

# FEMTOSECOND-LASER INDUCED DYNAMICS OF CO ON Ru(0001): NEW INSIGHTS FROM A HOT-ELECTRON, ELECTRONIC FRICTION MODEL INCLUDING SURFACE MOTION

Robert Scholz<sup>1,2</sup>, Gereon Floß<sup>1</sup>, Peter Saalfrank<sup>1</sup>, Gernot Fuchsels<sup>3</sup>, Ivor Lončarić<sup>4</sup>, and J. I. Juaristi<sup>4,5,6</sup>

<sup>1</sup>Institut für Chemie, Universität Potsdam, Karl-Liebknecht-Str. 24-25, D-14476 Potsdam, Germany

<sup>2</sup>Fritz-Haber-Institut der Max-Planck-Gesellschaft, Faradayweg 4-6, D-14195 Berlin, Germany

<sup>3</sup>Universiteit Leiden, Gorlaeus Laboratories, Einsteinweg 55, 2333 Leiden, The Netherlands

<sup>4</sup>Centro de Física de Materiales CFM/MPC (CSIC-UPV/EHU), Paseo Manuel de Lardizabal 5, 20018 Donostia-San Sebastián, Spain

<sup>5</sup>Departamento de Física de Materiales, Facultad de Químicas, Universidad del País Vasco (UPV/EHU), Apartado 1072, 20080 San Sebastián, Spain

<sup>6</sup>Donostia International Physics Center DIPC, P. Manuel de Lardizabal 4, 20018 San Sebastián, Spain

## Introduction

### Motivation

- research on small molecules adsorbed to metals is important for:
  - catalytic applications
  - fundamental understanding of bonding
- femtosecond(fs)-lasers are a valuable tool for such research as they
  - allow for investigations on small timescales
  - open up new processes compared to heating (femtochemistry)
  - may enable specific control over catalytic reactions (photocatalysis)
- specific motivation for system CO/Ru(0001)
  - experimentally well studied regarding fs-laser irradiation, e.g. [1, 2]
  - fulldimensional *ab-initio* potential recently developed in our group[3]
  - details of this indicate interpretation of experiment [2] may be wrong

### How does fs-laser-irradiation affect metal surfaces?

**fs-Laser** → **Desorption**  
**Diffusion (and possibly Reactions)**

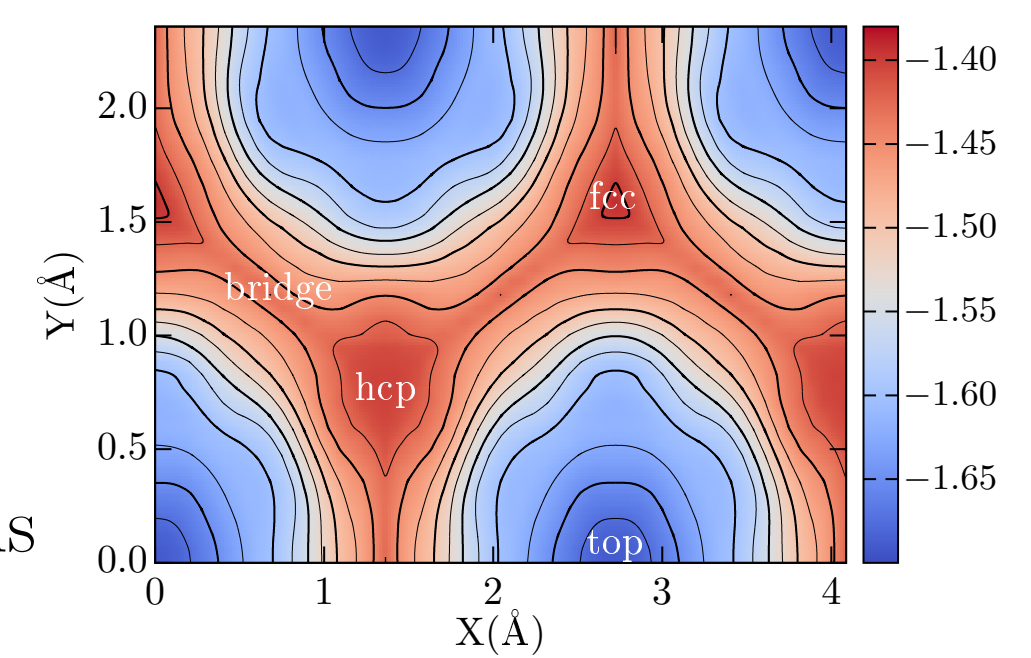
① **Electron-phonon coupling**  
② **Electronic friction**  
③ **Phonon-adsorbate interaction**

metals: ion lattice plus quasi-free electron gas  
visible light is absorbed only by the electrons  
produced electron-hole pairs thermalize quickly ⇒ “hot” Fermi-Dirac-distribution (after ~10 fs)  
electrons transfer part of energy to ion lattice, via ① **electron-phonon coupling** (phonons = lattice vibrations; quasi-particles)  
– electrons couple to phonons as their fast movement causes “shockwaves” in ion lattice  
– equilibration process completes after ~1 ps  
⇒ Thus, with fs-lasers, two different temperatures:  
–  $T_{el}$  - electron temperature  
–  $T_{ph}$  - phonon temperature  
• can be simulated using a Two-Temperature Model (2TM)[4] (see right)

## Models and Methods

### Six-dimensional Potential Energy Surface (6D PES)[3]

- Basis for dynamics: precomputed PES from DFT (rPBE + D2)
  - all 6 dimensions of the adsorbate
  - analytical PES and gradients ⇒ very fast
- ⇒ number and length of trajectories can be large
- downsides: – surface atoms frozen ⇒ no phonons  
– had to be constructed first



