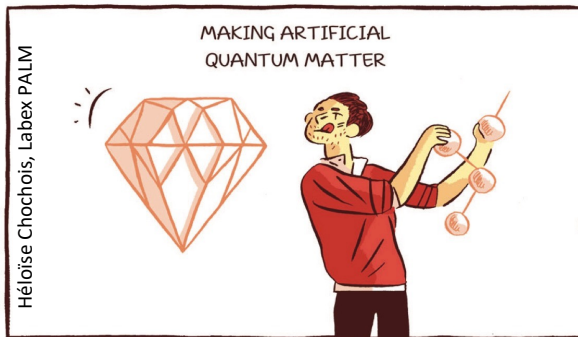


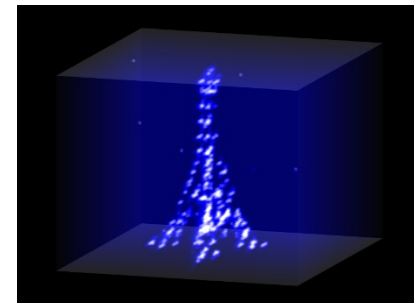
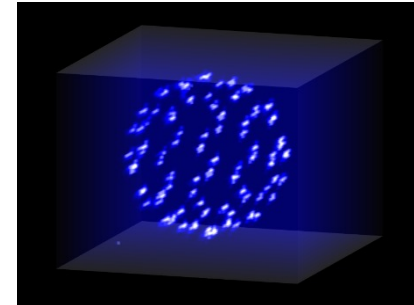
Introduction à la simulation quantique



Antoine Browaeys

*Laboratoire Charles Fabry,
Institut d'Optique, CNRS, FRANCE*

IMT Nantes, 10 novembre 2022



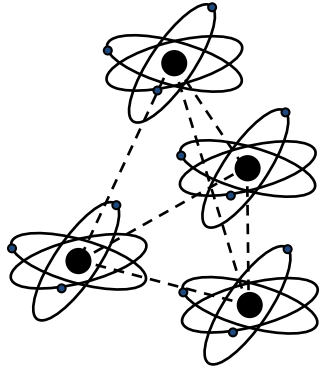
Le programme...

1. Physique à N-corps, simulation quantique et systèmes quantiques synthétiques
2. Exemples de simulateurs quantiques: ions et atomes
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 - A. Exploration du diagramme des phases
 - B. Dynamique hors-équilibre

Le problème à N-corps en physique

But: comprendre ensembles de **particules quantiques en interaction**

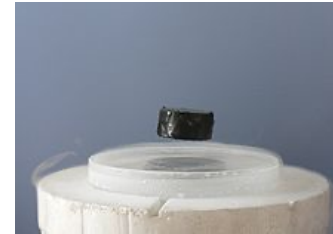
Microscopique



?



Macroscopique



Lois quantiques

Ce qu'il faut résoudre:

$$i\hbar \frac{\partial \Psi}{\partial t} = H_{\text{tot}} \Psi$$

$$H_{\text{tot}} = \sum_{i=1}^N -\frac{\hbar^2}{2m_i} \nabla_i^2 + \sum_{i=1}^N \sum_{j \neq i} \frac{q_i q_j}{r_{ij}} + \frac{\mu_B^2}{r_{ij}^3} \mathbf{s}_i \cdot \mathbf{s}_j$$

$N \approx 10^{23}$!!!

Complexité quantique: difficulté exponentielle

Fonction d'ondes à N -corps: $\Psi = \Psi(1, 2, \dots, N)$

« Taille » de la fonction d'onde ??

Exemple: particules à 2 d. de liberté (spin...) $\psi_i = \begin{pmatrix} a \\ b \end{pmatrix}$

$\Rightarrow \Psi$ nécessite 2^N composantes

Coder sur ordi. Ψ pour $N = 40 \Rightarrow 2^{50} \sim 10^{15} = \mathbf{1000\ To\ RAM\ !!}$

Record de calcul *ab-initio* (2022) $N \sim 50$

Des méthodes d'approximation...

$$H_{\text{tot}} = \sum_{i=1}^N \frac{p_i^2}{2m_i} - \frac{e^2}{r_i} + \sum_{i=1}^N \sum_{j \neq i} \frac{e^2}{r_{ij}}$$

Approx 1 : négliger interactions (mauvaise...)

$$H_{\text{tot}} = \sum_{i=1}^N H_i \Rightarrow \Psi(1, 2, \dots, N) = \psi_1 \psi_2 \dots \psi_N, \quad H_i \psi_i = E_i \psi_i$$

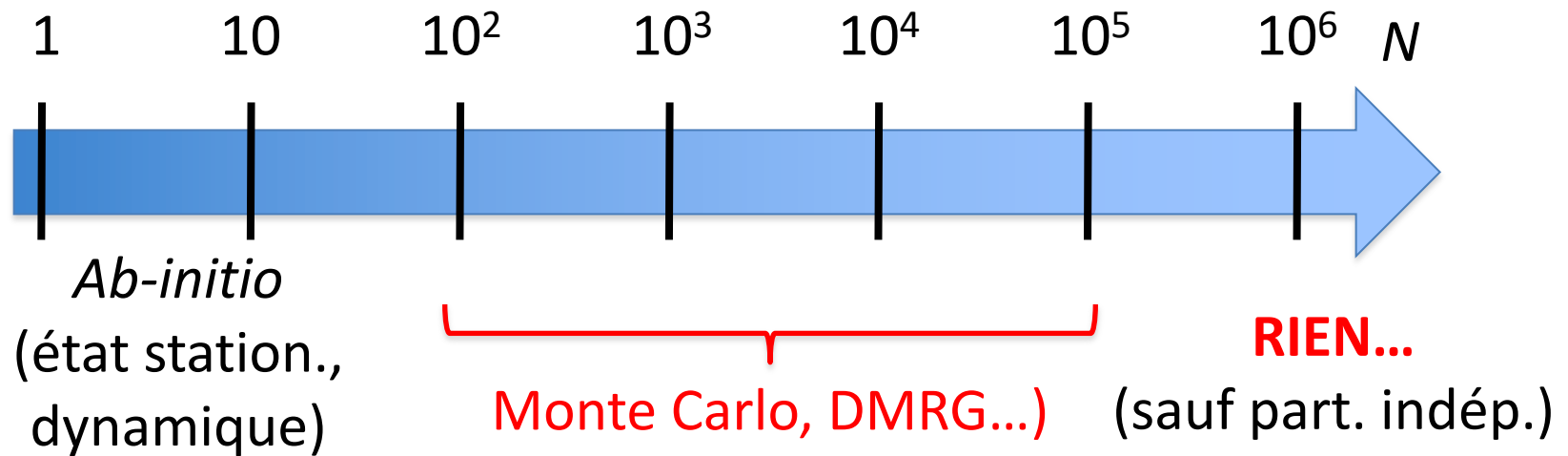
Approx 2 : champ moyen + perturbation (phys. atom, nucléaire, moléculaire, théorie des bandes phys. solide...)

$$H_{\text{tot}} = \sum_{i=1}^N \frac{p_i^2}{2m_i} + V(r_i) + \left[\sum_{i=1}^N \sum_{j \neq i} \frac{e^2}{r_{ij}} - \frac{e^2}{r_i} - V(r_i) \right]$$

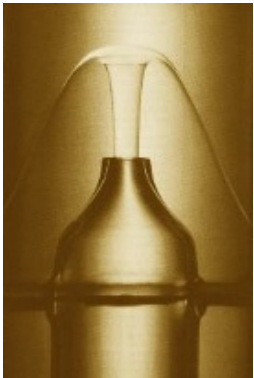
Méthode + raffinée : Monte Carlo Quantique, fonctionnel densité, DMRG, t-DMRG, Tensor Networks, Matrix Product states...

... qui ne marchent pas toujours !

Quand **interactions dominant** (systèmes **fortement corrélés**)



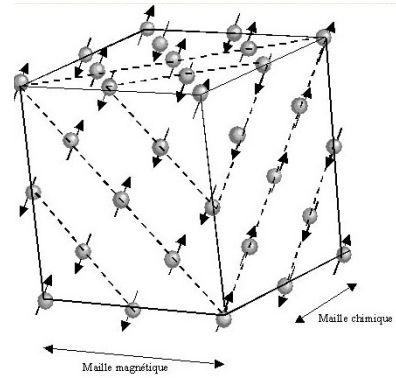
Exemples



superfluidité



supraconductivité

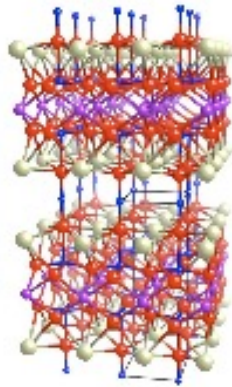
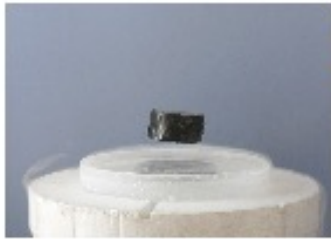


magnétisme



neutron star

Simulation quantique: exemple de démarche



Observation phénomène compliqué

Ex: supraconductivité haute T_c

Expérience sur
vrai système

simplification

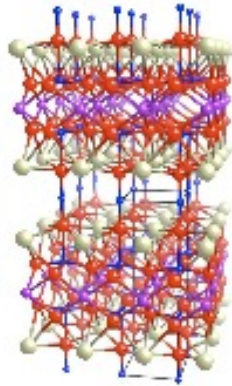
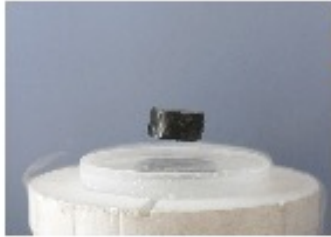


Calculs trop durs...

Hamiltonien modèle

$$H_{\text{model}} = - \sum_{i,j} J_{ij} a_i^\dagger a_j + \sum_i g (a_i^\dagger)^2 (a_i)^2$$

Simulation quantique: exemple de démarche



Observation phénomène compliqué

Ex: supraconductivité haute T_c

simplification

Expérience sur
vrai système



Hamiltonien modèle

“Préparable”

$$H_{\text{model}} = - \sum_{i,j} J_{ij} a_i^\dagger a_j + \sum_i g (a_i^\dagger)^2 (a_i)^2$$

$$|\psi(t)\rangle = e^{-\frac{i}{\hbar} H_{\text{model}} t} |\psi(0)\rangle$$

Lab...

Résultat mesure sur
simulateur: **état
fondamental = supra?**



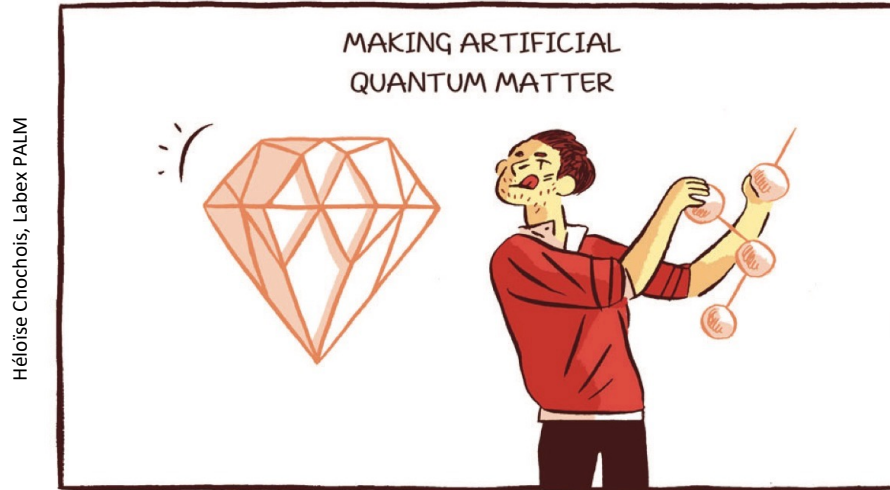
**Simulateur quantique =
Ingénierie d'atomes
gouvernés par H_{model}**

Problème à N-corps et systèmes quantiques synthétiques



R.P. Feynman

Simulating Physics with Computers, Int. J. Theo. Phys. **21** (1982)



i.e. systèmes quantiques **contrôlés** réalisant des **hamiltoniens à N-corps** (y compris artificiels...)

Plus grande programmabilité que systèmes réels
(géométrie, paramètres...)

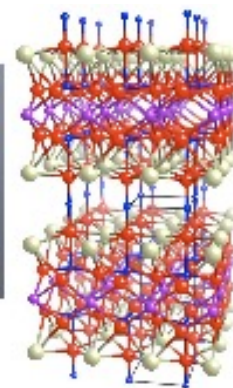
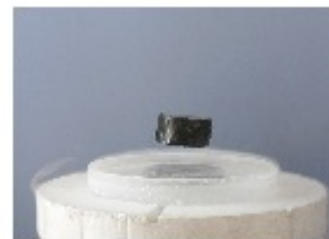
+

Nouvelles sondes, méthodes d'étude (ex.: dynamique...)

Qu'est ce que l'on peut simuler et à quoi ça peut servir...?

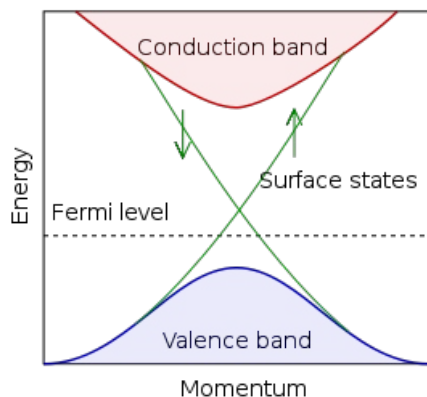
Supra-conductivité haute T_c

Vers le stockage de l'électricité?



Propriétés de conduction des métaux

influence du désordre et des interactions



Vers des métaux meilleurs conducteurs électriques?

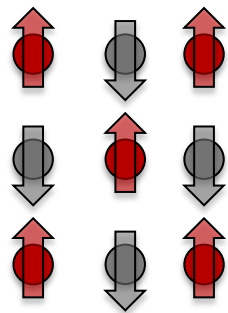
Magnétisme quantique

vers de nouveaux aimants pour moteurs électriques, stockage information...?



Les systèmes à N-corps les plus simples: spins en interaction

Particules de spin 1/2 sur un réseau:



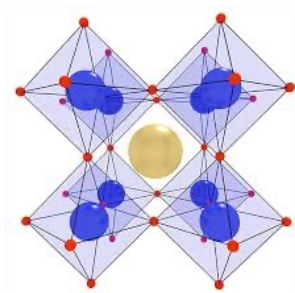
$$H \sim J \mathbf{S}_i \cdot \mathbf{S}_j$$

Ising $\hat{H} = \sum_{i \neq j} J_{ij} \hat{\sigma}_z^{(i)} \hat{\sigma}_z^{(j)}$

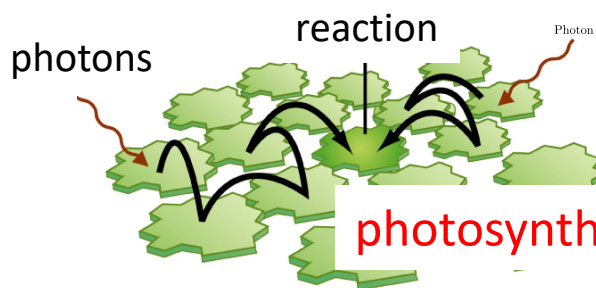
XY model $\hat{H} = \sum_{i \neq j} J_{ij} (\hat{\sigma}_i^+ \hat{\sigma}_j^- + \hat{\sigma}_i^- \hat{\sigma}_j^+)$

Magnétisme

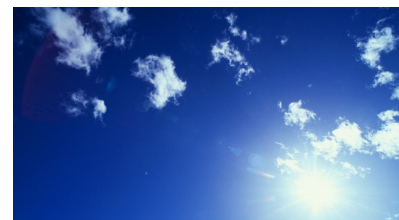
Propriétés de transport



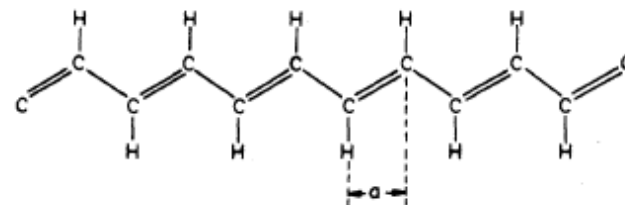
Perovskite
 $\text{Y}_2\text{Ti}_2\text{O}_7$



photosynthesis
excitons



Light scattering



Modèles génériques pour étude de transition de phase, dynamique...

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A. Ions piégés

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C. Atomes froids dans des pinces optiques

3. Exemples de simulation quantique

Manipulation de particules individuelles: Le point de vue d'un père fondateur...

..., nous ne faisons **jamais** d'expériences avec **juste un** électron, un atome ou une (petite) molécule. Dans des expériences de pensée, nous supposons parfois que nous le faisons ; cela conduit invariablement à des conséquences **ridicules**.



E. Schrödinger
British Journal of the Philosophy
of Science III (10), (1952)

60 ans plus tard...



