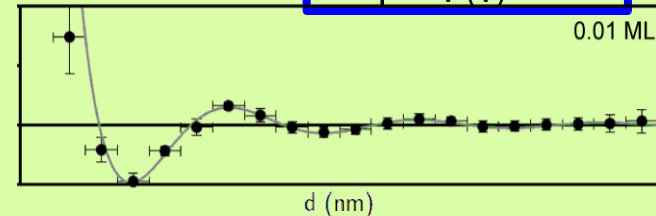
3) Quantum effects on surfaces measured by STS

Electron in a box II skip



4) Electron interference effects surface diffusion

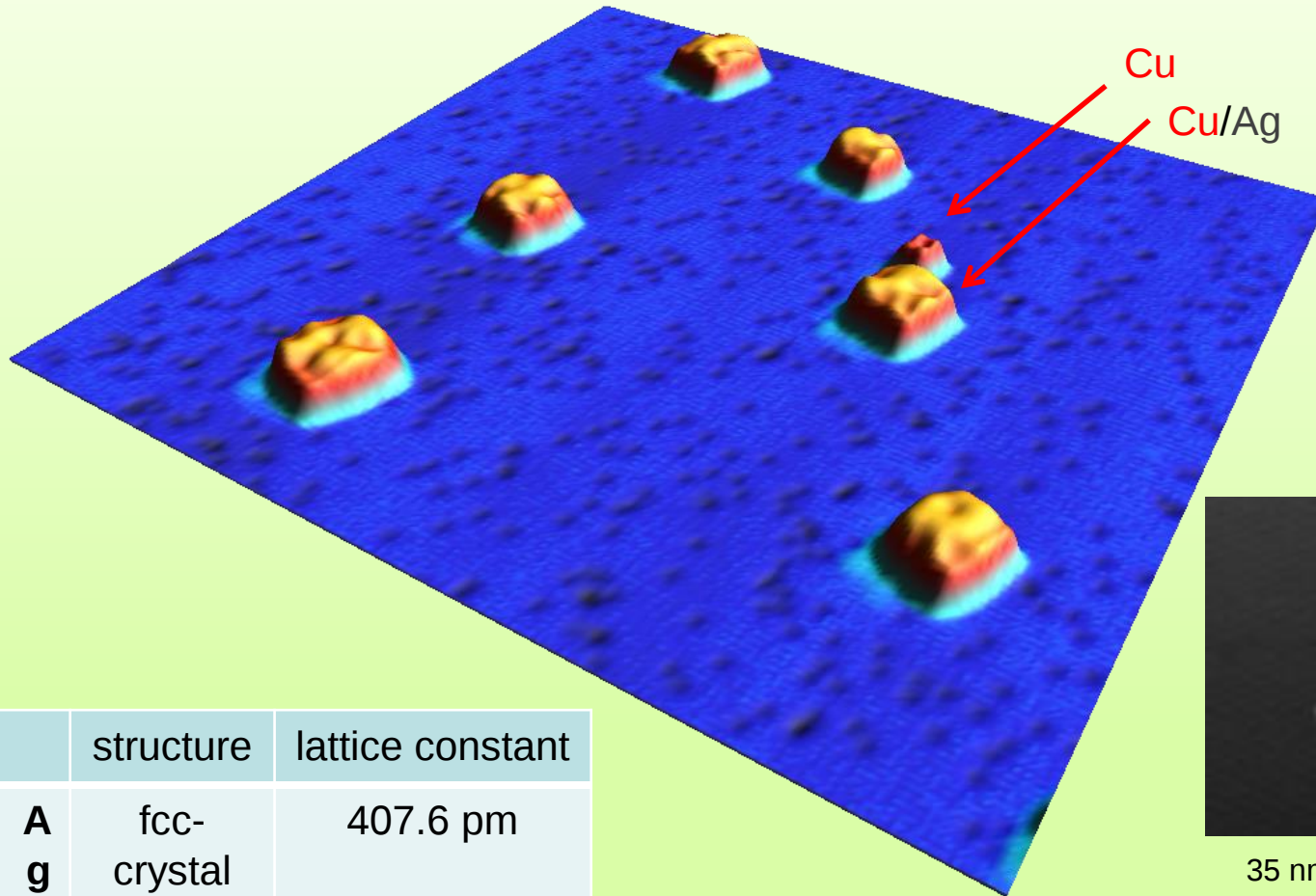
Influence of Friedel oscillations on diffusivity



Grain size stability of a complex surface: Cu on Ag(100)

RUB

PRL 107, 046101 (2011)
PRB 90, 165418 (2014)
PRB 92, 045422 (2015)



35 nm x 21 nm, 0.79 nA, 0.61 V, 300 K

	structure	lattice constant
A g	fcc-crystal	407.6 pm
C u	fcc-crystal	361.5 pm

13% lattice mismatch

- large strain
- quadratic islands of monatomic height

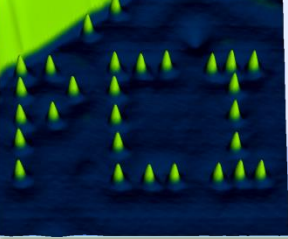
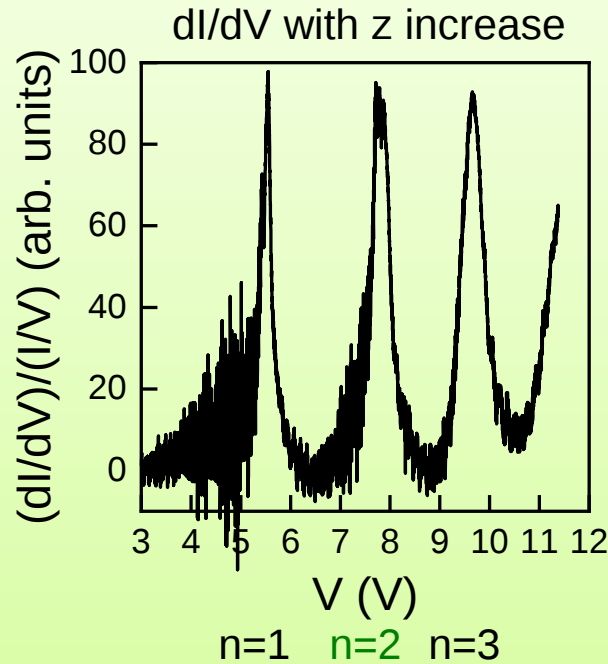