### Two-Temperature Model (2TM)[4]

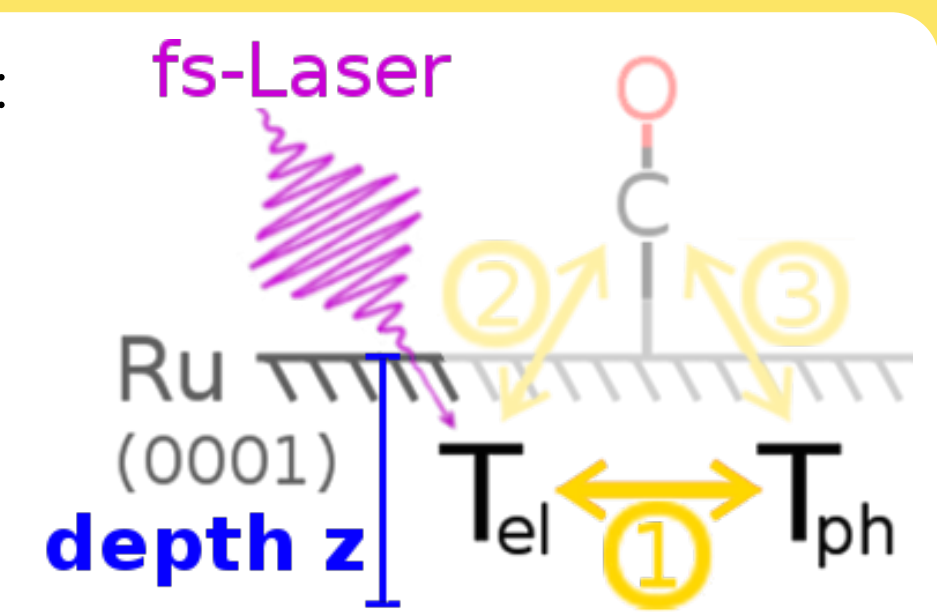
- describes interaction of metal with laser, using two differential equations:

$$C_{el} \frac{\partial T_{el}}{\partial t} = \frac{\partial}{\partial z} \kappa \frac{\partial T_{el}}{\partial z} - g(T_{el} - T_{ph}) + S(z, t),$$

$$C_{ph} \frac{\partial T_{ph}}{\partial t} = g(T_{el} - T_{ph}).$$

⇒ get  $T_{el}$  and  $T_{ph}$  as  $f(z, t)$  from laser parameters and material properties:

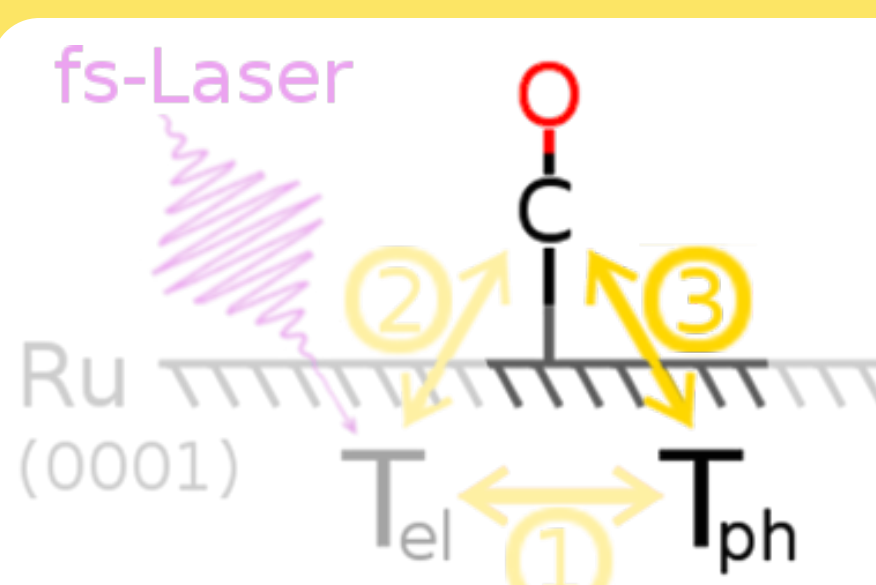
- laser wavelength  $\lambda$  (affects penetration depth into material)
- (effective) absorbed fluence  $F$  (energy/area)
- pulse duration  $\tau$  (all three appear in the “source term”  $S(z, t)$ )
- electron and phonon heat capacities  $C_{el}$  and  $C_{ph}$
- electron heat conductivity  $\kappa$
- electron-phonon coupling constant  $g$



### Electronic Friction: Langevin Dynamics[5] and Local Density Friction Approximation (LDFA)[6]

- Langevin equation of motion, a stochastic differential equation:
 
$$m_k \frac{d^2 r_k}{dt^2} = \underbrace{-\nabla_k V(r_1, r_2)}_{\text{Force on Atom } k} - \underbrace{\eta_{el,k}(r_k)}_{\text{Friction force slows movement}} \frac{dr_k}{dt} + \underbrace{R_{el,k}(t)}_{\text{Random force from e-h pairs}}$$
- describes movement of CO on the PES and interaction with electron-hole pairs (friction and excitation)
- Local Density Friction Approx. (LDFA): most simple model to calculate friction coefficients  $\eta_{el,k}$

### Inclusion of Phonons: Generalized Langevin Oscillator(GLO)-model[7, 8, 9]



## References

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