

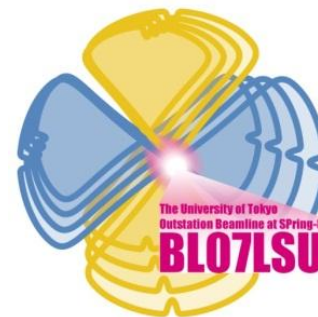
Carrier dynamics at semiconductor surfaces

Marie D'angelo

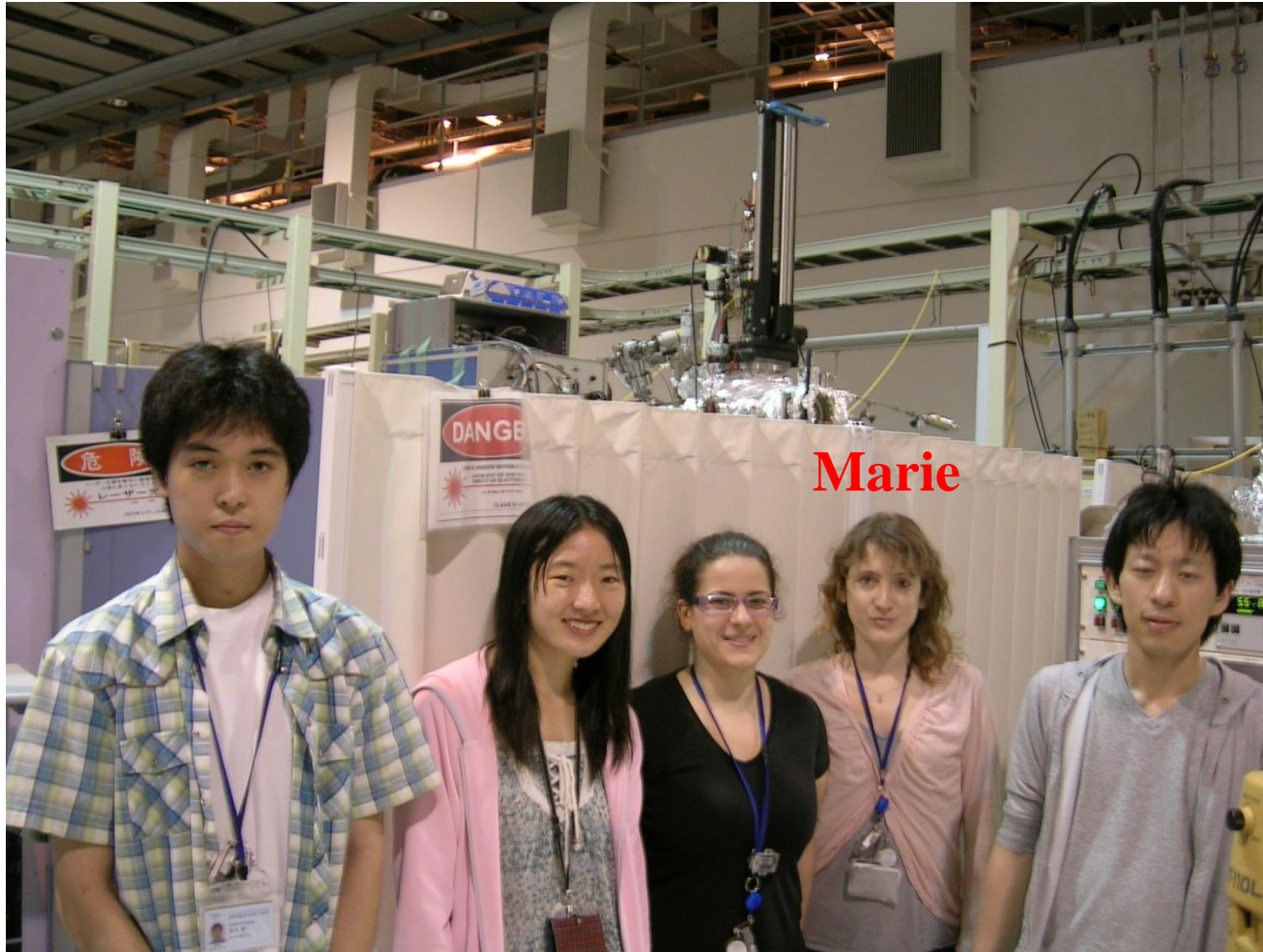
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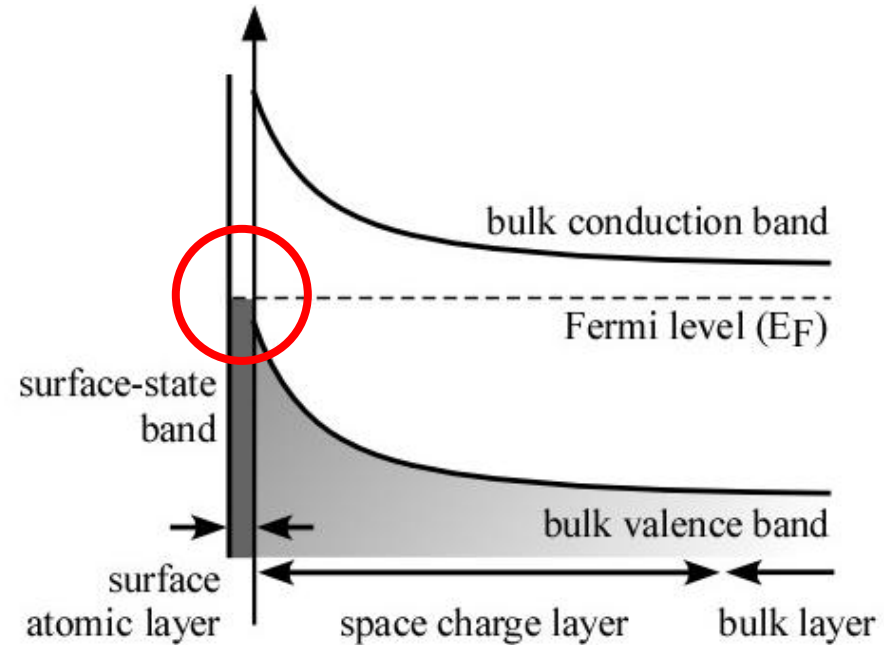
■ Low-dimensional physics at a surface