Image potential states

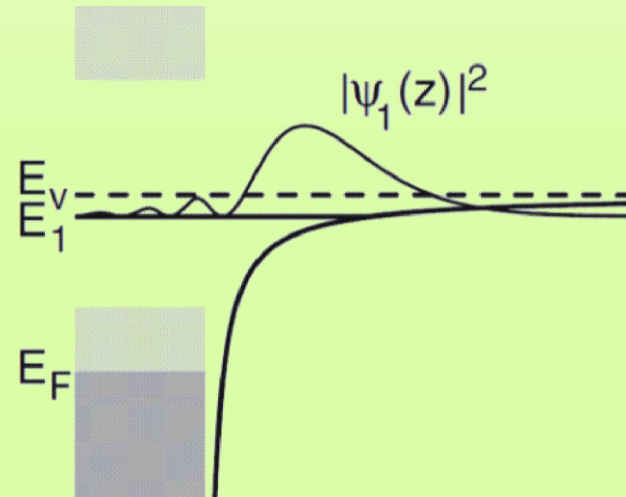
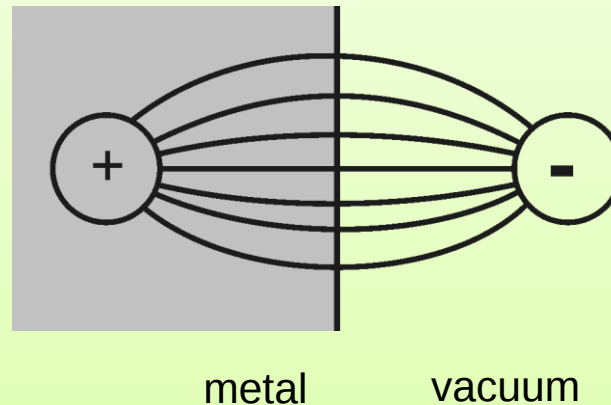
Cu/Ag(100)

Phys. Rev. B 90, 165418 (2014)



$$E_n = E_V - \frac{13.6\text{eV}}{[4(n+a)]^2}$$

Image potential



see: P.M. Enchenique, e.g. Chem. Rev. 106, 4160 (2006)

sketch from Crampin PRL 95 046801 (2005)

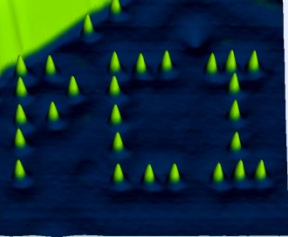
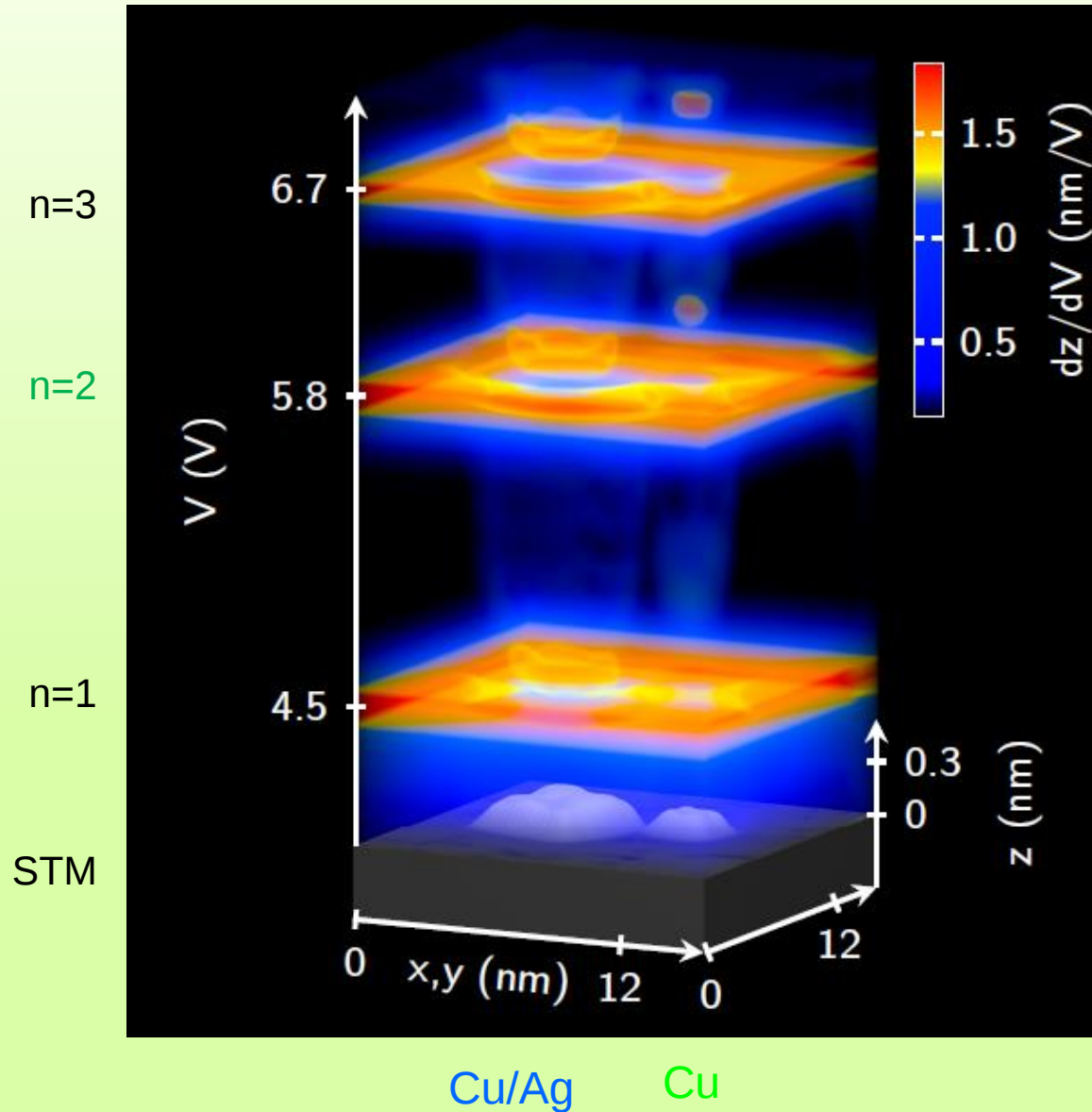


