

Probing the Electronic Structure of Complex Systems by State-of-the-Art ARPES

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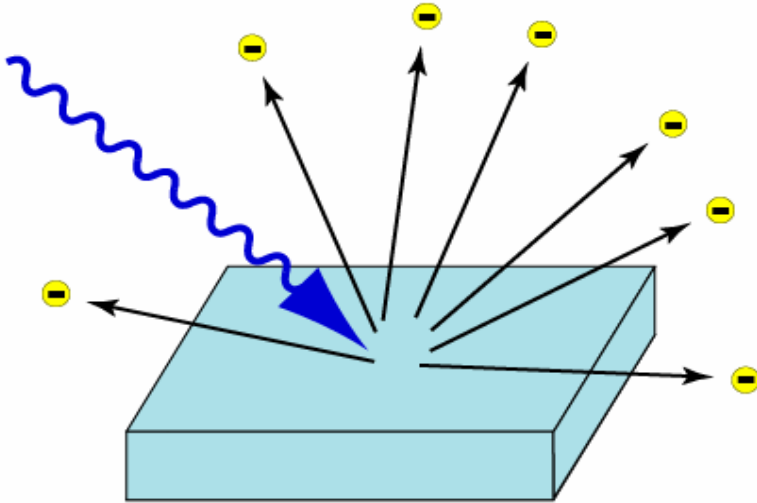
Outline: Part I

State-of-the-Art **ARPES**: the essentials

- ▶ **Motivation and potential**
- ▶ **Formal description**
 - One-step vs three-step model
 - The sudden approximation
 - Kinematics of photoemission
 - One-particle spectral function
- ▶ **Experimental**
 - State-of-art ARPES
 - Surface vs bulk sensitivity
- ▶ **Summary and discussion**

History of Photoemission

The Photoelectric Effect



- First experimental work performed by H. Hertz (1886), W. Hallwachs (1888), von Lenard (1900)
- Theoretical explanation by Einstein (1905)

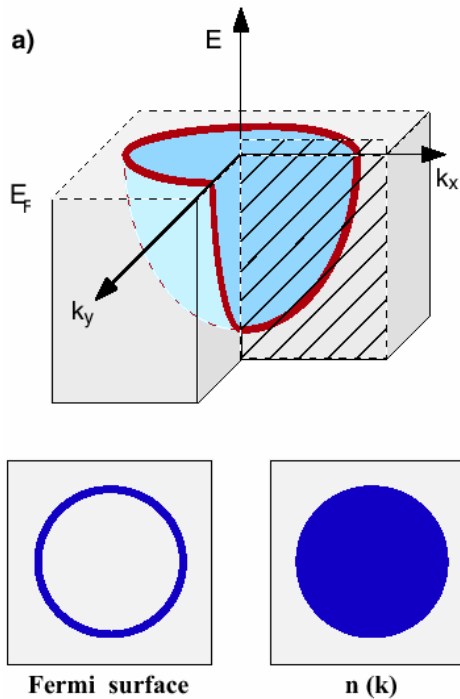
**FIRST EXPERIMENTAL EVIDENCE
FOR QUANTIZATION OF LIGHT!**

Is there anything else we can learn from the photoelectric effect?

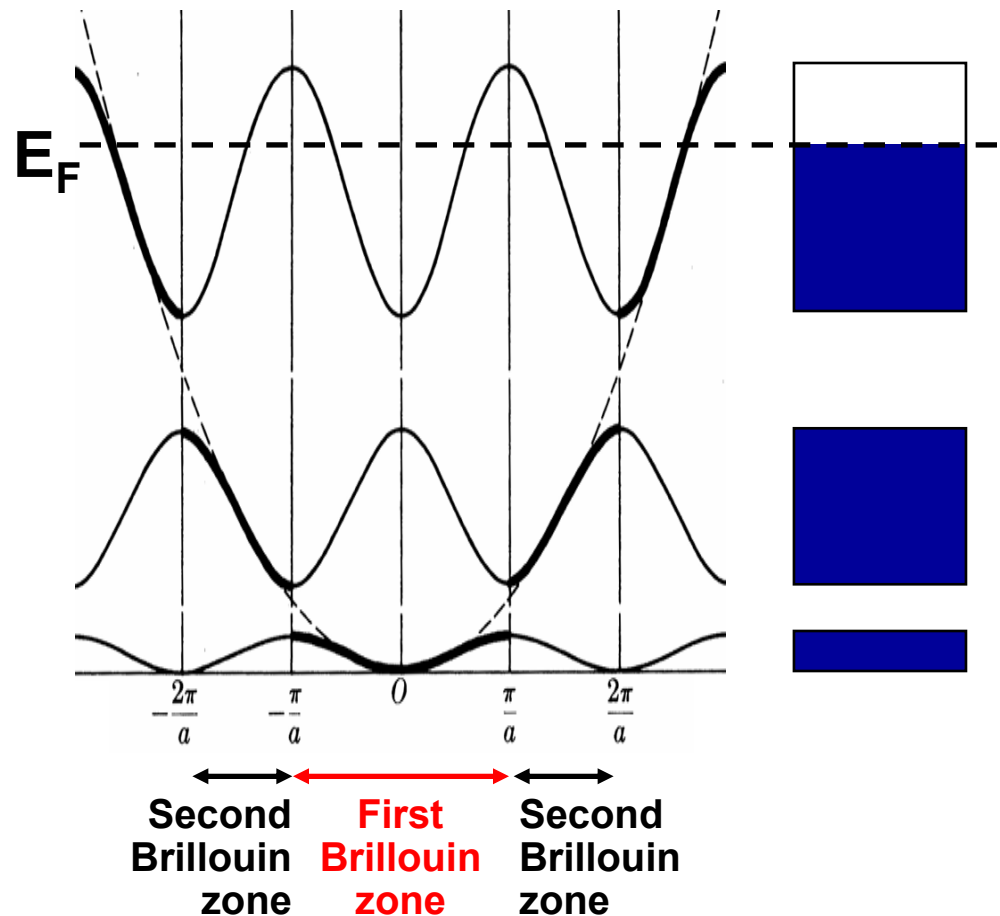
Insights into the solid-state!

Understanding the Solid State: Electrons in Reciprocal Space

Many **properties** of a solids are determined by **electrons near E_F** (**conductivity, magnetoresistance, superconductivity, magnetism**)



Allowed electronic states Repeated-zone scheme



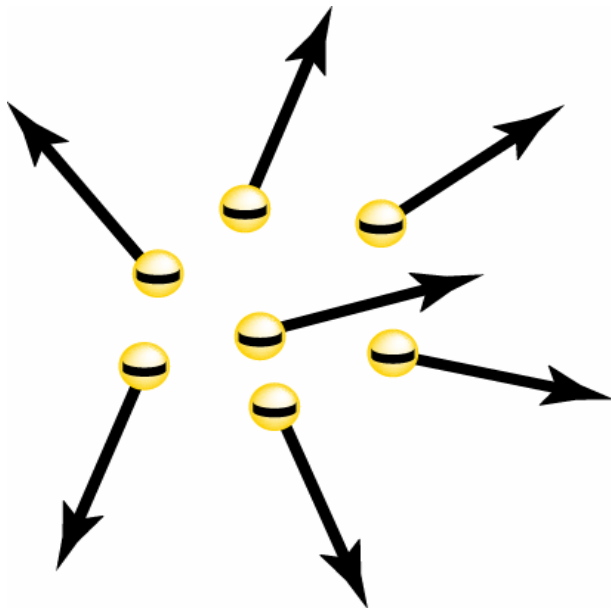
Only a **narrow energy slice** around E_F is relevant for these properties (**$kT=25$ meV** at room temperature)

Interaction Effects between Electrons : “Many-body Physics”

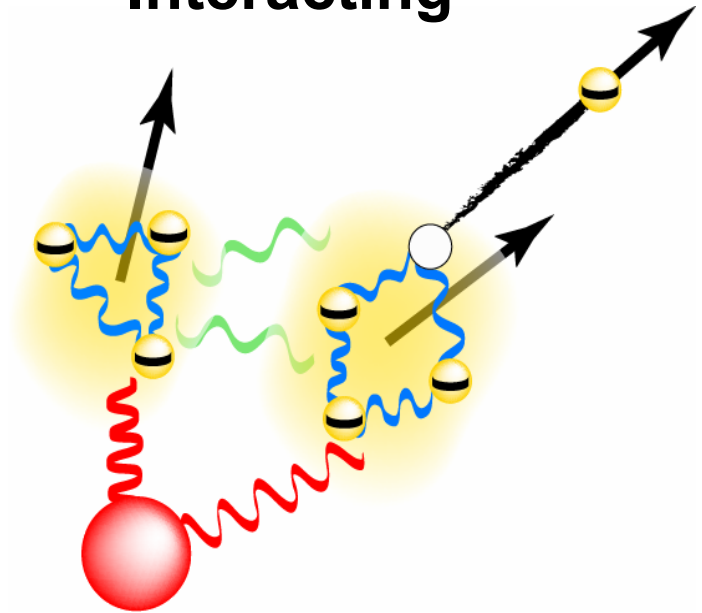
Many-body effects are due to the interactions between the **electrons** and **each other**, or with other **excitations inside the crystal** :

- 1) A “many-body” problem : intrinsically hard to calculate and understand
- 2) Responsible for many surprising phenomena :
Superconductivity, Magnetism, Density Waves,

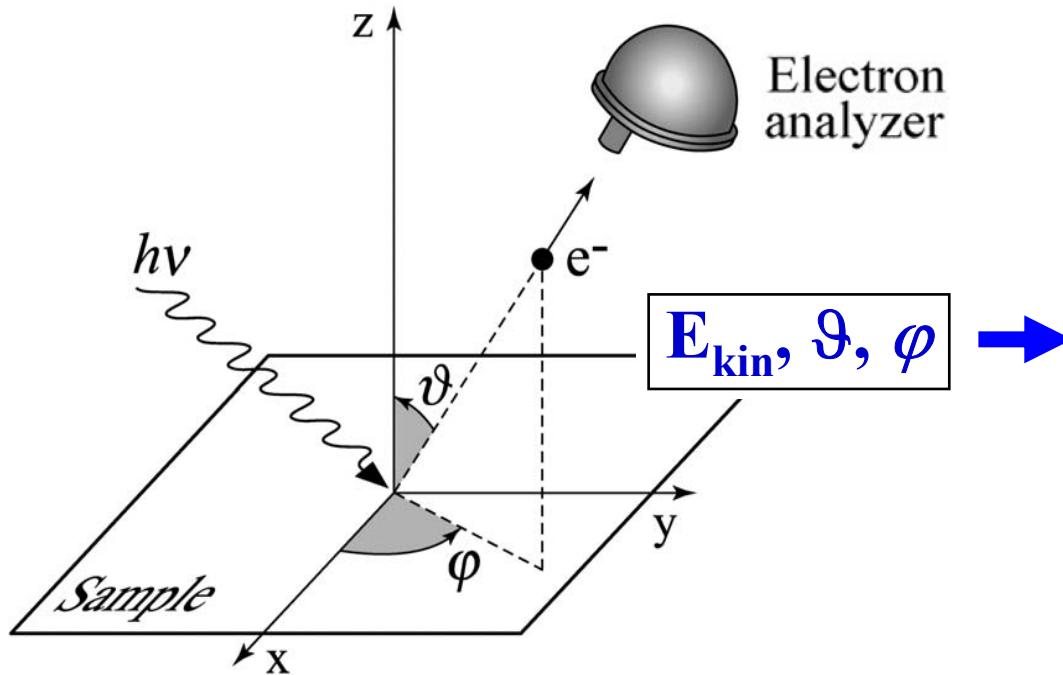
Non-Interacting



Interacting



Angle-Resolved Photoemission Spectroscopy



$$\mathbf{K} = \mathbf{p} / \hbar = \sqrt{2mE_{kin}} / \hbar$$

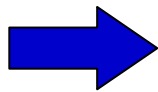
$$K_x = \frac{1}{\hbar} \sqrt{2mE_{kin}} \sin \vartheta \cos \varphi$$

$$K_y = \frac{1}{\hbar} \sqrt{2mE_{kin}} \sin \vartheta \sin \varphi$$

$$K_z = \frac{1}{\hbar} \sqrt{2mE_{kin}} \cos \vartheta$$

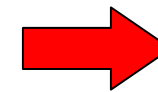
Vacuum

$$\begin{matrix} E_{kin} \\ \mathbf{K} \end{matrix}$$



Conservation laws

$$\begin{matrix} E_f - E_i = h\nu \\ \mathbf{k}_f - \mathbf{k}_i = \cancel{\mathbf{k}_{h\nu}} \end{matrix}$$



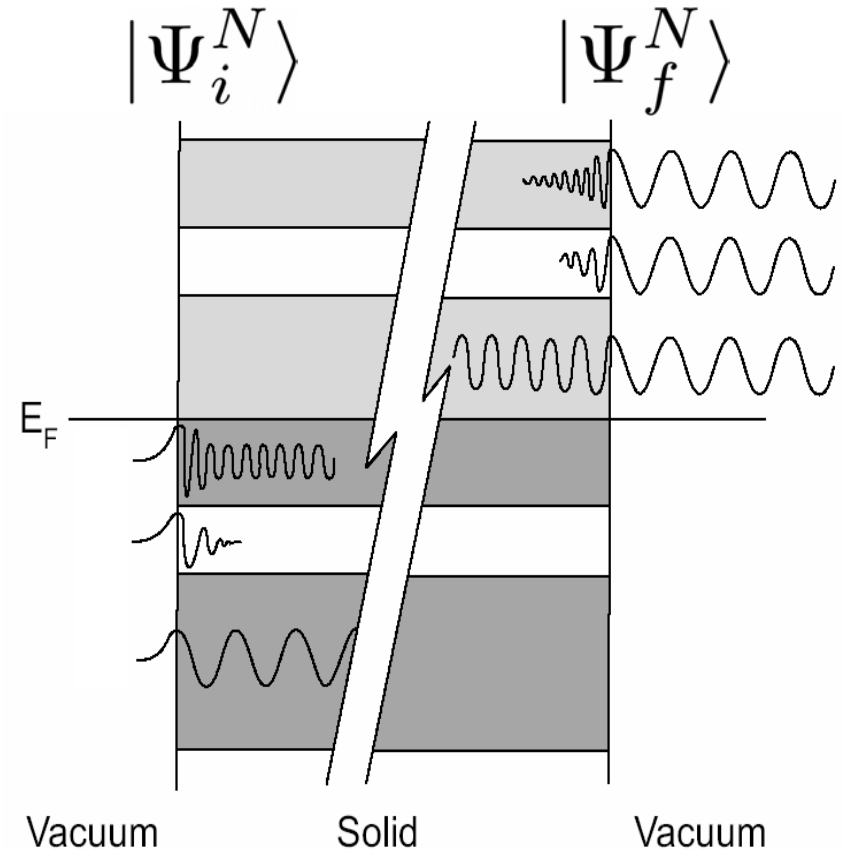
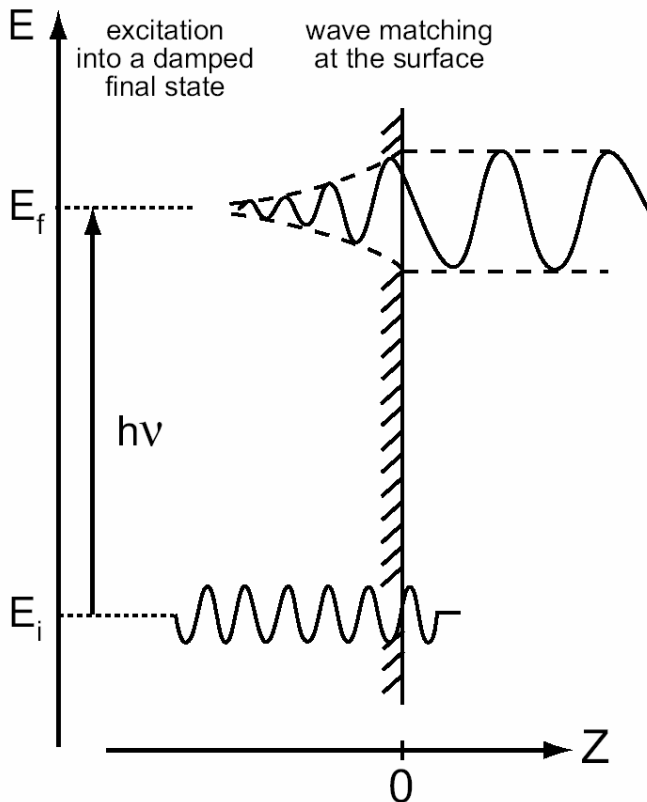
Solid

$$\begin{matrix} E_B \\ \mathbf{k} \end{matrix}$$

ARPES: One-Step vs Three-Step Model

$$\left. \begin{array}{l} \text{Photoemission} \\ \text{Intensity } I(k, \omega) \end{array} \right\} w_{fi} \propto |\langle \Psi_f^N | \mathbf{A} \cdot \mathbf{p} | \Psi_i^N \rangle|^2 \delta(E_f^N - E_i^N - h\nu)$$

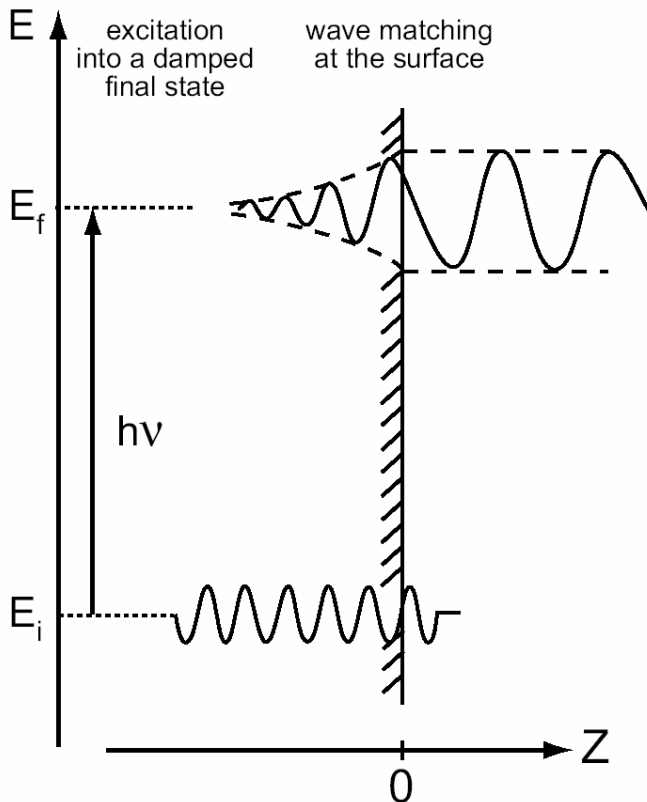
One-step model



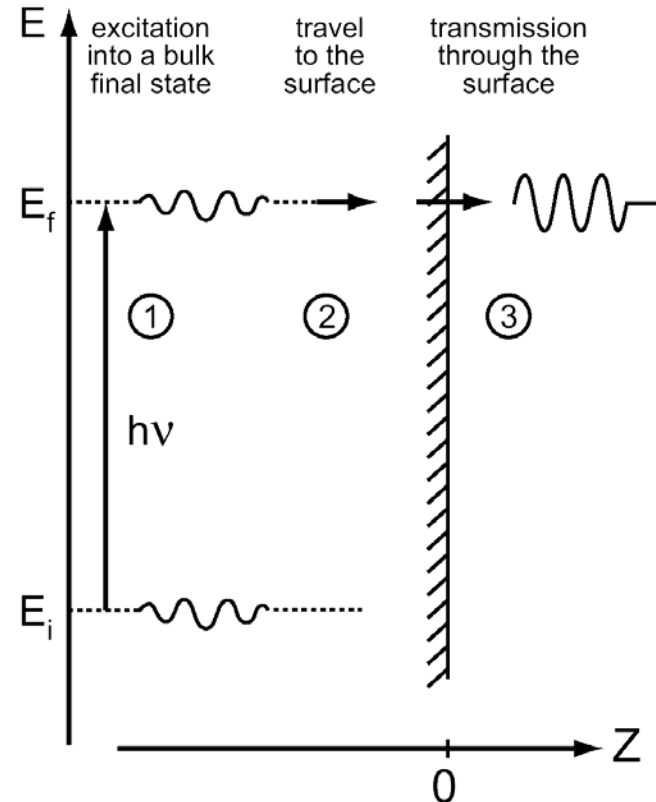
ARPES: One-Step vs Three-Step Model

$$\left. \begin{array}{l} \text{Photoemission} \\ \text{Intensity } I(k, \omega) \end{array} \right\} w_{fi} \propto |\langle \Psi_f^N | \mathbf{A} \cdot \mathbf{p} | \Psi_i^N \rangle|^2 \delta(E_f^N - E_i^N - h\nu)$$

