

Cooperative effects and photon localization in atomic gases :

Phase transition in non Hermitian random matrices

PRL **101**, 103602 (2008),
EPL **101**, (2013)
PR **A88**, (2013),
PR A **90**, 063822 (2014)

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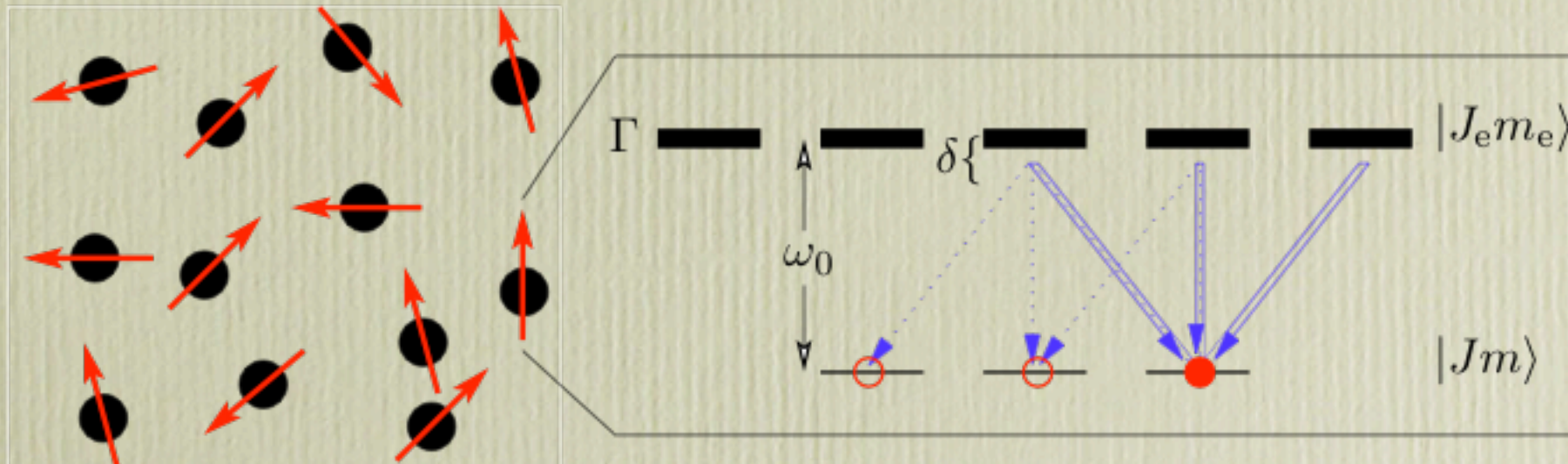
Cargese, September 28, 2016

What is it about ?

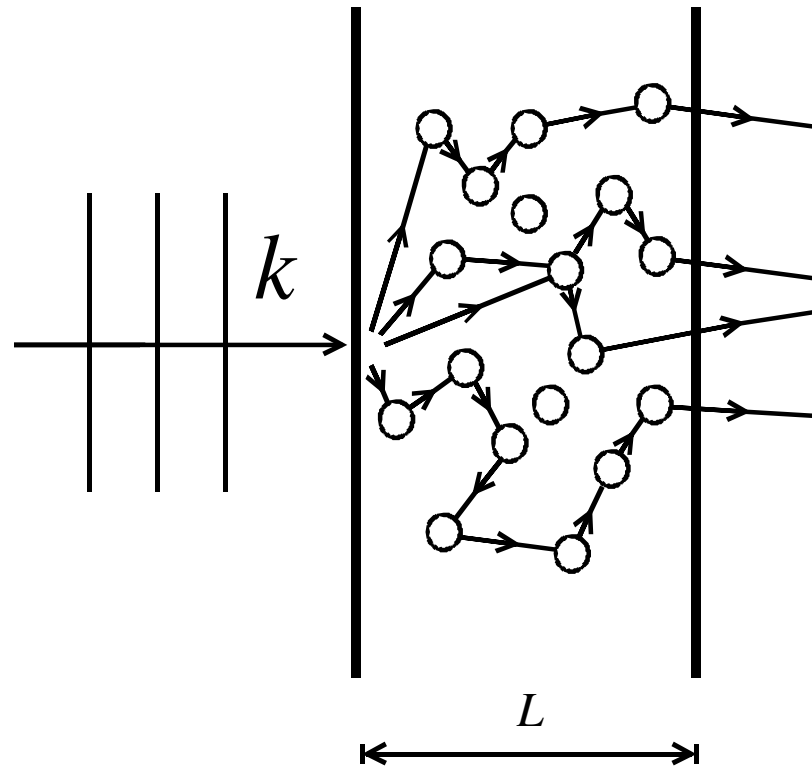
- Coherent multiple scattering of photons/waves
- Anderson photon localization : phase transition and scaling
- Cooperative effects and Dicke superradiance
- Photon escape rates : Competition between Anderson and Dicke mechanisms

