



Anderson transitions in Euclidean random matrix models

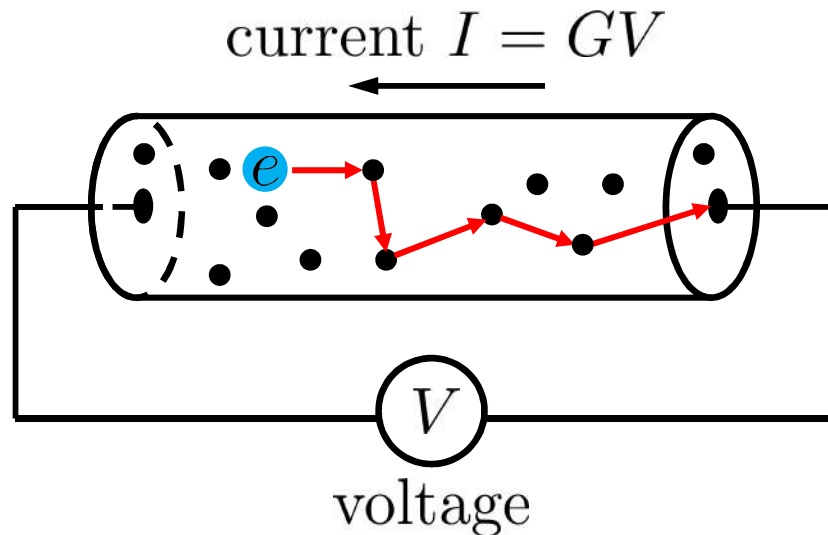
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Laboratoire de Physique et Modélisation des Milieux Condensés
CNRS and Université Grenoble Alpes, France



Anderson localization

Metal



Conductance $G = 1/R$

High T

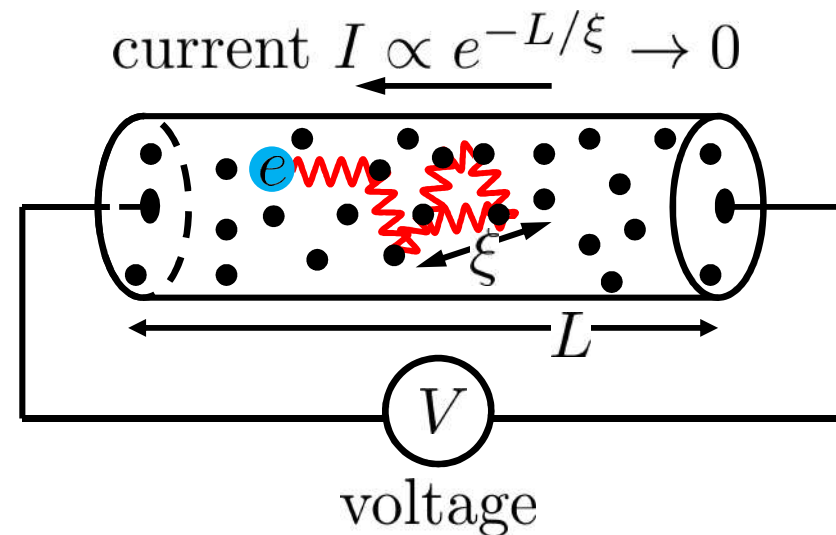
Few impurities

Metal-insulator
transition in 3D

$T \rightarrow 0$

Add impurities

Anderson insulator



$G \propto e^{-L/\xi} \rightarrow 0$

ξ — localization
length

Anderson, PR **109**, 1492 (1958)

Absence of transport versus localization

Green's function

$$\left\{ \nabla^2 + \frac{2m}{\hbar^2} [E - V(\mathbf{r})] \right\} G(\mathbf{r}, \mathbf{r}') = \delta(\mathbf{r} - \mathbf{r}')$$

Eigenmodes

$$\left\{ \nabla^2 + \frac{2m}{\hbar^2} [E_n - V(\mathbf{r})] \right\} \psi_n(\mathbf{r}) = 0$$

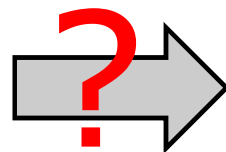
$$G(\mathbf{r}, \mathbf{r}') = \frac{\hbar^2}{2m} \sum_n \frac{\psi_n(\mathbf{r}) \psi_n^*(\mathbf{r}')}{E - E_n}$$

Transport

Localization

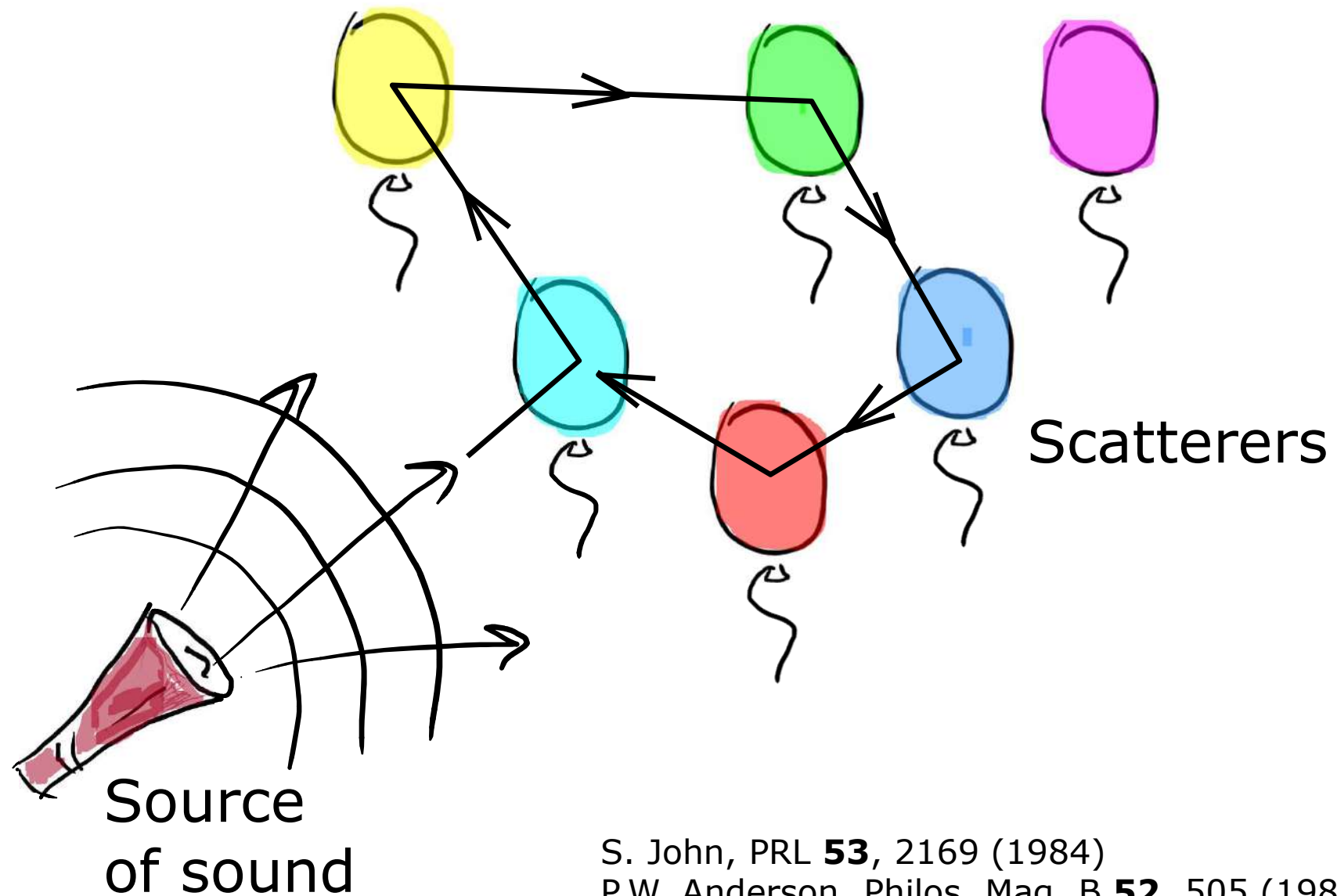
Localized modes do not contribute to transport

No diffusion



Localization

Can we do the same with “classical” waves?



S. John, PRL **53**, 2169 (1984)

P.W. Anderson, Philos. Mag. B **52**, 505 (1985)

Yes we can! ...in 1D and 2D

Quasi-1D: Chabanov et al., Nature **404**, 850 (2000)

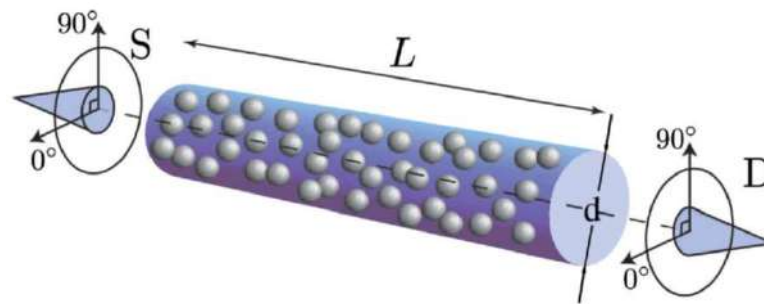
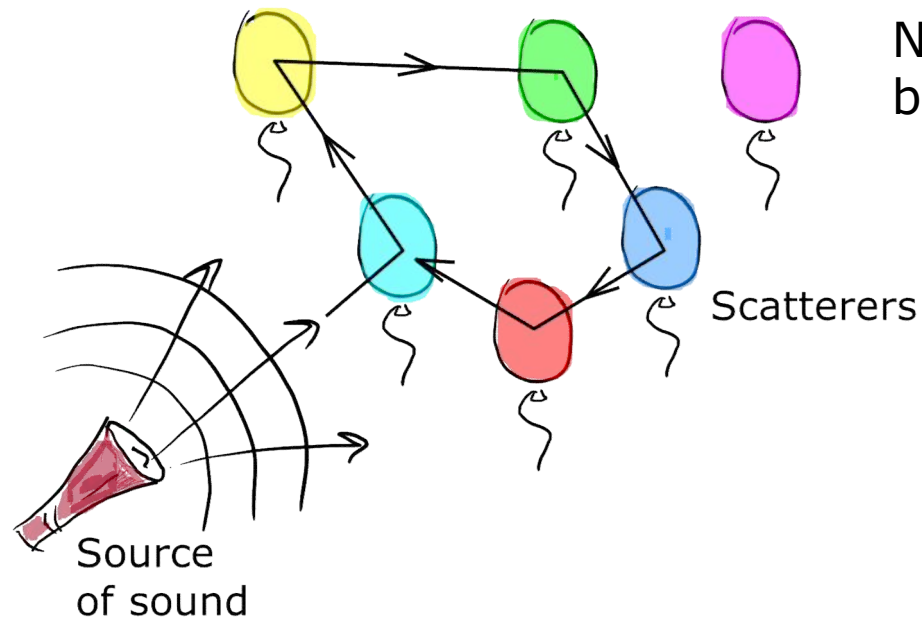


Figure from Cherroret et al., PRB **80**, 045118 (2009)

2D: Weaver, Wave Motion **12**, 129 (1990)
Dalichaouch et al., Nature **354**, 53 (1991)
Schwartz et al., Nature **446**, 52 (2007)

...

Yes we can! A real challenge in 3D



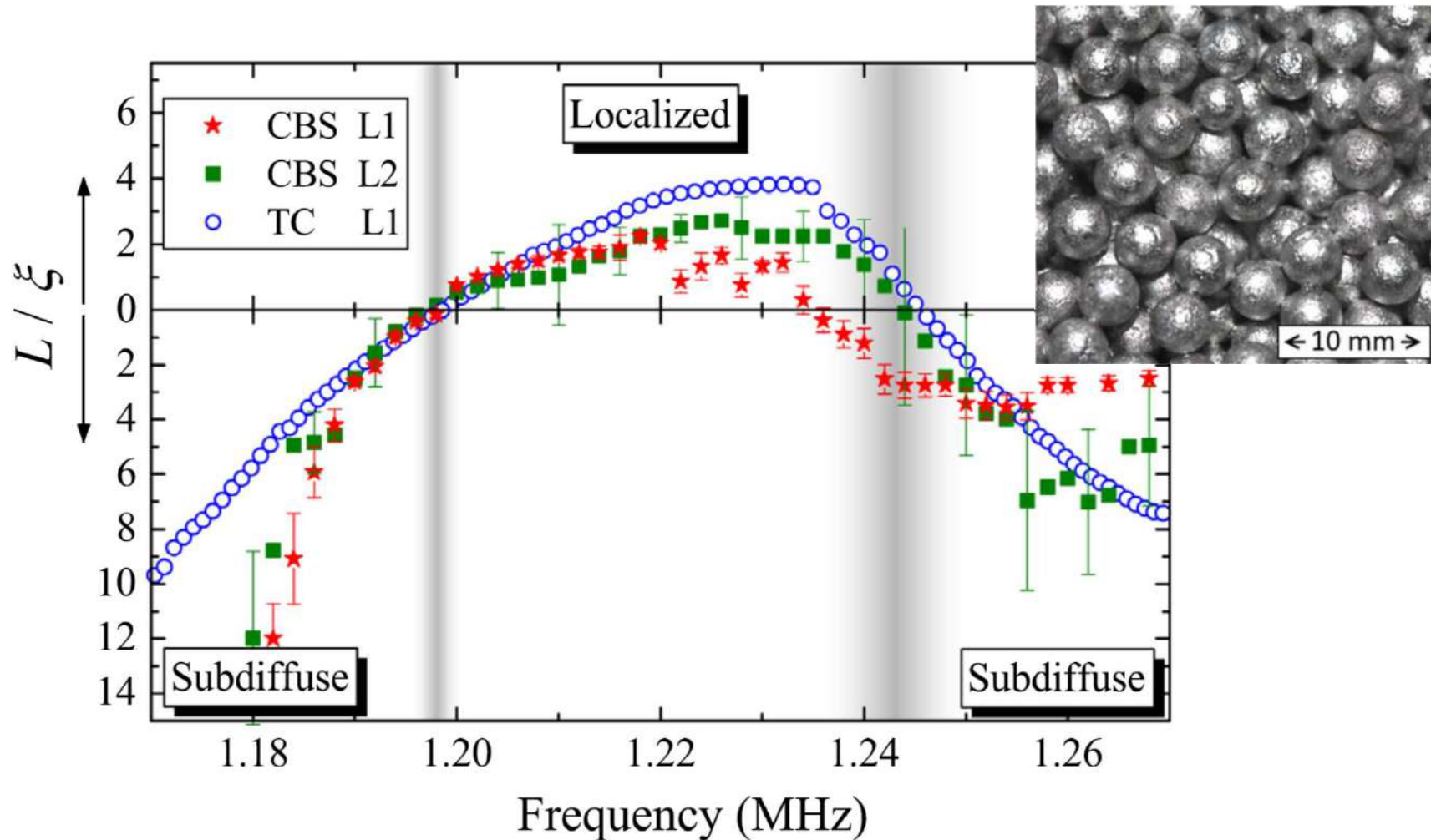
No Anderson localization of sound by balloons...

... but sound can be localized by **stronger** scattering objects (aluminum beads) that are **densely packed**



Hu et al., Nature Physics **4**, 945 (2008)

Anderson localization of elastic waves in 3D



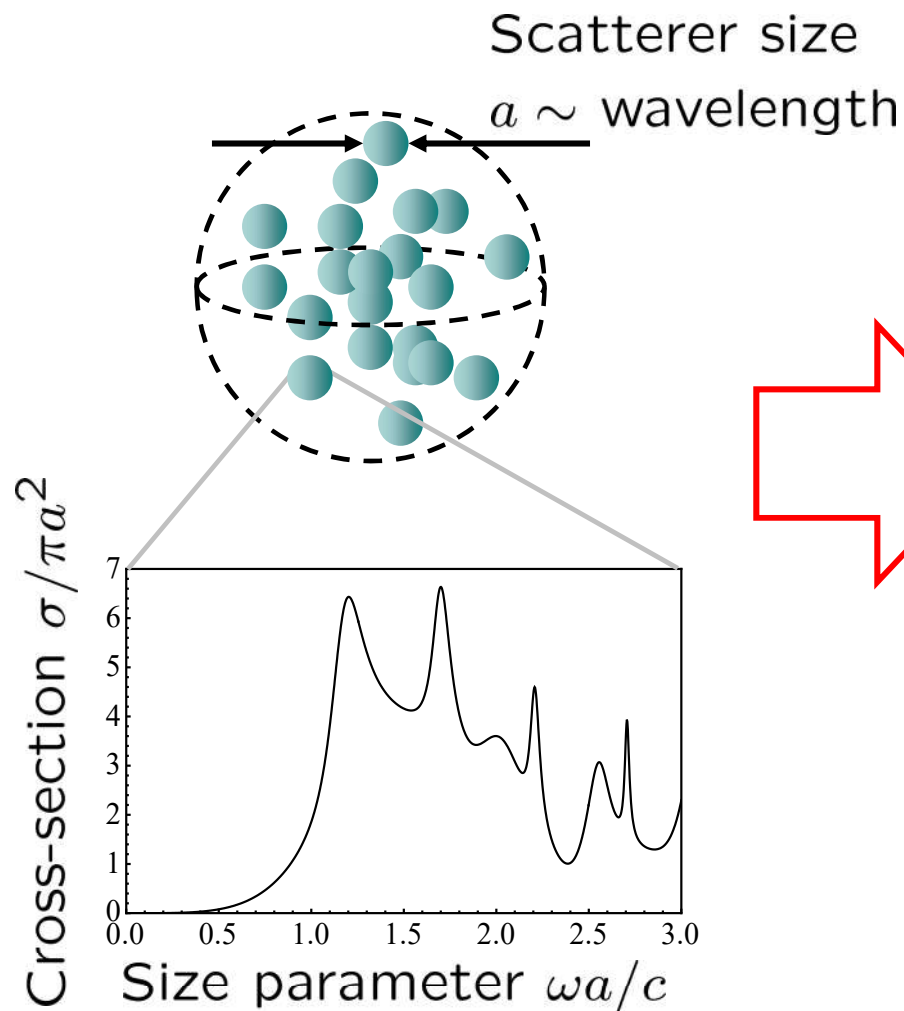
ξ — localization or correlation length

L — sample thickness

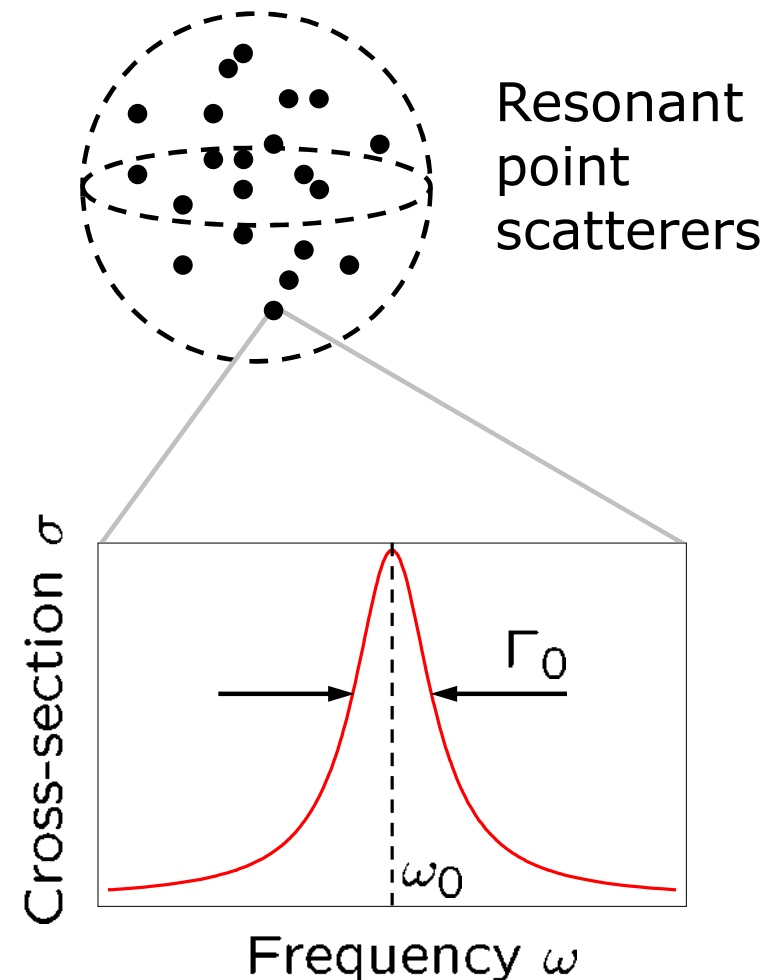
Cobus et al., PRL **116**, 193901 (2016)

A minimal model of disordered media

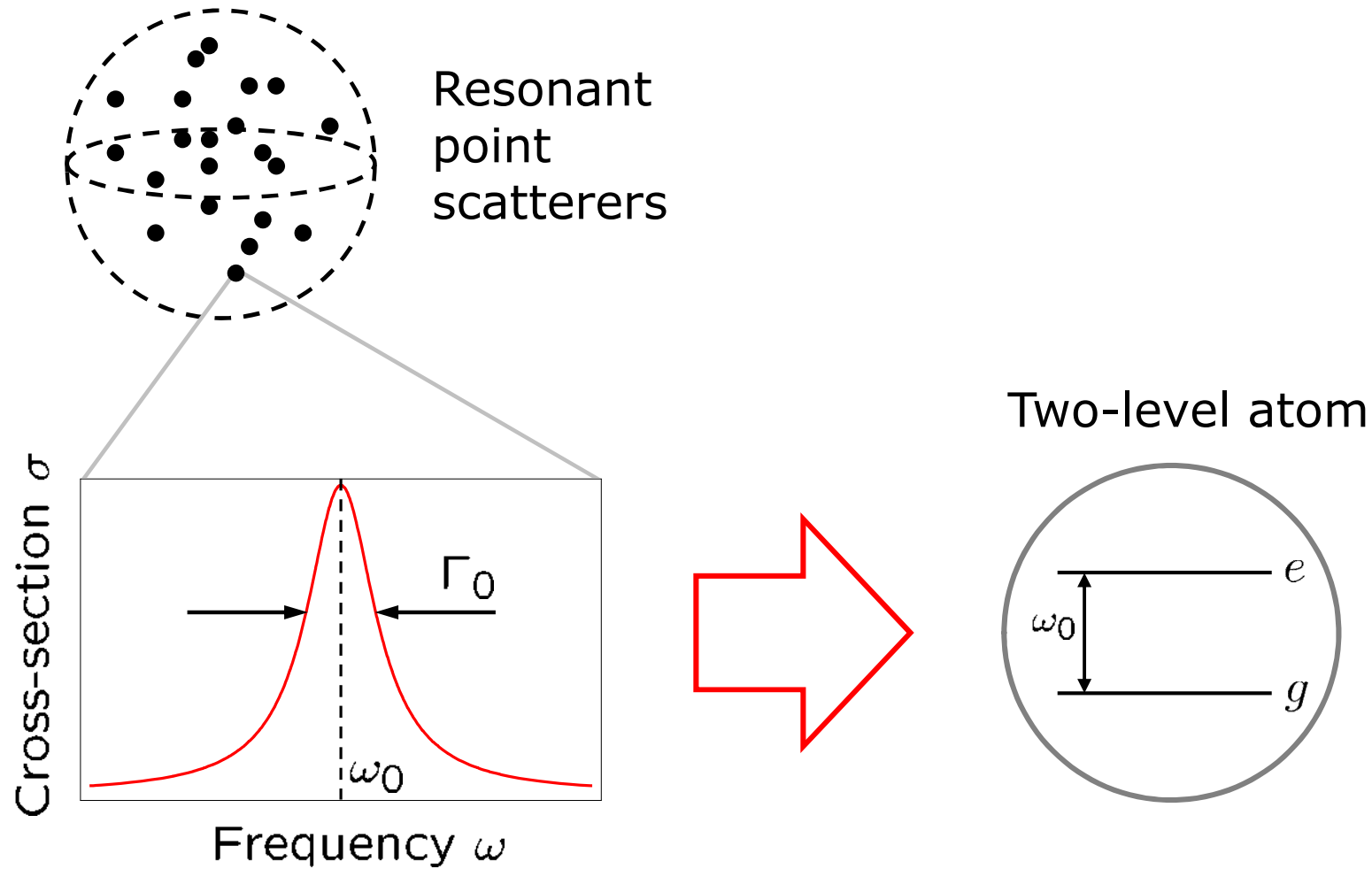
Real samples complex



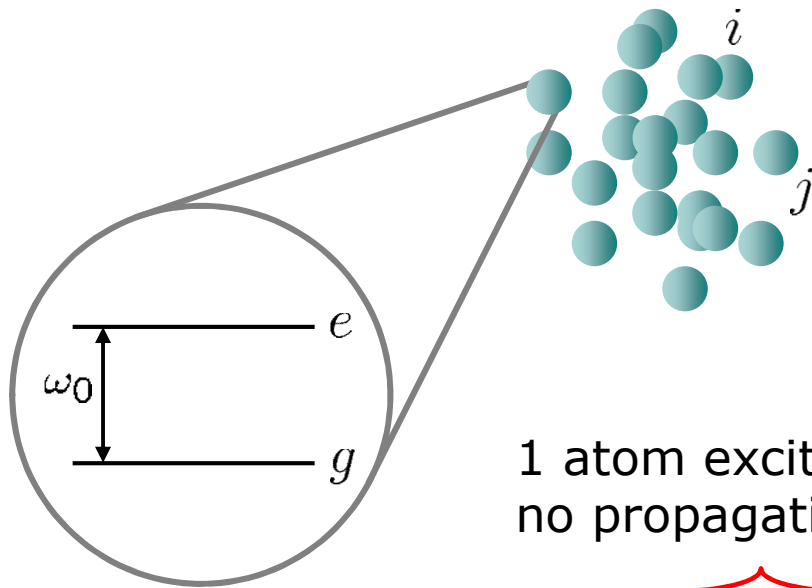
Our model simpler



Atom is a realistic resonant point scatterer



A scalar photon in a cloud of atoms



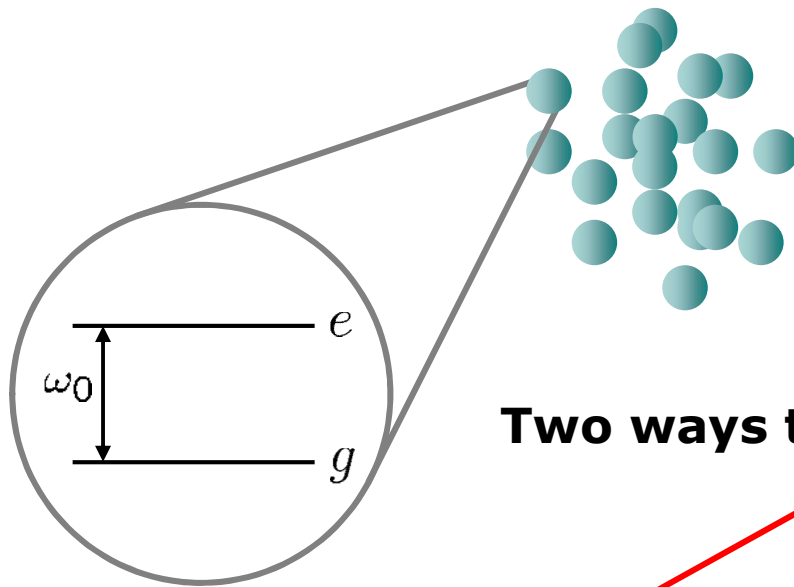
1 atom excited
no propagating photons

all atoms in the ground state
+ a photon in the mode \mathbf{k}

$$\begin{aligned}
 |\psi(t)\rangle &= \underbrace{\sum_{j=1}^N \beta_j(t) |(N-1) : g, j : e\rangle |0_R\rangle}_{\text{1 atom excited, no propagating photons}} + \underbrace{\sum_{\mathbf{k}} \gamma_{\mathbf{k}}(t) |N : g\rangle |\mathbf{k}\rangle}_{\text{all atoms in the ground state + a photon in the mode } \mathbf{k}} \\
 &+ \underbrace{\sum_{i < j}^N \sum_{\mathbf{k}} \alpha_{ij\mathbf{k}}(t) |(N-2) : g, i : e, j : e\rangle |\mathbf{k}\rangle}_{\text{2 atoms excited + a (virtual) photon in the mode } \mathbf{k}}
 \end{aligned}$$

