

# Recent progress in Casimir physics for atom-surface interactions

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# Collaborators

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# Outline of this talk

## ○ Part I

- Review of theory and experiment on Casimir atom-surface interactions

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- First measurement of thermal effect in Casimir physics (courtesy of Mauro Antezza, JILA-Trento collaboration)

## ○ Part III

- Casimir-Polder forces within scattering theory
- Cold atoms for probing lateral Casimir-Polder forces

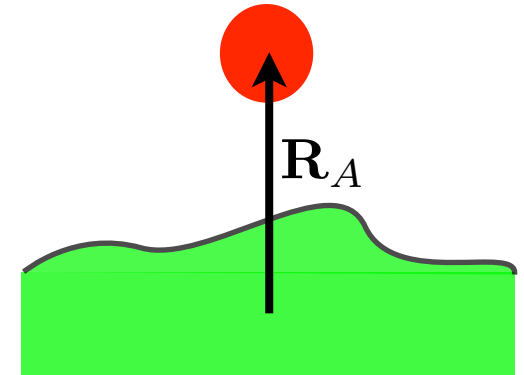
# Part I: Review

# The Casimir-Polder force

## ■ vdW - CP interaction Casimir and Polder (1948)

The interaction energy between a ground-state atom and a surface is given by

$$U_{\text{CP}}(\mathbf{R}_A) = \frac{\hbar}{c^2 \epsilon_0} \int_0^\infty \frac{d\xi}{2\pi} \xi^2 \alpha(i\xi) \text{Tr} \mathbf{G}(\mathbf{R}_A, \mathbf{R}_A, i\xi)$$



Atomic polarizability:  $\alpha(\omega) = \lim_{\epsilon \rightarrow 0} \frac{2}{3\hbar} \sum_k \frac{\omega_{k0} |\mathbf{d}_{0k}|^2}{\omega_{k0}^2 - \omega^2 - i\omega\epsilon}$

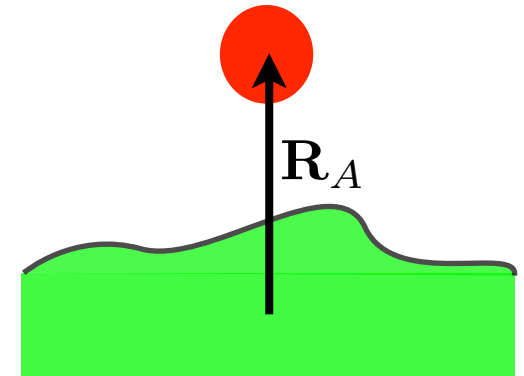
Scattering Green tensor:  $\left( \nabla \times \nabla \times - \frac{\omega^2}{c^2} \epsilon(\mathbf{r}, \omega) \right) \mathbf{G}(\mathbf{r}, \mathbf{r}', \omega) = \delta(\mathbf{r} - \mathbf{r}')$

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## ■ Eg: Ground-state atom near planar surface @ T=0

Non-retarded (vdW) limit  $z_A \ll \lambda_A$

$$U_{\text{vdW}}(z_A) = -\frac{\hbar}{8\pi\epsilon_0} \frac{1}{z_A^3} \int_0^\infty \frac{d\xi}{2\pi} \alpha(i\xi) \frac{\epsilon(i\xi) - 1}{\epsilon(i\xi) + 1}$$

Retarded (CP) limit  $z_A \gg \lambda_A$

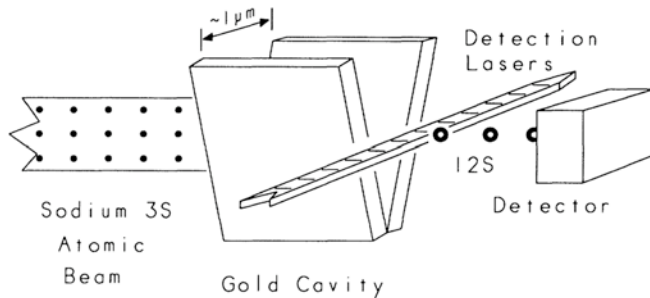
$$U_{\text{CP}}(z_A) = -\frac{3\hbar c \alpha(0)}{8\pi} \frac{1}{z_A^4} \frac{\epsilon_0 - 1}{\epsilon_0 + 1} \phi(\epsilon_0)$$



# Modern CP experiments

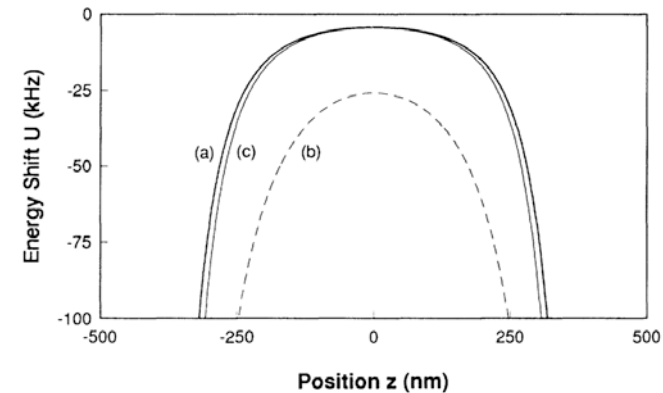
## Deflection of atoms

Hinds et al (1993)



$L = 0.7 - 1.2 \text{ } \mu\text{m}$

Exp-Th agreement @ 10%

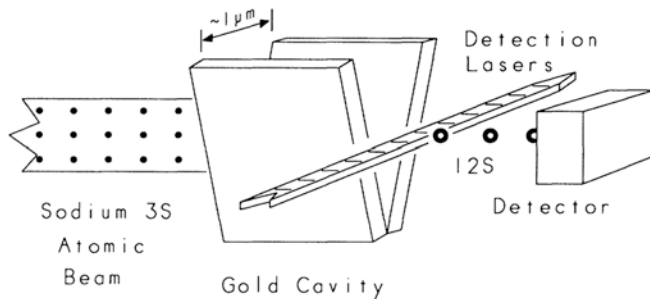


$$U_{CP} = -\frac{1}{4\pi\epsilon_0} \frac{\pi^3 \hbar c \alpha(0)}{L^4} \left[ \frac{3 - 2 \cos^2(\pi z/L)}{8 \cos^4(\pi z/L)} \right]$$

# Modern CP experiments

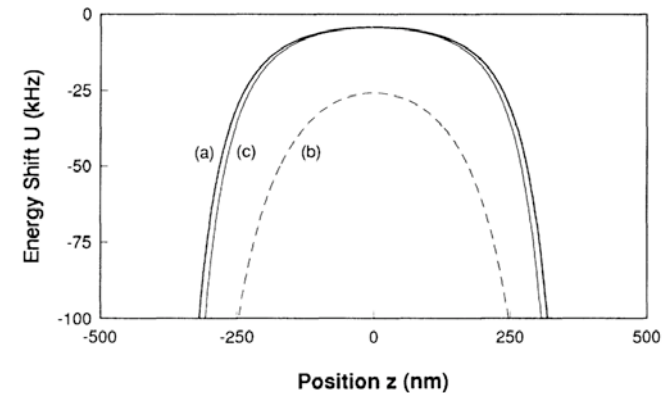
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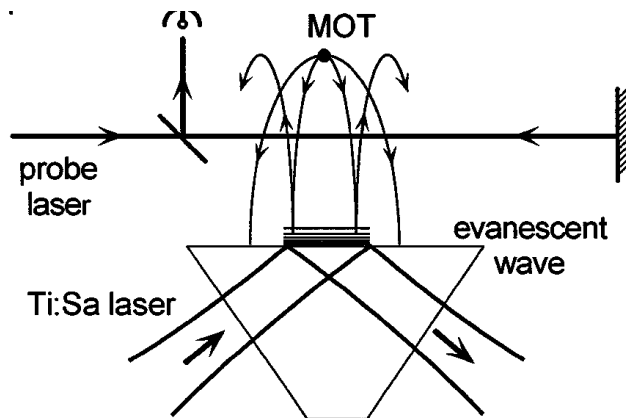
Exp-Th agreement @ 10%



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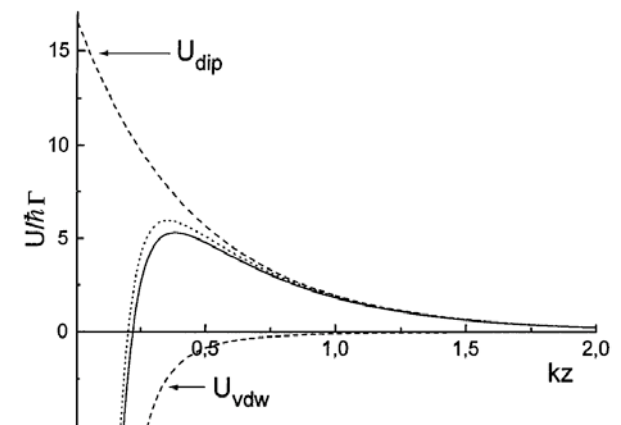
## Classical reflection on atomic mirror

Aspect et al (1996)



$$U_{\text{dip}} = \frac{\hbar}{4} \frac{\Omega^2}{\Delta} e^{-2kz}$$

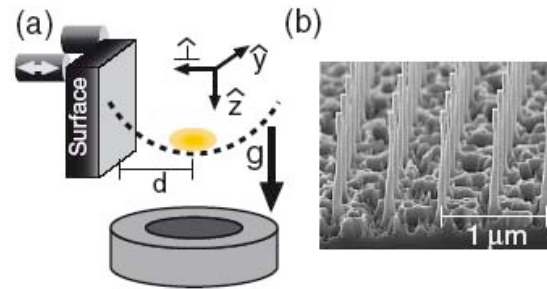
$$U_{\text{vdW}} = -\frac{\epsilon - 1}{\epsilon + 1} \frac{1}{48\pi\epsilon_0} \frac{D^2}{z^3}$$