2012



S. Haroche (France)

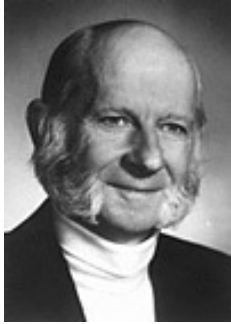


D. Wineland (USA)

**« pour le développements de méthodes
expérimentales permettant de mesurer et de
manipuler des systèmes quantiques individuels »**

Piégeage de particules uniques: les débuts...

H. Dehmelt



1989



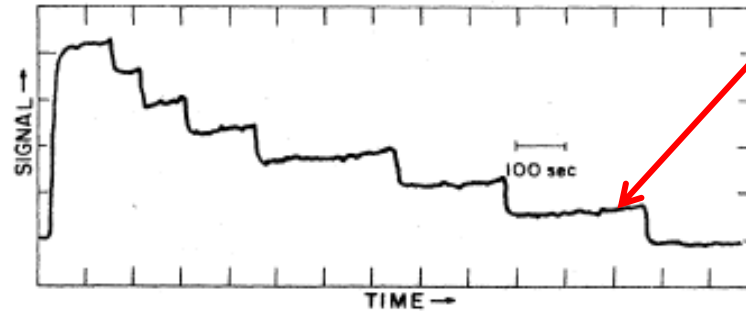
W. Paul: techniques de piégeage d'ions

1^{er} ion unique:

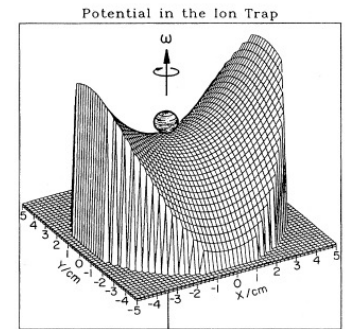
VOLUME 31 19 NOVEMBER 1973 NUMBER 21

Monoelectron Oscillator

D. Wineland, P. Ekstrom, and H. Dehmelt
Department of Physics, University of Washington, Seattle, Washington 98195
(Received 13 August 1973)



Electron unique



PHYSICAL REVIEW A

VOLUME 22, NUMBER 3

SEPTEMBER 1980

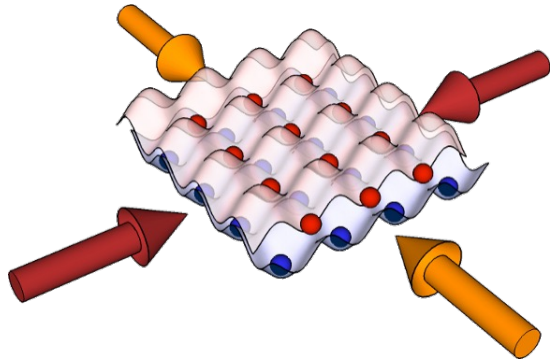
Localized visible Ba⁺ mono-ion oscillator

W. Neuhauser, M. Hohenstatt, and P. E. Toschek
Institut für Angewandte Physik I der Universität Heidelberg, D-69 Heidelberg, Federal Republic of Germany

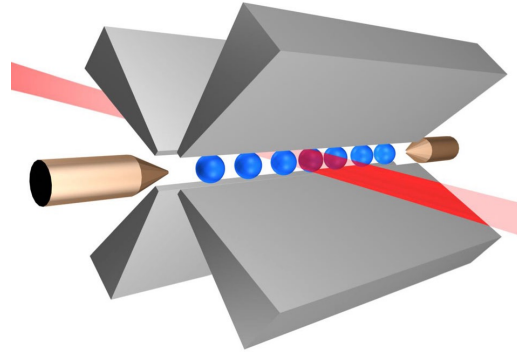
H. Dehmelt

Department of Physics, University of Washington, Seattle, Washington 98195
(Received 11 September 1979)

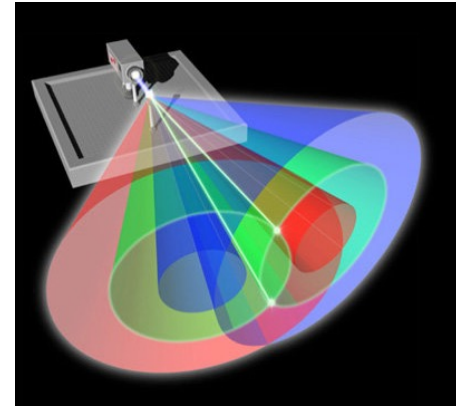
Ingénierie Quantique avec des systèmes *individuels*



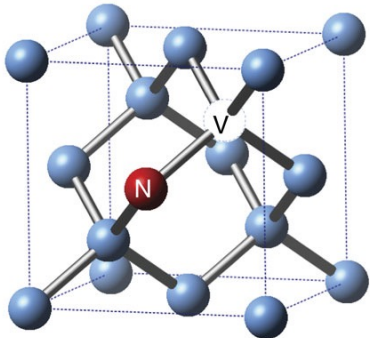
Atomes et molécules froides



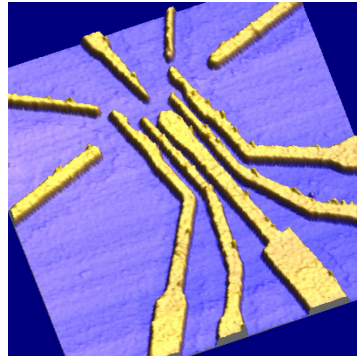
Ions piégés



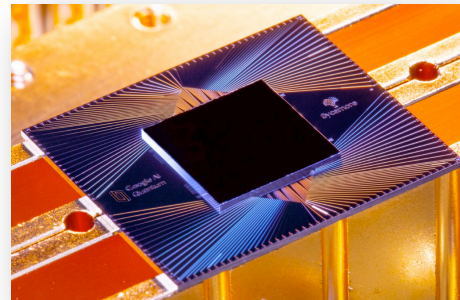
Photons



Centres NV

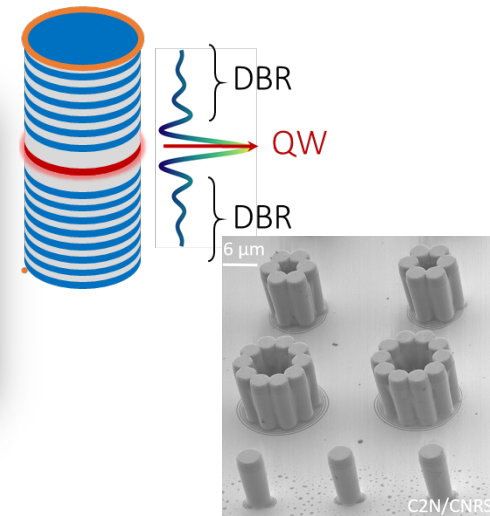


Spin électronique



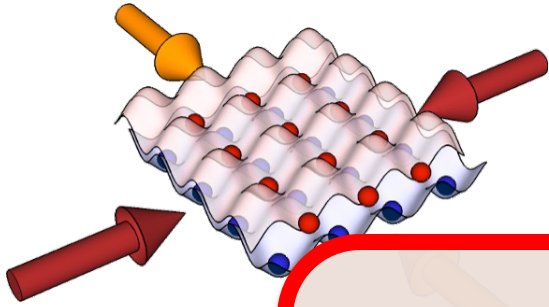
Sycamore processor, Google Inc.

Circuits supracond.

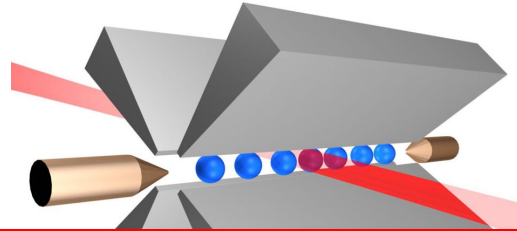


Polaritons (½ cond.)

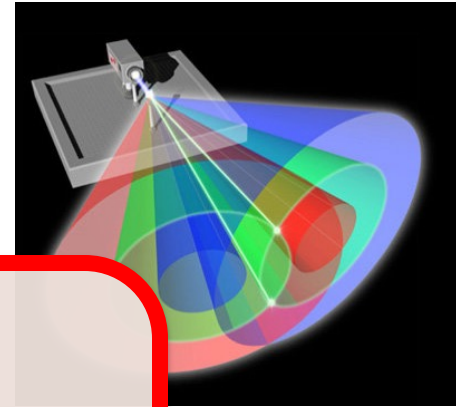
Ingénierie Quantique avec des systèmes *individuels*



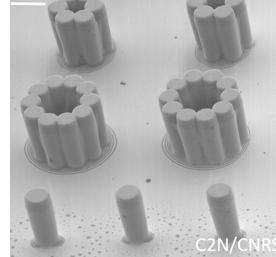
Atomes et molécules froides



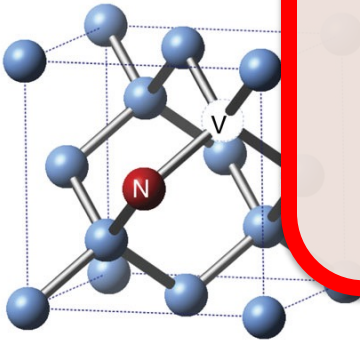
Circuits supracond.



Photons



Polaritons (½ cond.)



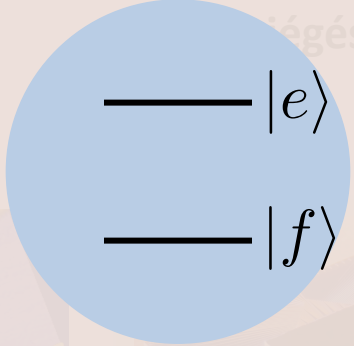
Centres NV



Spin électronique

Sycamore processor, Google Inc.

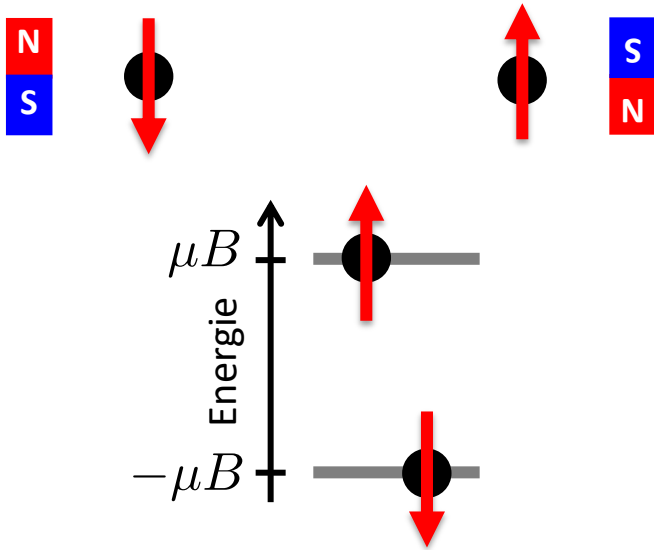
Systeme à 2 niveaux :



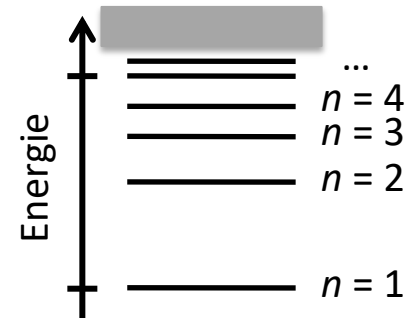
Adressable + interactions **controlées**

Correspondance entre système à 2 niveaux et spin 1/2

Spin 1/2 dans un champ B

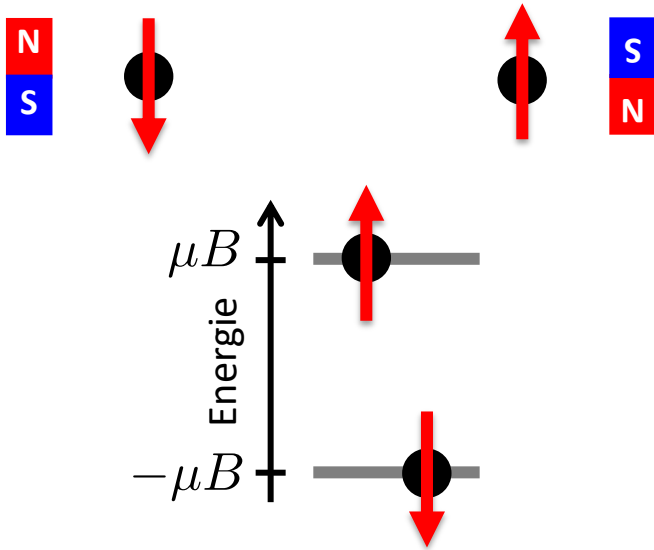


Atomes

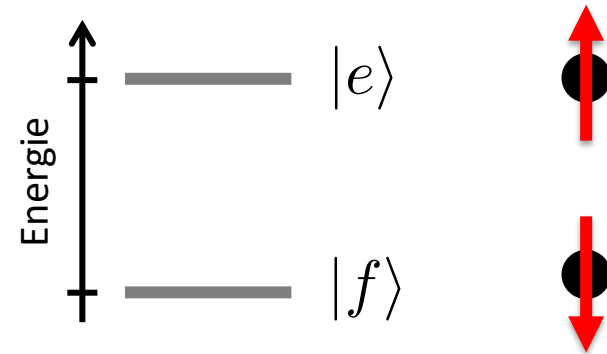


Correspondance entre système à 2 niveaux et spin 1/2

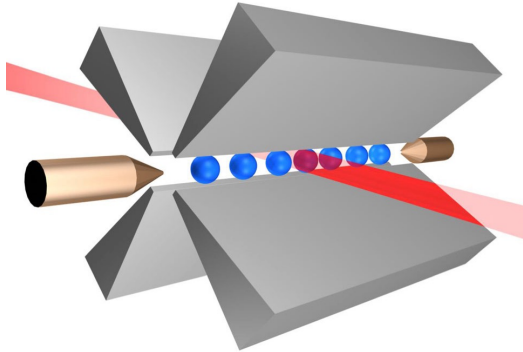
Spin 1/2 dans un champ B



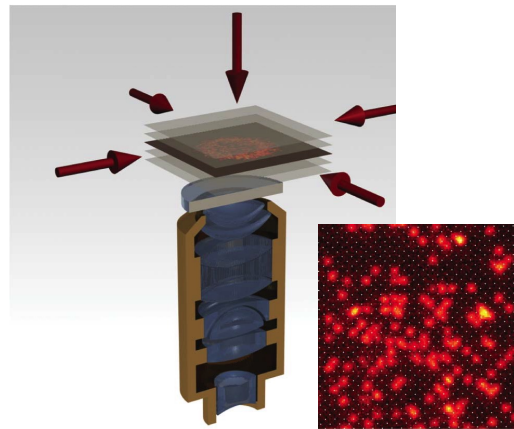
Atomes « à 2 niveaux »



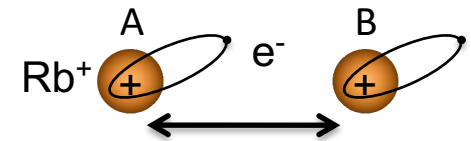
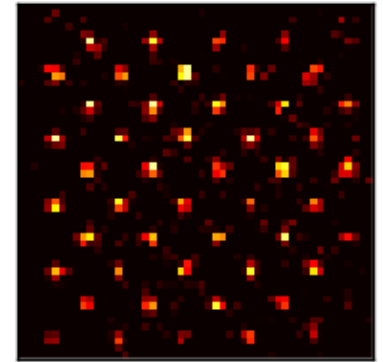
Exemples de réseaux d'atomes individuels



Ions piégés



Atomes dans réseaux optiques



Atomes dans pinces optiques

Scalable: plus de 100 atomes

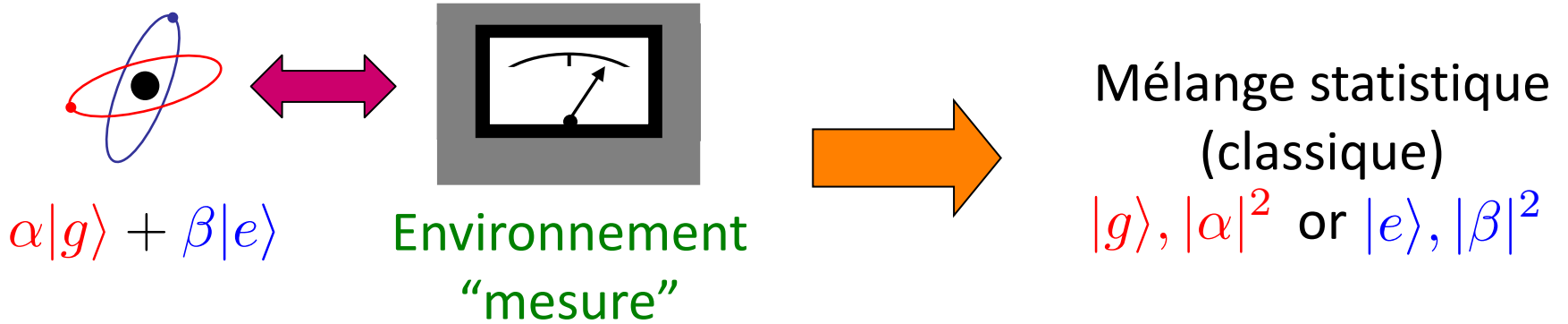
Adressable: manipulations et mesures locales

$$\langle \sigma_i^\alpha \rangle, \langle \sigma_i^\alpha \sigma_j^\beta \rangle, \dots$$

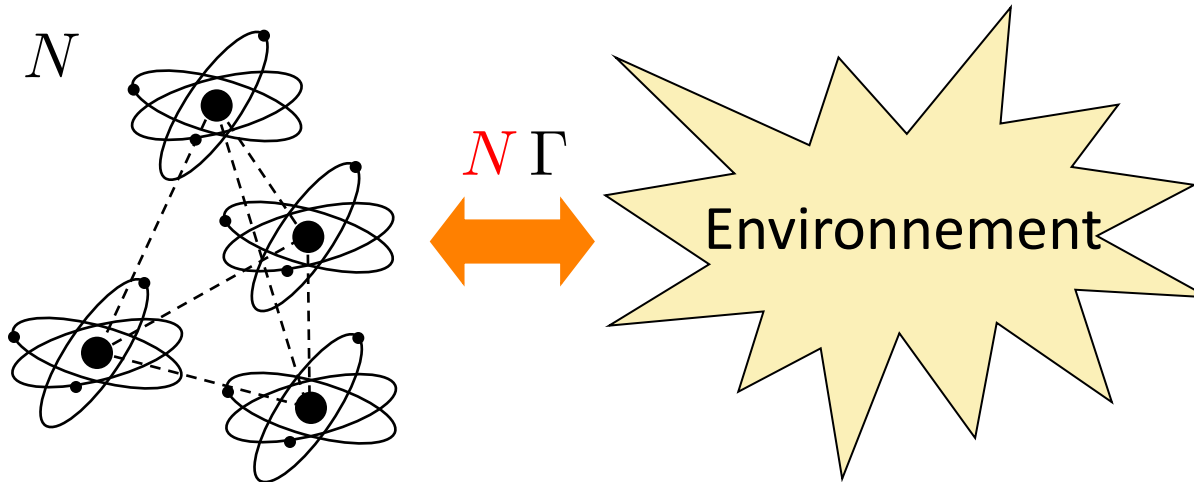
Programmable: contrôle géométrie, interactions...

