

A 3D rendering of a photonic crystal slab. The slab is blue and has a sawtooth-like top surface. A red beam of light is directed at the slab, and three gold-colored spheres representing atoms are positioned on the surface. The background is a light gray gradient.

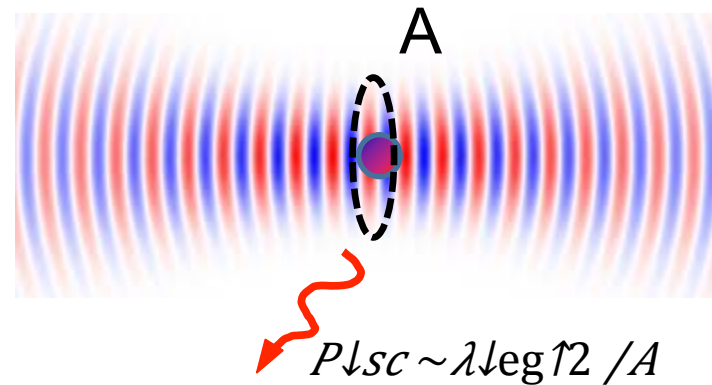
Enhancing atom-light interactions through subradiant dissipation

Darrick Chang
ICFO – The Institute of Photonic Sciences
Barcelona, Spain

KITP discussion
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Atom-photon interactions

- Problem: single atoms and photons don't like to talk to each other



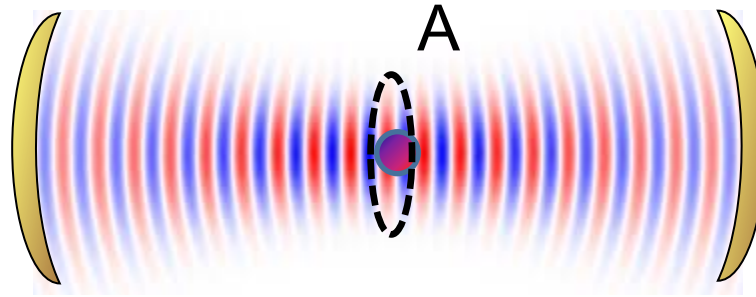
diffraction limit

$$A > \lambda^2$$

Atom-photon interactions

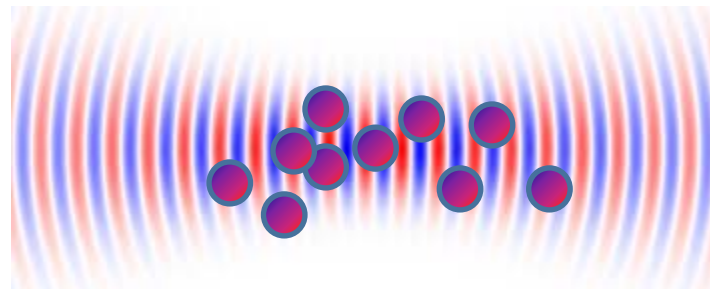
- Some fixes:

- Cavity QED



$$P_{sc} \sim \lambda \log^2 / A N_{\text{bounces}} \sim g^2 / \kappa \Gamma \sim \text{Cooperativity}$$

- Atomic ensembles



$$P_{sc} \sim \lambda \log^2 / A N_{\text{atoms}} \sim \text{Optical depth}$$

A fundamental limit

- Cooperativity or optical depth widely viewed as a fundamental limit to applications of atom-light interfaces

A quantum gate between a flying optical photon and a single trapped atom

Andreas Reiserer¹, Norbert Kalb¹, Gerhard Rempe¹ & Stephan Ritter¹

In principle, the gate mechanism presented in this work is deterministic. In our experimental implementation, the photon is not back-reflected¹⁸ from the coupled system $|\uparrow^a\uparrow^b\rangle$ with a probability of 34(2)% (due to the finite cooperativity $C = \frac{g^2}{2\kappa\gamma} = 3$) and in the uncoupled

All-Optical Switch and Transistor Gated by One Stored Photon

Wenlan Chen,¹ Kristin M. Beck,¹ Robert Bücker,^{1,2} Michael Gullans,³ Mikhail D. Lukin,³ Haruka Tanji-Suzuki,^{1,3,4} Vladan Vuletić^{1*}

stood in a simple cavity QED model: One atom in state $|s\rangle$ reduces the cavity transmission (I, \mathcal{A}) by a factor of $T = (1 + \eta)^{-2}$, where η is the single-atom cooperativity (30). In the strong-coupling

Universal Approach to Optimal Photon Storage in Atomic Media

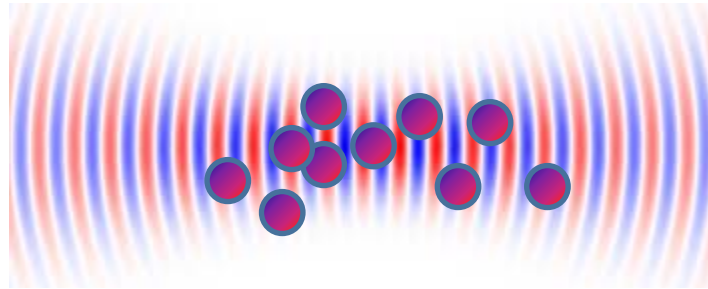
Alexey V. Gorshkov,¹ Axel André,¹ Michael Fleischhauer,² Anders S. Sørensen,³ and Mikhail D. Lukin¹

off-resonant Raman fields to photon-echo-based techniques. Furthermore, we derive an optimal control strategy for storage and retrieval of a photon wave packet of any given shape. All these approaches, when optimized, yield identical maximum efficiencies, which only depend on the optical depth of the medium.

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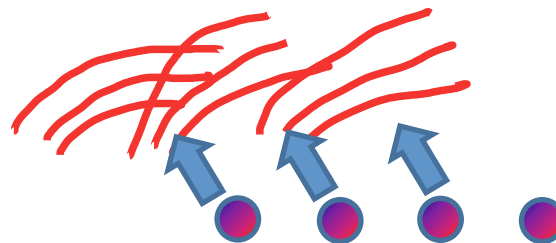
Is there a loophole?

- Atomic ensembles



$$P_{sc} \sim \lambda^4 \epsilon^2 / A N_{atoms} \sim \text{Optical depth}$$

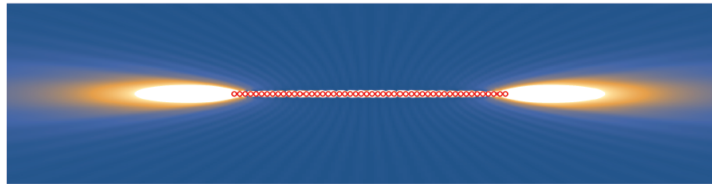
- Why this scaling?
 - Each atom has a probability of talking to the “good” mode (Gaussian) vs. the bad (free space)
 - Effectiveness is multiplied by # of attempts ($\propto N_{atoms}$)
- **Important assumption:** coupling to free space assumed to be independent
 - Cannot be true – atoms emit waves, which can interfere



Physics of super- and
sub-radiance

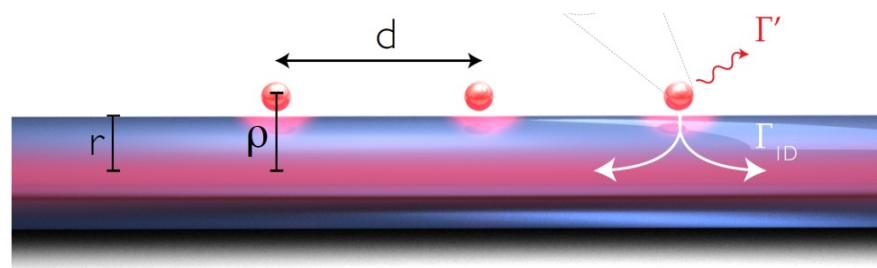
Outline

- **Big question:** can we exploit collective emission to enhance atom-light interfaces? **(Yes!)**
- Formalism to treat atom-light interfaces including subradiance
- Origin of subradiance in atomic arrays in free space



Subradiance = Guided modes

- “Selective” subradiance: atoms coupled to a nanofiber



Light-matter interactions as a spin model

Electromagnetic Green's function

- Formal definition:

$$\left[(\nabla \times \nabla \times) - \frac{\omega^2}{c^2} \epsilon(r, \omega) \right] G_{\alpha\beta}(r, r', \omega) = \delta(r - r') \cdot \mathbf{I}$$

- G describes the electric field at point r, of a (normalized) oscillating dipole at r'



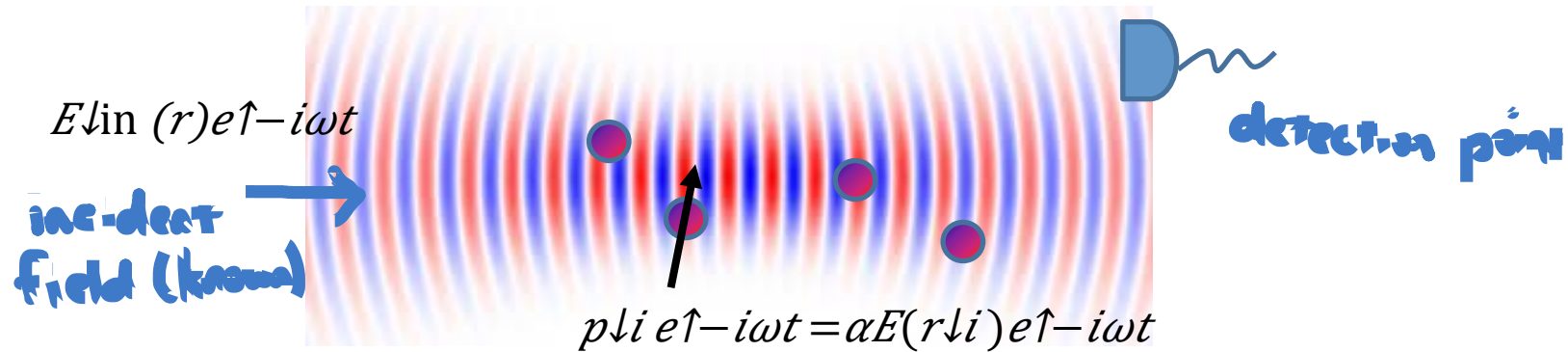
- Tensor quantity ($\alpha, \beta = x, y, z$): source dipole can have three orientations, and the electric field at r is a vector

- Simple case: free space

$$G(r, 0, \omega) = e^{ikr - i\omega t} \left[\frac{1}{r} (\hat{n} \times \hat{p}) \times \hat{n} + (3\hat{n}(\hat{n} \cdot \hat{p}) - \hat{p}) \left(\frac{1}{r^3} - \frac{ik}{r^2} \right) \right]$$

Getting rid of the light

- Classical scattering from polarizable particles



- Know the radiation pattern for a dipole
- Can calculate the total field

$$E(r, \omega) = E^{\text{lin}}(r, \omega) + \alpha \sum_i G(r, r_{li}, \omega) p_{li}(\omega)$$

Becomes convolution in time domain

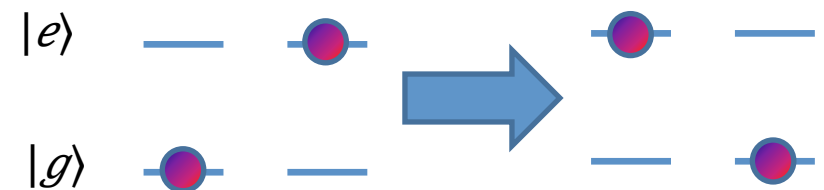
- Classical and quantum fields propagate the same way
- Generalized “input-output” equation *in time* for atoms

$$E(r, t) = E^{\text{lin}}(r, t) + (3\pi\hbar c \Gamma / d \int \omega \text{leg}) \sum_i G(r, r_{li}, \omega \text{leg}) \sigma_{lg} e^{i\omega t} \quad \text{Field encoded in atoms!}$$

What about the atoms?