Image potential states

Phys. Rev. B 90, 165418 (2014)



Cu/Ag(100) islands

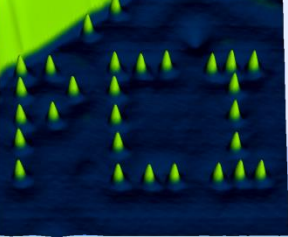
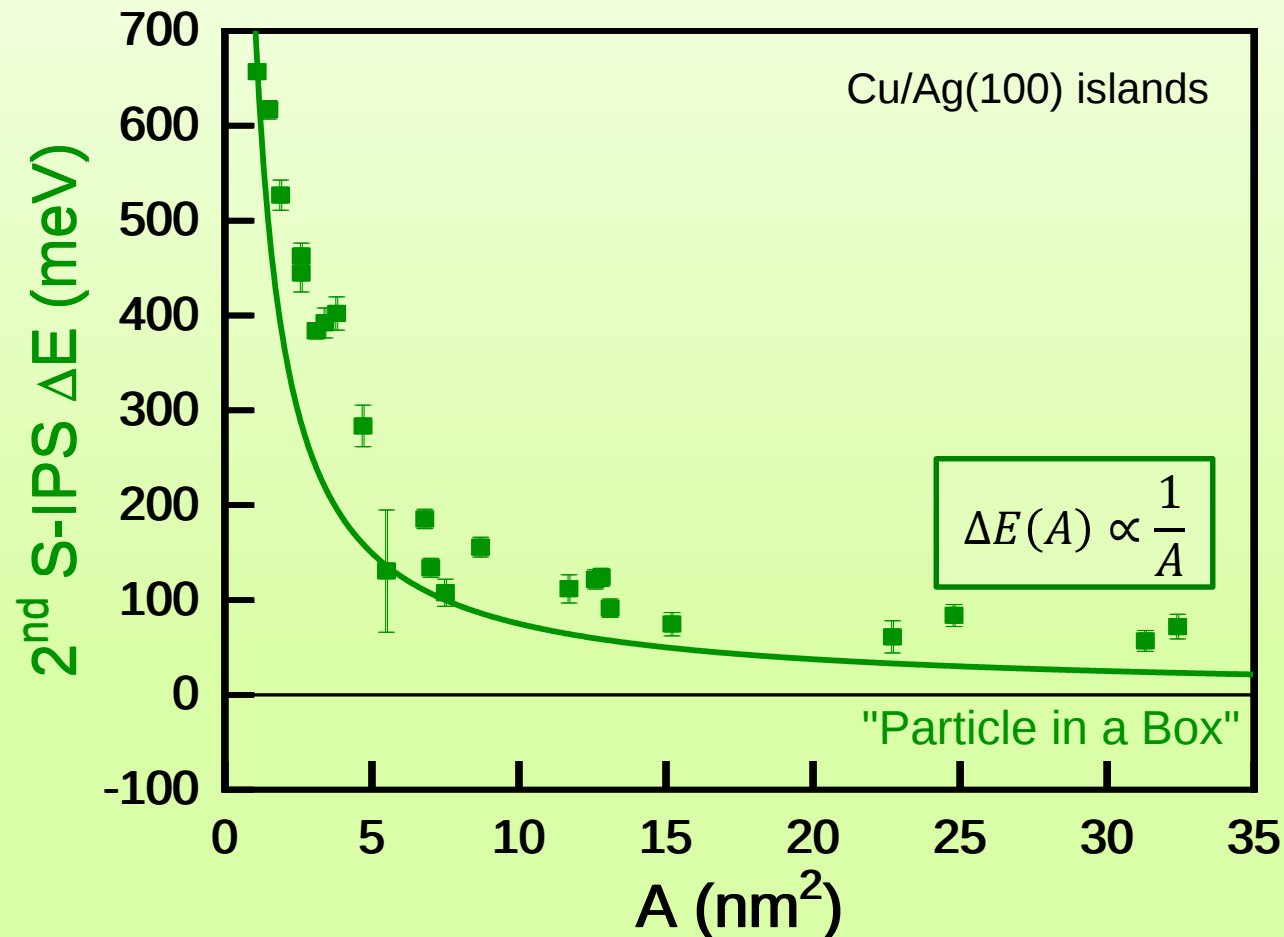
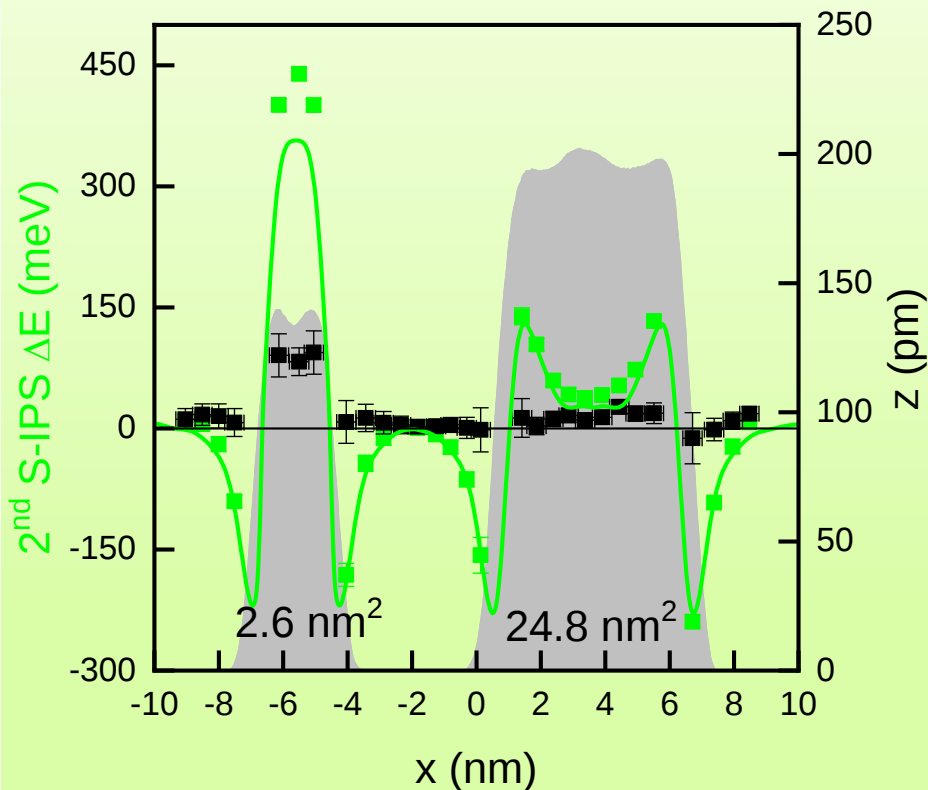