governed by **surface electrons**
at the **Fermi level**

The regulation



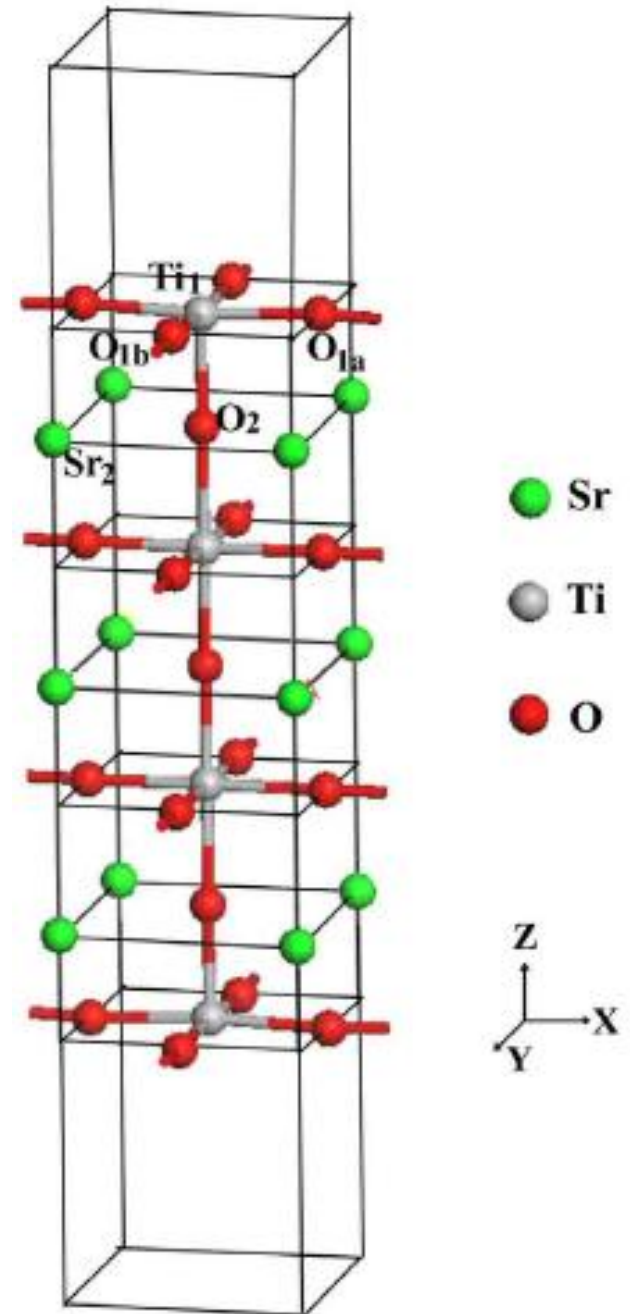
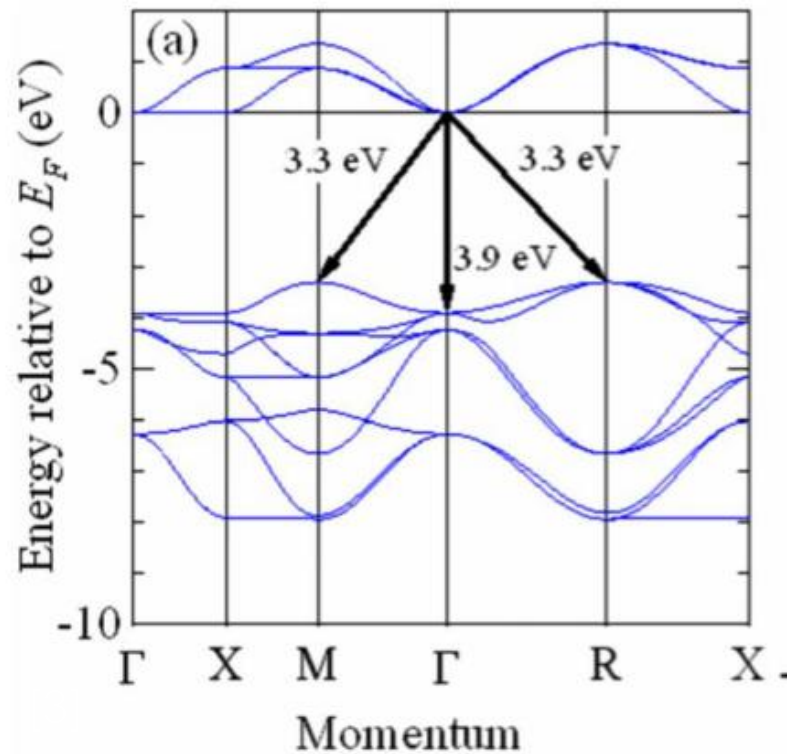
Induction of intriguing electronic states at a surface



- Formation of long-ranged ordered phase (surface superstructure)
- Insulator-to-metal surface phase transition
- Etc

SrTiO₃ (001)

a playground for studying correlated electrons
a model for catalytic reaction
a candidate of future device (oxide electronics)



H/SrTiO₃(001)

Hydrogen-induced metallicity of SrTiO₃ (001) surfaces

Feng Lin *et al*, Phys. Rev. B **79**, 035311 (2009).

- SrTiO₃: d⁰ configuration.

Electron filling into the Ti 3d band

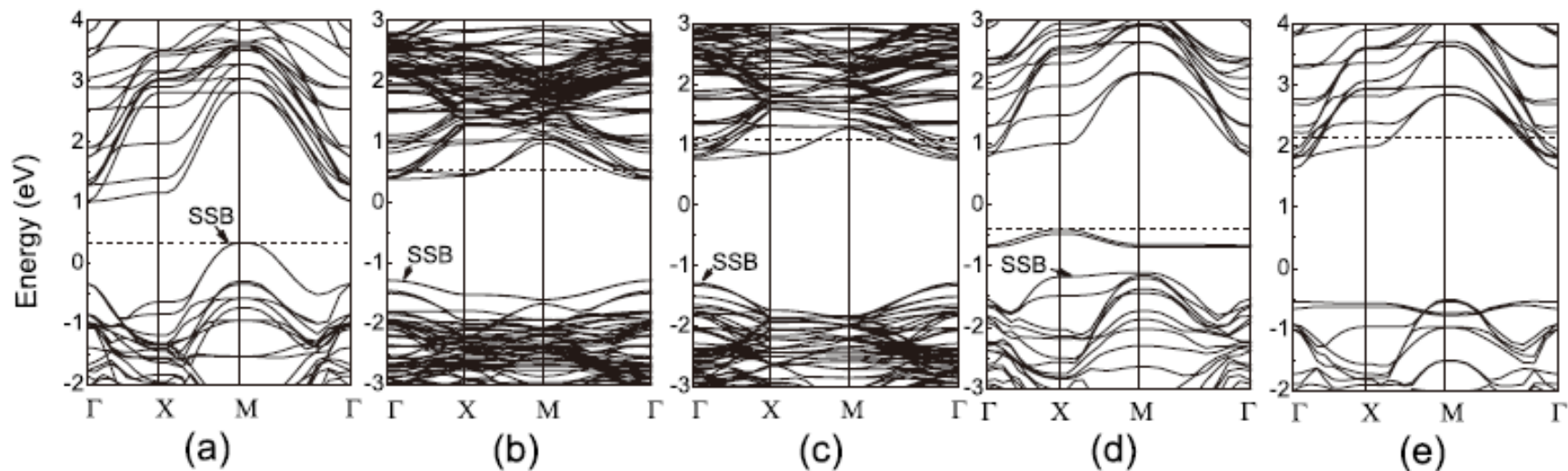
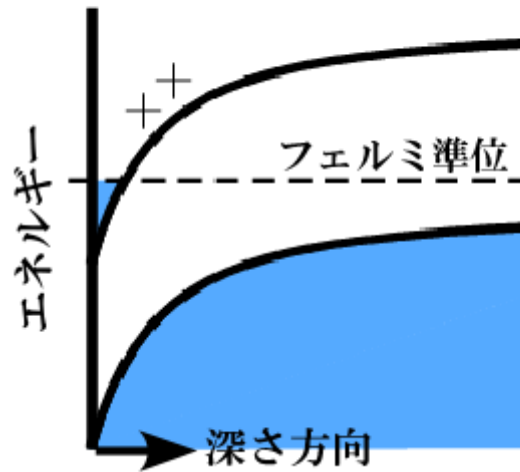