Framework

Multiple scattering of photons/waves by a cold atomic gas.



Multiple scattering



2 characteristic lengths:

Wavelength: λ_0

Elastic mean free path: $l = \frac{1}{n_i \sigma} \gg n_i^{-1/d}$

density of scatterers

scattering cross section $\sigma \propto \lambda^2$

Disorder strength :

$$W = \frac{1}{k_0 l} = \frac{\pi}{2} \frac{\lambda}{L} \frac{N}{N_{\perp}}$$

$$\lambda = 2\pi / k_0$$

Elastic mean free path

$$l = \frac{1}{n\sigma} = \frac{L^3}{N\lambda^2}$$

Number of transverse channels $(d = 3)$

$$N_{\perp} = (k_0 L)^2 / 4$$

Weak disorder limit $W \ll 1$

Numerical calculations on the Anderson Hamiltonian

$d = 2$

$d = 3$

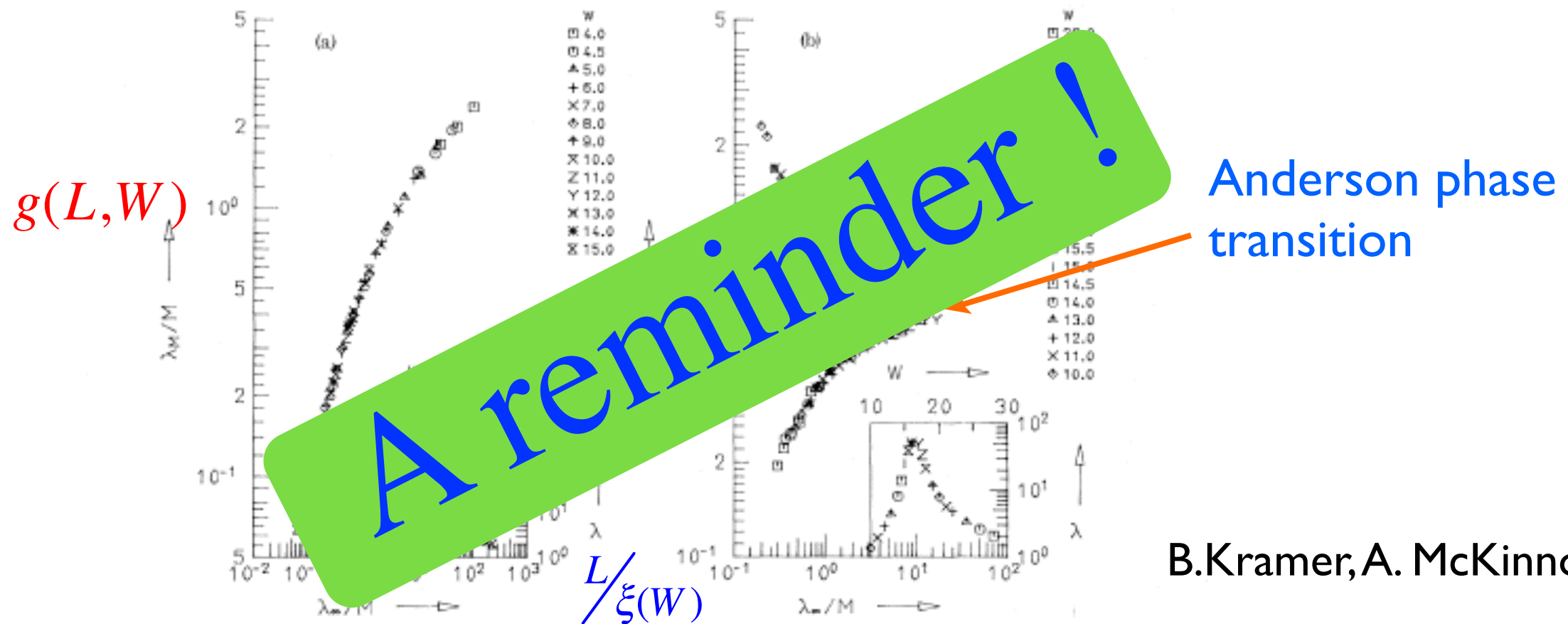


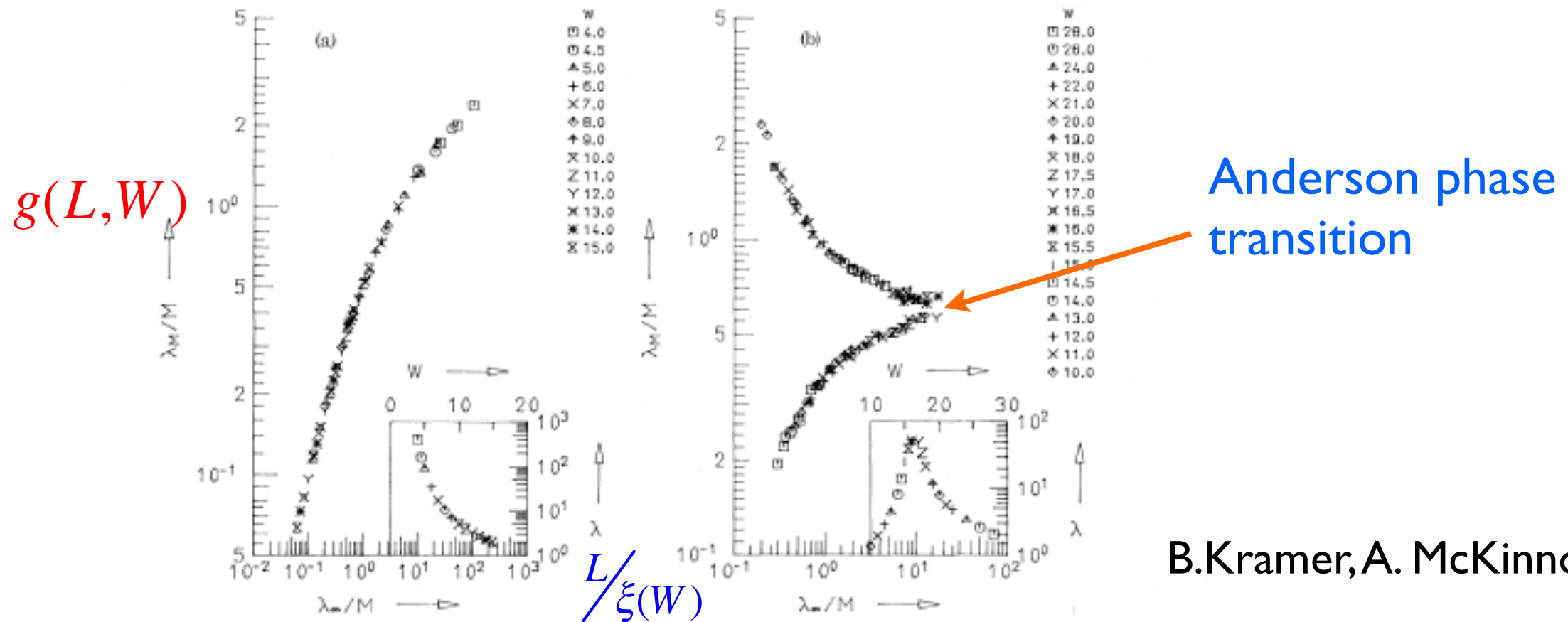
FIG. 1. Scaling function λ_m/M vs λ_m/M for the localization length λ_m of a system of thickness M for (a) $d=2$ ($M \geq 4$) and (b) $d=3$ ($M \geq 3$). Insets show the scaling parameter λ_m as a function of the disorder W .