Image potential states



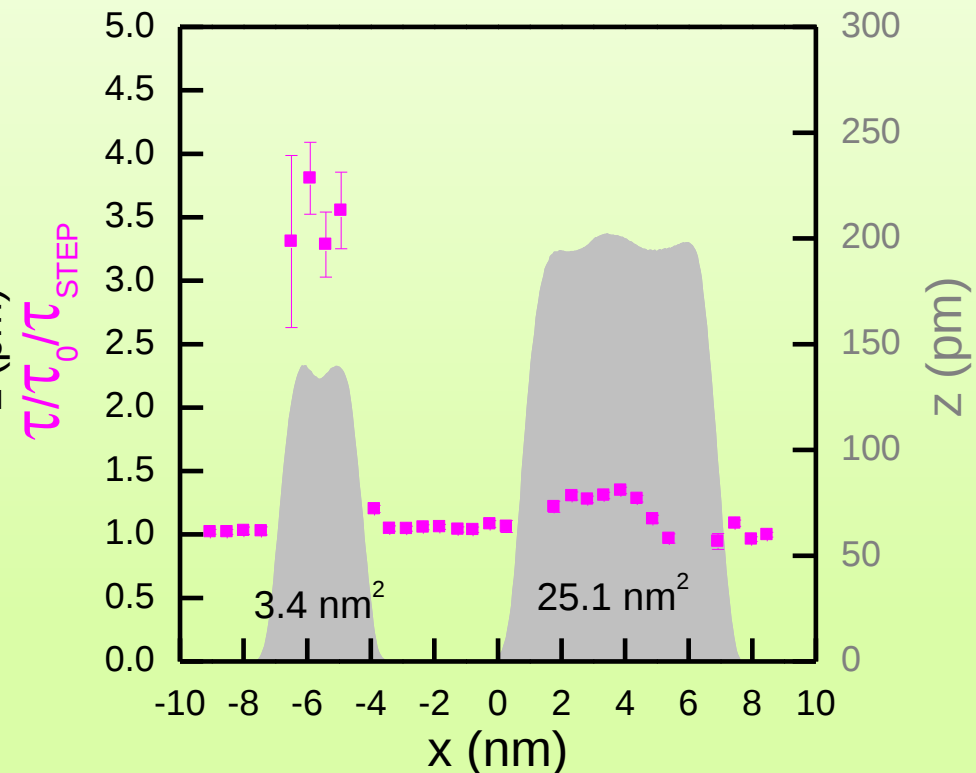
Local dependence of image potential states

Energy shift



⇒ quantum interference effect
at step edge

Change in life time

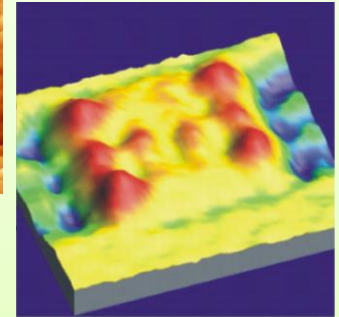
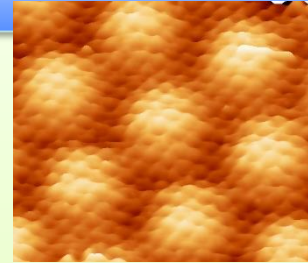