One-step model



Three-step model

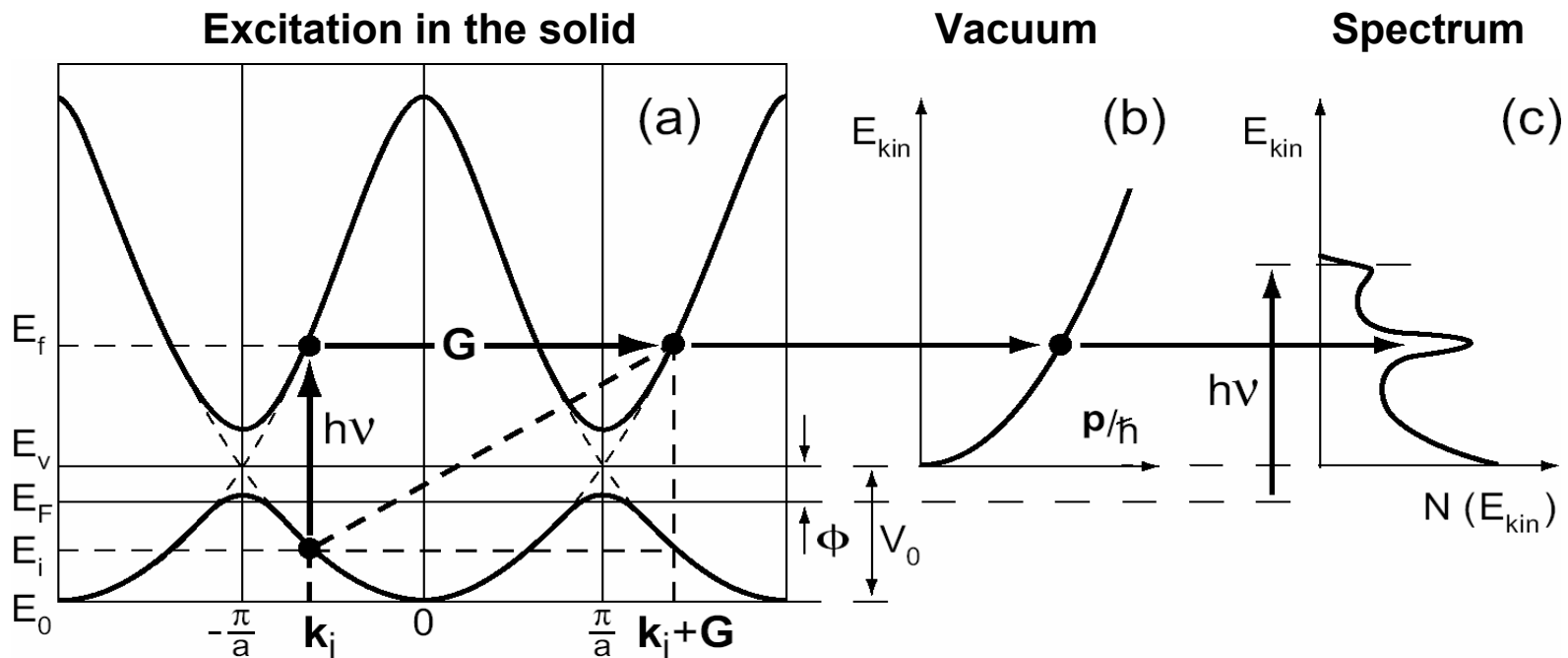


ARPES: The Sudden Approximation

$$\left. \begin{array}{l} \text{Photoemission} \\ \text{Intensity } I(\mathbf{k}, \omega) \end{array} \right\} w_{fi} \propto |\langle \phi_f^{\mathbf{k}} | \mathbf{A} \cdot \mathbf{p} | \phi_i^{\mathbf{k}} \rangle \langle \Psi_m^{N-1} | \Psi_i^{N-1} \rangle|^2 \delta(\omega - h\nu)$$

$$\Psi_f^N: \text{Sudden approximation} \rightarrow \Psi_f^N = \mathcal{A} \phi_f^{\mathbf{k}} \Psi_f^{N-1}$$

$$\Psi_i^N: \text{One Slater determinant} \rightarrow \Psi_i^N = \mathcal{A} \phi_i^{\mathbf{k}} \Psi_i^{N-1}$$

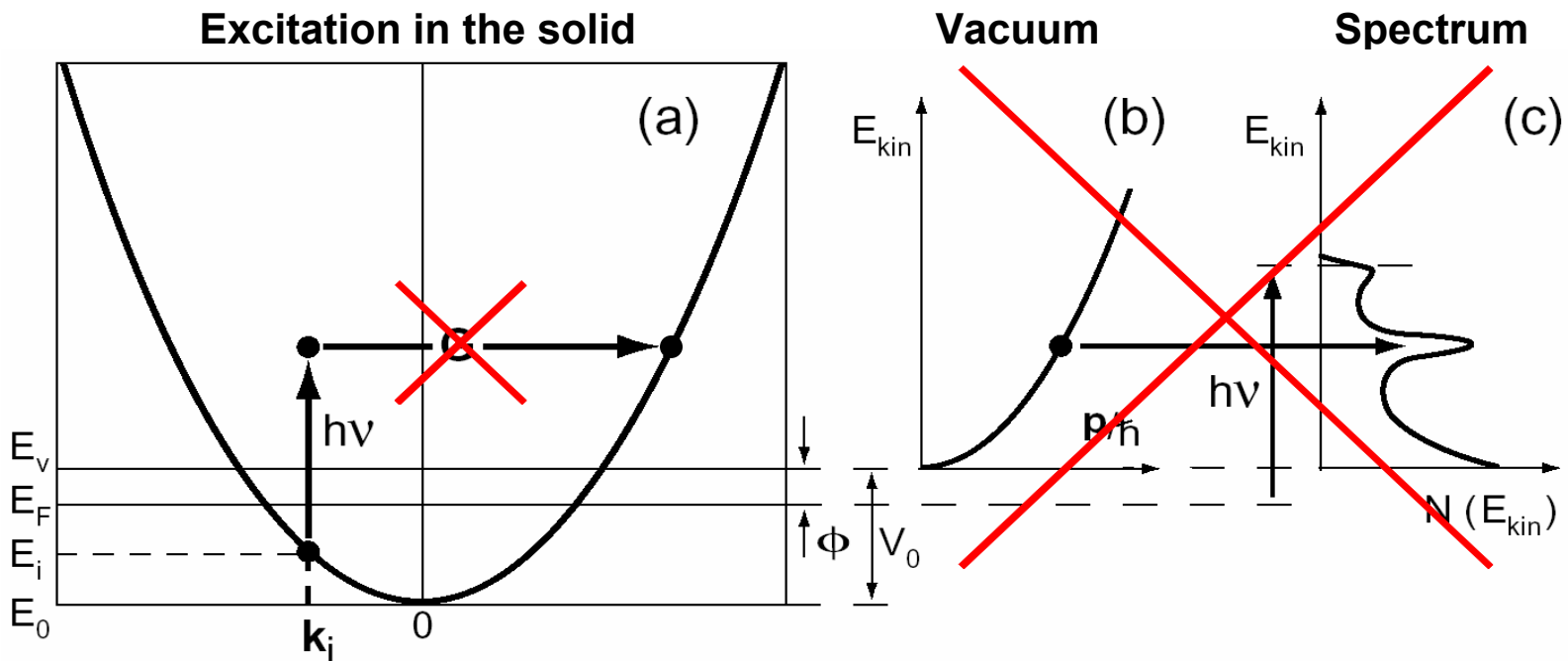


ARPES: Role of the Crystal Potential

Photoemission Intensity $I(k, \omega)$ } $w_{fi} \propto |\langle \phi_f^{\mathbf{k}} | \underline{\mathbf{A}} \cdot \nabla V | \phi_i^{\mathbf{k}} \rangle \langle \Psi_m^{N-1} | \Psi_i^{N-1} \rangle|^2 \delta(\omega - h\nu)$

Ψ_f^N : **Sudden approximation** $\rightarrow \Psi_f^N = \mathcal{A} \phi_f^{\mathbf{k}} \Psi_f^{N-1}$

Ψ_i^N : **One Slater determinant** $\rightarrow \Psi_i^N = \mathcal{A} \phi_i^{\mathbf{k}} \Psi_i^{N-1}$

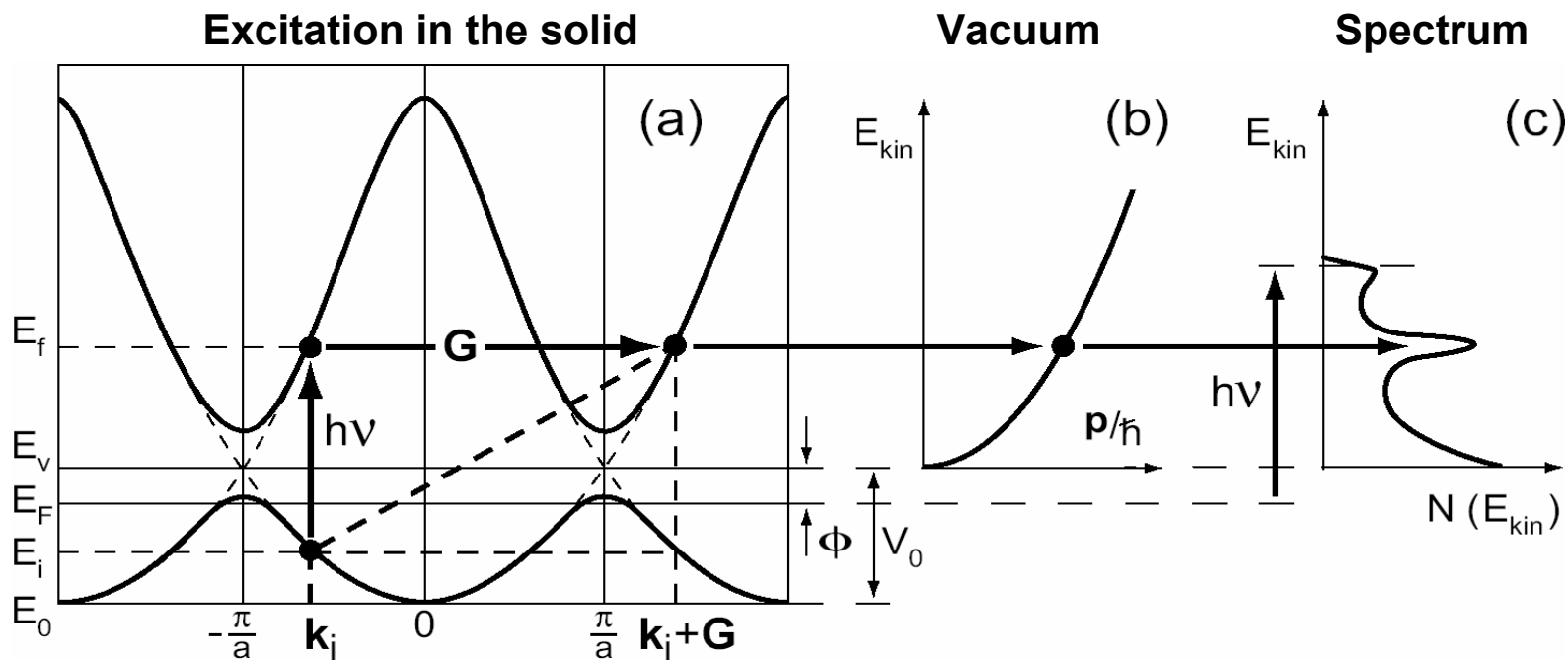


ARPES: Inner Potential and Determination of k_z

Free-electron final state $E_f(\mathbf{k}) = \frac{\hbar^2 \mathbf{k}^2}{2m} - |E_0| = \frac{\hbar^2 (\mathbf{k}_{\parallel}^2 + \mathbf{k}_{\perp}^2)}{2m} - |E_0|$

because $\frac{\hbar^2 \mathbf{k}_{\parallel}^2}{2m} = E_{kin} \sin^2 \vartheta$ $E_f = E_{kin} + \phi$ $V_0 = |E_0| + \phi$

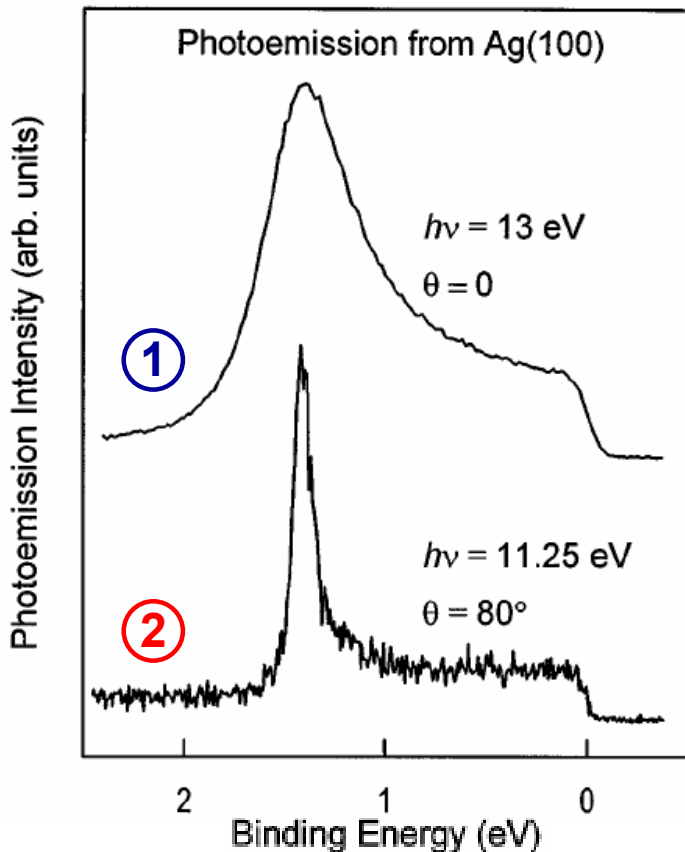
➔
$$\mathbf{k}_{\perp} = \frac{1}{\hbar} \sqrt{2m(E_{kin} \cos^2 \vartheta + V_0)}$$



ARPES: FWHM and Inverse Lifetime

FWHM of an ARPES peak }
$$\Gamma = \frac{\frac{\Gamma_i}{|v_{i\perp}|} + \frac{\Gamma_f}{|v_{f\perp}|}}{\left| \frac{1}{v_{i\perp}} \left[1 - \frac{mv_{i\parallel} \sin^2 \vartheta}{\hbar k_{\parallel}} \right] - \frac{1}{v_{f\perp}} \left[1 - \frac{mv_{f\parallel} \sin^2 \vartheta}{\hbar k_{\parallel}} \right] \right|}$$

Hansen *et al.*, PRL **80**, 1766 (1998)



