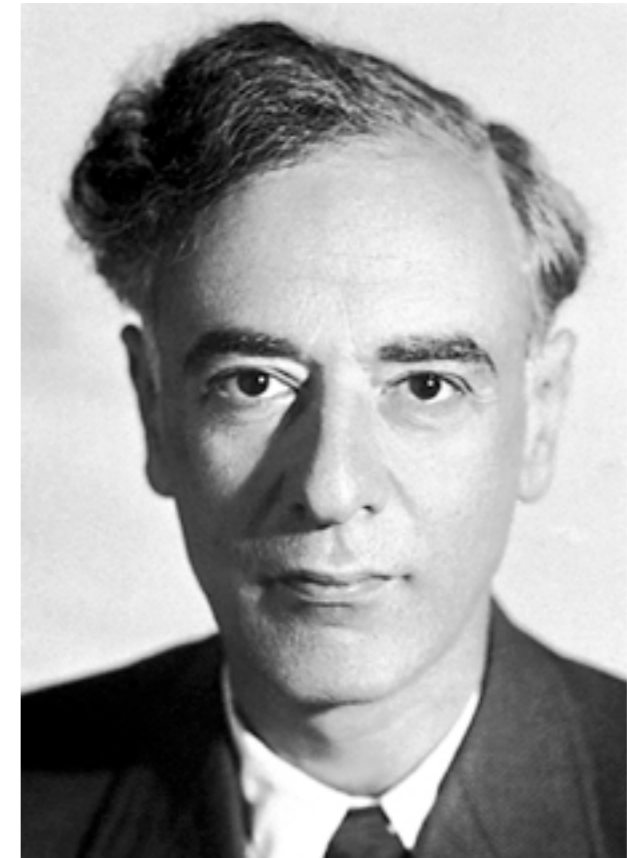


Quantum Field Theory and Condensed Matter Physics: making the vacuum concrete

Fabian Essler (Oxford)

Oxford, June 2013

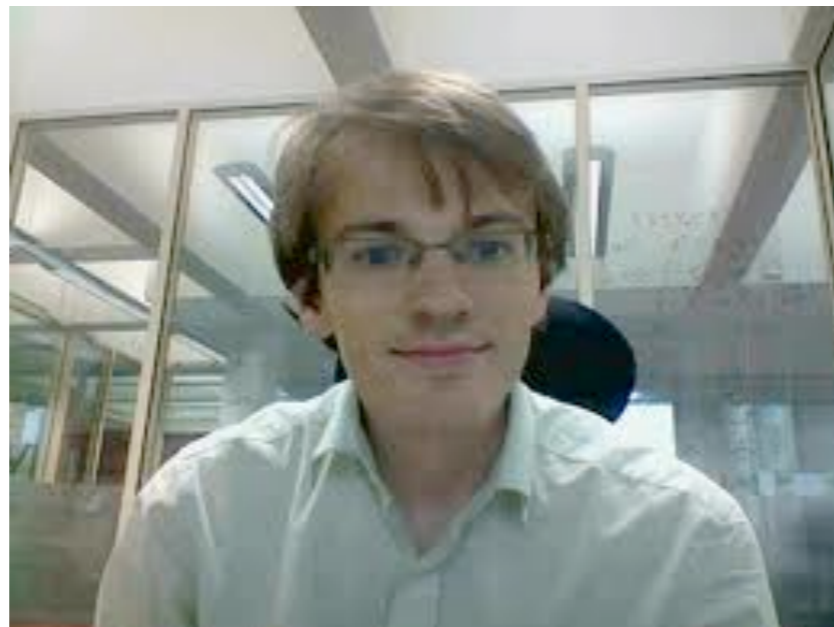


Lev Landau

“This work contains many things which are new and interesting. Unfortunately, everything that is new is not interesting, and everything which is interesting, is not new.”



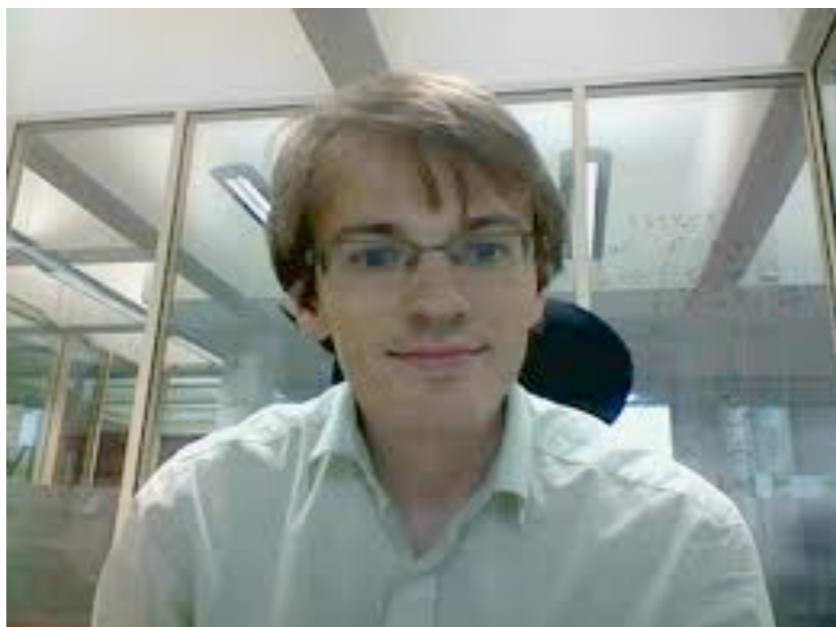
The Field
is real!



$\hbar > 0!$



The Field
is real!



$\hbar > 0!$



Here:

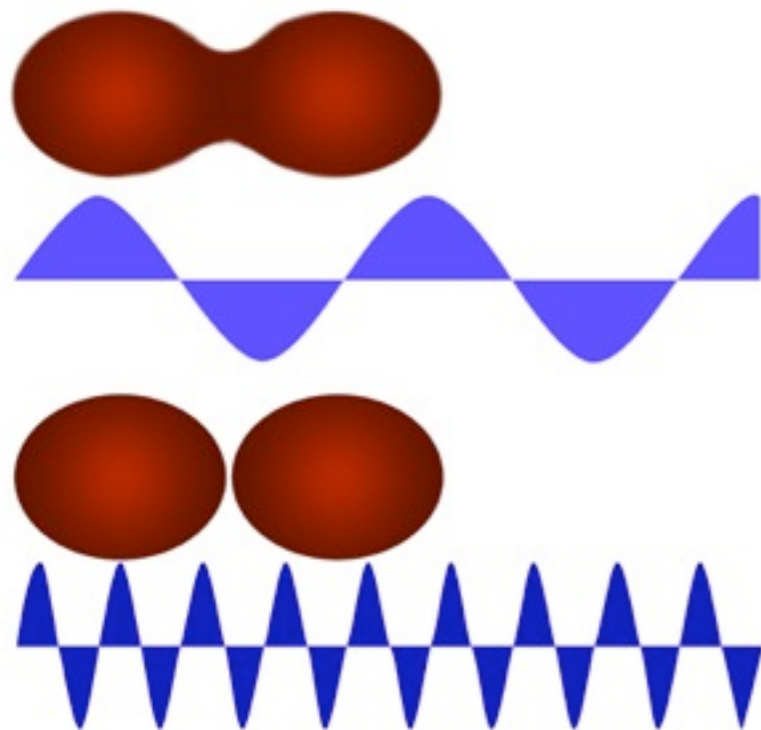
Relativistic QFT as an emergent
(rather than fundamental) phenomenon.

Length/Energy Scales: Organizing Principle of Physics

high energies \Leftrightarrow short distances

low energies \Leftrightarrow large distances

Think of this in terms of probing a physical system by e.g. light:



to resolve what happens at length scale a ,
the wavelength λ must be smaller than a .

short distances \rightarrow small $\lambda \rightarrow$ high energies as $E=hc/\lambda$

Length



← **Classical Mechanics**

← **Quantum Mechanics**

← **Standard Model**

← **ToE?**

A Reductionist View of the World

Quantum Field Theory



Quantum Mechanics



Chemistry



“If we understand
QFT/ToE, we understand
everything.”

A Reductionist View of the World

Quantum Field Theory



Quantum Mechanics



Chemistry



“If we understand
QFT/ToE, we understand
everything.”

Depends on what
“understand” means...