- Atoms interact with fields, but fields are dependent on the atoms themselves
- Effective “spin” Hamiltonian involving atoms alone

$$H_{\text{eff}} = - \left(\frac{3\pi\hbar\Gamma}{4} \frac{c}{\omega_{eg}} \right) \sum_{i,j} G(r_j, r_i, \omega_{eg}) \sigma_{eg}^{\uparrow i} \sigma_{ge}^{\uparrow j}$$



What about the atoms?

- Atoms interact with fields, but fields are dependent on the atoms themselves
- Effective “spin” Hamiltonian involving atoms alone

$$H_{\text{eff}} = - (3\pi\hbar\Gamma \downarrow 0 \ c/\omega \downarrow eg \) \sum_{i,j} G(r \downarrow j, r \downarrow i, \omega \downarrow eg) \sigma \downarrow eg \uparrow i \sigma \downarrow ge \uparrow j$$



- Strength depends on how field propagates from j to i
- Non-Hermitian Hamiltonian – describes both coherent interactions and (collective) spontaneous emission

A "trivial" example

- Must work for a single atom in free-space too

$$H_{\text{eff}} = - (3\pi\hbar\Gamma \downarrow 0 \ c/\omega_{\text{leg}}) \sum_{i,j} G(r_{\downarrow j}, r_{\downarrow i}, \omega_{\text{leg}}) \sigma_{\text{leg}}^{\uparrow i} \sigma_{\text{leg}}^{\uparrow j}$$



$$H_{\text{eff}} = - (3\pi\hbar\Gamma \downarrow 0 \ c/\omega_{\text{leg}}) G(r_{\downarrow \text{atom}}, r_{\downarrow \text{atom}}, \omega_{\text{leg}}) \sigma_{\text{leg}}$$



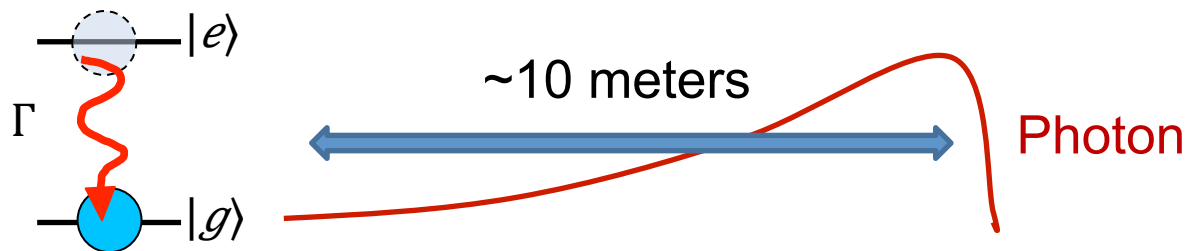
$$H_{\text{eff}} = - (3\pi\hbar\Gamma \downarrow 0 \ c/\omega_{\text{leg}}) i\omega_{\text{leg}} / 6\pi c \sigma_{\text{leg}}^{\uparrow} = -i\hbar\Gamma \downarrow 0 / 2 \sigma_{\text{leg}}$$

- Describes single-atom spontaneous emission at rate $\Gamma \downarrow 0$!

Spin model

$$H_{\text{eff}} = -\mu \sum_i d_i \sum_j \omega_{eg} \sum_{\sigma} G(r_j, r_i, \omega_{eg}) \sigma_{eg}^{\dagger i} \sigma_{ge}^{\dagger j}$$

- Limits of validity
 - No strong coupling effects (e.g. vacuum Rabi oscillations)
 - Ignores time retardation



- Equally captures **any system** of atoms interacting with light, and treats them on equal footing
 - Cavity QED, free-space atomic ensembles, nanophotonic systems, ...

1D chain of atoms in free space

Green's function for 1D atomic chain

- Spin model in general:

$$H_{\text{eff}} = -\mu_0 \hbar \omega \sum_{i,j} G(r_{ij}, r_{li}, \omega) \sigma_{i\uparrow} \sigma_{j\uparrow}$$

- Simple case: atoms in free space

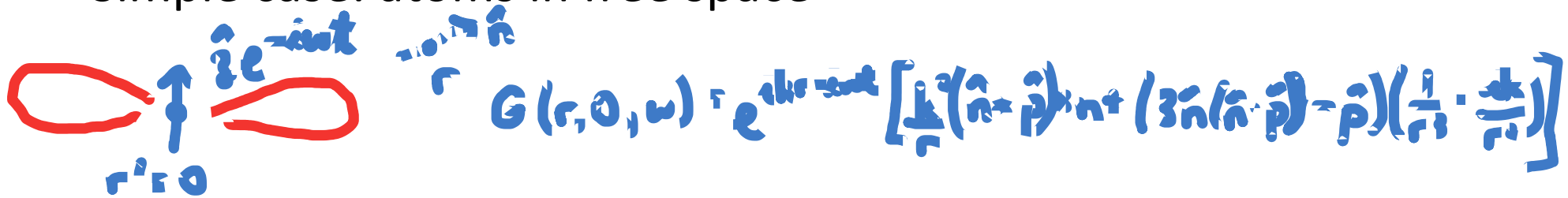
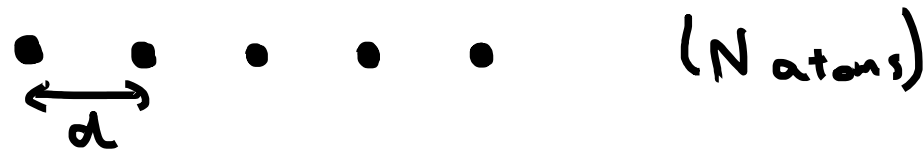


Diagram showing a dipole in free space with a red loop and an upward arrow, and a handwritten equation for the Green's function $G(r, 0, \omega) = e^{ikr} \left[\frac{1}{r} (\hat{n} \cdot \hat{p})^2 + (3\hat{n}(\hat{n} \cdot \hat{p}) - \hat{p}) \cdot \left(\frac{1}{r^3} - \frac{ik}{r^2} \right) \right]$

- Must specify a geometry: 1D chain of atoms (but living in 3D)



- Considering just one excitation, there are N possible states

$$|eggg\dots\rangle, |gegg\dots\rangle, |ggeg\dots\rangle, \dots$$

- Hamiltonian becomes $N \times N$ matrix in this subspace, $H_{ij} \propto G(r_{ij}, r_{li})$

Exact diagonalization

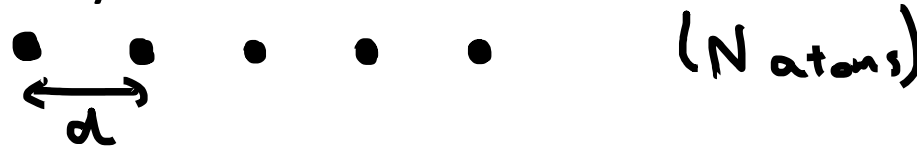
- Numerically diagonalize H
 - In general, each eigenstate is a superposition of excitations sitting on different atoms
 - Eigenvalues give the frequency shift and decay rate of each collective state

$$(j=1, \dots, N) \quad \omega_j = \underbrace{(\text{Re } \omega_j)}_{\text{real energy shift}} + i \underbrace{(\text{Im } \omega_j)}_{\text{decay rate}}$$

- Energies and decay rates modified from single-atom values, $\omega \downarrow eg$ and $\Gamma \downarrow 0$, due to photon-mediated interactions

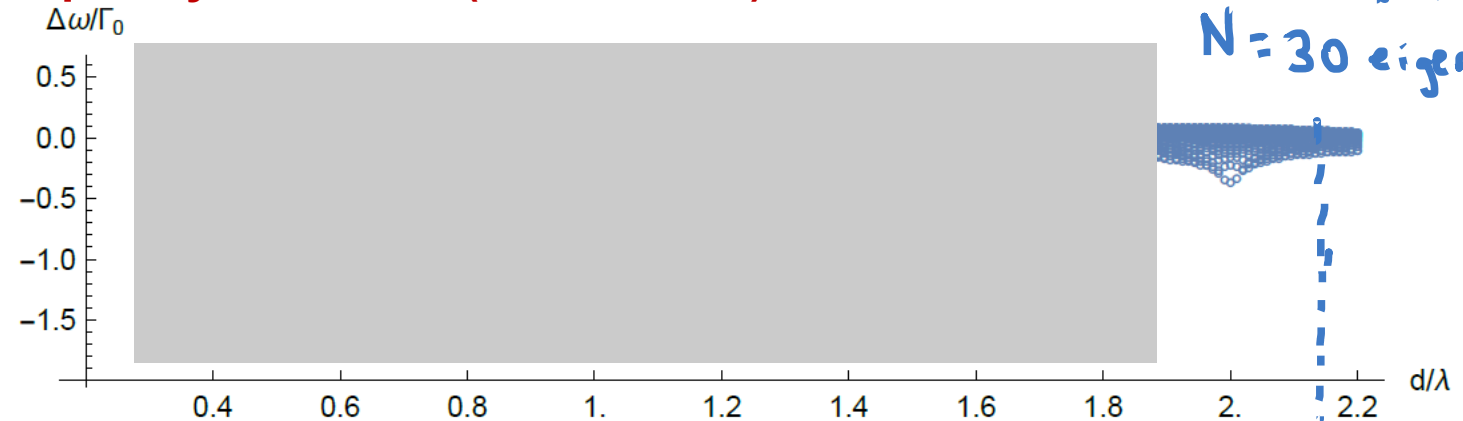
Energy shifts and decay rates

- Shifts and decay rates vs. lattice constant



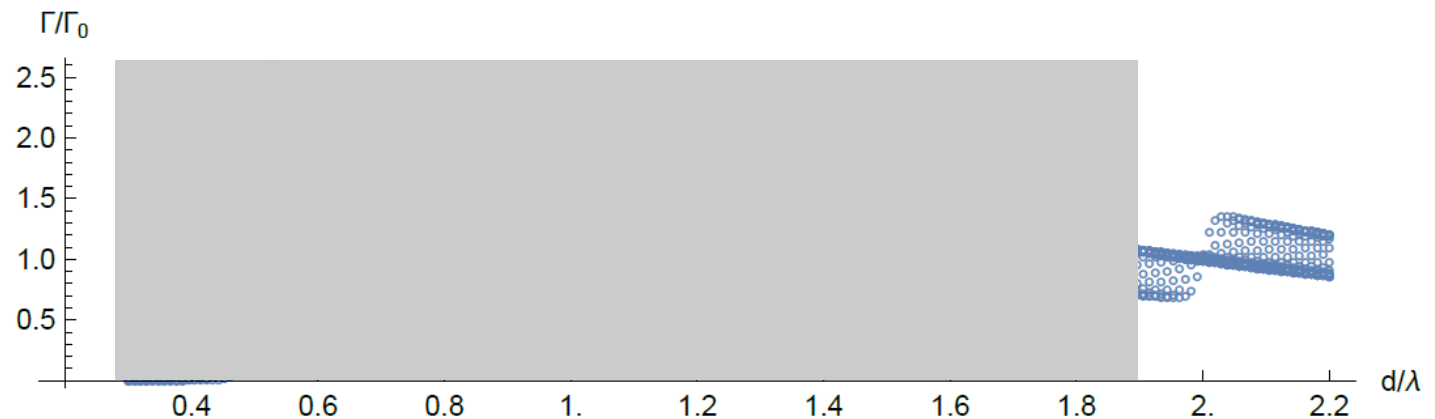
Frequency shift vs. d (N=30 atoms)