⇒ size-dependent lifetime of
electrons in image potential states

Outline

1) STM, STS, IETS, and manipulation

An extended introduction

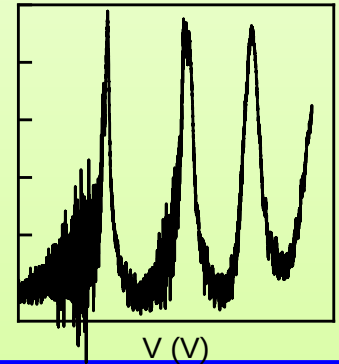


2) Quantum effects on surface imaged by STM

Electron in a box I

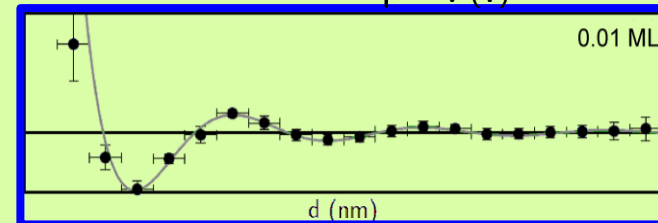
3) Quantum effects on surfaces measured by STS

Electron in a box II



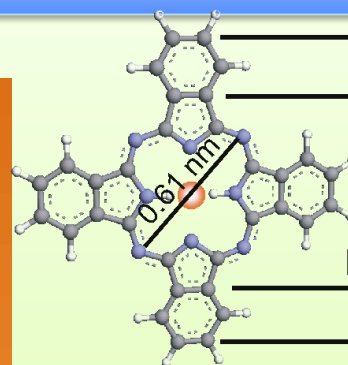
4) Electron interference effects surface diffusion

Influence of Friedel oscillations on diffusivity skip



Diffusion of molecules on surfaces

Phtalocyanine / Ag(100)



Diffusion studies (KM et al.):

Phys. Stat. Sol. B 242, 773 (2005)

Nano Lett. 23, 4793 (2023)

Angew. Chem. Int. Ed. 61 (2022)

e202212245

Nano Lett. 22, 340 (2022)

Nano Lett. 19, 710 (2019)

Nano Lett. 16, 3001 (2016)

J. Phys. Chem. Lett. 6, 4165 (2015)

[J. Am. Chem. Soc. 137, 14920 \(2015\)](#)

ACS Nano 9, 3572 (2015)

Phys. Rev. Lett. 114, 146104 (2015)

Phys. Rev. Lett. 107, 046101 (2011)

Phys. Rev. Lett. 104, 076101 (2010)

Phys. Rev. Lett. 100, 116104 (2008)

Phys. Rev. Lett. 94, 166104 (2005)

Phys. Rev. Lett. 93, 056102 (2004)

Phys. Rev. Lett. 86, 5739 (2001)

Phys. Rev. Lett. 83, 1613 (1999)

Phys. Rev. Lett. 80, 556 (1998)

Phys. Rev. Lett. 76, 2113 (1996)

Phys. Rev. Lett. 74, 2058 (1995)