Following 30 minutes:

Quantum Field Theory



Quantum Mechanics



Quantum Field Theory

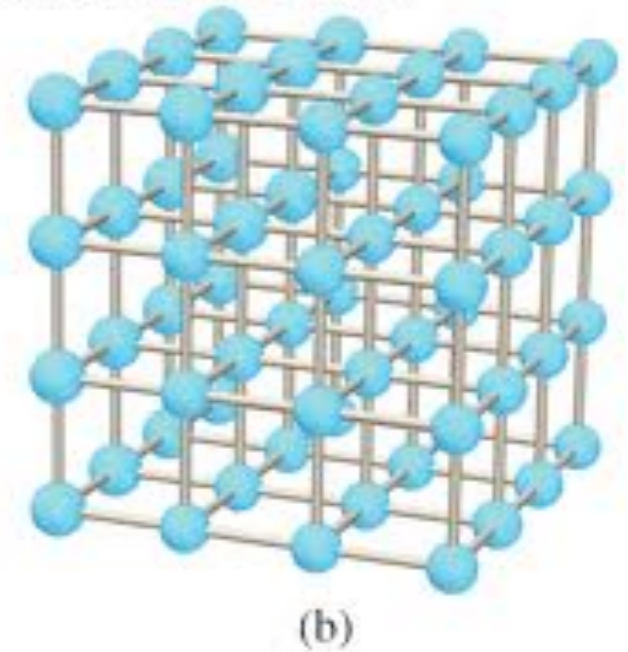


interesting perspective on QFT (and perhaps the reductionist view)

Example: Lattice Vibrations in Crystals

Crystal: atoms in a periodic array

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Interaction between the atoms (e.g. Coulomb) $V(\vec{R}_1, \vec{R}_2, \dots, \vec{R}_N)$

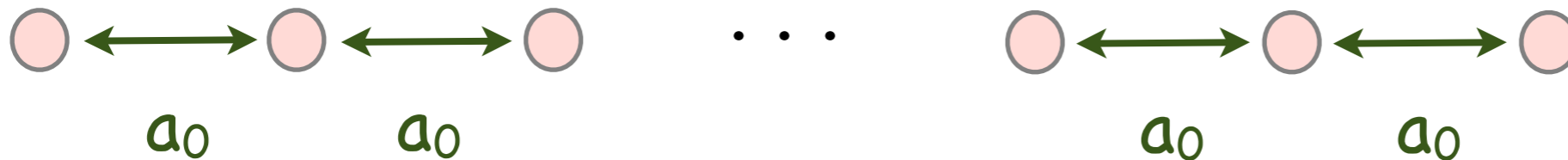
→ causes oscillations of atoms

Crystal ↔ mean ("equilibrium") positions

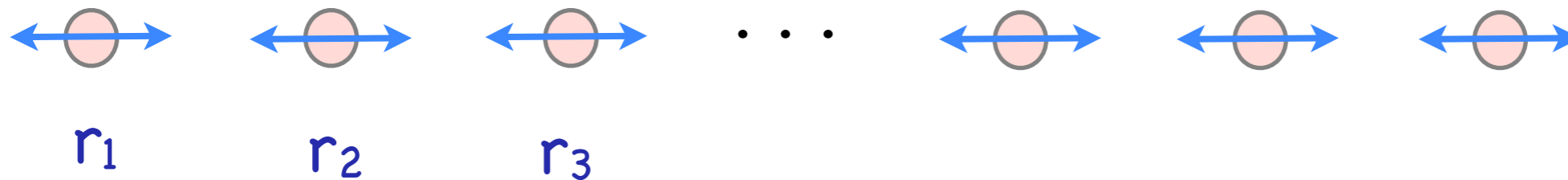
$$\vec{R}_1^{(0)}, \vec{R}_2^{(0)}, \dots$$

Lattice Vibrations of a Linear Chain

equilibrium positions: $R_j^{(0)} = ja_0$



oscillations: $r_j = R_j - R_j^{(0)}$



deviations from equilibrium positions

- Main interactions between nearest neighbours
- Oscillations around equilibrium positions typically **small**:

$$|r_j| \ll a_0$$

$$V(R_1, R_2, \dots, R_N) = V_0 + \frac{\kappa}{2} \sum_{j=1}^{N-1} (r_j - r_{j+1})^2 + \dots$$

- ◇ $V_0 = V(R_1^{(0)}, R_2^{(0)}, \dots, R_N^{(0)})$
- ◇ The linear terms add up to zero \leftrightarrow equilibrium positions
- ◇ ... \rightarrow small cubic, quartic etc "anharmonic" terms

$$H = \underbrace{\sum_{l=1}^N \frac{p_l^2}{2m}}_{\text{Kinetic energy}} + \frac{\kappa}{2} \underbrace{\sum_{l=1}^{N-1} (r_l - r_{l+1})^2}_{\text{Potential energy}} + \dots$$

Kinetic energy

Potential energy

→ N coupled harmonic oscillators!



Quantum mechanically:

$$p_l = -i\hbar \frac{\partial}{\partial r_l}$$

Solution of the Classical Problem

Newton's
equations:

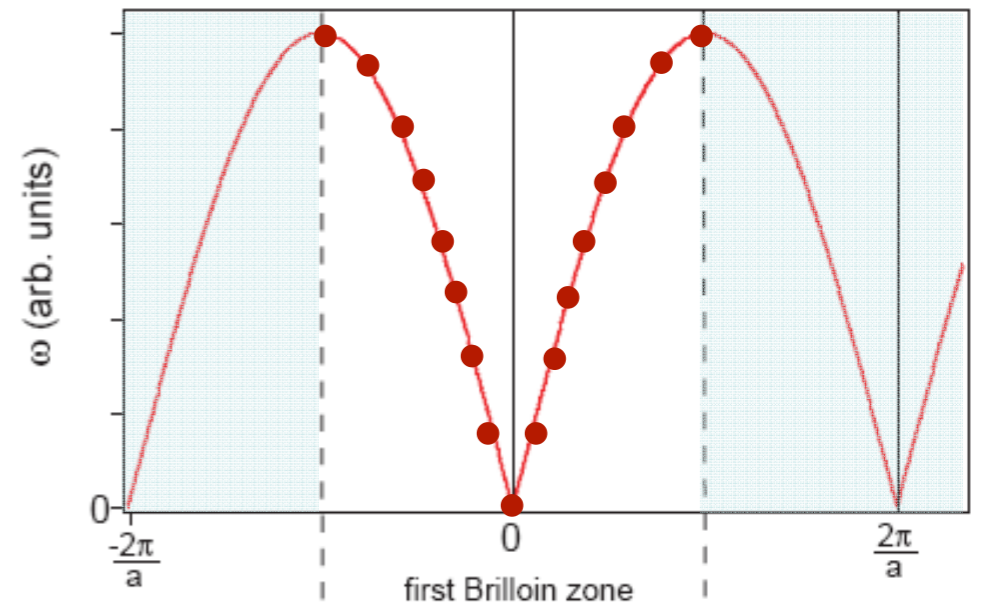
$$m \frac{\partial^2 r_l}{\partial t^2} = \kappa(r_{l+1} - r_l) - \kappa(r_l - r_{l-1}) \quad l = 1, 2, \dots, N$$

Ansatz:

$$r_l(t) = A \cos(kla_0 - \omega t + \delta)$$

Works if

$$\omega(k_l) = \sqrt{\frac{2\kappa}{m} [1 - \cos(k_l a_0)]}$$



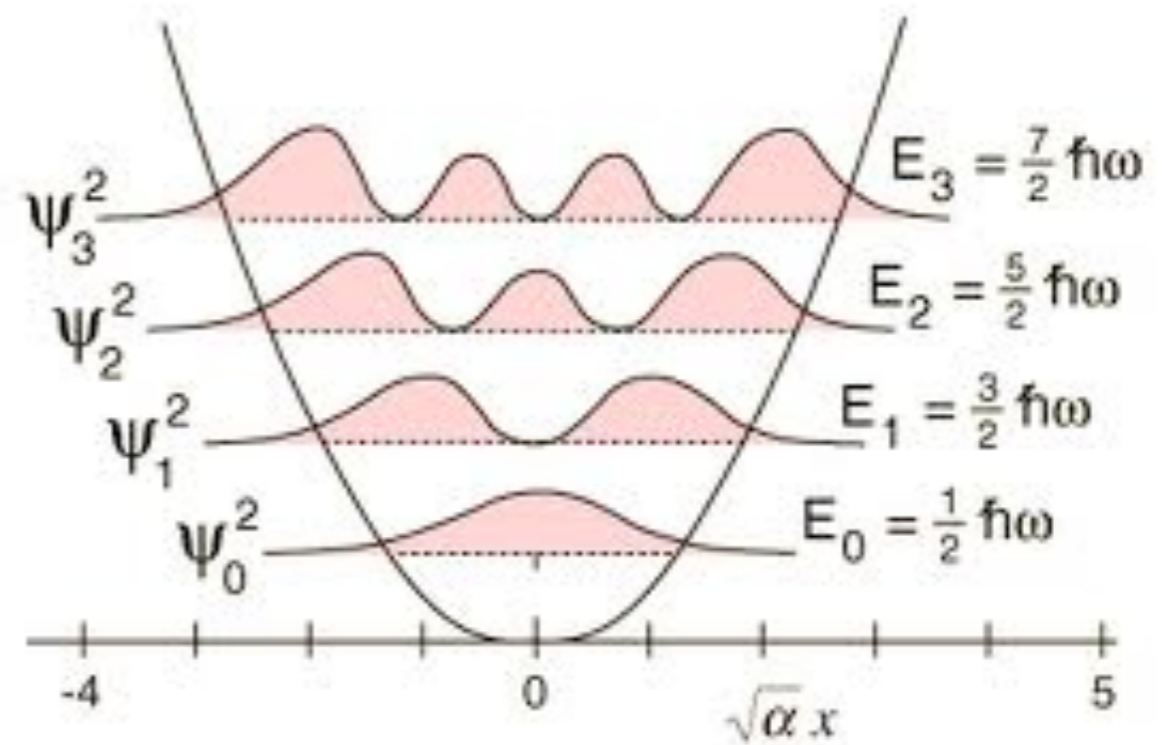
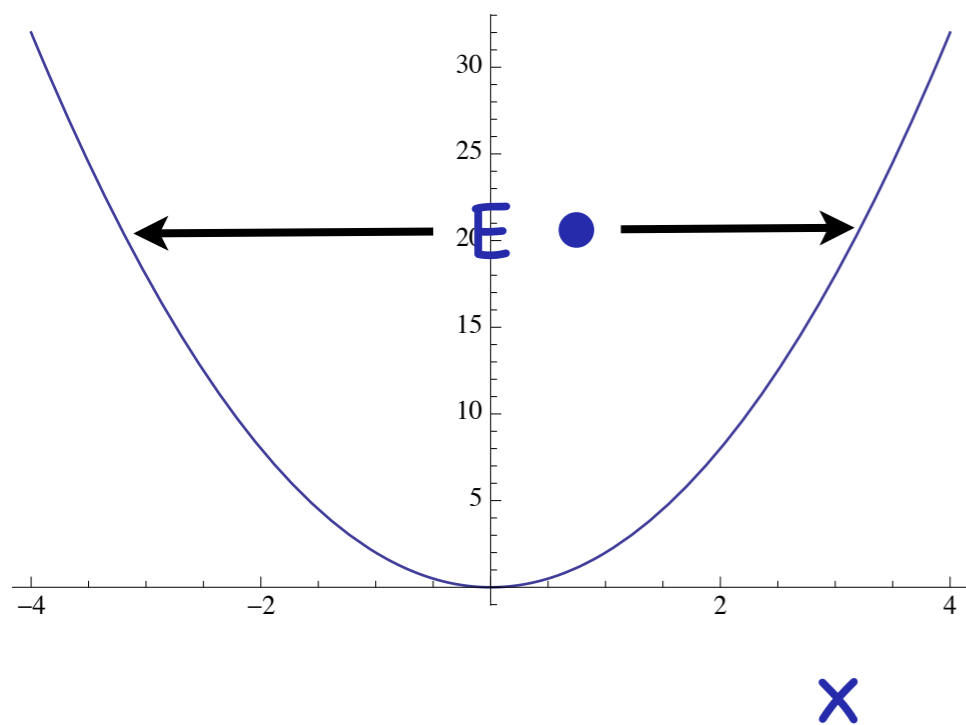
periodicity in $k \leftrightarrow$ periodicity of the lattice

finite number of atoms \rightarrow finite number of **normal modes**



Quantum Mechanics

Each $\omega(k)$ gives a simple harmonic oscillator
(in an appropriate coordinate) \rightarrow



quantized energies:

$$E_n(k_l) = \hbar\omega(k_l) \left[n + \frac{1}{2} \right], \quad l = 1, \dots, N$$

Ground State ("Vacuum"):

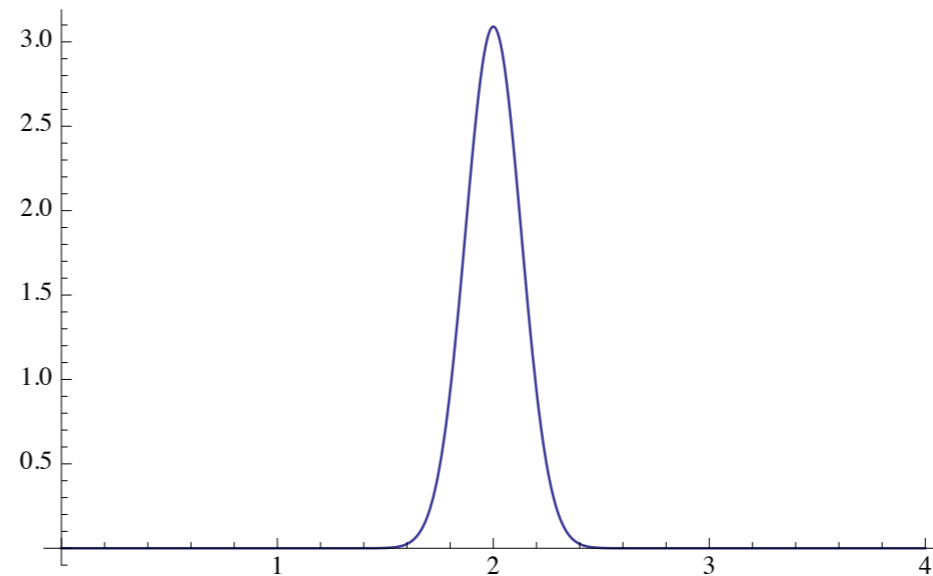
zero-point energy:

$$E_{\text{GS}} = \sum_{l=1}^N \frac{\hbar\omega(k_l)}{2}$$

wave function:

$$\Psi(r_1, \dots, r_N) \propto \exp\left(-\frac{1}{2} \sum_{j,k} r_j M_{jk} r_k\right)$$

Probability distr.
of 2nd atom



Excited states:

Single particles
("Phonons")

$$E = E_{\text{GS}} + \hbar\omega(k_l)$$

wave function:

$$\Psi(r_1, \dots, r_N) \propto \left[\sum_{n=1}^N \cos(k_l n) r_n \right] \Psi_{\text{GS}}$$

Two phonons:

$$E = E_{\text{GS}} + \hbar\omega(k_l) + \hbar\omega(k_n)$$

Quantum Field Theory

Recall that

$$H = \sum_{l=1}^N \frac{p_l^2}{2m} + \frac{\kappa}{2} \sum_{l=1}^{N-1} (r_l - r_{l+1})^2 + \dots$$

$$p_l = -i\hbar \frac{\partial}{\partial r_l}$$

Quantum Field Theory

Define

$$\Phi(la_0) = \frac{1}{\sqrt{a_0}} r_l$$

Hamiltonian becomes

$$H = a_0 \sum_{l=1}^N \frac{m}{2} \left[\frac{\partial}{\partial t} \Phi(la_0) \right]^2 + \frac{\kappa}{2} \left[\Phi([l+1]a_0) - \Phi(la_0) \right]^2$$

Now consider the limit

$$N \rightarrow \infty, \quad a_0 \rightarrow 0, \quad \kappa \rightarrow \infty$$

$$L = Na_0 \text{ and } \bar{\kappa} = \kappa a_0^2 \text{ fixed}$$

↑
volume

↑
vibrations
remain $\ll a_0$

$$\begin{aligned} la_0 &\longrightarrow x \\ \Phi(la_0) &\longrightarrow \Phi(x) \\ a_0 \sum_l &\longrightarrow \int_0^L dx \end{aligned}$$

corresponding
Lagrangian density

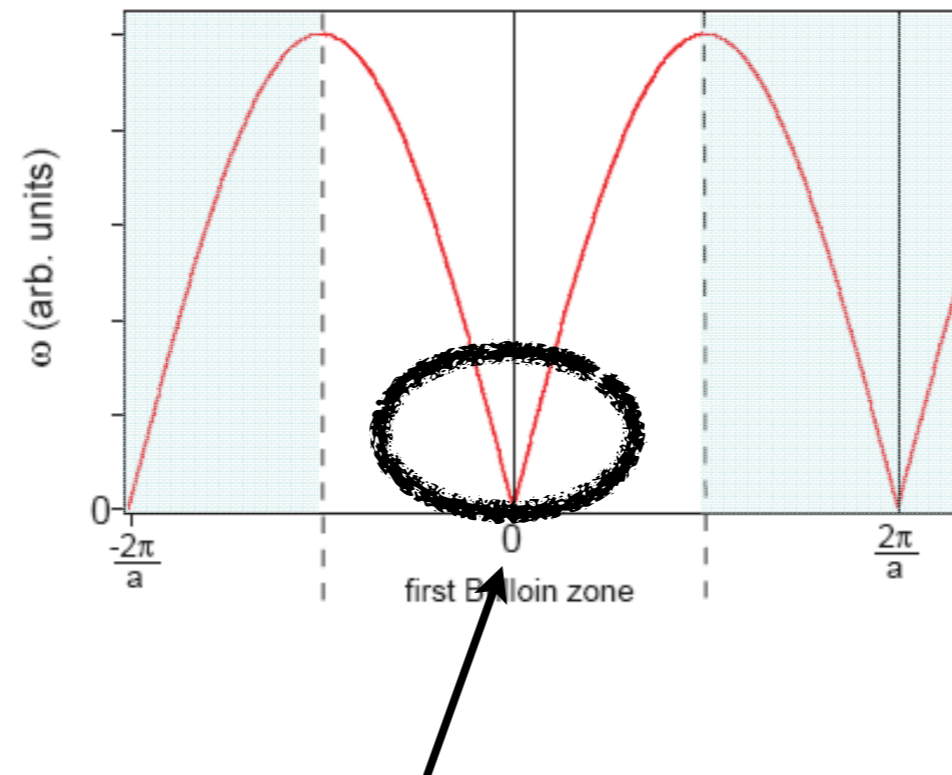
$$H \longrightarrow \int_0^L dx \left[\frac{m}{2} \left(\frac{\partial \Phi}{\partial t} \right)^2 + \frac{\bar{\kappa}}{2} \left(\frac{\partial \Phi}{\partial x} \right)^2 \right]$$