Shift normalized by free-space linewidth ($\Delta\omega_j/\Gamma_0$)



Decay rate vs. d (N=30 atoms)

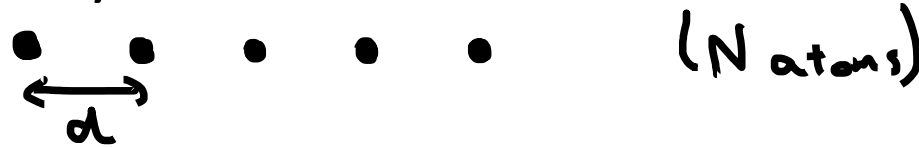
Decay rate normalized by free-space linewidth (Γ_j/Γ_0)



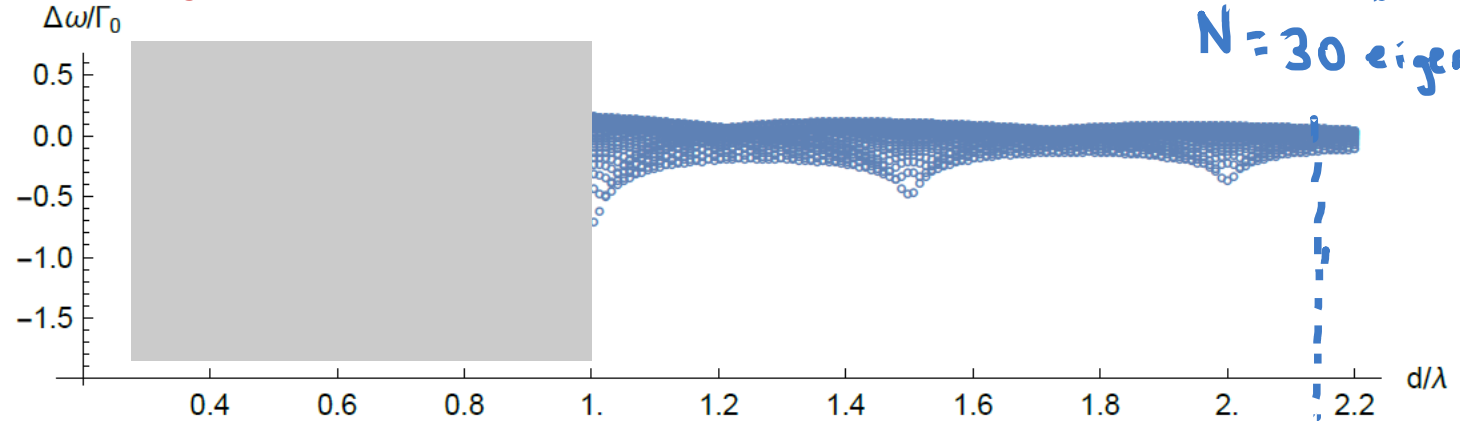
- Minimal effect of interactions at large distances

Energy shifts and decay rates

- Shifts and decay rates vs. lattice constant

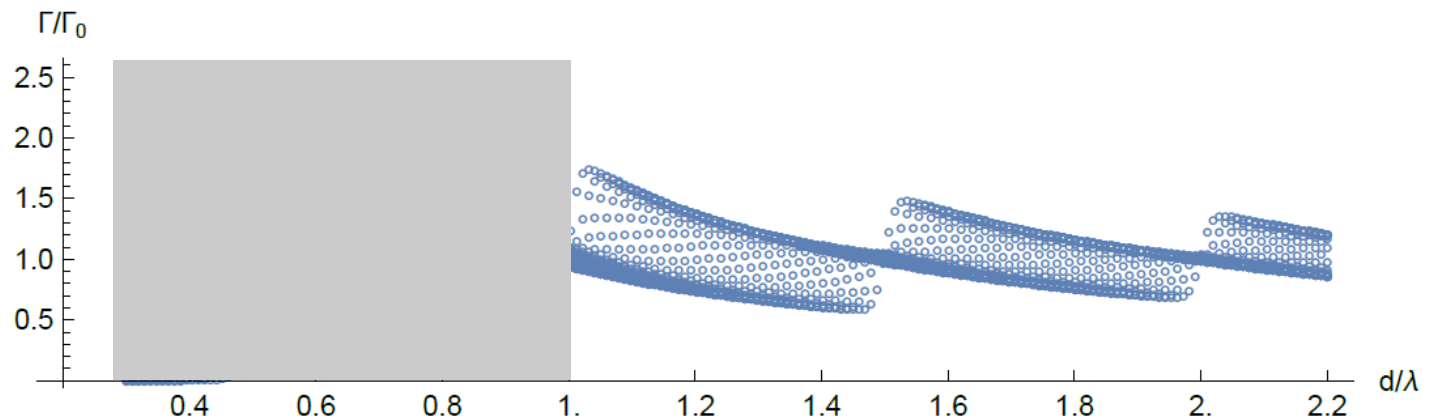


Frequency shift vs. d (N=30 atoms)



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Decay rate vs. d (N=30 atoms)

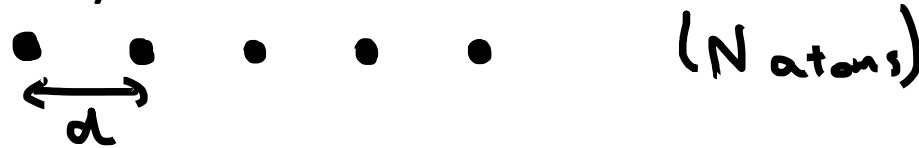


Decay rate normalized by free-space linewidth (Γ_j/Γ_0)

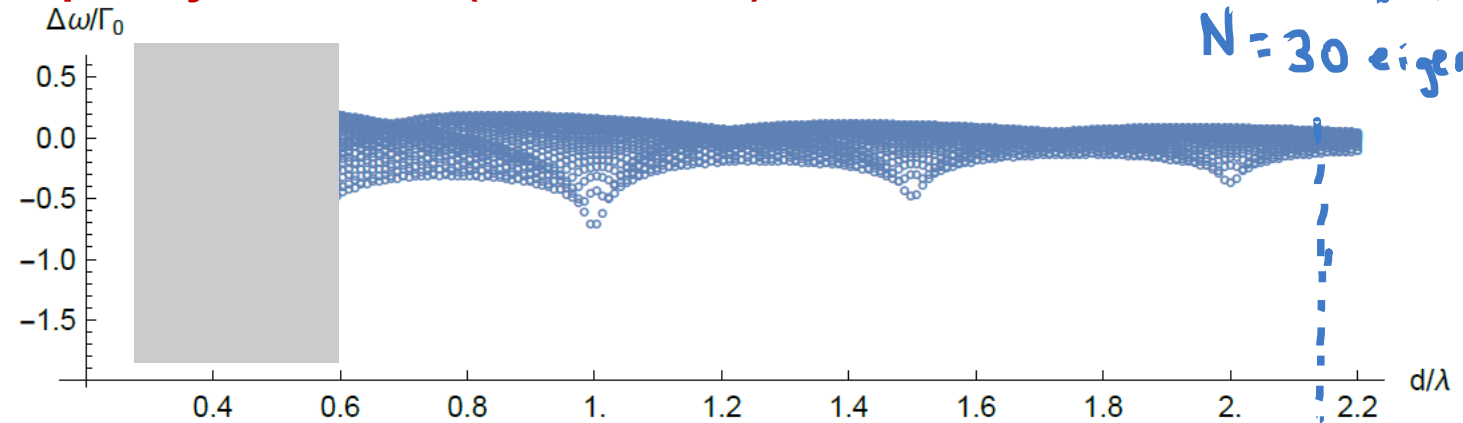
- Noticeable corrections at $d \sim \lambda$ (>50% modification of decay)

Energy shifts and decay rates

- Shifts and decay rates vs. lattice constant

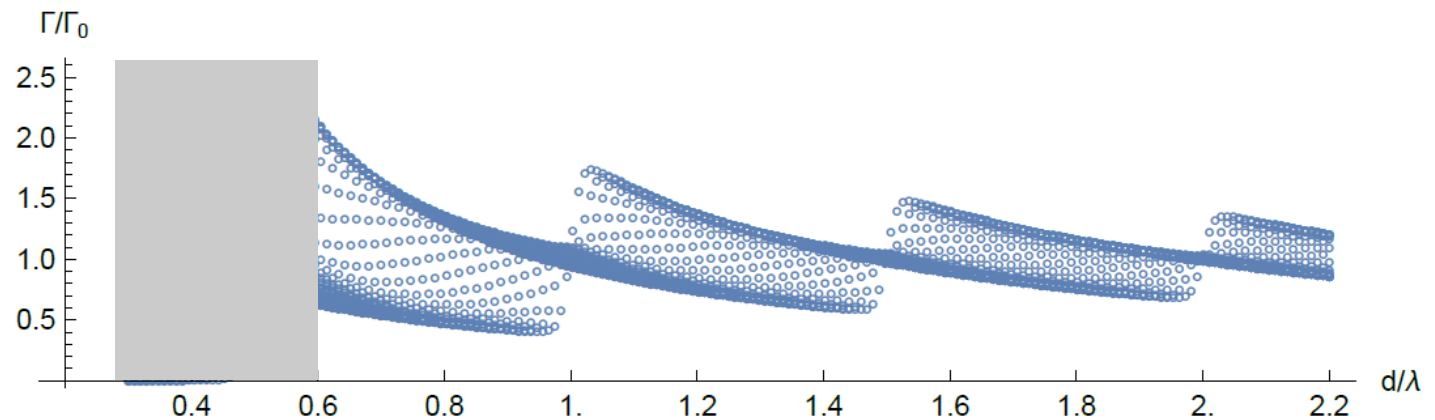


Frequency shift vs. d ($N=30$ atoms)