2 atoms excited + a (virtual) photon in the mode \mathbf{k}

A scalar photon in a cloud of atoms



Two ways to deal with the problem

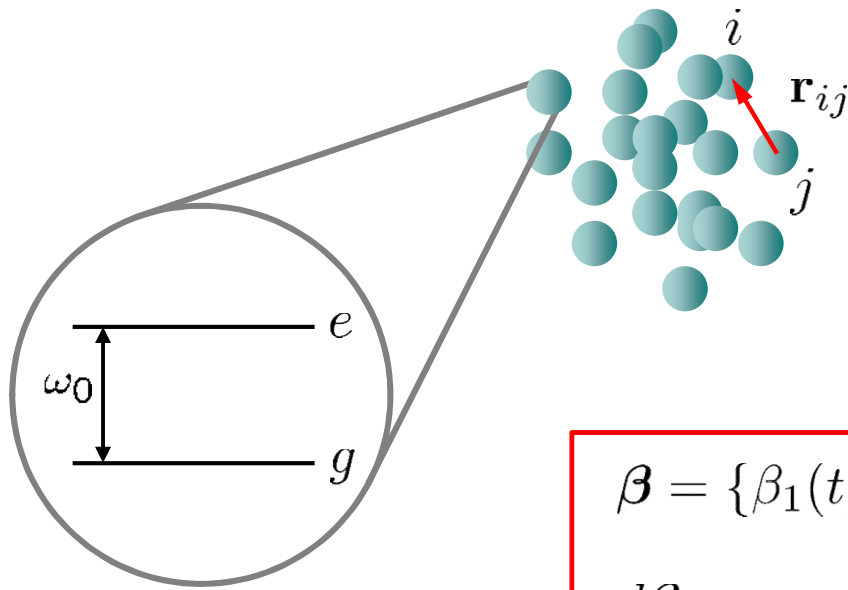
Waves (light) scattered
by the scatterers (atoms)

Equations to analyze:
wave equation with
random material properties
(dielectric constant)

Scatterers (atoms) coupled
by waves (light)

Equations to analyze:
dynamic equations for
identical nodes with
random couplings

Green's matrix



Green's matrix G describes propagation of light between pairs of atoms
 $r_{ij} = |\mathbf{r}_i - \mathbf{r}_j|$

Euclidean random matrix

$$\boldsymbol{\beta} = \{\beta_1(t), \beta_2(t), \dots, \beta_N(t)\}$$

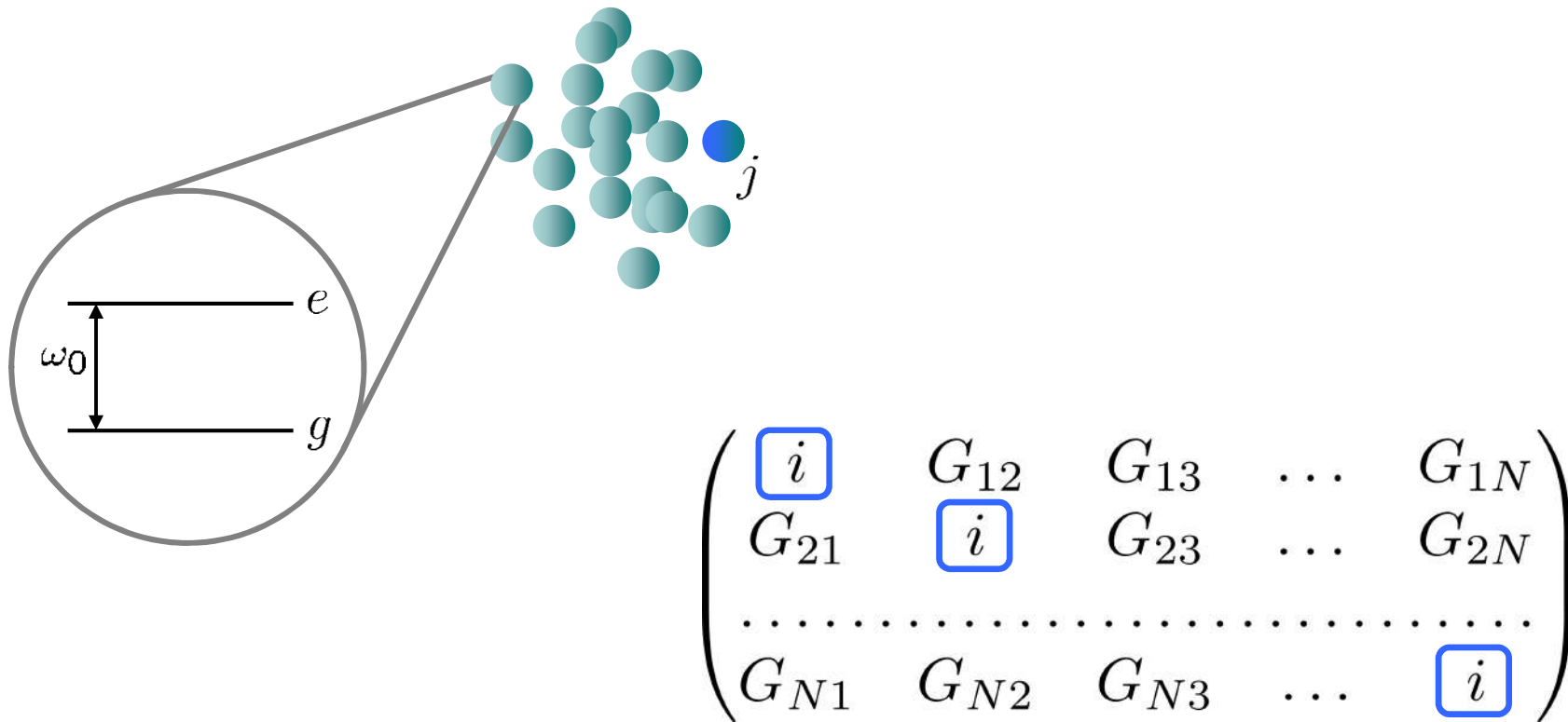
$$\frac{d\boldsymbol{\beta}}{dt} = iG\boldsymbol{\beta}(t)$$

$$G_{ij} = i\delta_{ij} + (1 - \delta_{ij}) \frac{e^{ik_0 r_{ij}}}{k_0 r_{ij}}$$

$$k_0 = \frac{\omega_0}{c}$$

Prasad & Glauber, PRA **31**, 1583 (1985)
Svidzinsky & Chang, PRA **77**, 043833 (2008)
Friedberg & Manassah, Phys. Lett. A **372**, 2514 (2008)

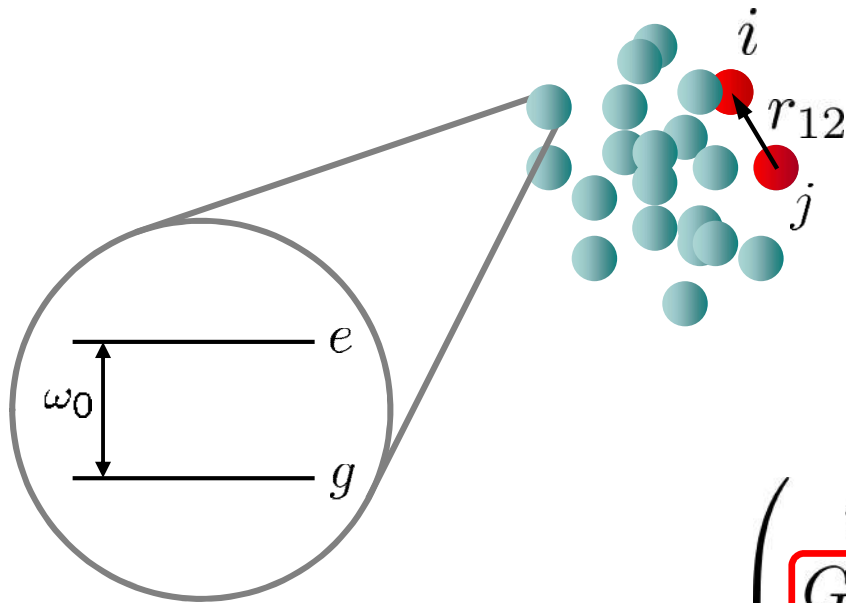
Structure of the Green's matrix



One-atom dynamics:

$G_{jj} = i$ describes the decay $e^{-\Gamma_0 t}$ of the excitation of an isolated excited atom

Structure of the Green's matrix

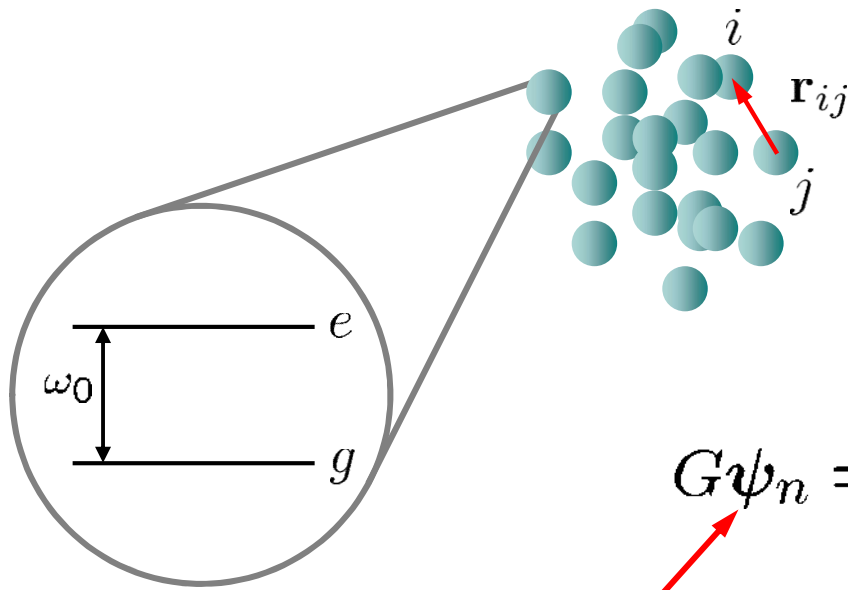


$$\begin{pmatrix} i & G_{12} & G_{13} & \dots & G_{1N} \\ G_{21} & i & G_{23} & \dots & G_{2N} \\ \dots & \dots & \dots & \dots & \dots \\ G_{N1} & G_{N2} & G_{N3} & \dots & i \end{pmatrix}$$

Pairwise coupling between atoms 1 & 2:

$G_{12} = e^{ik_0 r_{12}} / k_0 r_{12}$ is the field at position 2 due to a source at position 1

Quasi-modes of the system



Green's matrix G describes propagation of light between pairs of atoms
 $r_{ij} = |\mathbf{r}_i - \mathbf{r}_j|$

$$G\psi_n = \Lambda_n\psi_n$$

Eigenvectors:

$$\psi_n = \{\psi_n^1, \psi_n^2, \dots, \psi_n^N\}$$

Eigenvalues:

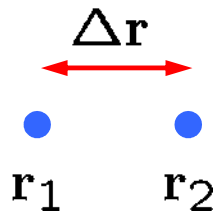
$$\Lambda_n = \text{Re}\Lambda_n + i\text{Im}\Lambda_n$$

Frequency
of the mode

Decay rate
of the mode

Green's matrix for $N = 2$

$$G_{ij} = i\delta_{ij} + (1 - \delta_{ij}) \frac{e^{ik_0|\mathbf{r}_i - \mathbf{r}_j|}}{k_0|\mathbf{r}_i - \mathbf{r}_j|}$$
$$\hat{G}\psi_n = \Lambda_n\psi_n$$



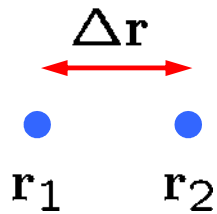
$$\hat{G} = \begin{pmatrix} i & \frac{e^{ik_0\Delta r}}{k_0\Delta r} \\ \frac{e^{ik_0\Delta r}}{k_0\Delta r} & i \end{pmatrix}$$

$$\Lambda_{1,2} = i \pm G_{12}$$
$$\psi_{1,2} = (\pm 1, 1)$$

Green's matrix for $N = 2$

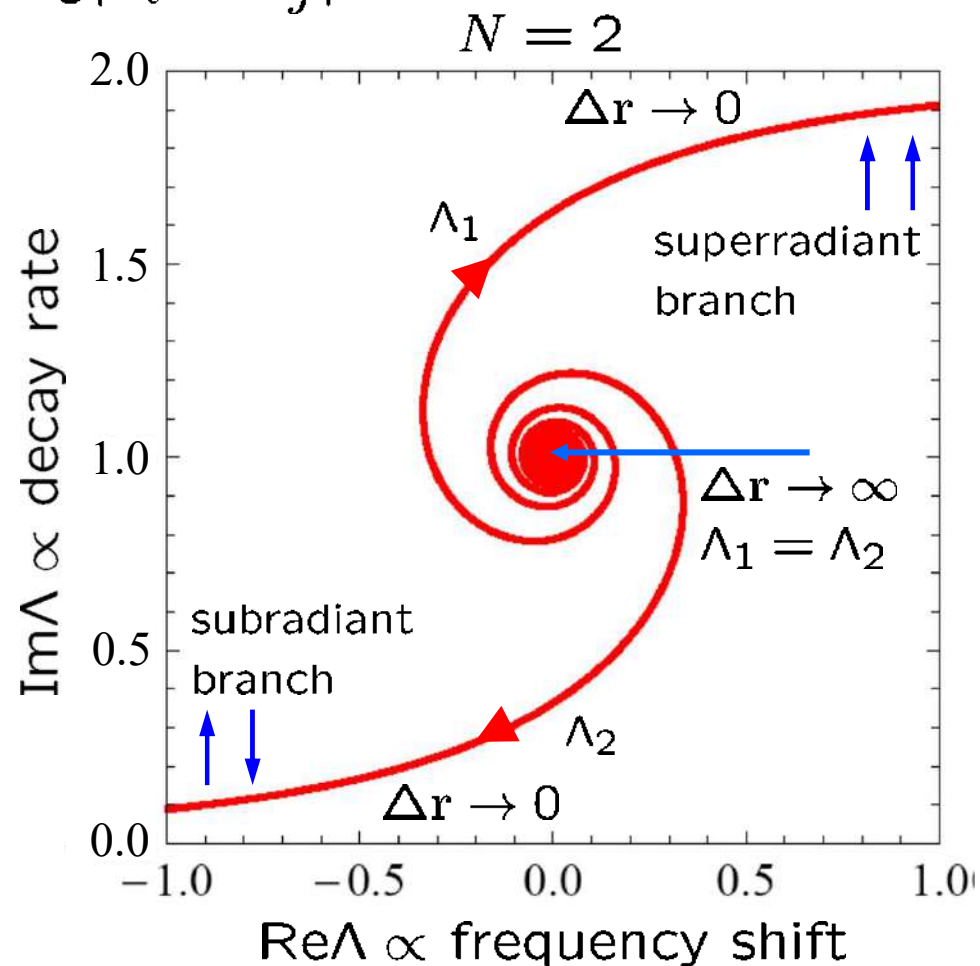
$$G_{ij} = i\delta_{ij} + (1 - \delta_{ij}) \frac{e^{ik_0|\mathbf{r}_i - \mathbf{r}_j|}}{k_0|\mathbf{r}_i - \mathbf{r}_j|}$$