① if $E_i \simeq E_F$

→ $\Gamma_i \longrightarrow 0$ → $\Gamma \propto \Gamma_f$

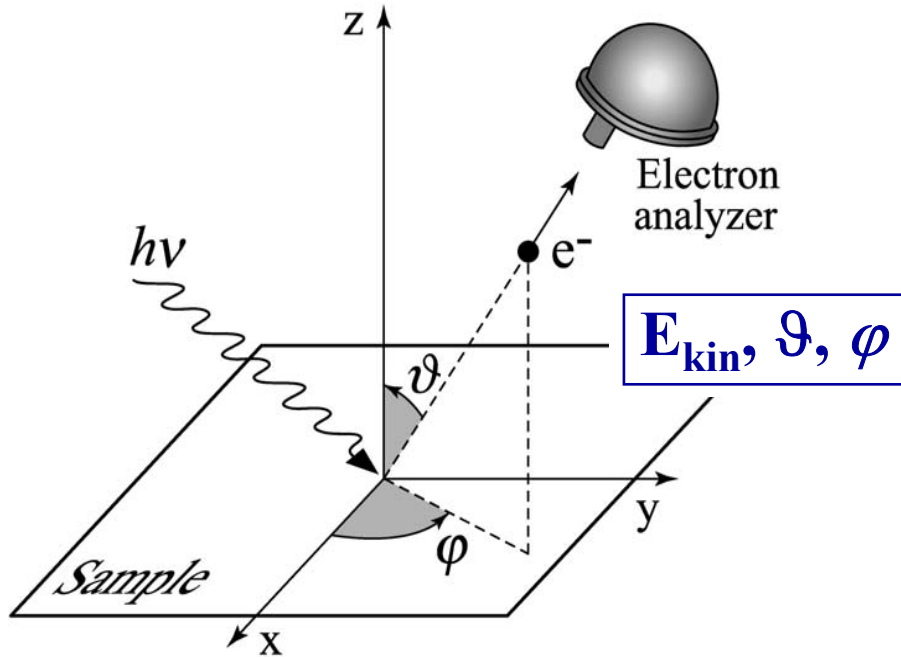
② if $|v_{i\perp}| \simeq 0$

→ $\Gamma = \frac{\Gamma_i}{\left| 1 - \frac{mv_{i\parallel} \sin^2 \vartheta}{\hbar k_{\parallel}} \right|} \equiv C \Gamma_i$

if $v_{i\parallel} < 0$, large; θ large; k_{\parallel} small

→ $C < 1$, and $\Gamma < \Gamma_i$

ARPES: Energetics and Kinematics

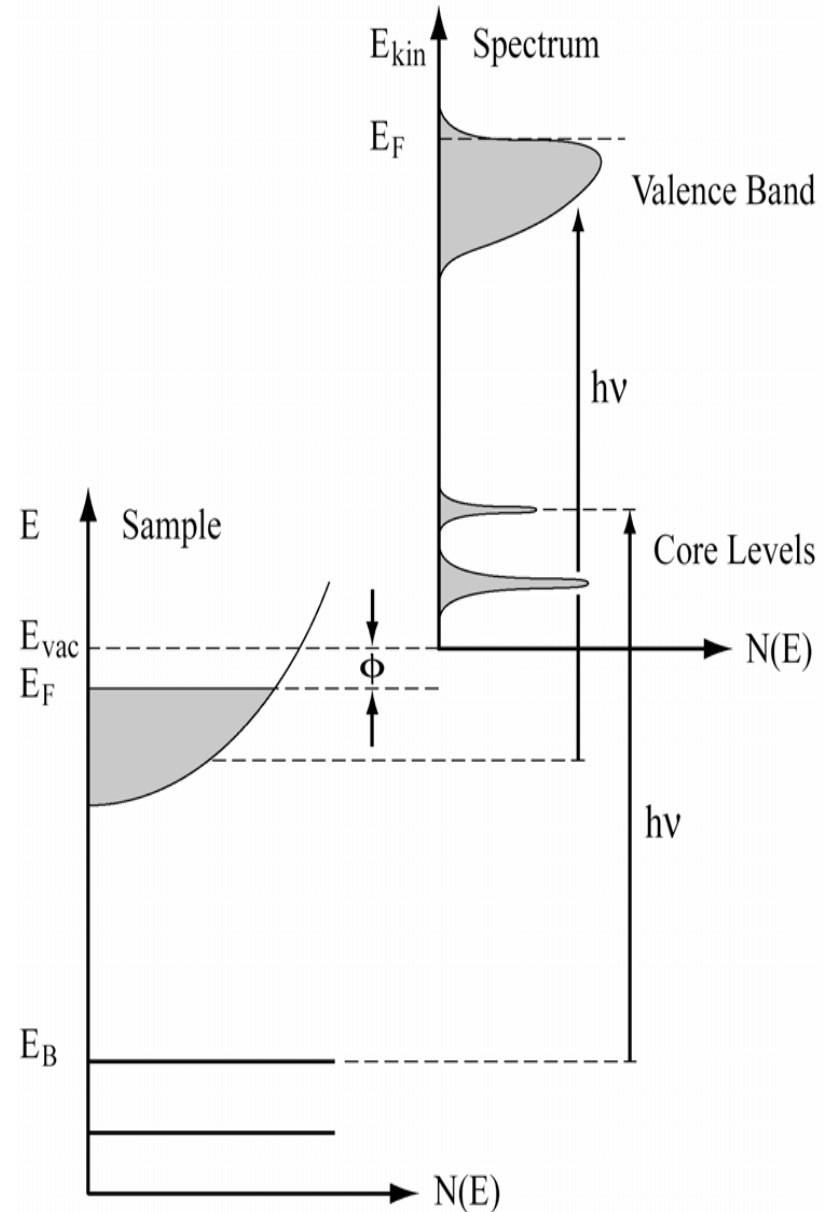


Energy Conservation

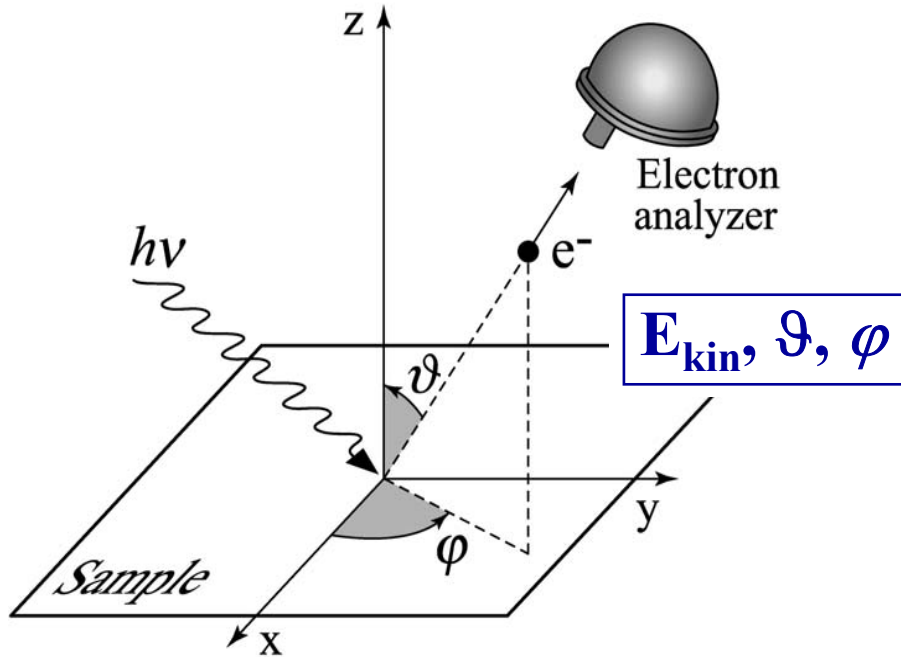
$$E_{kin} = h\nu - \phi - |E_B|$$

Momentum Conservation

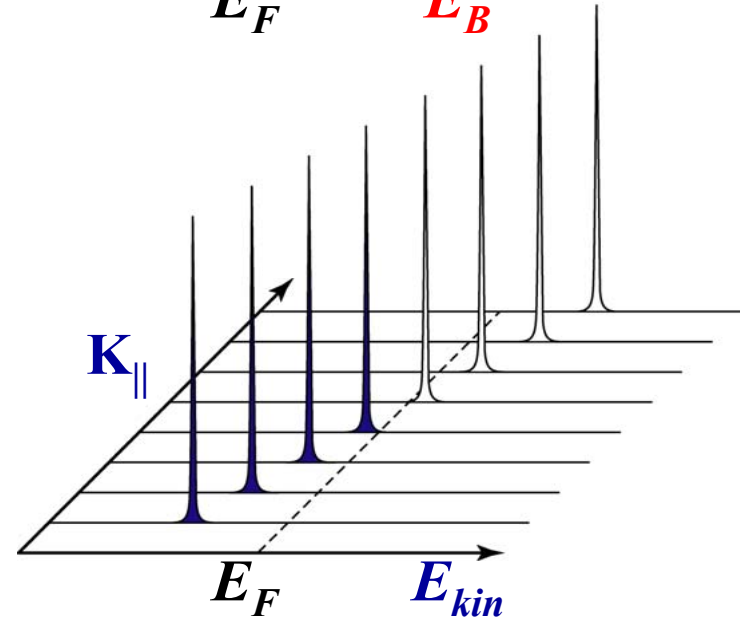
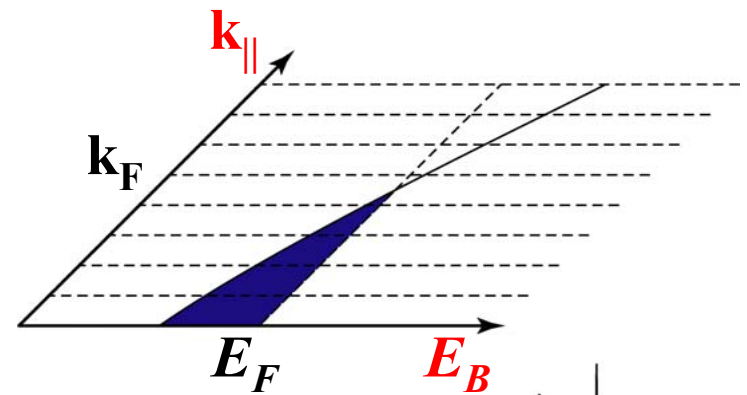
$$\hbar \mathbf{k}_{||} = \hbar \mathbf{K}_{||} = \sqrt{2m E_{kin}} \cdot \sin \theta$$



ARPES: Energetics and Kinematics



Electrons in Reciprocal Space



Energy Conservation

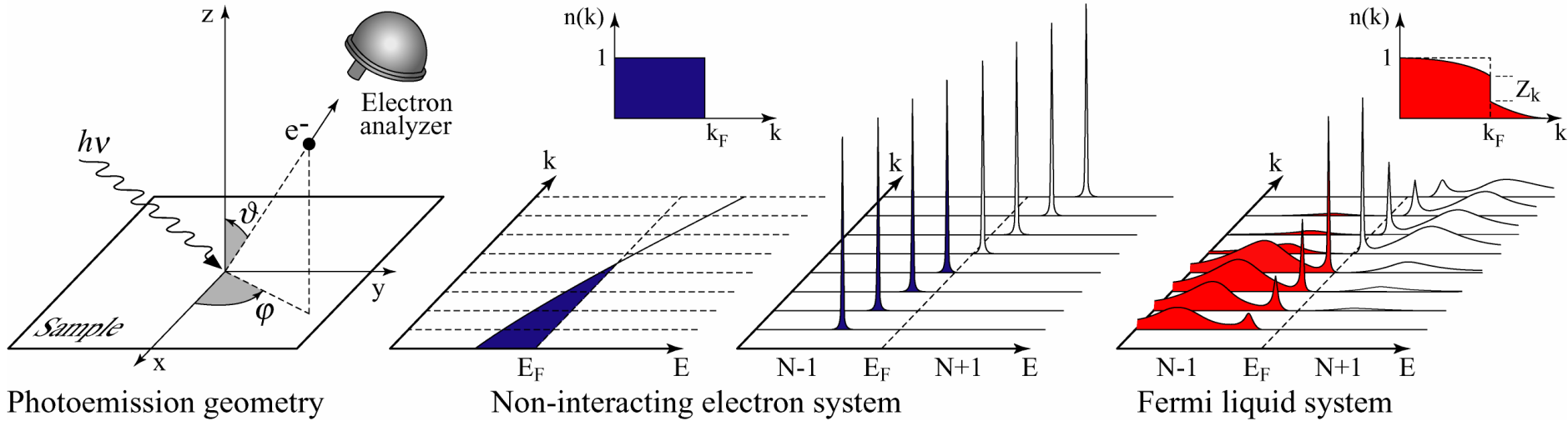
$$E_{kin} = h\nu - \phi - |E_B|$$

Momentum Conservation

$$\hbar \mathbf{k}_{\parallel} = \hbar \mathbf{K}_{\parallel} = \sqrt{2m E_{kin}} \cdot \sin \theta$$

ARPES: Interacting Systems

A. Damascelli, Z. Hussain, Z.-X Shen, Rev. Mod. Phys. **75**, 473 (2003)



Photoemission intensity: $I(\mathbf{k}, E_{kin}) = \sum_{f,i} w_{f,i}$

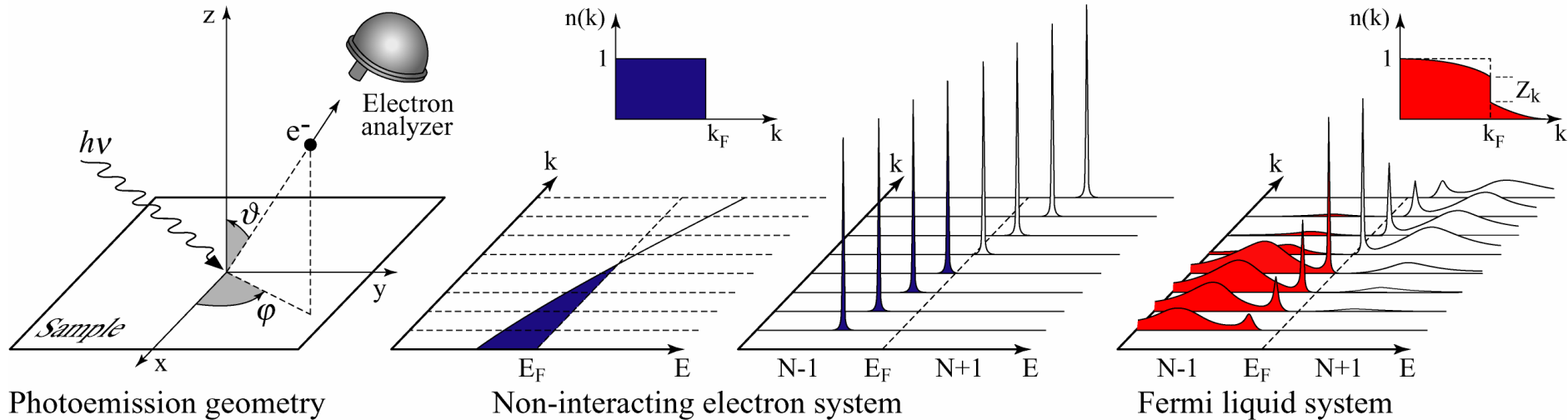
$$I(\mathbf{k}, E_{kin}) \propto \sum_{f,i} |M_{f,i}^{\mathbf{k}}|^2 \sum_m |c_{m,i}|^2 \delta(E_{kin} + E_m^{N-1} - E_i^N - h\nu)$$

$$|M_{f,i}^{\mathbf{k}}|^2 \equiv |\langle \phi_f^{\mathbf{k}} | \mathbf{A} \cdot \mathbf{p} | \phi_i^{\mathbf{k}} \rangle|^2 \quad |c_{m,i}|^2 = |\langle \Psi_m^{N-1} | \Psi_i^{N-1} \rangle|^2$$

In general $\Psi_i^{N-1} = c_{\mathbf{k}} \Psi_i^N$ **NOT orthogonal** Ψ_m^{N-1}

ARPES: The One-Particle Spectral Function

A. Damascelli, Z. Hussain, Z.-X Shen, Rev. Mod. Phys. **75**, 473 (2003)



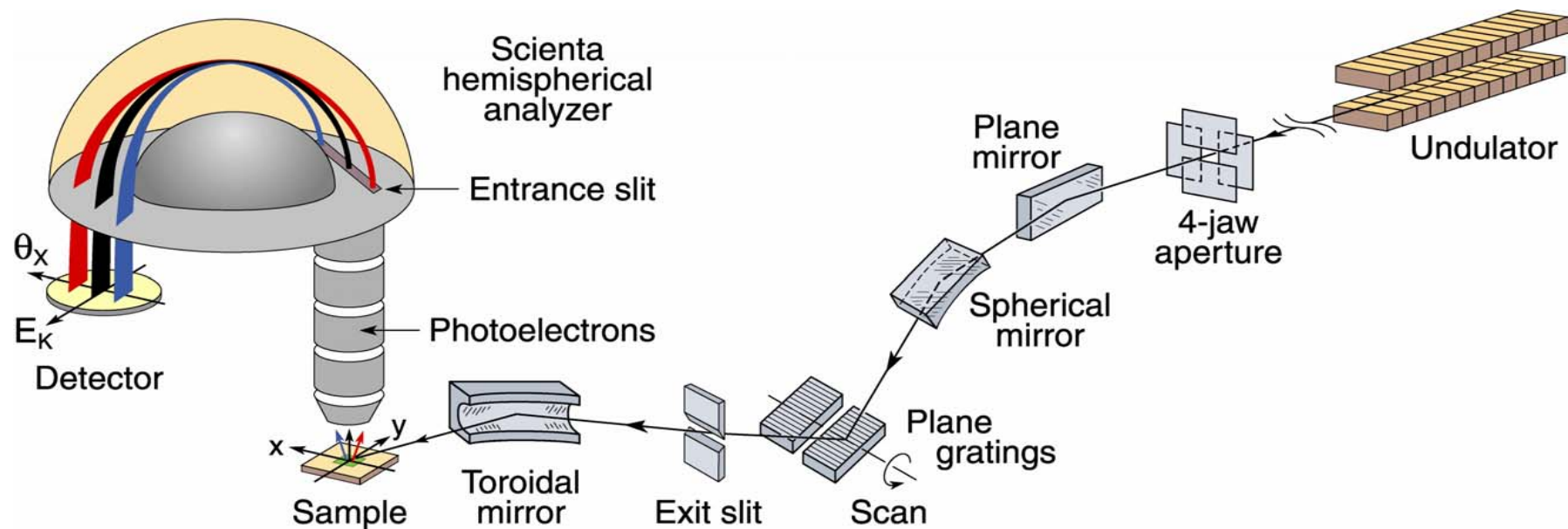
Photoemission intensity: $I(\mathbf{k}, \omega) = I_0 |M(\mathbf{k}, \omega)|^2 f(\omega) A(\mathbf{k}, \omega)$

Single-particle spectral function

$$A(\mathbf{k}, \omega) = -\frac{1}{\pi} \frac{\Sigma''(\mathbf{k}, \omega)}{[\omega - \epsilon_{\mathbf{k}} - \Sigma'(\mathbf{k}, \omega)]^2 + [\Sigma''(\mathbf{k}, \omega)]^2}$$

$S(\mathbf{k}, \omega)$: the “self-energy” captures the effects of interactions

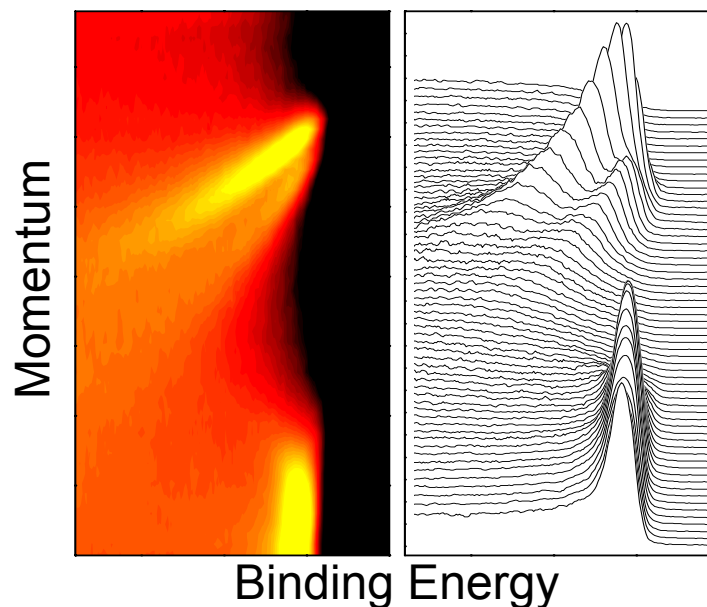
Angle-Resolved Photoemission Spectroscopy



Parallel multi-angle recording

- Improved **energy resolution**
- Improved **momentum resolution**
- Improved **data-acquisition efficiency**