Shift normalized by
 free-space linewidth ($\Delta\omega_j/\Gamma_0$)

Decay rate vs. d ($N=30$ atoms)

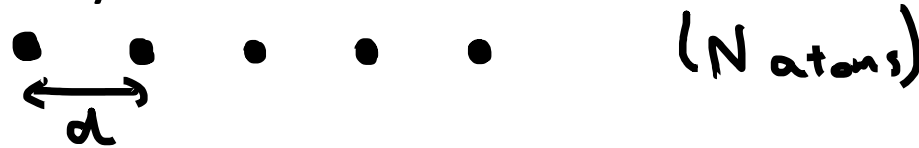


Decay rate normalized
 by free-space linewidth
 (Γ_j/Γ_0)

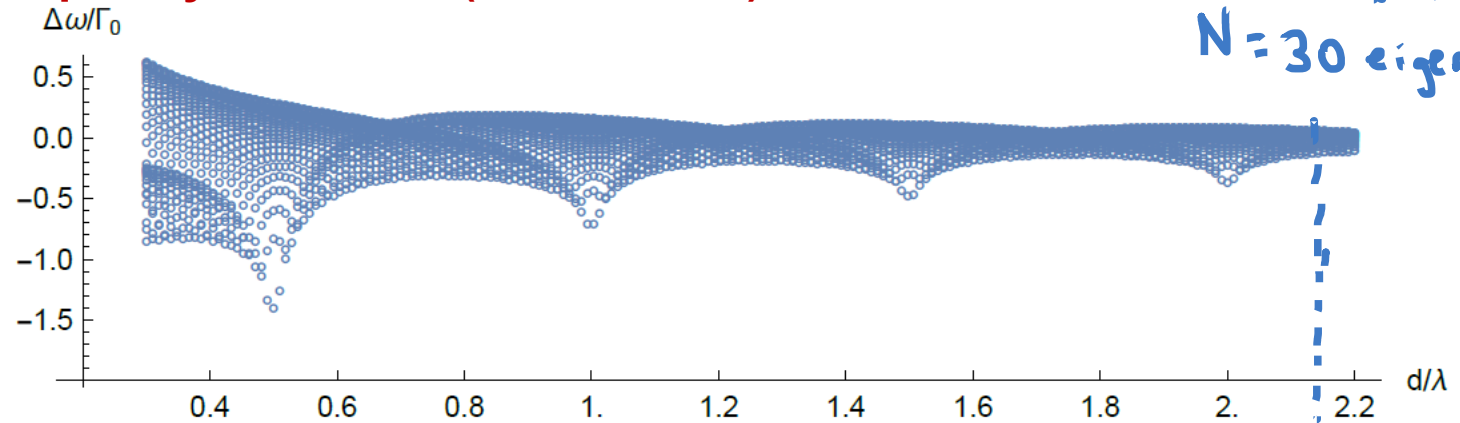
- One state doubles in decay rate as $d \sim \lambda/2$

Energy shifts and decay rates

- Shifts and decay rates vs. lattice constant



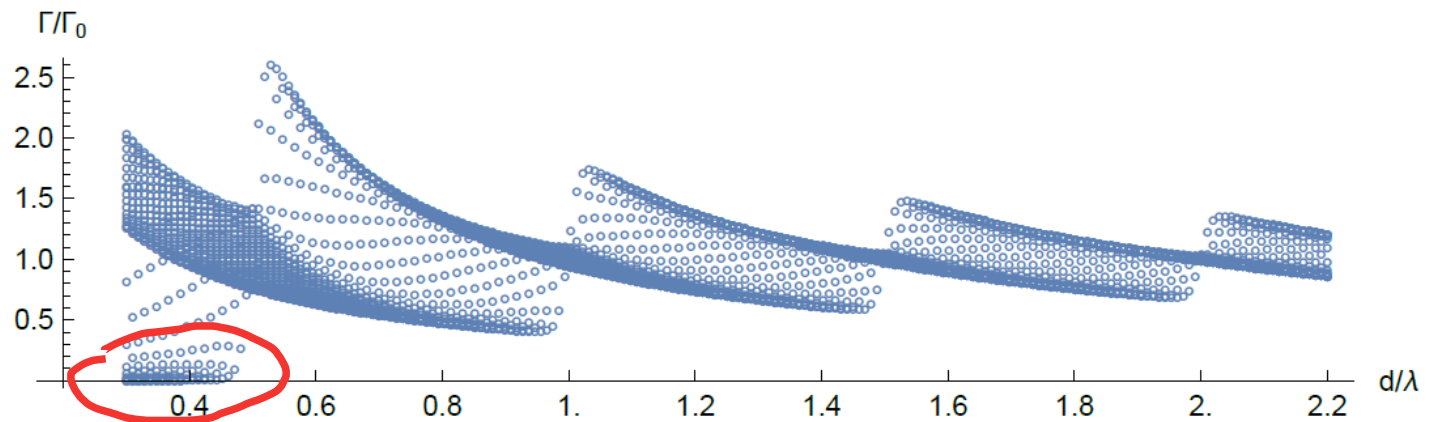
Frequency shift vs. d ($N=30$ atoms)



Shift normalized by
free-space linewidth ($\Delta\omega \downarrow j / \Gamma \downarrow 0$)

For fixed d ,
 $N=30$ eigenvalues

Decay rate vs. d ($N=30$ atoms)

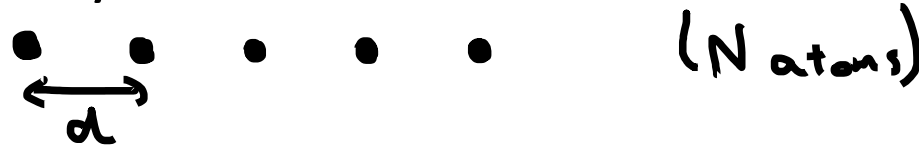


Decay rate normalized by
free-space linewidth ($\Gamma \downarrow j / \Gamma \downarrow 0$)

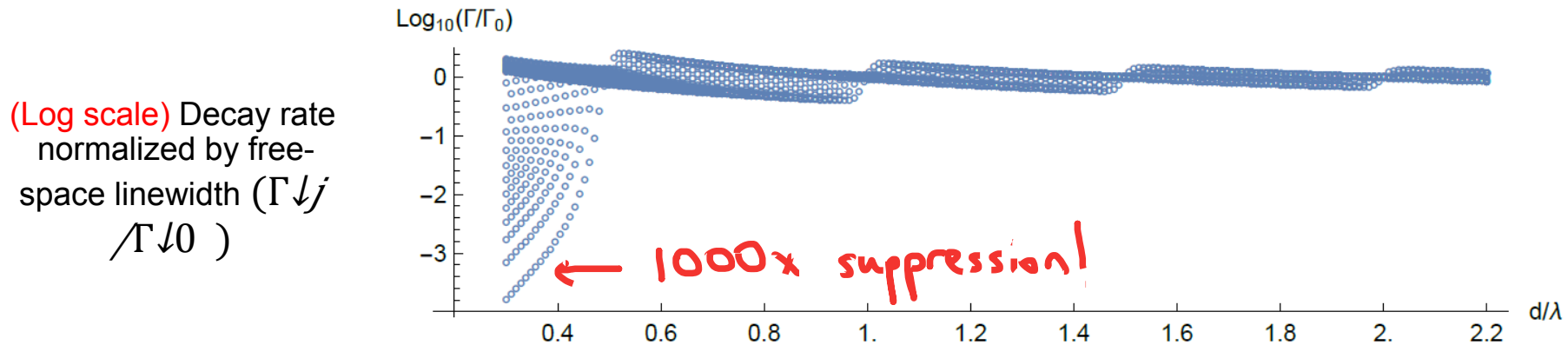
- States with **nearly zero decay rate** for $d < \lambda/2$!

Energy shifts and decay rates

- Shifts and decay rates vs. lattice constant



Decay rate (log scale) vs. d ($N=30$ atoms)



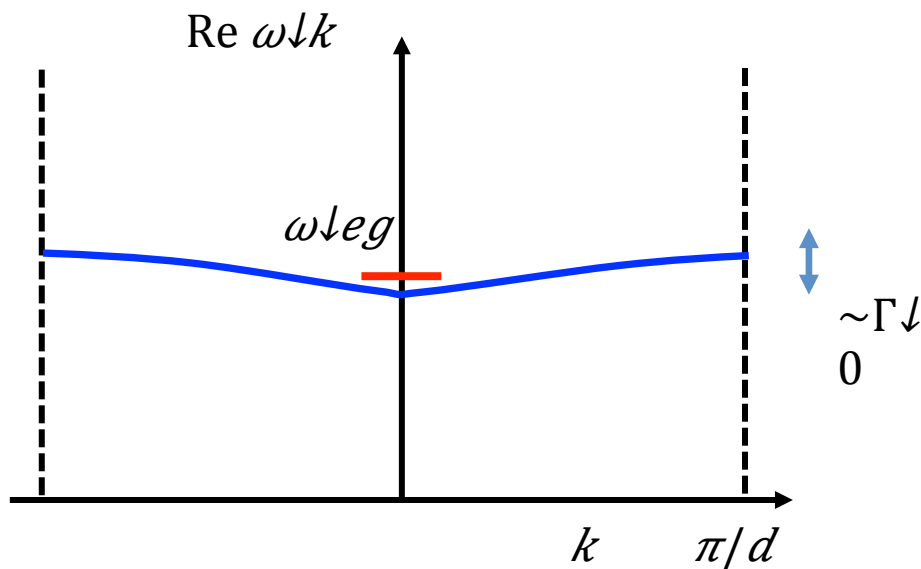
- Numerical analysis: smallest decay rates go like $\Gamma_m \sim \Gamma_0 \frac{m^2}{N^3}$
 $m=1,2,3,\dots$
- Interference of wave emission really matters at close distances!