$$\mathcal{L}(t, x) = \left[\frac{m}{2} \left(\frac{\partial \Phi(t, x)}{\partial t} \right)^2 - \frac{\bar{\kappa}}{2} \left(\frac{\partial \Phi(t, x)}{\partial x} \right)^2 \right]$$

Massless (relativistic) Scalar Field

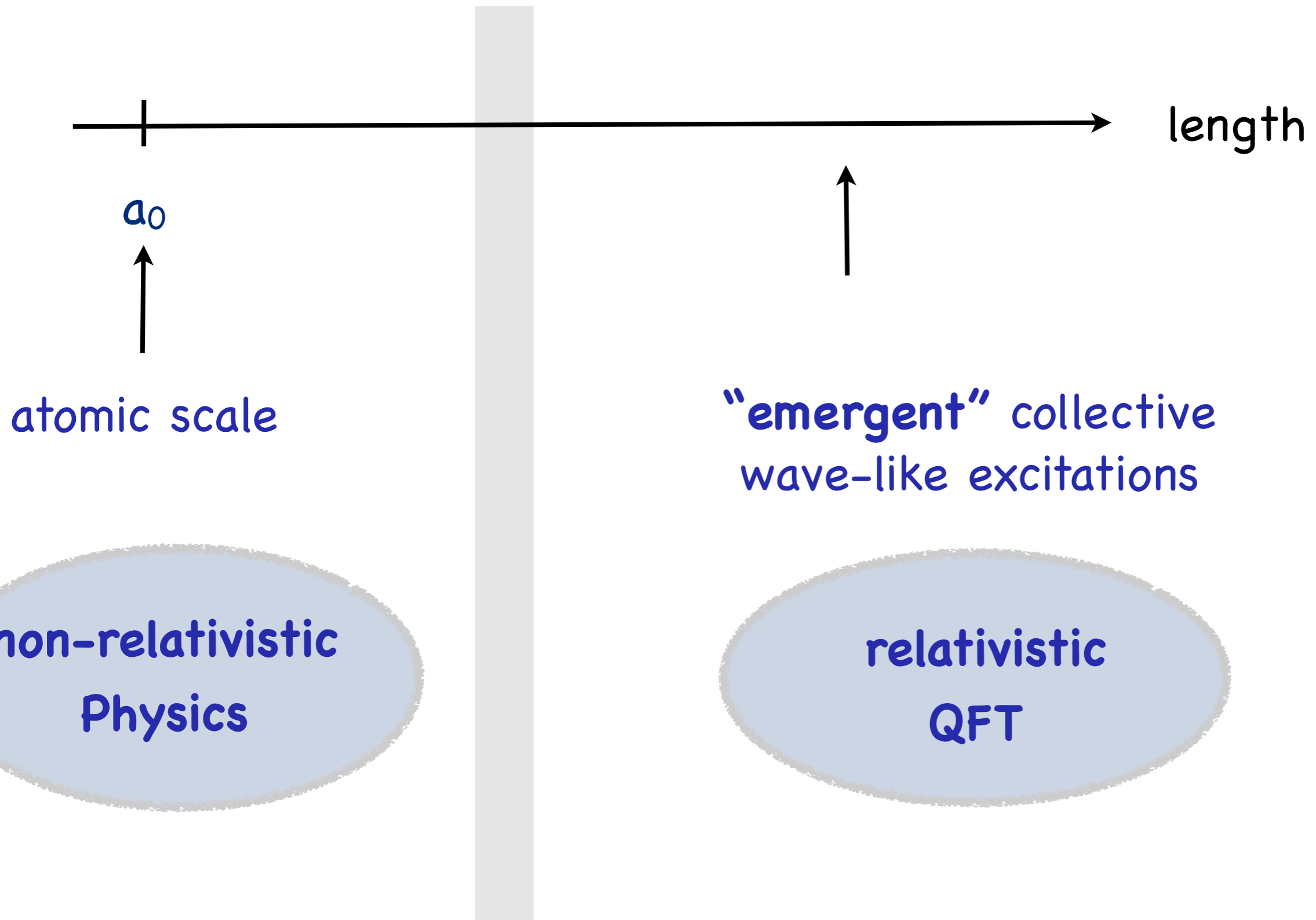
Which part of the physics does QFT describe?

- **Large distances** compared to “lattice spacing” a_0
- Small frequencies $\omega(k)$, i.e. **low energies**