	ΔE (meV)	$\Delta\theta$
past	20-40	2°
now	2-10	0.2°



SSRL Beamline 5-4 : NIM / Scienta System

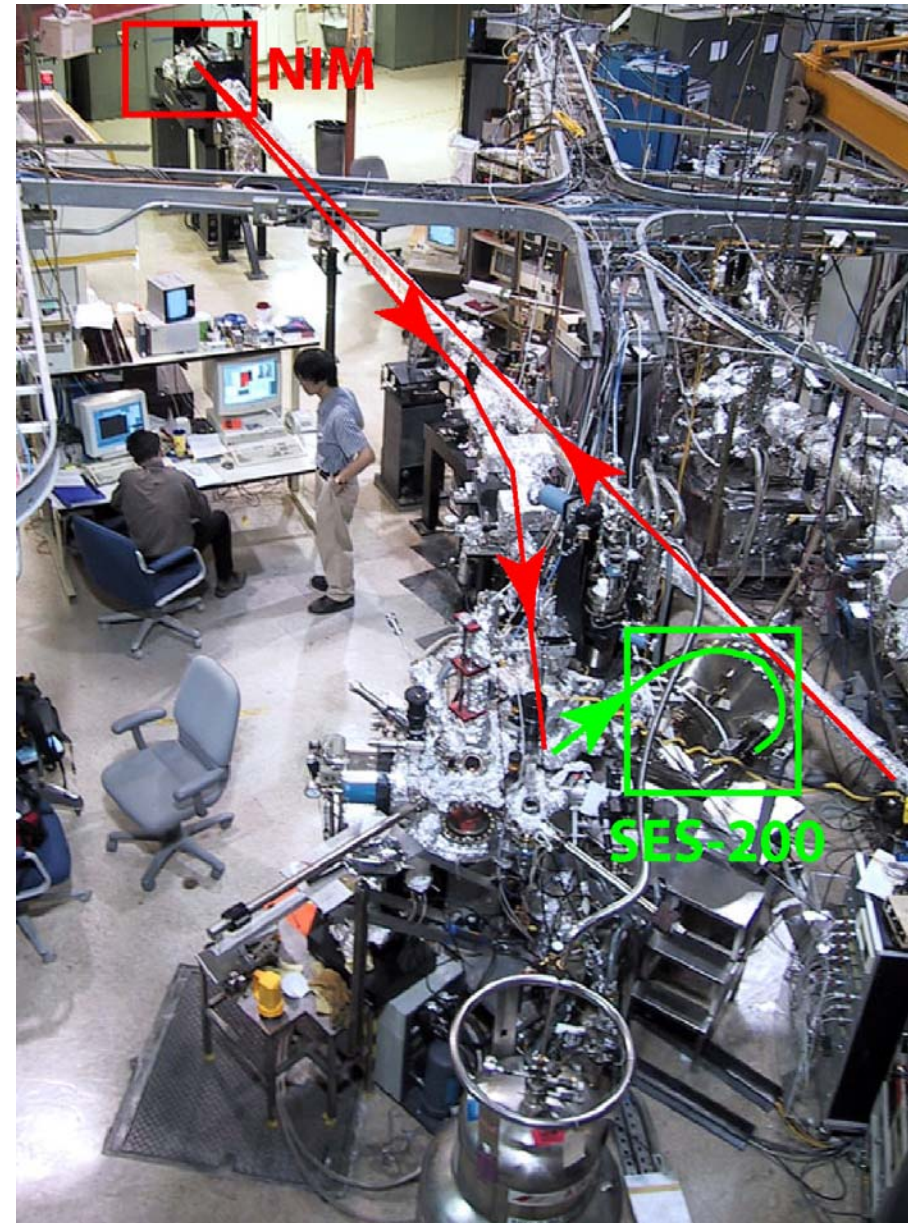
STANFORD SYNCHROTRON RADIATION LABORATORY



- High resolution

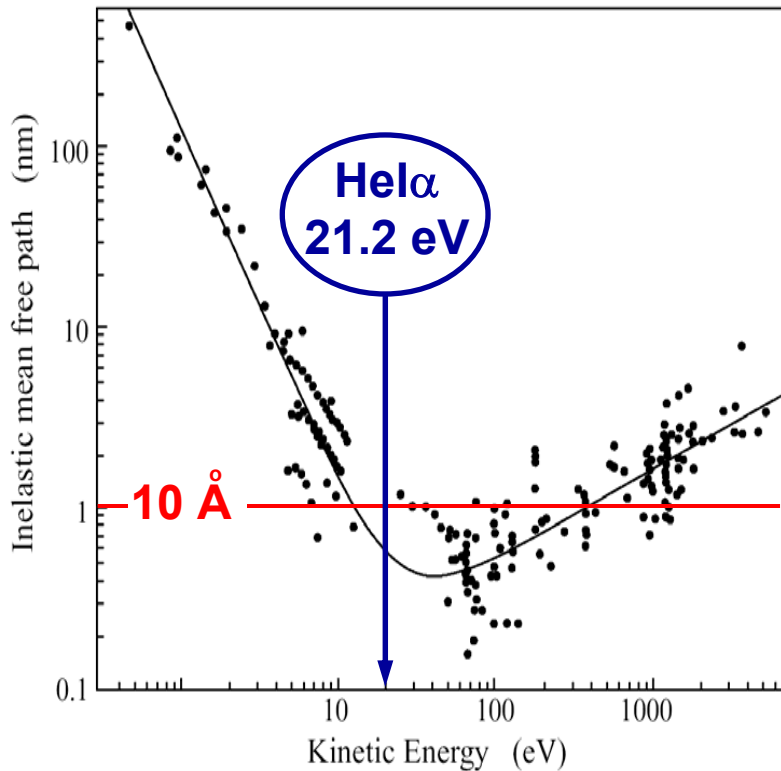
ΔE (meV)	$\Delta\theta$
2-10	0.2°

- Ultra-high vacuum ($\sim 10^{-11}$ torr)
- High angular precision ($\pm 0.1^\circ$)
- Low base temperature (< 10 K)
- Wide temperature range (10-350 K)
- Variable photon energies (12-30 eV)
- Multiple light sources (He lamp)
- Control of light polarization
- Single crystal cleaving tools
- Sample surface preparation & cleaning
- Low-Energy Electron Diffraction (LEED)

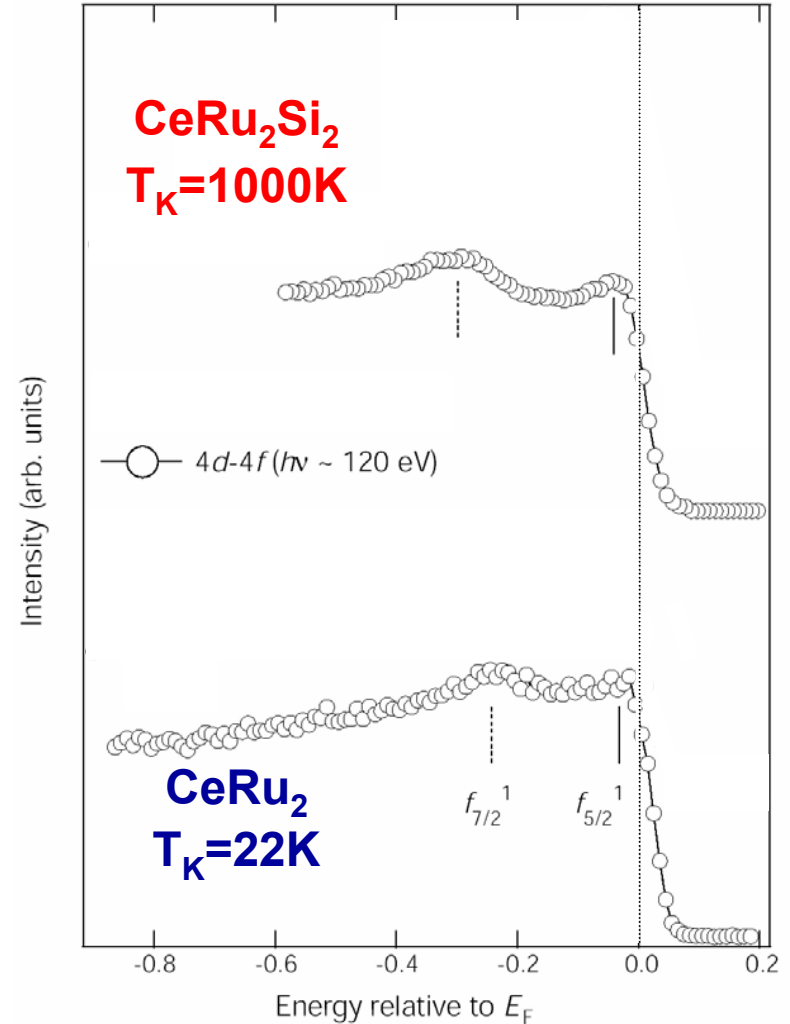


ARPES: Surface vs Bulk Sensitivity

Mean-free path for excited electrons



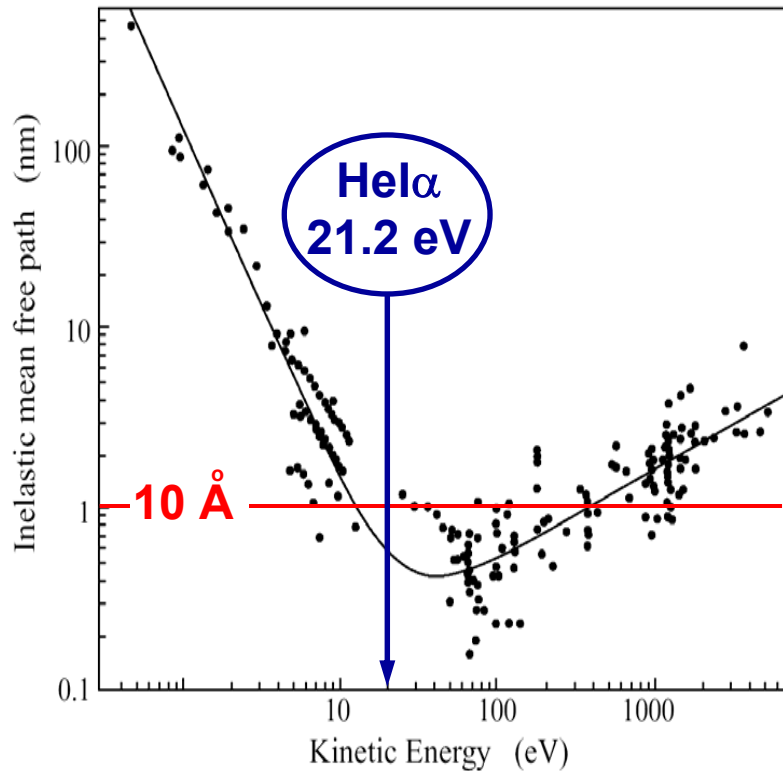
Seah, Dench *et al.*, SIA 1, 2 (1979)



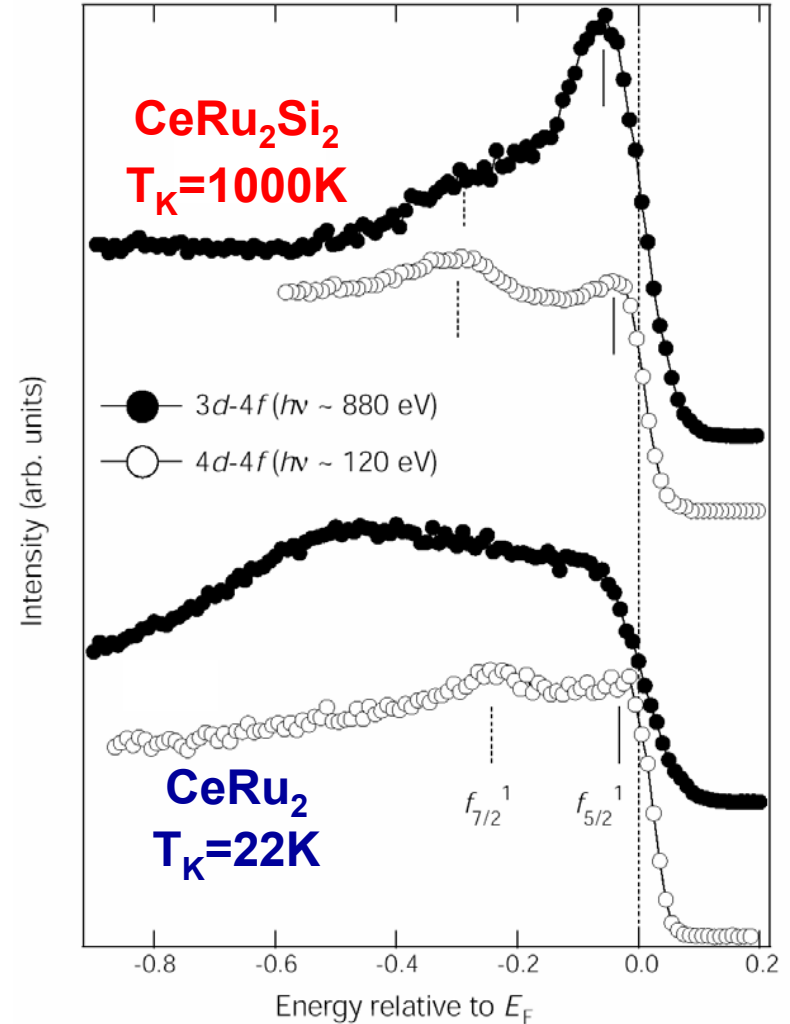
Sekiyama *et al.*, Nature **403**, 396 (2000)

ARPES: Surface vs Bulk Sensitivity

Mean-free path for excited electrons



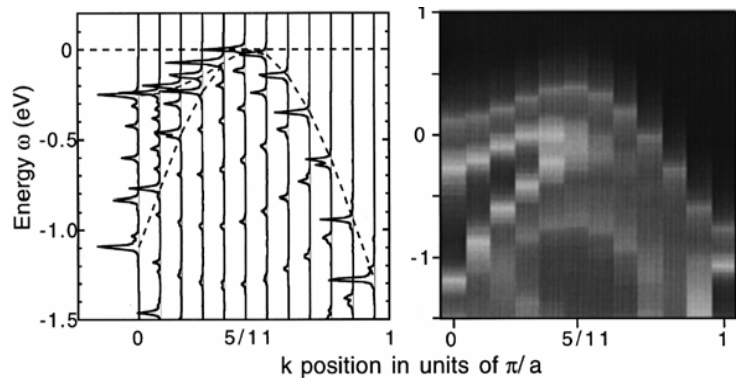
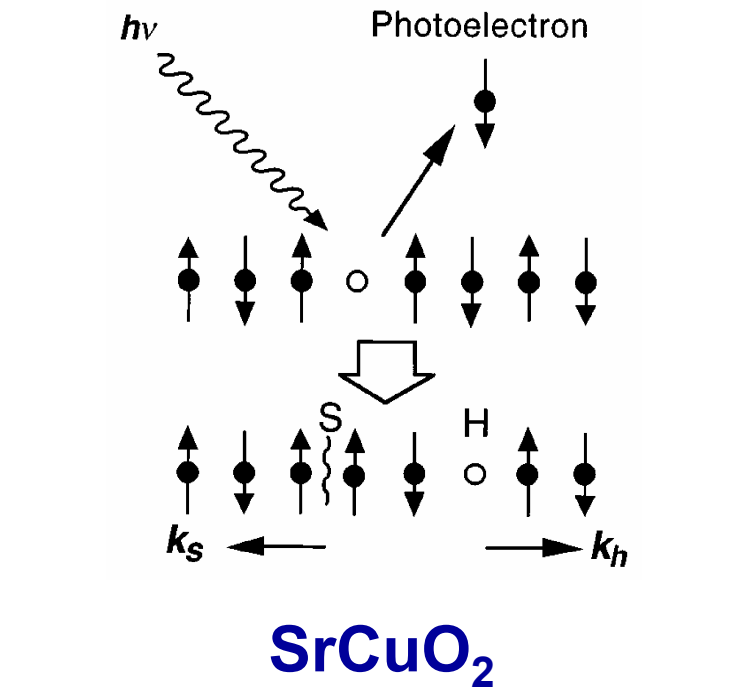
Seah, Dench *et al.*, SIA 1, 2 (1979)



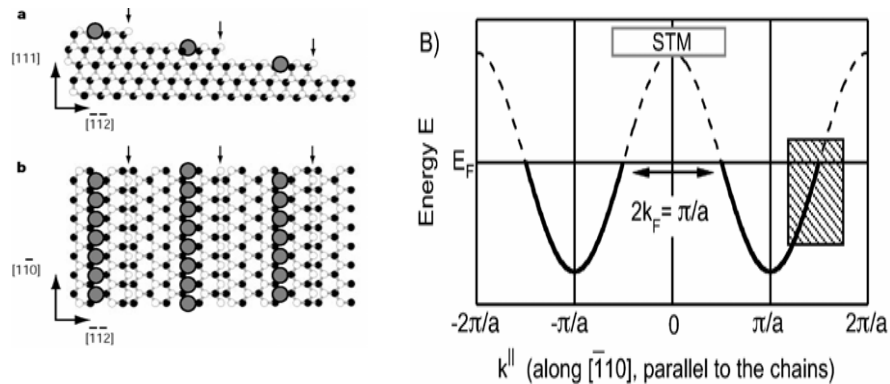
Sekiyama *et al.*, Nature **403**, 396 (2000)

ARPES on 1D Nanostructures: Spin-Charge Separation?