FIG. 4. Band structures of (a) clean and (b)–(e) hydrogen-adsorbed TiO₂-terminated SrTiO₃ surface. Herein, (b) and (c), respectively, correspond to the adsorption coverages of 1/12 and 1/6 ML, which are calculated using 2 × 2 supercell, while, (a), (d), and (e), respectively, correspond to the clean surface, Ti-2/3 ML adsorption, and 1 ML adsorption, which are calculated using 1 × 1 supercell. The Fermi level is denoted by the dashed line. Surface-state bands (SSBs) are denoted by arrows.

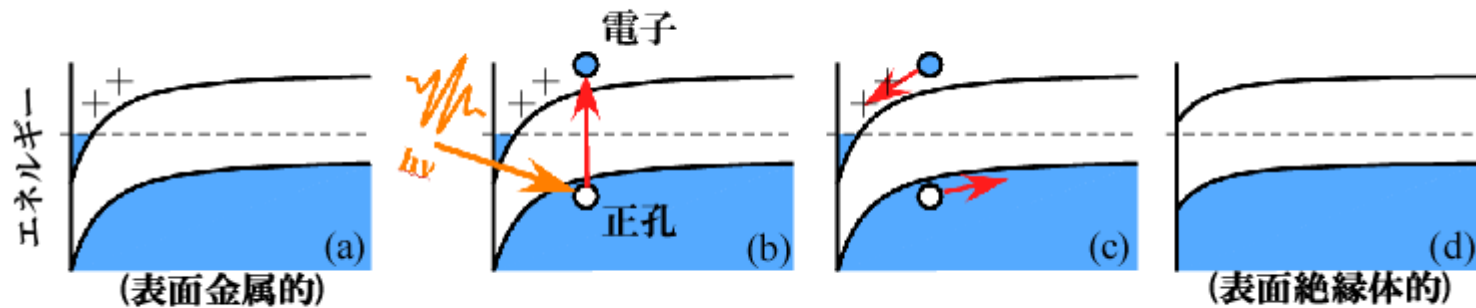
H/SrTiO₃(001)

Motivation

Surface metallization?

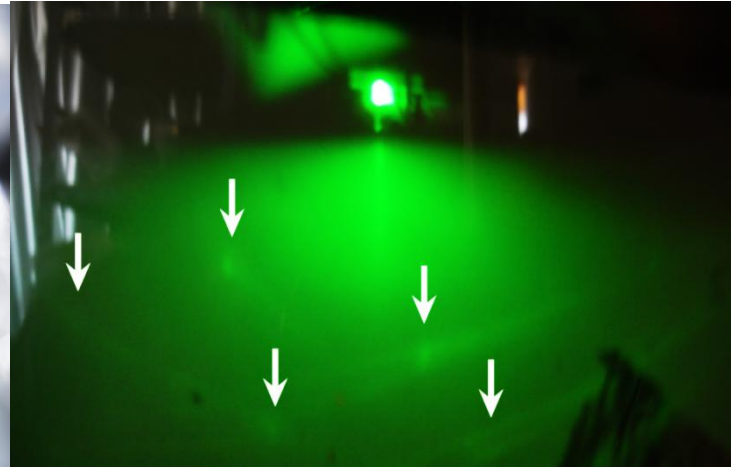
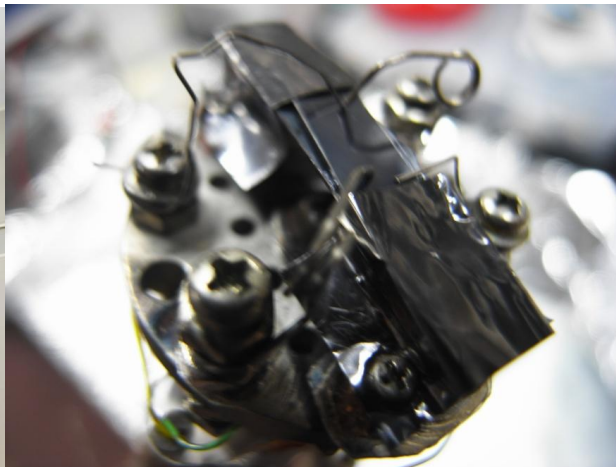


Optically-induced surface metal-to-insulator transition?



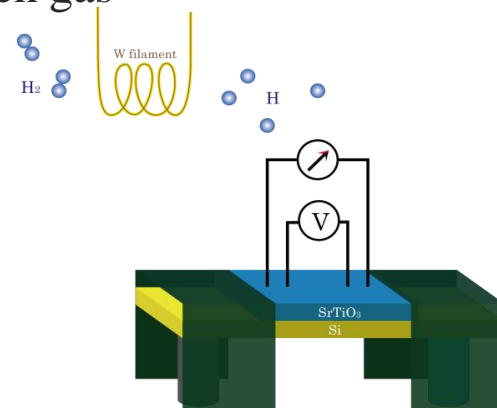
H/SrTiO₃(001)

Surface transport measurements

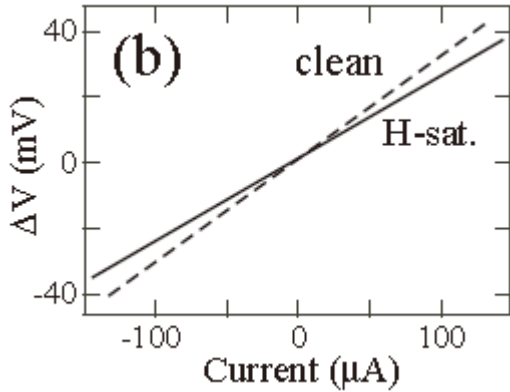


SrTiO₃(001)