Exp-Th agreement @ 30%

## Quantum reflection

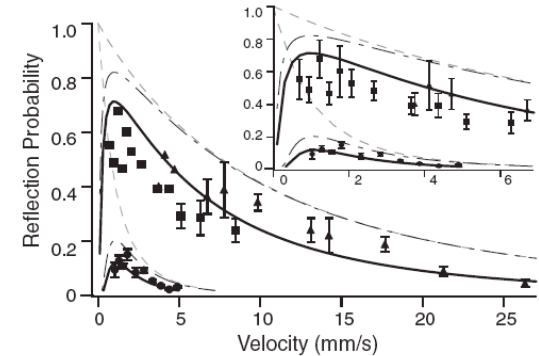
Wave-nature of atoms implies that slow atoms can reflect from purely attractive potentials



$$k = \sqrt{k_0^2 - 2mU/\hbar^2} \quad \phi = \frac{1}{k^2} \frac{dk}{dr} > 1$$

$$U = -C_n/r^n \quad (n > 2)$$

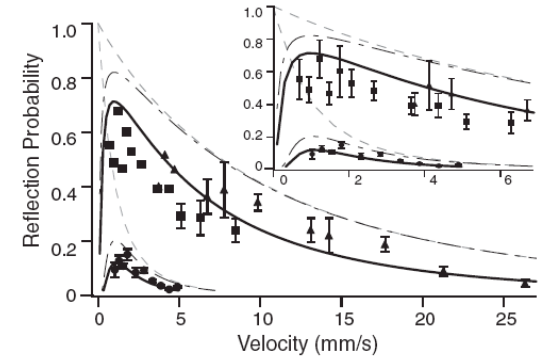
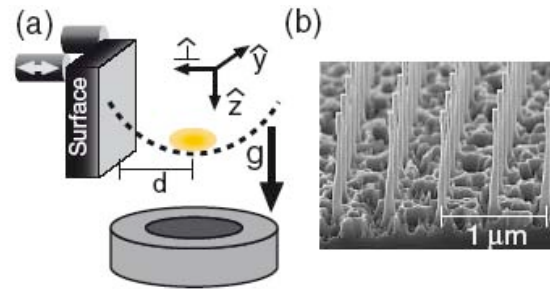
Shimizu (2001)    Ketterle et al (2006)  
DeKieviet et al (2003)



# Modern experiments (cont'd)

## Quantum reflection

Wave-nature of atoms implies that slow atoms can reflect from purely attractive potentials



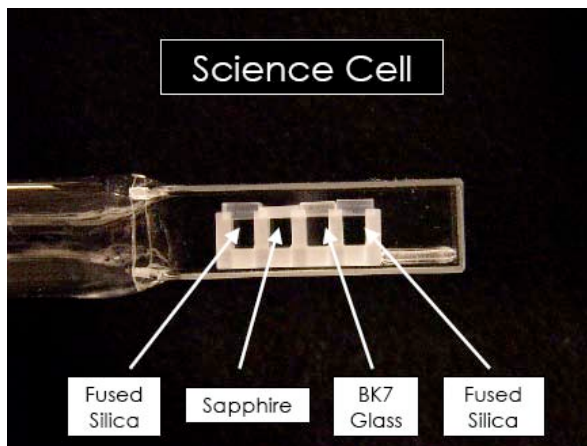
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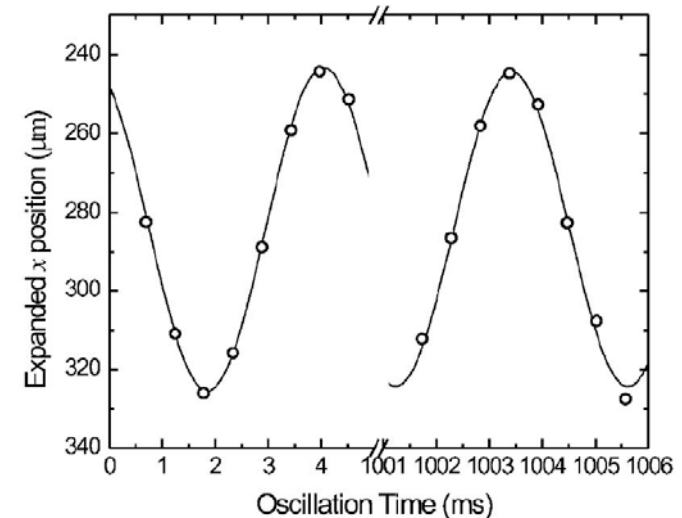
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## BEC oscillator

Cornell et al (2007)

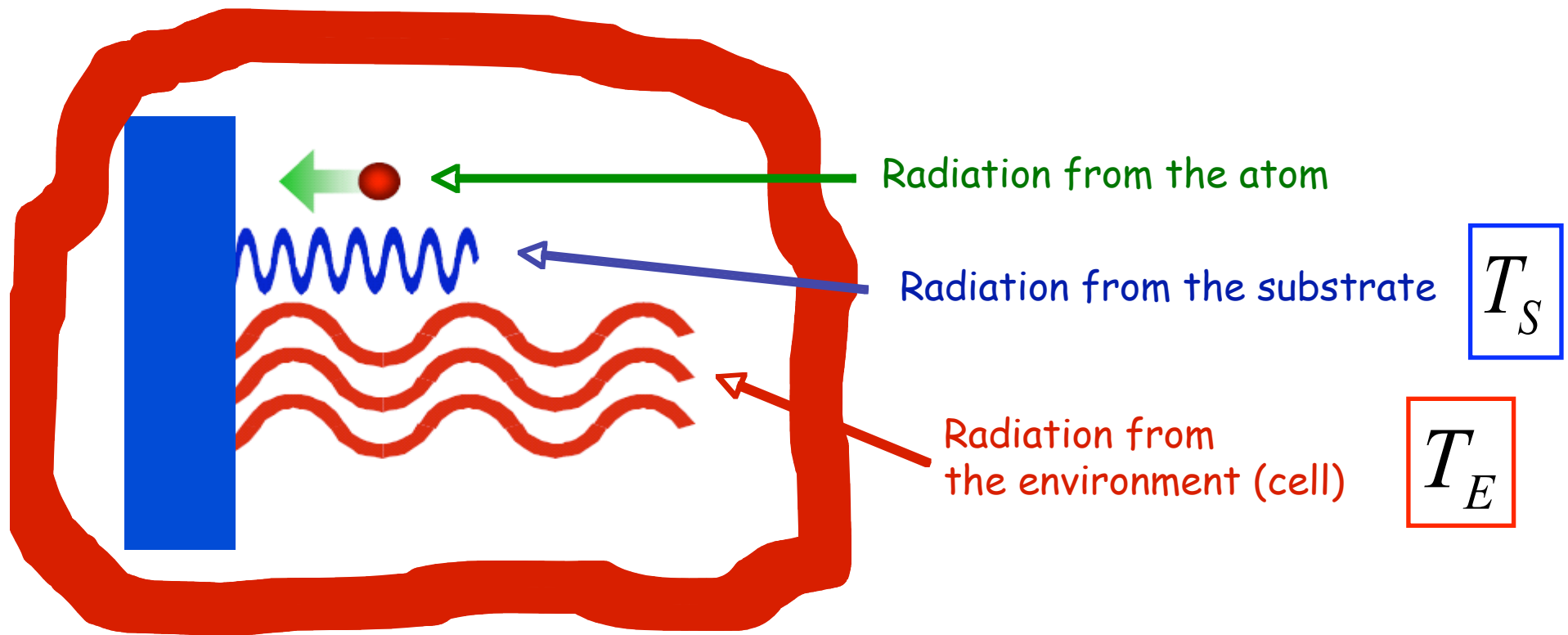


$$\gamma_x \equiv \frac{\omega_x - \omega'_x}{\omega_x} \approx -\frac{1}{2\omega_x^2 m} \partial_x^2 U^*$$



# Part II: Thermal CP force

## Surface-atom force



$$\vec{F}(\vec{r}) = \left\langle d_i^{tot}(t) \vec{\nabla} E_i^{tot}(\vec{r}, t) \right\rangle \approx \left\langle d_i^{ind}(t) \vec{\nabla} E_i^{fl}(\vec{r}, t) \right\rangle + \left\langle d_i^{fl}(t) \vec{\nabla} E_i^{ind}(\vec{r}, t) \right\rangle$$

Force includes **zero-point (or vacuum)** fluctuations effects + **thermal (or radiation)** fluctuations effects (**crucial at large distance!**)



## Large distance asymptotic behaviours

System at equilibrium

$$F^{eq} = -\frac{3k_B T \alpha_0 (\epsilon_0 - 1)}{4z^4 (\epsilon_0 + 1)}$$

E.M. Lifshitz, Dokl. Akad. Nauk. **100**, 879 (1955)

System out of equilibrium

substrate  environment 

$$F^{neq} = -\frac{\pi \alpha_0 k_B^2 (T_S^2 - T_E^2)}{6 z^3 c \hbar} \frac{\epsilon_0 + 1}{\sqrt{\epsilon_0 - 1}}$$

M.Antezza, L.P.Pitaevskii and S.Stringari, PRL. **95**, 093202 (2005)

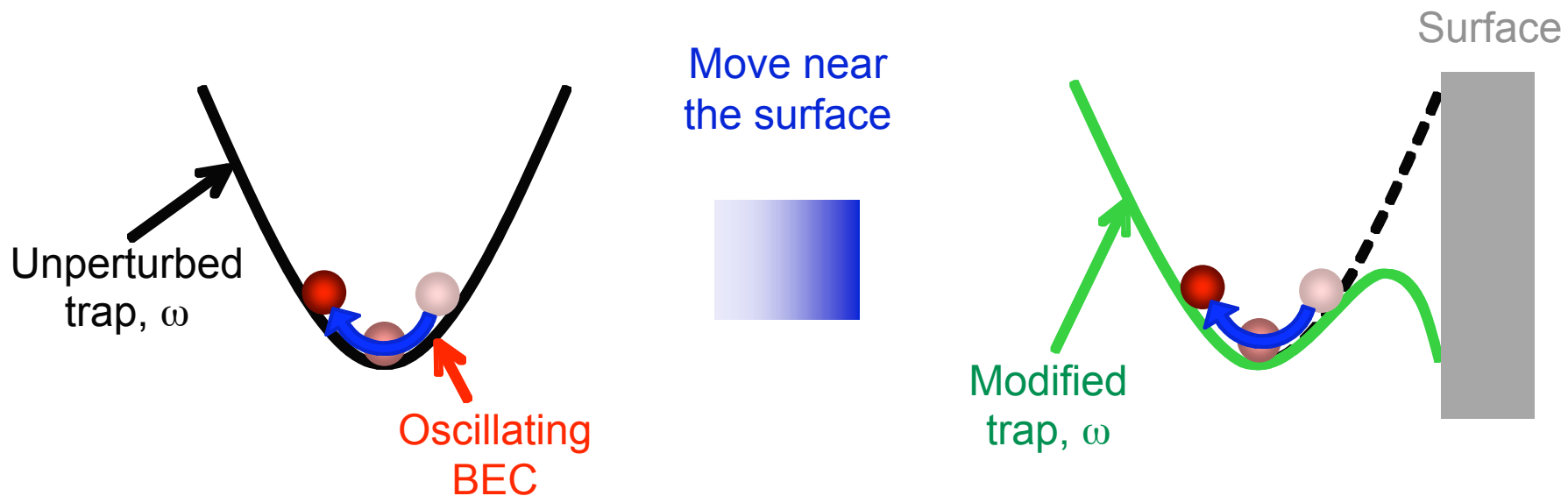
**NEW**

- ✓ force **decays slower than** at thermal **equilibrium**:
- ✓ force depends on **temperature** more **strongly** than at equilibrium
- ✓ force can be **attractive** or **repulsive** depending on relative temperatures of substrate and environment
- ✓ simple extension to **metals** (Drude model  $\epsilon'' = 4\pi\sigma / \omega$ )

# Measuring atom-surface interactions: dipolar oscillations of a BEC

**Use trapped BEC as a mechanical oscillator:**

Measure changes in oscillation frequency



Total force on the BEC is the sum of the forces on individual atoms. **Role of BEC coherence/superfluidity not central.** A BEC is convenient since it is a spatially compact collection of large number of particles, well controlled.