Band structure of infinite lattice

- Infinite chain: single-excitation eigenstates are Bloch modes

$$|\psi_{\downarrow k}\rangle = \sum_j e^{ikz_{\downarrow j}} |e_{\downarrow j}\rangle$$

- Diagonalize H, represent spectrum by band structure

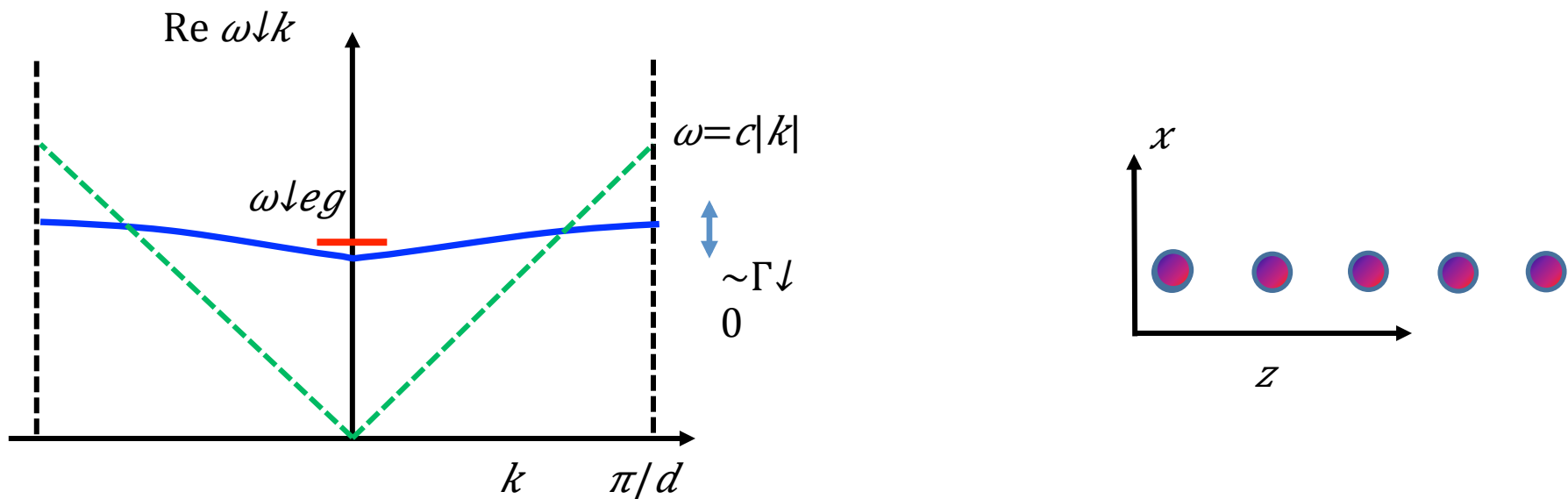


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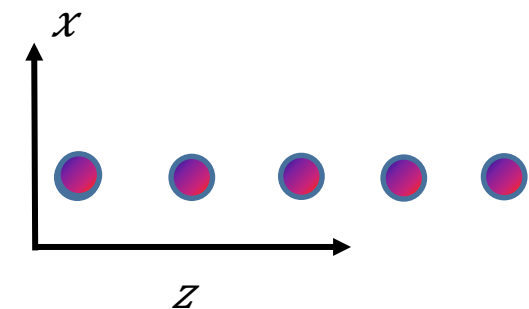
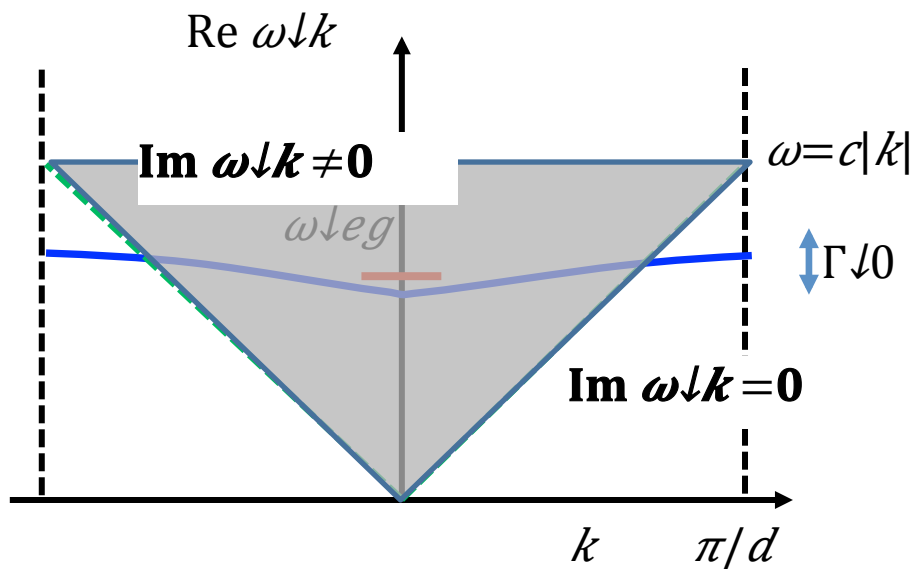
- Emitted light also has same wavevector: $E(r) \sim e^{ikz + ik_{\perp}x}$
 - Dispersion relation $k_{\parallel}^2 + k_{\perp}^2 = (\omega/c)^2$
 - $|k| > \omega/c$ implies $k_{\perp} \in \text{Im}$ (evanescent or guided mode)

Band structure of infinite lattice

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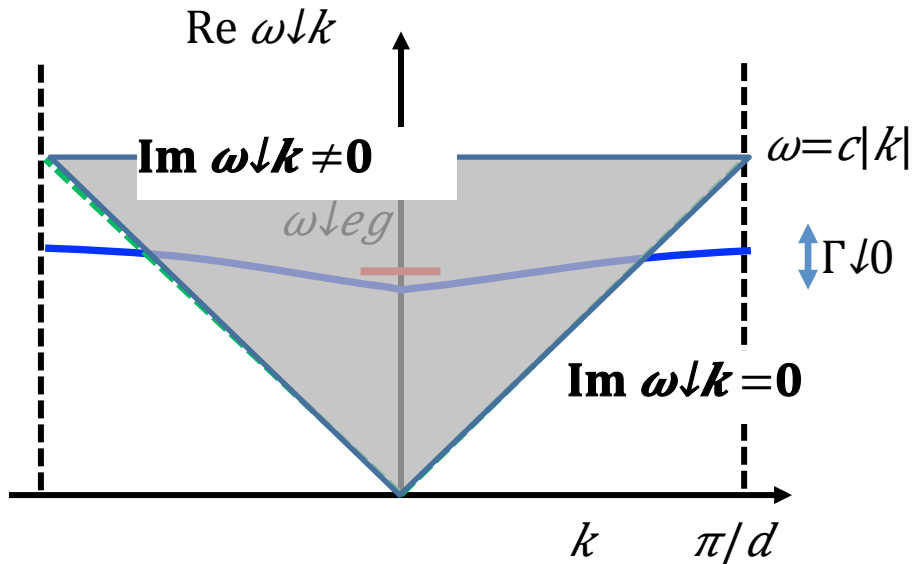
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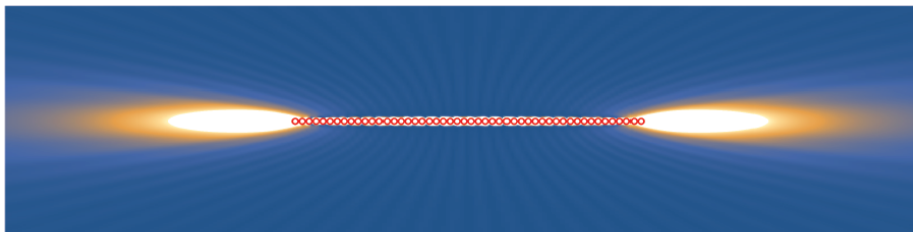
Subradiant states as guided modes

- An infinite chain supports states with *zero* decay rate, which are simply guided “fiber” modes of the chain!



- Condition: $\omega=c/|k|$ intersects Brillouin zone above $\omega\downarrow eg$
- Equivalent to $d < \lambda\downarrow eg / 2$

- Finite chain: finite decay rate $\Gamma\downarrow m \propto \Gamma\downarrow 0 \ m\uparrow 2 / N\uparrow 3$ due to end-fire emission off the “fiber” ends



Emission pattern for most sub-radiant state, N=30 atoms

Multiple excitations

- Single-excitation physics is classical
- Can we encode many excitations in sub-radiant states?

- Most sub-radiant single excitation $\Gamma \downarrow 1 \propto \Gamma \downarrow 0 / N \uparrow 3$

$$|\psi\rangle = S \downarrow 1 \uparrow \uparrow |g\rangle \uparrow \otimes N = \sum_j c_j |e \downarrow j\rangle$$

- Example: two excitations

- What if we create same excitation twice: $|\psi\rangle = (S \downarrow 1 \uparrow \uparrow) \uparrow 2 |g\rangle \uparrow \otimes N$

Eigenstate?

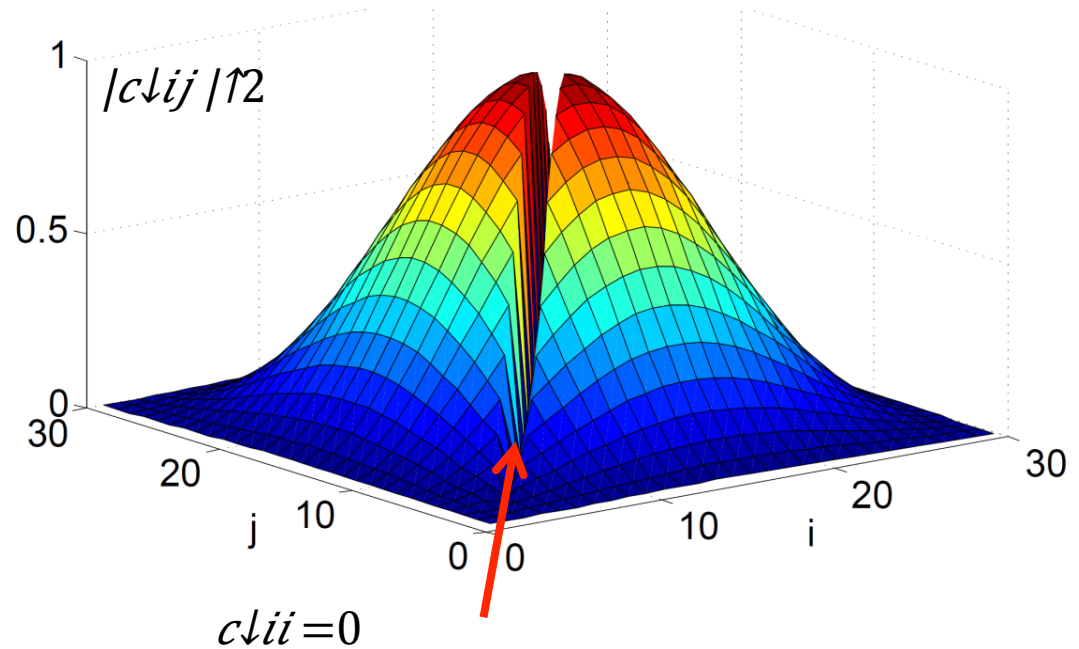
- Bosons: yes! Decay rate $2\Gamma \downarrow 1$
- Spins: no!