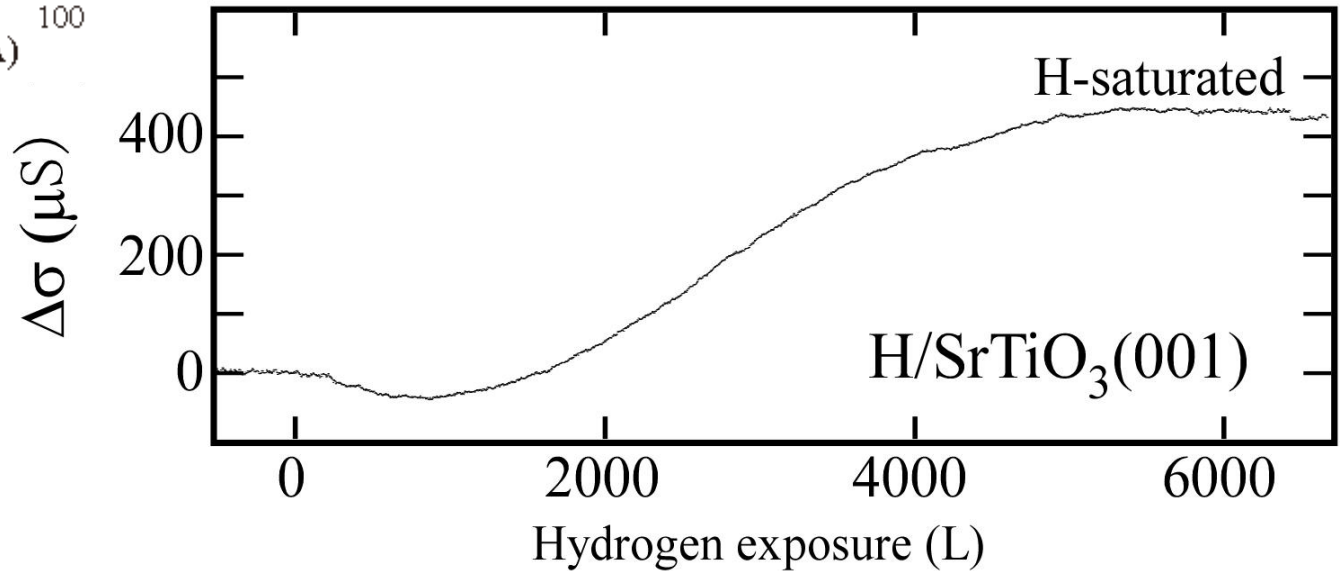
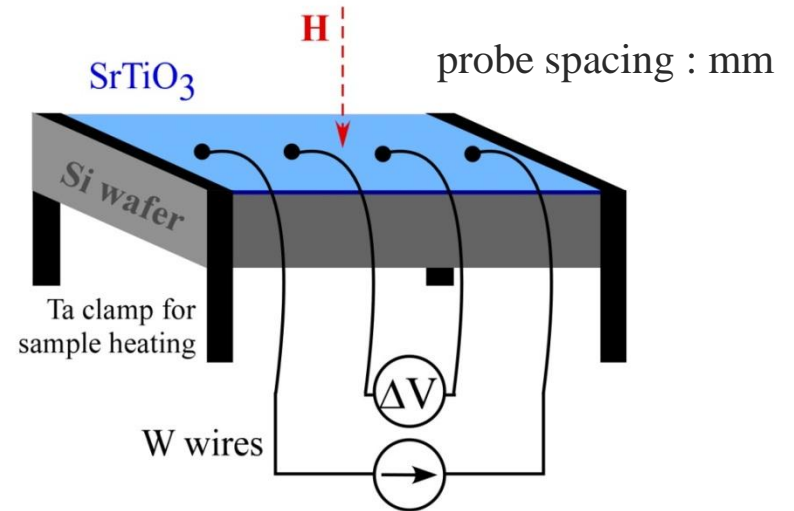
- 0.05wt %Nb-doped (n-type)
- TiO₂-terminated (HF treatment)
- Heat treatment in oxygen gas (free-oxygen vacancy)
- H atoms by a hot filament in hydrogen gas



H/SrTiO₃(001)



(a)



$$\Delta\sigma = \sigma_{\text{H/STO}} = 440 \mu\text{S} > 39 \mu\text{S} \quad (\text{2-D Ioffe-Regel limit})$$