## Frequency shift of collective oscillations of a BEC

In M. Antezza, L.P. Pitaevskii and S. Stringari, PRA 70, 053619 (2004), the surface-atom force has been calculated and used to predict the frequency shift of the center of mass oscillation of a trapped Bose-Einstein condensate, including:

- **Effects of finite size** of the condensate
- **Non harmonic effects** due to the finite amplitude of the oscillations
- Dipole (center of mass) and quadrupole (long living mode) **frequency shifts**

In the presence of harmonic potential + surface-atom force frequency of center of mass motion is given by

$$V_{ho}(\vec{r}) = \frac{m}{2}\omega_x^2 x^2 + \frac{m}{2}\omega_y^2 y^2 + \frac{m}{2}\omega_z^2 z^2$$

$$\omega_{cm}^2 - \omega_z^2 = \frac{1}{m} \int n_0(\vec{r}) \partial_z^2 V_{surf-at}(z) d\vec{r} + \frac{a^2}{8m} \int n_0(\vec{r}) \partial_z^4 V_{surf-at}(z) d\vec{r}$$

Linear approximation

First non-linear correction

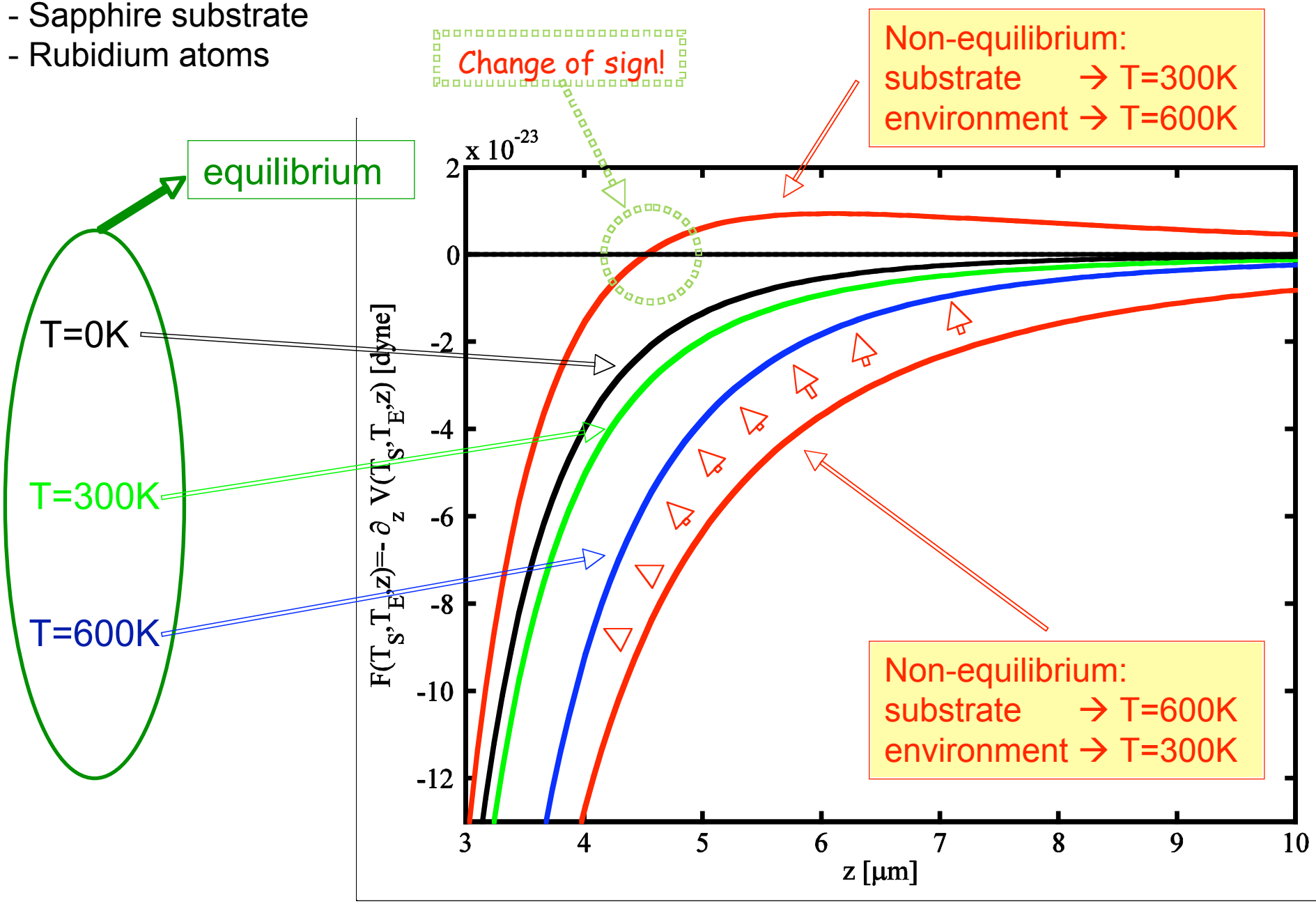
$a$  = amplitude of c.m. oscillation

$$Z_{cm} = Z_0 + a \cos(\omega t)$$

$n_0(r) \equiv$  Thomas-Fermi inverted parabola

# Thermal effects on the surface-atom force

- Sapphire substrate
- Rubidium atoms



equilibrium

Non-equilibrium:  
substrate  $\rightarrow T=300\text{K}$   
environment  $\rightarrow T=600\text{K}$

Non-equilibrium:  
substrate  $\rightarrow T=600\text{K}$   
environment  $\rightarrow T=300\text{K}$

Change of sign!

$T=0\text{K}$

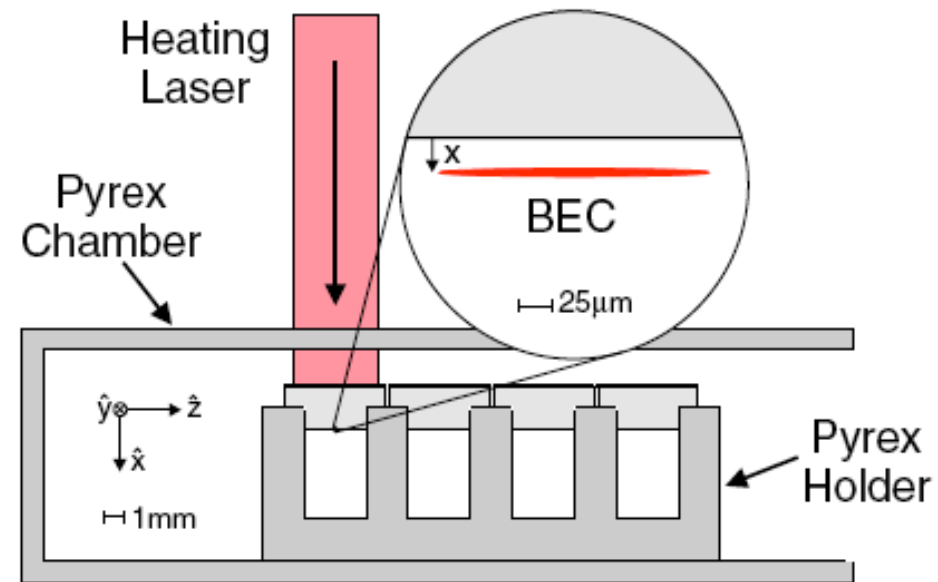
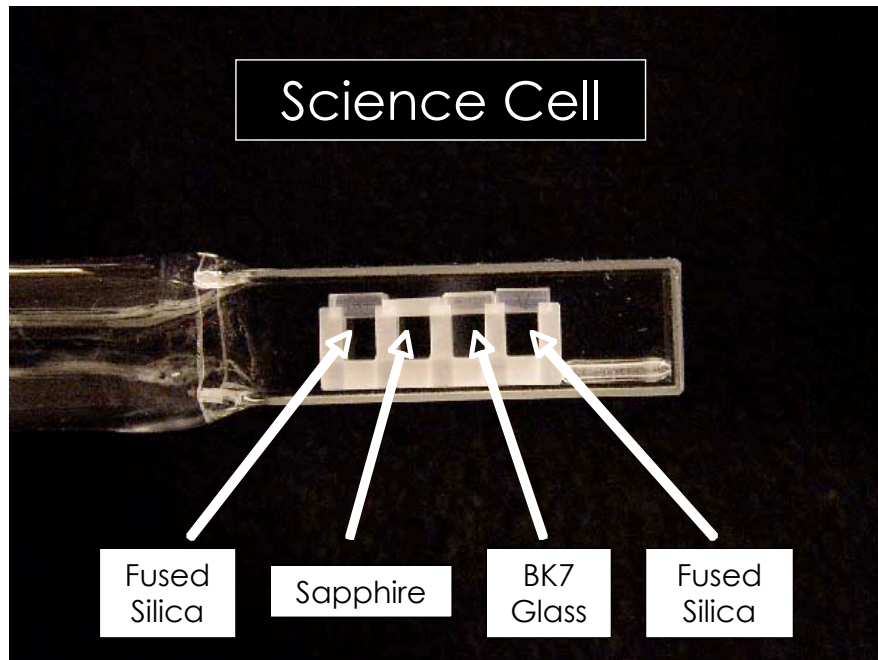
$T=300\text{K}$