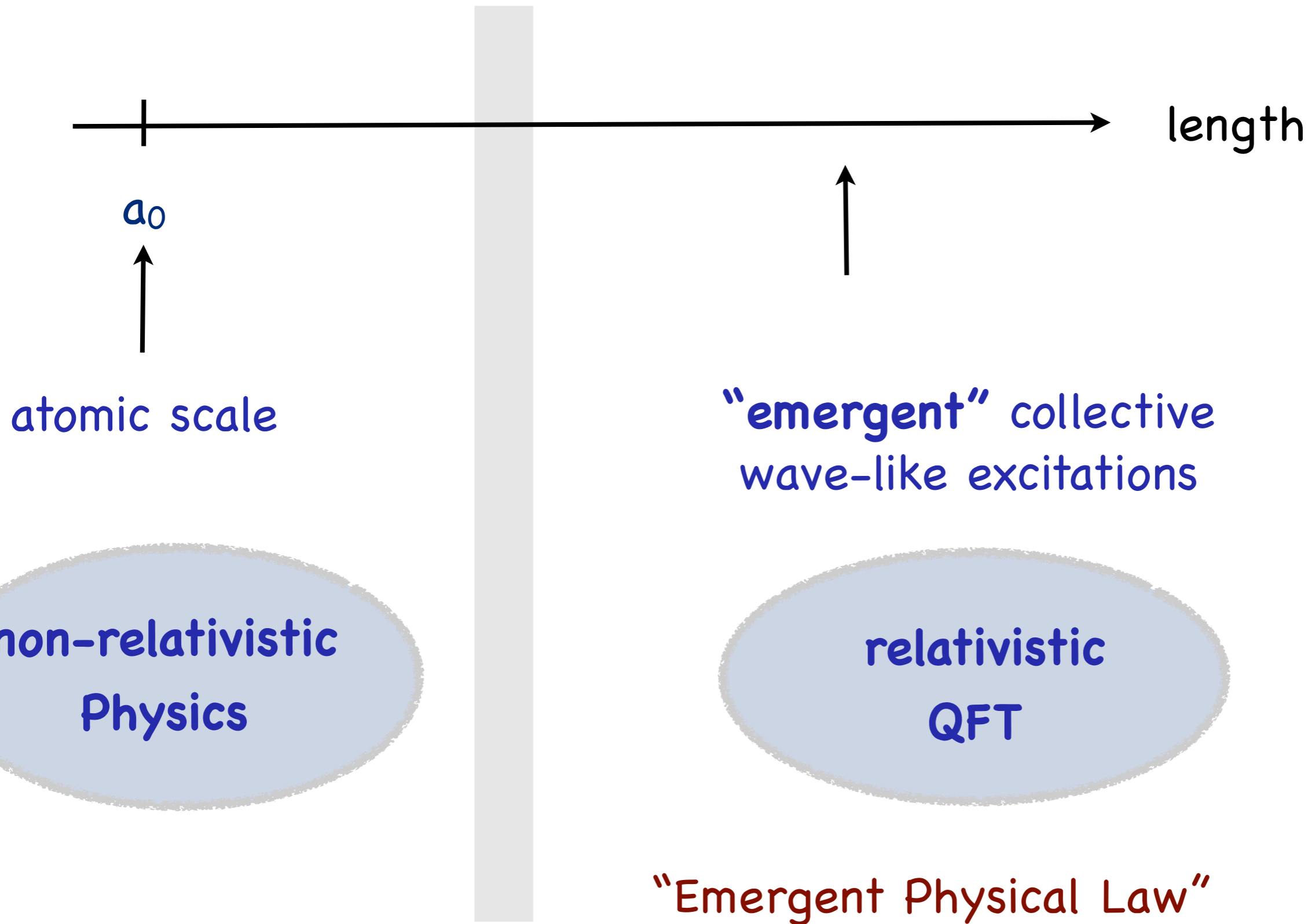


i.e. normal modes in **this** region

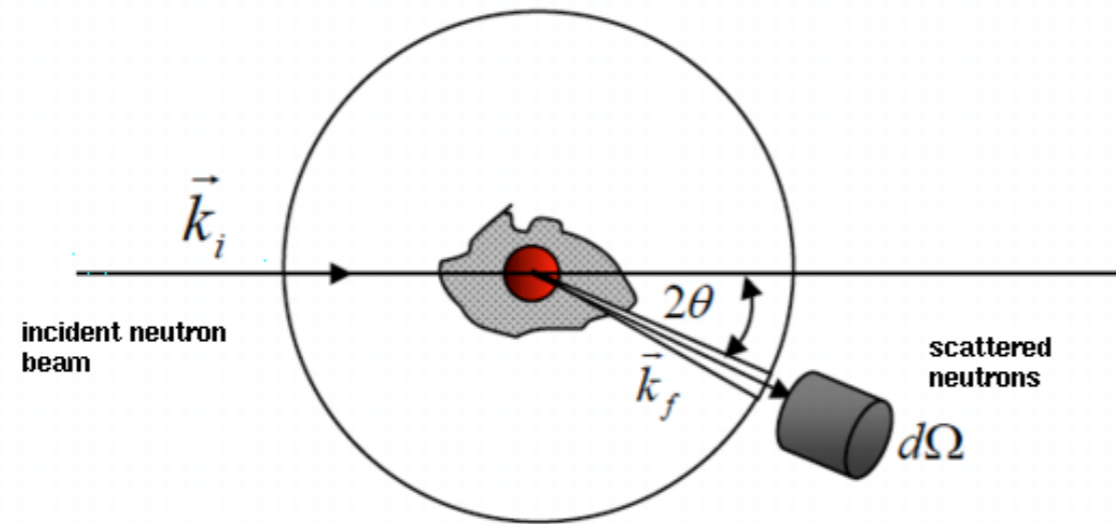
Which part of the physics does QFT describe?



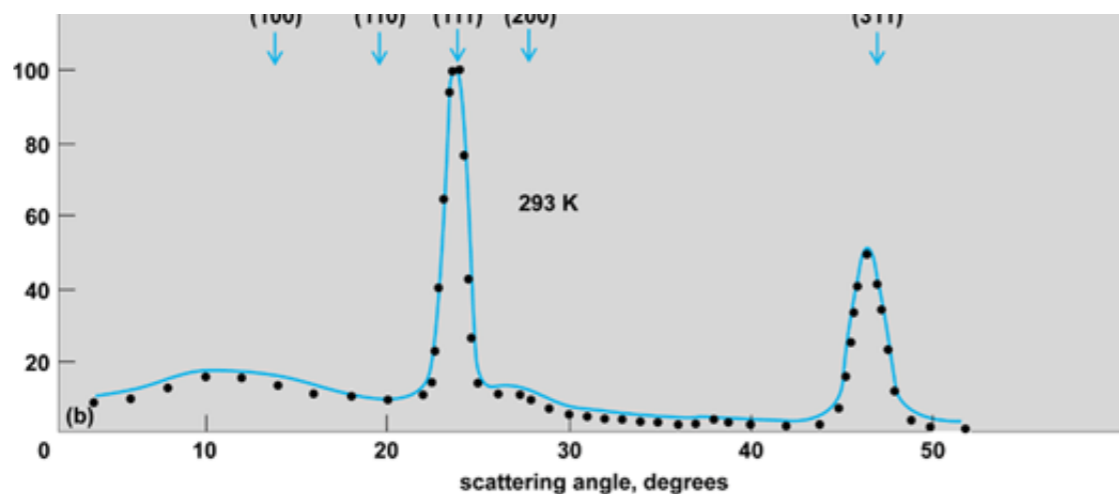
Which part of the physics does QFT describe?



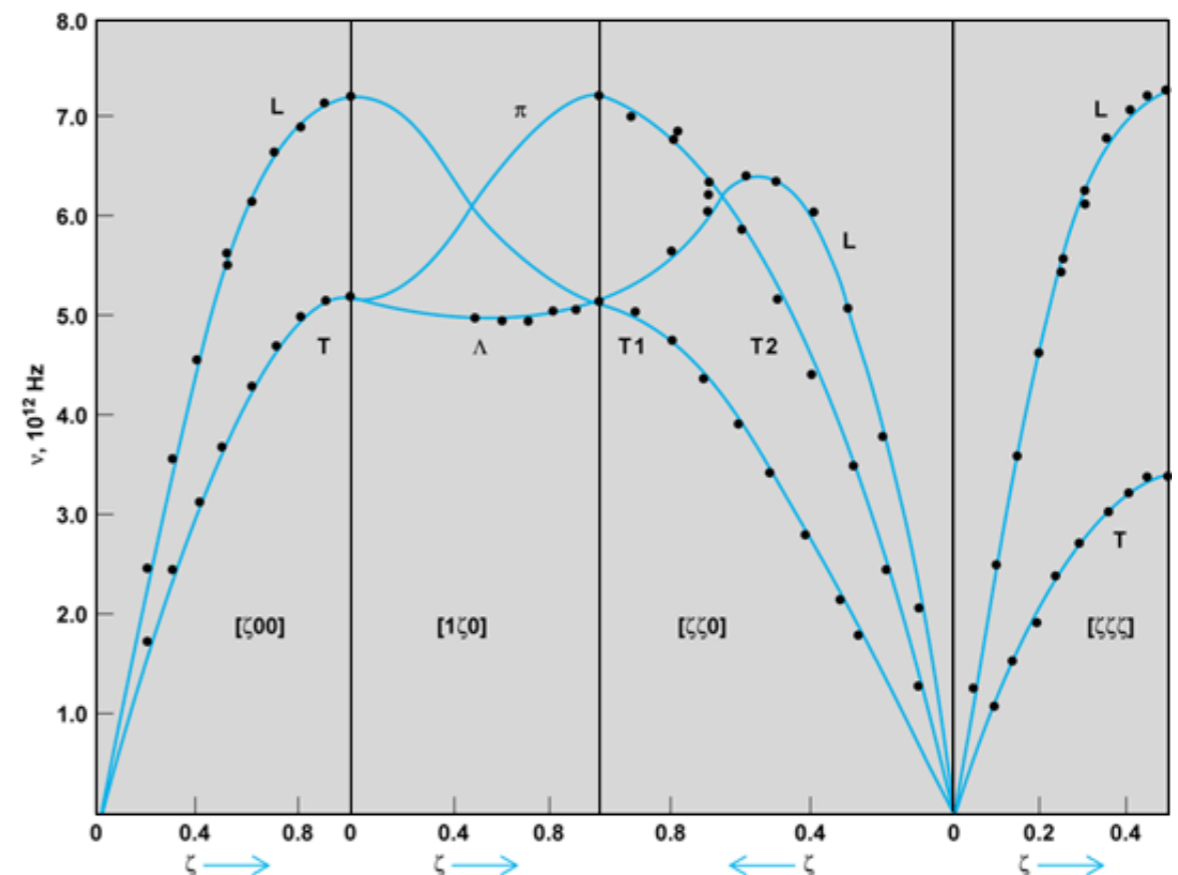
Measuring the collective modes: inelastic neutron scattering