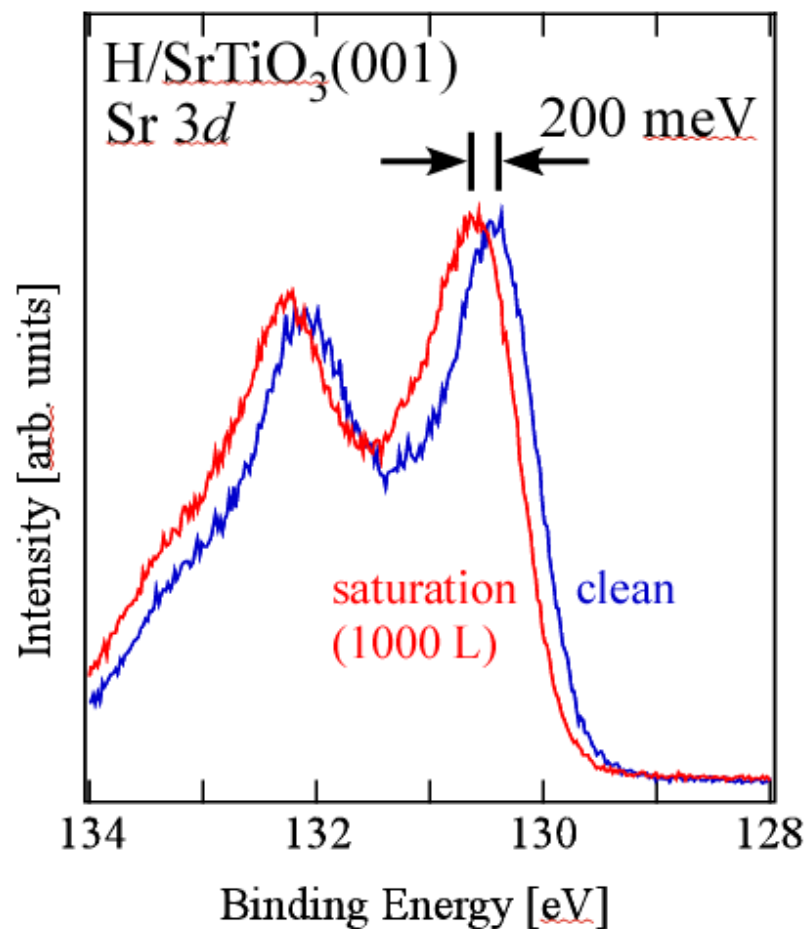
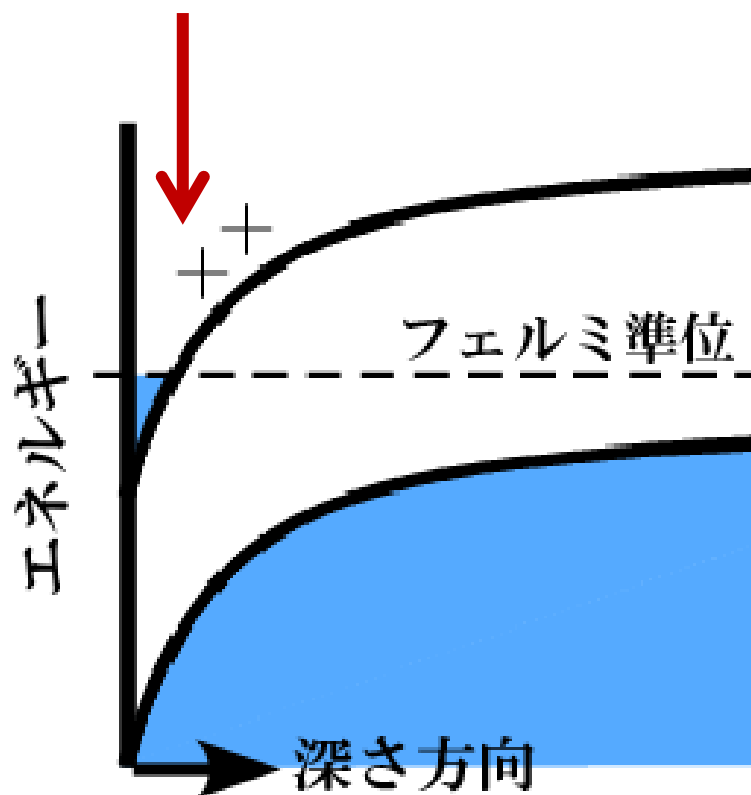
Band transport regime

H/SrTiO₃(001)

Photoemission Experiments

- SPring-8 BL07LSU

**Downward energy shift
for the surface metallization**



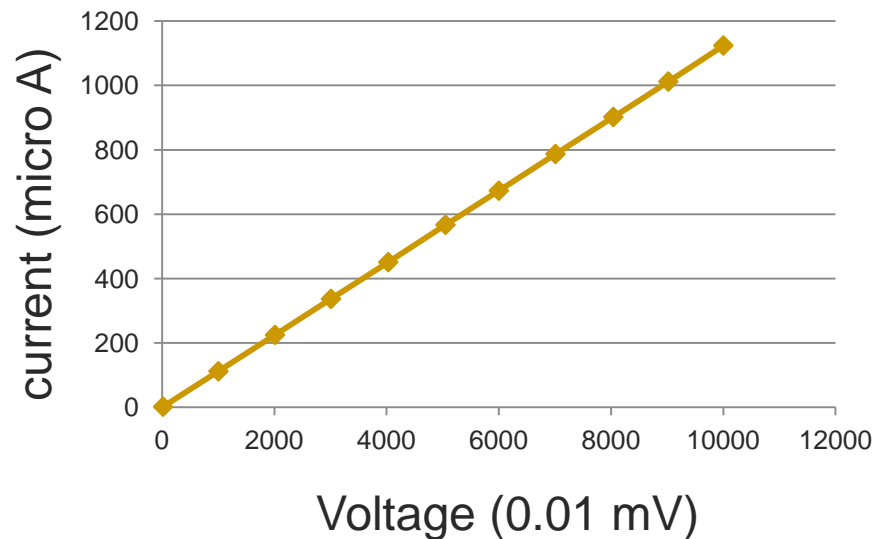
H/SrTiO₃(001)

Note:

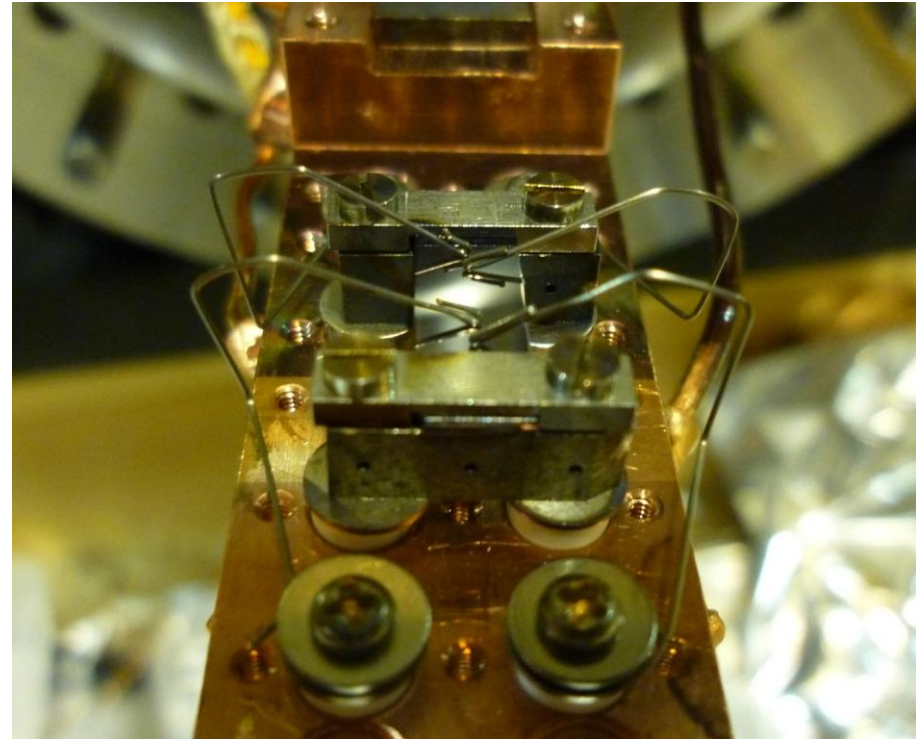
Photoemission and Transport Experiments

- SPring-8 BL07LSU

異なる手法の同時測定が可能に



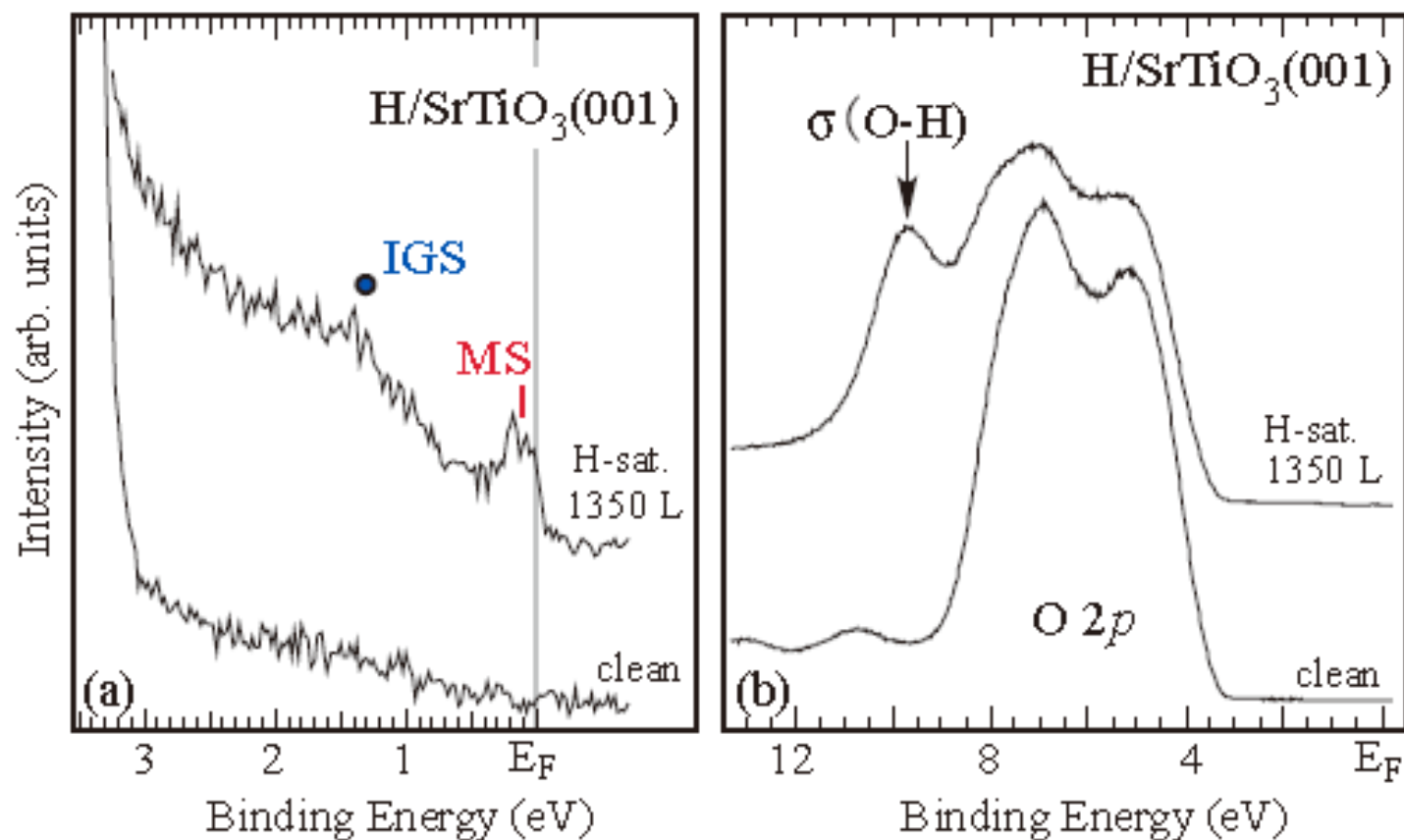
Photoemission and transport measurements
on the same sample



H/SrTiO₃(001)