$T=600\text{K}$

$$F(T_S, T_E, z) = -\partial_z V(T_S, T_E, z) \text{ [dyne]}$$

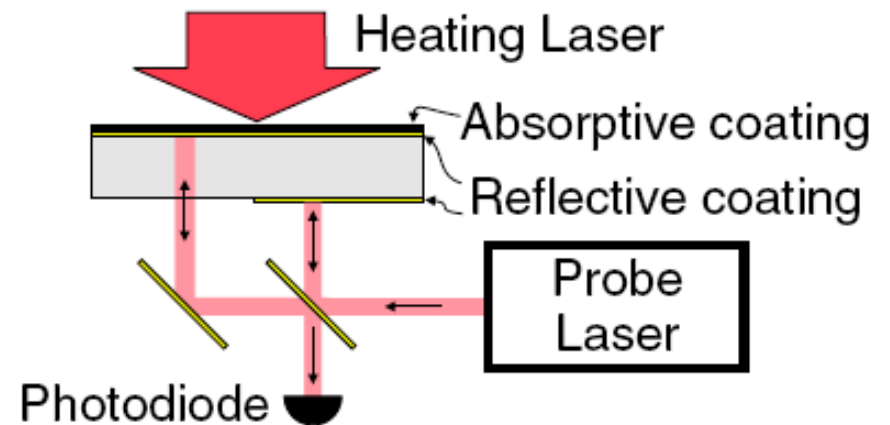
$z \text{ [}\mu\text{m]}$

## The experimental apparatus



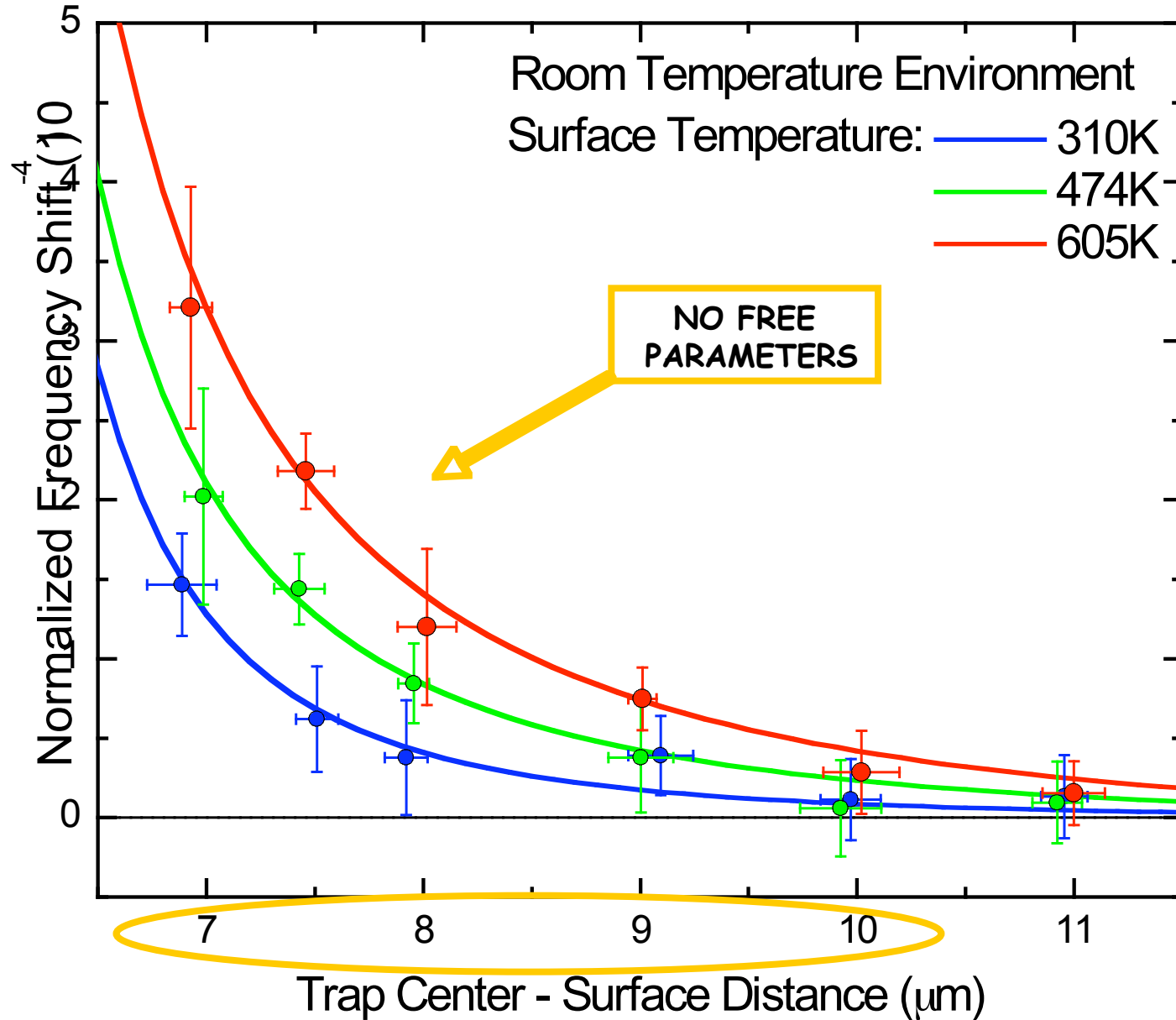
- ✓ Multiple dielectric surfaces! Amorphous glass, crystalline sapphire.
- ✓ No conducting objects near atoms!

Interferometric measurement  
of temperature of substrate:



Experimental results from JILA

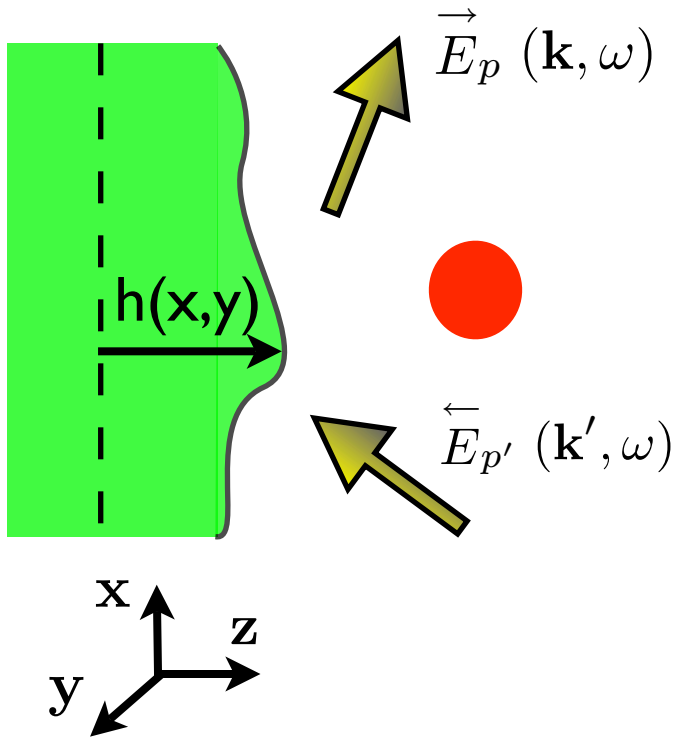
OUT OF EQUILIBRIUM



# Part III: CP and scattering theory



# CP within scattering theory



Output fields:  $\vec{\mathbf{E}}(\mathbf{R}, \omega) = \int \frac{d^2\mathbf{k}}{(2\pi)^2} e^{i\mathbf{k}\cdot\mathbf{r}} \vec{\mathbf{E}}(\mathbf{k}, z, \omega)$

$$\vec{\mathbf{E}}(\mathbf{k}, z, \omega) = [\vec{E}_{\text{TE}}(\mathbf{k}, \omega) \hat{\epsilon}_{\text{TE}}^+(\mathbf{k}) + \vec{E}_{\text{TM}}(\mathbf{k}, \omega) \hat{\epsilon}_{\text{TM}}^+(\mathbf{k})] e^{ik_z z}$$

$$\hat{\epsilon}_{\text{TE}}^+(\mathbf{k}) = \mathbf{z} \times \mathbf{k} \quad \hat{\epsilon}_{\text{TM}}^+(\mathbf{k}) = \hat{\epsilon}_{\text{TE}}^+(\mathbf{k}) \times \mathbf{K} \quad (\mathbf{K} = \mathbf{k} + k_z \mathbf{z})$$

Input fields: idem with  $k_z \rightarrow -k_z$

Input-output fields related via reflection operators for the surface

$$\vec{E}_p(\mathbf{k}, \omega) = \int \frac{d^2\mathbf{k}'}{(2\pi)^2} \sum_{p'} \langle \mathbf{k}, p | \mathcal{R}(\omega) | \mathbf{k}', p' \rangle \vec{E}_{p'}(\mathbf{k}', \omega)$$

$$\mathcal{E} = \hbar \int_0^\infty \frac{d\xi}{2\pi} \text{Tr} \log(1 - \mathcal{R} e^{-\mathcal{K}z_A} \mathcal{R}_{\text{at}} e^{-\mathcal{K}z_A})$$