Anderson localisation phase transition occurs in $d > 2$

Numerical calculations on the Anderson Hamiltonian

$d = 2$

$d = 3$



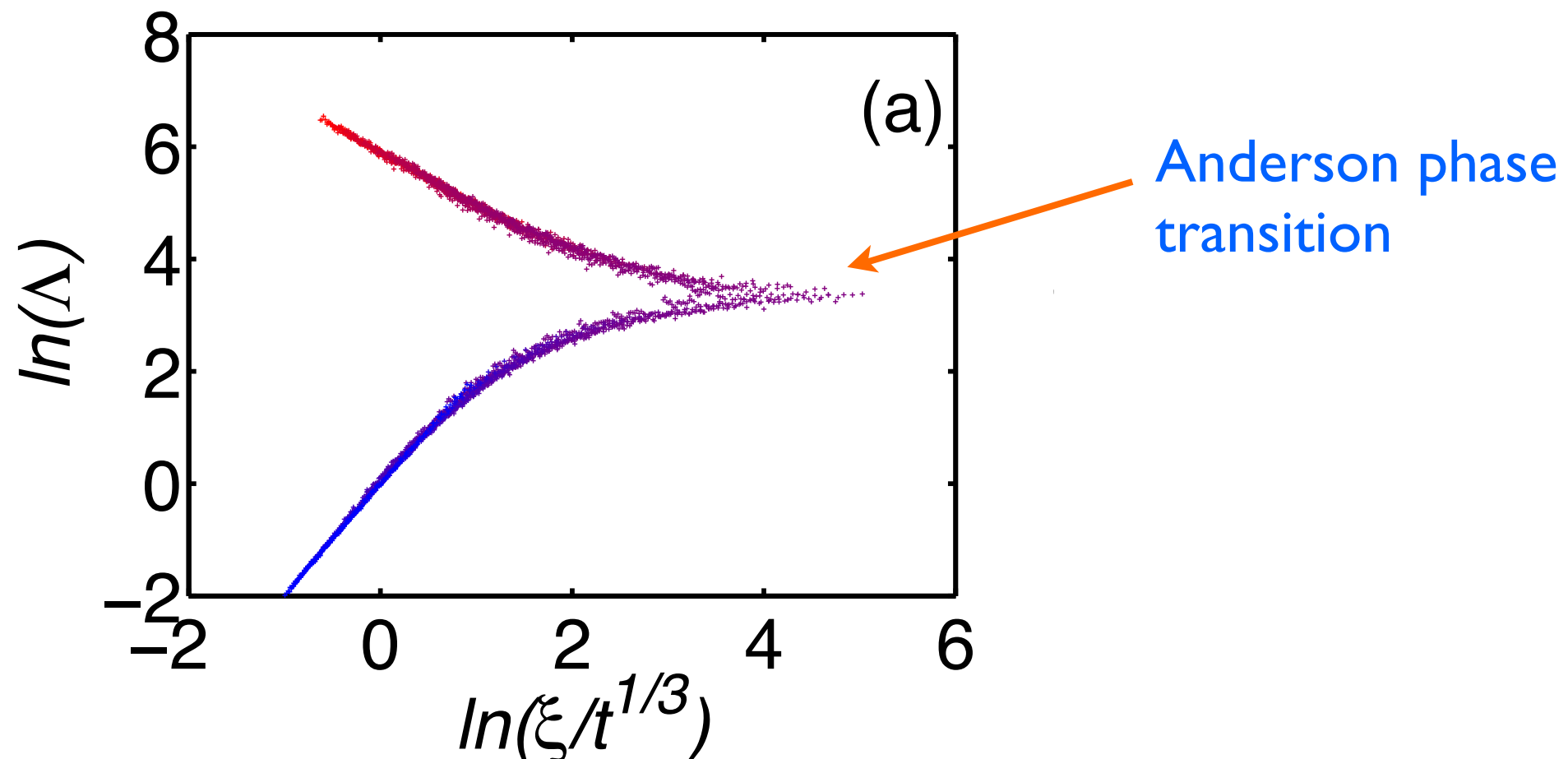
B.Kramer, A. McKinnon, 1981

FIG. 1. Scaling function λ_M/M vs λ_w/M for the localization length λ_M of a system of thickness M for (a) $d=2$ ($M \geq 4$) and (b) $d=3$ ($M \geq 3$). Insets show the scaling parameter λ_w as a function of the disorder W .