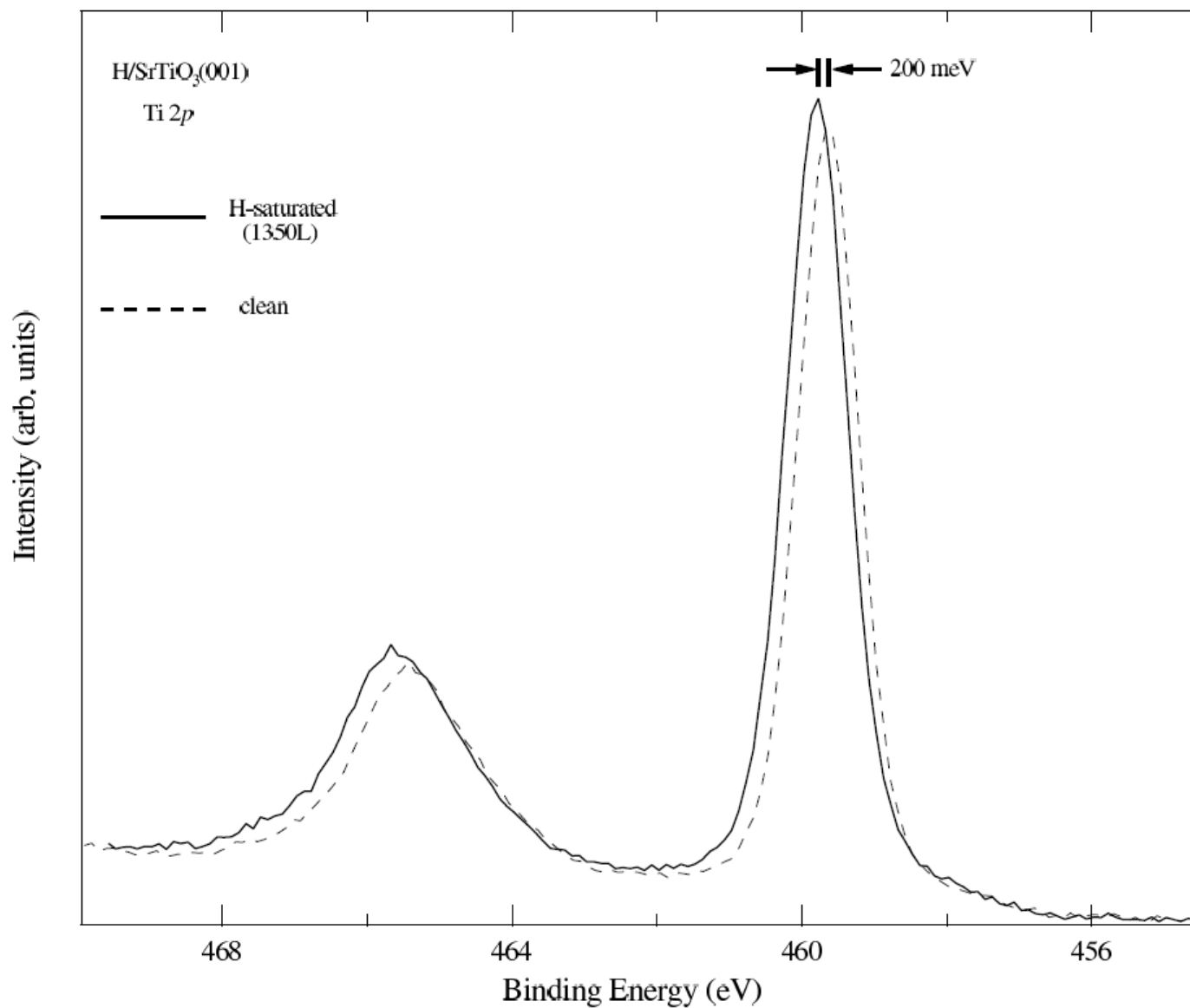
Photoemission Experiments

- SOLEIL TEMPO

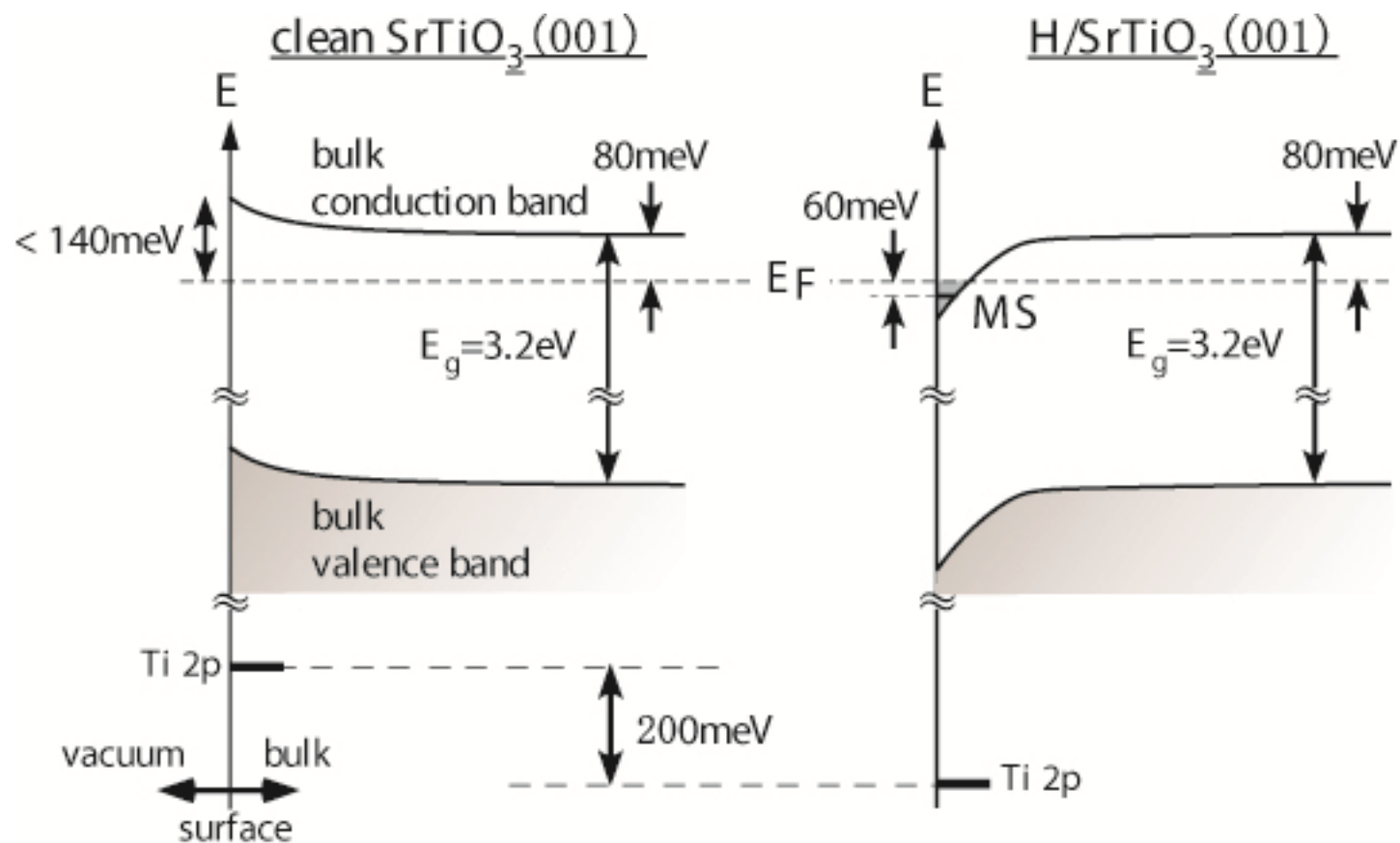


$h\nu=81$ eV, tuned to match with the minimum of Ti 3d bulk conduction band

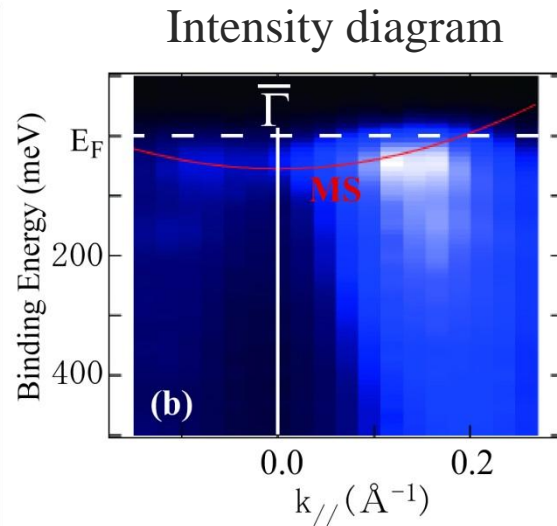
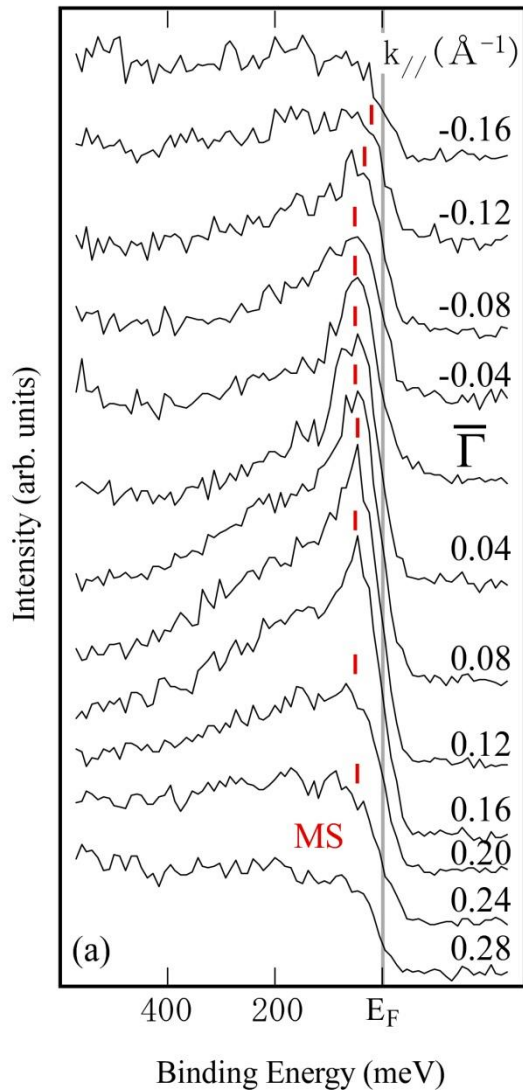
H/SrTiO₃(001)