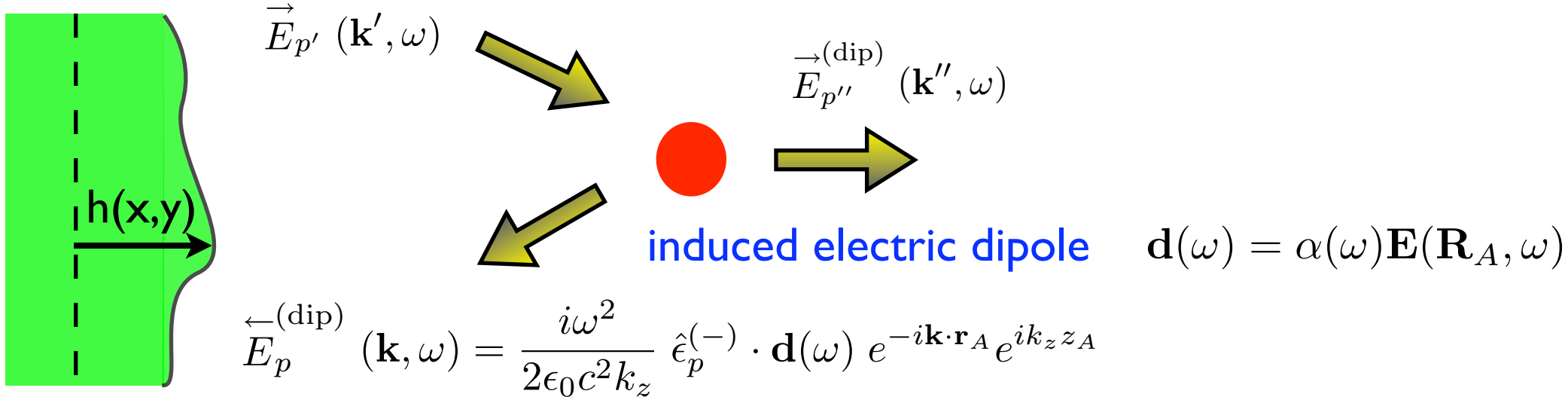
aka "TGTG" formula

$$\langle \mathbf{k}, p | \mathcal{K} | \mathbf{k}', p' \rangle = \delta^2(\mathbf{k} - \mathbf{k}') \delta_{pp'} \sqrt{k^2 + \xi^2/c^2}$$

$\mathcal{R}_{\text{at}}$  reflection operator for the atom ( $z_A = 0$ )

# Atom reflection operator

Calculating reflection operator for the atom:  $(\forall \mathbf{R}_A)$



$$\overset{\leftarrow}{E}_p^{(\text{dip})}(\mathbf{k}, \omega) = \frac{i\omega^2}{2\epsilon_0 c^2 k_z} \hat{\epsilon}_p^{(-)} \cdot \mathbf{d}(\omega) e^{-i\mathbf{k}\cdot\mathbf{r}_A} e^{ik_z z_A}$$

$$\overset{\leftarrow}{E}_p^{(\text{dip})}(\mathbf{k}, \omega) = \frac{i\omega^2 \alpha(\omega)}{2\epsilon_0 c^2 k_z} \int \frac{d^2 \mathbf{k}'}{(2\pi)^2} e^{-i(\mathbf{k}-\mathbf{k}')\cdot\mathbf{r}_A} e^{i(k_z+k'_z)z_A} \sum_{p'} \hat{\epsilon}_p^{(-)}(\mathbf{k}) \cdot \hat{\epsilon}_{p'}^{(+)}(\mathbf{k}') \overset{\rightarrow}{E}_{p'}(\mathbf{k}', \omega)$$

$$\langle \mathbf{k}, p | \mathcal{R}_{\text{at}}(\omega) | \mathbf{k}', p' \rangle = \frac{i\omega^2 \alpha(\omega)}{2\epsilon_0 c^2 k_z} e^{-i(\mathbf{k}-\mathbf{k}')\cdot\mathbf{r}_A} e^{i(k_z+k'_z)z_A} \hat{\epsilon}_p^{(-)}(\mathbf{k}) \cdot \hat{\epsilon}_{p'}^{(+)}(\mathbf{k}')$$

# From TGTG to GTG

Exact expression for the atom-surface  
interaction energy:

$$U_{\text{CP}}(\mathbf{R}_A) = \frac{\hbar}{c^2 \epsilon_0} \int_0^\infty \frac{d\xi}{2\pi} \xi^2 \alpha(i\xi) \int \frac{d^2 \mathbf{k}}{(2\pi)^2} \int \frac{d^2 \mathbf{k}'}{(2\pi)^2} e^{i(\mathbf{k}-\mathbf{k}') \cdot \mathbf{r}_A} e^{-(\kappa+\kappa')z_A} \frac{1}{2\kappa'} \sum_{p,p'} \hat{\epsilon}_p^+(\mathbf{k}) \cdot \hat{\epsilon}_{p'}^-(\mathbf{k}') R_{p,p'}(\mathbf{k}, \mathbf{k}')$$

with  $\kappa \equiv \sqrt{\xi^2/c^2 + k^2}$  and  $R_{p,p'}(\mathbf{k}, \mathbf{k}')$  dependent on material properties at freq.  $i\xi$

■ Analogous to known expressions based on Green function formalism

$$U_{\text{CP}}(\mathbf{R}_A) = \frac{\hbar}{c^2 \epsilon_0} \int_0^\infty \frac{d\xi}{2\pi} \xi^2 \alpha(i\xi) \text{Tr} \mathbf{G}(\mathbf{R}_A, \mathbf{R}_A, i\xi)$$

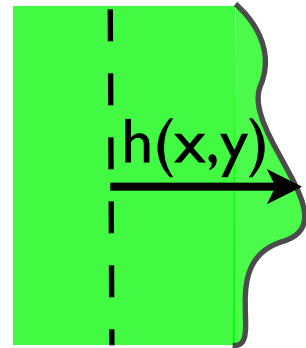
■ Remaining difficulty: calculation of the surface reflection matrix



# Perturbation theory

In order to treat a general rough or corrugated surface, we make a perturbative expansion in powers of  $h(x,y)$

$$\mathcal{R} = \mathcal{R}^{(0)} + \mathcal{R}^{(1)} + \dots$$



## □ Specular reflection:

$$\langle \mathbf{k}, p | \mathcal{R}^{(0)} | \mathbf{k}', p' \rangle = (2\pi)^2 \delta^{(2)}(\mathbf{k} - \mathbf{k}') \delta_{p,p'} r_p(\mathbf{k}, \xi)$$

Fresnel coefficients  $r_{\text{TE}} = \frac{\kappa - \kappa_t}{\kappa + \kappa_t}$   $r_{\text{TM}} = \frac{\epsilon(i\xi)\kappa - \kappa_t}{\epsilon(i\xi)\kappa + \kappa_t}$  ( $\kappa_t = \sqrt{\epsilon(i\xi)\xi^2/c^2 + k^2}$ )

## □ Non-specular reflection:

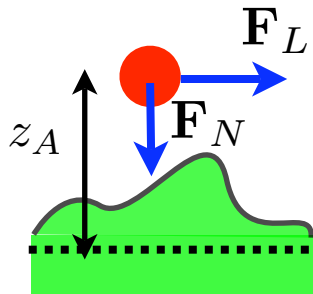
$$\langle \mathbf{k}, p | \mathcal{R}^{(1)} | \mathbf{k}', p' \rangle = R_{p,p'}(\mathbf{k}, \mathbf{k}') H(\mathbf{k} - \mathbf{k}') \leftarrow \text{Fourier transform of } h(x,y)$$



The non-specular reflection matrices depend on the geometry and material properties.

Greffet (1988), Reynaud et al (2005)

# Lateral Casimir-Polder force



$$U_{\text{CP}} = U_{\text{CP}}^{(0)}(z_A) + U_{\text{CP}}^{(1)}(z_A, x_A)$$

■ Normal CP force:

$$U_{\text{CP}}^{(0)}(z_A) = \frac{\hbar}{c^2 \epsilon_0} \int_0^\infty \frac{d\xi}{2\pi} \xi^2 \alpha(i\xi) \int \frac{d^2 \mathbf{k}}{(2\pi)^2} \frac{1}{2\kappa} \sum_p \hat{\epsilon}_p^+ \cdot \hat{\epsilon}_p^- r_p(\mathbf{k}, \xi) e^{-2\kappa z_A}$$

■ Lateral CP force:

$$U_{\text{CP}}^{(1)}(z_A, x_A) = \int \frac{d^2 \mathbf{k}}{(2\pi)^2} e^{i\mathbf{k} \cdot \mathbf{r}_A} g(\mathbf{k}, z_A) H(\mathbf{k})$$

Response function  $g$ :

$$g(\mathbf{k}, z_A) = \frac{\hbar}{c^2 \epsilon_0} \int_0^\infty \frac{d\xi}{2\pi} \xi^2 \alpha(i\xi) \int \frac{d^2 \mathbf{k}'}{(2\pi)^2} a_{\mathbf{k}', \mathbf{k}' - \mathbf{k}}(z_A, \xi)$$

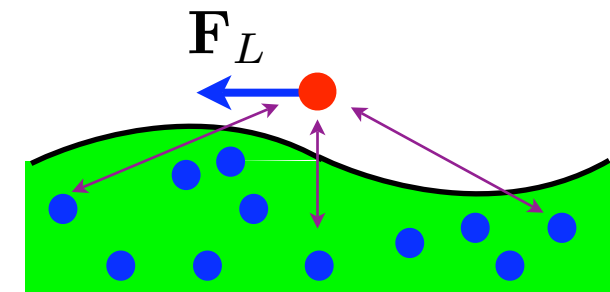
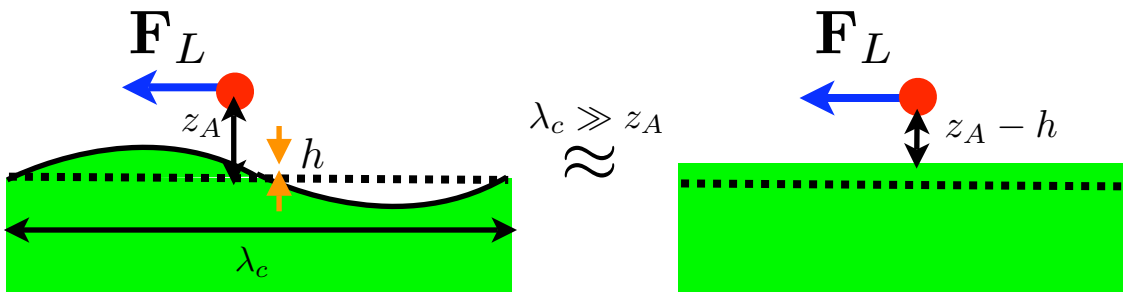
$$a_{\mathbf{k}', \mathbf{k}''} = \sum_{p', p''} \hat{\epsilon}_{p'}^+ \cdot \hat{\epsilon}_{p''}^- \frac{e^{-(\kappa' + \kappa'')z_A}}{2\kappa''} R_{p', p''}(\mathbf{k}', \mathbf{k}'')$$

Our approach is perturbative in  $h(x, y)$ , which should be the smallest length scale in the problem  $h \ll z_A, \lambda_c, \lambda_A, \lambda_0$