$$\hat{G}\psi_n = \Lambda_n\psi_n$$

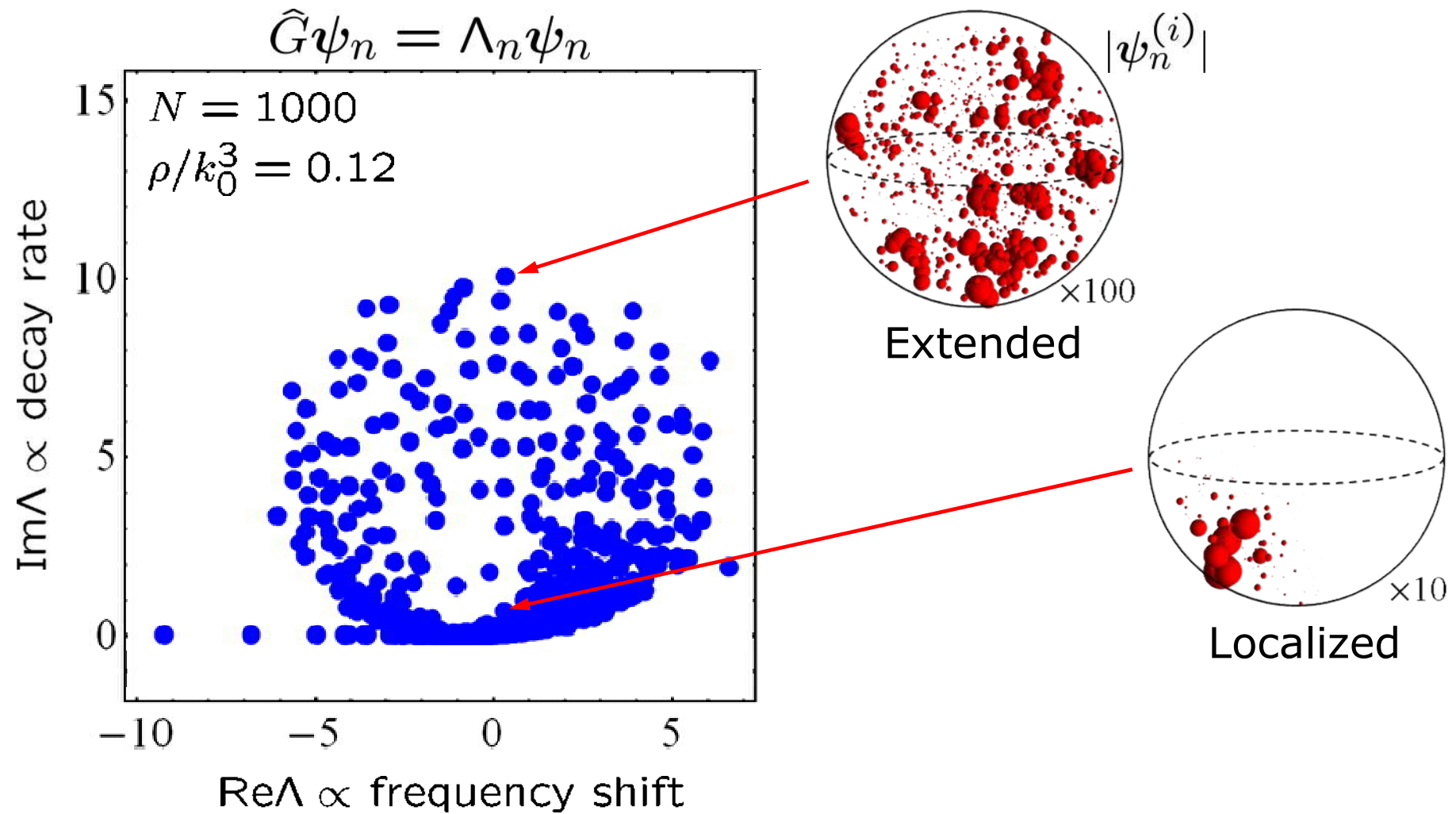


$$\Lambda_{1,2} = i \pm G_{12}$$

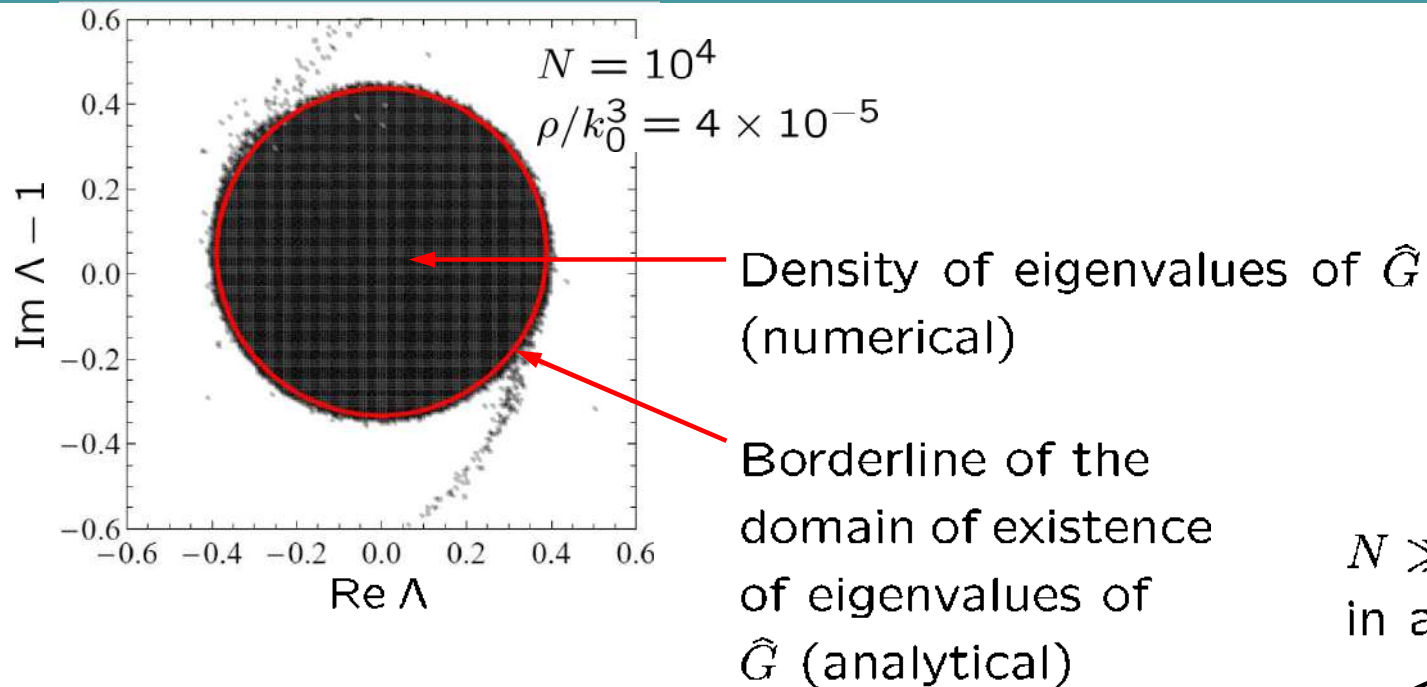
$$\psi_{1,2} = (\pm 1, 1)$$



Green's matrix for $N \gg 1$



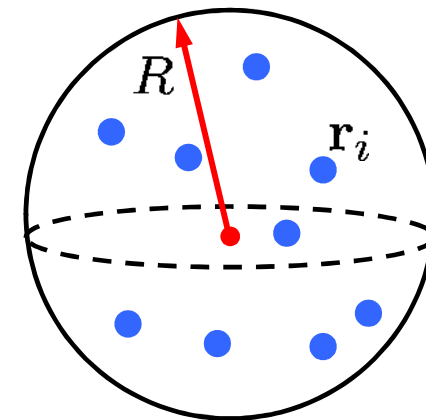
Eigenvalue density for $N \gg 1$



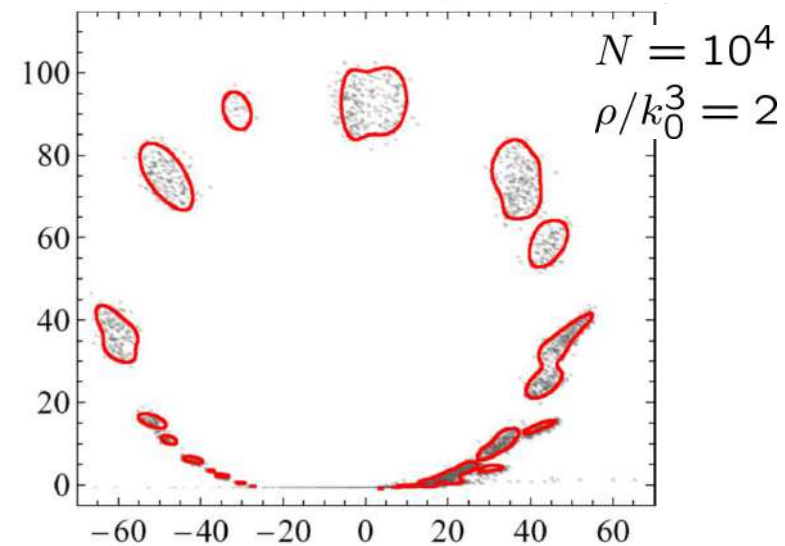
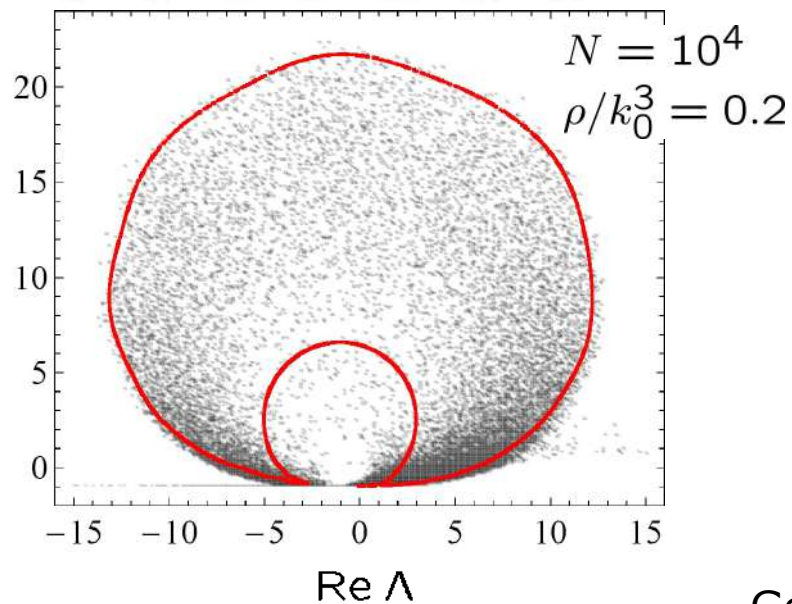
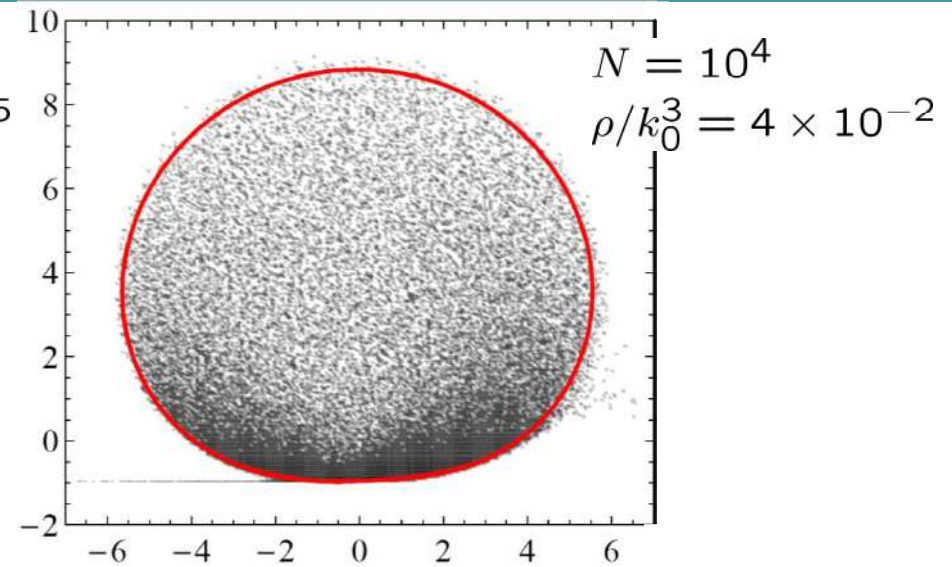
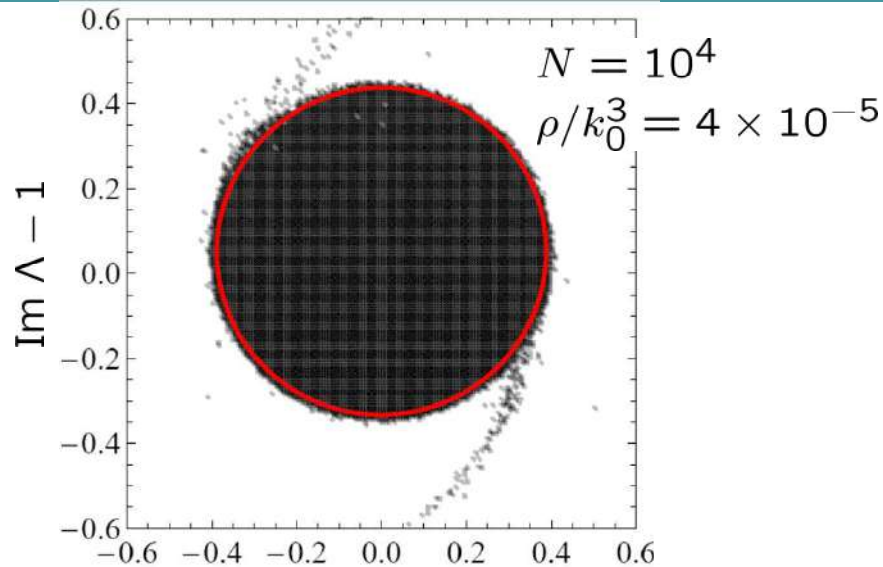
$$G_{ij} = i\delta_{ij} + (1 - \delta_{ij}) \frac{e^{ik_0|\mathbf{r}_i - \mathbf{r}_j|}}{k_0|\mathbf{r}_i - \mathbf{r}_j|}$$

$$\hat{G}\psi_n = \Lambda_n\psi_n$$

$N \gg 1$ points
in a sphere:



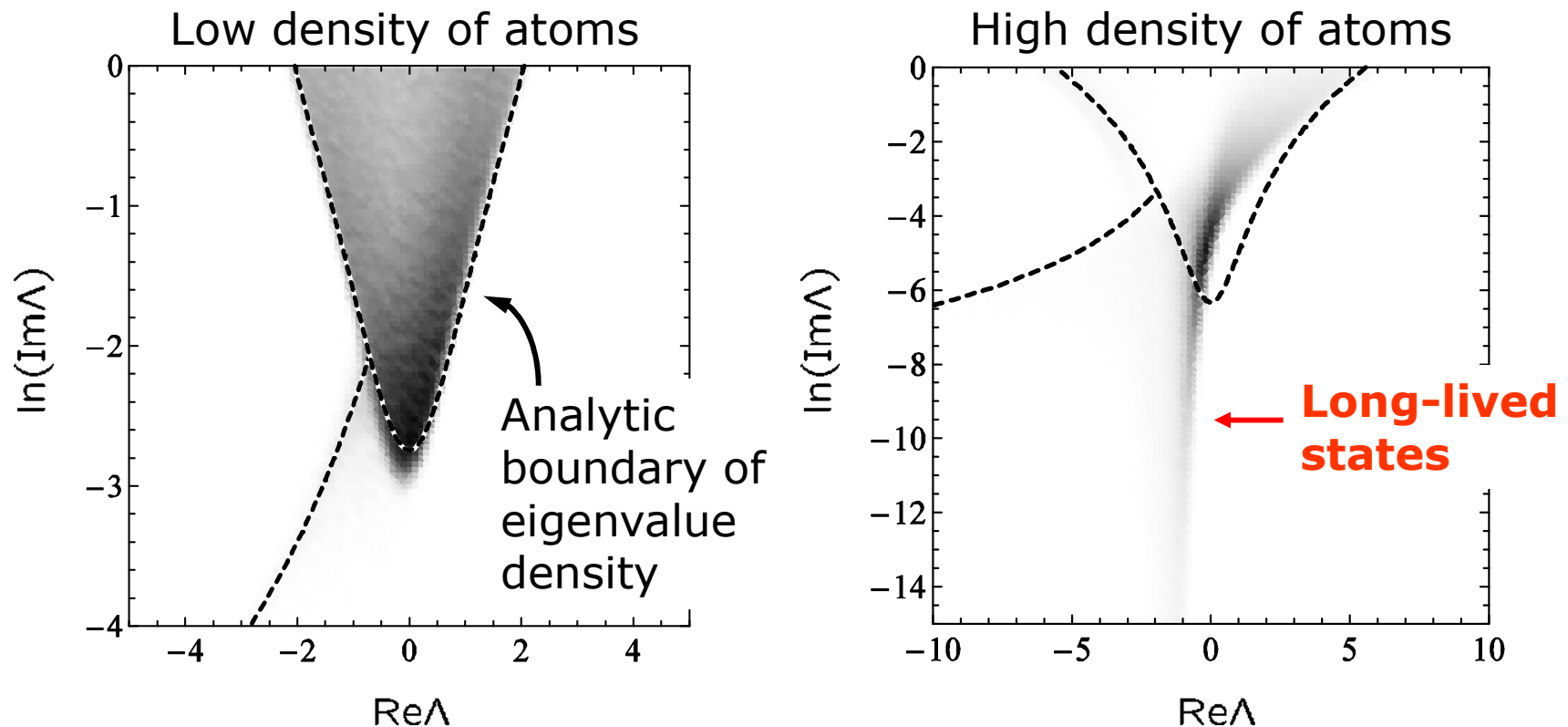
Eigenvalue density for $N \gg 1$



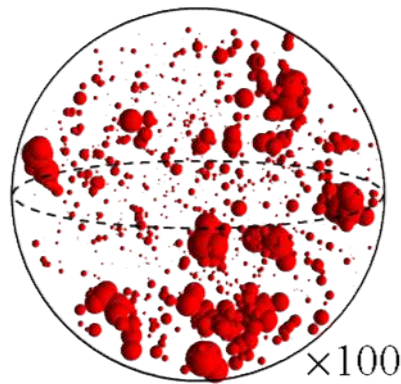
Goetschy & Skipetrov, PRE **84**, 011150 (2011)

Breakdown of analytic theory for $\text{Im } \Lambda \ll 1$

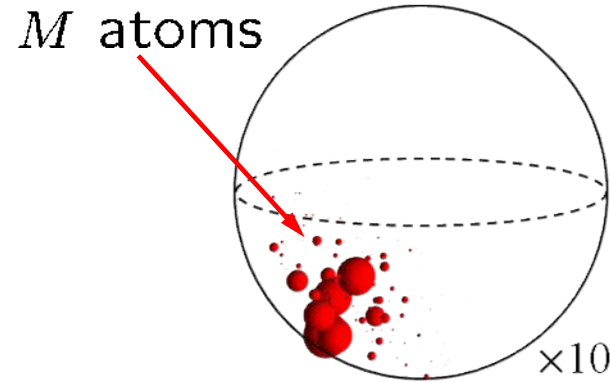
Density of eigenvalues



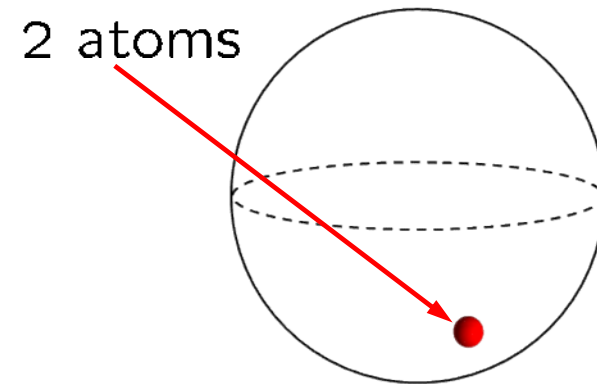
Inverse participation ratio (IPR)



Small IPR
 $\sim 1/N \ll 1$



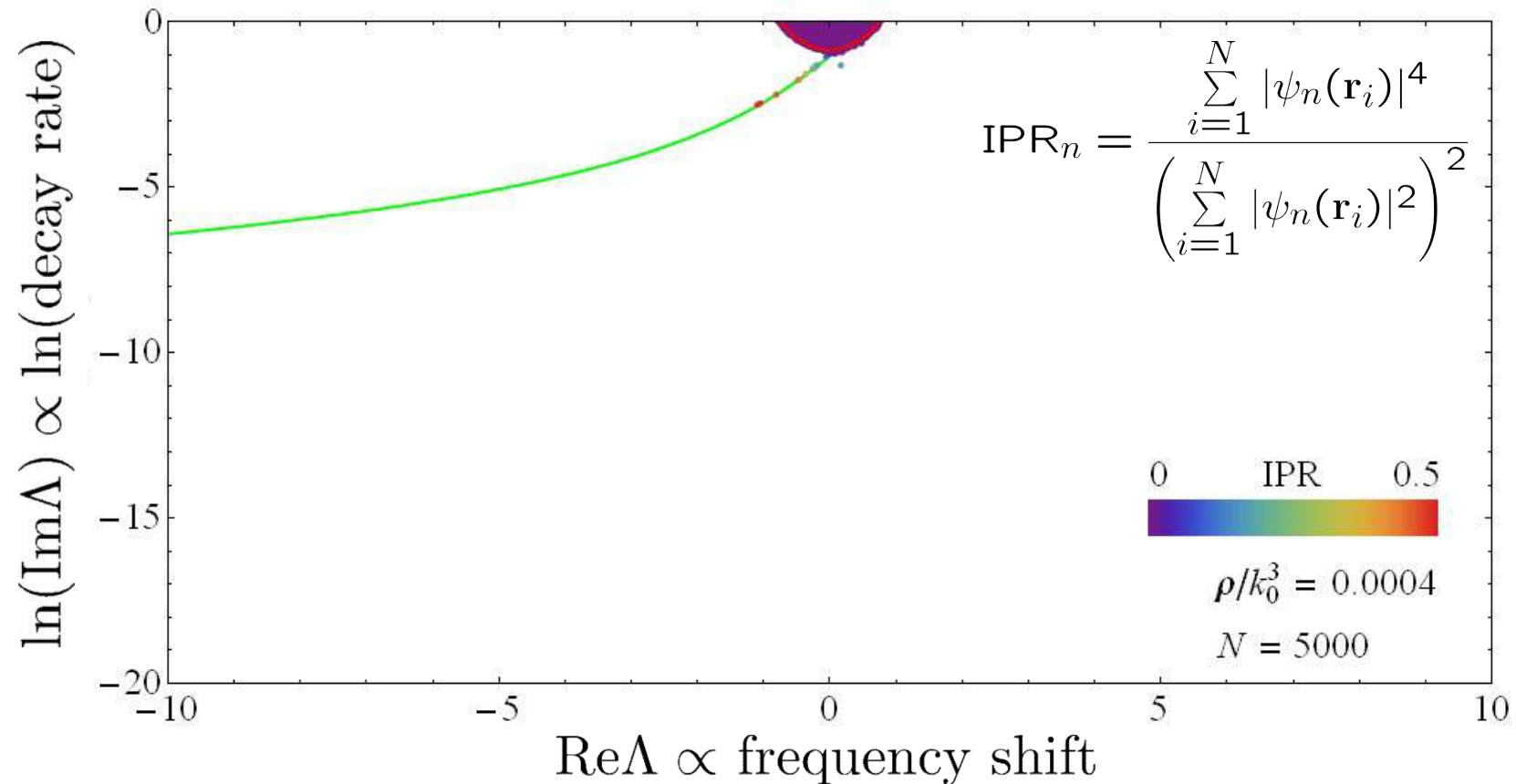
Appreciable IPR
 $\sim 1/M \gg 1/N$



Large IPR
 $\sim 1/2$

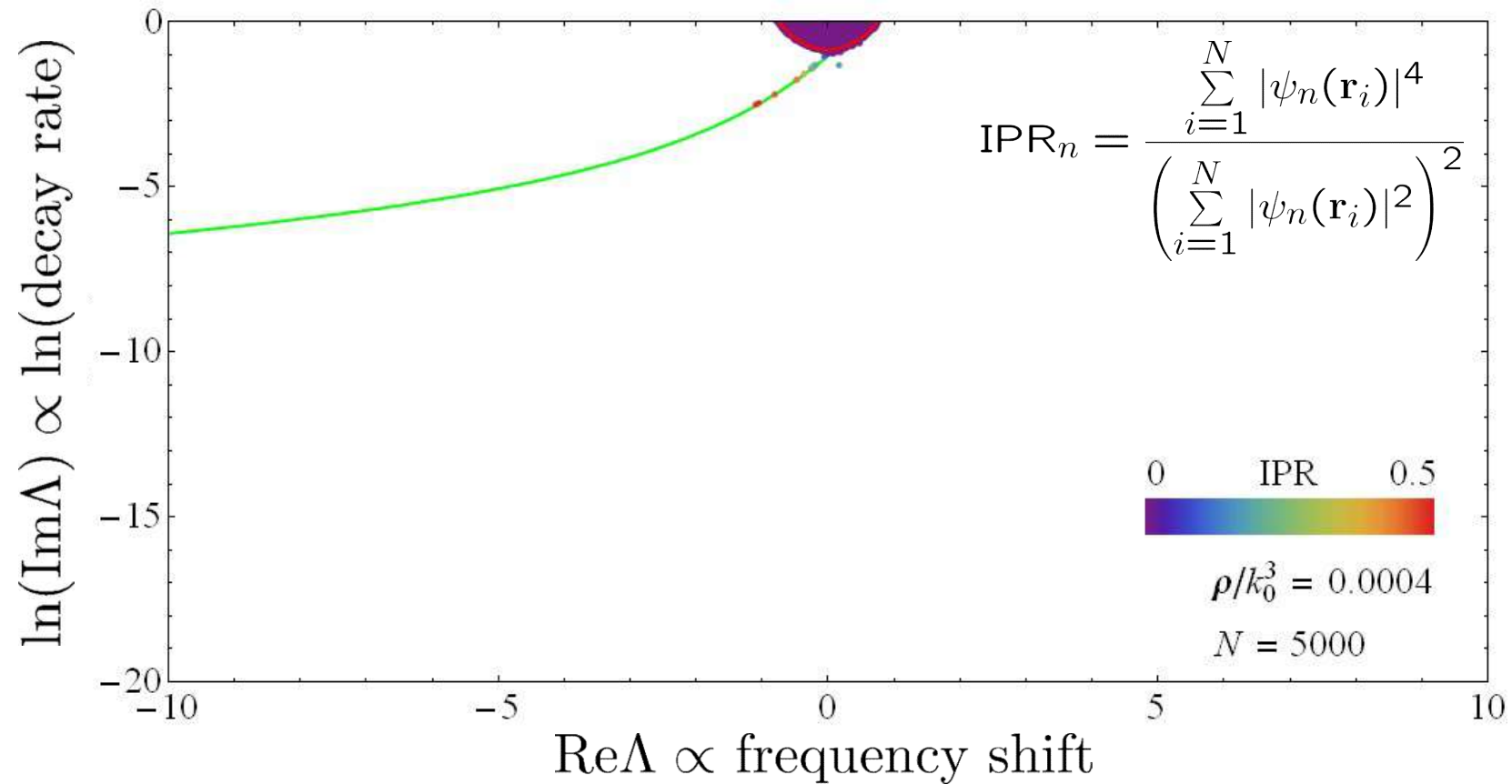
$$\text{IPR}_n = \frac{\sum_{i=1}^N |\psi_n(\mathbf{r}_i)|^4}{\left(\sum_{i=1}^N |\psi_n(\mathbf{r}_i)|^2 \right)^2} \sim \frac{1}{M} \quad \text{for a state (mode) localized on } M \text{ atoms}$$

Evolution of IPR with increasing density



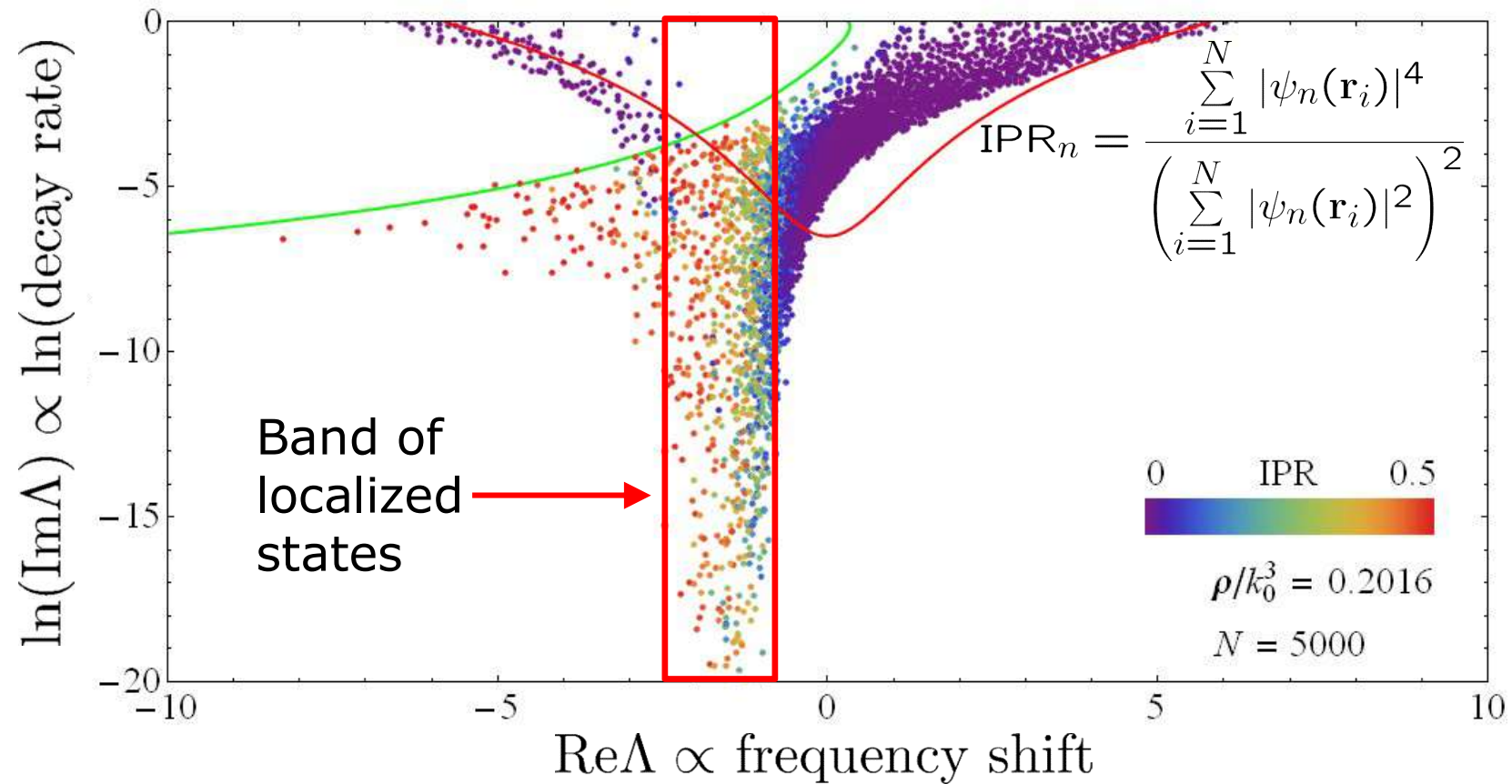
- Eigenvalue domain boundary from the diffusion theory
- Subradiant states localized on 2 closely located atoms

Evolution of IPR with increasing density



- Eigenvalue domain boundary from the diffusion theory
- Subradiant states localized on 2 closely located atoms