Pourquoi ingénierie quantique est-elle si difficile ?

Superposition quantique = **très fragile!!** Décohérence



Plus c'est gros, plus ça « décohère »...



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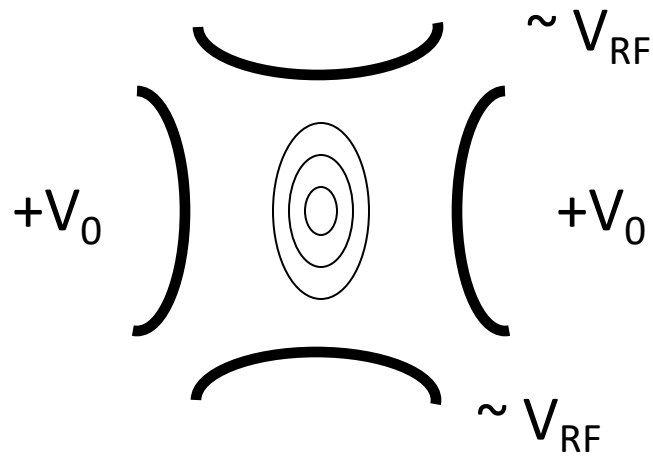
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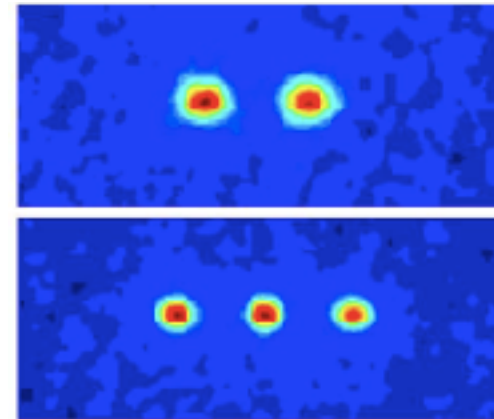
Piégeage d'ions refroidis par laser

Piège de Paul (RF + statique)

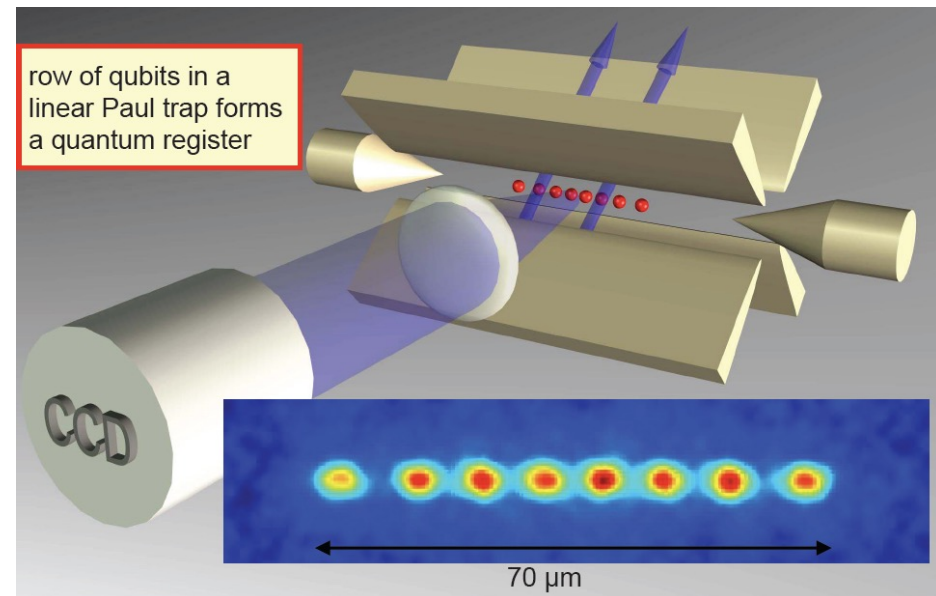
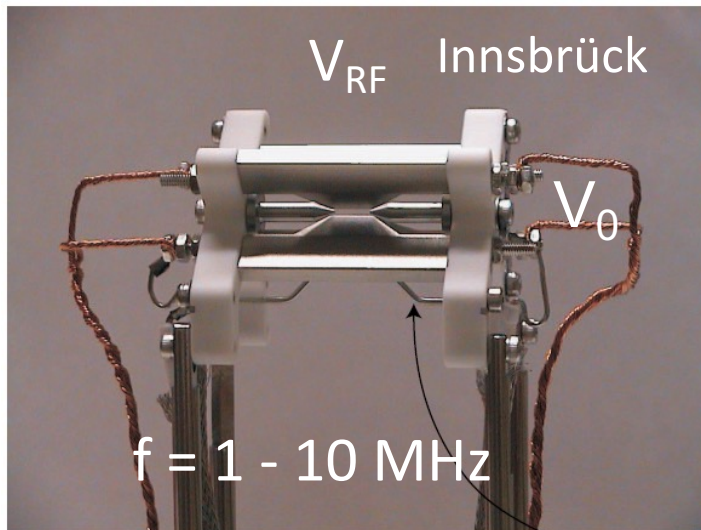


Répulsion de Coulomb

\Rightarrow Cristal ionique



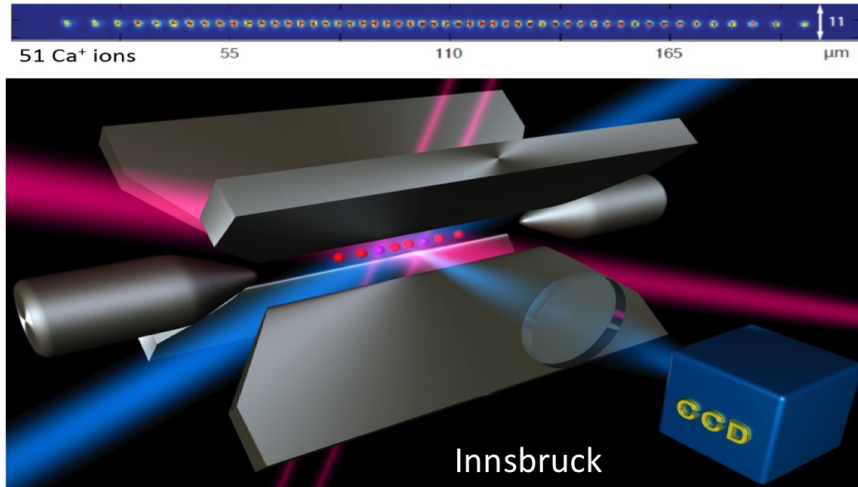
Piège de Paul linéaire



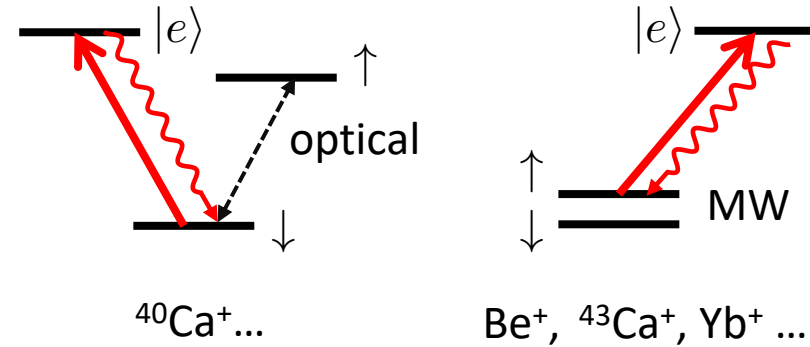
Innsbrück

Simulateur à ions piégés: modèles de spin

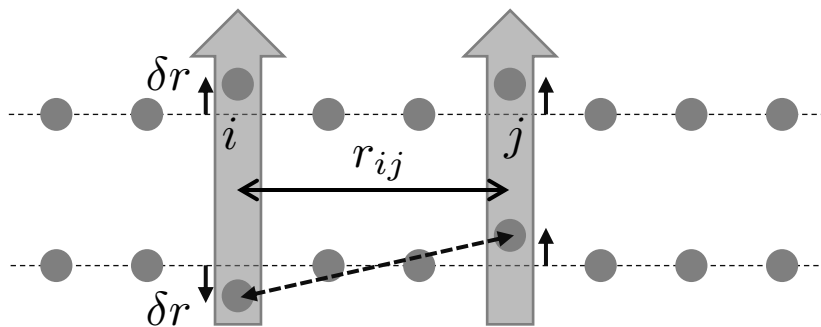
Cristal de Coulomb



« Coder » le spin 1/2



Force lumineuse dépendant de l'état



Les Hamiltoniens

$$H_{XY} = \sum_{i,j} J_{ij} (\hat{\sigma}_i^+ \hat{\sigma}_j^- + \hat{\sigma}_i^- \hat{\sigma}_j^+)$$

$$H_{\text{Ising}} = \sum_{i,j} J_{ij} \hat{\sigma}_i^x \hat{\sigma}_j^x + \frac{\delta}{2} \sum_i \hat{\sigma}_i^z$$

$$J_{ij} \approx \frac{J_0}{|i-j|^\alpha}, \quad \alpha = 0.5 - 2.5$$

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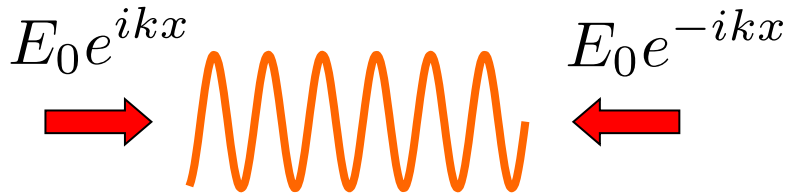
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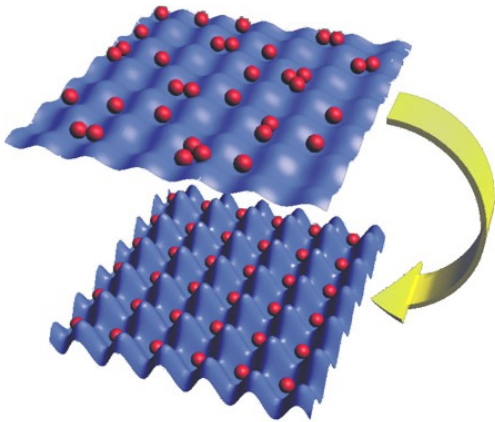
Atomes ultra-froids dans les réseaux optiques

Force dipolaire: $\mathbf{F} \propto -\nabla I(\mathbf{r})$

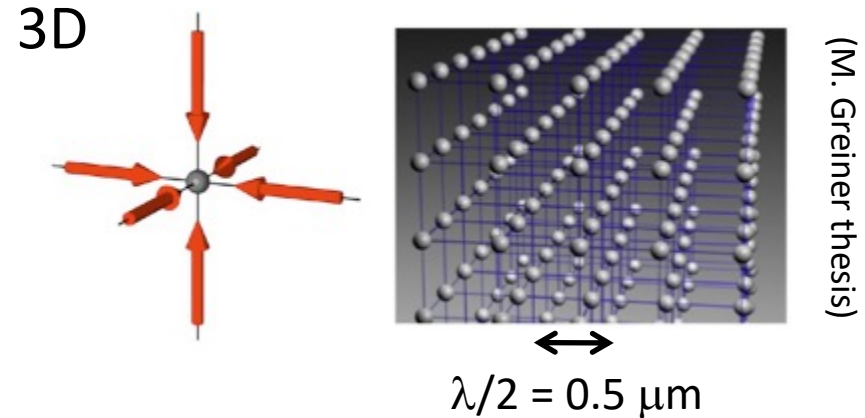


$$I(x) = 2E_0^2(1 + \cos 2kx)$$

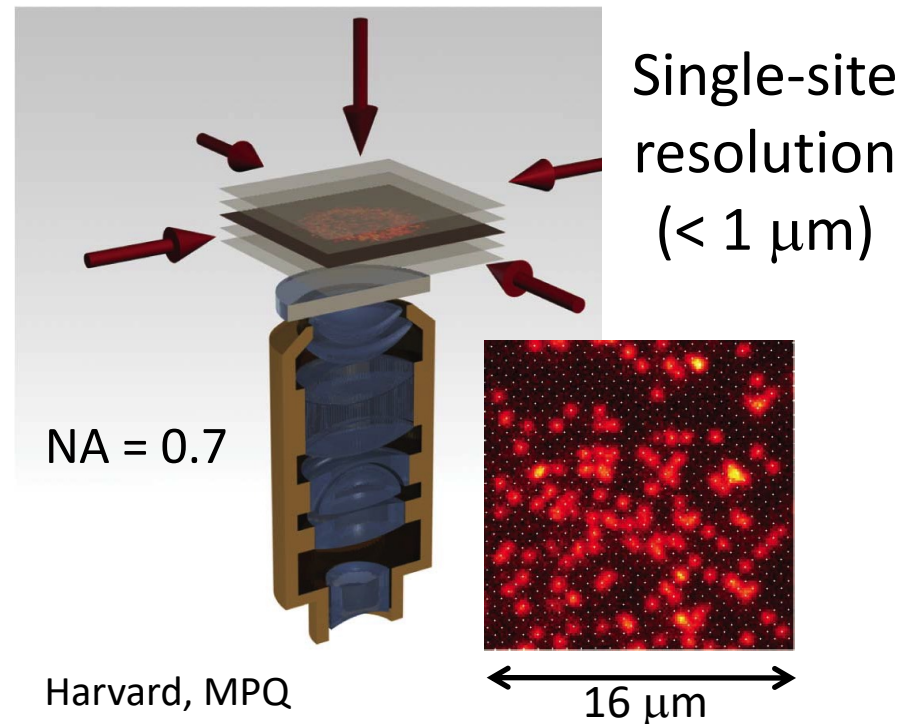
Chaque site contient 1 atome !



Boson (Rb, Na, ^7Li , ^{39}K , $^4\text{He}^*$),
Fermion (^6Li , ^{40}K),
Magnetic atoms (Cr, Dy...)

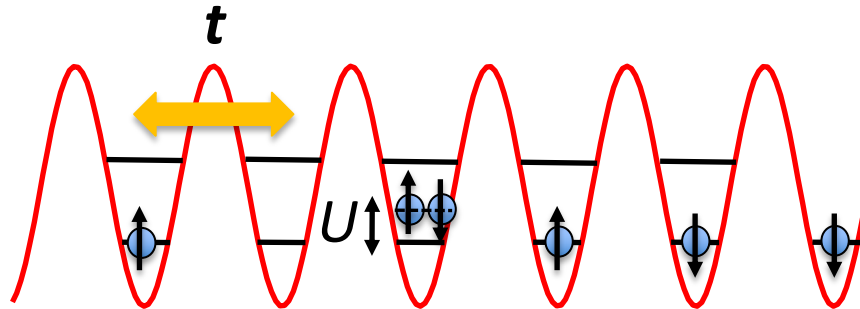


Quantum gas microscope



Réseaux optiques: modèles de Hubbard et de spin

Compétition effet tunnel / interaction



U contrôlable par B!!

Modèles d'Hubbard

$$H = -t \sum_{\langle i,j \rangle} (a_i^\dagger a_j + \text{h.c.}) + \frac{U}{2} \sum_i n_i (n_i - 1)$$

Modèle le plus simple:
conductivité / interaction

Le programme...