# Approx. methods: PFA & PWS

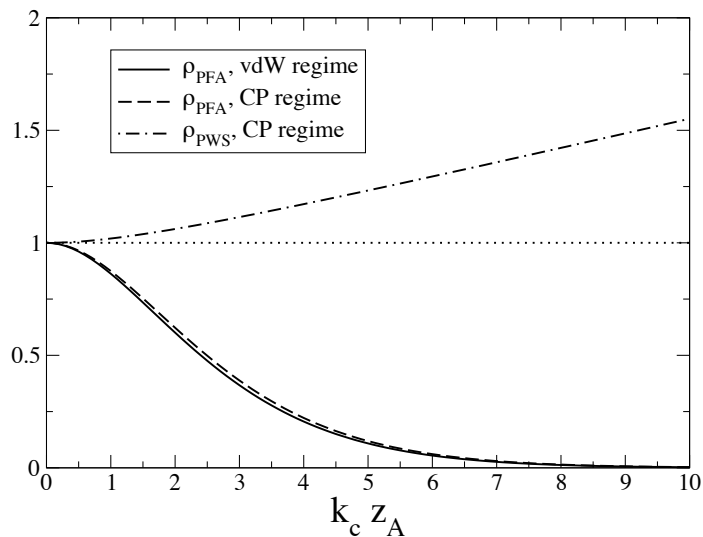
## Proximity Force Approximation (PFA)

## Pair-wise Summation (PWS)



## Deviations from PFA and PWS

$$\rho_{\text{PFA}} = \frac{g(k_c, z_A)}{g(0, z_A)} \quad \rho_{\text{PWS}} \equiv \frac{g(k_c, z_A)}{g_{\text{PWS}}(k_c, z_A)}$$



### Example:

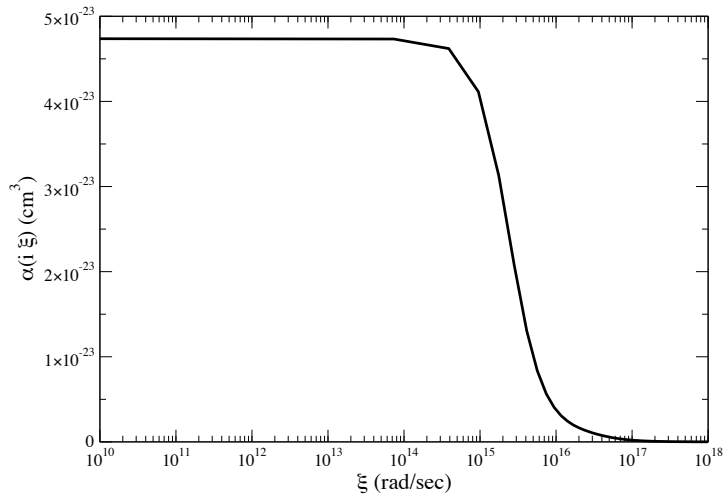
atom-surface distance  $z_A = 2\mu\text{m} \gg \lambda_A$

corrugation wavelength  $\lambda_c = 3.5\mu\text{m}$

➔  $\rho_{\text{PFA}} \approx 30\% \quad \rho_{\text{PWS}} \approx 115\%$

PFA largely overestimates the lateral CP force  
PWS underestimates the lateral CP force

**Dynamic polarizability of Rb**  
Babb et al (1999)



**Calculation of  $R_{p,p'}^{(1)}(\mathbf{k}, \mathbf{k}', \xi)$  in terms of  $\epsilon(i\xi)$  of bulk materials** Reynaud et al (2005)

$$R_{\text{TE,TE}}^{(1)}(\mathbf{k}, \mathbf{k}'; \xi) = 2\kappa C h_{\text{TE,TE}}(\mathbf{k}, \mathbf{k}', \xi)$$

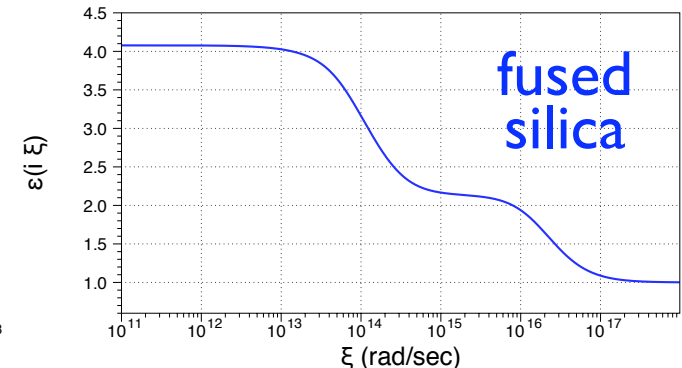
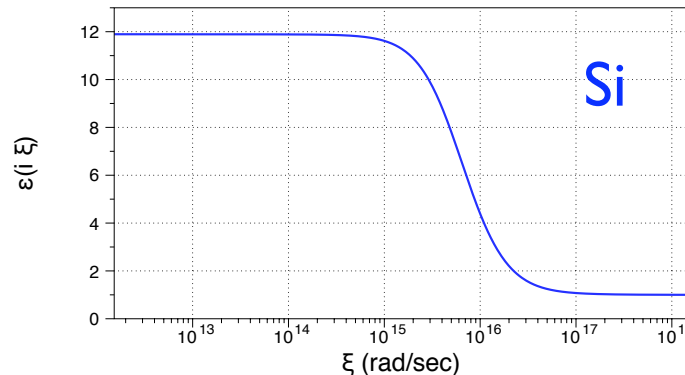
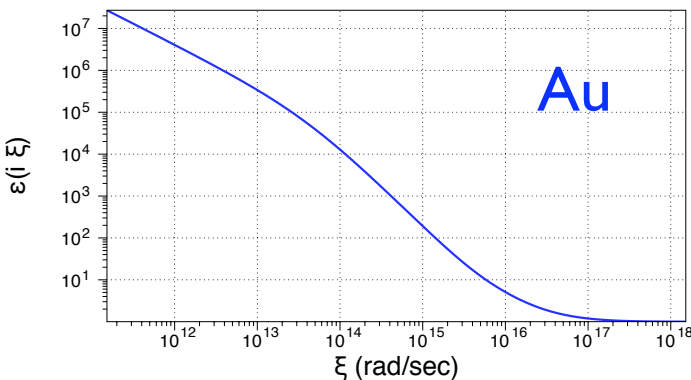
$$R_{\text{TE,TM}}^{(1)}(\mathbf{k}, \mathbf{k}'; \xi) = 2\kappa S \frac{c\kappa'_t}{\sqrt{\epsilon}\xi} h_{\text{TE,TM}}(\mathbf{k}, \mathbf{k}', \xi)$$

$$R_{\text{TM,TE}}^{(1)}(\mathbf{k}, \mathbf{k}'; \xi) = \frac{2\sqrt{\epsilon}\kappa\kappa'_t \frac{\xi}{c} S}{\left(\frac{\xi}{c}\right)^2 - (\epsilon + 1)\kappa^2} h_{\text{TM,TE}}(\mathbf{k}, \mathbf{k}', \xi)$$

$$R_{\text{TM,TM}}^{(1)}(\mathbf{k}, \mathbf{k}'; \xi) = -2\kappa \frac{\epsilon k k' + \kappa_t \kappa'_t C}{\left(\frac{\xi}{c}\right)^2 - (\epsilon + 1)\kappa^2} h_{\text{TM,TM}}(\mathbf{k}, \mathbf{k}', \xi)$$

$$h_{pp'}(\mathbf{k}, \mathbf{k}') = \frac{r^p(\mathbf{k}) t^{p'}(\mathbf{k}')}{t^p(\mathbf{k})}$$

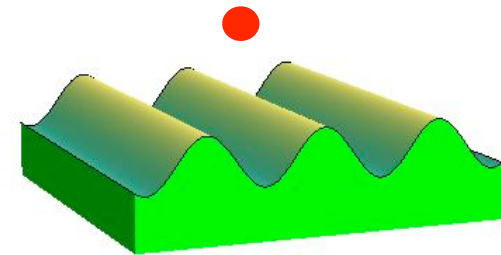
**Optical data + Kramers-Kronig relations**



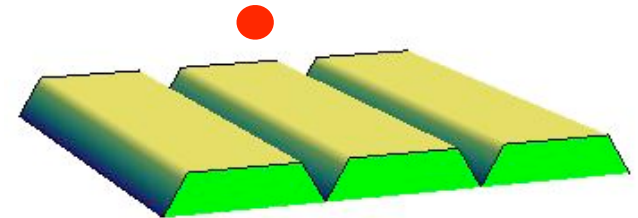
# Atoms as local probes

In contrast to the case of the lateral Casimir force between corrugated surfaces, an atom is a **local probe** of the lateral Casimir-Polder force. **Deviations from the PFA can be much larger than for the force between two surfaces!**

☑ **Deviations** from PFA/PWS can be obtained for a **sinusoidal corrugated surface**.



☑ **Even larger deviations** from PFA/PWS can be obtained for a **periodically grooved surface**.



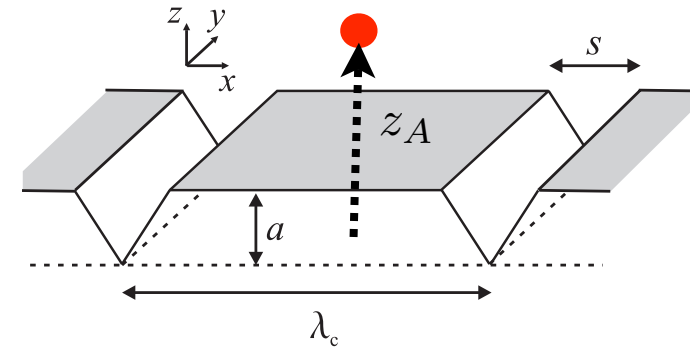
🔍 If the atom is located above one plateau, the PFA predicts that the lateral Casimir-Polder force should vanish. A non-vanishing force appearing when the atom is moved above the plateau thus clearly signals a deviation from PFA!

🔍 A lateral force appears for PWS, but it should be much smaller than the exact result.