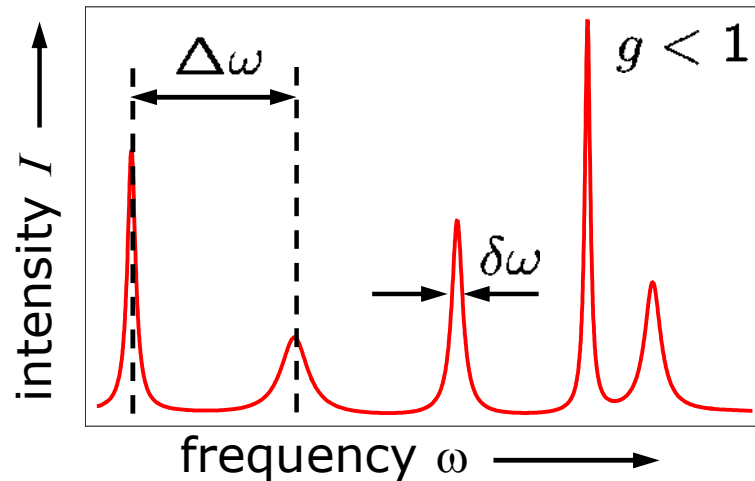
Evolution of IPR with increasing density



- Eigenvalue domain boundary from the diffusion theory
- Subradiant states localized on 2 closely located atoms

Dimensionless conductance

Spectrum of a disordered system

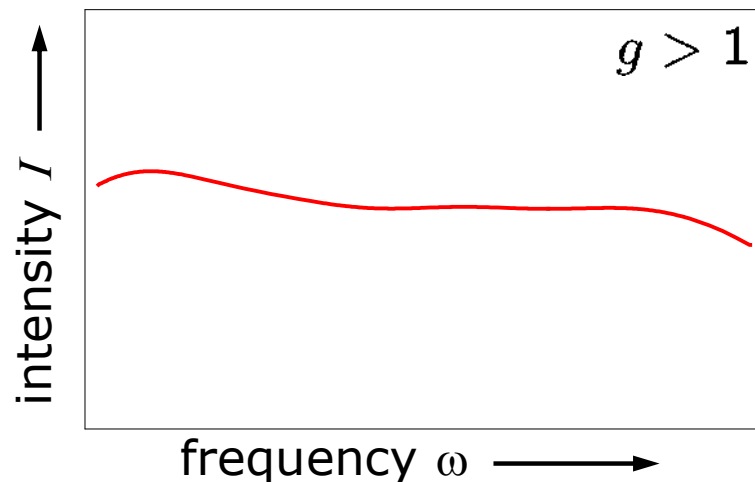


$\delta\omega \sim \text{Im}\Lambda$ — “mode width”

$\Delta\omega \sim \text{Re}\Lambda_{n+1} - \text{Re}\Lambda_n$ — “mode spacing”

$g = \frac{\delta\omega}{\Delta\omega}$ - Thouless parameter
(dimensionless conductance)

The same with mode widths x 10



Thouless criterion
of Anderson localization:

$$g < 1$$

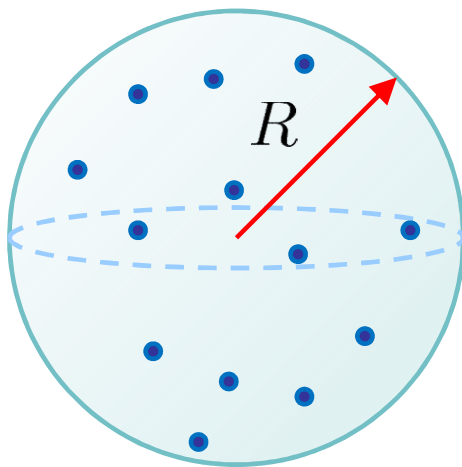
Scaling theory of Anderson localization

Main idea: Study how g evolves with sample size R

If the modes are extended, g grows with R

If the modes are localized, g decreases with R

At the critical point $g = g_c$ is independent of R



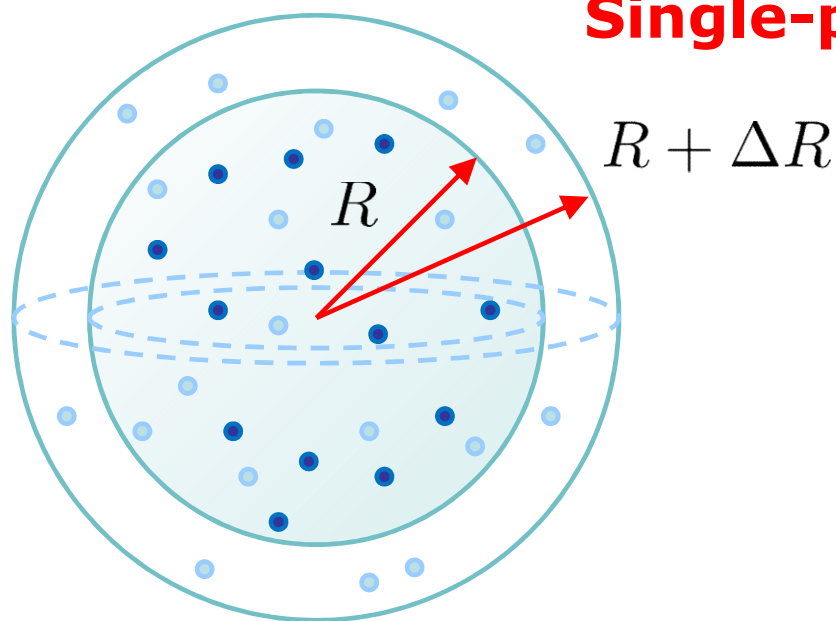
Scaling theory of Anderson localization

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At the critical point $g = g_c$ is independent of R



Single-parameter scaling hypothesis

$$\beta(g) = \frac{\partial \ln g}{\partial \ln k_0 R}$$

> 0 if extended eigenstates

< 0 if localized eigenstates

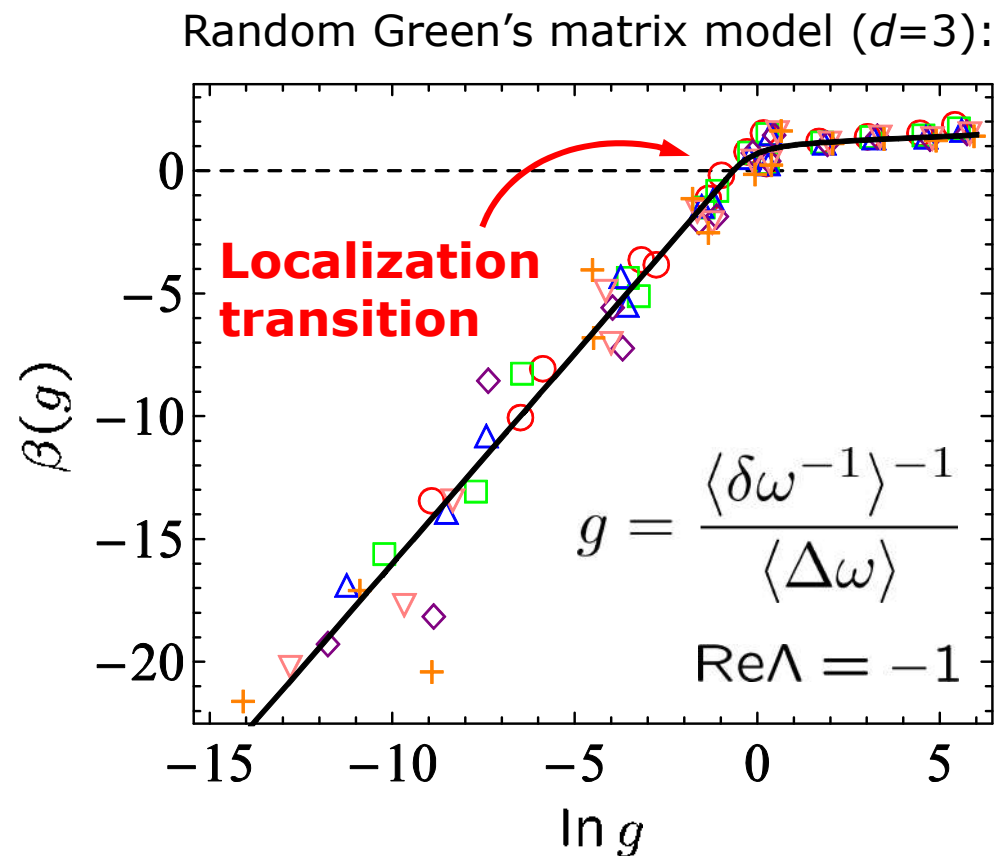
Abrahams et al., PRL **42**, 673 (1979)

Scaling theory of Anderson localization

Abrahams et al., PRL **42**, 673 (1979):

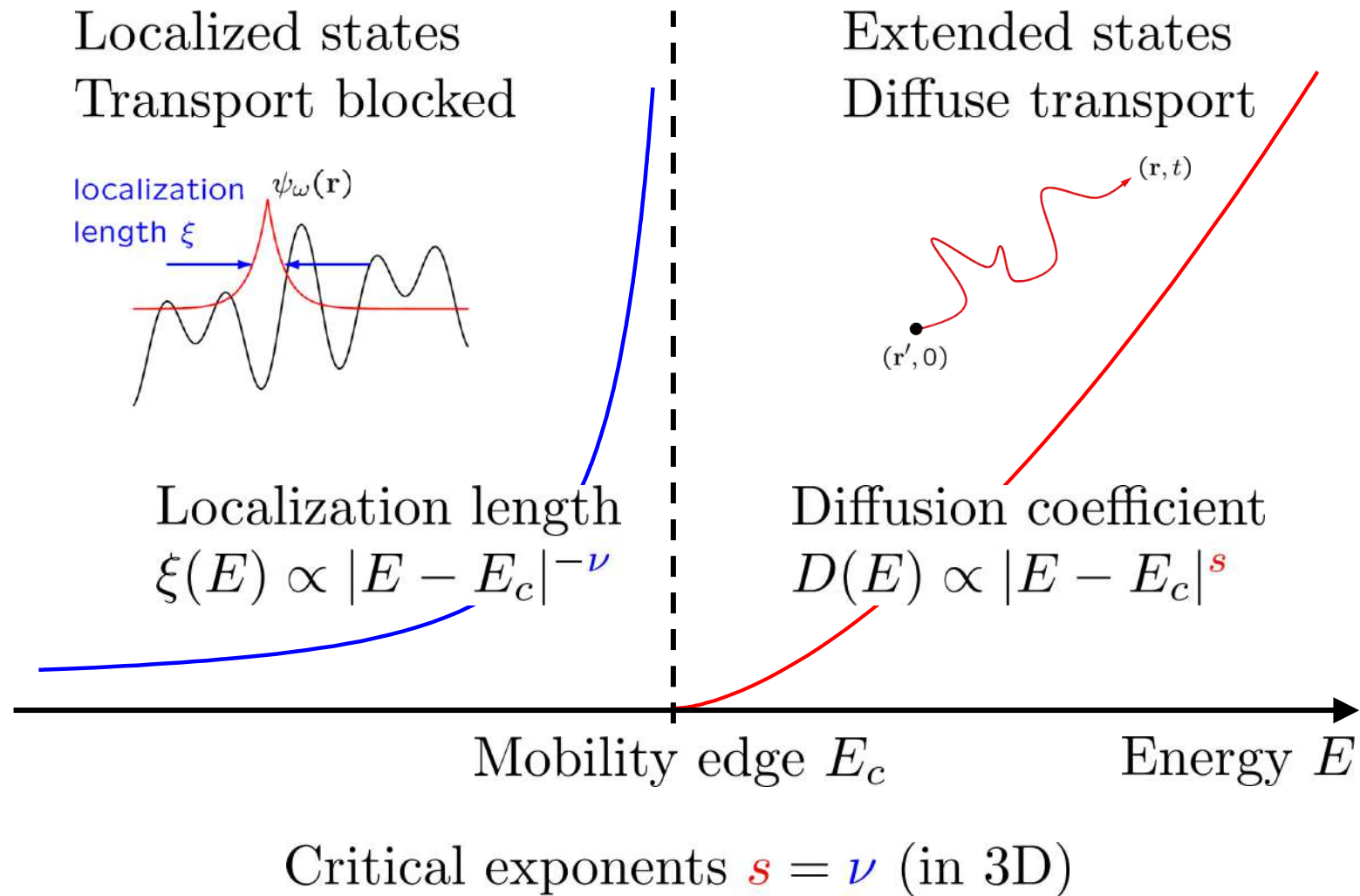
$$\beta(g) = \frac{\partial \ln g}{\partial \ln k_0 R}$$

- > 0 if extended eigenstates
- < 0 if localized eigenstates
- = 0 at the critical point



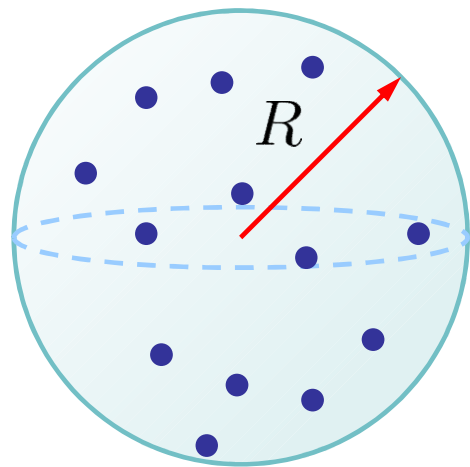
Skipetrov & Sokolov, PRL **112**, 023905 (2014)

Critical behavior around the mobility edge



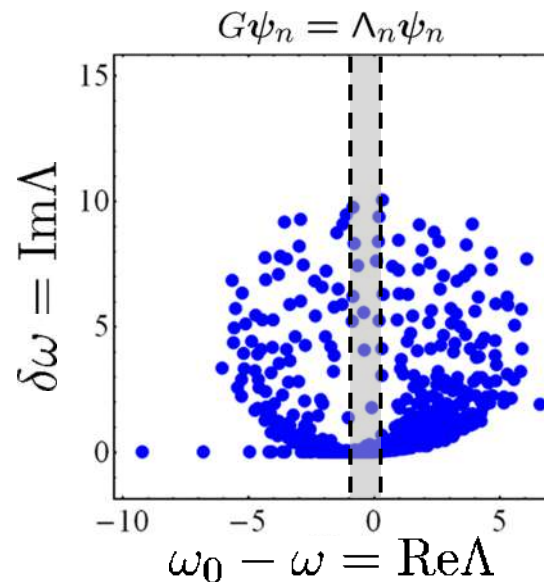
Thouless conductance & scaling

N scatterers at density ρ



$$G_{ij} = i\delta_{ij} + (1 - \delta_{ij}) \frac{e^{ik_0 r_{ij}}}{k_0 r_{ij}}$$

Eigenvalues of
the Green's matrix



Thouless conductance

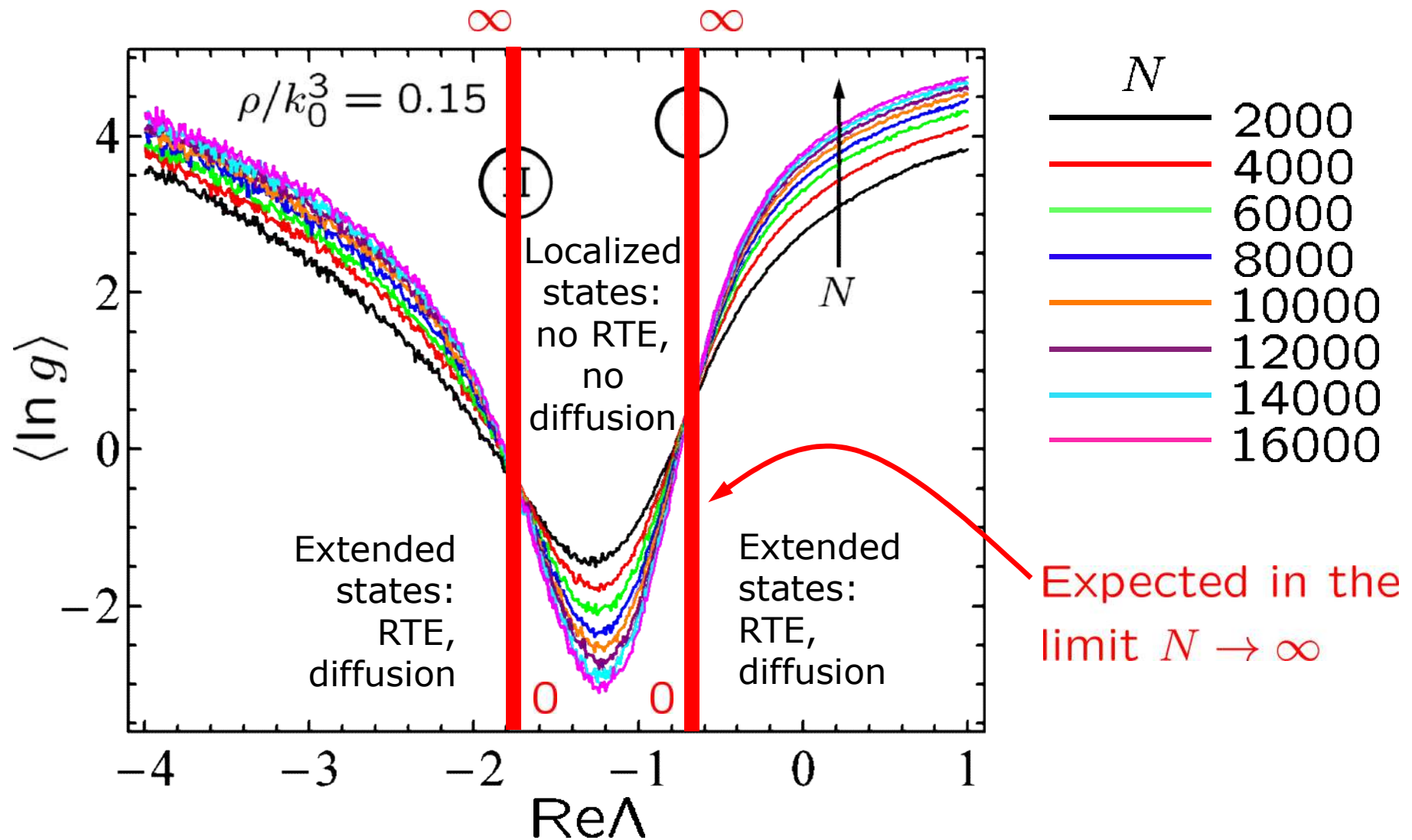
$$\delta\omega = \text{Im}\Lambda$$

$$g(\omega) = \frac{\delta\omega}{\langle \Delta\omega \rangle}$$

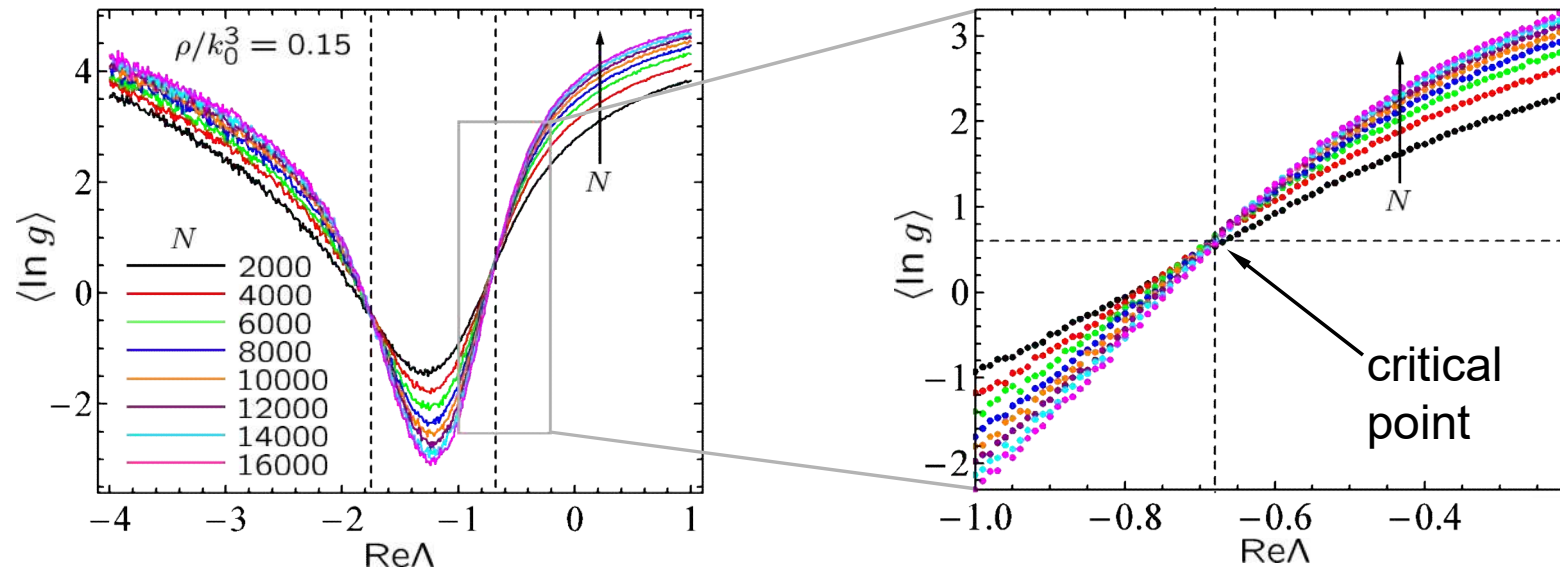
$$\langle \Delta\omega \rangle = \langle \omega_{n+1} - \omega_n \rangle$$

Now we are going to study statistical properties of $g(\omega)$ at high $\rho = 0.15k_0^3$ at which localized states are expected

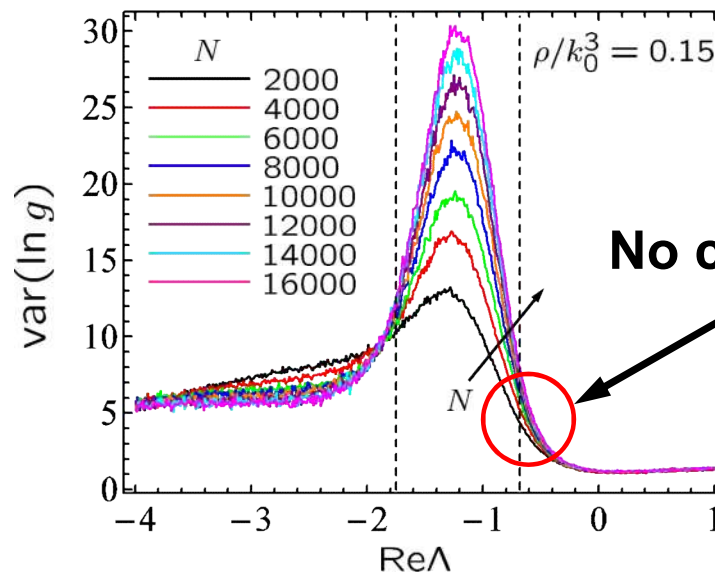
Scaling of moments of $\ln g$



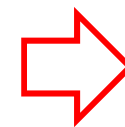
Scaling of moments of $\ln g$



But...