1. Physique à N-corps, simulation quantique et systèmes quantique synthétiques

2. Exemples de simulateurs quantiques

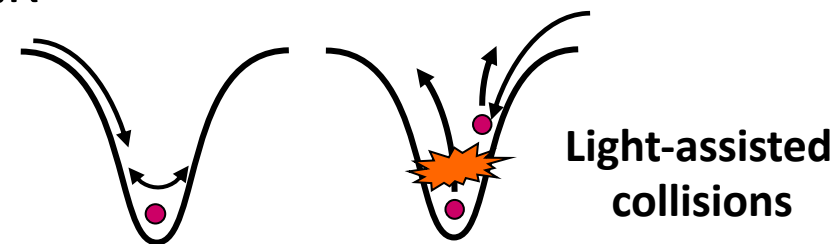
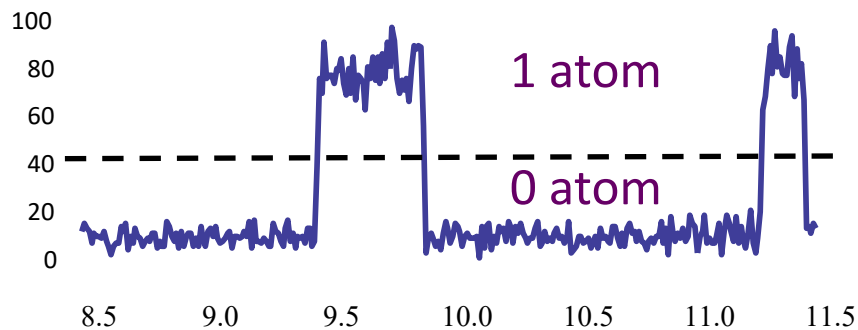
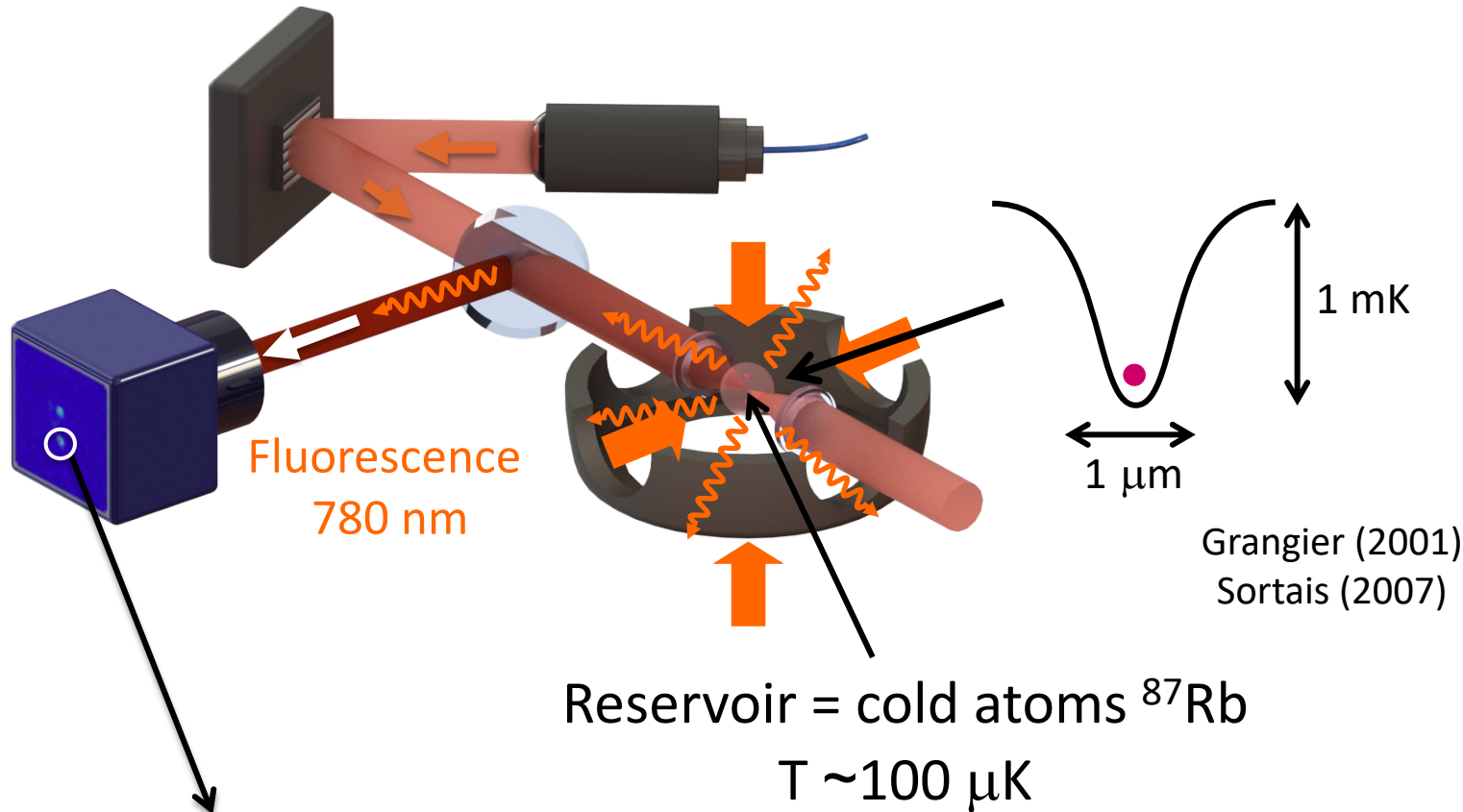
A. Ions piégés

B. Atomes ultra-froids dans des réseaux optiques

C. Atomes froids dans des pinces optiques

3. Exemples de simulation quantique

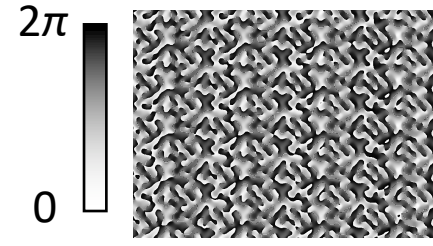
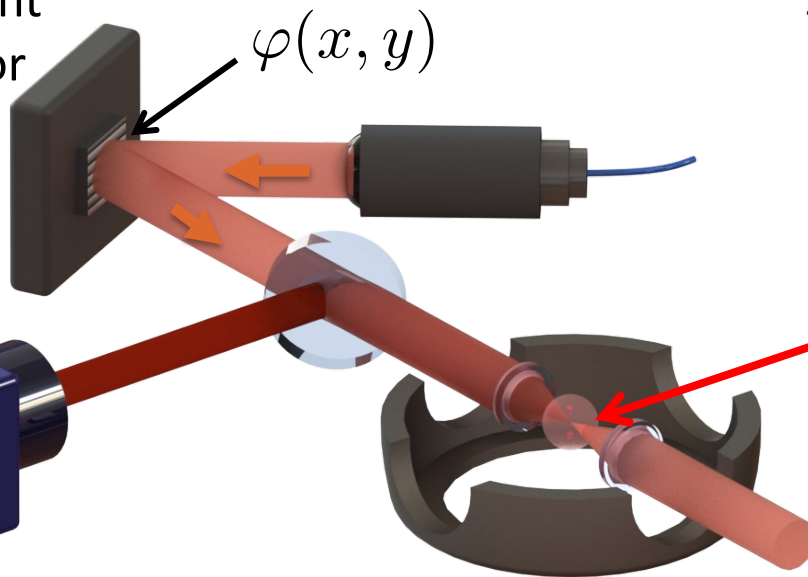
Atomes dans des matrices de pinces optiques



Non-deterministic
single-atom source

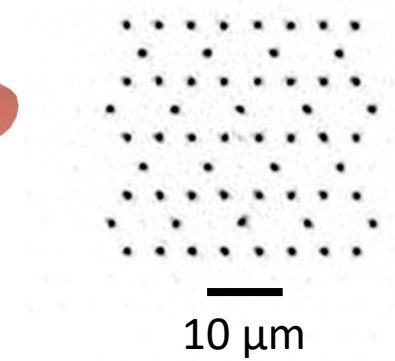
Atomes dans des matrices de pinces optiques

Spatial Light Modulator



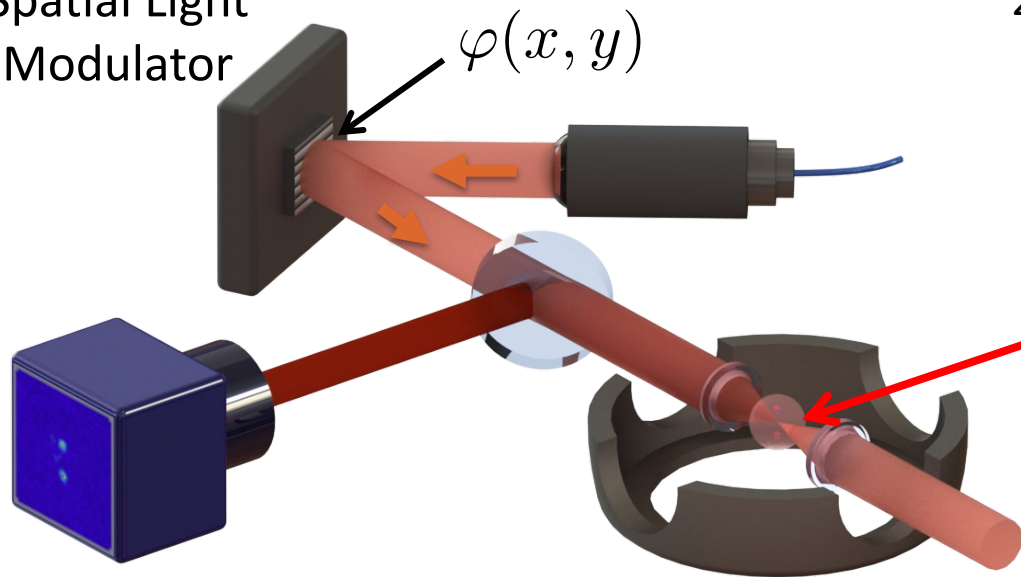
Nogrette, PRX (2014)

$$\left| \text{FT}[e^{i\varphi(x,y)}] \right|^2$$

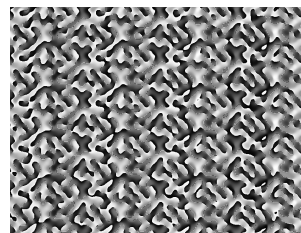


Atomes dans des matrices de pinces optiques

Spatial Light Modulator



2π



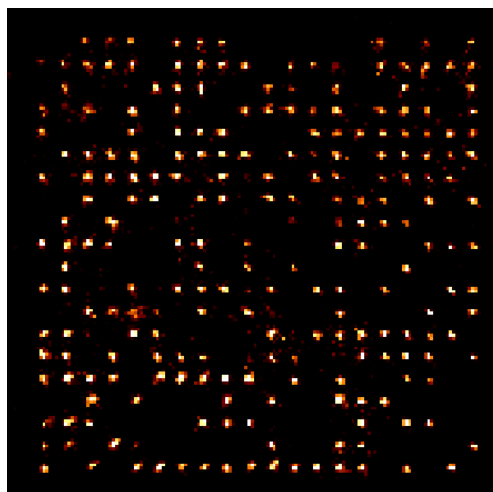
SLM pattern

0

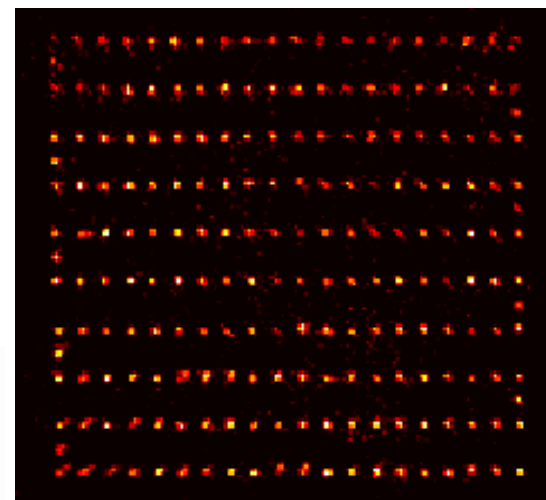
Nogrette, PRX (2014)

$$\left| \text{FT}[e^{i\varphi(x,y)}] \right|^2$$

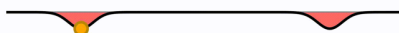
Initial configuration



Assembled configuration

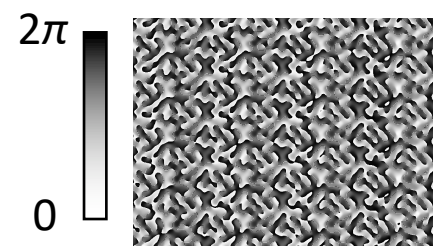
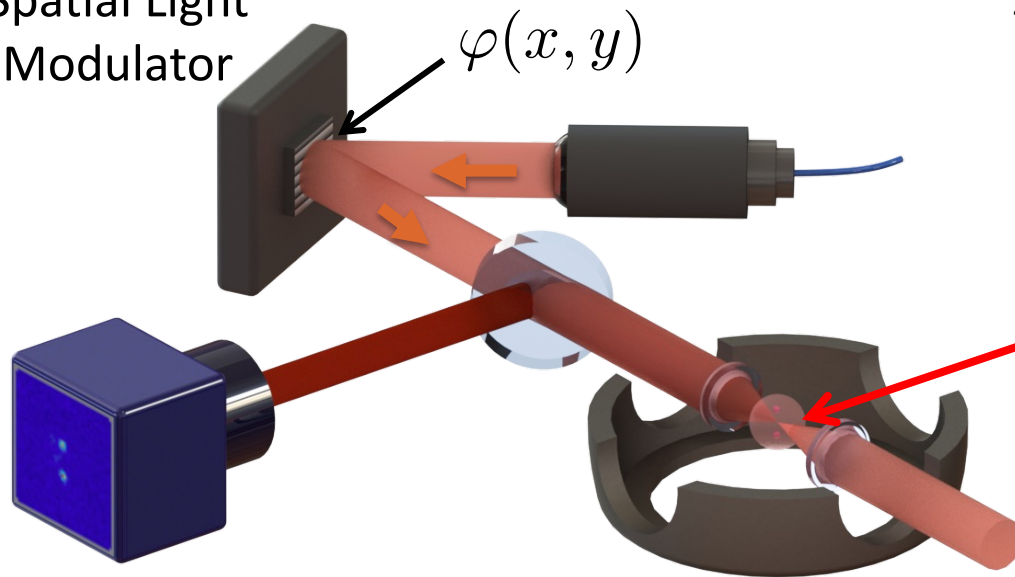


Assembling process



Atomes dans des matrices de pinces optiques

Spatial Light Modulator

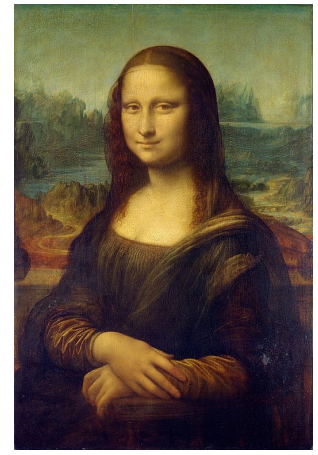


SLM pattern

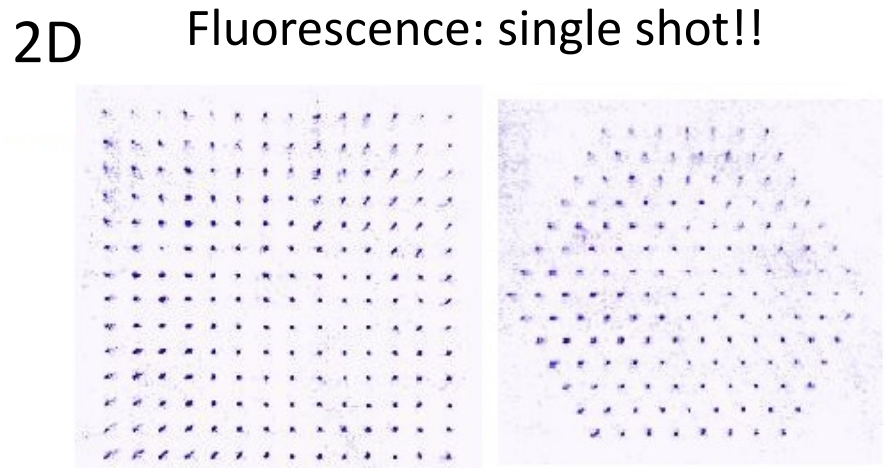
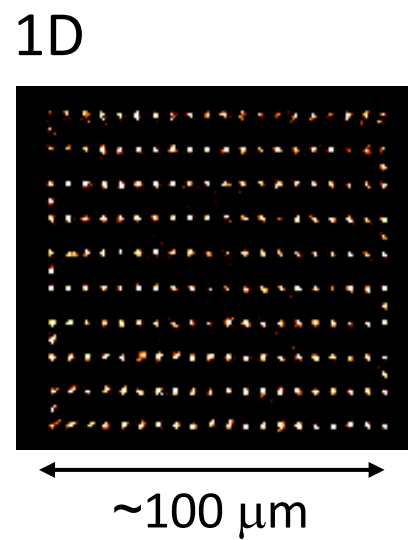
Nogrette, PRX (2014)

$$\left| \text{FT}[e^{i\varphi(x,y)}] \right|^2$$

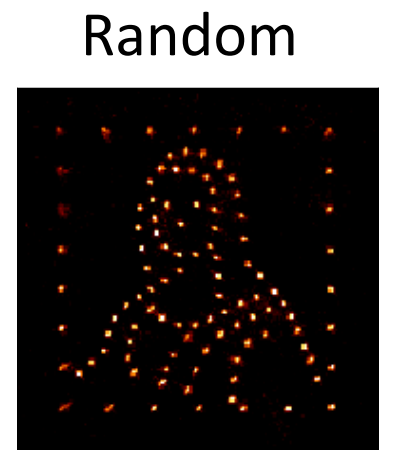
L. da Vinci



Assembled arrays of individual atoms ($N \sim 200$)