H/SrTiO₃(001)

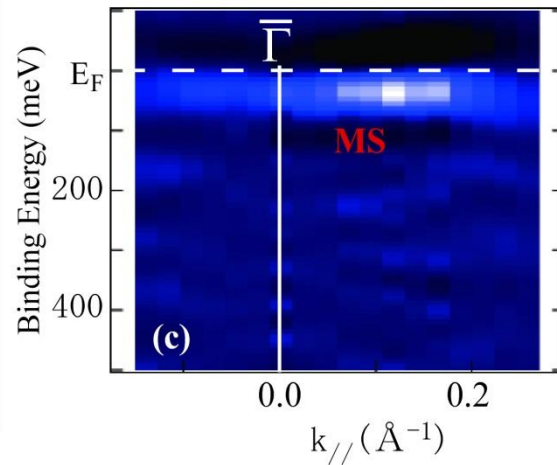


H/SrTiO₃(001)



$$k_F = 0.2 \text{ \AA}^{-1}$$

Mean free path
 $l = 2 \text{ nm (RT)}$



Broad spectral feature,
appeared at high binding energy

The second derivative

H/SrTiO₃(001)

As expected, not only the MS peak but also peaks, IGS and “broad feature”, were found.

“broad feature”

- Polaron formation due to the electron-phonon interaction [Y.J.Chang *et al.*, PRB 81, 235109(2010).]
- Electron correlation effect (Fermi liquid) [W. Meevasana *et al.*, Nature Materials, **10**, 114 (2011).]

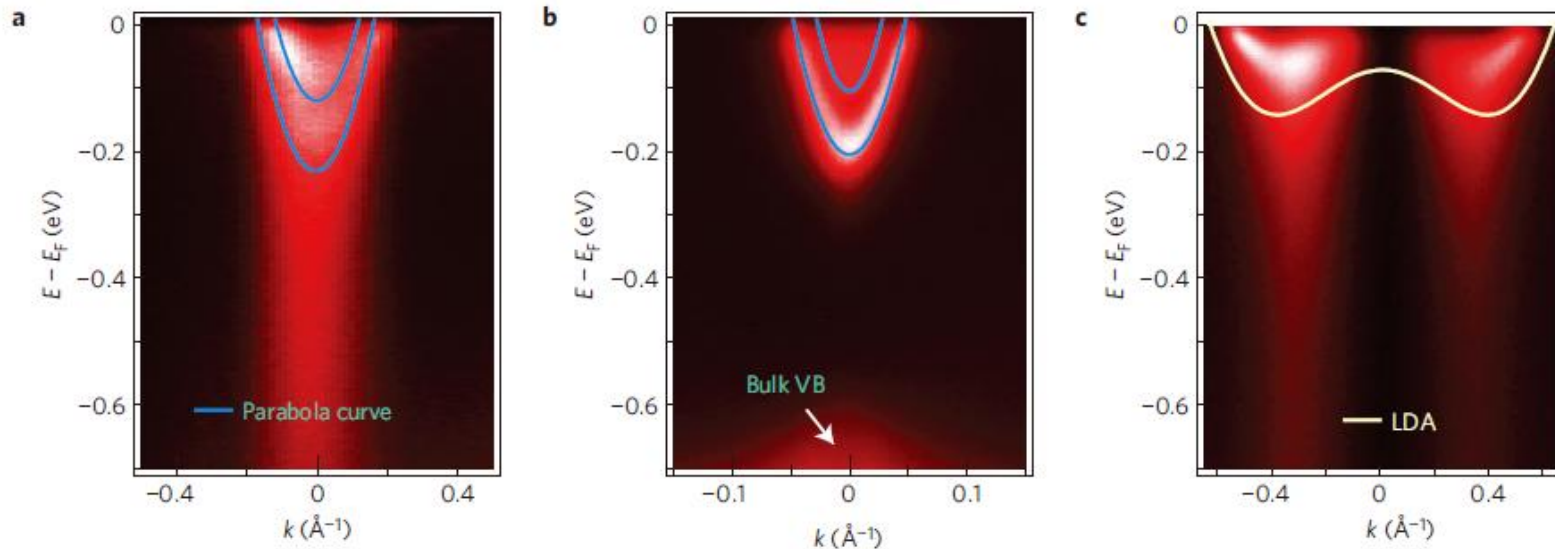
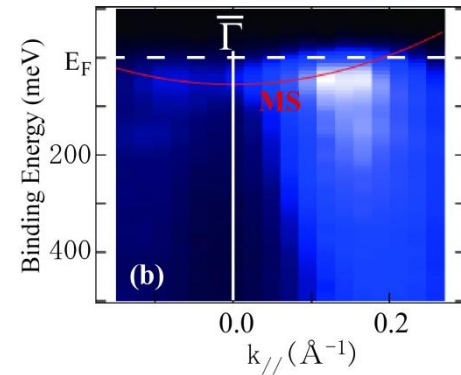


Figure 4 | Comparison of ARPES data from SrTiO₃, InAs and Bi₂Sr₂CuO₆ samples. a,b, 2DEG states at the surface of SrTiO₃ (a) and InAs (b). **c**, ARPES data from the single-layer cuprate Bi₂Sr₂CuO₆ (Bi2201) along $(\pi, 0)$ to $(0, \pi)$. The lines in **a** and **b** are parabolic dispersion relations, to guide the eye, and the result of a local density approximation band-structure calculation in **c**.

Summary

- A solid surface is an interesting playground to study low-dimensional electronic system.
- Surface insulator-to-metal transition or surface metallization was found by hydrogen adsorption on a SrTiO_3 surface.

Spectral features, likely related to the many body effects, were observed.

M. D'angelo *et al.*, Phys. Rev. Lett. accepted

- The VUV-soft X-ray photoemission research is now extending to the time-resolved measurement.

- **The optically-induced surface metal-to-insulator transition**

- fs-ns time: dynamics of the correlated carriers
- ps- ms time: a model for photo-chemistry