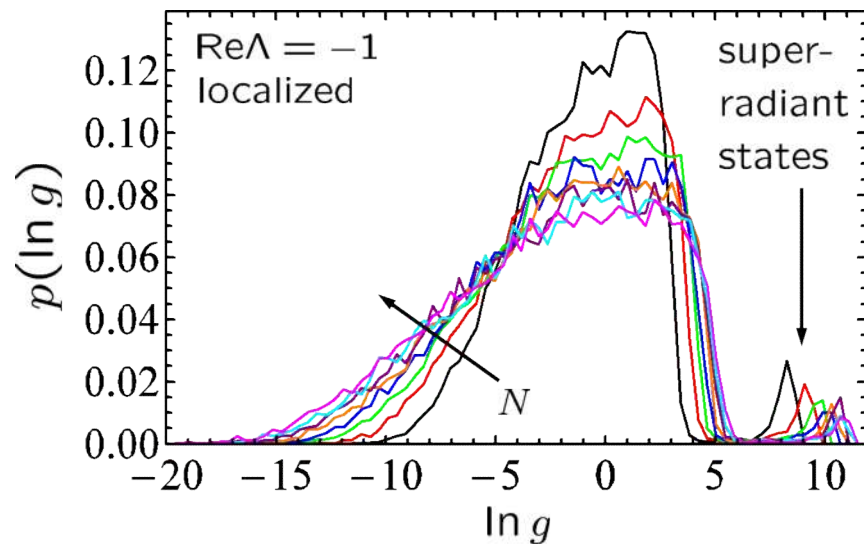
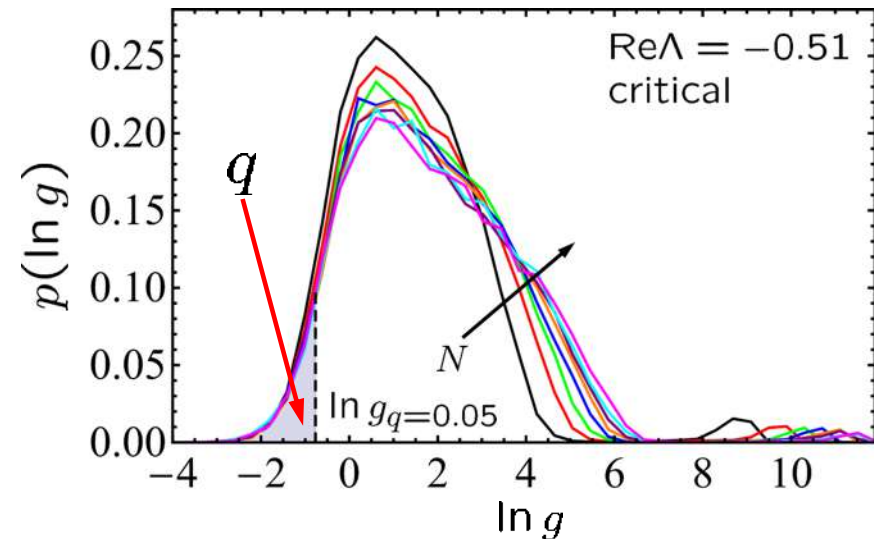
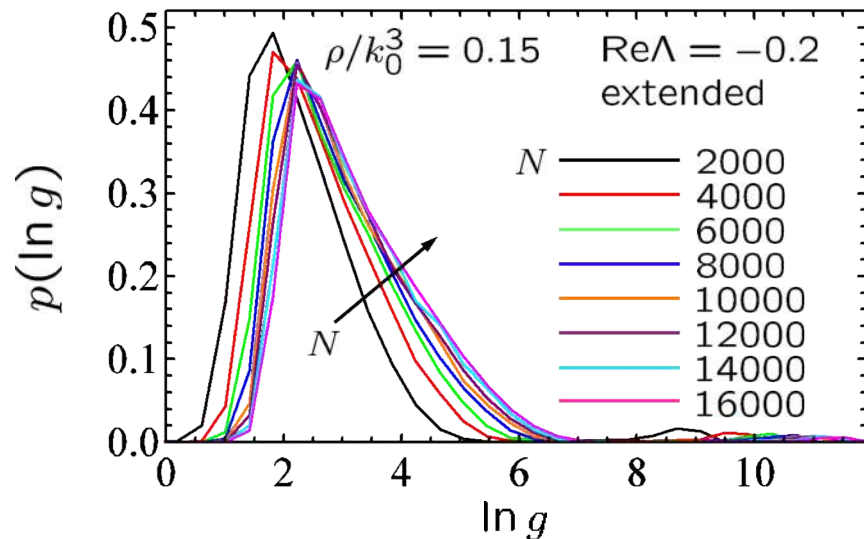


No crossing



No single-parameter scaling

Distribution of conductance



Percentile g_q :

$$q = \int_0^{g_q} p(g) dg$$

Single-parameter scaling

R — system size

 $\ln g_q = F_q(R/\xi)$

 $\xi \propto \frac{1}{(\text{Re}\Lambda - \text{Re}\Lambda_c)^\nu}$
 — localization length

$$\begin{aligned}
 \ln g_q &= F_q(R/\xi) = F_q[R(\text{Re}\Lambda - \text{Re}\Lambda_c)^\nu] \\
 &= F_q[R^{1/\nu}(\text{Re}\Lambda - \text{Re}\Lambda_c)] \longrightarrow F_q(\psi, \phi)
 \end{aligned}$$

Relevant scaling variable:

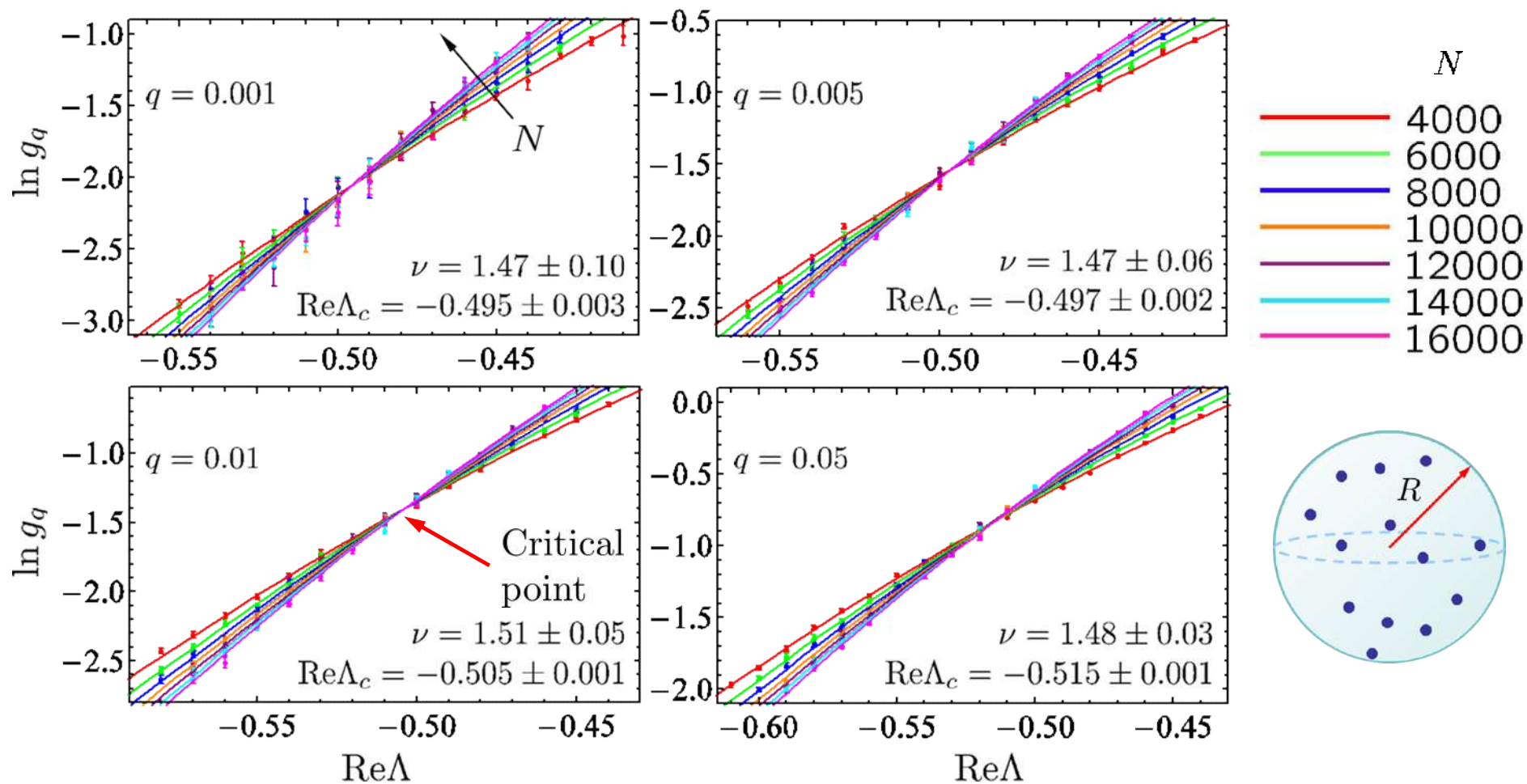
$$\psi = R^{1/\nu} u(\text{Re}\Lambda - \text{Re}\Lambda_c), \quad u(x) = u_1 x + u_2 x^2 + \dots$$

Irrelevant scaling variable:

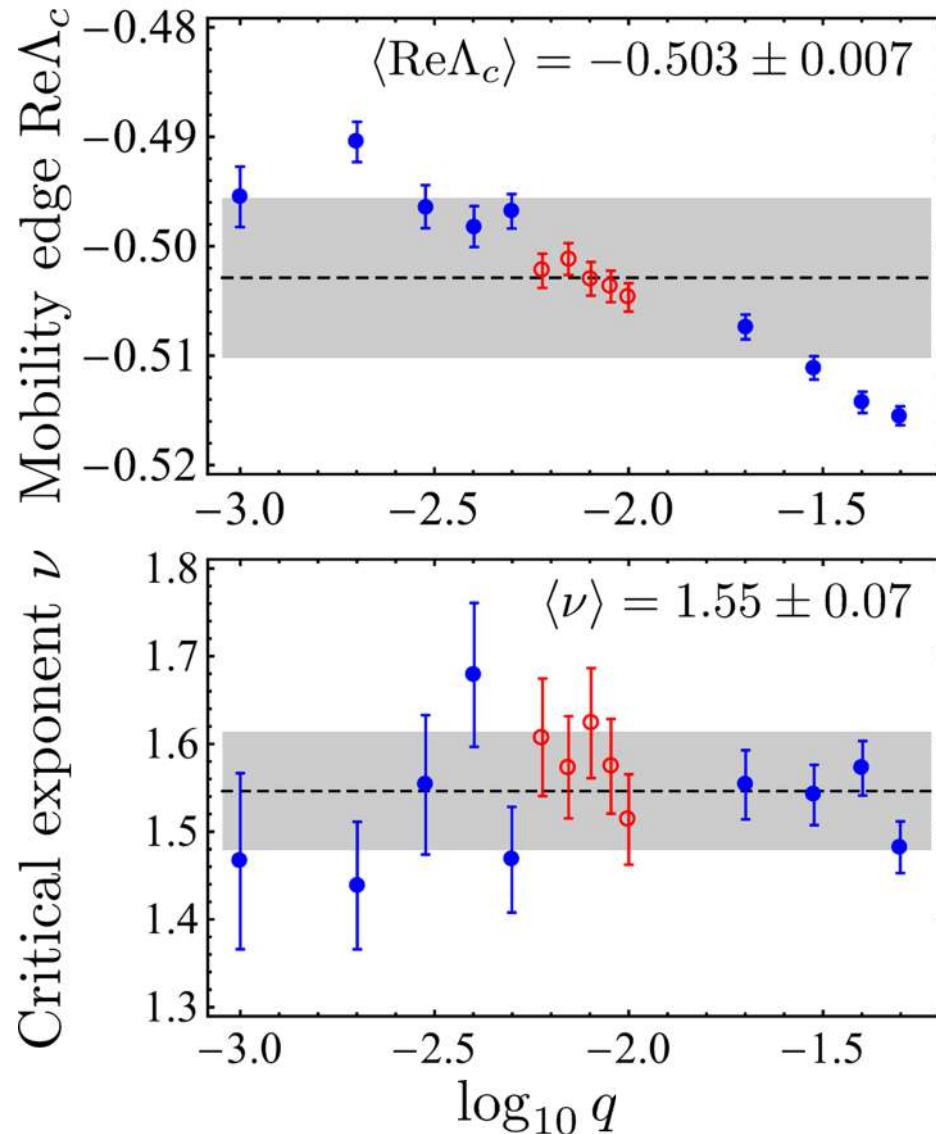
$$\phi = R^{-y} v(\text{Re}\Lambda - \text{Re}\Lambda_c), \quad v(x) = v_0 + v_1 x + v_2 x^2 + \dots$$

Finite-size scaling of percentiles

$$\ln g_q = F_q(R/\xi), \quad \xi \propto (\text{Re}\Lambda - \text{Re}\Lambda_c)^{-\nu}$$



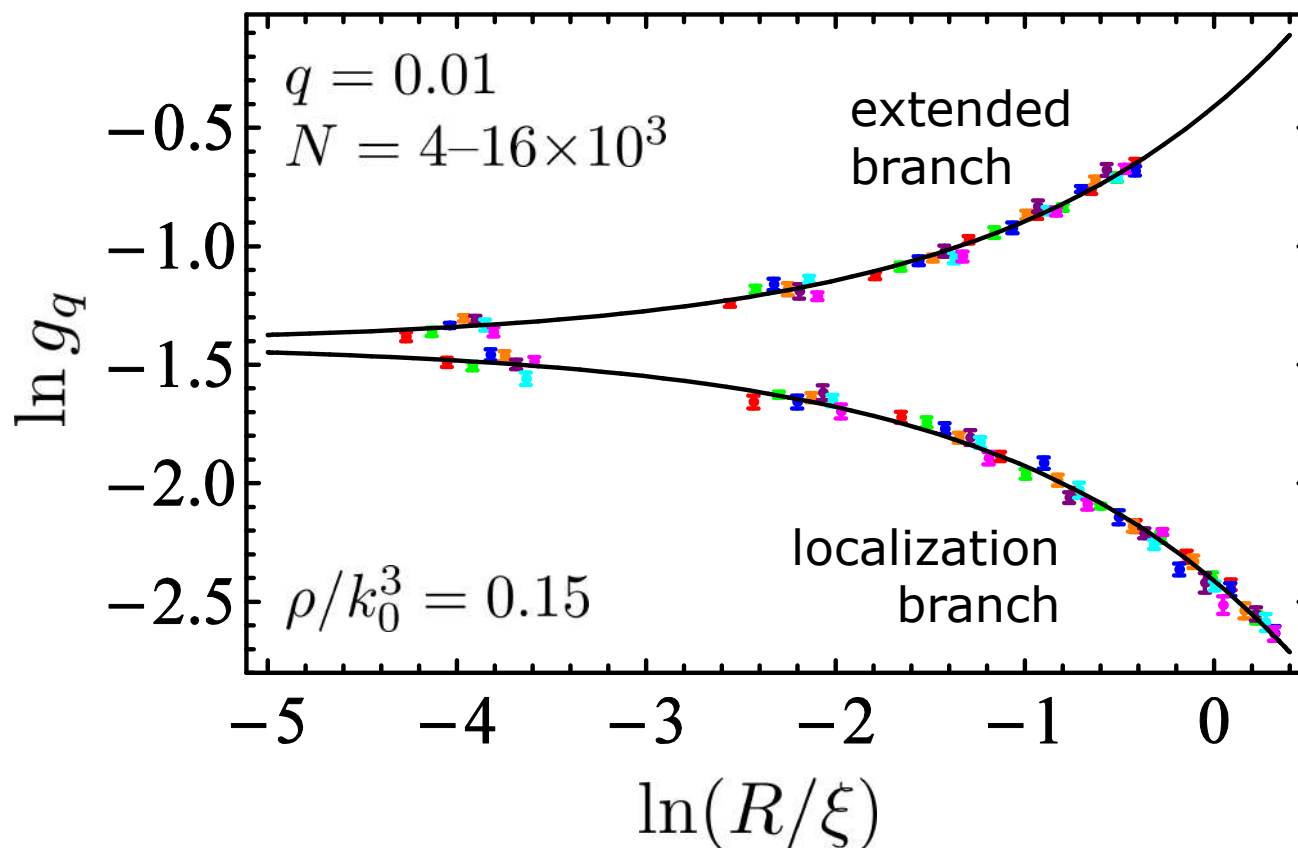
Best-fit parameters



The value of critical exponent following from the fits is close to $\nu \simeq 1.57$ expected for the 3D orthogonal symmetry class.

We conclude that the observed transition is likely to belong to the same symmetry class as the **Anderson transition in a system of spinless electrons**.

Single-parameter scaling: a posteriori check

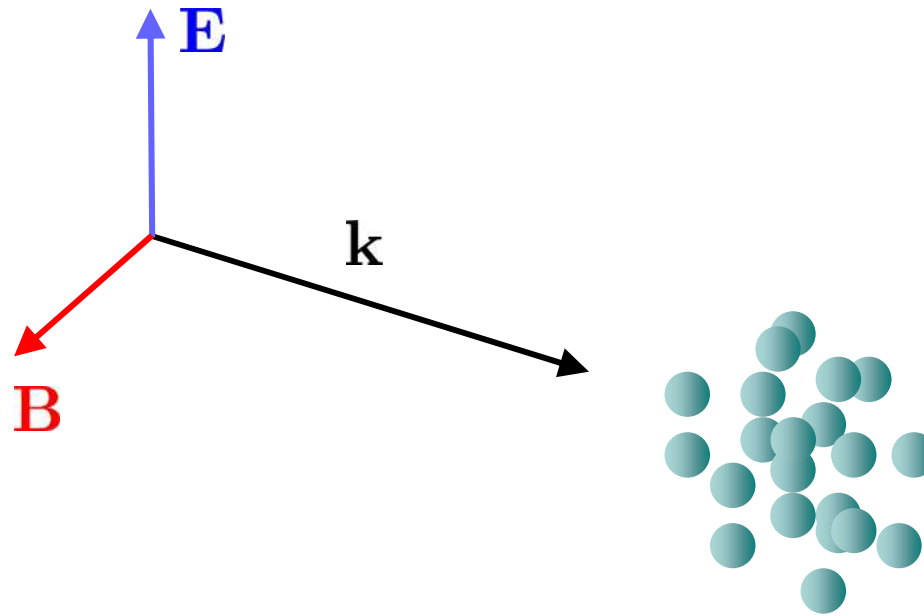


Data for different N all fall on a single line $\ln g_q = F_q(R/\xi)$

R — sample size

ξ — localization (correlation) length

Anderson localization of light



Pioneering theoretical works: John, PRL **53**, 2169 (1984)

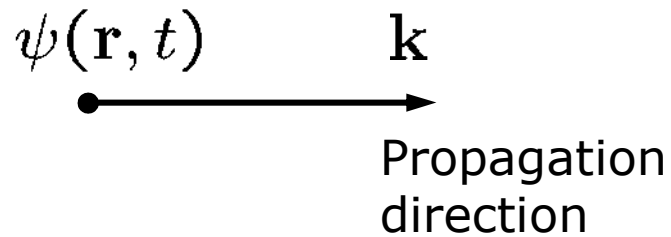
Anderson, Philos. Mag. B **52**, 505 (1985)

Experiments inconclusive: Wiersma et al., Nature **390**, 671 (1997)

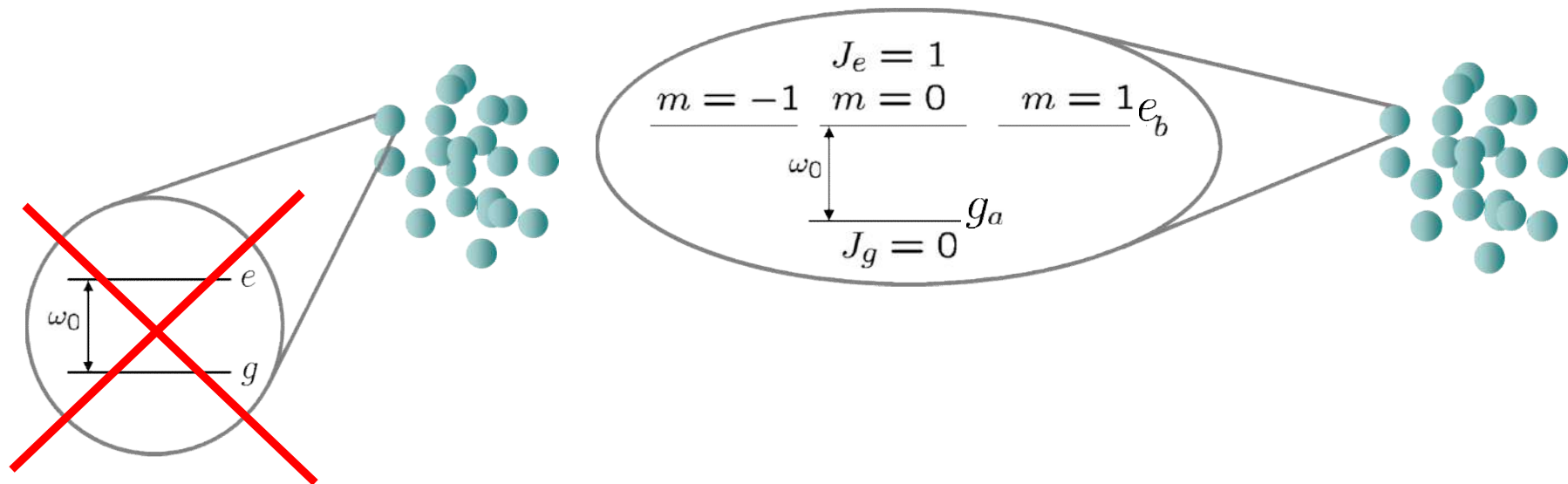
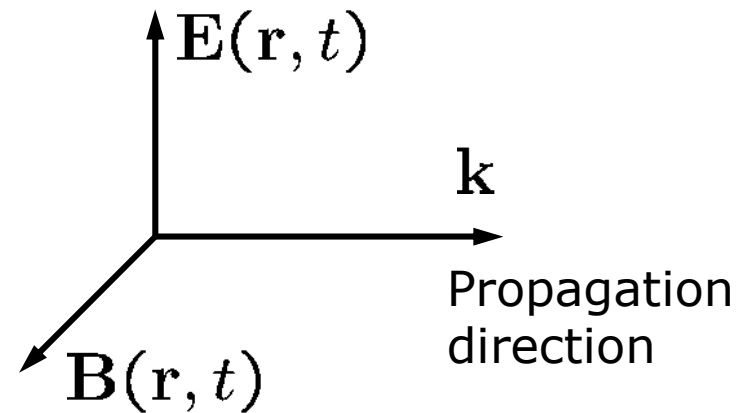
Sperling et al., Nat. Photonics **7**, 48 (2013)

Light is a vector wave

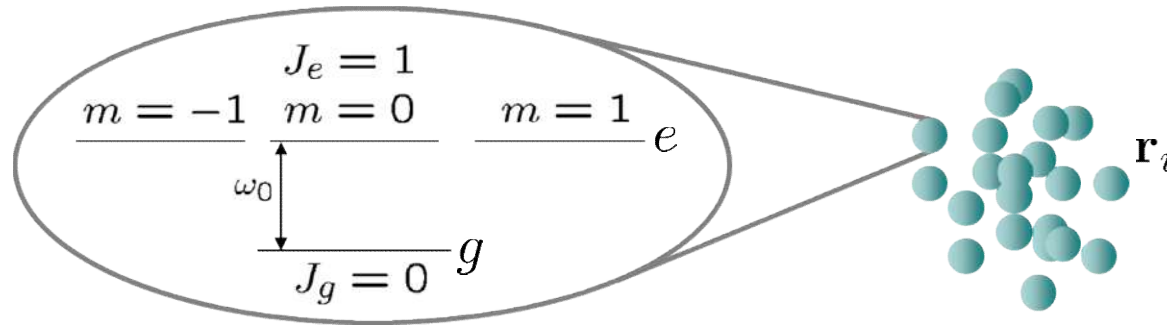
**“Schrödinger” waves
or sound**



Electromagnetic waves



Familiar textbook Hamiltonian...



isolated atoms

free electromagnetic field

$$\hat{H} = \underbrace{\sum_{i=1}^N \sum_{m=-1}^1 \hbar \omega_0 |e_{im}\rangle \langle e_{im}|}_{\text{isolated atoms}} + \underbrace{\sum_{\mathbf{s} \perp \mathbf{k}} \hbar c k \left(\hat{a}_{\mathbf{k}\mathbf{s}}^\dagger \hat{a}_{\mathbf{k}\mathbf{s}} + \frac{1}{2} \right)}_{\text{free electromagnetic field}}$$

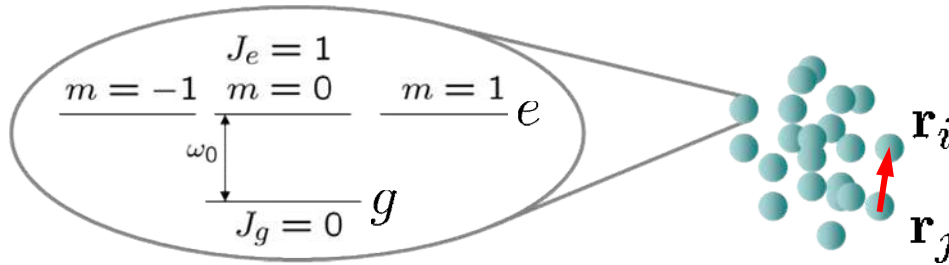
$$- \underbrace{\sum_{i=1}^N \hat{\mathbf{D}}_i \cdot \hat{\mathbf{E}}(\mathbf{r}_i)}_{\text{atom-field interaction}} + \underbrace{\frac{1}{2\epsilon_0} \sum_{i \neq j}^N \hat{\mathbf{D}}_i \cdot \hat{\mathbf{D}}_j \delta(\mathbf{r}_i - \mathbf{r}_j)}_{\text{contact term}}$$

atom-field interaction
in the dipole approximation

contact term

Cohen-Tannoudji, Dupont-Roc & Grynberg,
Photons and Atoms: Introduction to Quantum Electrodynamics (1992)
Morice, Castin & Dalibard, PRA **51**, 3896 (1995)

Green's matrix for light



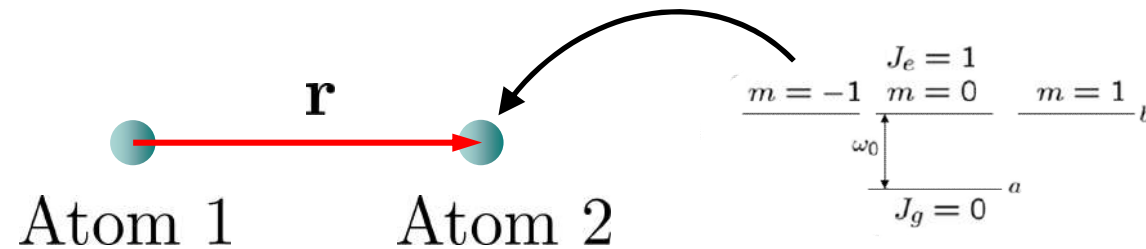
Green's matrix G describes propagation of light between pairs of atoms
 $r_{ij} = |\mathbf{r}_i - \mathbf{r}_j|$

$$G_{e_{im}e_{jm'}} = i\delta_{e_{im}e_{jm'}} - \frac{2}{\Gamma_0}(1 - \delta_{e_{im}e_{jm'}}) \sum_{\mu,\nu} d_{e_{im}g_i}^\mu d_{g_j e_{jm'}}^\nu \frac{e^{ik_0 r_{ij}}}{\hbar r_{ij}^3} \\ \times \left\{ \delta_{\mu\nu} \left[1 - ik_0 r_{ij} - (k_0 r_{ij})^2 \right] - \frac{r_{ij}^\mu r_{ij}^\nu}{r_{ij}^2} \left[3 - 3ik_0 r_{ij} - (k_0 r_{ij})^2 \right] \right\}$$

$$\mathbf{d}_{e_{im}g_i} = \langle J_e m | \hat{\mathbf{D}}_i | J_g 0 \rangle$$

($3N \times 3N$ matrix)