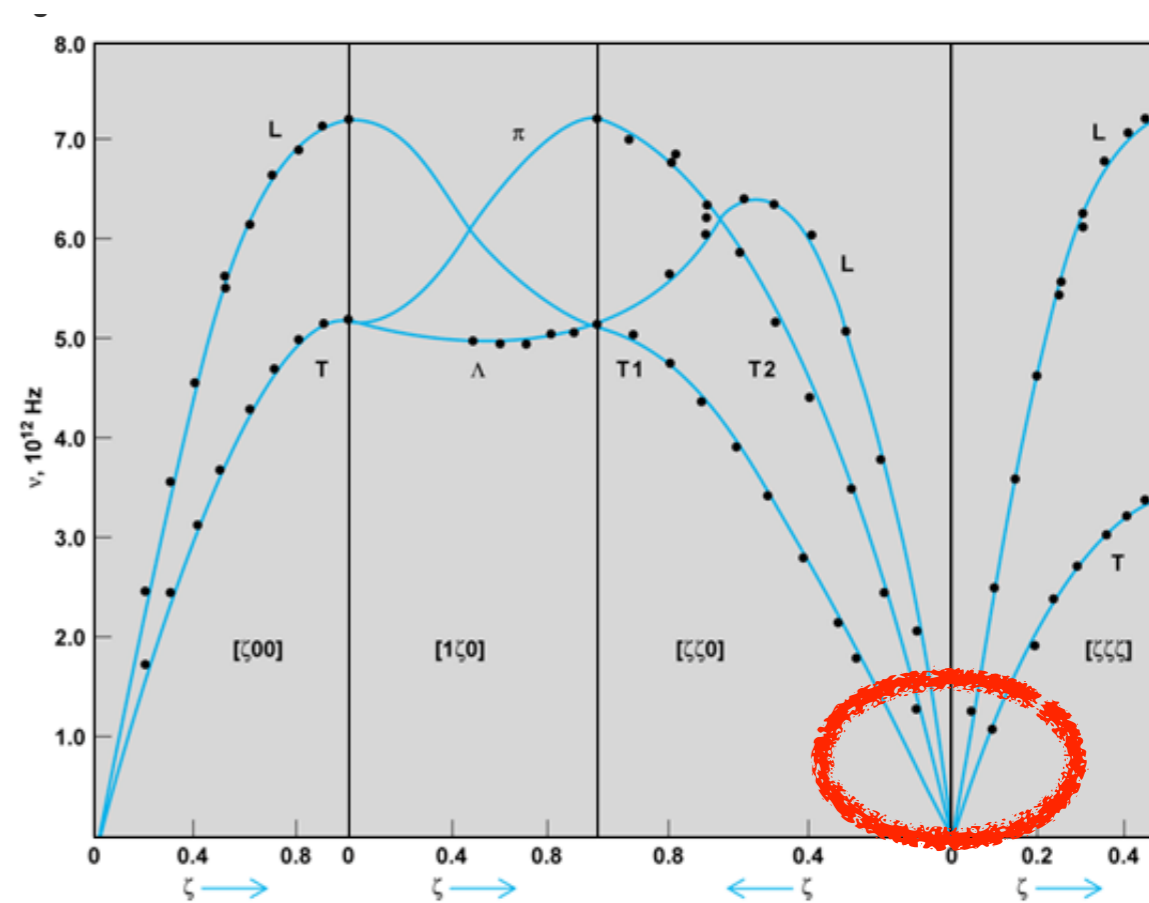
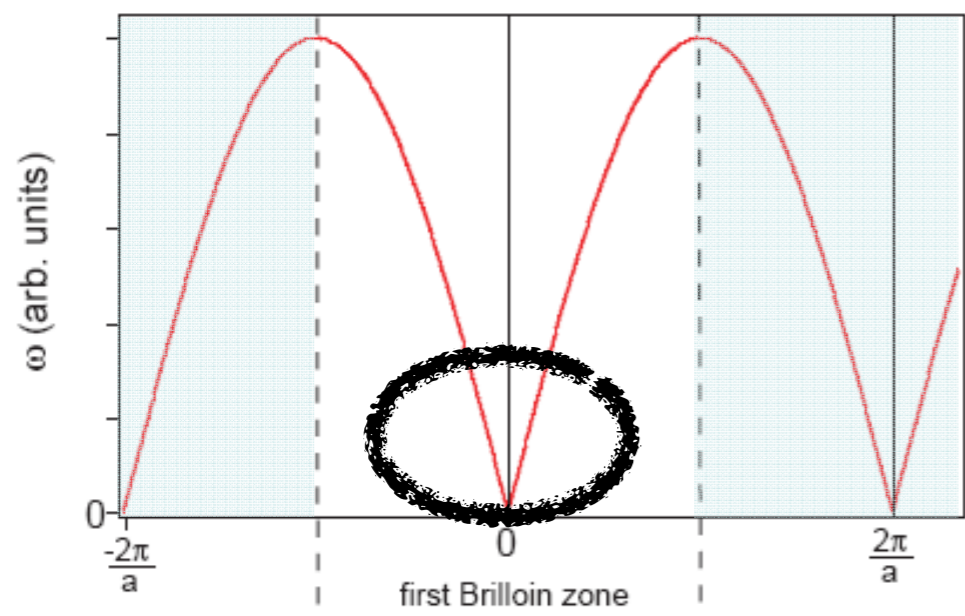


neutron gives energy ω and momentum k to the crystal
 → excites a **single** normal mode if ω and k “match”



Phonon spectrum of Cu





Main Points so far

- Quantum mechanics of atoms in solids can give rise to **collective excitations**, which are described by a QFT
- The lowest energy state (“vacuum”) has a non-trivial wave function.

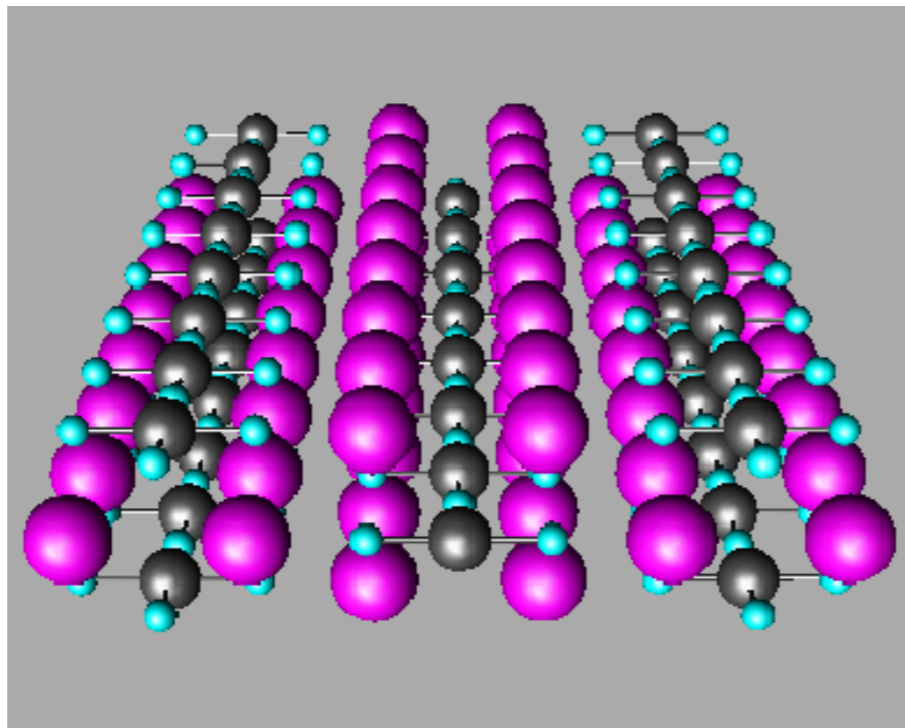
Main Points so far

- Quantum mechanics of atoms in solids can give rise to **collective excitations**, which are described by a QFT
- The lowest energy state (“vacuum”) has a non-trivial wave function.

Much more exotic physics and QFTs can arise in this way!

"Splitting the Electron"

Consider electronic degrees of freedom in "quasi-1D crystals"

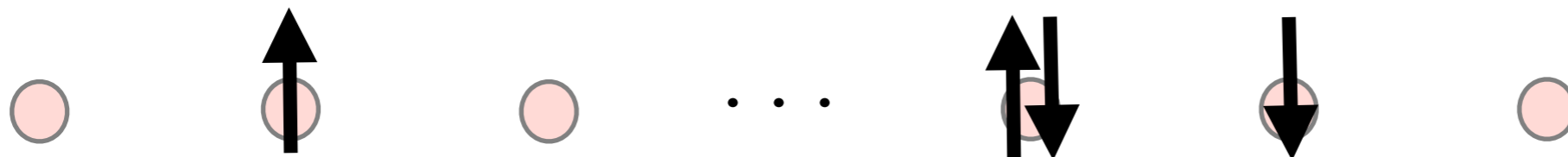


low energy electronic physics due to outer electrons on Cu atoms (black)

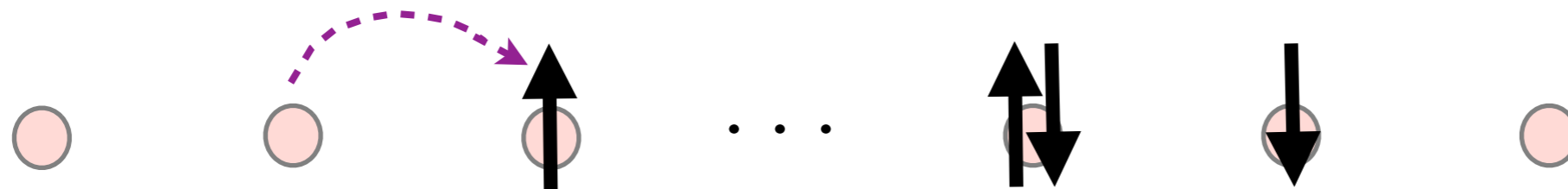
anisotropy \rightarrow e^- move essentially only along 1D chains

Basic model:

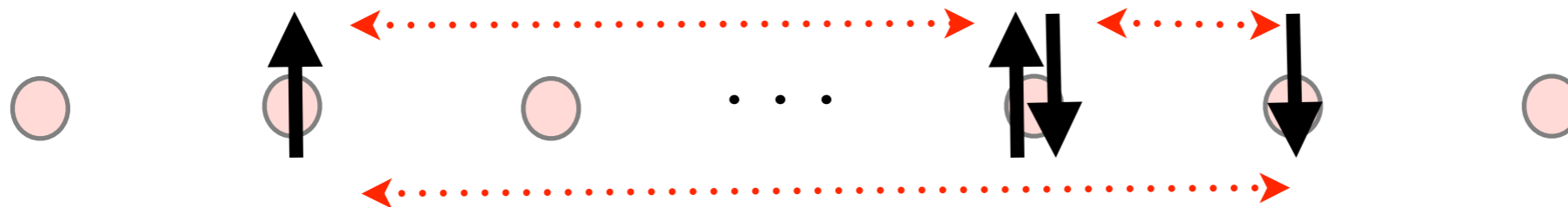
Lattice: on each site either 0, 1 or 2 electrons (spin!)