# CP energy for grooved surface

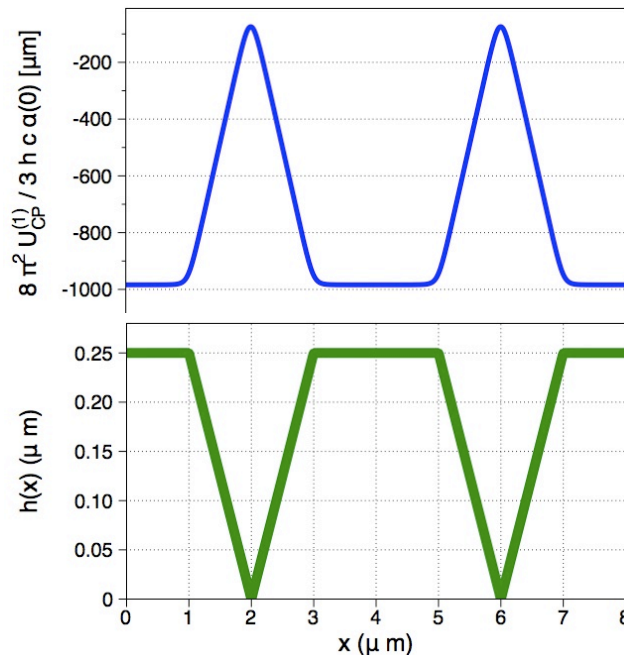
Surface profile for periodical grooved corrugation

$$h(x) = a \left( 1 - \frac{s}{2\lambda_c} \right) + \frac{2a\lambda_c}{\pi^2 s} \sum_{n=1}^{\infty} (-1)^{n+1} \frac{1 - \cos(n\pi s/\lambda_c)}{n^2} \cos\left(\frac{2\pi n x}{\lambda_c}\right)$$

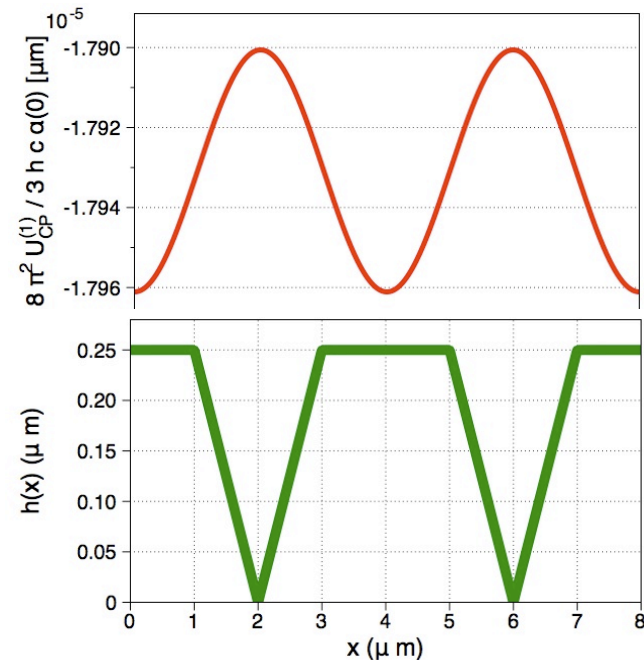


Single-atom lateral CP energy: it can be easily calculated using that the first order lateral CP energy  $U_{CP}^{(1)}(\mathbf{R}_A) = \int \frac{d^2\mathbf{k}}{(2\pi)^2} e^{i\mathbf{k}\cdot\mathbf{r}_A} g(\mathbf{k}, z_A) H(\mathbf{k})$  is linear in  $H(\mathbf{k})$

$$k_c z_A = 0.3$$



$$k_c z_A = 10$$

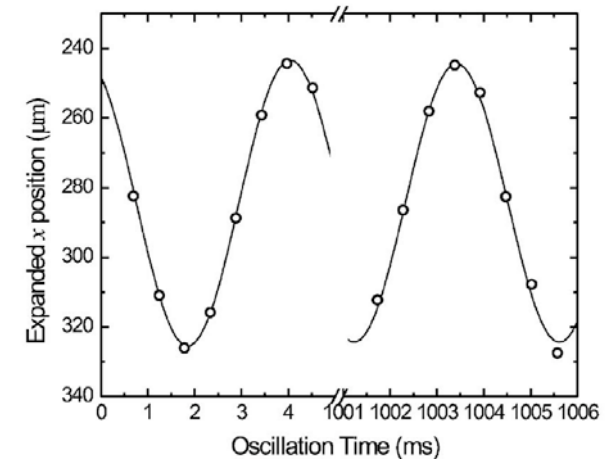


# BEC as a field sensor

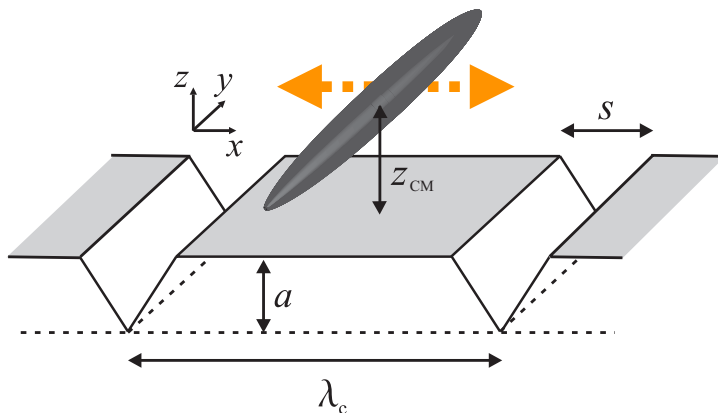
## ■ BEC oscillator

● The normal component of Casimir-Polder force  $U_{CP}^{(0)}(z)$  shifts the **normal dipolar oscillation frequency** of a BEC trapped above a surface

Antezza et al (2004) Cornell et al (2005, 2007)



● In order to measure the lateral component  $U_{CP}^{(1)}(x, z)$ , a cigar-shaped BEC could be trapped parallel to the corrugation lines, and the **lateral dipolar oscillation** measured as a function of time



$$V(\mathbf{r}) = V_{ho}(\mathbf{r}) + U_{CP}(\mathbf{r})$$

$$V_{ho}(\mathbf{r}) = \frac{m}{2}(\omega_x^2 x^2 + \omega_y^2 y^2 + \omega_z^2 z^2) \quad \omega_y \ll \omega_x = \omega_z$$

**Lateral frequency shift:**

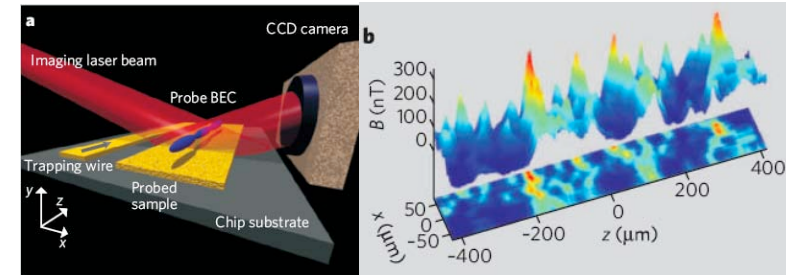
$$\omega_{x,CM}^2 = \omega_x^2 + \frac{1}{m} \int dx dz n_0(x, z) \frac{\partial^2}{\partial x^2} U_{CP}^{(1)}(x, z)$$

# BEC as a field sensor (cont'd)

## Density variations of a BEC above an atom chip

For a quasi one-dimensional BEC, the potential is related to the 1D density profile as

$$V_{\text{ho}}(x) + U_{\text{CP}}(x) = -\hbar\omega_x \sqrt{1 + 4a_{\text{scat}}n_{1d}(x)}$$



Measurement of the magnetic field variations along a current-carrying wire

Schmiedmayer et al (2005)



# BEC as a field sensor (cont'd)

## Density variations of a BEC above an atom chip

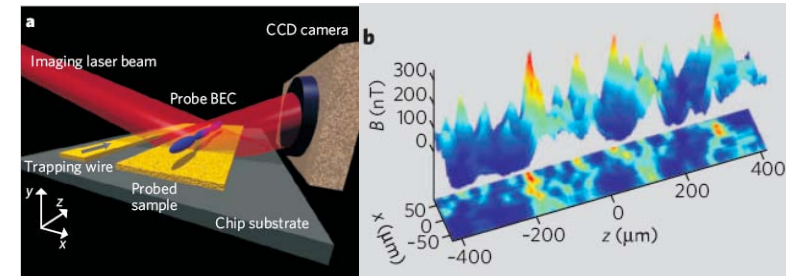
- For a quasi one-dimensional BEC, the potential is related to the 1D density profile as

$$V_{\text{ho}}(x) + U_{\text{CP}}(x) = -\hbar\omega_x \sqrt{1 + 4a_{\text{scat}}n_{1d}(x)}$$

For the lateral CP force, perfect conductor, sinusoidal corrugation ( $a = 100\text{nm}$ ), distance  $z_A = 2\mu\text{m}$ , PFA limit ( $k_c z_A \ll 1$ )

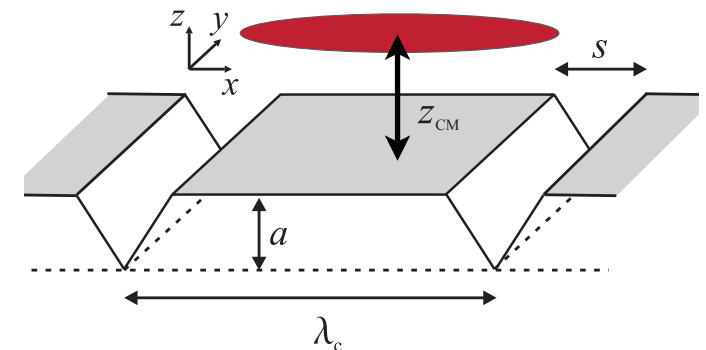
$$\Delta U_{\text{CP}}^{(1)} \simeq 10^{-14} \text{ eV}$$

- To measure the lateral CP force, the elongated BEC should be aligned along the x-direction, and a **density modulation** along this direction **above the plateau** would be a signature of a nontrivial (beyond-PFA) geometry effect.