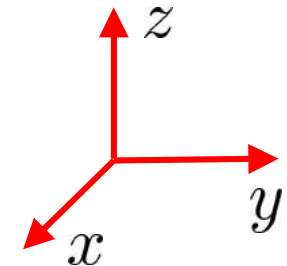
Green's function in different bases



$$G_{\mu\nu}(\mathbf{r}) = \frac{3}{2} \frac{e^{ik_0 r}}{k_0 r} \left[P(ik_0 r) \delta_{\mu\nu} + Q(ik_0 r) \frac{r_\mu r_\nu}{r^2} \right] \quad \text{linear basis}$$

$$\mu, \nu = x, y, z$$

$$P(x) = 1 - 1/x + 1/x^2, \quad Q(x) = -1 + 3/x - 3/x^2$$

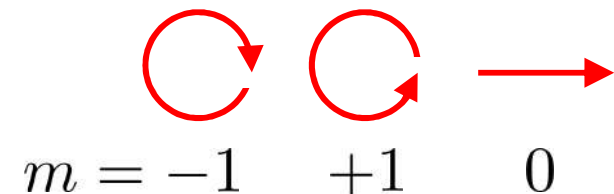


$$G_{mm'}(\mathbf{r}) = -\frac{4k_0^3}{3\hbar\Gamma_0} \sum_{\mu,\nu} d_{m\mu} G_{\mu\nu}(\mathbf{r}) d_{m'\nu}$$

$$m, m' = -1, 0, 1$$

$$\mathbf{d}_m = \langle J_e m | \hat{\mathbf{D}} | J_g 0 \rangle$$

spiral basis



Structure of the Green's matrix

$$\begin{pmatrix} \boxed{\begin{matrix} i & 0 & 0 \\ 0 & i & 0 \\ 0 & 0 & i \end{matrix}} & \begin{matrix} G_{12}^{xx} & G_{12}^{xy} & G_{12}^{xz} \\ G_{12}^{yx} & G_{12}^{yy} & G_{12}^{yz} \\ G_{12}^{zx} & G_{12}^{zy} & G_{12}^{zz} \end{matrix} & \dots & \begin{matrix} G_{1N}^{xx} & G_{1N}^{xy} & G_{1N}^{xz} \\ G_{1N}^{yx} & G_{1N}^{yy} & G_{1N}^{yz} \\ G_{1N}^{zx} & G_{1N}^{zy} & G_{1N}^{zz} \end{matrix} \\ \begin{matrix} G_{21}^{xx} & G_{21}^{xy} & G_{21}^{xz} \\ G_{21}^{yx} & G_{21}^{yy} & G_{21}^{yz} \\ G_{21}^{zx} & G_{21}^{zy} & G_{21}^{zz} \end{matrix} & \boxed{\begin{matrix} i & 0 & 0 \\ 0 & i & 0 \\ 0 & 0 & i \end{matrix}} & \dots & \dots \\ \dots & \dots & \dots & \dots \\ \dots & \dots & \dots & \dots \\ \begin{matrix} G_{N1}^{xx} & G_{N1}^{xy} & G_{N1}^{xz} \\ G_{N1}^{yx} & G_{N1}^{yy} & G_{N1}^{yz} \\ G_{N1}^{zx} & G_{N1}^{zy} & G_{N1}^{zz} \end{matrix} & \dots & \dots & \boxed{\begin{matrix} i & 0 & 0 \\ 0 & i & 0 \\ 0 & 0 & i \end{matrix}} \end{pmatrix}$$

One-atom dynamics:

Excitation of an isolated excited atom decays as $e^{-\Gamma_0 t}$

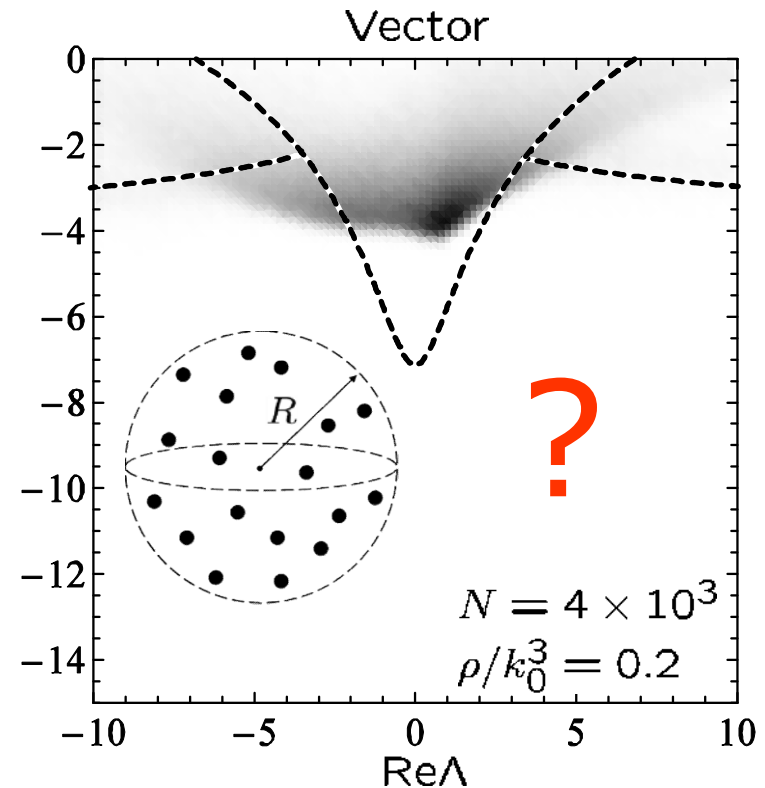
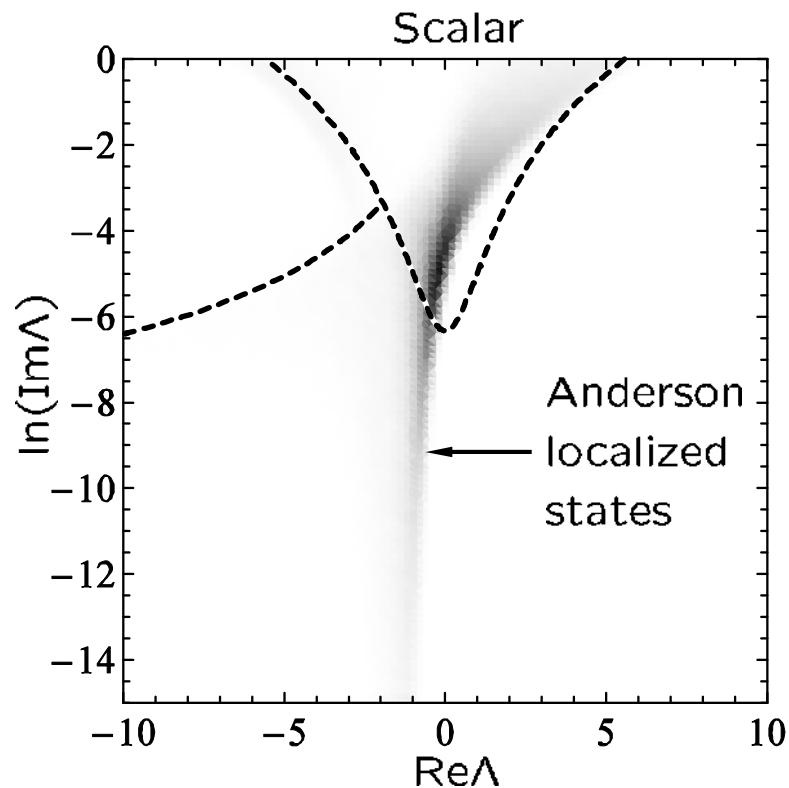
Structure of the Green's matrix

$$\begin{pmatrix}
 i & 0 & 0 & G_{12}^{xx} & G_{12}^{xy} & G_{12}^{xz} & \dots & G_{1N}^{xx} & G_{1N}^{xy} & G_{1N}^{xz} \\
 0 & i & 0 & G_{12}^{yx} & G_{12}^{yy} & G_{12}^{yz} & \dots & G_{1N}^{yx} & G_{1N}^{yy} & G_{1N}^{yz} \\
 0 & 0 & i & G_{12}^{zx} & G_{12}^{zy} & G_{12}^{zz} & \dots & G_{1N}^{zx} & G_{1N}^{zy} & G_{1N}^{zz} \\
 G_{21}^{xx} & G_{21}^{xy} & G_{21}^{xz} & i & 0 & 0 & \dots & \dots & \dots & \dots \\
 G_{21}^{yx} & G_{21}^{yy} & G_{21}^{yz} & 0 & i & 0 & \dots & \dots & \dots & \dots \\
 G_{21}^{zx} & G_{21}^{zy} & G_{21}^{zz} & 0 & 0 & i & \dots & \dots & \dots & \dots \\
 \dots & \dots & \dots & \dots & \dots & \dots & \dots & \dots & \dots & \dots \\
 G_{N1}^{xx} & G_{N1}^{xy} & G_{N1}^{xz} & \dots & \dots & \dots & \dots & i & 0 & 0 \\
 G_{N1}^{yx} & G_{N1}^{yy} & G_{N1}^{yz} & \dots & \dots & \dots & \dots & 0 & i & 0 \\
 G_{N1}^{zx} & G_{N1}^{zy} & G_{N1}^{zz} & \dots & \dots & \dots & \dots & 0 & 0 & i
 \end{pmatrix}$$

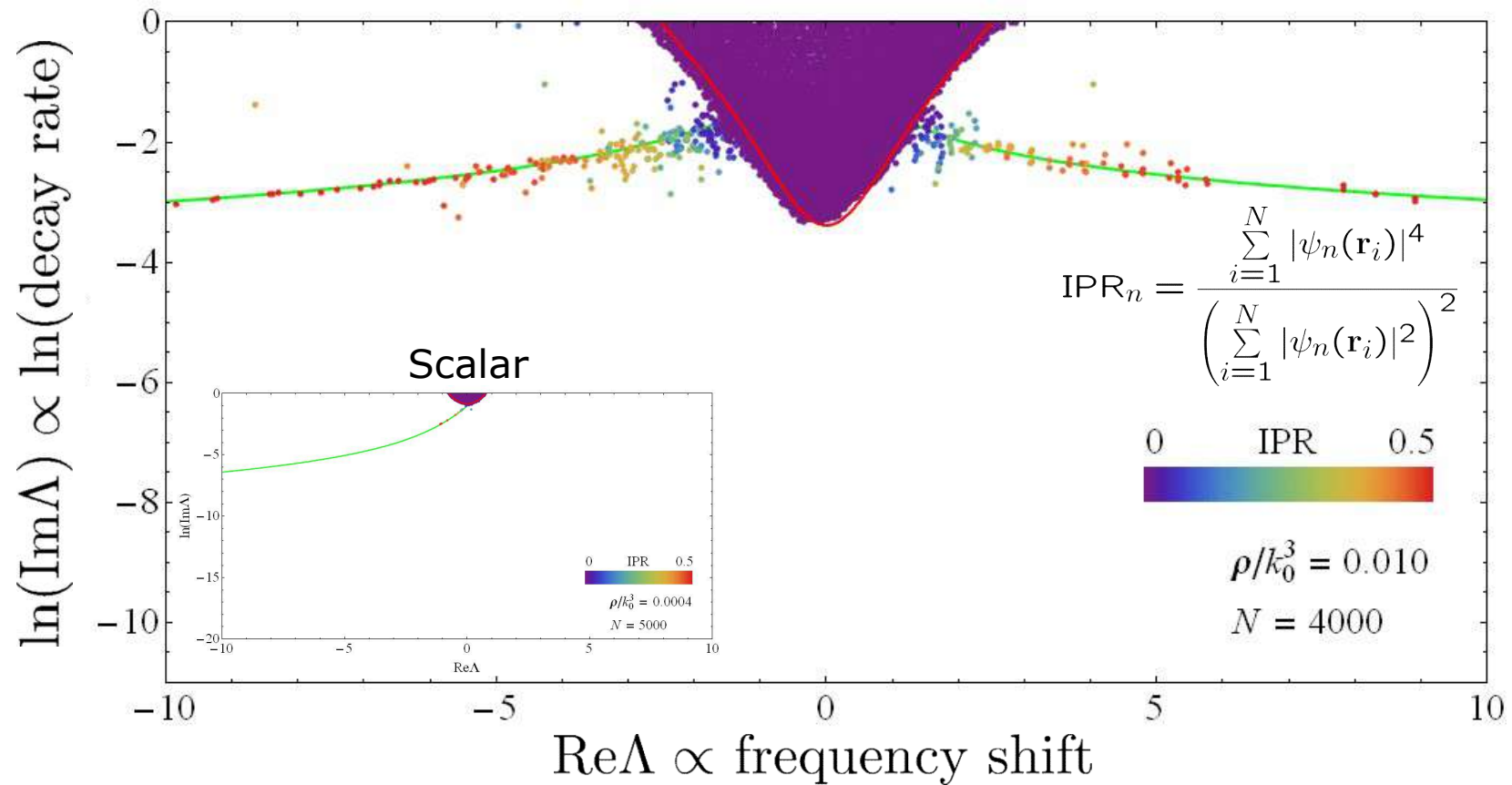
Pairwise coupling between atoms 1 & 2:

G_{12}^{xy} is the y component of the field at position 2 due to a dipole oscillating along x at position 1

Probability density of eigenvalues

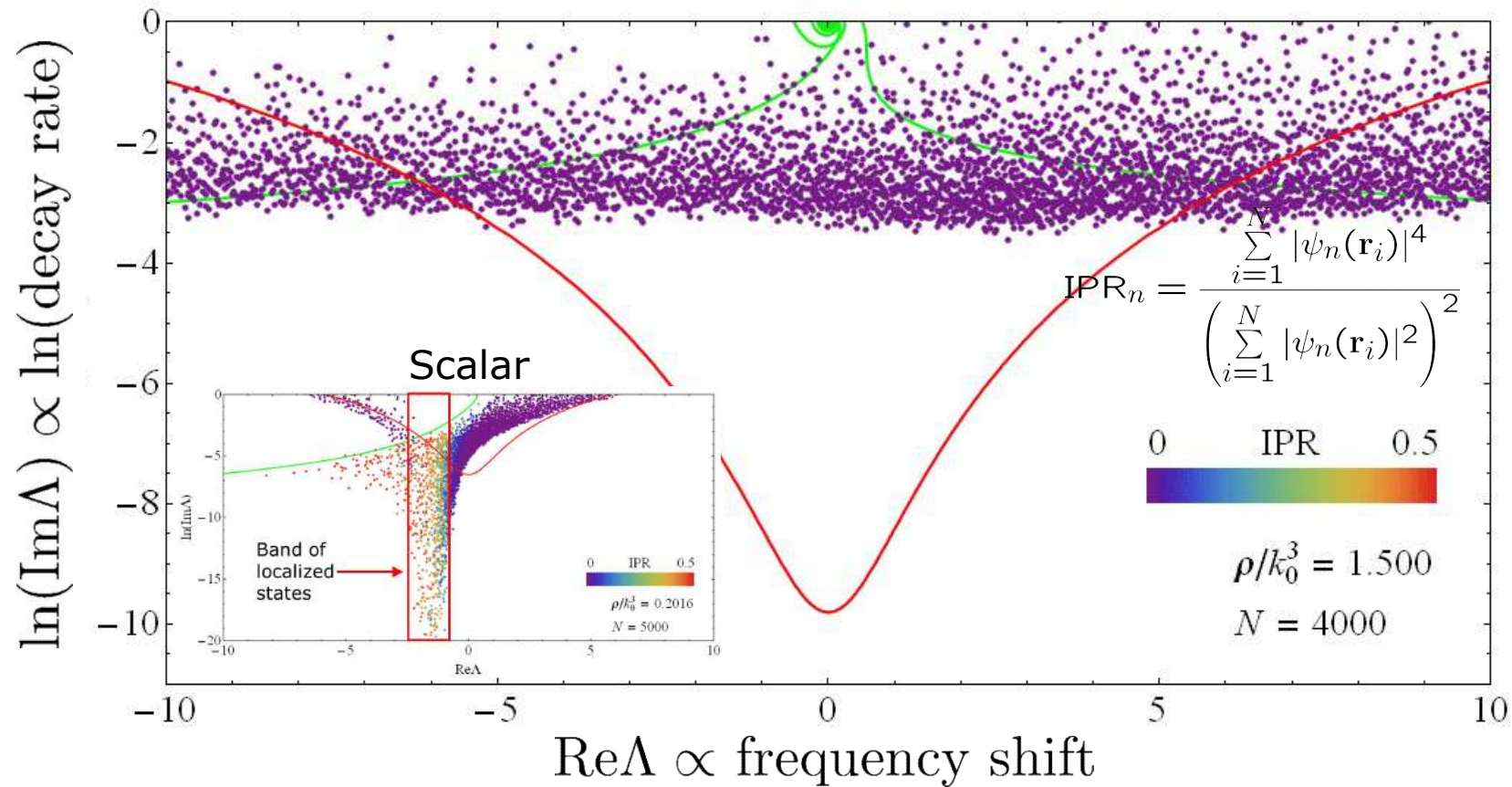


IPR for light



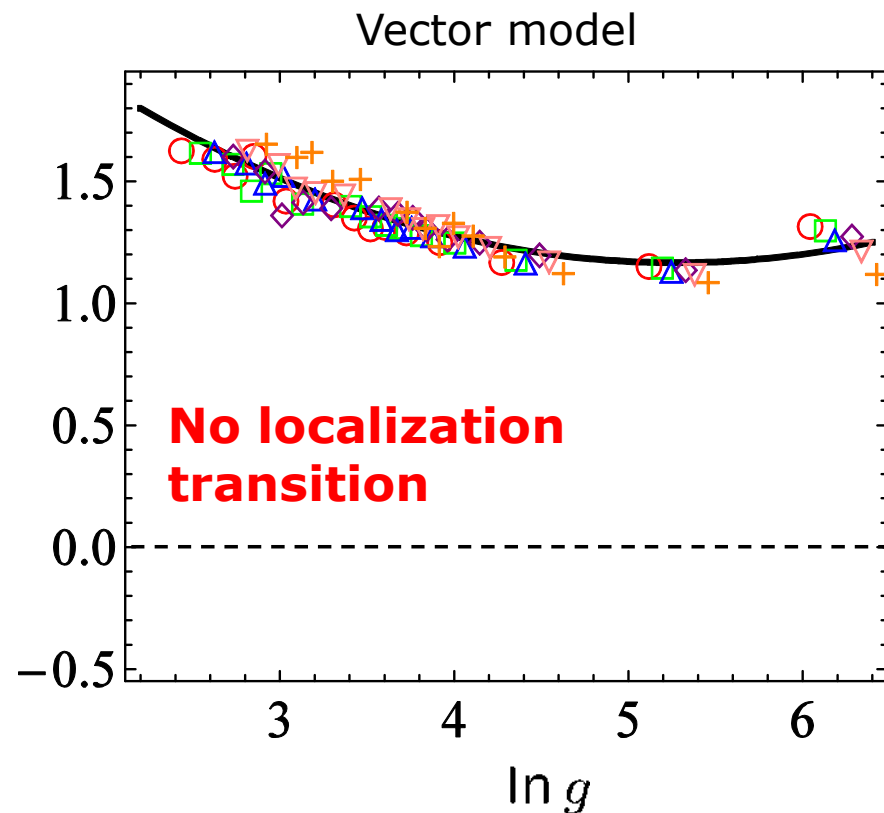
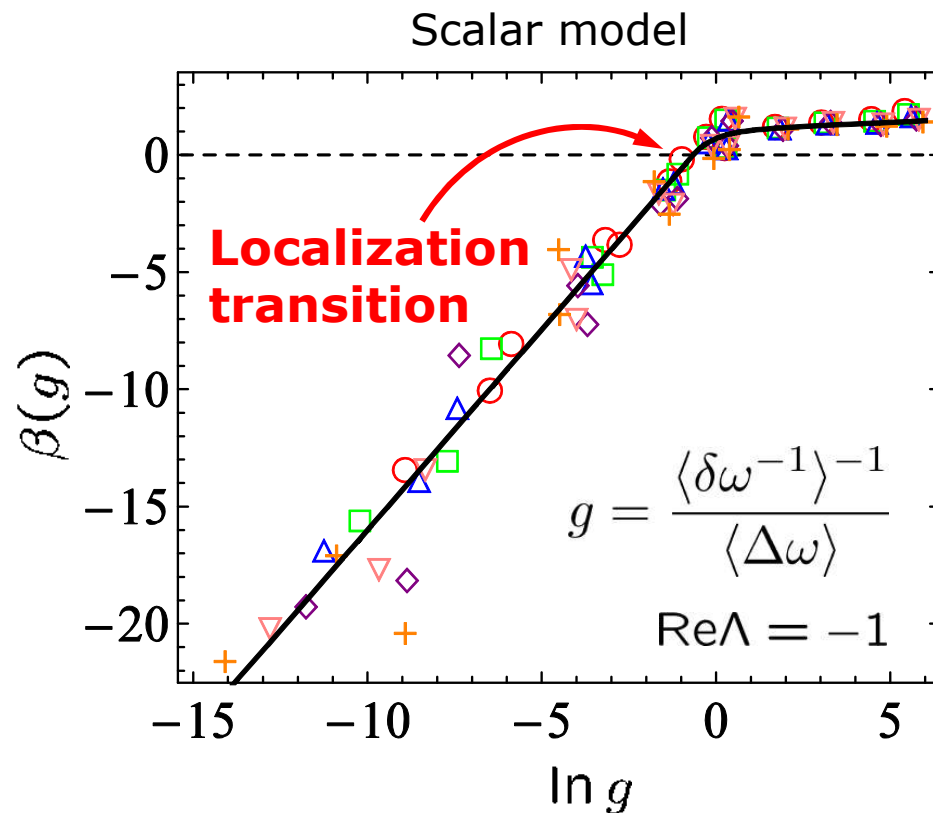
- Eigenvalue domain boundary from the diffusion theory
- Subradiant states localized on 2 closely located atoms

IPR for light



- Eigenvalue domain boundary from the diffusion theory
- Subradiant states localized on 2 closely located atoms

No Anderson localization for light in 3D

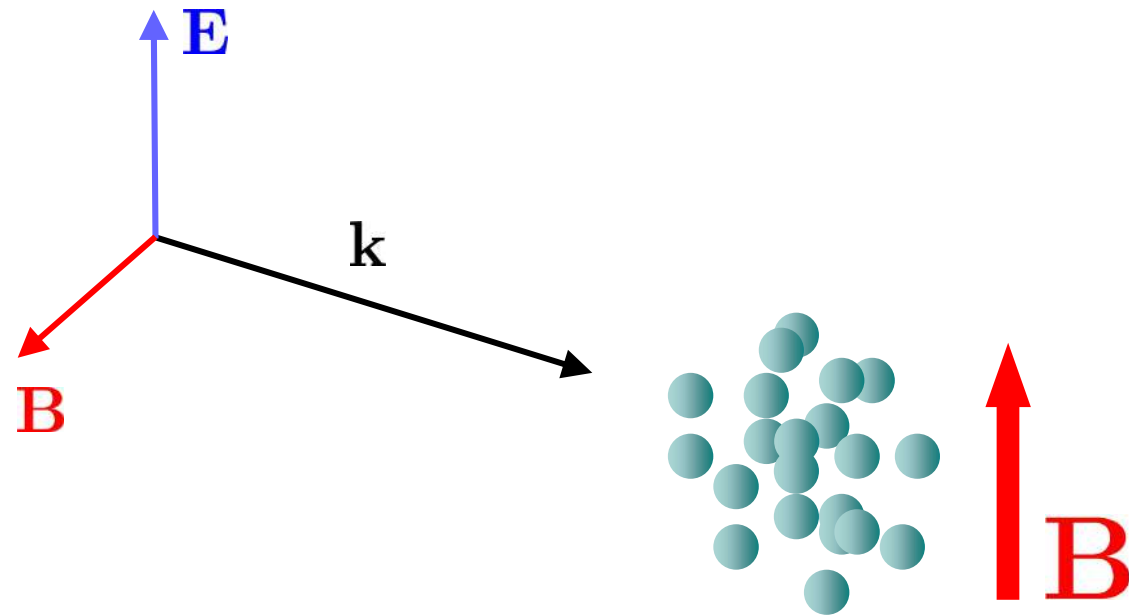


Explanation stems from near-field effects:

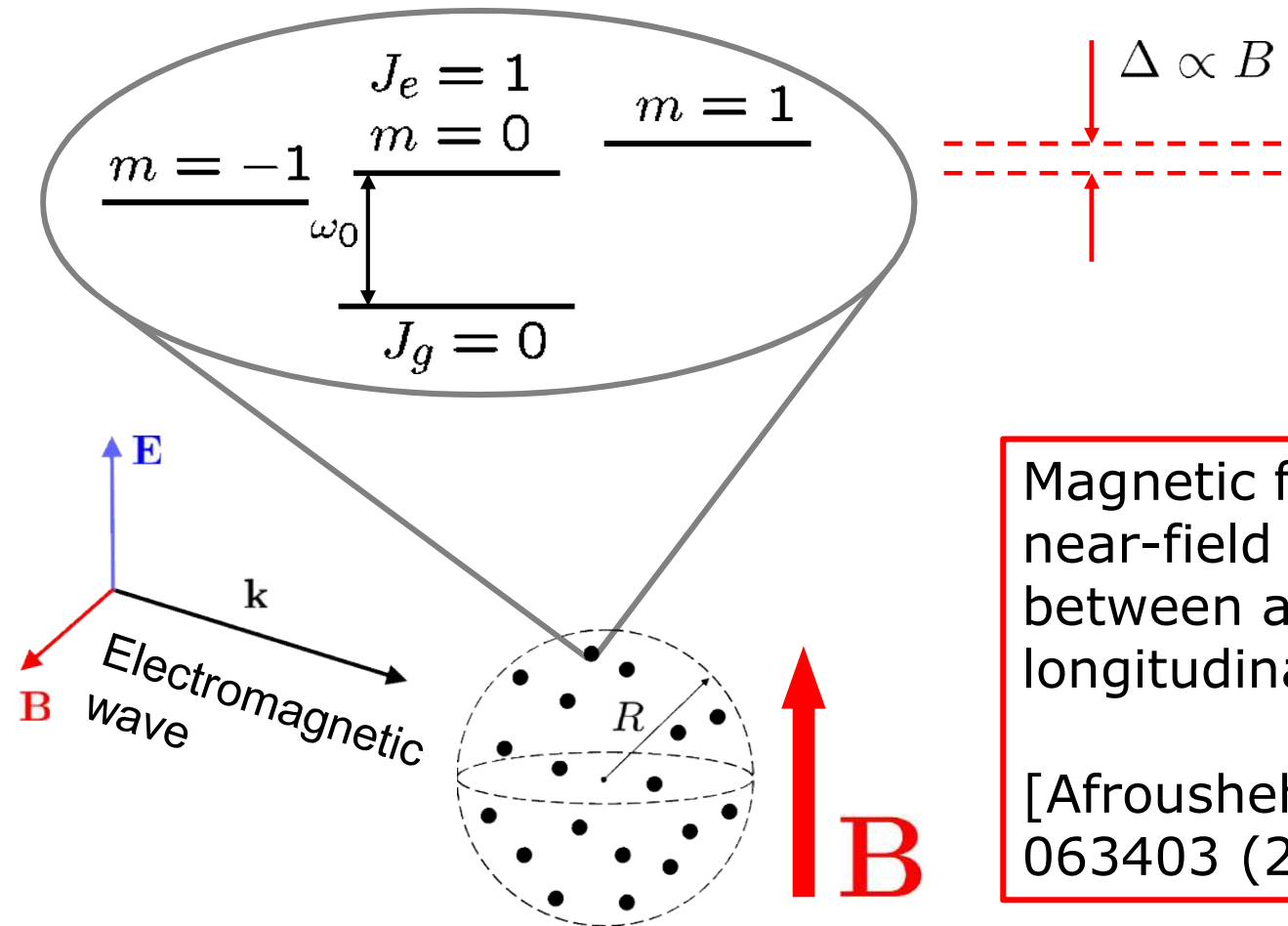
$$G_{\text{scalar}}(\mathbf{r})|_{r \rightarrow 0} \propto \frac{1}{r}$$

$$\hat{G}_{\text{EM}}(\mathbf{r})|_{r \rightarrow 0} \propto \frac{1}{r^3}$$

Anderson localization of light in a magnetic field



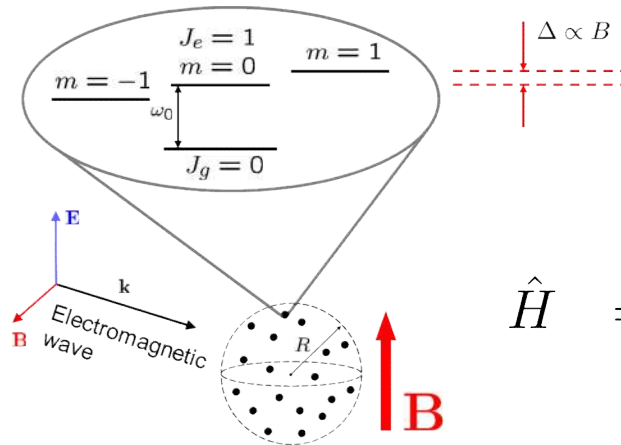
Zeeman effect



Magnetic field suppresses near-field coupling between atoms by longitudinal fields

[Afrousheh et al., PRA **73**, 063403 (2006)]

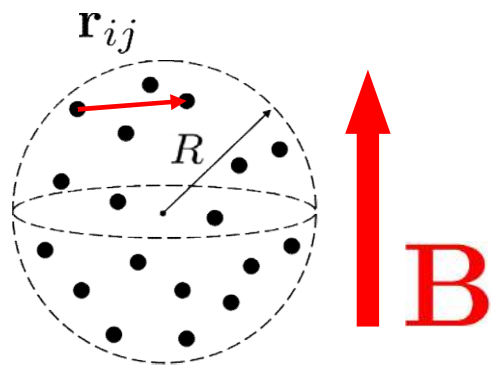
The Hamiltonian



$$\begin{aligned}
 \hat{H} = & \sum_{i=1}^N \sum_{m=-1}^1 \hbar \omega_0 |e_{im}\rangle \langle e_{im}| + \sum_{\mathbf{s} \perp \mathbf{k}} \hbar c k \left(\hat{a}_{\mathbf{k}\mathbf{s}}^\dagger \hat{a}_{\mathbf{k}\mathbf{s}} + \frac{1}{2} \right) \\
 & - \sum_{i=1}^N \hat{\mathbf{D}}_i \cdot \hat{\mathbf{E}}(\mathbf{r}_i) + \frac{1}{2\epsilon_0} \sum_{i \neq j}^N \hat{\mathbf{D}}_i \cdot \hat{\mathbf{D}}_j \delta(\mathbf{r}_i - \mathbf{r}_j) \\
 & + \underbrace{g_e \mu_B B \sum_{i=1}^N \sum_{m=-1}^1 m |e_{im}\rangle \langle e_{im}|}_{\text{Coupling of atoms with a magnetic field}}
 \end{aligned}$$

Coupling of atoms with a magnetic field

Green's matrix in a magnetic field

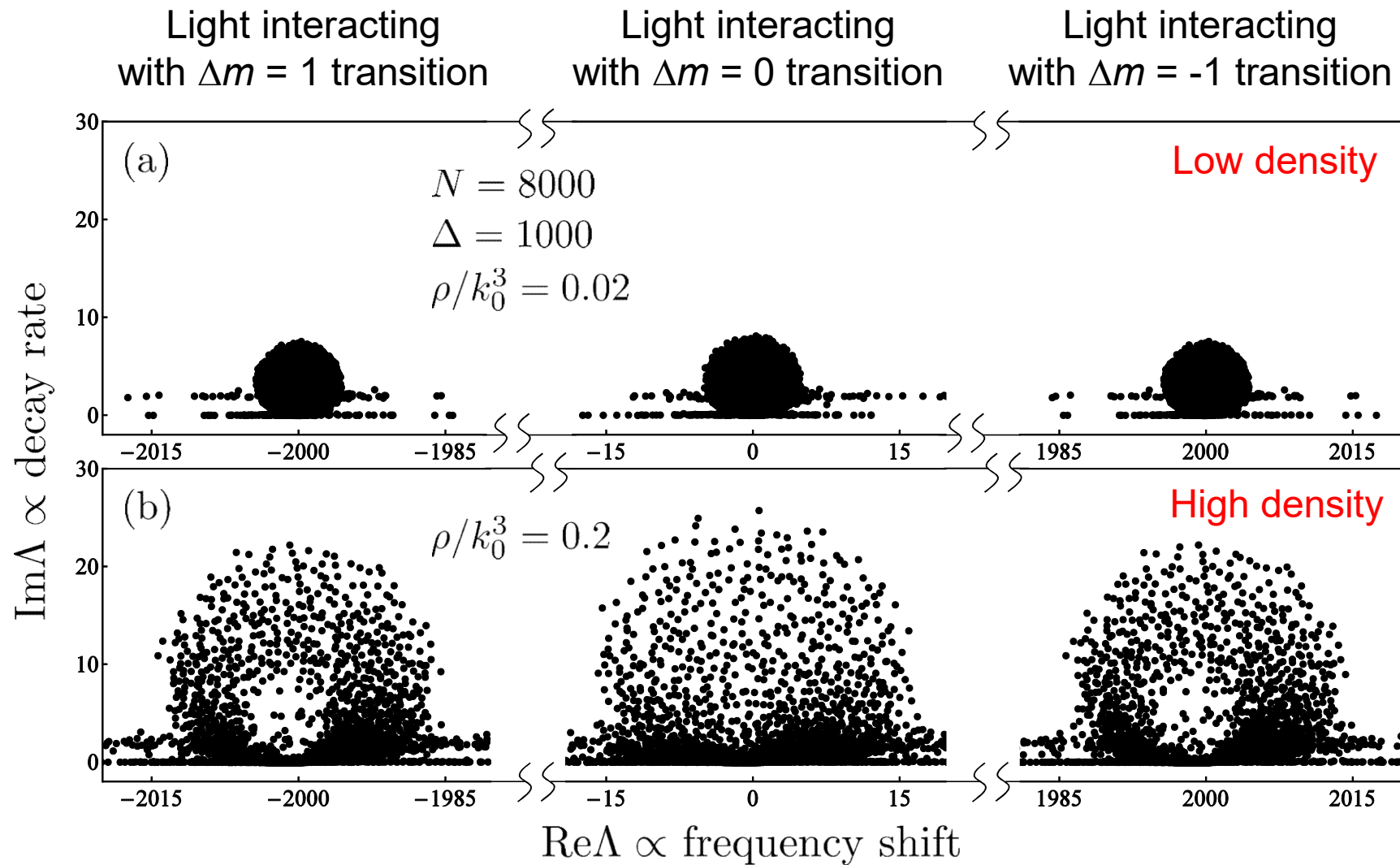
$$\begin{aligned}
 G_{e_{im}e_{jm'}} &= (i - 2m\Delta) \delta_{e_{im}e_{jm'}} - \frac{2}{\hbar\Gamma_0} (1 - \delta_{e_{im}e_{jm'}}) \\
 &\times \sum_{\mu,\nu} d_{e_{im}g_i}^\mu d_{g_j e_{jm'}}^\nu \frac{e^{ik_0 r_{ij}}}{r_{ij}^3} \\
 &\times \left\{ \delta_{\mu\nu} [1 - ik_0 r_{ij} - (k_0 r_{ij})^2] \right. \\
 &\quad \left. - \frac{r_{ij}^\mu r_{ij}^\nu}{r_{ij}^2} [3 - 3ik_0 r_{ij} - (k_0 r_{ij})^2] \right\}
 \end{aligned}$$


$$\Delta = g_e \mu_B B / \hbar \Gamma_0$$

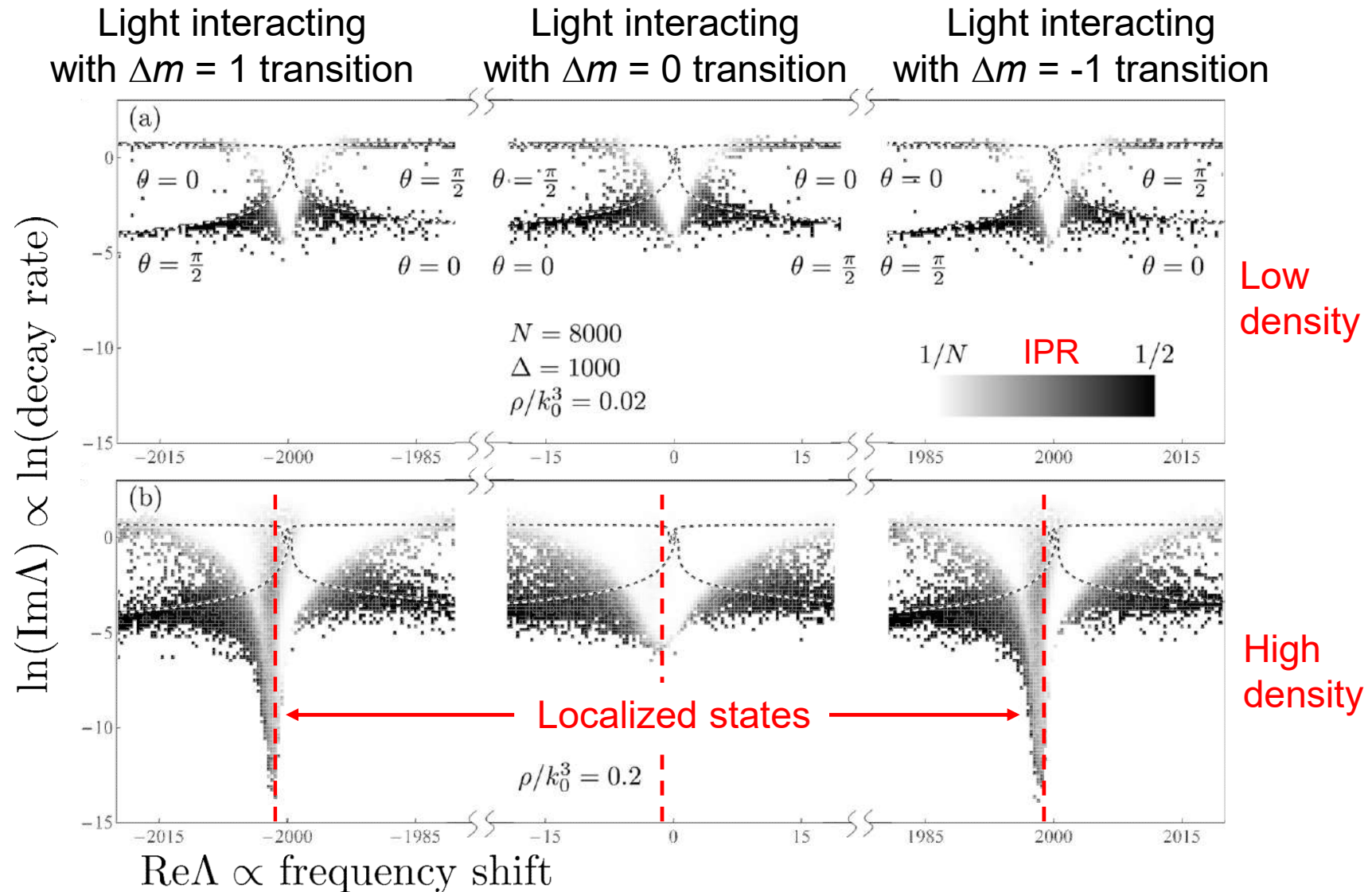
$$d_{e_{im}g_i} = \langle J_e m | \hat{\mathbf{D}}_i | J_g 0 \rangle$$

see also Pinheiro et al., Acta. Phys. Pol. A **105**, 339 (2004)

Eigenvalues in a strong magnetic field



Average inverse participation ratio

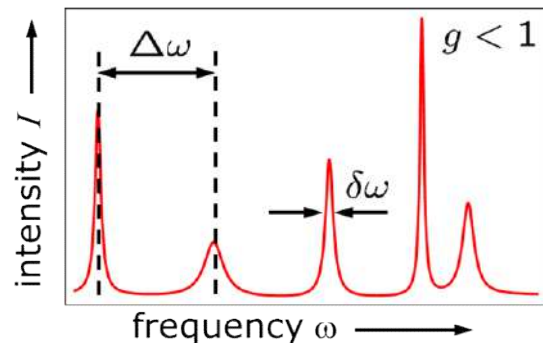
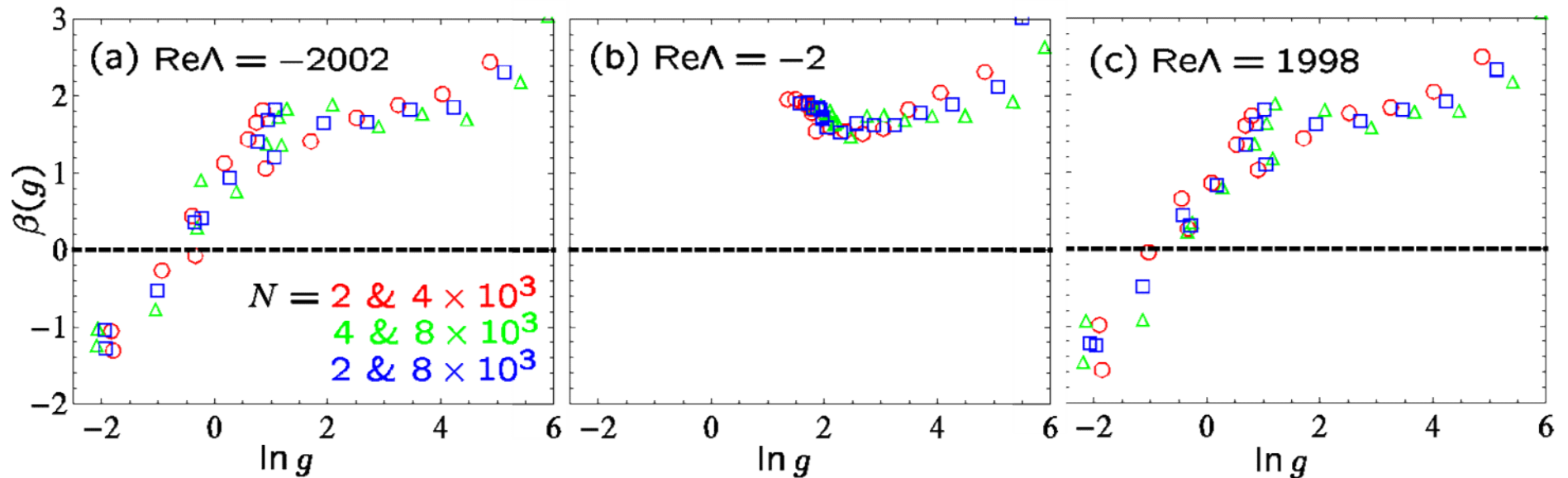


Scaling with sample size

Light interacting
with $\Delta m = 1$ transition

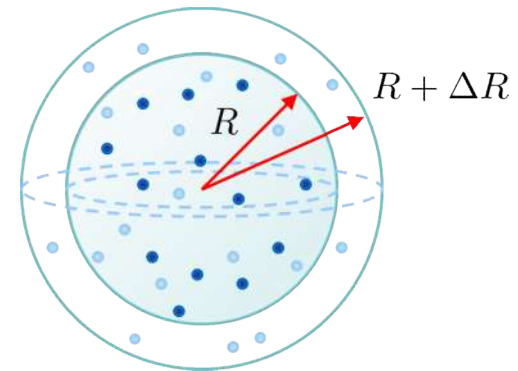
Light interacting
with $\Delta m = 0$ transition

Light interacting
with $\Delta m = -1$ transition



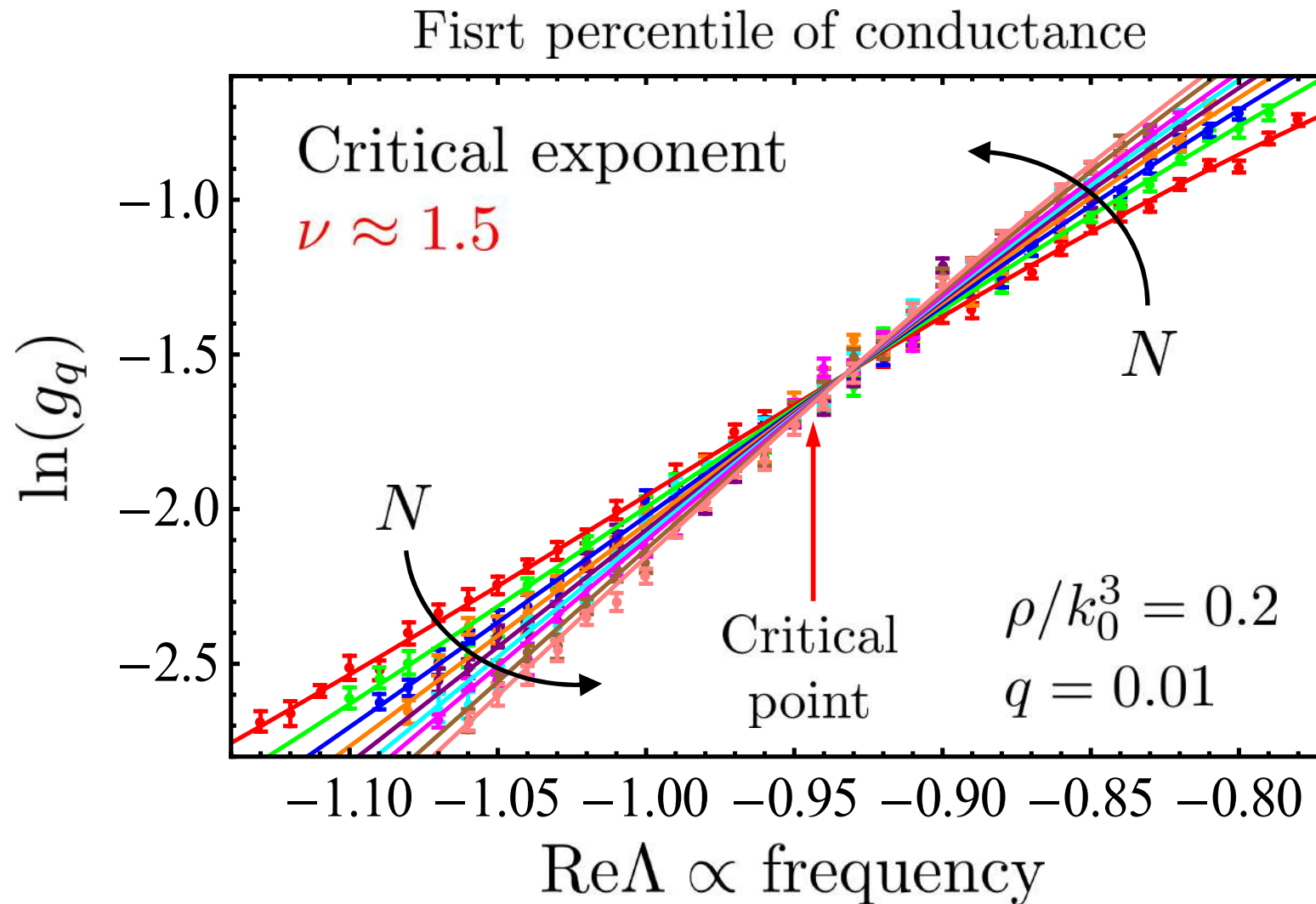
$$g = \frac{\delta\omega}{\Delta\omega}$$

$$\beta(g) = \frac{\partial \ln g}{\partial \ln k_0 R}$$



Skipetrov & Sokolov, PRL **114**, 053902 (2015)

Finite-size scaling of percentiles



Work in progress...