Electrons can hop to neighbouring sites



Electrons repel through Coulomb interaction



Field Theory Limit: description in terms of



$$\Psi_1(t, x) = \begin{pmatrix} R_{\uparrow}(t, x) \\ L_{\uparrow}(t, x) \end{pmatrix}, \quad \Psi_2(t, x) = \begin{pmatrix} R_{\downarrow}(t, x) \\ L_{\downarrow}(t, x) \end{pmatrix}$$

fermionic quantum fields



2 species of 2-dimensional Dirac spinors

$$\mathcal{L}(t, x) = \sum_{a=1}^2 \bar{\Psi}_a(t, x) \left[i\gamma^0 \frac{\partial}{\partial t} - i\gamma^1 \frac{\partial}{\partial x} \right] \Psi_a(t, x) + g \sum_{\alpha=1}^3 J_1^{\alpha}(t, x) J_1^{\alpha}(t, x) - J_0^{\alpha}(t, x) J_0^{\alpha}(t, x)$$

$$\bar{\Psi}(t, x) = \Psi^{\dagger}(t, x) \gamma^0$$

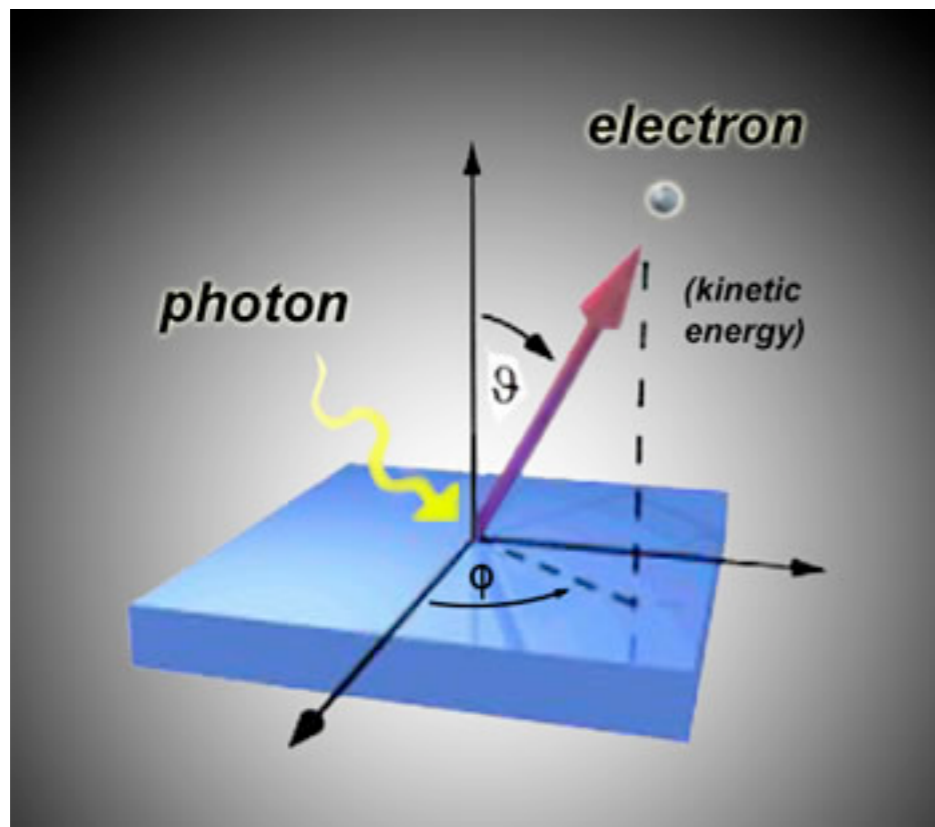
$$J_{\mu}^{\alpha} = \frac{1}{2} \bar{\Psi}_a \gamma_{\mu} \sigma_{ab}^{\alpha} \Psi_b, \quad \mu = 0, 1$$

$$\gamma^0 = -\gamma_0 = i\sigma^y = \begin{pmatrix} 0 & 1 \\ -1 & 0 \end{pmatrix}$$

$$\gamma^1 = \gamma_1 = \sigma^x = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}$$

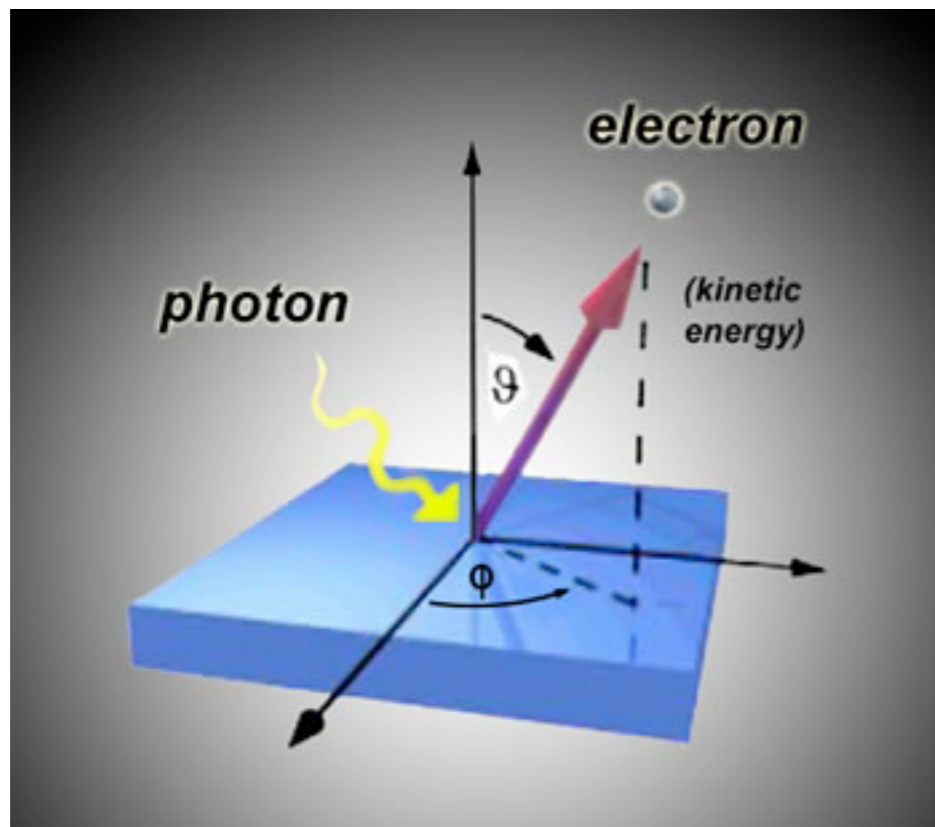
“SU(2) Thirring Model”

Measure excitations by scattering photons (Angle Resolved Photo Emission Spectroscopy)

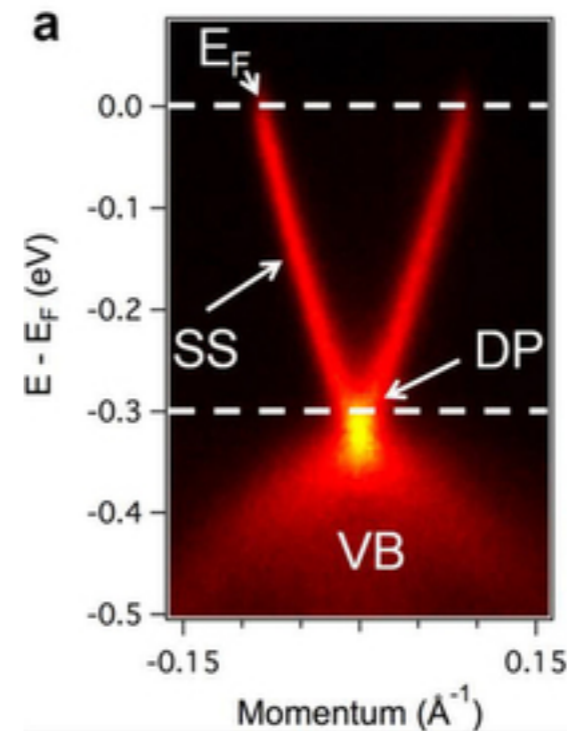


Count emitted electrons
with given energy and
momentum (\leftrightarrow angles)