Two-excitation wavefunction

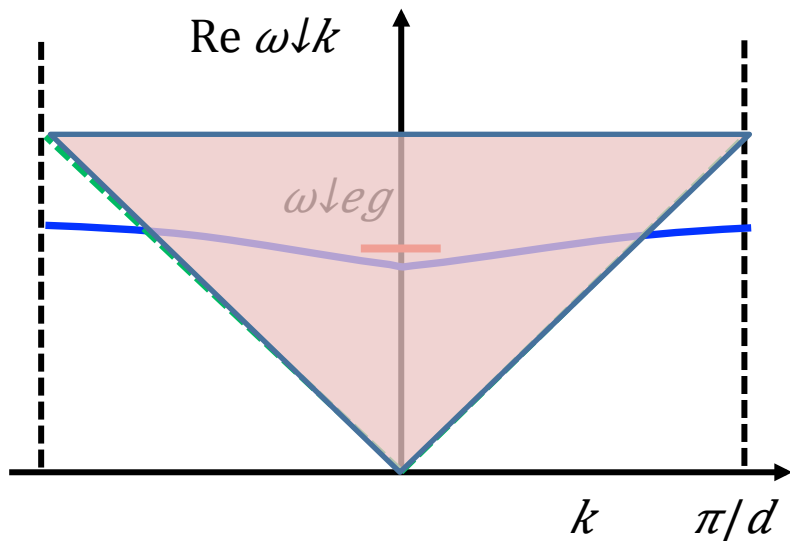
- Wave function:

$$|\psi\rangle = (S \downarrow 1 \uparrow \uparrow) \uparrow 2 |g\rangle \uparrow \otimes N$$

$$= \sum_{i,j} c_{ij} |e_i e_j\rangle$$



- State contains many momentum components within light cone



- Only weakly subradiant

$$\langle \Gamma \rangle \sim \Gamma_0 / N$$

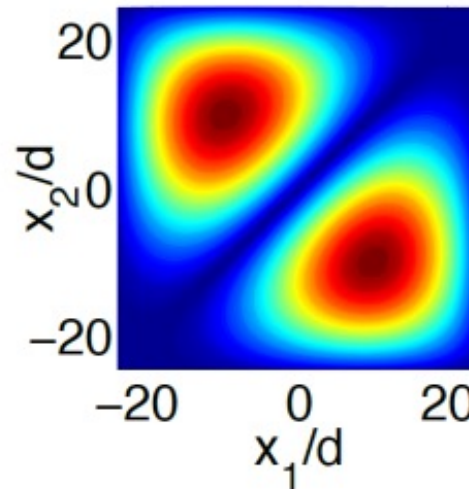
Multiple excitations

- How to construct many-excitation subradiant states?
- Single-particle eigenstates
 $\Gamma \downarrow m \propto m \uparrow^2 \Gamma \downarrow 0 / N \uparrow^3$, $|\psi \downarrow m \rangle = \sum_{j \uparrow} c \downarrow j \uparrow(m) |e \downarrow j \rangle$
- Sub-radiant states should satisfy:
 - Wavevectors beyond the light line, *and* spatially non-overlapping
- Create a *fermionized* wave function

- Example: two excitations

$$|\psi \rangle = \sum_{i \uparrow} \sum_{j \uparrow} c \downarrow ij |e \downarrow i e \downarrow j \rangle$$

$$c \downarrow ij = c \downarrow i \uparrow(1) c \downarrow j \uparrow(2) - c \downarrow i \uparrow(2) c \downarrow j \uparrow(1)$$



$$\Gamma \sim \Gamma \downarrow 1 + \Gamma \downarrow 2 \propto 1 / N \uparrow^3$$

Multiple excitations

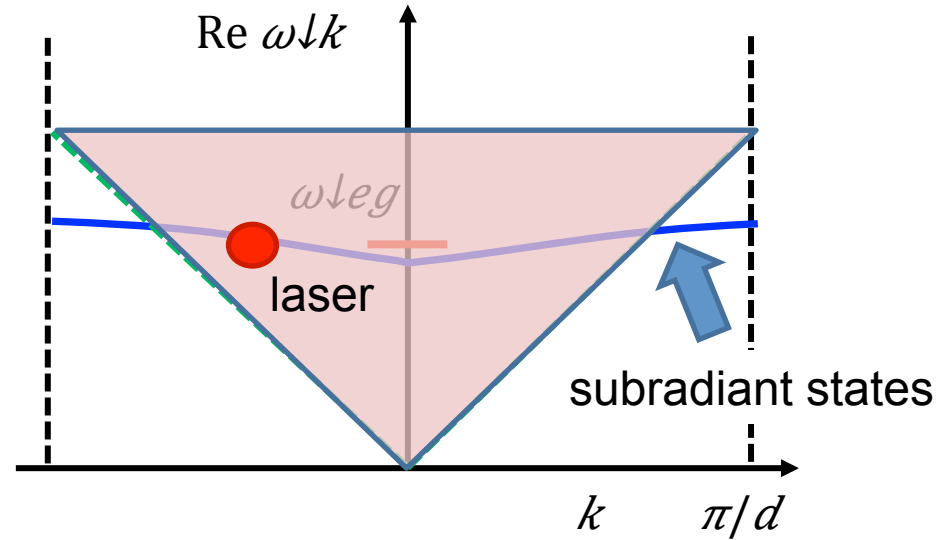
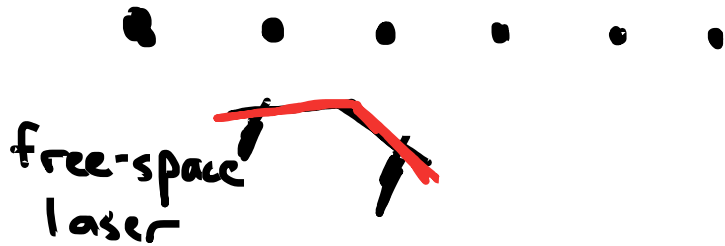
- Can extend to many-body limit
- N atoms, m excitations

$$\Gamma/\Gamma_0 \sim \sum_{j=1}^m \Gamma_j^2 / N^3 \sim (m/N)^3$$

- System can support a low density of excitations in subradiant manifold

Accessing subradiant states

- Hard to couple to subradiant states from free space

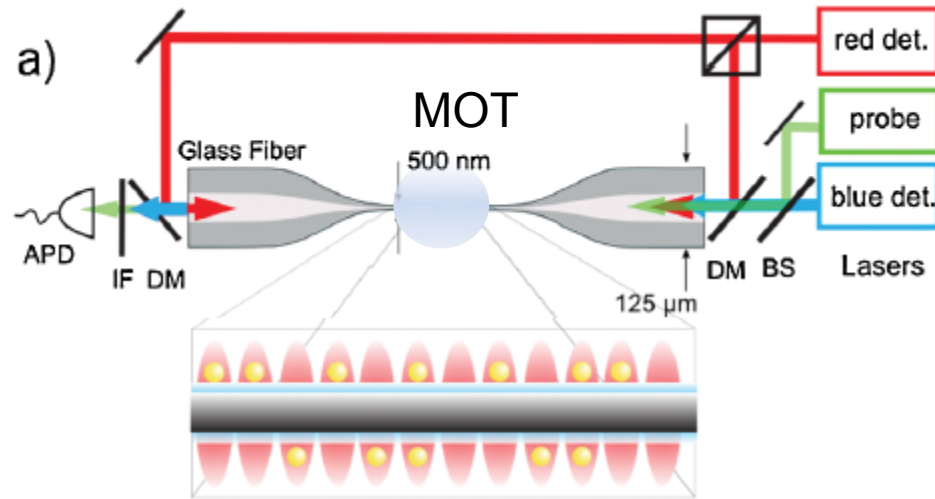


- Excite with another set of guided modes (optical nanofiber!)

Atom-nanofiber interfaces

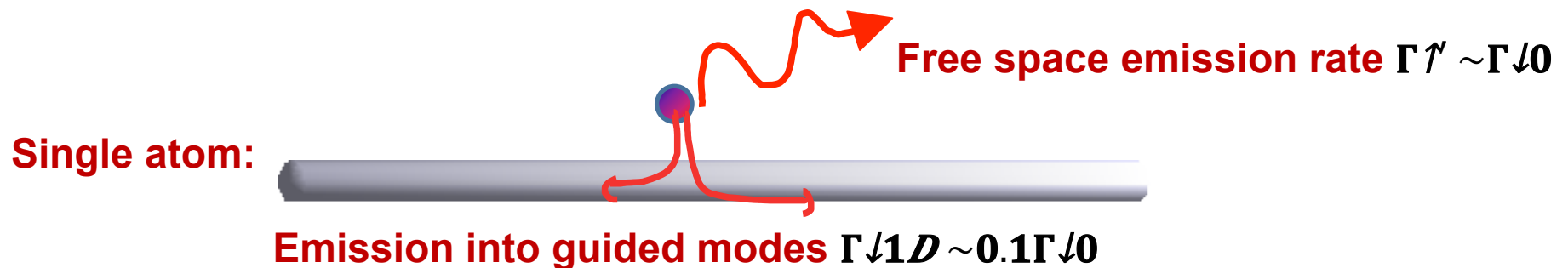
Atom-nanofiber interface

- Trap atoms in lattice with far off-resonant guided modes



Vetsch et al, PRL 104, 203603 (2010)

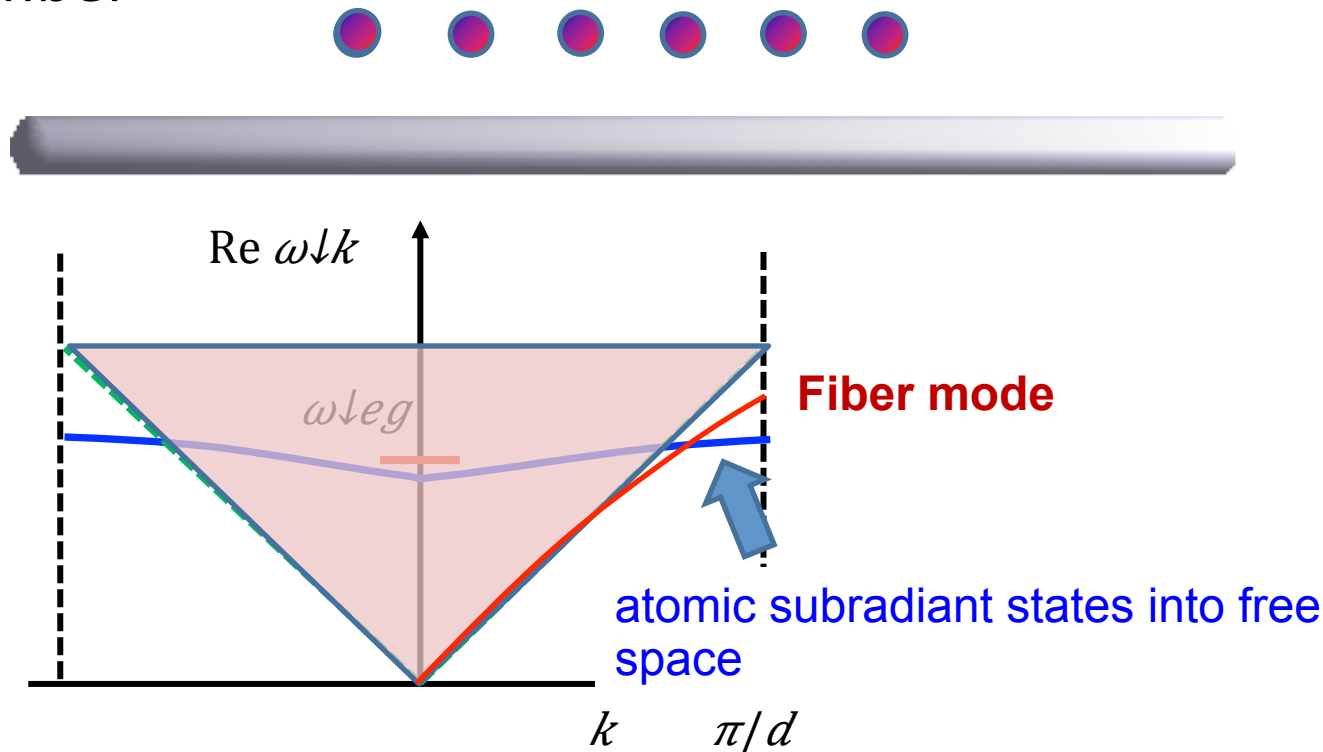
- Efficient coupling of atoms with near-resonant photons



- Expts: Rauschenbeutel (TU Vienna), Kimble (Caltech), Orozco (JQI), Laurat (Paris), Polzik (Niels Bohr Institute), ...