C. Kim *et al.*, PRL **77**, 4054 (1996)

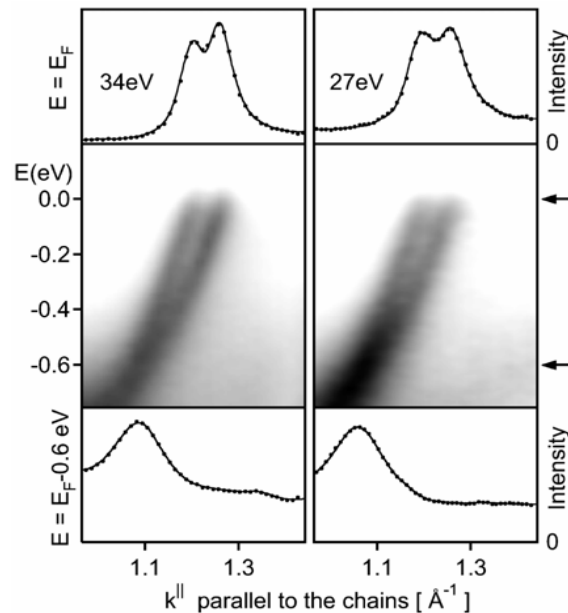


P. Segovia *et al.*, Nature **402**, 504 (1999)



Au chains on Si(557)

Losio *et al.*, PRL **86**, 4632 (2001)

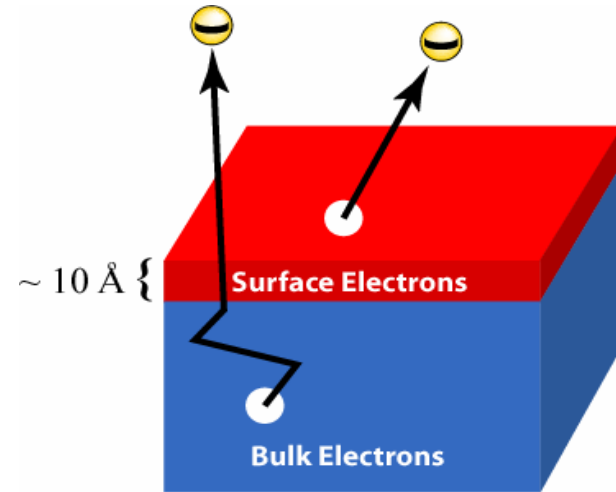


ARPES: Advantages and Limitations

Advantages

- **Direct information about the electronic states!**
- Straightforward comparison with theory - little or no modeling.
- High-resolution information about **BOTH energy and momentum**
- **Surface-sensitive probe**
- Sensitive to “**many-body**” effects
- Can be applied to small samples (100 μm x 100 μm x 10 nm)

Limitations



- **Not bulk sensitive**
- Requires clean, atomically flat surfaces in **ultra-high vacuum**
- Cannot be studied as a function of pressure or magnetic field

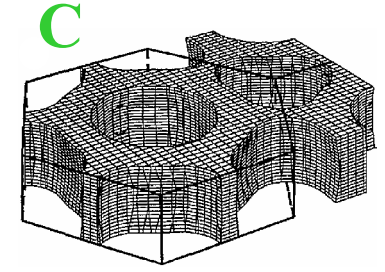
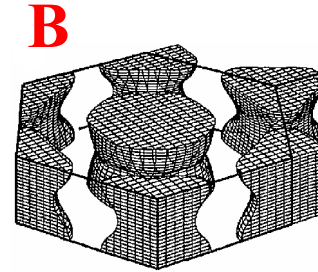
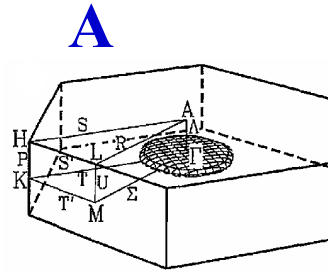
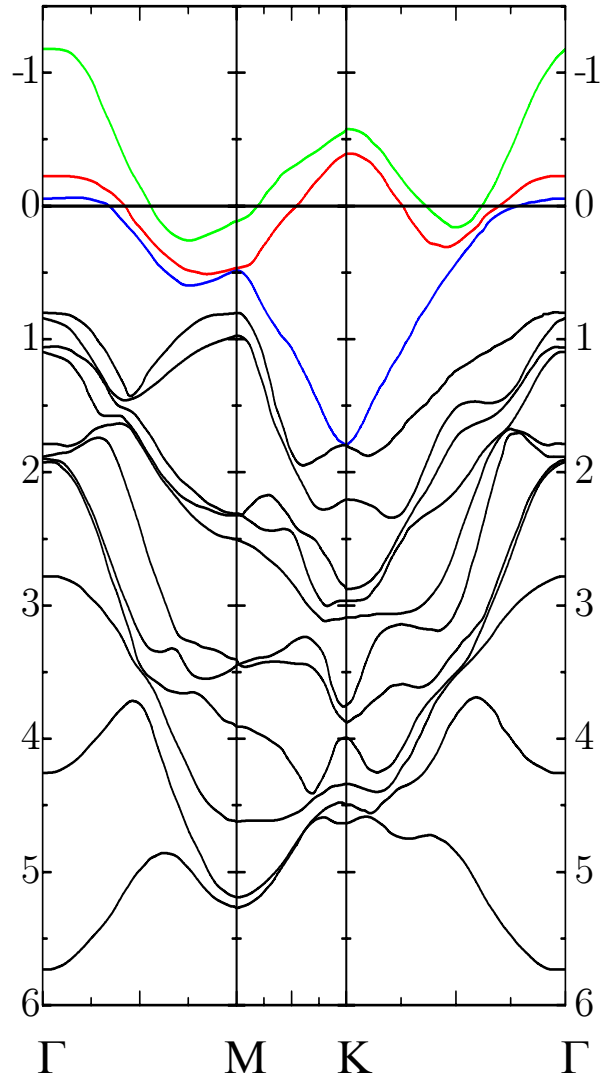
Outline: Part II

Electronic structure of complex systems

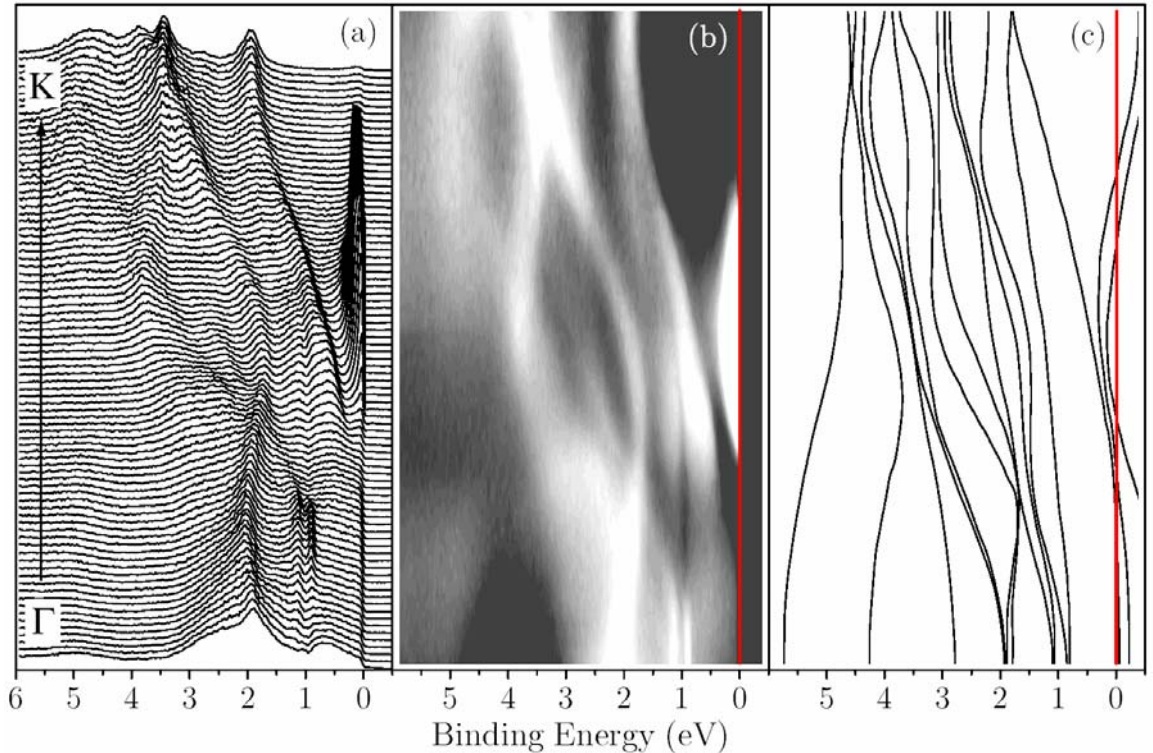
- ▶ **2H-NbSe₂ - Sr₂RuO₄**
 - Detecting bands and Fermi surface
 - Bulk & surface electronic structure
- ▶ **Nb - Bi₂Sr₂CaCu₂O_{8+δ}**
 - Superconducting gap: s-wave vs. d-wave
 - Bogoliubov quasiparticles in high-T_c cuprates
- ▶ **Be(0001) - Mo(110) - Bi₂Sr₂CaCu₂O_{8+δ}**
 - Many-body effects in the quasiparticle dispersion
- ▶ Conclusions and discussion

2H-NbSe₂: Normal State Electronic Structure

Corcoran *et al.*, JPCM **6**, 4479 (1994)

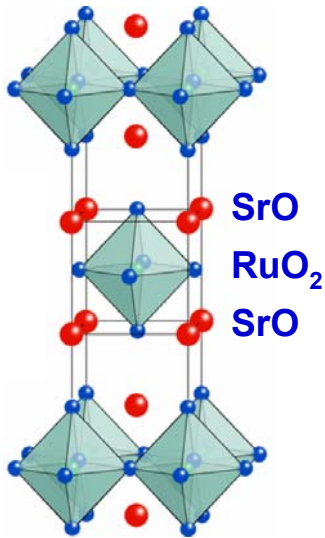


Damascelli *et al.* (2000)



Sr₂RuO₄: basic properties

2D perovskite

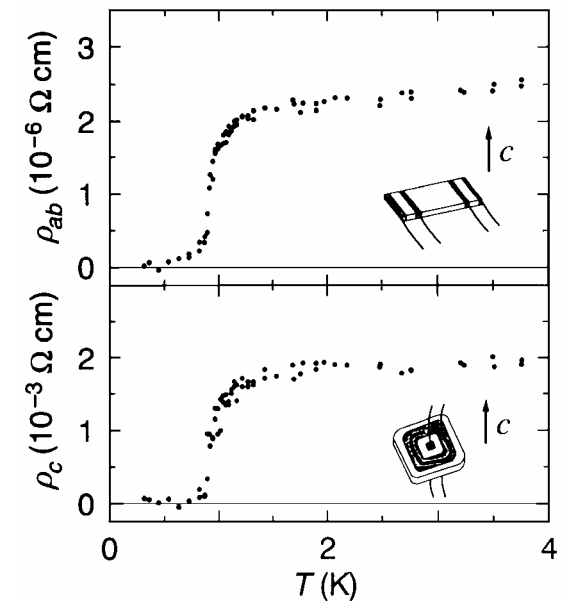


Unconventional superconductivity

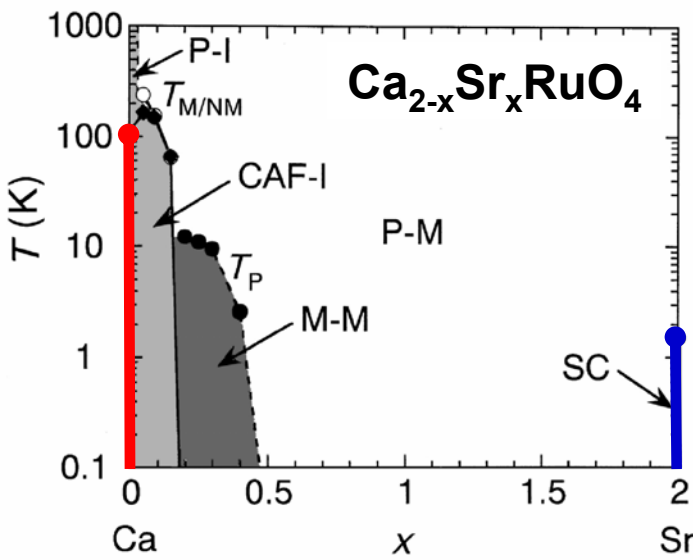
- Pairing mechanism ?
- Order parameter ?
- FM-AF fluctuations ?

Rice & Sigrist, JPCM 7, L643 (1995)

Maeno *et al.*, Nature 372, 532 (1994)



Nakatsuji & Maeno, PRL 84, 2666 (2000)



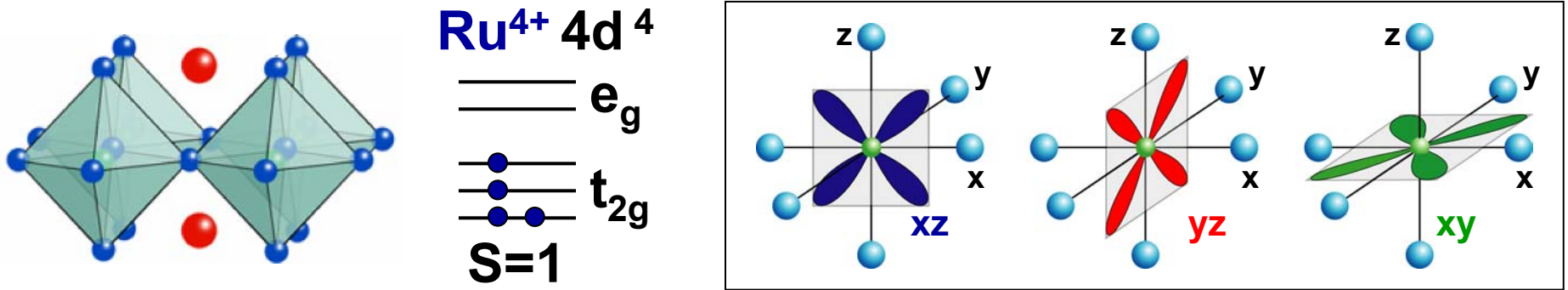
Lattice-magnetism interplay Orbital degrees of freedom

Sr₂RuO₄ : 2D Fermi Liquid ($\rho_c/\rho_{ab}=850$)

Ca₂RuO₄ : insulating **Anti-Ferromagnet**

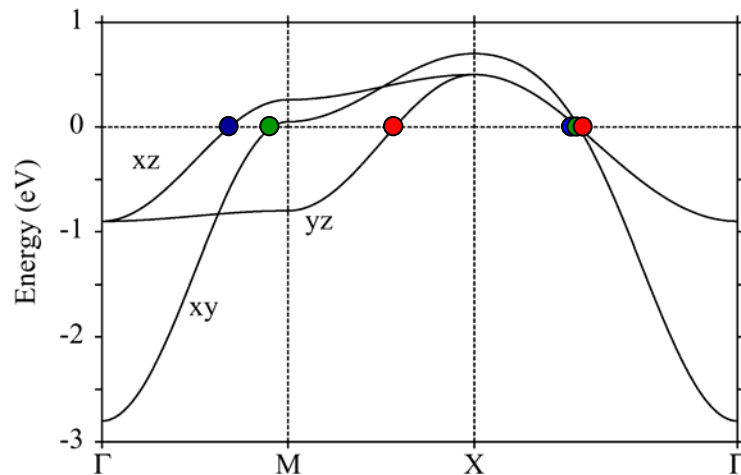
SrRuO₃ : metallic **Ferromagnet**

Low-Energy Electronic structure of Sr_2RuO_4

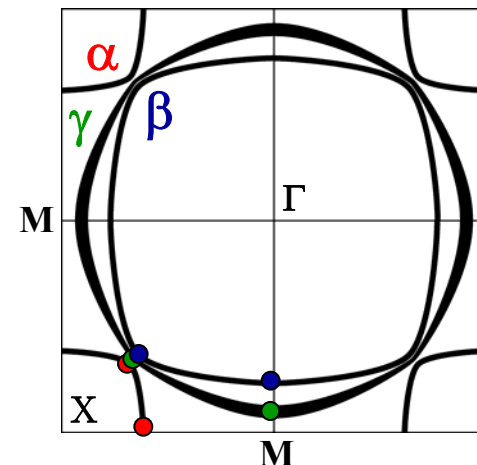


► Band structure calculation: **3** t_{2g} bands crossing E_F

→ 3 sheets of FS $\left\{ \begin{array}{l} \alpha \text{ (hole-like)} \\ \beta \text{ and } \gamma \text{ (electron-like)} \end{array} \right.$



A. Liebsch *et al*, PRL **84**, 1591 (2000)

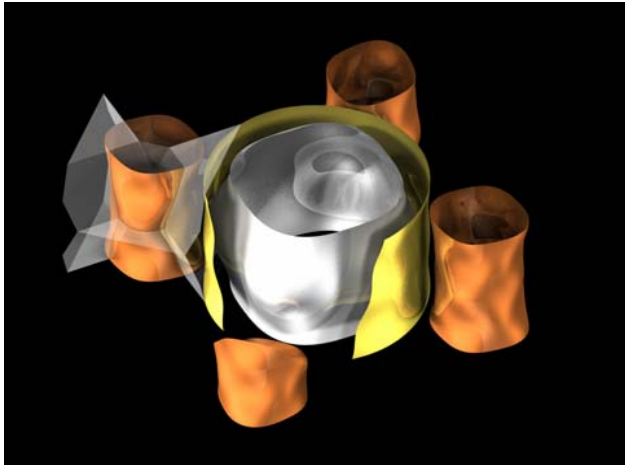


I.I. Mazin *et al*, PRL **79**, 733 (1997)

Fermi Surface Topology of Sr_2RuO_4

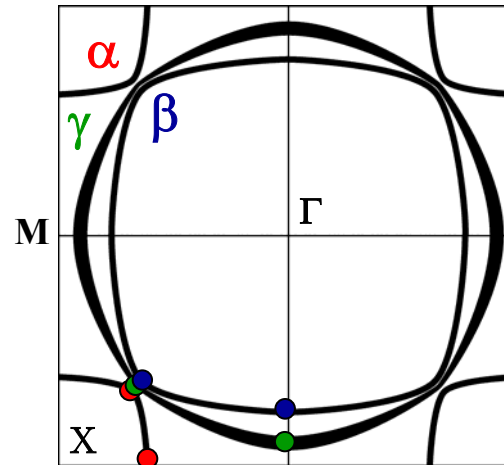
► Early ARPES results gave a different topology

de Haas-van Alphen



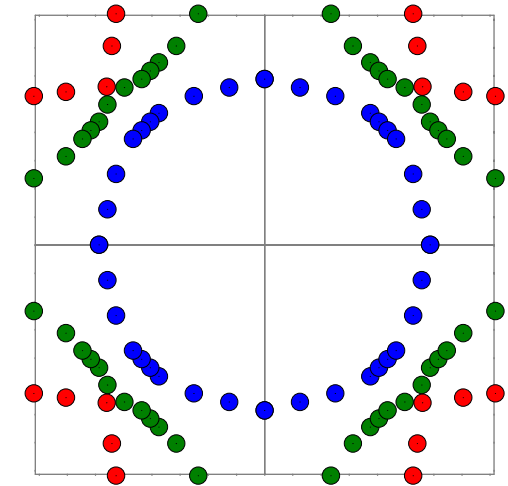
A.P. Mackenzie *et al.*, PRL **76**, 3786 (1996)
C. Bergemann *et al.*, PRL **84**, 2662 (2000)

LDA



I.I. Mazin *et al.*, PRL **79**, 733 (1997)

ARPES



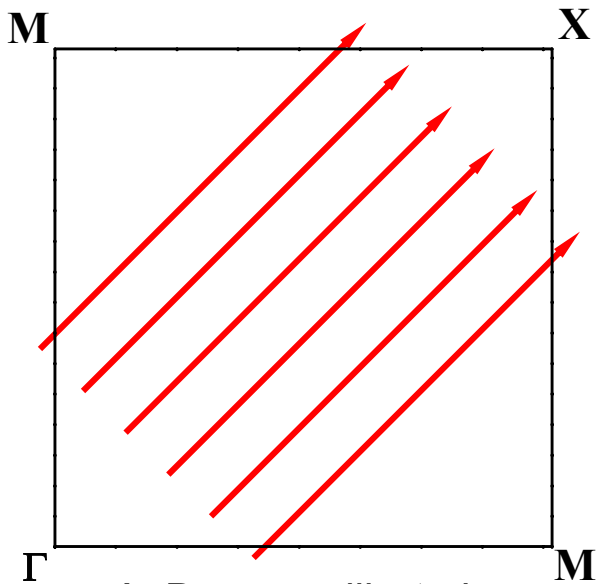
T.Yokoya *et al.*, PRB **54**, 13311 (1996)
D.H. Lu *et al.*, PRL **76**, 4845 (1996)