Anderson localization of elastic waves

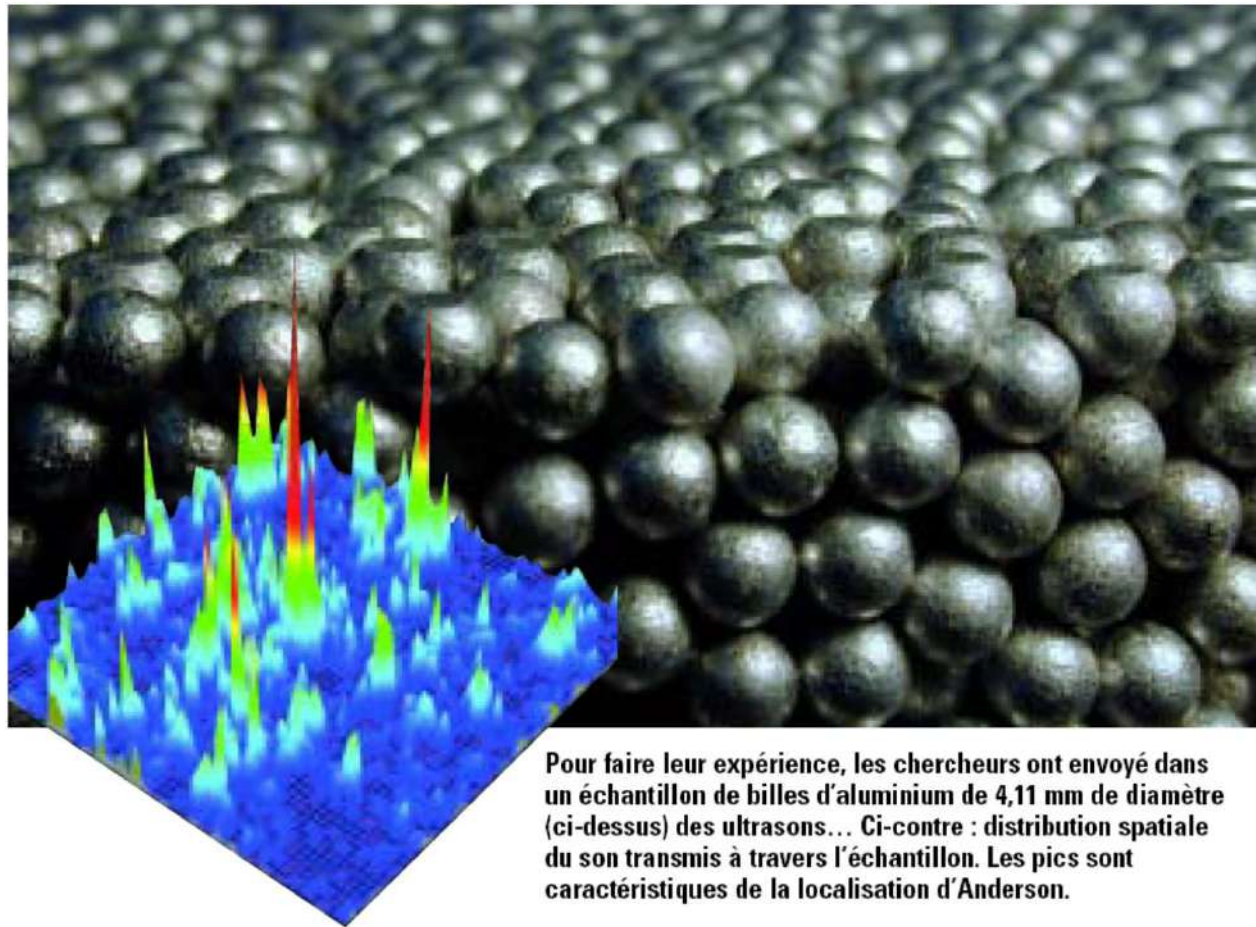


Image from *Le Journal du CNRS* (December 2008)

Elastic wave equation

$$\rho\omega^2 u_i + \frac{\partial}{\partial x_j} \left(c_{ijkl} \frac{\partial}{\partial x_k} u_l \right) = -f_i$$

mechanical
displacement
in the direction i

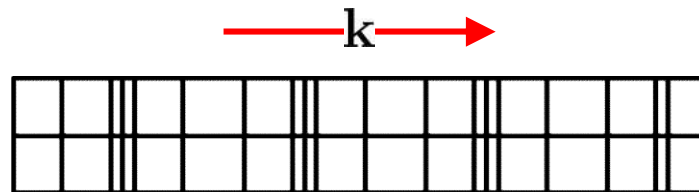
elasticity
tensor

excitation

$$c_{ijkl} = \lambda \delta_{ij} \delta_{kl} + \mu (\delta_{ik} \delta_{jl} + \delta_{il} \delta_{jk})$$

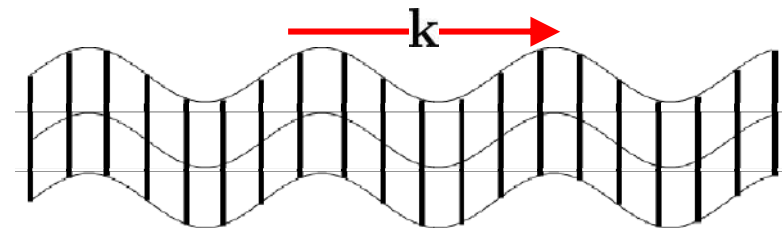
Lamé parameters

compressional waves:



velocity $\alpha = \sqrt{\frac{\lambda + 2\mu}{\rho}}$

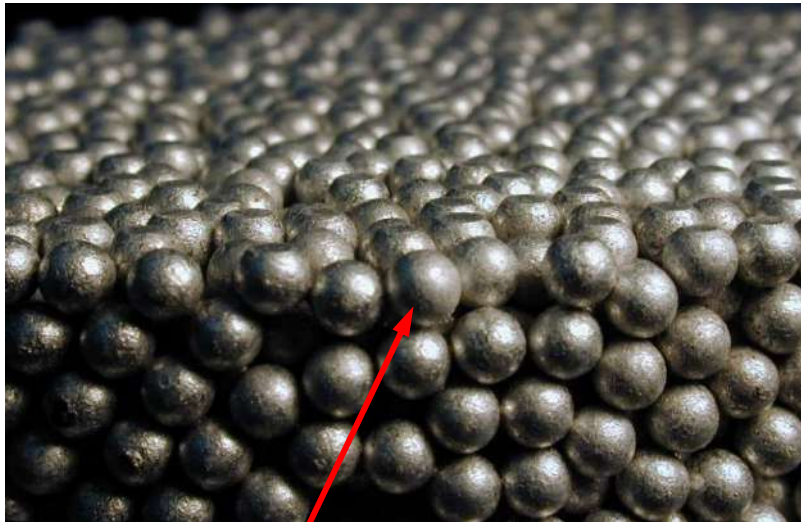
shear waves:



velocity $\beta = \sqrt{\frac{\mu}{\rho}}$

Point-scatterer model

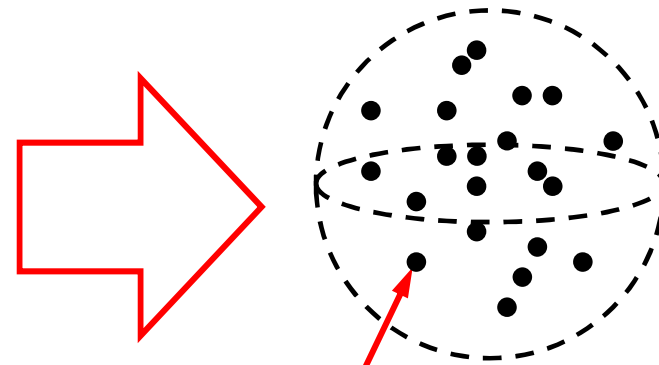
Real sample



Hu et al., Nature Physics **4**, 945 (2008)

Identical aluminum beads
with many resonances

Model



Identical point scatterers
with a single resonance

Elastic Green's function

$$\begin{aligned}\hat{G}(\mathbf{r}) = & \frac{k_\alpha}{4\pi(\lambda + 2\mu)} \times \left\{ \frac{e^{ik_\alpha r}}{3k_\alpha r} [\mathbb{1} + (\mathbb{1} - 3\hat{r} \otimes \hat{r})] \left(-1 - \frac{3i}{k_\alpha r} + \frac{3}{(k_\alpha r)^2} \right) \right. \\ & \left. - \left(\frac{\alpha}{\beta} \right)^3 \frac{e^{ik_\beta r}}{3k_\beta r} [-2\mathbb{1} + (\mathbb{1} - 3\hat{r} \otimes \hat{r})] \left(-1 - \frac{3i}{k_\beta r} + \frac{3}{(k_\beta r)^2} \right) \right\}\end{aligned}$$

$$k_\alpha = \frac{\omega}{\alpha}, \quad k_\beta = \frac{\omega}{\beta}, \quad \hat{r} = \frac{\mathbf{r}}{r}$$

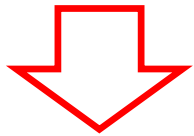
Typically, $\alpha > \beta$ ($\alpha/\beta \simeq 2$ for aluminium)

Equipartition principle:

$$\frac{\langle \text{Energy of shear waves} \rangle}{\langle \text{Energy of compressional waves} \rangle} = 2 \left(\frac{\alpha}{\beta} \right)^3 > 1$$

Elastic Green's function in the near field

$$\begin{aligned}\hat{G}(\mathbf{r}) = & \frac{k_\alpha}{4\pi(\lambda + 2\mu)} \times \left\{ \frac{e^{ik_\alpha r}}{3k_\alpha r} [\mathbb{1} + (\mathbb{1} - 3\hat{\mathbf{r}} \otimes \hat{\mathbf{r}})] \left(-1 - \frac{3i}{k_\alpha r} + \frac{3}{(k_\alpha r)^2} \right) \right. \\ & \left. - \left(\frac{\alpha}{\beta} \right)^3 \frac{e^{ik_\beta r}}{3k_\beta r} [-2\mathbb{1} + (\mathbb{1} - 3\hat{\mathbf{r}} \otimes \hat{\mathbf{r}})] \left(-1 - \frac{3i}{k_\beta r} + \frac{3}{(k_\beta r)^2} \right) \right\}\end{aligned}$$



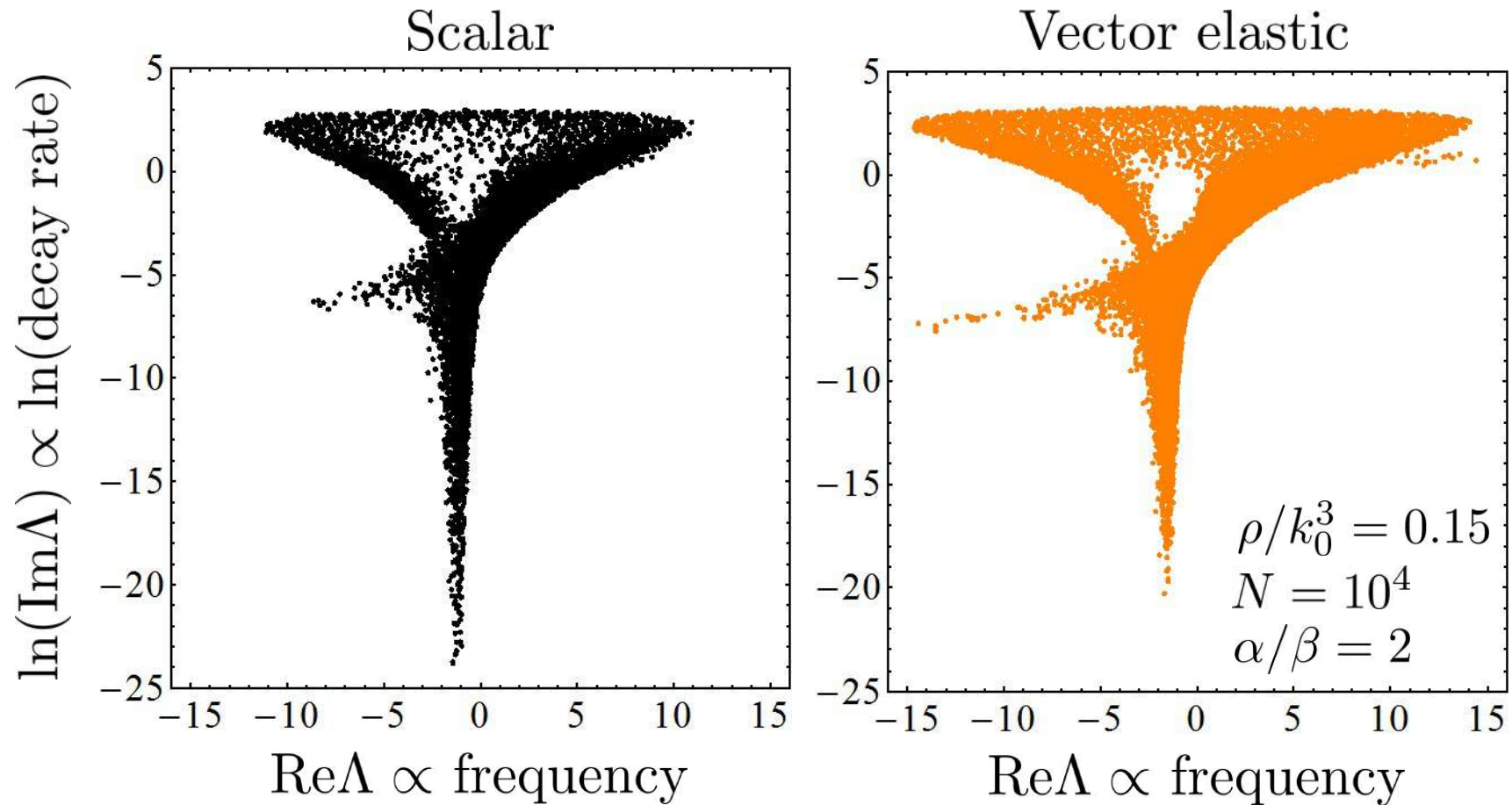
Near-field behavior:

$$\hat{G}(\mathbf{r})|_{r \rightarrow 0} = -\frac{1}{8\pi\mu} \left(1 - \frac{\beta^2}{\alpha^2} \right) \frac{\mathbb{1} - 3\hat{\mathbf{r}} \otimes \hat{\mathbf{r}}}{r} \propto \frac{1}{r}$$

Similar to the scalar case and different from the electromagnetic one:

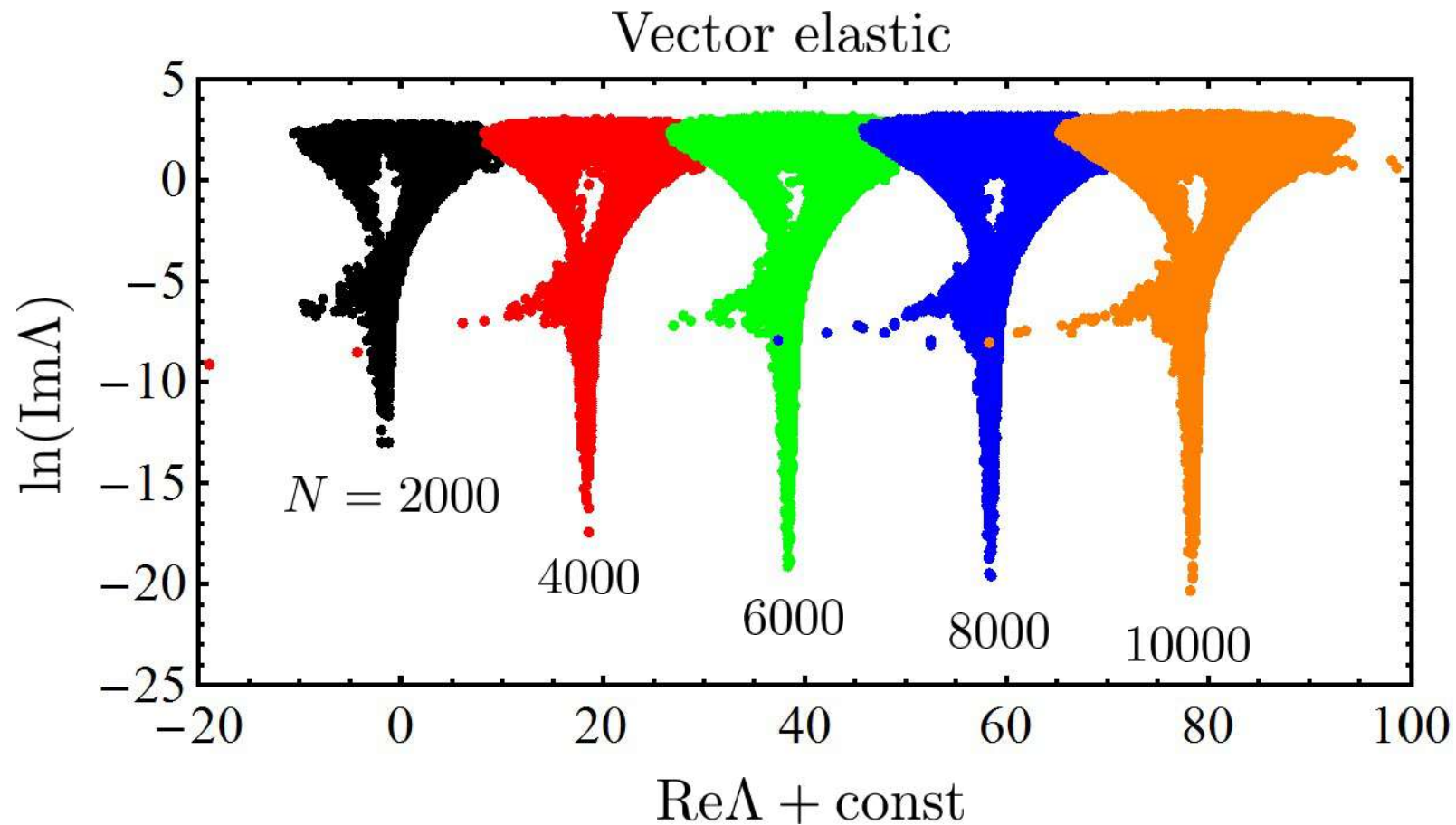
$$\hat{G}_{\text{EM}}(\mathbf{r})|_{r \rightarrow 0} \propto \frac{1}{r^3}$$

Eigenvalues: scalar vs elastic



Work in progress...

Eigenvalues: evolution with sample size



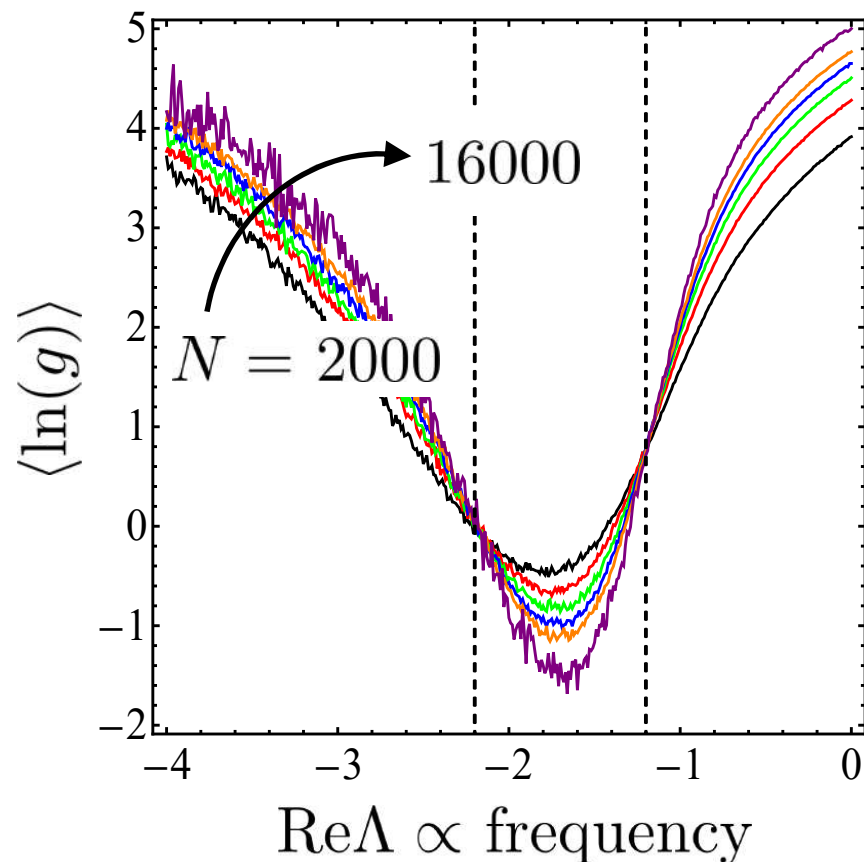
$$\rho/k_0^3 = 0.15$$

$$\alpha/\beta = 2$$

Work in progress...

Scaling for elastic waves

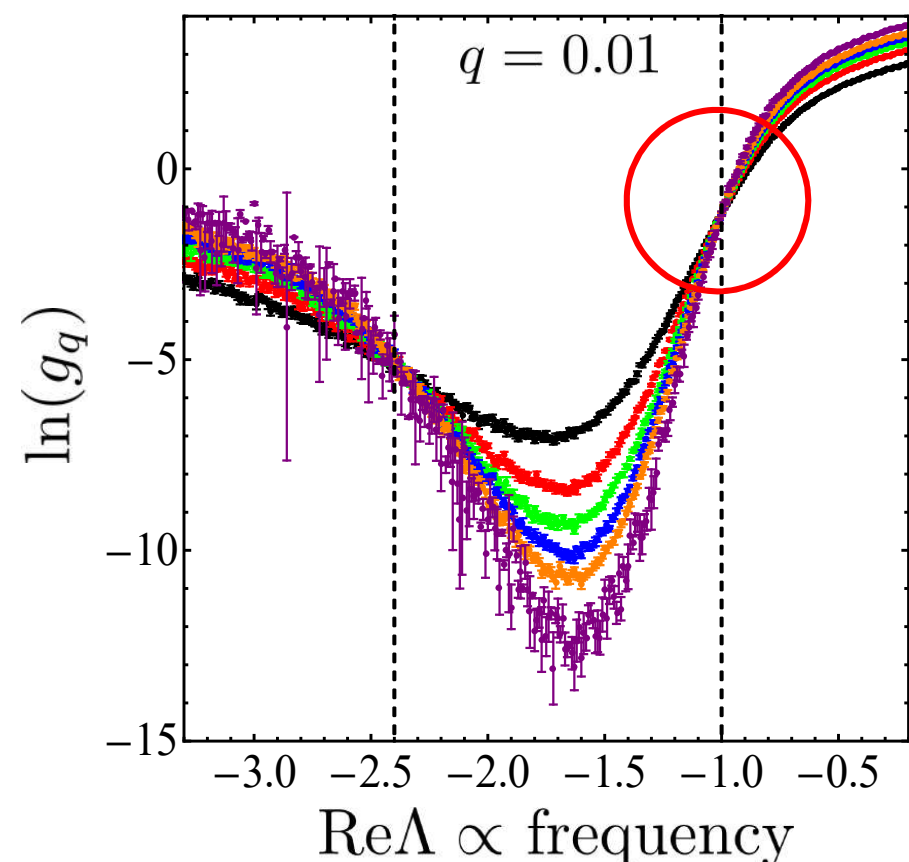
Average ln of conductance



$$\rho/k_0^3 = 0.15$$

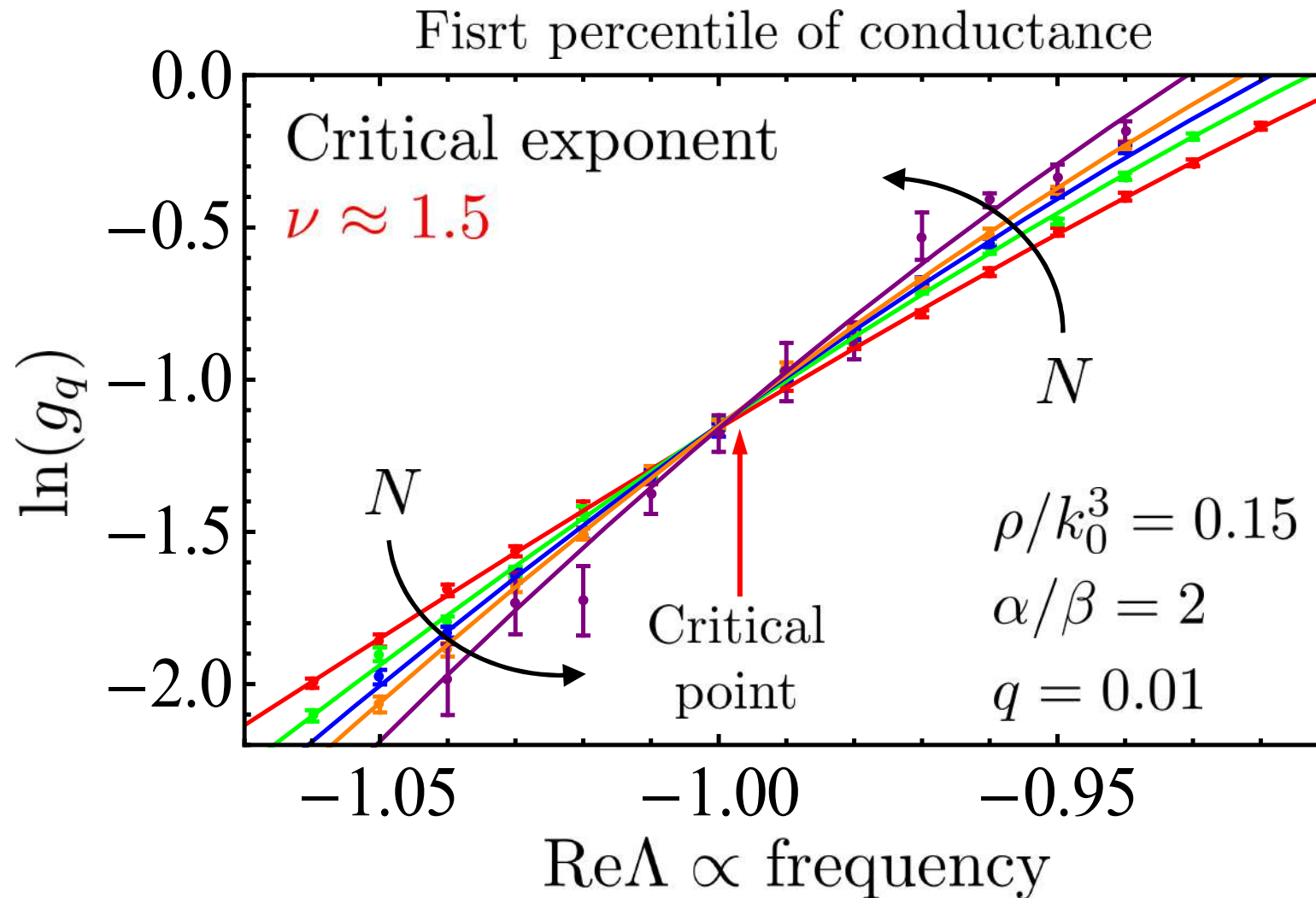
$$\alpha/\beta = 2$$

First percentile of conductance



Work in progress...

Finite-size scaling of percentiles



Work in progress...

Conclusions

- Wave scattering by random ensembles of resonant point scatterers is naturally described by **Euclidean random matrix (ERM) models**
- ERM models can be used to study **Anderson localization transitions** with account for peculiarities specific for a particular type of waves: scalar vs vector waves, light vs elastic waves, etc.
- **Anderson localization of light by atoms** is possible only in the presence of a strong magnetic field and that the localization transition for elastic waves is similar to that of scalar waves
- More developments of **analytic approaches to ERM** are necessary for further progress

Thank you for your attention!