T = 75 K

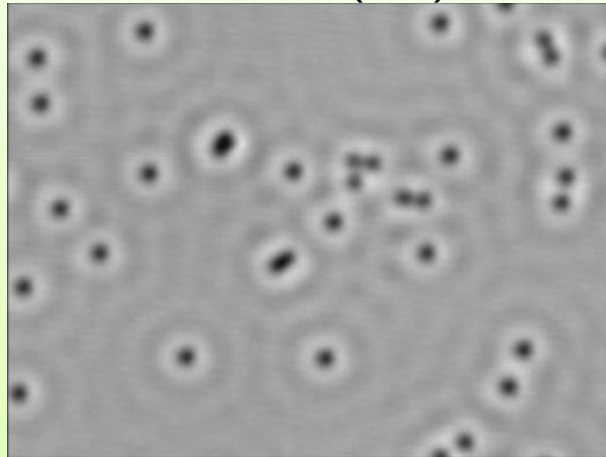
$\Delta t = 135$ s

V = 289 mV

I = 51 pA

Diffusion of molecules on surfaces

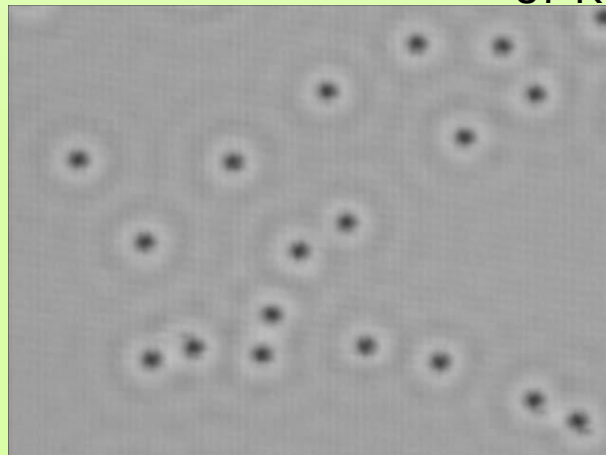
CO / Cu(111) 32 K



44pA, 200 mV, 180 s

5 nm

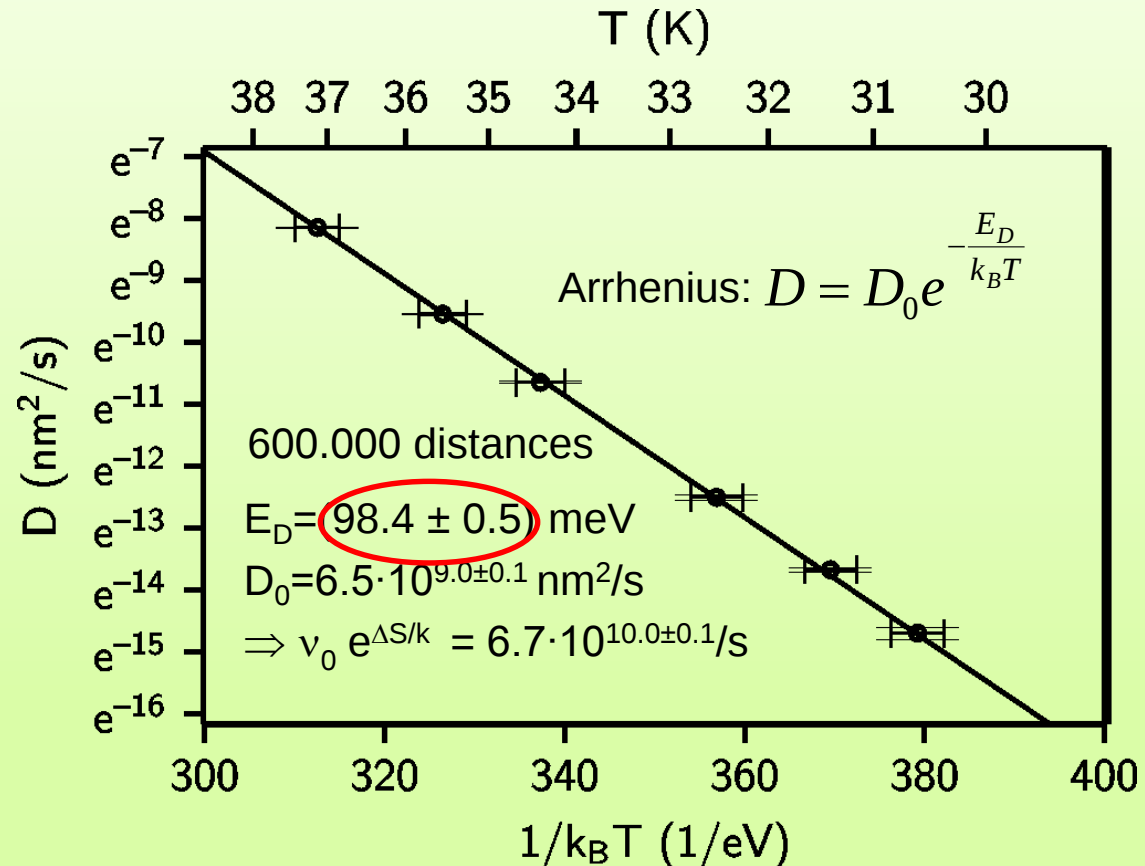
37 K



44pA, 200 mV, 180 s

5 nm

Phys. Rev. Lett.114, 146104 (2015)



Helium-Spin Echo spectroscopy:

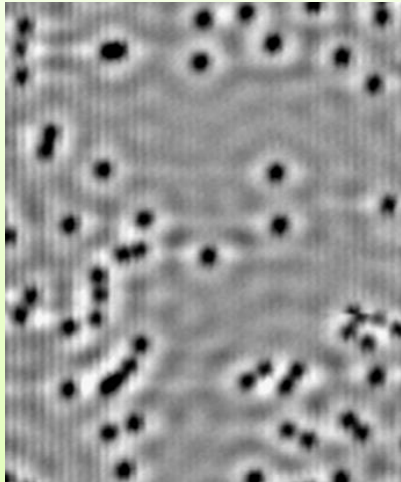
$$E_D = 98 \pm 5 \text{ meV}$$

P. R. Kole et al., J. Phys. Condens. Matter 24, 104016 (2012).

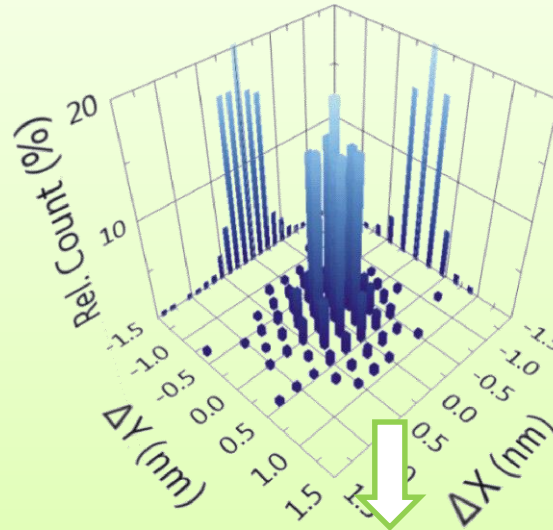
Influence of neighbors on diffusion

T = 42 K

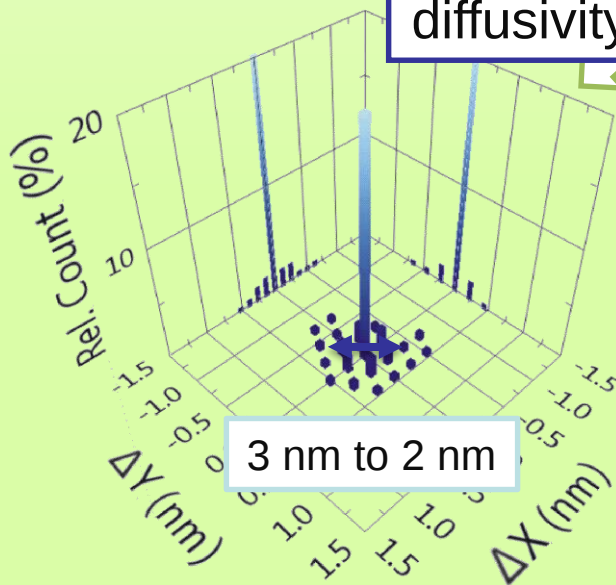
Phys. Rev. Lett. 114, 146104 (2015)
CO / Cu(111)