“Selective” subradiance

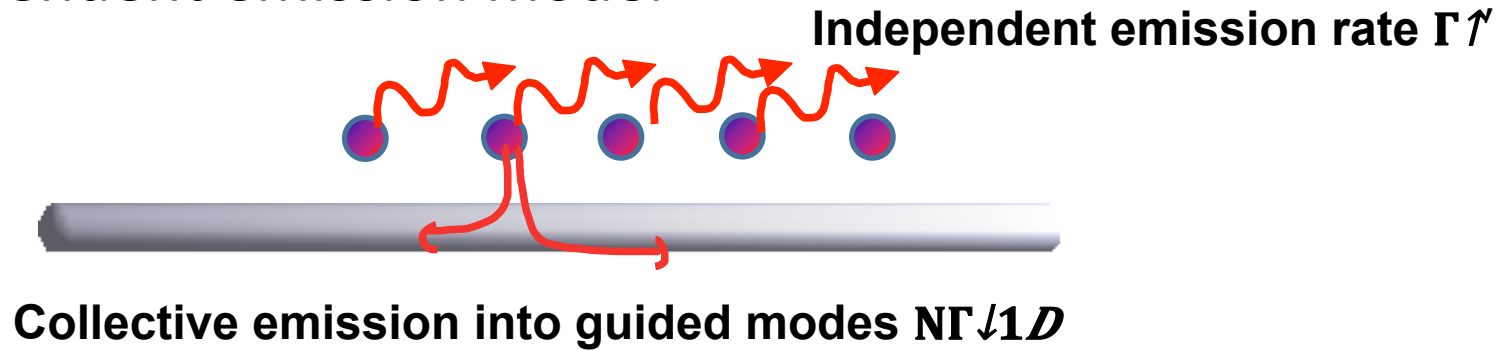
- Previously subradiant states can now couple to guided modes of the fiber



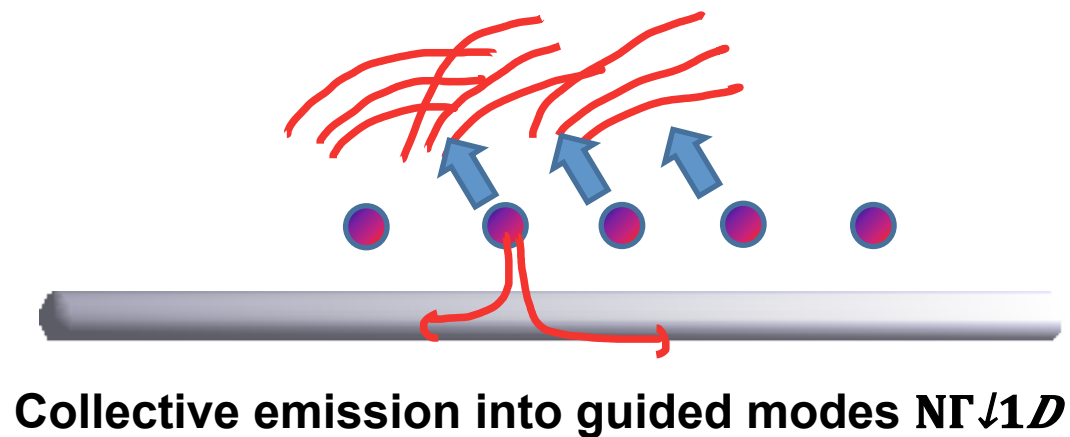
- Not subradiant, but “selectively” subradiant... **perfect!**
- Efficient atom-light interface: atoms talk very well to modes of interest, rather than undesired modes (free space)
 - Use this to beat the “optical depth” limit

Model of atom-light interactions

- “Independent emission model”

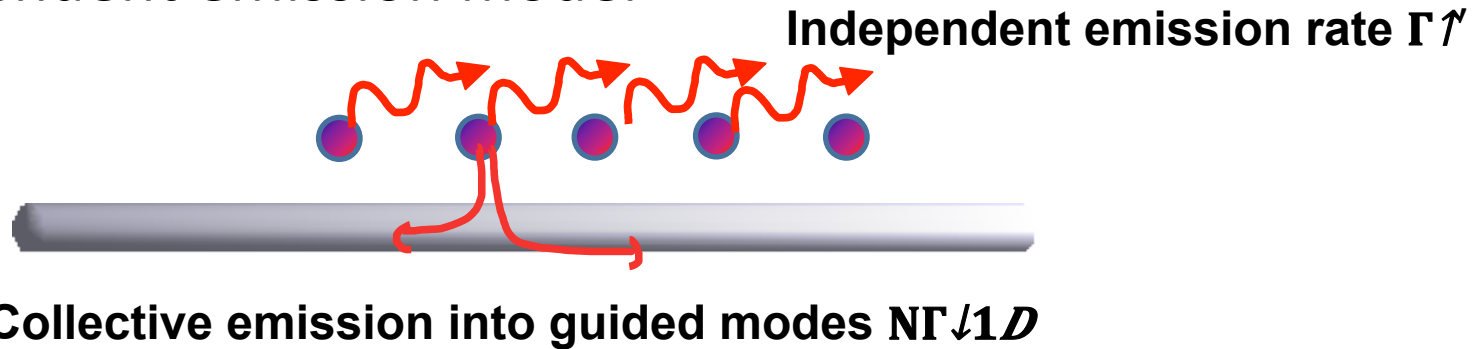


- Collective emission model



Model of atom-light interactions

- “Independent emission model”



- Collective emission model
- Use our “spin model”

$$H_{\text{eff}} = -\mu \sum_{i,j} \frac{d_{ij}}{d} \frac{\omega_{\text{leg}}}{\omega_{\text{leg}} - \omega} \sum_{i,j} G(r_j, r_i, \omega_{\text{leg}}) \sigma_{\text{leg}}^{\uparrow i} \sigma_{\text{leg}}^{\uparrow j}$$

Exact for cylindrical fiber

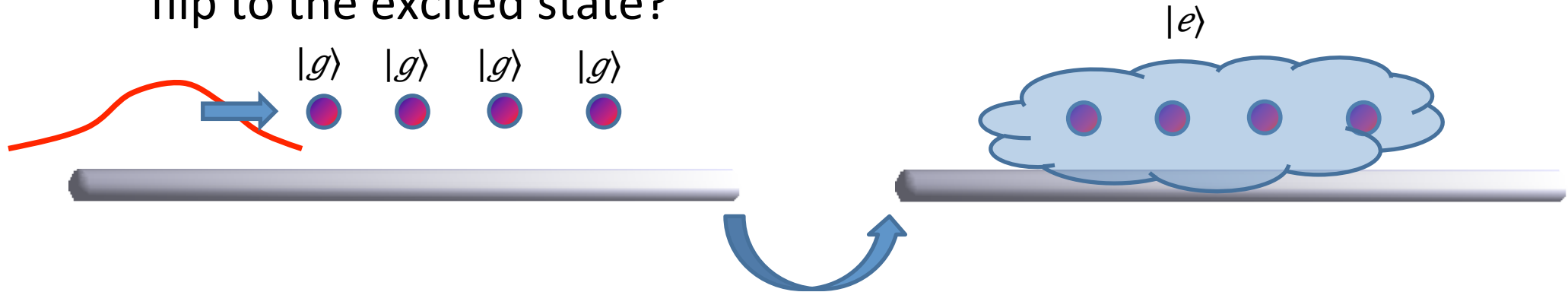
- Field propagation

$$E(r,t) = E_{\text{in}}(r,t) + (3\pi\hbar c \Gamma / d \omega_{\text{leg}}) \sum_i G(r, r_i, \omega_{\text{leg}}) \sigma_{\text{leg}}^{\uparrow i}(t)$$

Project into guided mode of fiber

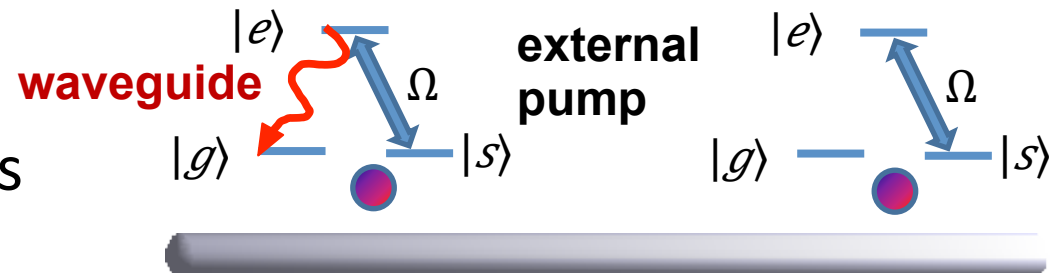
Photon storage

- How well can one store an incoming photon as a collective spin flip to the excited state?