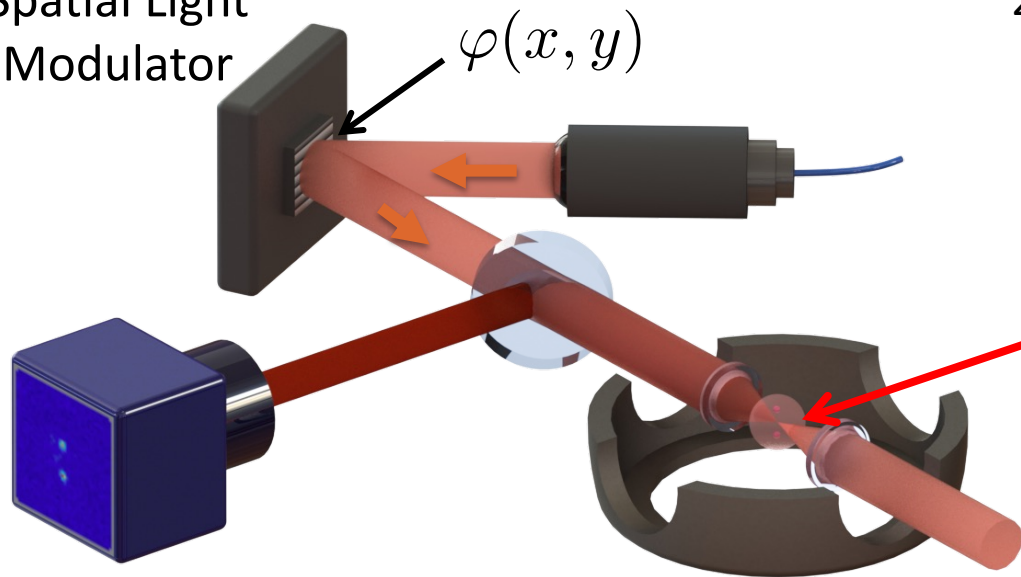


Barredo, Nature 2016 ; Schymik, PRA 2020

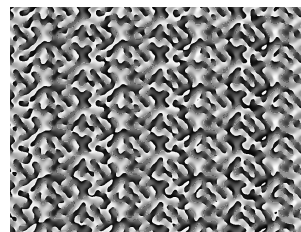


Atomes dans des matrices de pinces optiques

Spatial Light Modulator



2π



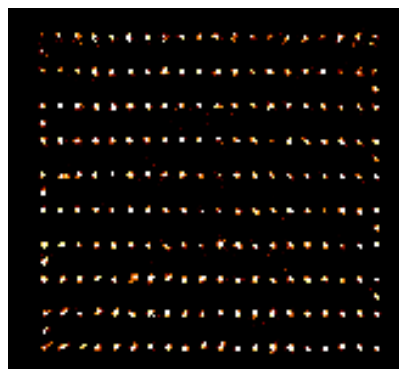
SLM pattern

Nogrette, PRX (2014)

$$\left| \text{FT}[e^{i\varphi(x,y)}] \right|^2$$

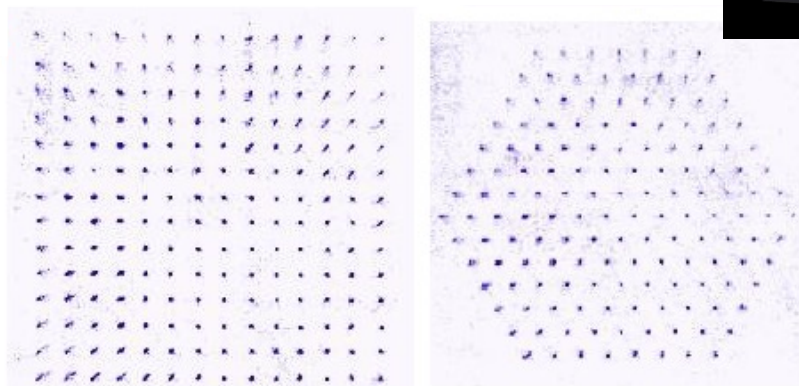
Assembled arrays of individual atoms ($N \sim 200$)

1D



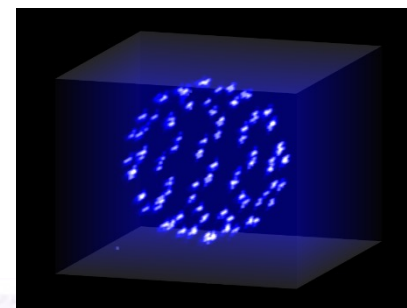
$\sim 100 \mu\text{m}$

2D

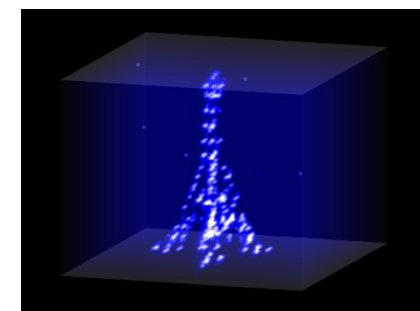


Barredo, Nature 2016 ; Schymik, PRA 2020

Fluorescence: single shot!!



(averaged)

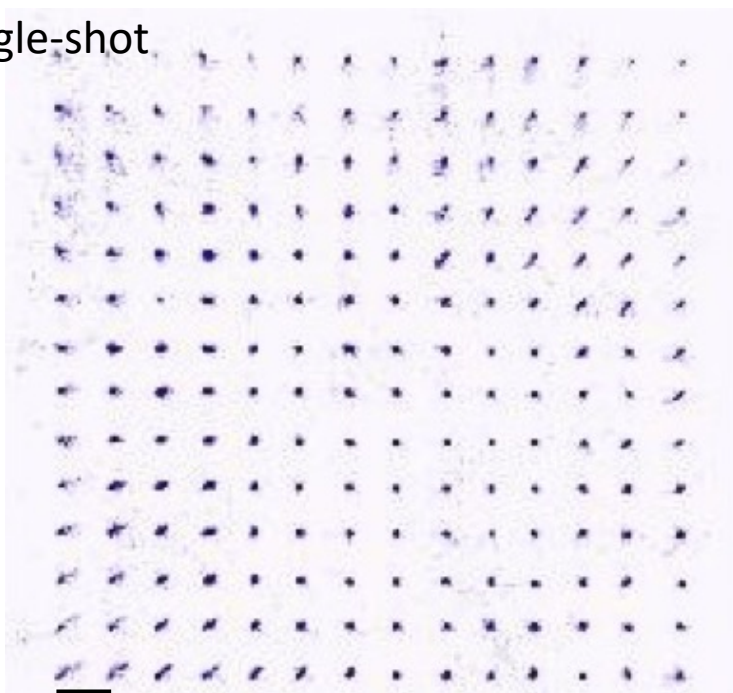


Barredo, Nature (2018)

Arrays of interacting Rydberg atoms

Arrays of atoms

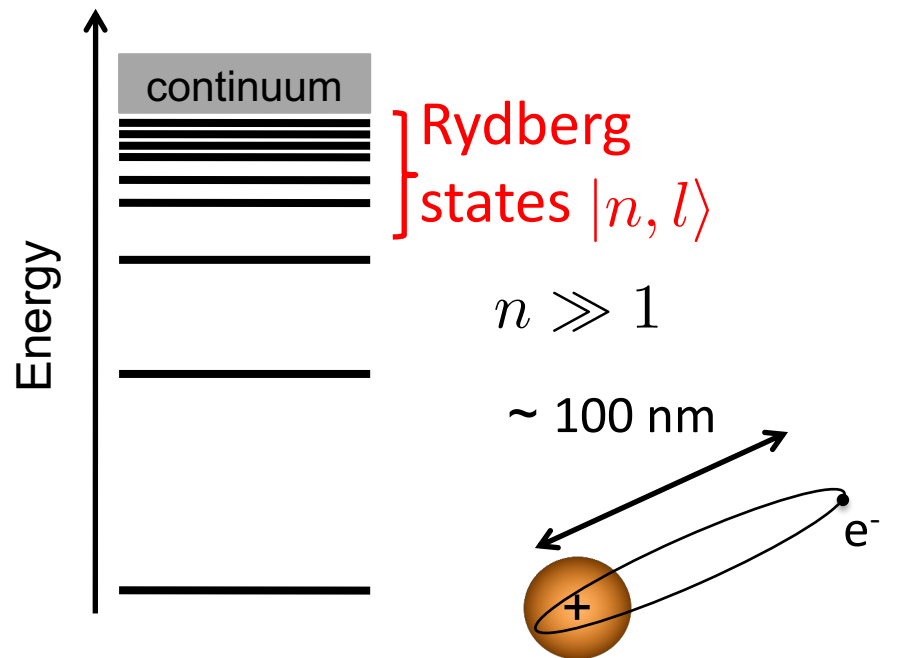
Single-shot



5 μm

Addressable!!

Rydberg atoms



Lifetime $> 100 \mu\text{s}$

Transition dipole: $d_{n, n\pm 1} \sim n^2 e a_0$

\Rightarrow Large dipole-dipole interactions

$$R = 10 \mu\text{m} \Rightarrow V_{\text{int}}/h \sim 1 - 10 \text{ MHz}$$

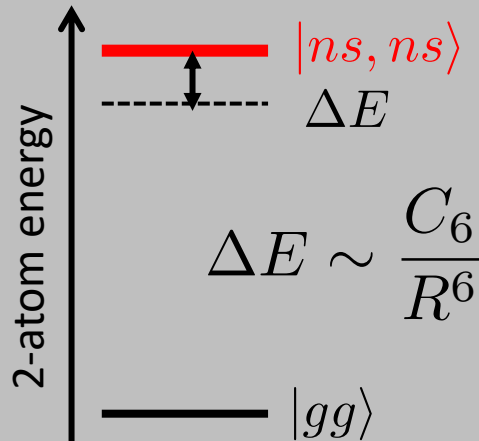
\Rightarrow timescales $< \mu\text{sec}$

Interactions between Rydberg atoms and spin models

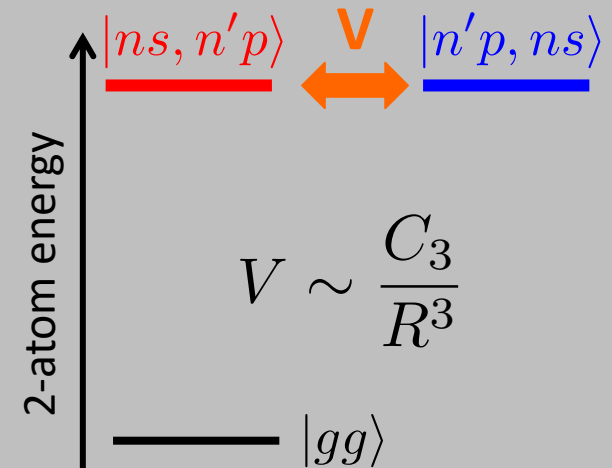


Browaeys & Lahaye, Nat.Phys. (2020)

van der Waals



Resonant dipole



Quantum Ising

$$\hat{H} = \sum_{i \neq j} J_{ij} \hat{\sigma}_z^{(i)} \hat{\sigma}_z^{(j)}$$

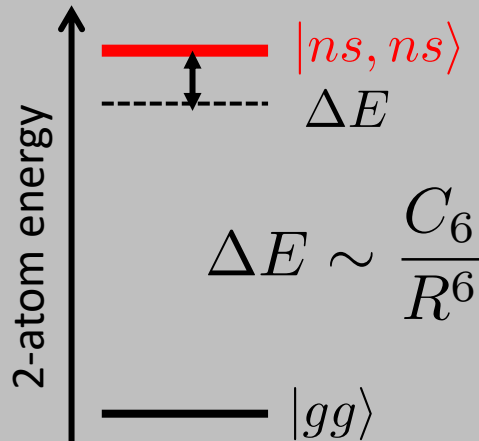
XY model

$$\hat{H} = \sum_{i \neq j} J_{ij} (\hat{\sigma}_i^+ \hat{\sigma}_j^- + \hat{\sigma}_i^- \hat{\sigma}_j^+)$$

From van der Waals interactions to Ising model...



van der Waals



$$C_6 \propto n^{11} \Rightarrow \text{switchable interaction}$$

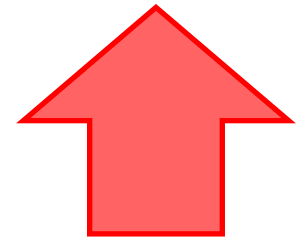
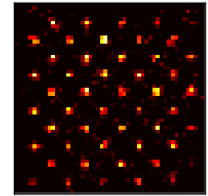
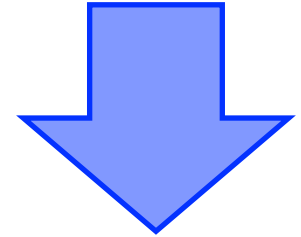
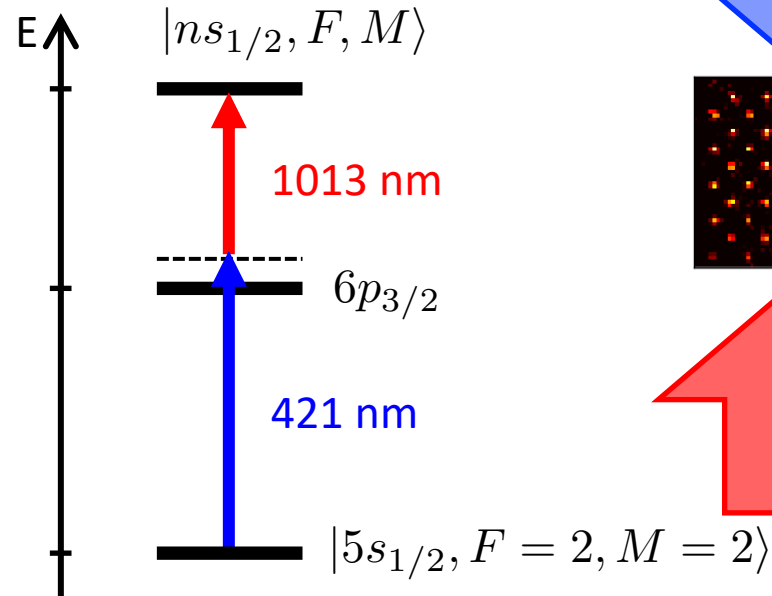
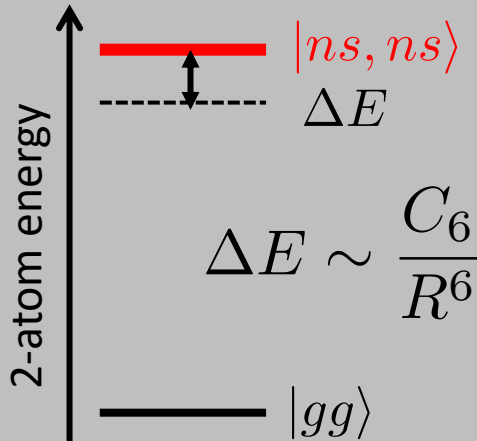
$$\hat{H}_{\text{int}} = \frac{C_6}{R^6} \hat{n}_1 \hat{n}_2 \sim J \hat{\sigma}_1^z \hat{\sigma}_2^z$$

Rydberg occupation
number

From van der Waals interactions to Ising model...