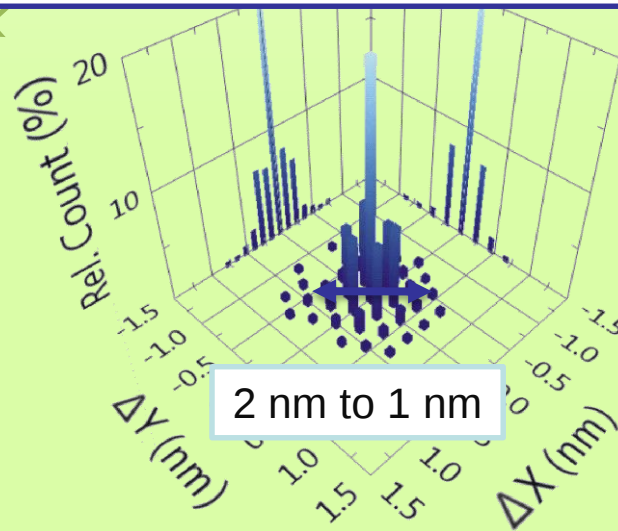
16 nm x 13 nm, 50 pA, 230 mV
 $\Delta t = 140$ s



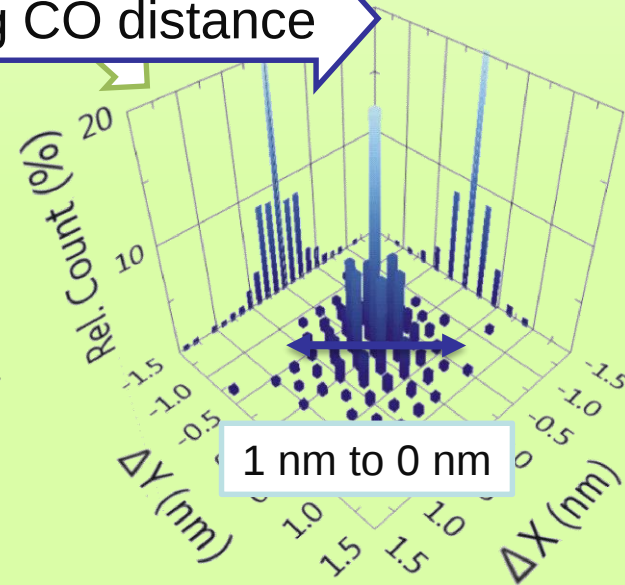
diffusivity increases with decreasing CO distance



3 nm to 2 nm



2 nm to 1 nm

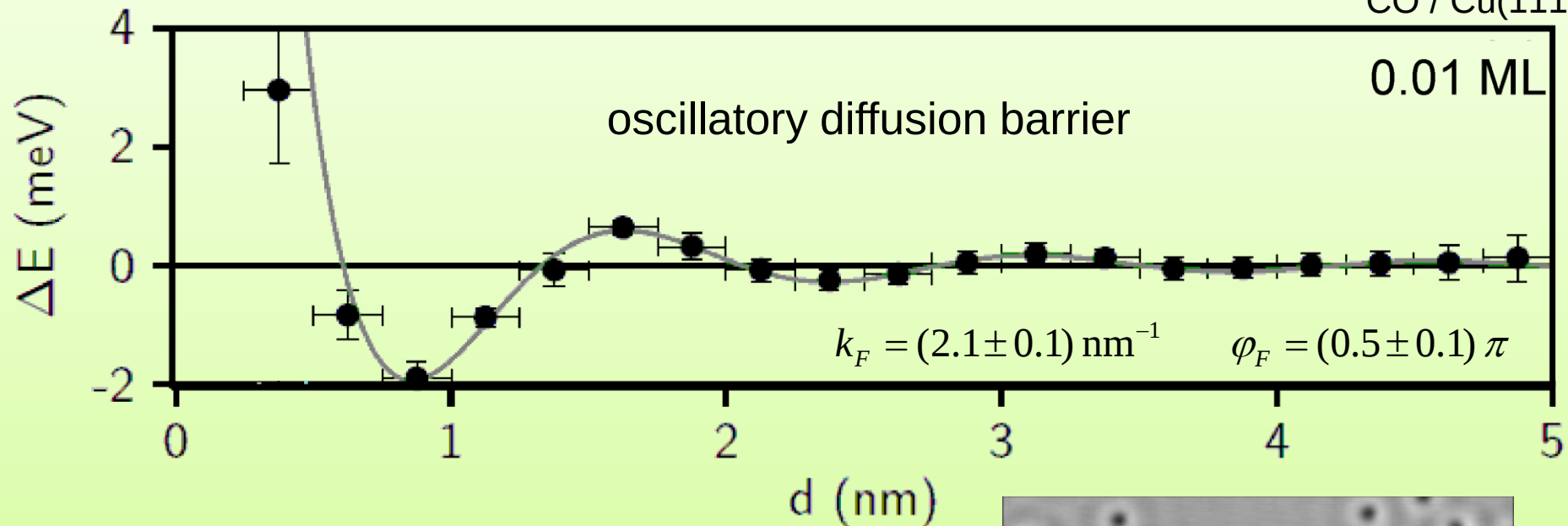


1 nm to 0 nm

Diffusion of CO on Cu(111)

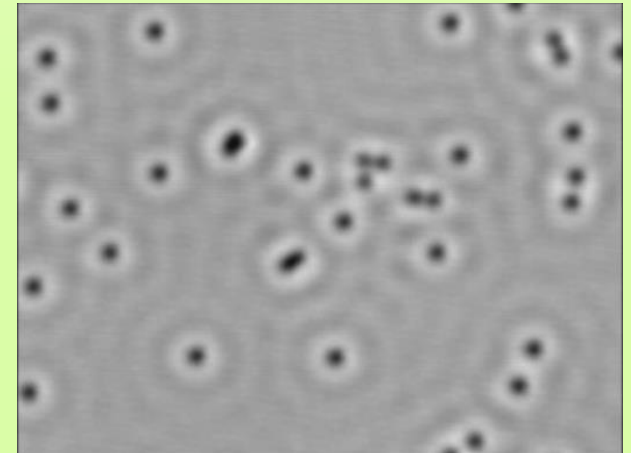
Phys. Rev. Lett. 114, 146104 (2015)