Measure excitations by scattering photons (Angle Resolved Photo Emission Spectroscopy)



If the collective excitations are electrons, we expect to see something like this

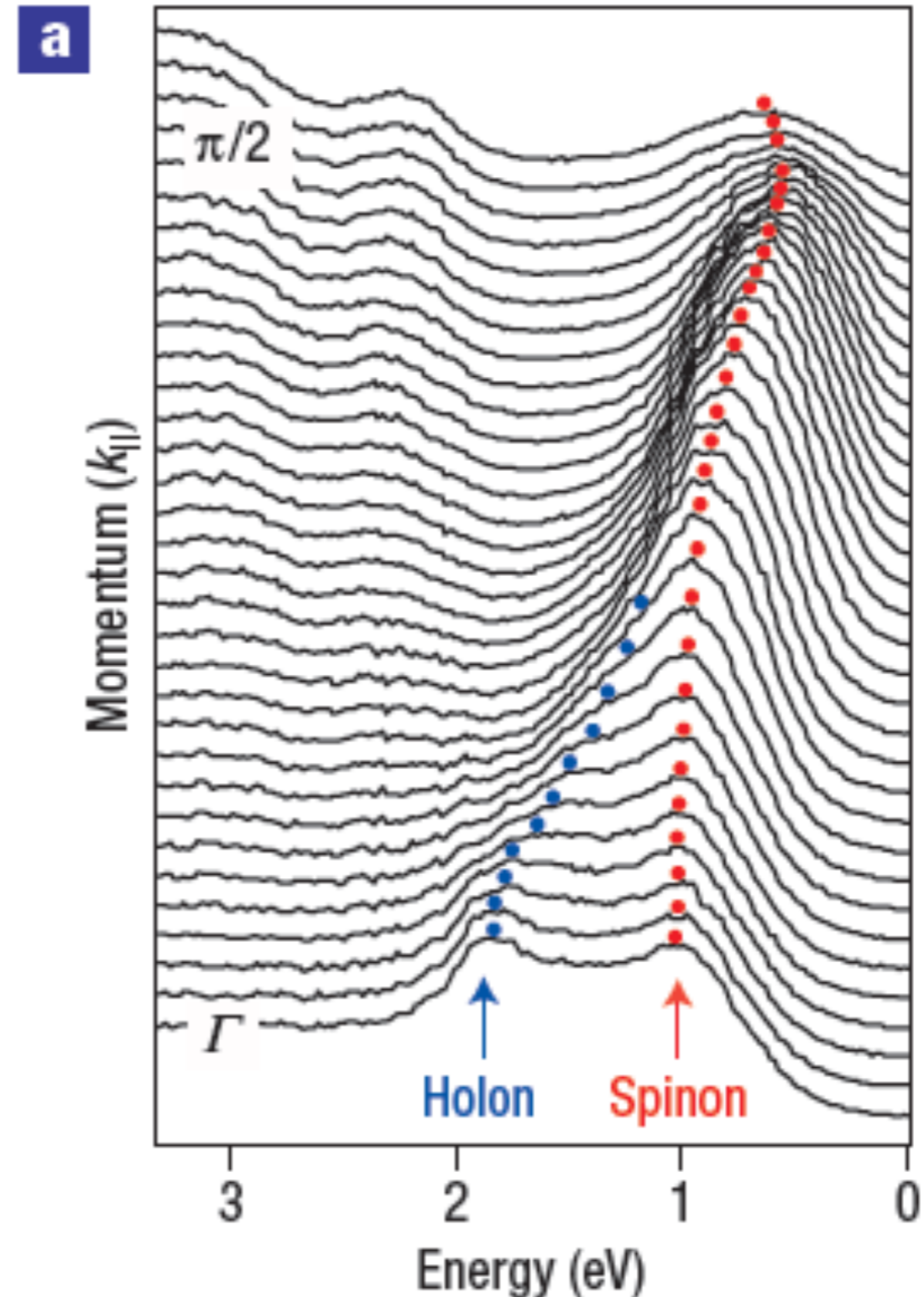


Spectral Function : ARPES on SrCuO₂

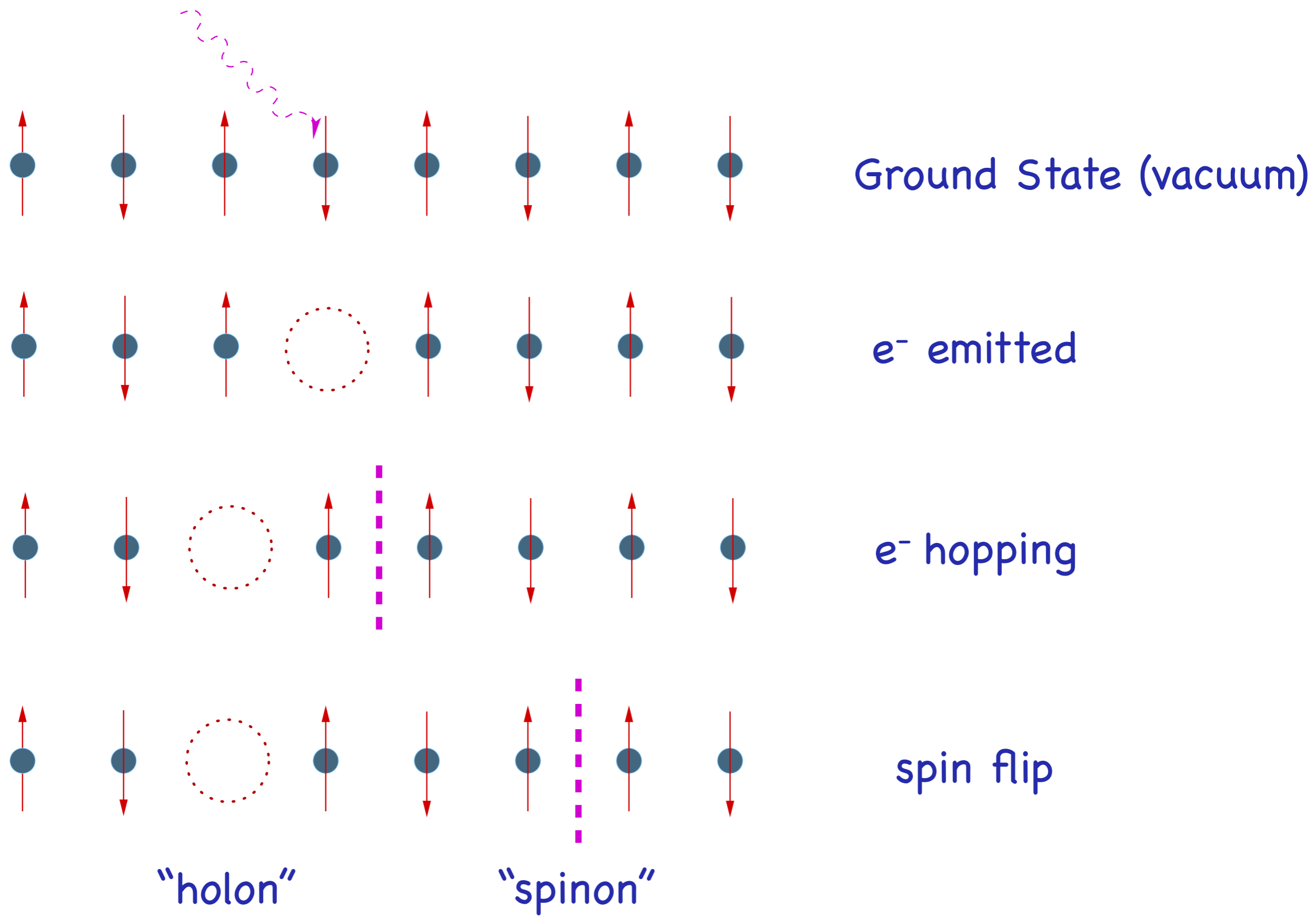
(Kim et al '06)

Instead, for quasi-1D systems
one observes

**The electron has fallen apart!
("spin-charge separation")**



How to understand this? In Sr_2CuO_3 we have 1 electron per site



The origin for this “splitting of the electron” is the highly non-trivial nature of the ground state (vacuum)!!

Summary

- Quantum Field Theories often describe **collective** properties of solids at large distances/low energies.
- Lorentz covariance can be an **emergent** feature in this regime.
- Low-energy physics can be very exotic because the “vacuum” is highly non-trivial.
- In the Cond. Mat. context it is possible to vary the spatial dimensionality $D=1,2,3$ (anisotropy!).