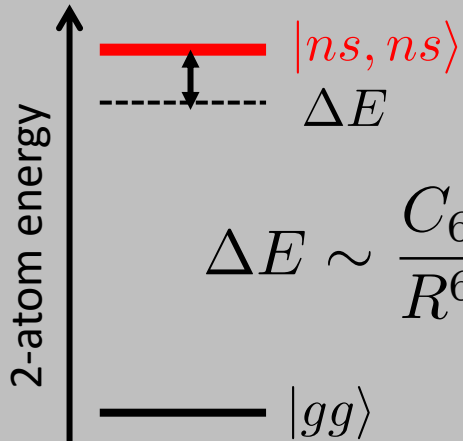
van der Waals



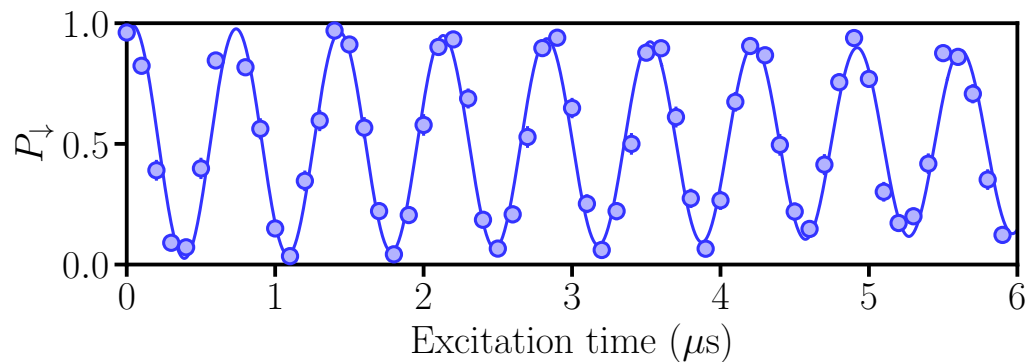
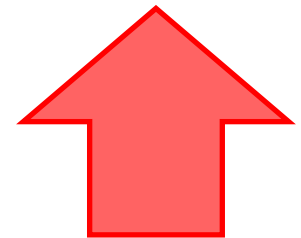
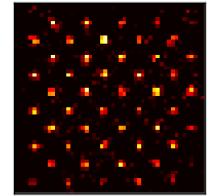
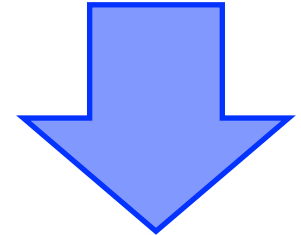
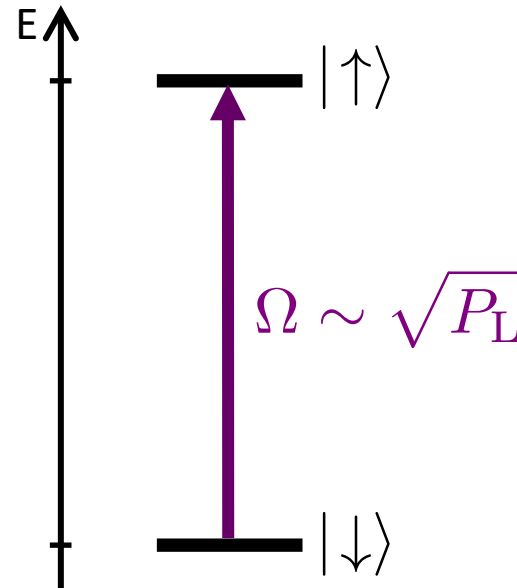
From van der Waals interactions to Ising model...



van der Waals

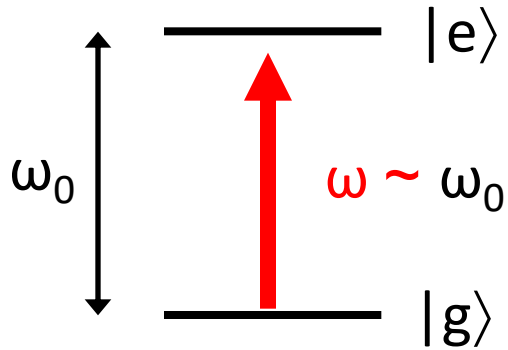


$$\Delta E \sim \frac{C_6}{R^6}$$



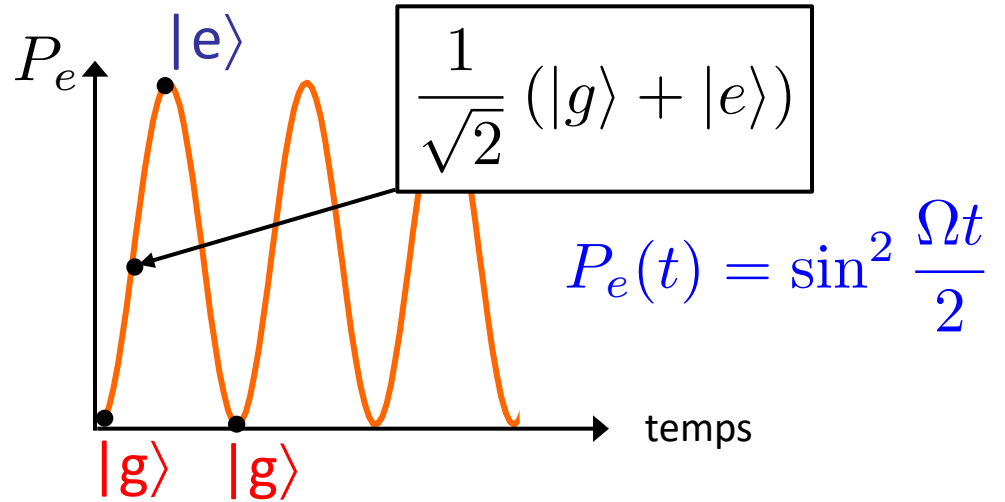
Les oscillations de Rabi à un atome

Oscillations de Rabi (1930s!): émission stimulée

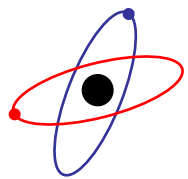


Ω = fréquence de Rabi

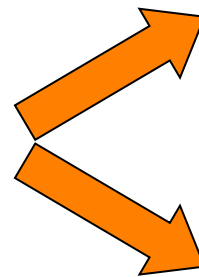
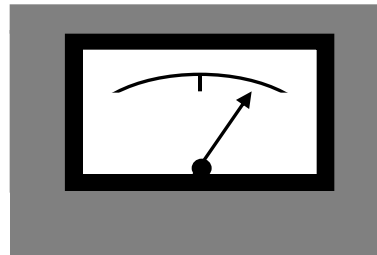
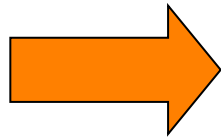
$$\propto \sqrt{\text{Intensity}}$$



Mesure sur un système individuel



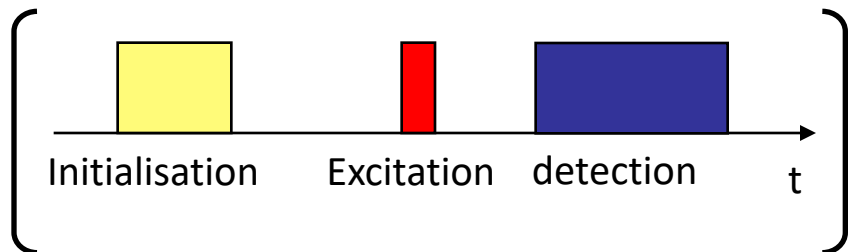
$$\alpha|g\rangle + \beta|e\rangle$$



$$|g\rangle, |\alpha|^2$$

OU

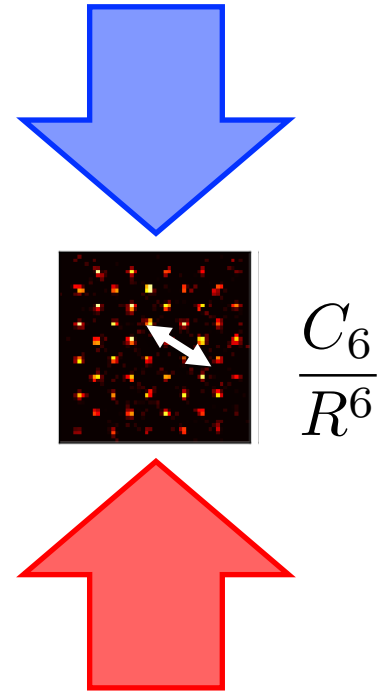
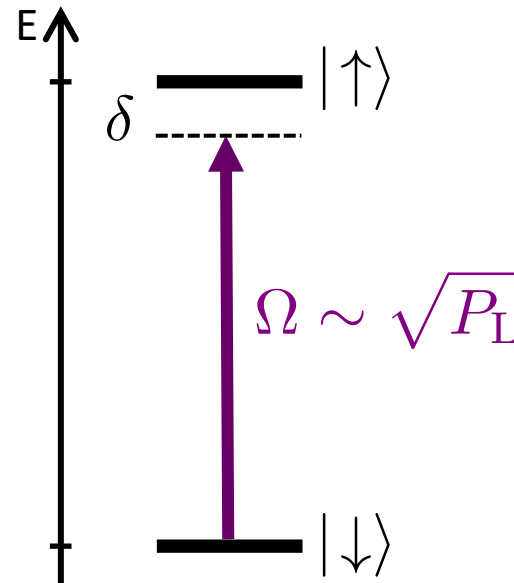
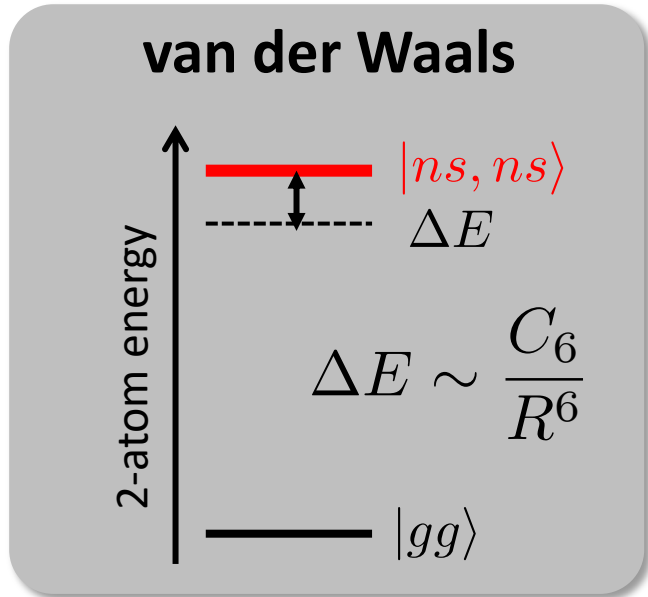
$$|e\rangle, |\beta|^2$$



$\times 100$

$|g\rangle, |e\rangle, |e\rangle, |g\rangle, \dots, |e\rangle, |g\rangle$
 \Rightarrow Probabilité $P(g)$

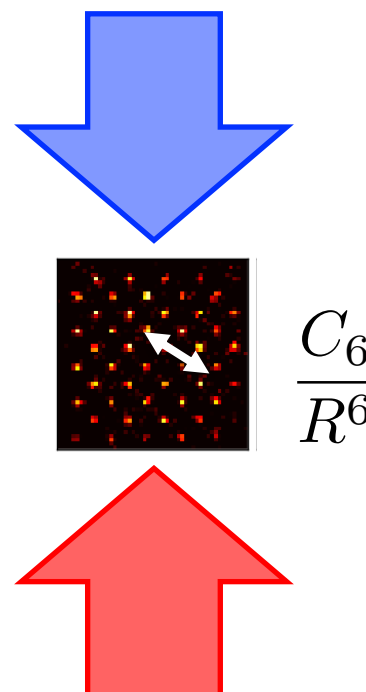
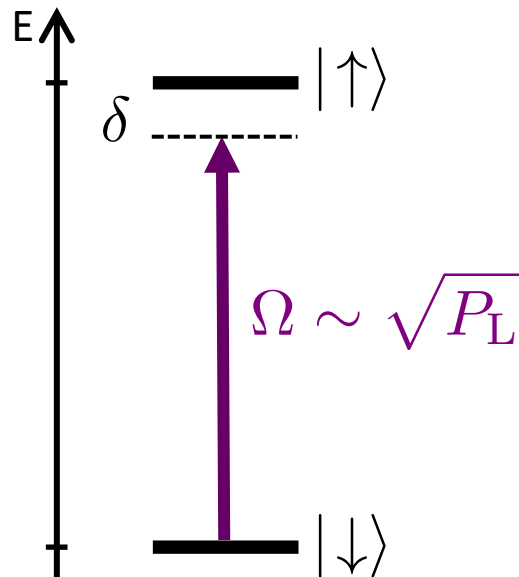
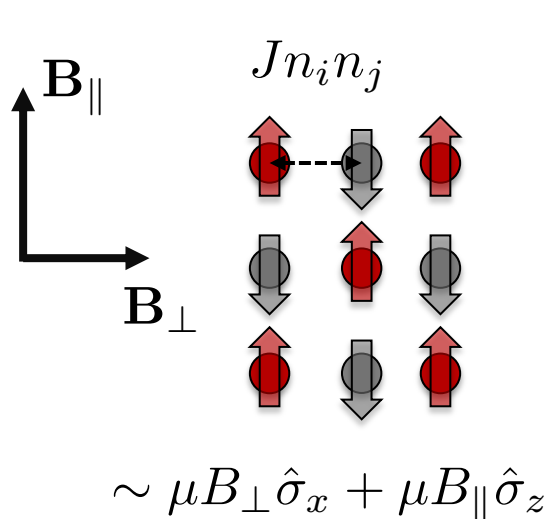
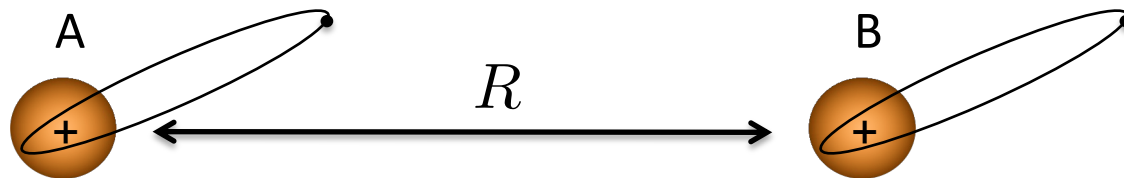
From van der Waals interactions to Ising model...



Quantum Ising-like model ($s=1/2$):

$$H = \frac{\hbar\Omega}{2} \sum_i \sigma_x^i + \hbar\delta \sum_i \hat{n}_i + \sum_{i<j} \frac{C_6}{R_{ij}^6} \hat{n}_i \hat{n}_j$$

From van der Waals interactions to Ising model...



Quantum Ising-like model ($s=1/2$):

$$H = \frac{\hbar\Omega}{2} \sum_i \sigma_x^i + \hbar\delta \sum_i \hat{n}_i + \sum_{i<j} \frac{C_6}{R_{ij}^6} \hat{n}_i \hat{n}_j$$

Experiment^t.

$$\frac{C_6/a^6}{\Omega} = [0 - 20]$$

Laser: B_{\perp}

B_{\parallel}

Spin-spin interaction

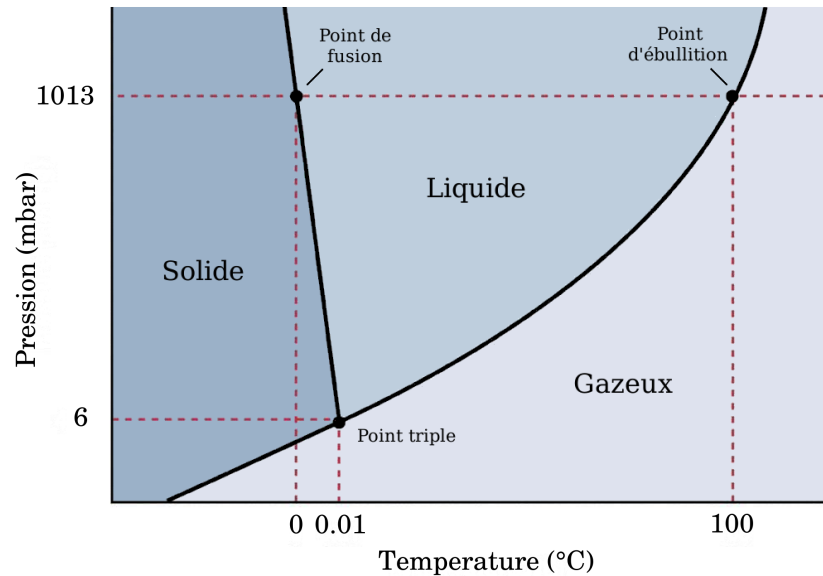
Le programme...

1. Physique à N-corps, simulation quantique et systèmes quantique synthétiques
2. Exemples de simulateurs quantiques: ions et atomes
3. Exemples de simulation quantique
 - A. Exploration du diagramme des phases
 - B. Dynamique hors-équilibre

Diagramme des phases

Etat fondamental d'un système en fonction de paramètres de contrôle

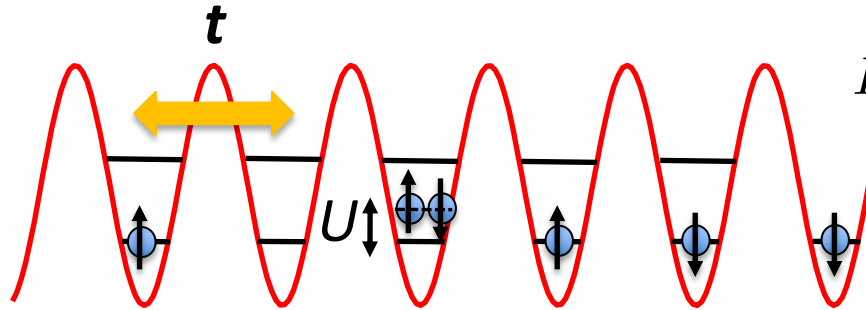
Diagramme des phases de l'eau (P, T)



Système quantique: $\langle \Psi_f | O | \Psi_f \rangle (T, B, M \dots)$

Atoms in optical lattices implements Hubbard models

Competition tunneling / interaction

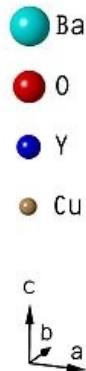
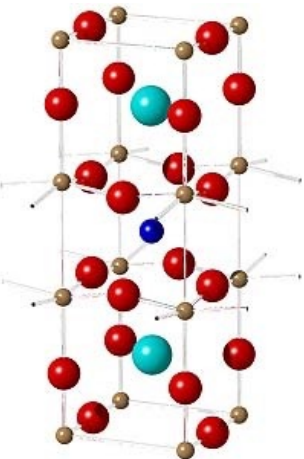


$$H = -t \sum_{\langle i,j \rangle, \sigma} (c_{i\sigma}^\dagger c_{j\sigma} + \text{h.c.}) + U \sum_i n_{i\downarrow} n_{i\uparrow}$$

Fermi-Hubbard models

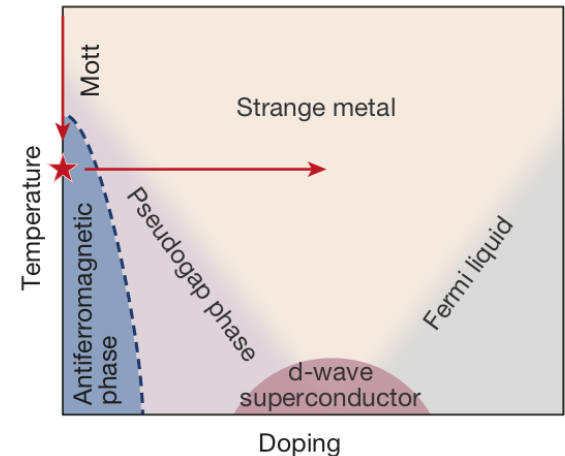
U tunable by B-field!!