ARPES: additional information

dHvA: limited set of systems

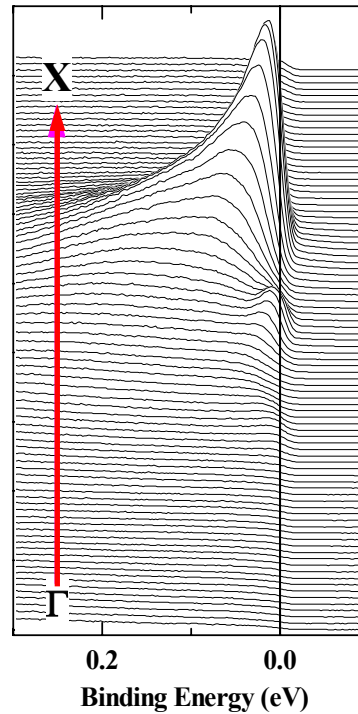
Fermi Surface Topology of Sr_2RuO_4

ARPES : present day

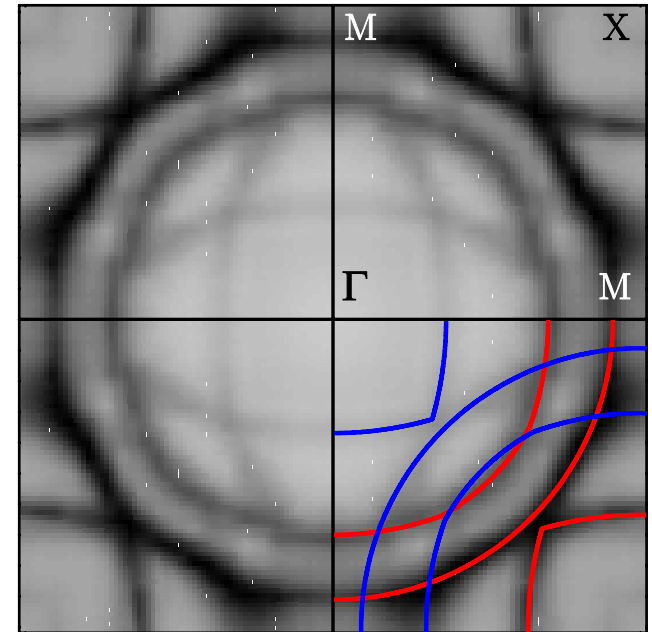


A. Damascelli *et al.*,
PRL **85**, 5194 (2000)

ARPES Spectra

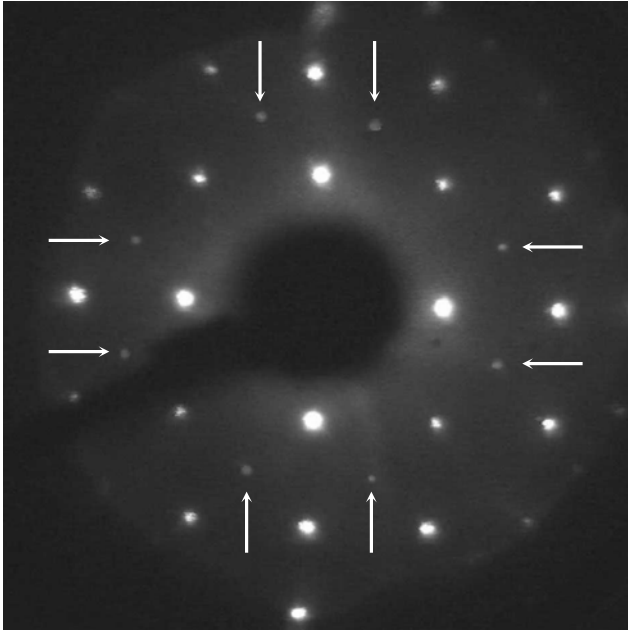


Intensity at E_F



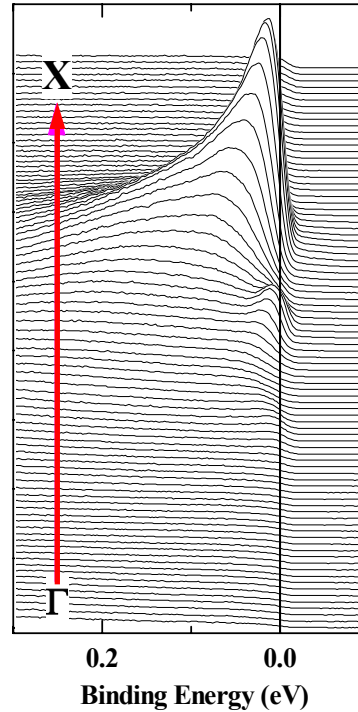
Fermi Surface Topology of Sr_2RuO_4

LEED Pattern

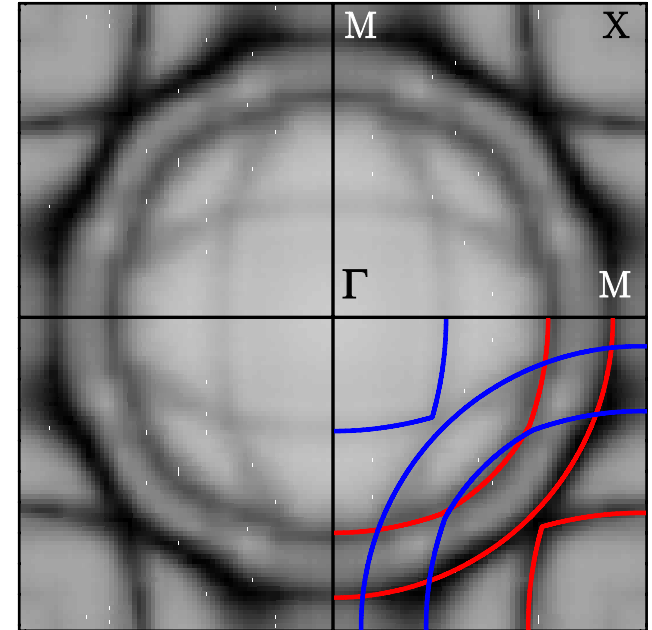


A. Damascelli *et al.*,
PRL **85**, 5194 (2000)

ARPES Spectra



Intensity at E_F



Surface instability



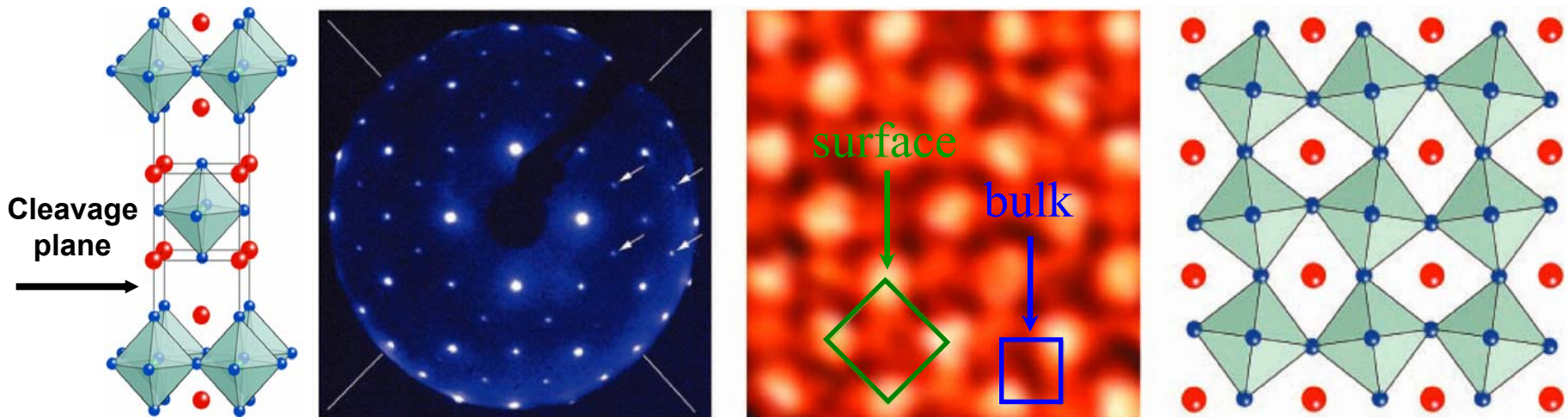
Band folding

Surface reconstruction of cleaved Sr_2RuO_4

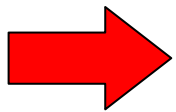
LEED

STM

a-b plane



R. Matzdorf *et al.*, Science **289**, 746 (2000)

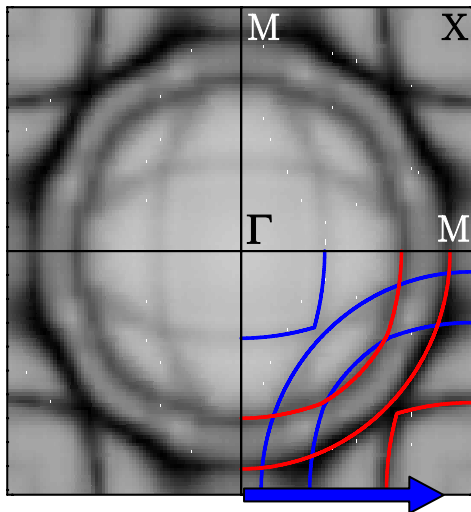


**Rotation of the RuO_6 octahedra
around the c axis (9°)**

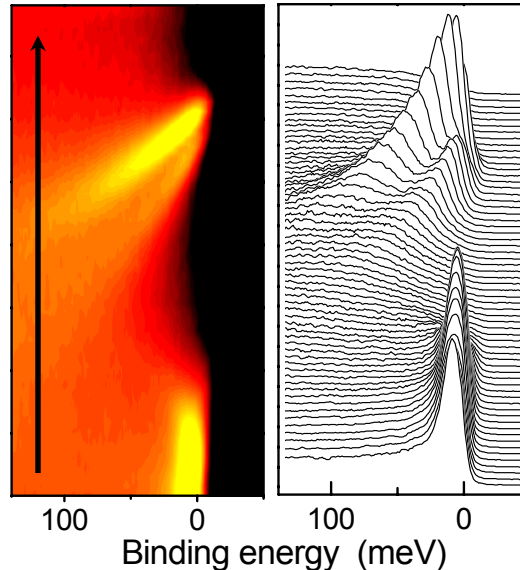
Surface electronic structure of Sr_2RuO_4

On samples cleaved at **180 K**
the **surface**-related features are
suppressed

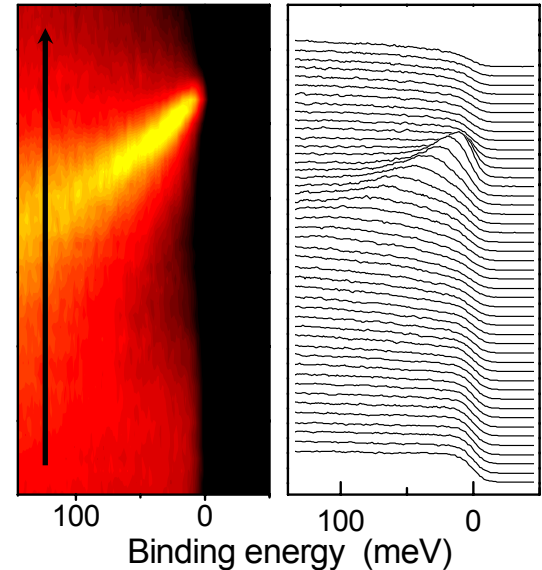
E_F mapping
 ± 10 meV



Cold cleave
 $T=10$ K



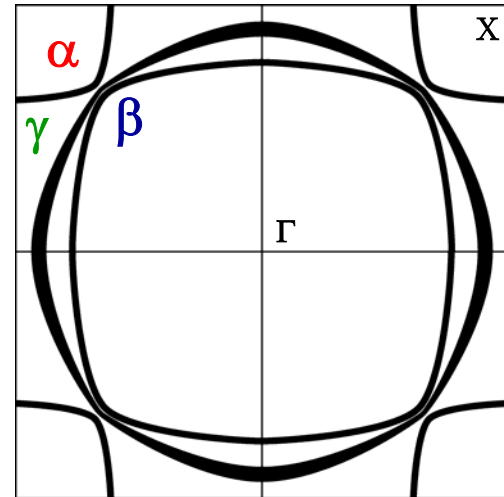
Hot cleave
 $T=180$ K



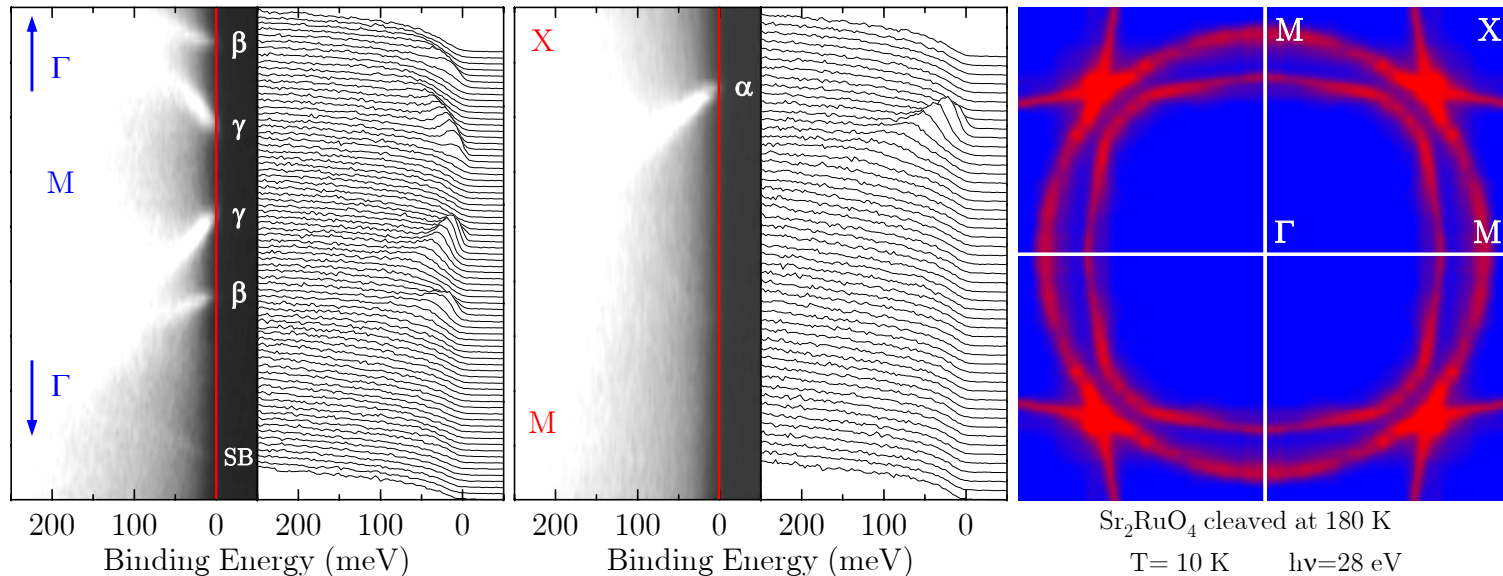
Bulk electronic structure of Sr_2RuO_4

What do we learn about the **bulk electronic structure**?

- Fermi Surface
- Fermi velocity
- Effective mass

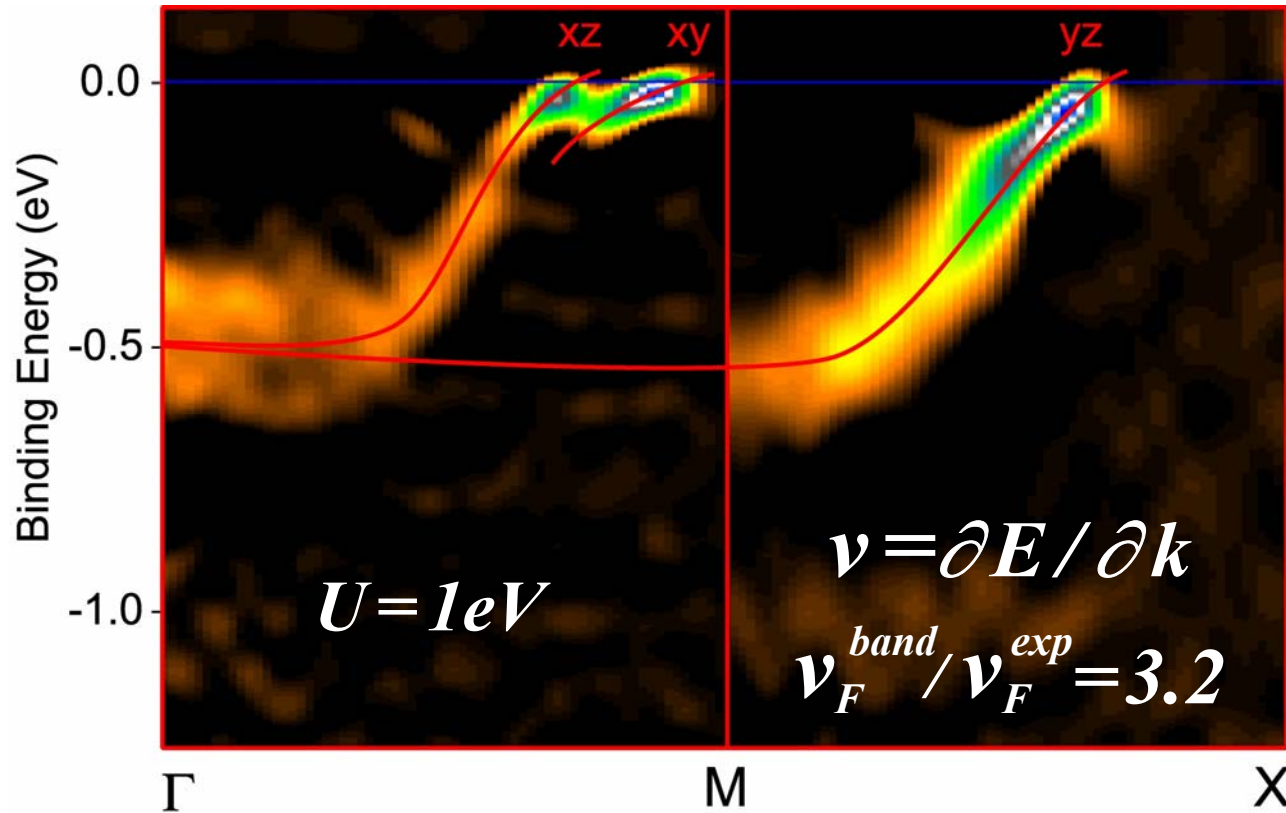


I.I. Mazin *et al.*, PRL **79**, 733 (1997)