Anderson localisation phase transition occurs in $d > 2$

Realisations of the Anderson Hamiltonian

Quantum evolution of the atomic kicked rotor (localisation of the momentum in phase space (d=3))

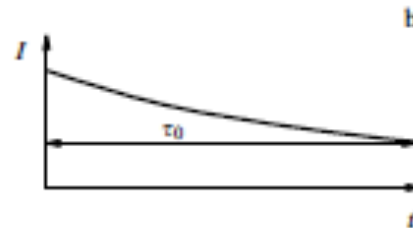
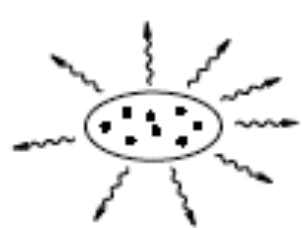


(P. Szriftgiser et al. 2010, for the experiment, theory : Casati, Chirikov, ('79)
Fishman, Grempel, Prange, ('84), Guarneri et al. ('89),

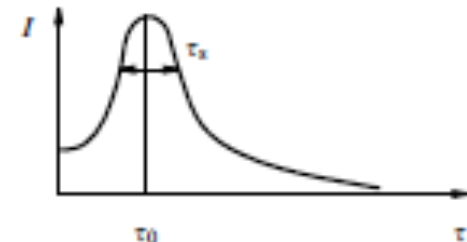
Cooperative effects (superradiance-subradiance)

Cooperative spontaneous radiation (**Superradiance**) results from *quantum phase correlations* induced between atoms by dipole-dipole interactions.

Superradiant emission can be summarised by



$$I \approx N$$



$$I \approx N^2$$

But the dependence $I \approx N^2$ *does not constitute the main distinguishing feature of superradiance.*

It is rather the mechanism leading to *coherent phasing of atoms.*

Superradiant emission : all atoms must see (in phase) the same
electromagnetic field.

- small volumes (*Dicke limit*)
- large systems : *Anderson localization* may play a role :

photon modes are spatially localized in volumes ξ^d

only a fraction $N \left(\frac{\xi}{L} \right)^d$ of atoms are coherent so that the

emission time τ_s becomes large:

$$\tau_s \approx \left(\frac{L}{\xi} \right)^d \frac{\ln N}{N} \gg \frac{\ln N}{N}$$

Superradiant emission : all atoms must see (in phase) the same electromagnetic field.

→ small volumes (*Dicke limit*)

→ large systems : *Anderson localization* plays a role :

photon modes are scattered in large volumes ξ^d

only a fraction of atoms are coherent so that the

emission becomes large:

$$\tau_s \approx \left(\frac{L}{\xi} \right)^d \frac{\ln N}{N} \gg \frac{\ln N}{N}$$

Model

N identical two-level atoms located at random positions \vec{r}_i (uniform distribution) with electric dipole moments \vec{d}_i in the quantum radiation field \vec{E}

- Total Hamiltonian

$$H = H_0 + U$$

- Non-interacting Hamiltonian

$$H_0 = \hbar\omega_0 \sum_{i=1}^N |e_i\rangle\langle e_i| + \sum_{\vec{k}\epsilon} \hbar\omega_k a_{\vec{k}\epsilon}^+ a_{\vec{k}\epsilon}$$