The simplest model to describe interplay conductivity / interactions



YBa2Cu3O7
 $d \sim 0.4 \text{ nm}$
 $T_c = 92 \text{ K}$

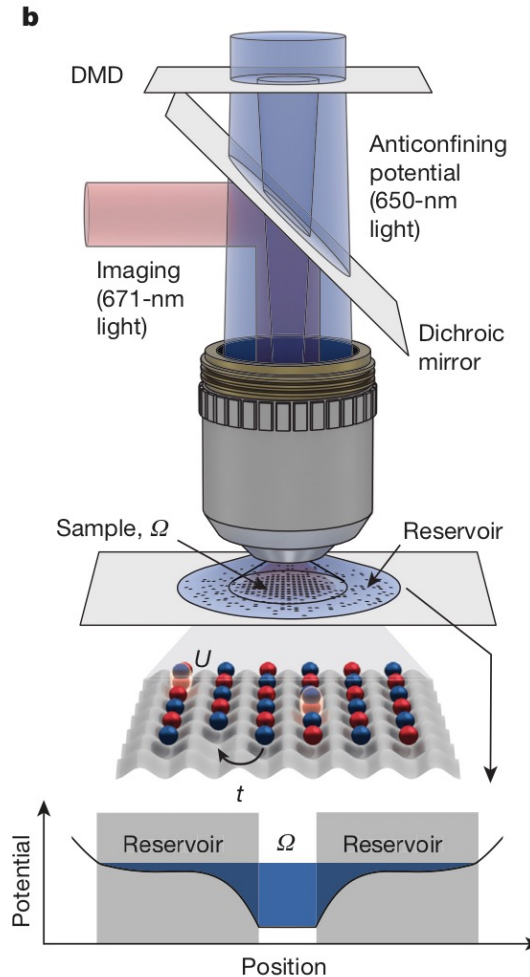
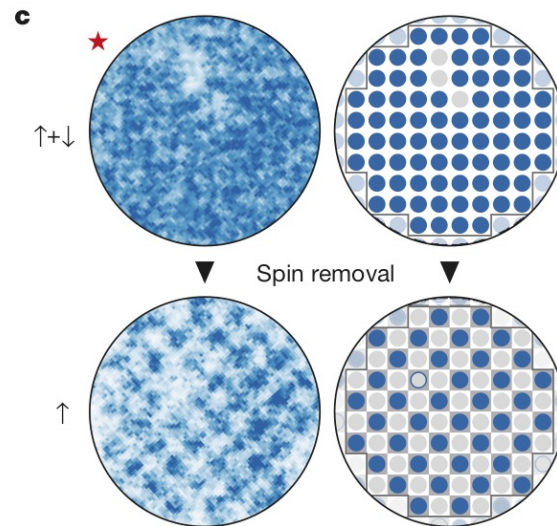
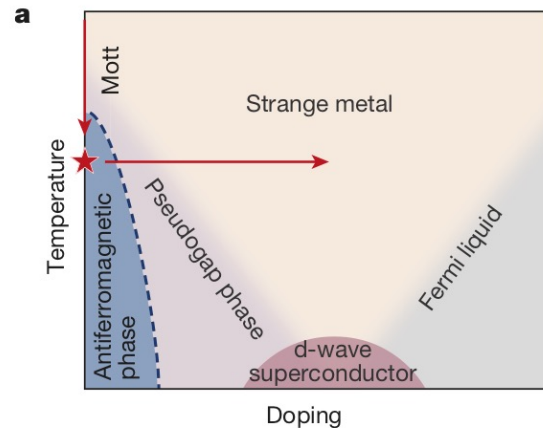
YET: not solved!



Believed to explain high- T_c superconductivity

Accessing the AF phase using a quantum simulator

Greiner, Nature 2017

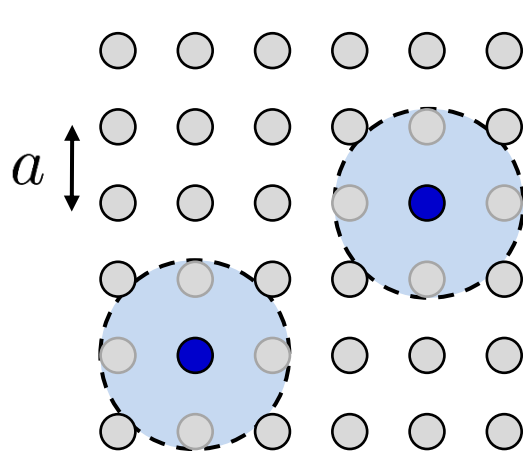


Also: Bloch (MPQ), M. Koehl (Bonn), Kuhr (Glasgow), Zwierlein (MIT), Thywissen (Toronto)...

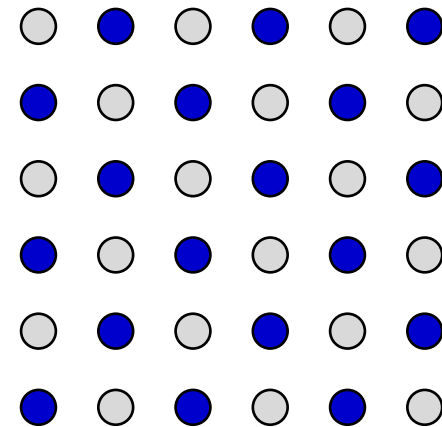
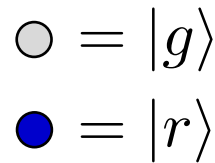
2D Ising anti-ferromagnet on a square

Nearest neighb. interaction

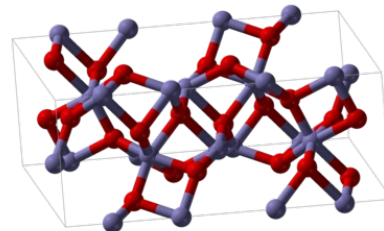
Anti-ferromagnetic ground state



$$\frac{C_6}{a^6} \sim \Omega$$



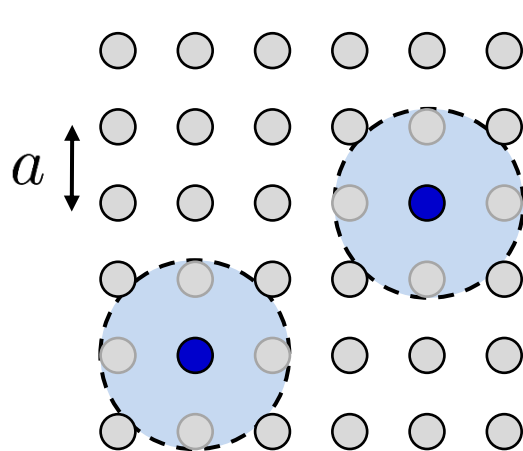
Ex of antiferromagnets: MnO, FeO, CoO, NiO, FeCl₂...



2D Ising anti-ferromagnet on a square

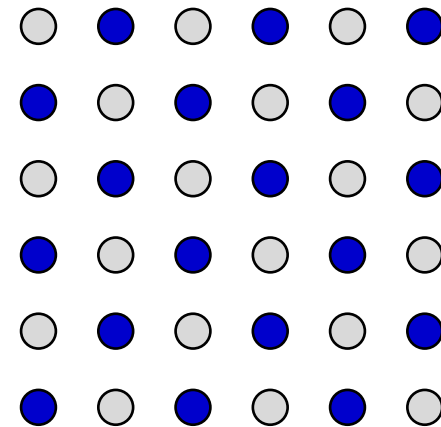
Nearest neighb. interaction

Anti-ferromagnetic ground state

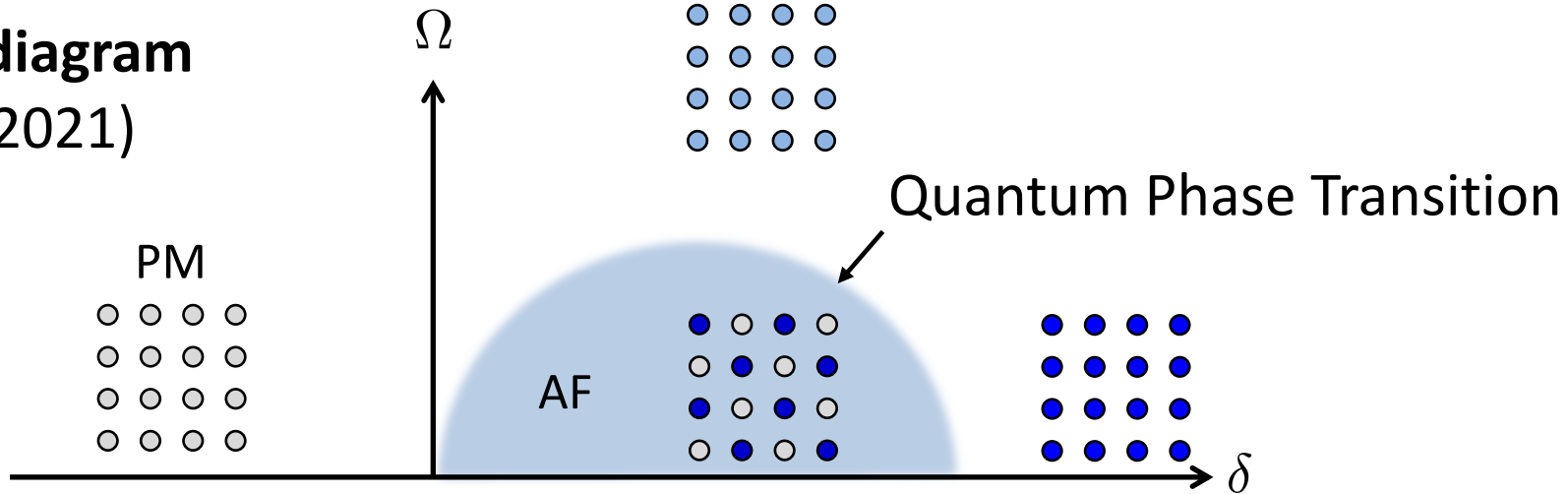


$$\frac{C_6}{a^6} \sim \Omega$$

○ = $|g\rangle$
● = $|r\rangle$



2D phase diagram
(1970 - 2021)



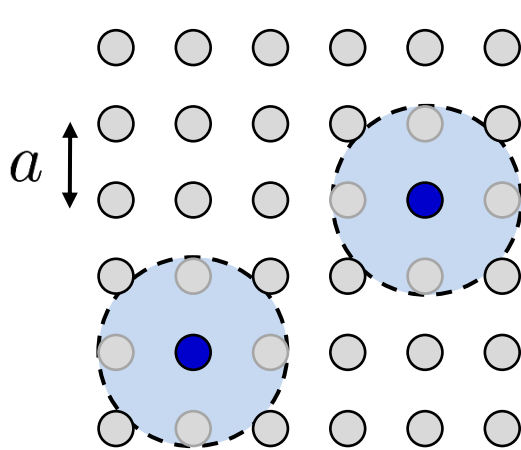
Known by Quantum Monte-Carlo

Never measured in 2D...!!!

2D Ising anti-ferromagnet on a square

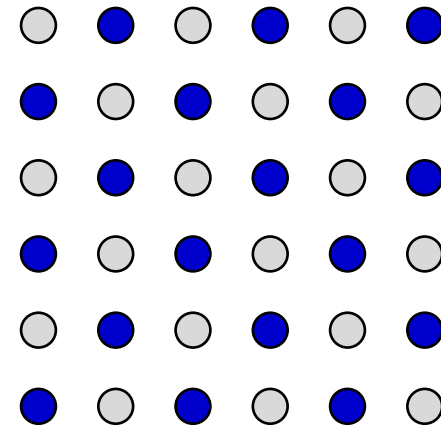
Nearest neighb. interaction

Anti-ferromagnetic ground state

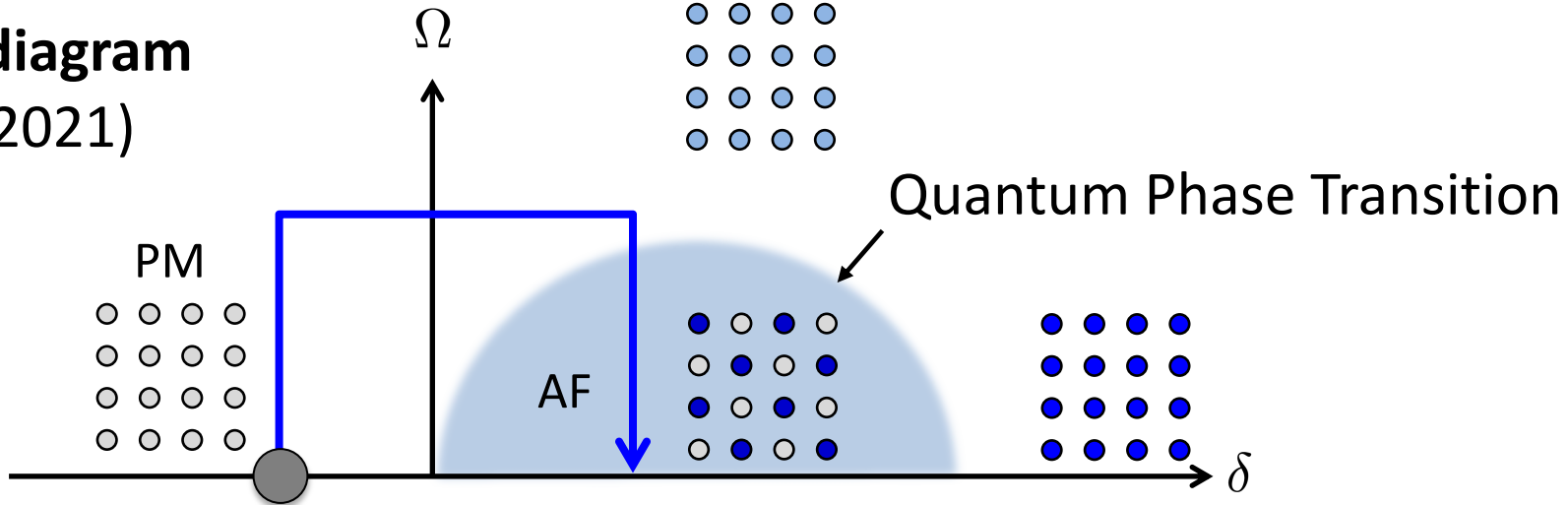


$$\frac{C_6}{a^6} \sim \Omega$$

$$\begin{aligned} \circ &= |g\rangle \\ \bullet &= |r\rangle \end{aligned}$$



2D phase diagram
(1970 - 2021)



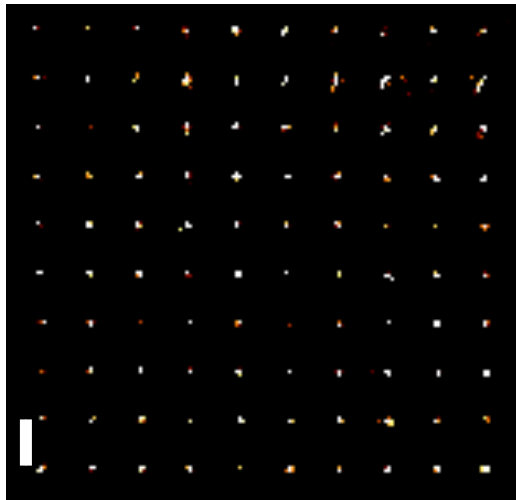
$$H = \sum_i \left(\frac{\hbar\Omega(t)}{2} \sigma_x^i - \hbar\delta(t) \hat{n}_i \right) + \sum_{i<j} \frac{C_6}{R_{ij}^6} \hat{n}_i \hat{n}_j$$

Preparation of a 2D Ising anti-ferromagnet on a square

10 × 10 square array



Scholl et al. Nature (2021)

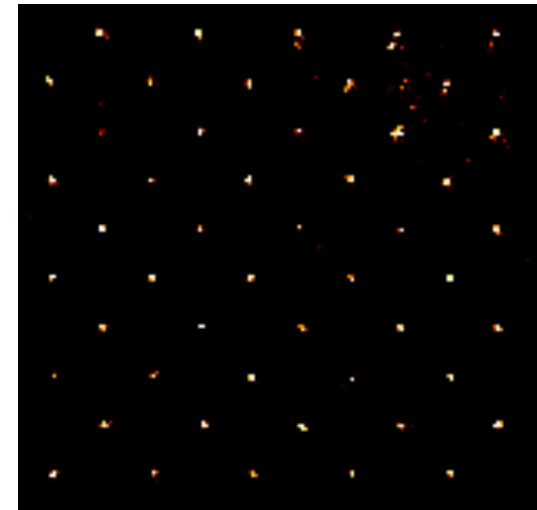
10 μm



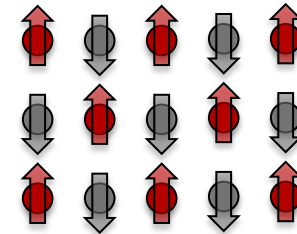
$\Omega(t), \delta(t)$

sweep

 = $|g\rangle$ bright
 = $|r\rangle$ dark



Perfect AF (Néel) ordering!