A. Damascelli *et al.*, PRL **85**, 5194 (2000)

Dispersion of the bulk electronic bands

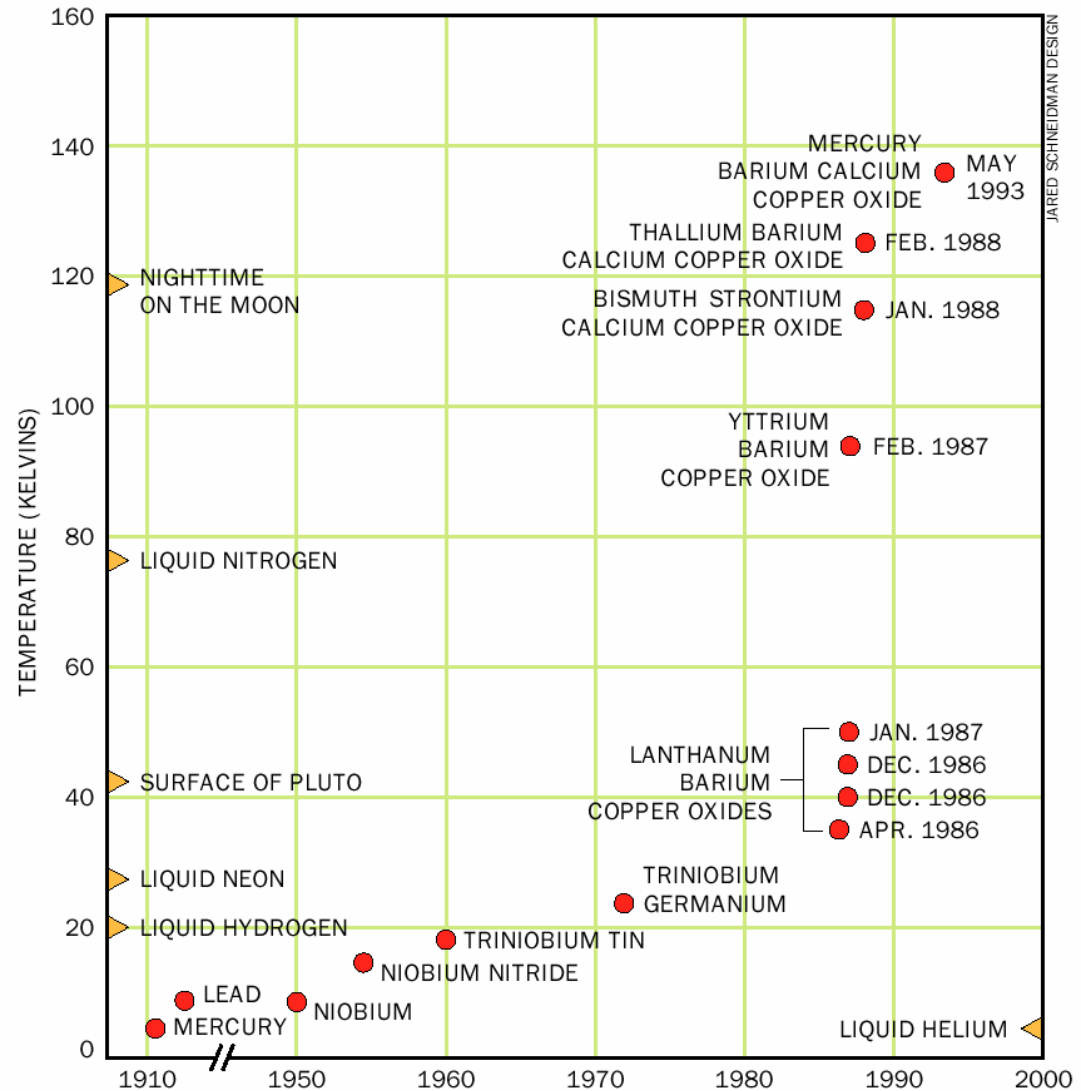
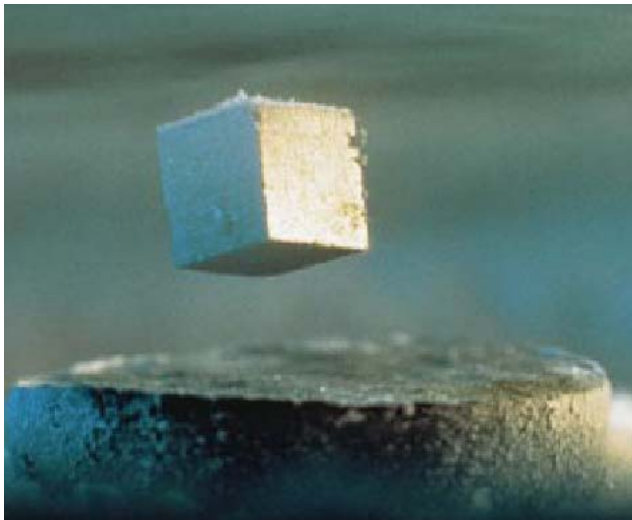
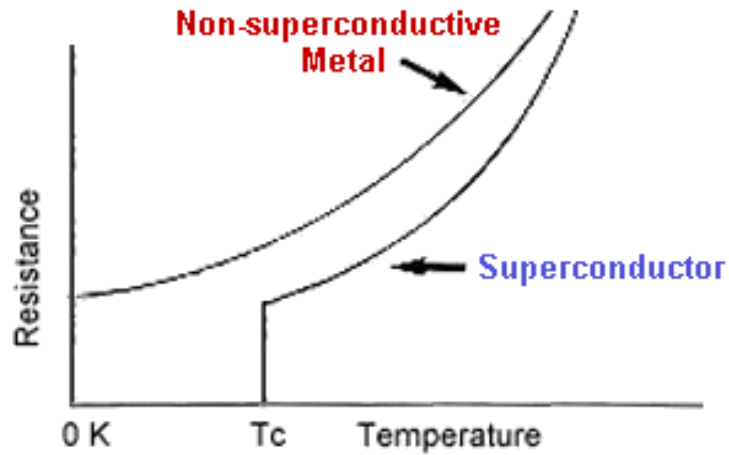


Experiment compares well with **LDA+U** calculations

A. Damascelli *et al.*, PRL **85**, 5194 (2000)

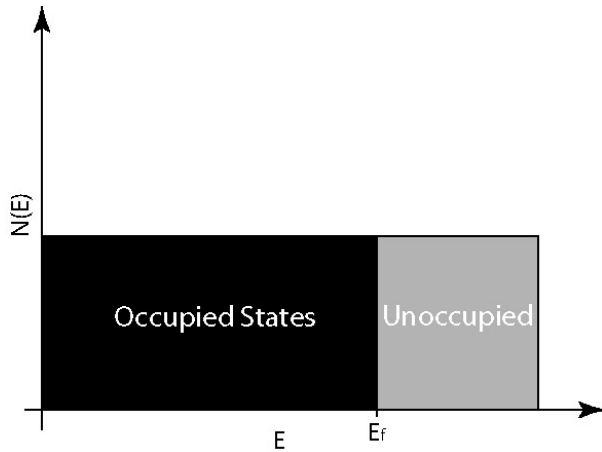
A. Liebsch & A. Lichtenstein, PRL **84**, 1591 (2000)

Superconductivity

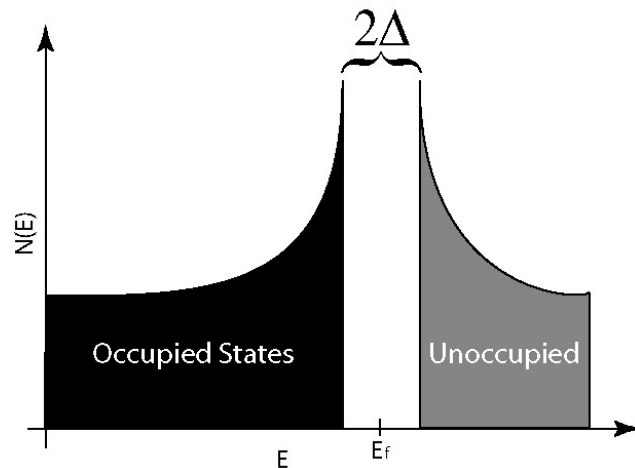


“Classic Low-temperature” Superconductors

Metallic Density of States



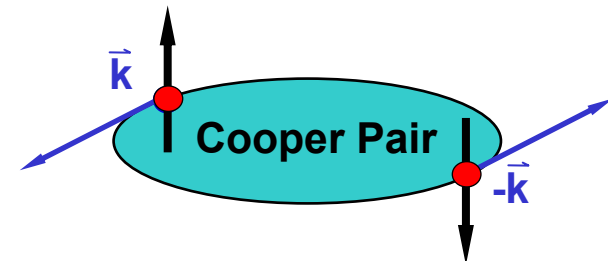
Superconducting Density of States



Superconductivity can only be seen on low energy scales and needs **high resolution!**

Superconductivity

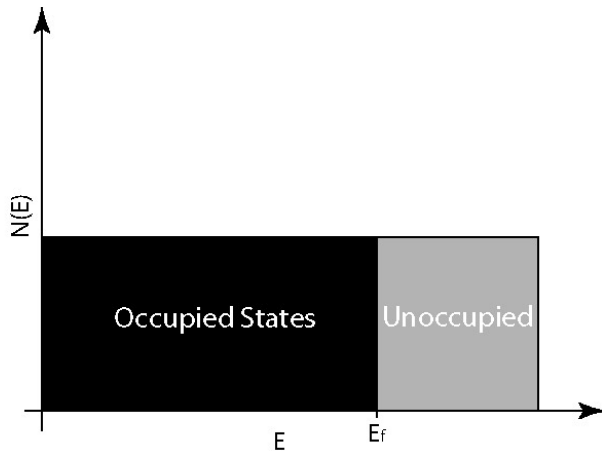
2-electron
bound state



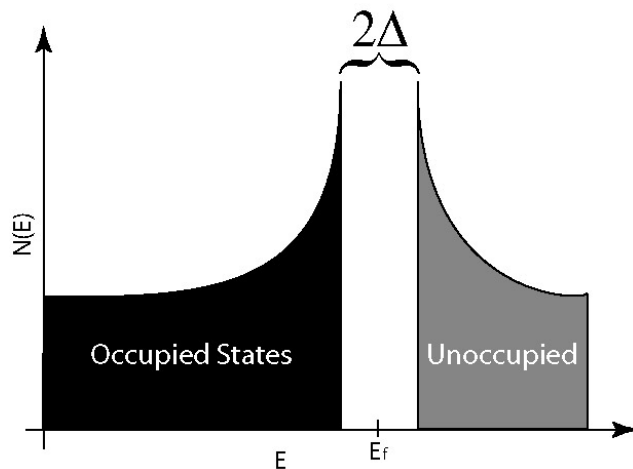
spin-singlet pairing

“Classic Low-temperature” Superconductors

Metallic Density of States

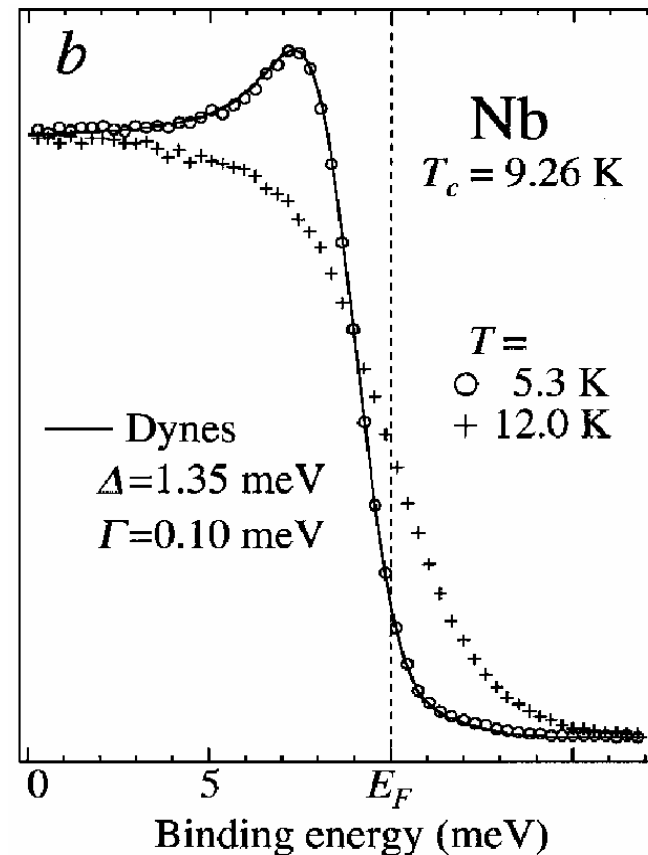


Superconducting Density of States



Superconductivity can only be seen on low energy scales and needs **high resolution!**

A. Chainani et al., PRL **85** (2000)



High-Temperature Superconductors: Bi2212

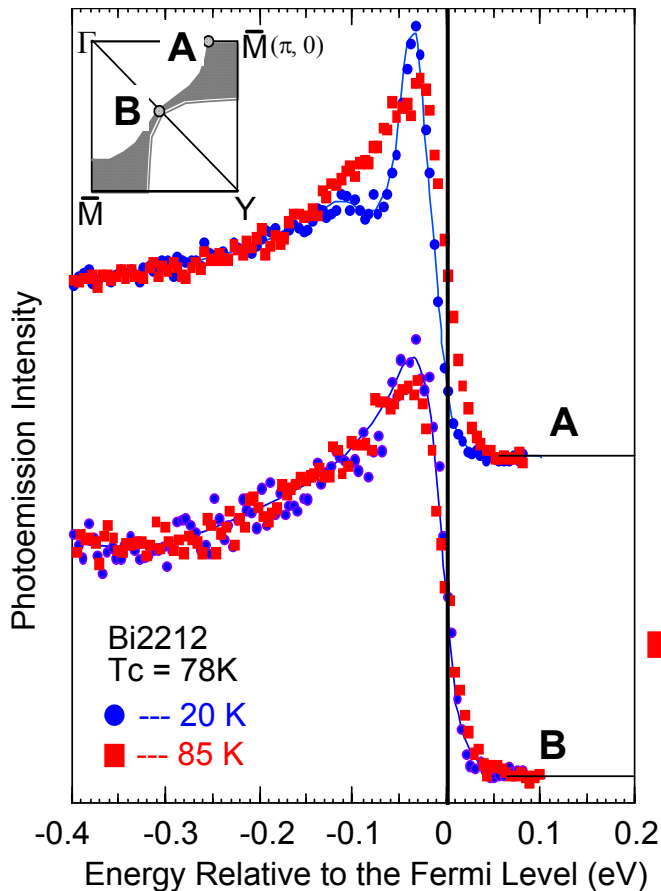
VOLUME 70, NUMBER 10

PHYSICAL REVIEW LETTERS

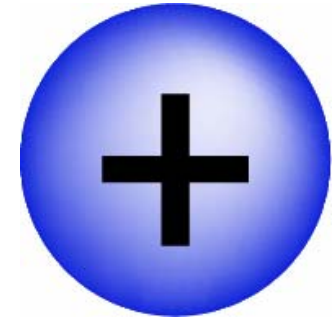
8 MARCH 1993

Anomalous Large Gap Anisotropy in the a - b Plane of $\text{Bi}_2\text{Sr}_2\text{CaCu}_2\text{O}_{8+\delta}$

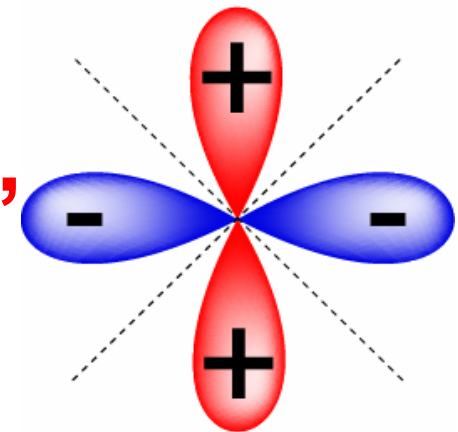
Z.-X. Shen,^{(1),(2)} D. S. Dessau,^{(1),(2)} B. O. Wells,^{(1),(2),(a)} D. M. King,⁽²⁾ W. E. Spicer,⁽²⁾ A. J. Arko,⁽³⁾
D. Marshall,⁽²⁾ L. W. Lombardo,⁽¹⁾ A. Kapitulnik,⁽¹⁾ P. Dickinson,⁽¹⁾ S. Doniach,⁽¹⁾ J. DiCarlo,^{(1),(2)}
A. G. Loeser,^{(1),(2)} and C. H. Park^{(1),(2)}



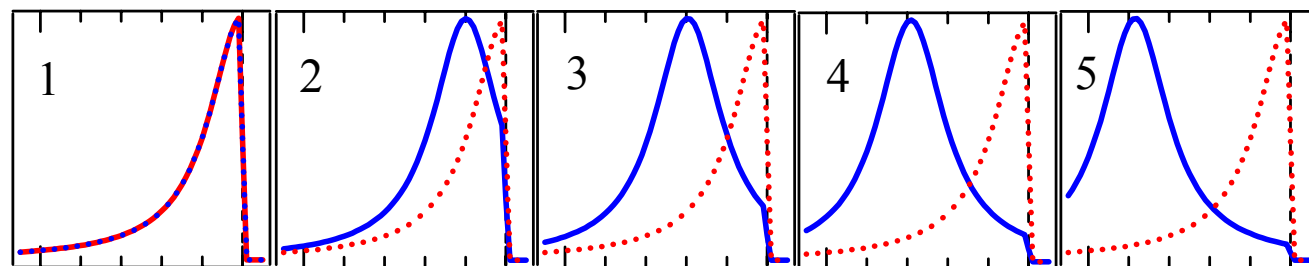
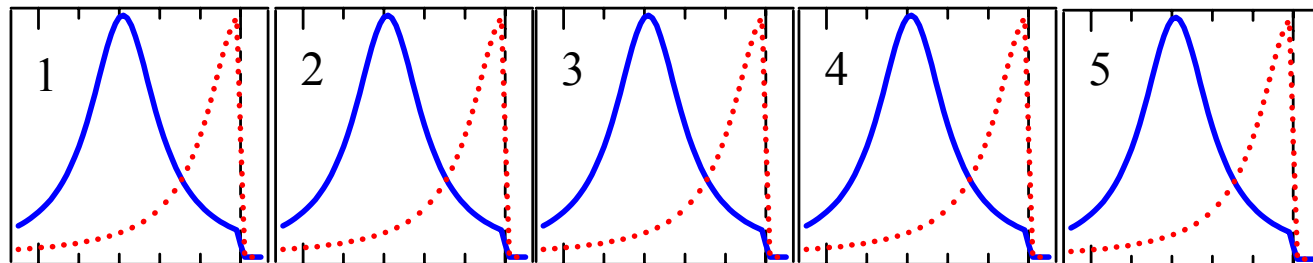
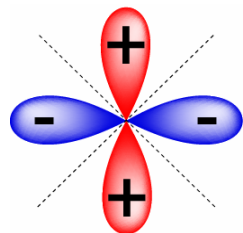
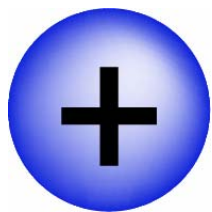
“s-wave”



“d-wave”