CO / Cu(111)



$$\Delta E(R) = -\alpha \cdot \frac{\sin(2k_F R + 2\phi)}{(k_F R)^2}$$

- ⇒ Friedel oscillations in diffusivity
- ⇒ long-range interaction mediated by surface state electrons



Thanks to

Experiments:

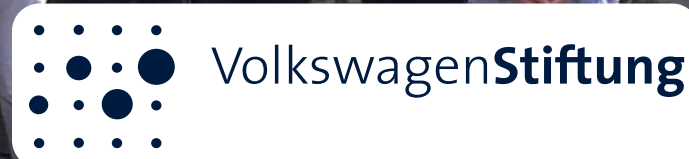
Grazyna Antzcak
Yunjun Cao
Heiko Gawronski
Sarah Heidorn
Jörg Henzl
Ting-Chieh Hung
Karsten Lucht
Michael Mehlhorn
Violeta Simic-Vilosevic
Kai Volgmann
Christopher Zaum

Theory:

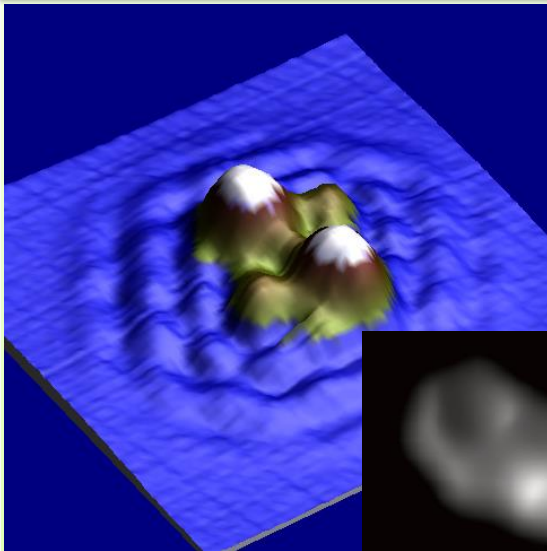
Javier Carrasco, Angelos Michaelides, UCL, UK
Galina Rusina, Svetlana Borisova, Tomsk, Russia
Evgueni Chulkov, San Sebastian
Joel Mieres-Perez, Elsa Sanchez-Garcia, UDE, TUD, D
Dave Austin, Eric Switzer, Talat Raman, UCF, USA

Molecules:

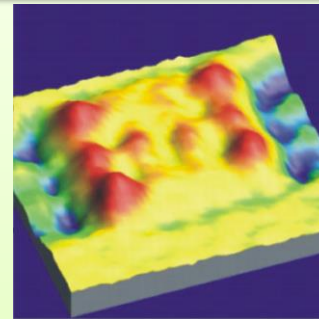
Julien Rowen, Wolfram Sander, R.U.B., D



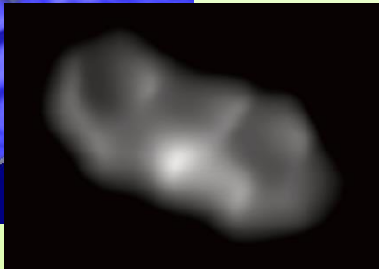
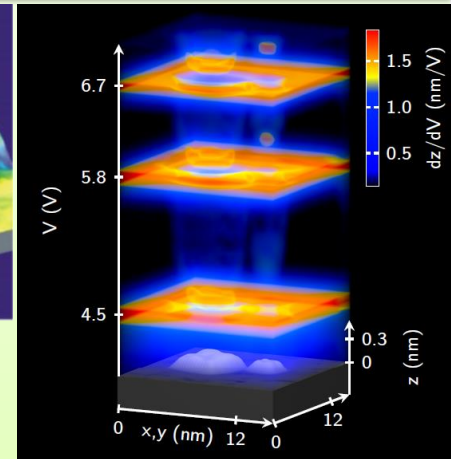
Summary



STM images present a convolution of topography, electronic density of states, and Pauli repulsion

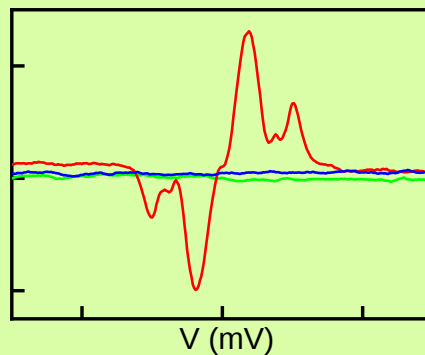
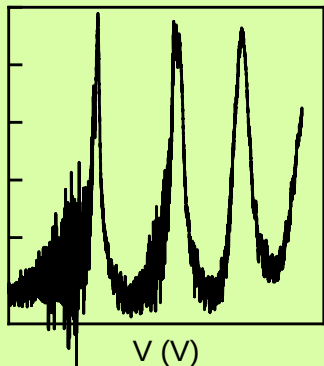


Particles in a box

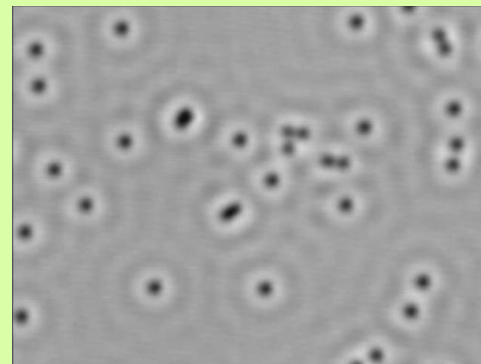


STS measures electronic states

IETS measures phonons and vibrations



⇒ fascinating view of the nanoworld by quantum mechanics based STM



Friedel-like oscillations influence surface processes