- By time reversal symmetry, can study the reverse problem of mapping an initial spin excitation into a photon

- Use three-level systems

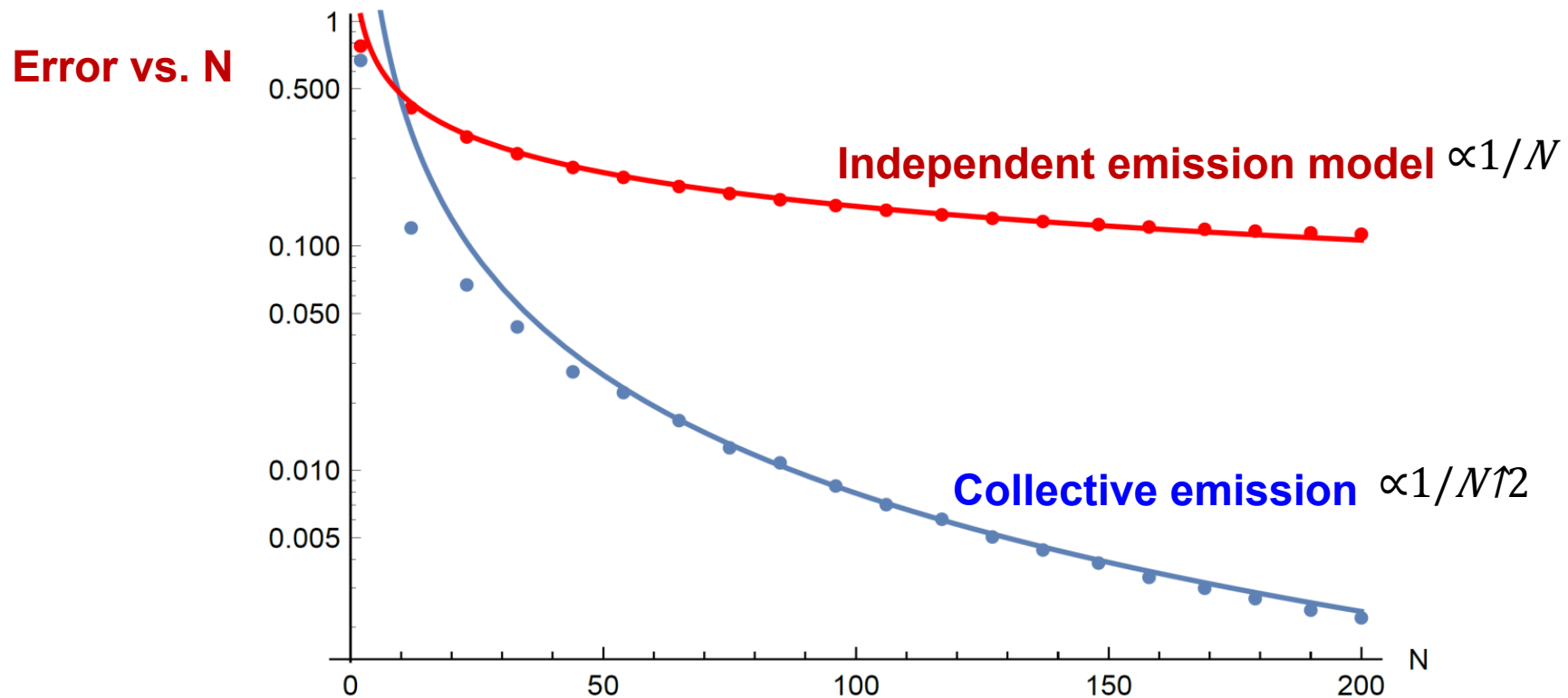


- Independent emission model: error $\sim \Gamma \uparrow / M \downarrow 1D = 1/OD$

Same result as free space: AV Gorshkov et al, PRL 98, 123601 (2007)

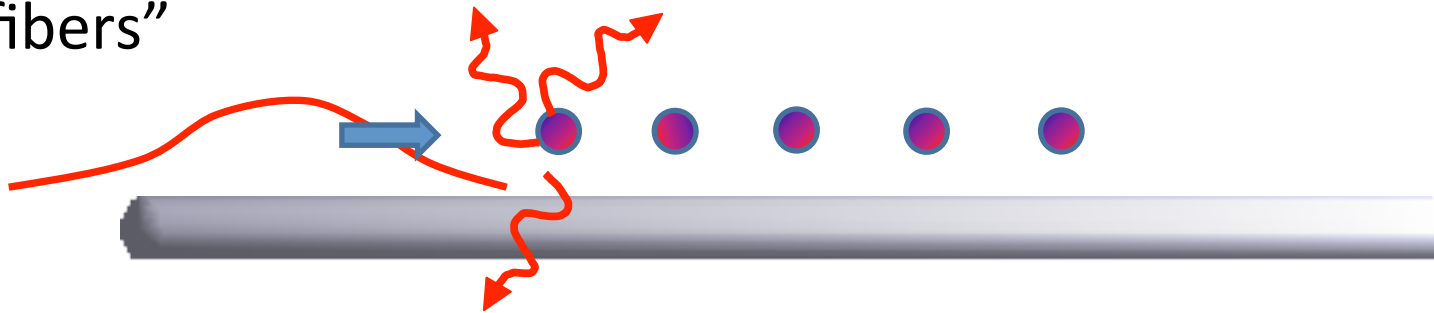
Photon storage/generation

- Collective emission:
 - Numerically diagonalize H_{eff} within manifold of single $|s\rangle$ excitation
 - Look for eigenstate with highest emission probability into waveguide



What limits the process?

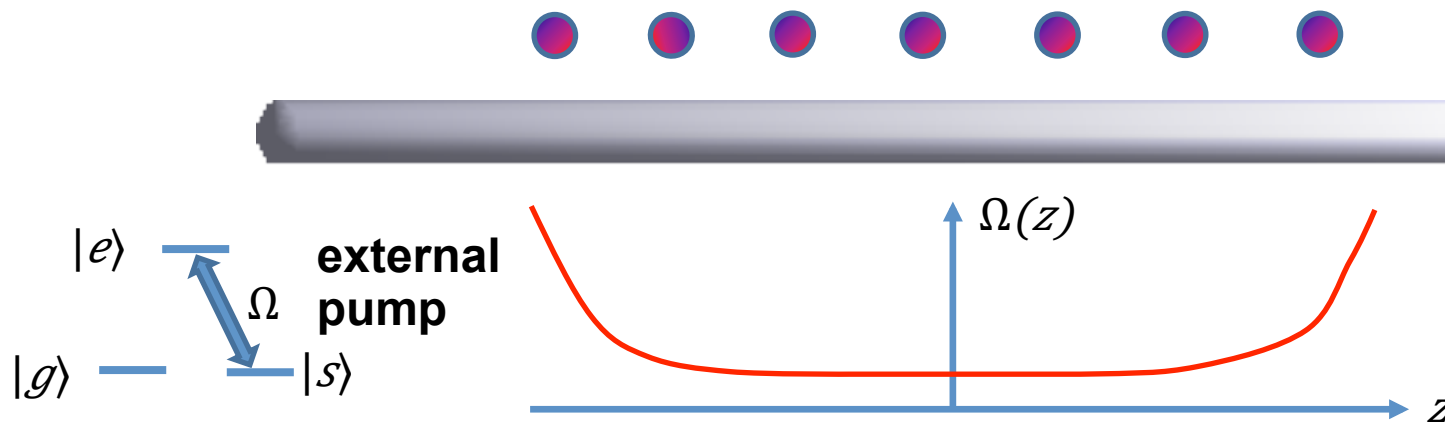
- $1/N^2$ error seems less than ideal
 - $1/N$ from collective enhancement into guided mode
 - “Only” $1/N$ suppression of emission into free space
- Origin: scattering losses at interface between two different “fibers”



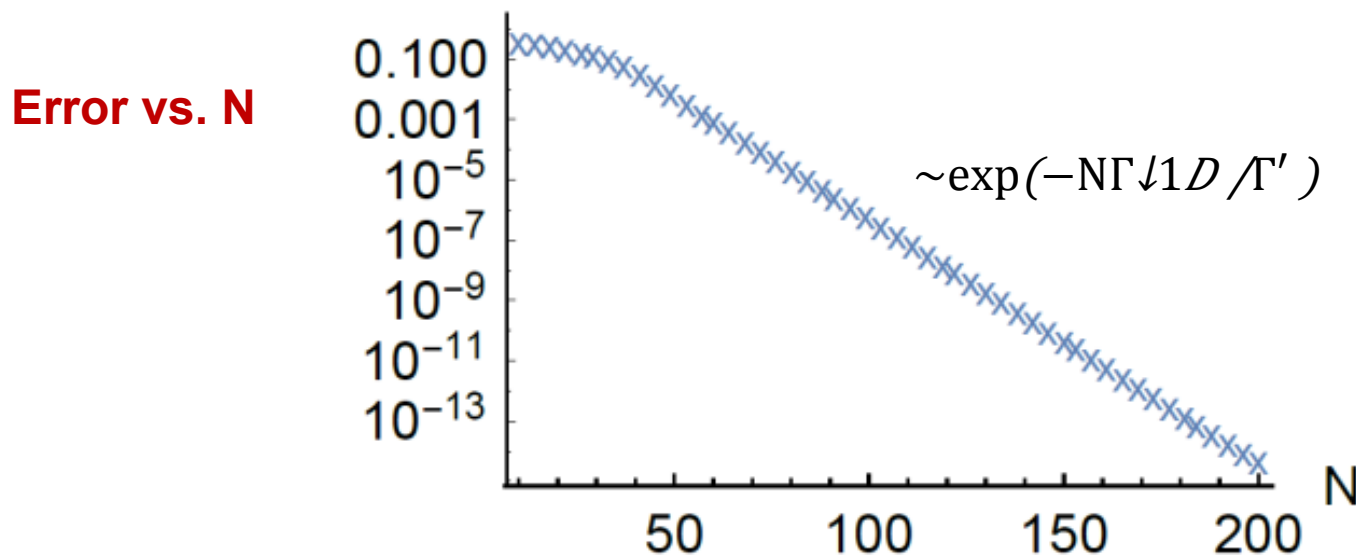
- Need a fiber “connector!”

Photon storage with impedance matching

- Spatially vary the pump field profile, so that atoms at the ends “gradually” couple to fiber

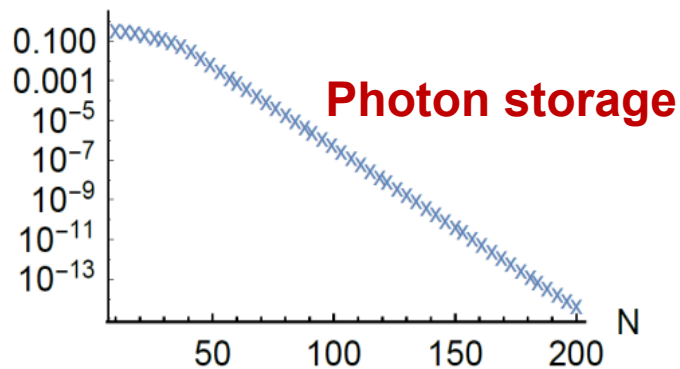


- Exponential suppression of error!



Outlook

- Spontaneous emission represents a fundamental barrier for atom-light interactions
- What is the new ceiling if one exploits subradiance?



**Nonlinear optics,
Photon gates,
Spin physics,
Lattice clocks, ...**

- Atom-light interactions as a quantum spin model

$$H_{\text{eff}} = - (3\pi\hbar\Gamma / 4c/\omega_{\text{leg}}) \sum_{i,j} G(r_{\downarrow j}, r_{\downarrow i}, \omega_{\text{leg}}) \sigma_{\text{leg}\uparrow i} \sigma_{\text{leg}\uparrow j}$$

$$E(r,t) = E_{\text{lin}}(r,t) + (3\pi\hbar c\Gamma / 4d / \omega_{\text{leg}}) \sum_i \hat{G}(r, r_{\downarrow i}, \omega_{\text{leg}}) \sigma_{\text{leg}\uparrow i}(t)$$

- New insights or numerical tools for AMO physics?

Matrix product states