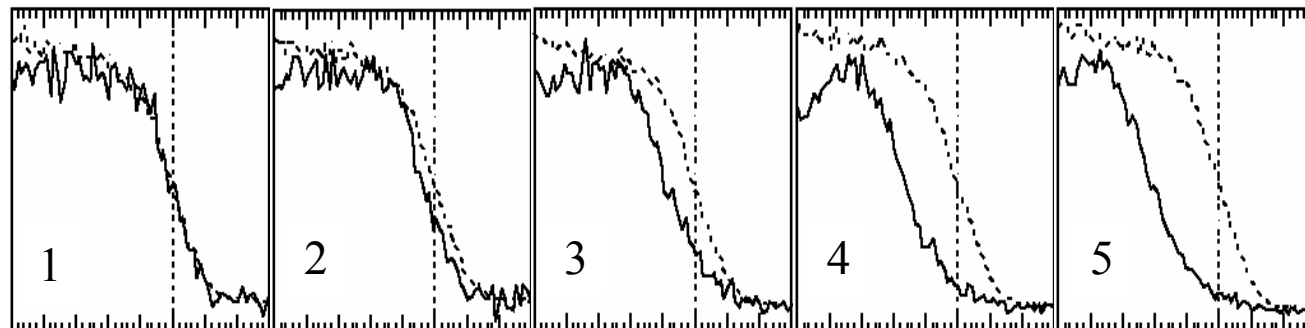
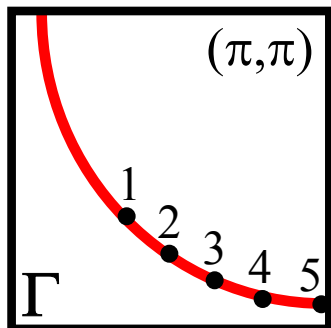


High- T_c Superconductors: s-wave vs. d-wave gap



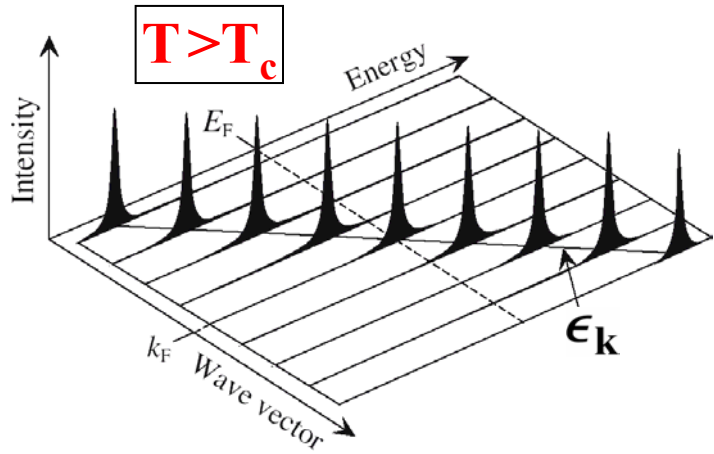
Binding Energy (meV)

Ding et al. PRB **54**, R9678 (1996)

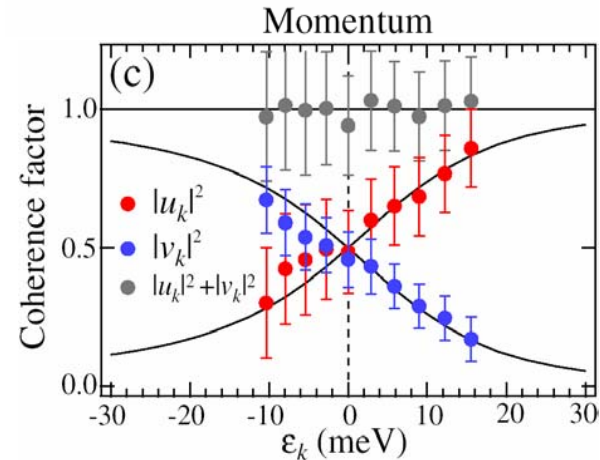
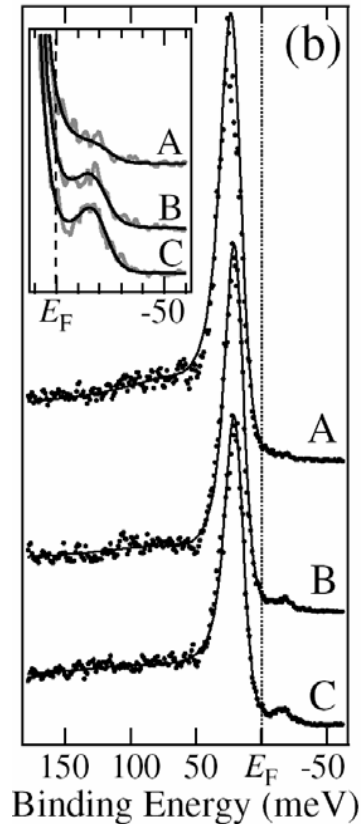
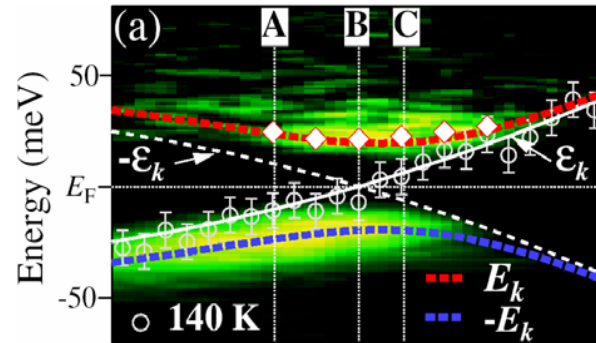


Binding Energy (meV)

High- T_c Superconductors: Bogoliubov QP in Bi2223

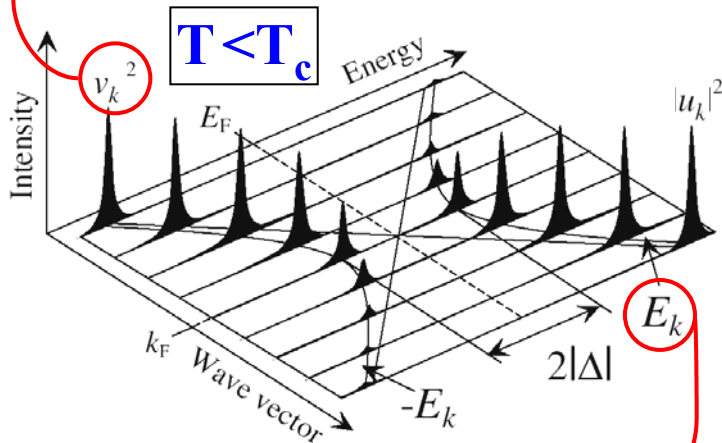


$$A_{\text{BCS}}(k, \omega) = \frac{1}{\pi} \left\{ \frac{|u_{\mathbf{k}}|^2 \Gamma}{(\omega - E_{\mathbf{k}})^2 + \Gamma^2} + \frac{|v_{\mathbf{k}}|^2 \Gamma}{(\omega + E_{\mathbf{k}})^2 + \Gamma^2} \right\}$$



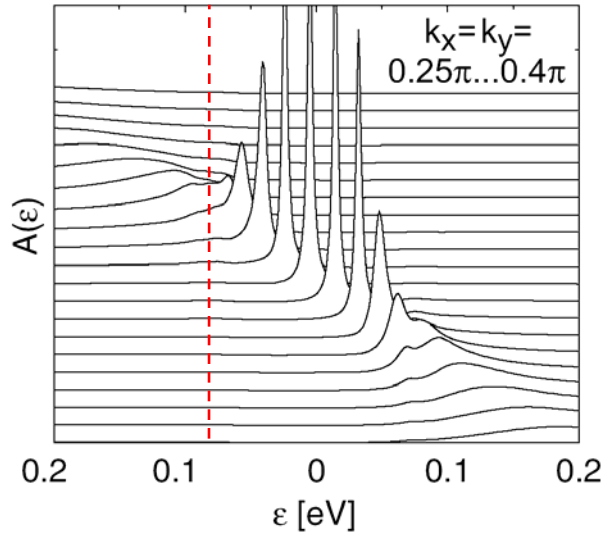
Matsui *et al.*, PRL **90**, 217002 (2003)

$$v_{\mathbf{k}}^2 = 1 - u_{\mathbf{k}}^2 = \frac{1}{2} \left(1 - \epsilon_{\mathbf{k}} / E_{\mathbf{k}} \right)$$



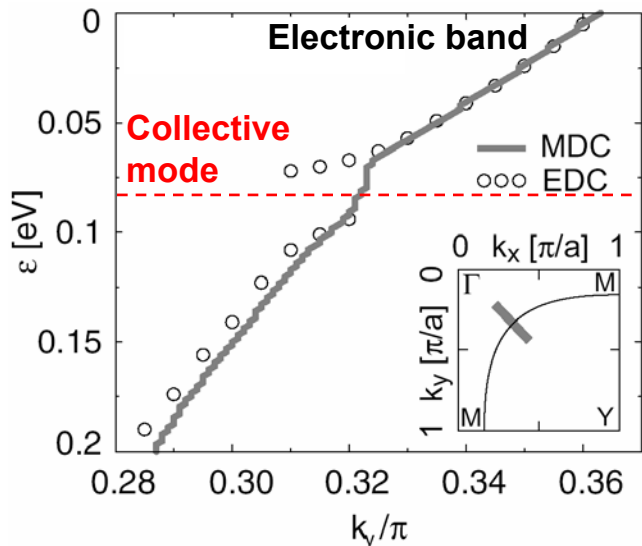
$$E_{\mathbf{k}} = \sqrt{\epsilon_{\mathbf{k}}^2 + |\Delta(\mathbf{k})|^2}$$

Many-Body effects: Electron-Phonon Coupling

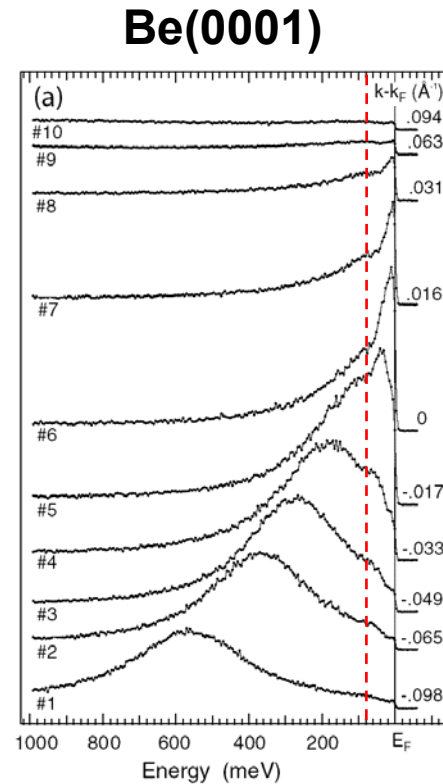


Single-particle spectral function

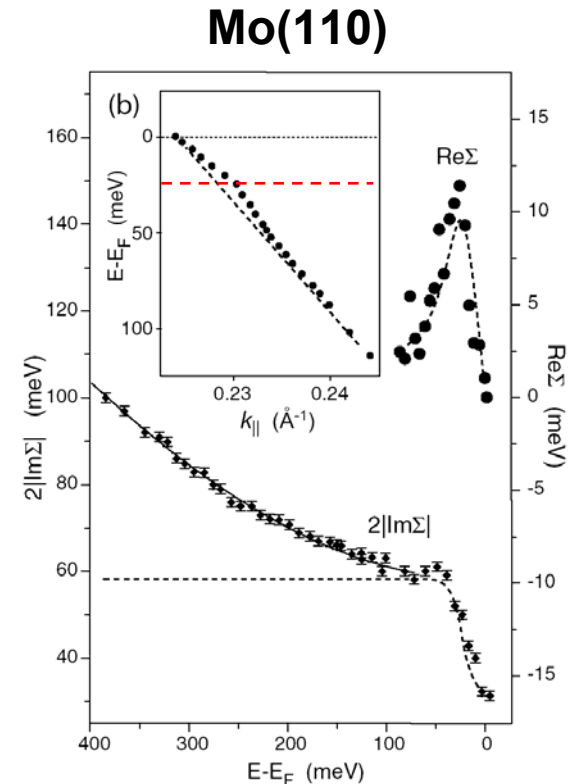
$$A(\mathbf{k}, \omega) = -\frac{1}{\pi} \frac{\Sigma''(\mathbf{k}, \omega)}{[\omega - \epsilon_{\mathbf{k}} - \Sigma'(\mathbf{k}, \omega)]^2 + [\Sigma''(\mathbf{k}, \omega)]^2}$$



Eschrig, Norman, PRB **67**, 144503 (2003)



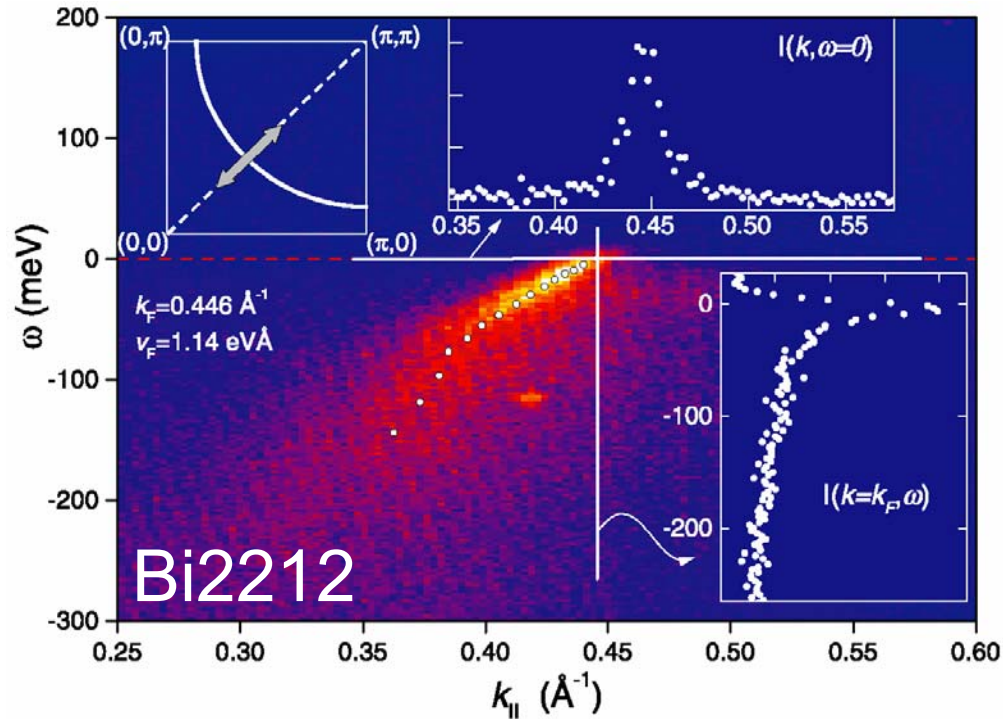
Hengsberger *et al.*, PRL **83**, 592 (1999)



Valla *et al.*, PRL **83**, 2085 (1999)

Many-Body effects in the High- T_c Cuprates

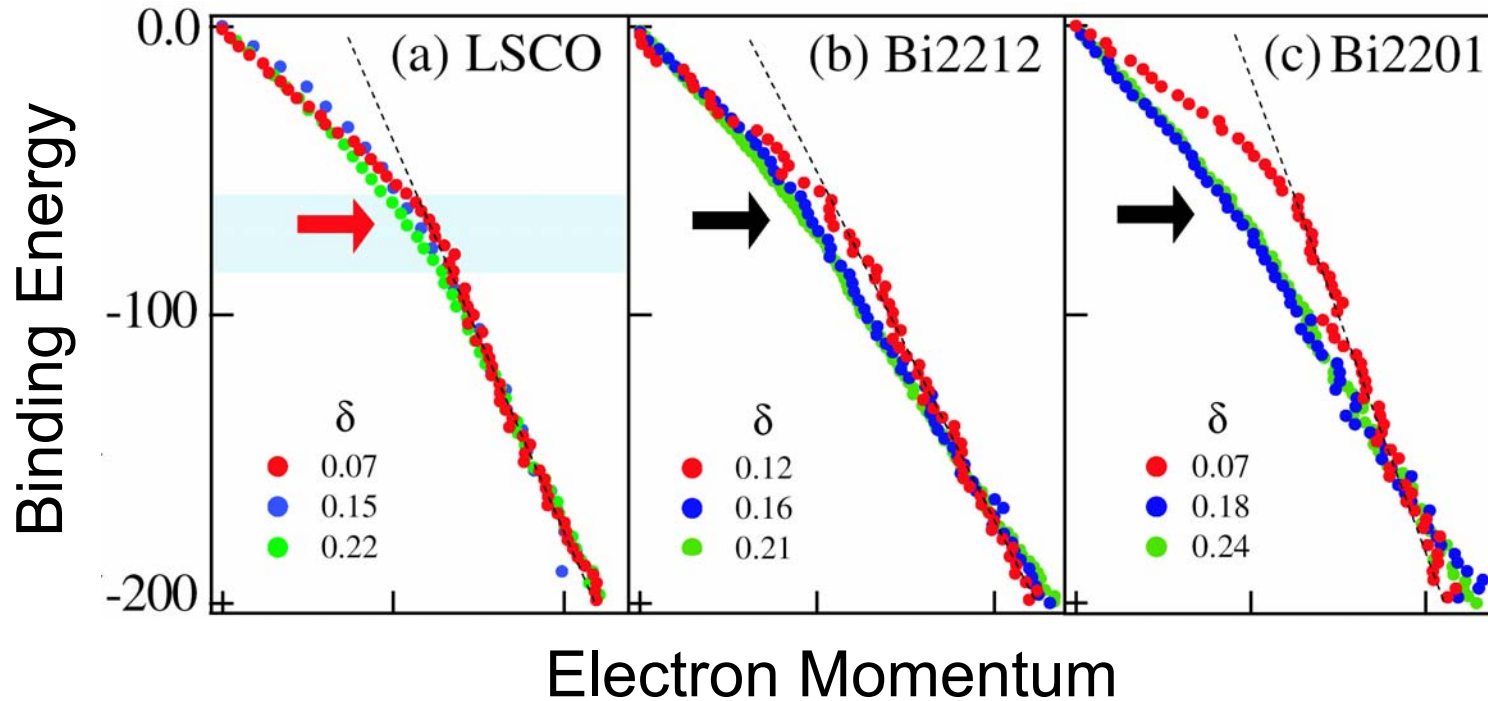
Valla *et al.*, Science **285**, 2110 (1999)



Mechanism for High- T_c { **Magnetic fluctuations ?**
Electron-phonon coupling ?

Many Body effects in the Quasiparticle Dispersion

Lanzara *et al.*, Nature **412**, 510

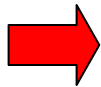


Mechanism for High- T_c { **Magnetic fluctuations ?**
Electron-phonon coupling ?

Conclusions

ARPES results from complex systems

- **Bands** and **FS** in unprecedented detail
- Fermi **velocity** and **effective mass**
- Superconducting (d-wave) **gap**
- **Many-body effects** in the QP dispersion
- Nanostructured materials (**surface FM**)



ARPES is a **powerful tool** for the study of the electronic structure of complex systems

For a review article see:

A. Damascelli, Z. Hussain, and Z.-X Shen, Rev. Mod. Phys. **75**, 473 (2003)

For additional material see:

<http://www.physics.ubc.ca/~damascel/>