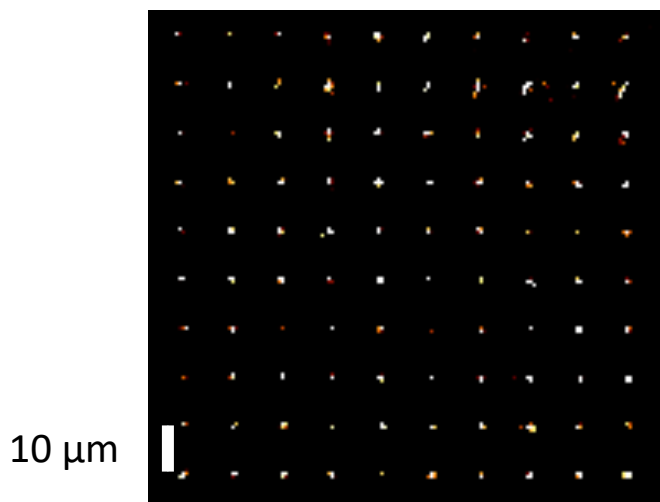
1D: Pohl PRL 2010; Bloch Science 2015; Lukin Nature 2017, 2019;

2D: Lienhard PRX 2018, Bakr PRX 2018; Lukin Nature 2021

Preparation of a 2D Ising anti-ferromagnet on a square

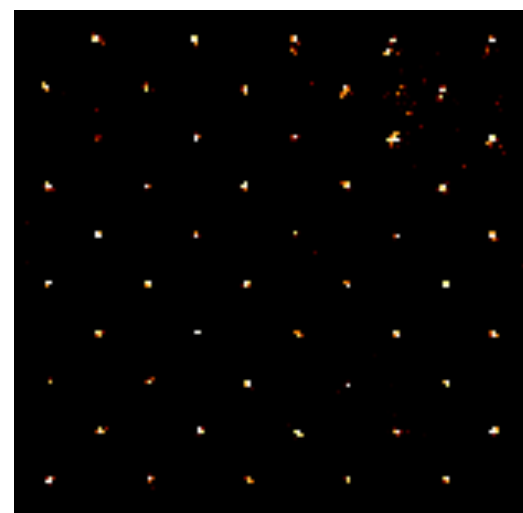
Scholl et al. Nature (2021)

10 × 10 square array

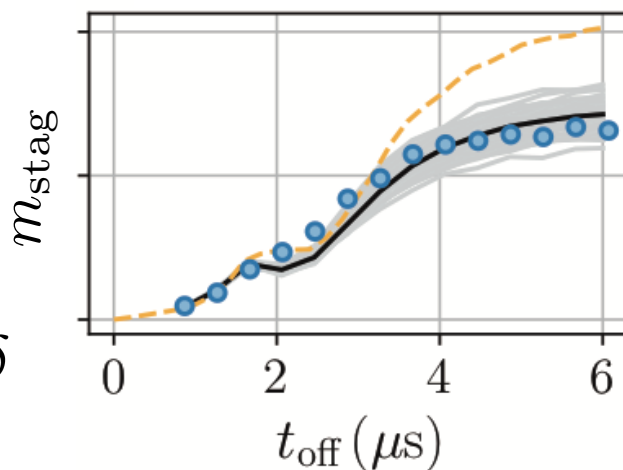
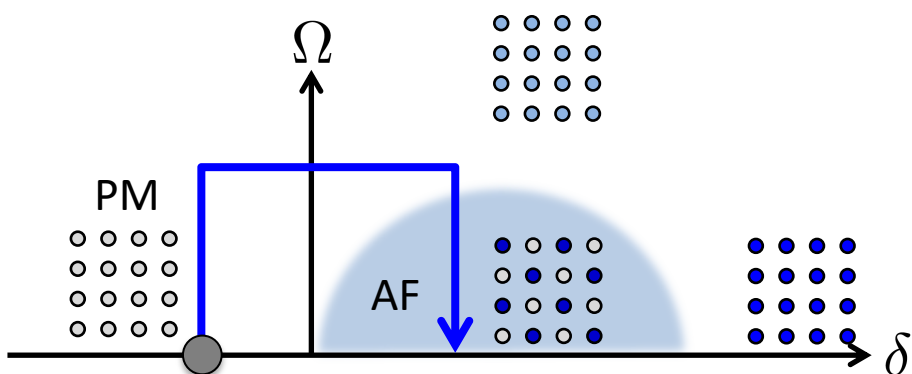


$\Omega(t), \delta(t)$

sweep



Staggered magnetization: $m_{\text{stag}} = \langle |n_A - n_B| \rangle$



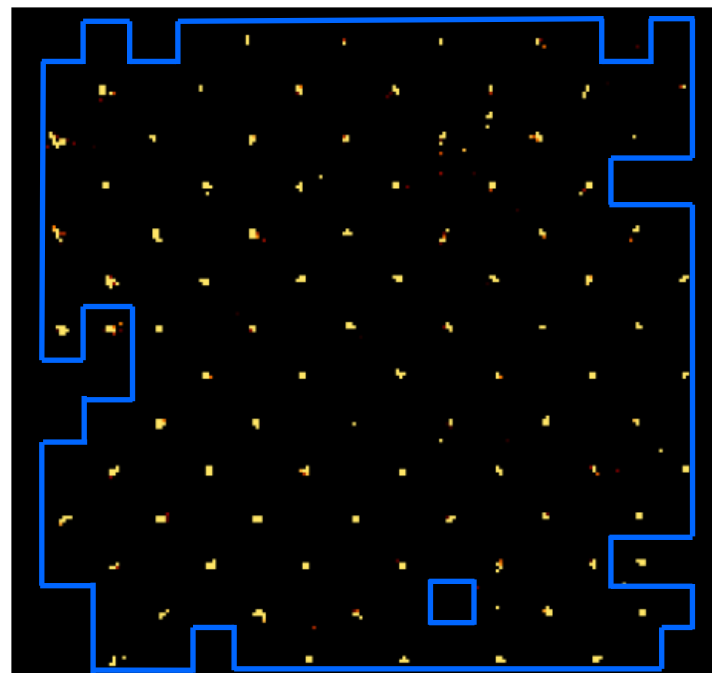
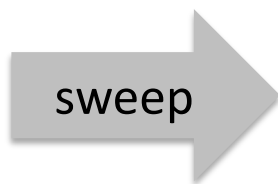
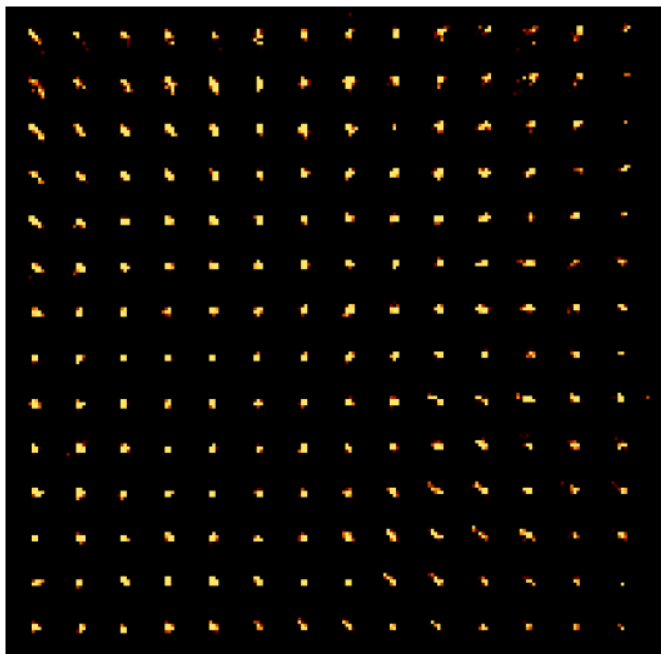
Accurate
MPS limited
to 10 x 10
(14 days!!)

Including experimental imperfections: $U_{ij}, \Omega_i, \delta_i$, real ramp...

Preparation of a 2D Ising anti-ferromagnet on a square

Scholl et *al.* Nature (2021)

14x14 square array

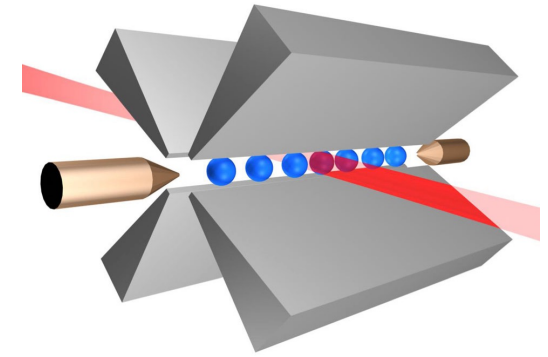
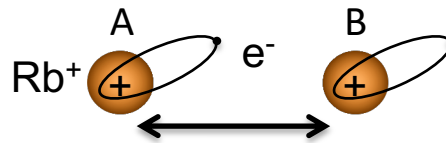
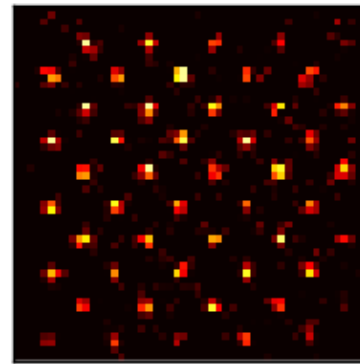
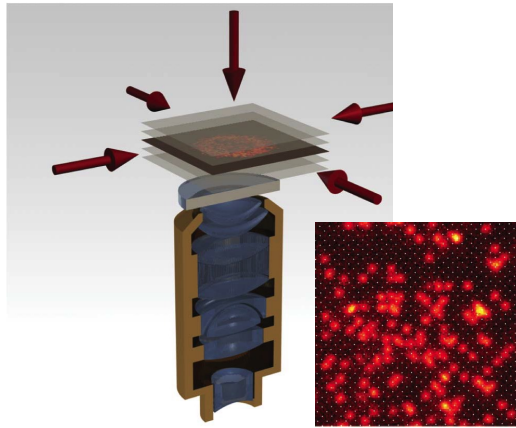


182-atom antiferromagnetic cluster!

2022: Data beyond $N > 100$ to test of new tensor network methods...!!

Conclusion

Useful quantum simulators already exist, and challenge theory!!



A very active field internationally

Programmable Atomic Large-Scale Quantum Simulation
(10 groups + 5 industries, coord.: A. Browaeys & I. Bloch)



Startups recently created to develop industry graded simulators



Applications: scientific computing, optimization in finance and industry...

Conclusion: Quantum Simulators can do more...!

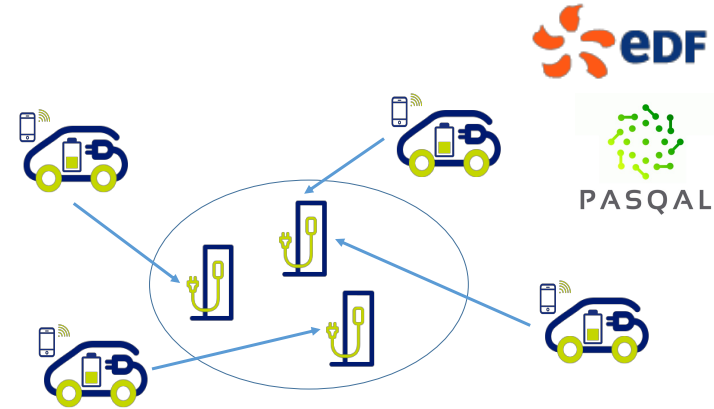
Quant. sim. = machine to prepare quantum states

Optimization problems

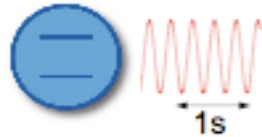
Mapped onto Ising model

$$H = \sum_i h_i n_i + \sum_{i < j} J_{ij} n_i n_j, \quad n_i = 0, 1$$

Solution = ground-state

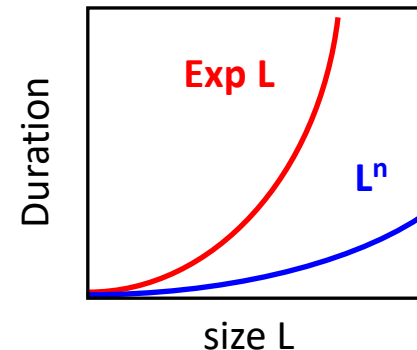


Quantum metrology



$$\frac{\Delta\nu}{\nu_0} \propto \frac{1}{\sqrt{N}} \xrightarrow{\text{Entangled states}} \frac{\Delta\nu}{\nu_0} \propto \frac{1}{N}$$

Quantum computer (long-term...)

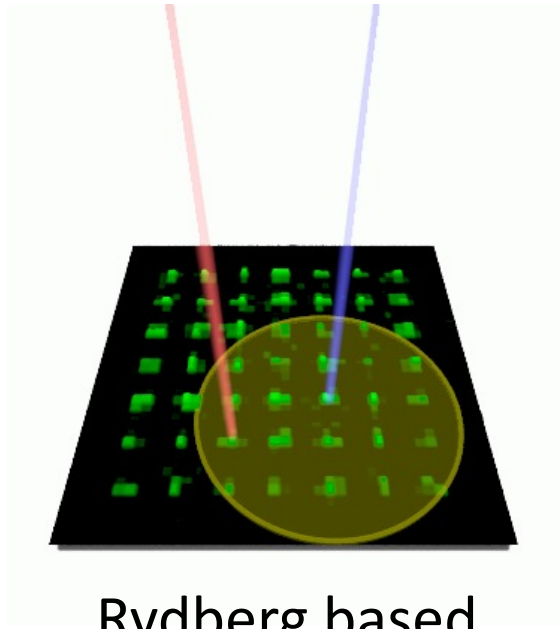


Outlook: towards Industrial Simulators



PASQAL

www.pasqal.io



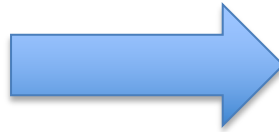
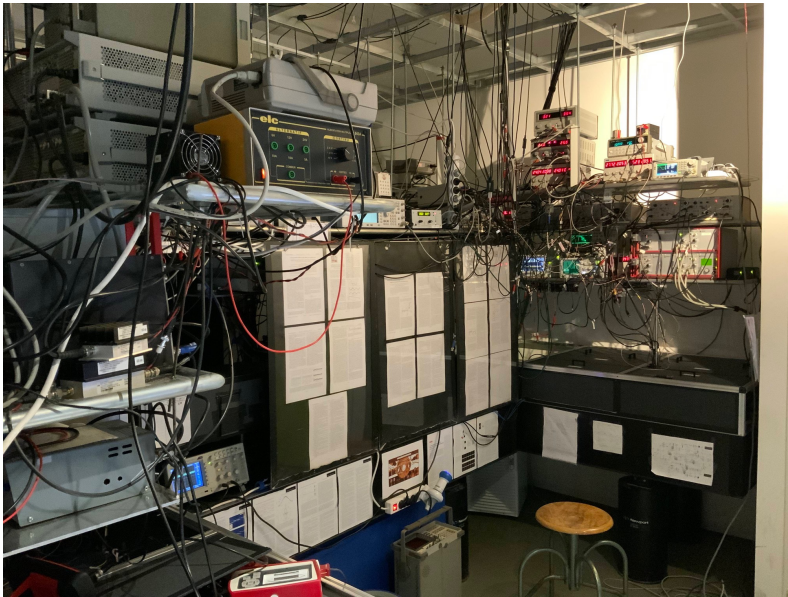
Rydberg based
technology

- Founded april 2019 (90 staff sept. 2022)
- **Applications:** scientific computing, optimisation in finance, industries...
- Software et hardware

USA



From a lab experiment to an industrial machine...



Jobs opportunities !!

References:

“Quantum Simulation”, I.M. Georgescu *et al.*, Rev. Mod. Phys. **49**, 153 (2014)

“Quantum Simulation with ultra-cold atoms in optical lattices”, C. Gross and I. Bloch., Science **357**, 995 (2017)

“Many-Body Physics with individually controlled Rydberg Atoms”, A. Browaeys and T. Lahaye, Nat. Phys. **16**, 132 (2020)

“Programmable quantum simulation of spin systems with trapped ions”, C. Monroe *et al.*, arXiv:1912.07845