Measurement of the magnetic field variations along a current-carrying wire

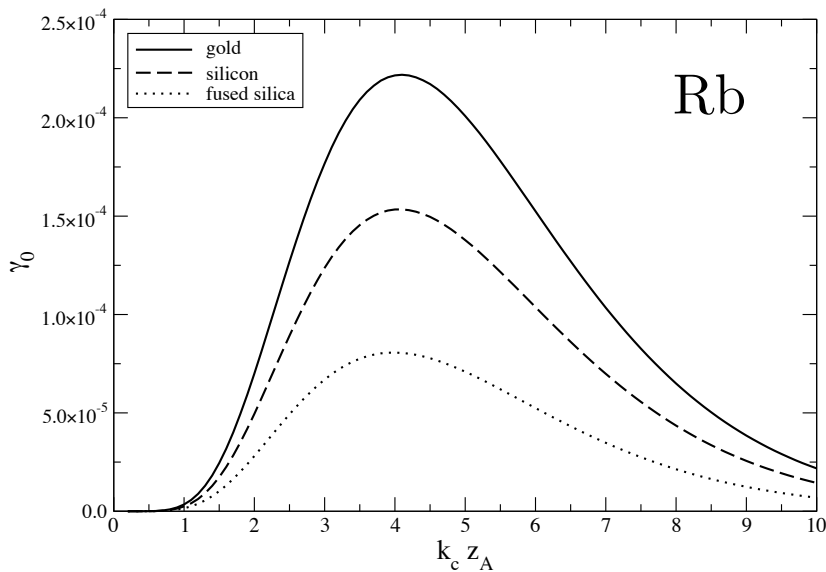
Schmiedmayer et al (2005)



# Single-atom/BEC frequency shift

$$\gamma_0 \equiv \frac{\omega_{x,CM} - \omega_x}{\omega_x}$$

## Single-atom lateral freq. shift



$$\omega_x/2\pi = 229 \text{ Hz}$$

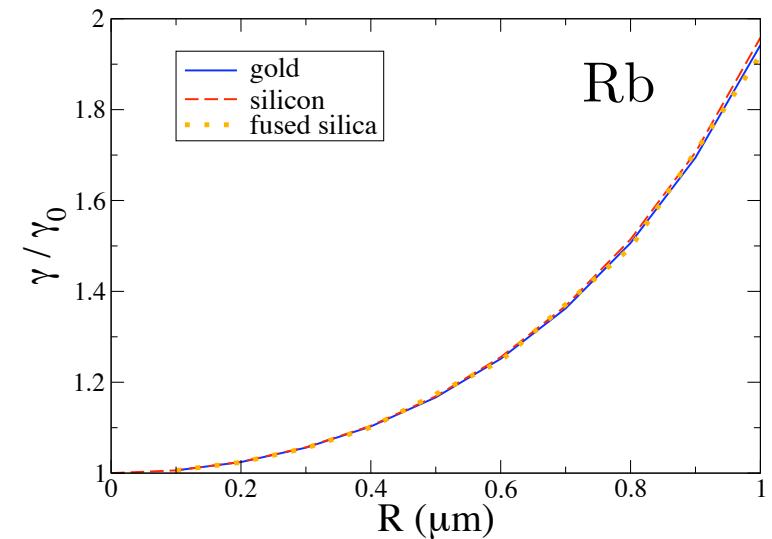
$$z_{CM} = 2 \mu\text{m}$$

$$\lambda_c = 4 \mu\text{m}$$

$$a = 250 \text{ nm}$$

$$s = \lambda_c/2$$

## Single-atom / BEC comparison



Given the **reported sensitivity**  $\gamma = 10^{-5} - 10^{-4}$  for relative frequency shifts from E. Cornell's experiment, we expect that **beyond-PFA lateral CP forces on a BEC above a plateau of a periodically grooved silicon surface should be detectable** for distances  $z_{CM} < 3 \mu\text{m}$ , groove period  $\lambda_c = 4 \mu\text{m}$ , groove amplitude  $a = 250 \text{ nm}$ , and a BEC radius of, say,  $R \approx 1 \mu\text{m}$

# Towards an experiment

Surfaces are being fabricated by Matt Blain



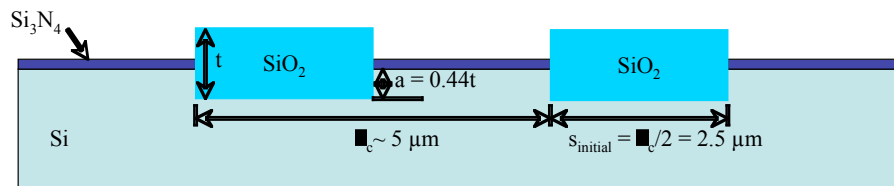
CP force measurements with BEC will be done by Malcolm Boshier



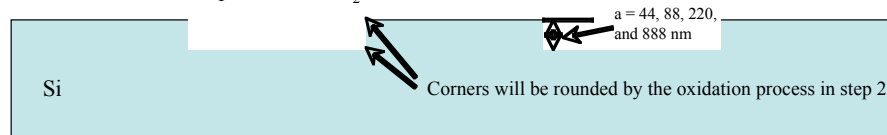
Number of periods, 100/set  
 • Length of “grooves”, 2 mm  
 • see next slide

**Process sequence**

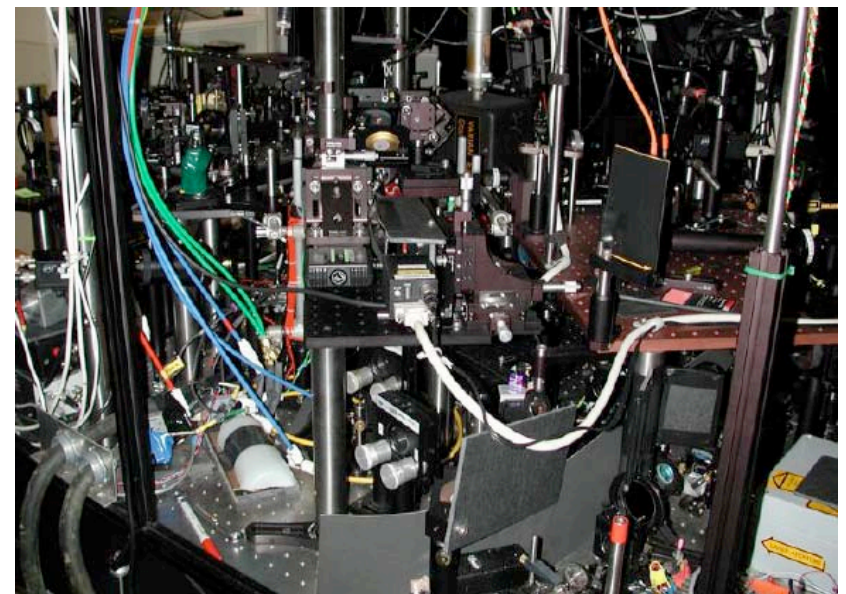
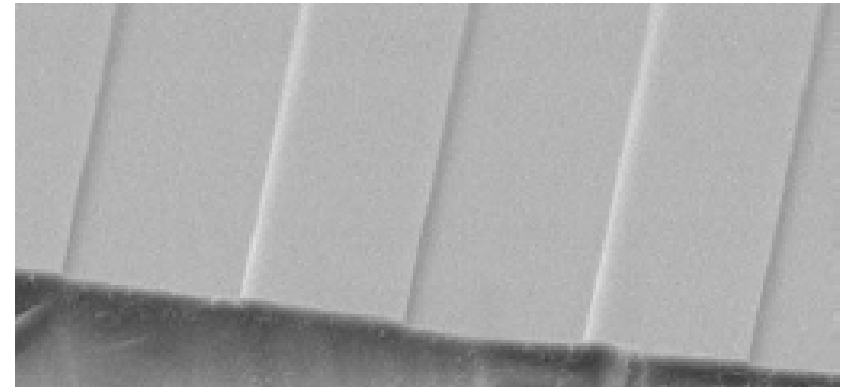
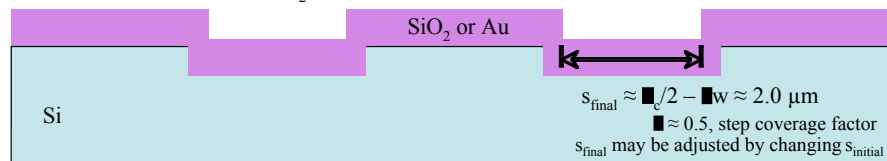
1. Deposit/pattern SiN oxidation mask
2. Grow SiO<sub>2</sub> in exposed Si to thicknesses of  $t = 100, 200, 500$  and  $2000$  nm



3. Strip SiN and SiO<sub>2</sub>

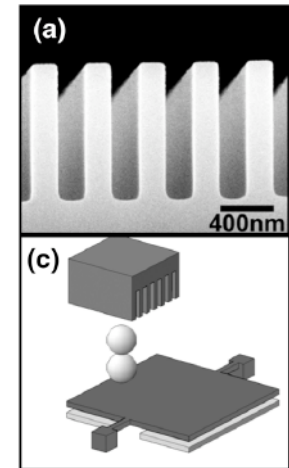


4. Deposit SiO<sub>2</sub> or Au to a thickness of  $w = 1 \mu\text{m}$



# Exact analytical results

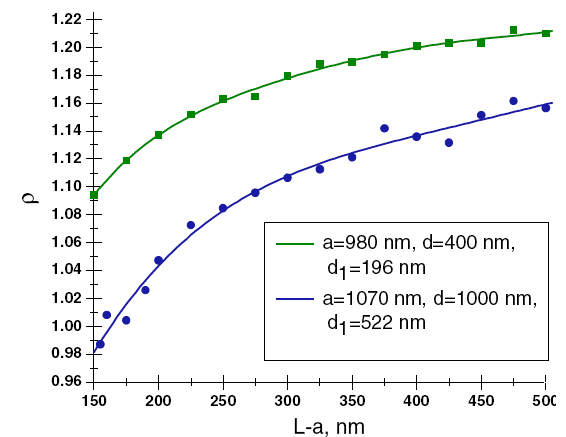
■ Large corrections to PFA/PWS have been recently observed in the Casimir force between a Au sphere and a Si uniaxial corrugated surface Chan et al (2008)



■ Exact analytical expressions for the Casimir force for uniaxial corrugations have been derived:

● Ideal reflectors Buscher+Emig (2004)

● Real materials Lambrecht+Marachevsky (2008)



■ We have already computed the exact CP force for an atom above uniaxial corrugated surfaces taking into account optical data of the materials Dalvit, Maia Neto, Lambrecht, Marachevsky (to be submitted)

# Summary part III

- Novel cold atoms techniques open a promising way of investigating nontrivial geometrical effects on quantum vacuum
- Important feature of atoms: they can be used as local probes of quantum vacuum fluctuations
- Non-trivial, beyond-PFA/PWS effects should be measurable using a BEC as a vacuum field sensor with available technology

For more details see:

Dalvit, Maia Neto, Lambrecht, and Reynaud,  
Phys. Rev. Lett. 100, 040405 (2008)  
J. Phys.A 41, 164028 (2008)