- Electric dipole representation of the interaction

$$U = - \sum_{i=1}^N \vec{d}_i \cdot \vec{E}(\vec{r}_i)$$

Model

- Effective Hamiltonian
 - Tracing over the EM field degrees of freedom

$$H_e = \left(\hbar\omega_0 - i\frac{\hbar\Gamma_0}{2} \right) \sum_{i=1}^N |e_i\rangle\langle e_i| + \frac{\hbar\Gamma_0}{2} \sum_{i \neq j} V_{ij} \Delta_i^+ \Delta_j^-$$

- Atomic raising and lowering operators

$$\Delta_i^+ = |e_i\rangle\langle g_i| \quad \Delta_j^- = |g_j\rangle\langle e_j|$$

Model

- Effective Hamiltonian
 - Tracing over the EM field degrees of freedom

$$H_e = \left(\hbar\omega_0 - i\frac{\hbar\Gamma_0}{2} \right) \sum_{i=1}^N |e_i\rangle\langle e_i| + \frac{\hbar\Gamma_0}{2} \sum_{i \neq j} V_{ij} \Delta_i^+ \Delta_j^-$$

- Atomic raising and lowering operators

$V_{ij} = \beta_{ij} - i\gamma_{ij}$ is random and complex valued

- Real part : interaction potential

$$\beta_{ij} = \frac{3}{2} \left[-p \frac{\cos k_0 r_{ij}}{k_0 r_{ij}} + q \left(\frac{\cos k_0 r_{ij}}{(k_0 r_{ij})^3} + \frac{\sin k_0 r_{ij}}{(k_0 r_{ij})^2} \right) \right]$$

- Imaginary part : **photon escape rate**

$$\gamma_{ij} = \frac{3}{2} \left[p \frac{\sin k_0 r_{ij}}{k_0 r_{ij}} - q \left(\frac{\sin k_0 r_{ij}}{(k_0 r_{ij})^3} - \frac{\cos k_0 r_{ij}}{(k_0 r_{ij})^2} \right) \right]$$

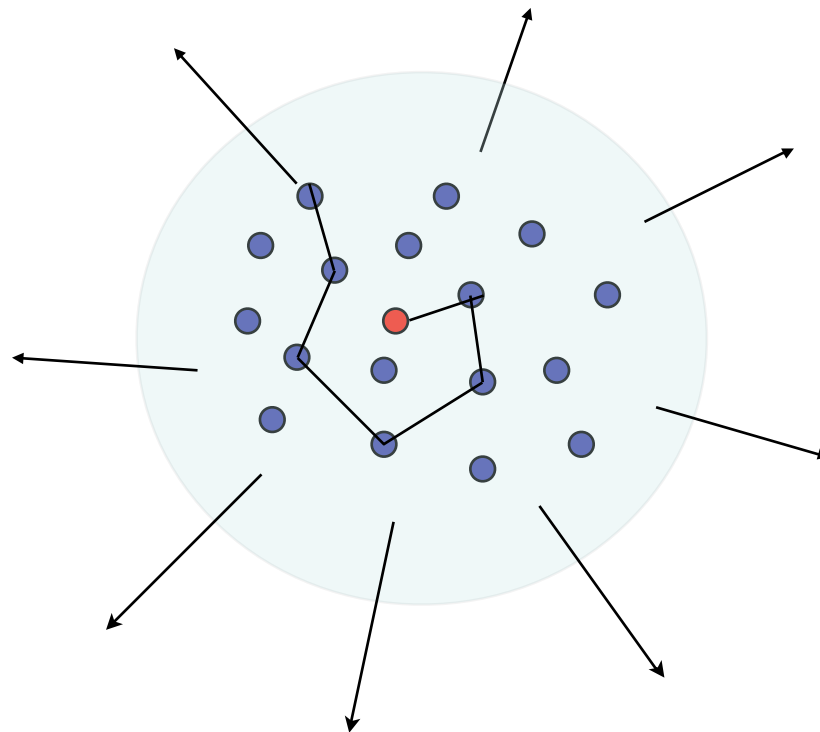
- **For a scalar wave:**

$$\beta_{ij} = -\frac{\cos(k_0 r_{ij})}{k_0 r_{ij}}$$

$$\gamma_{ij} = \frac{\sin k_0 r_{ij}}{k_0 r_{ij}}$$

Which quantity to study ?

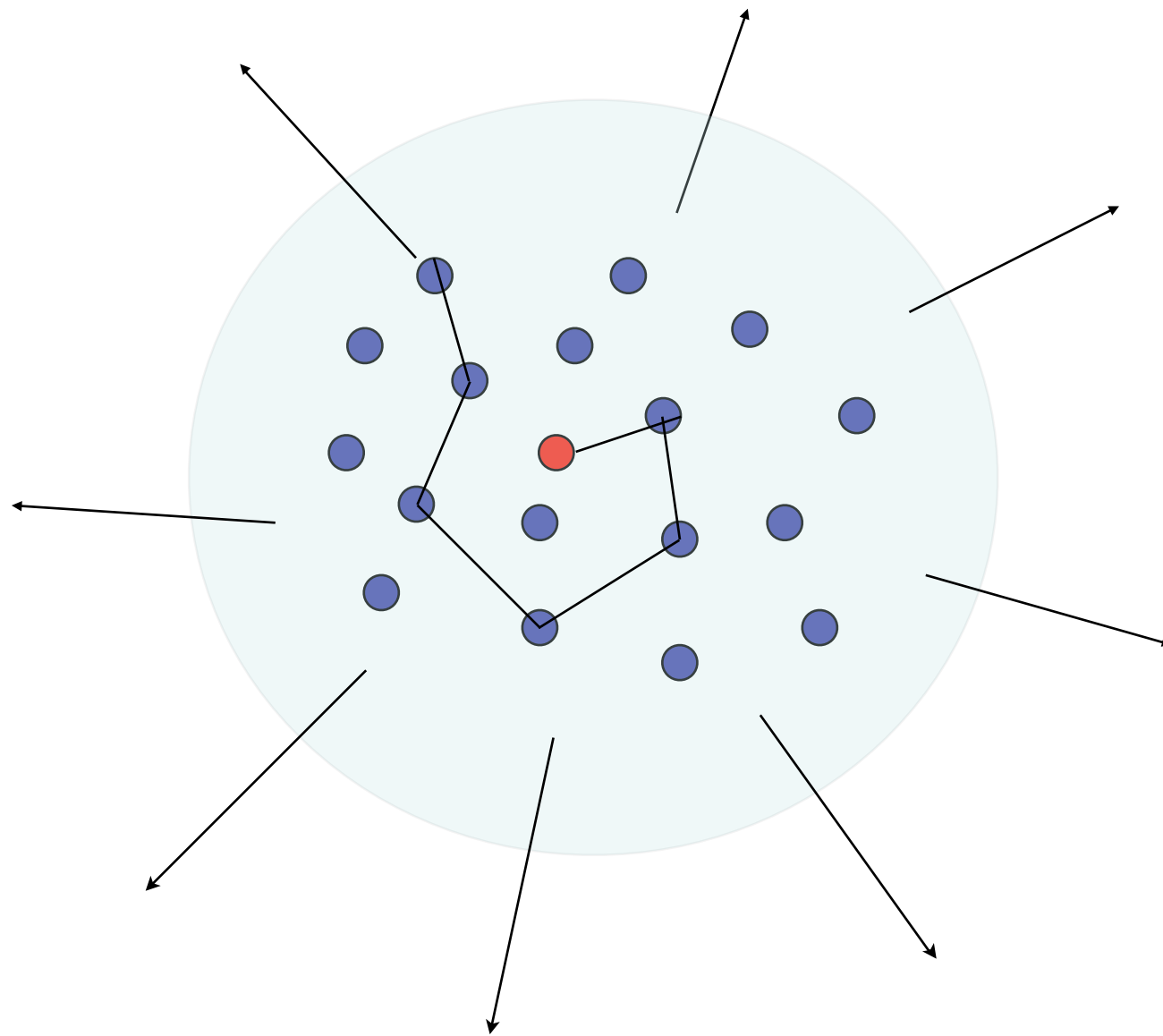
- The radiation pattern/intensity of the atomic cloud with a single excited atom



$$\Psi = \sum_{j=1}^N \beta_j(t) |b_1 b_2 \cdots a_j \cdots b_N\rangle |0\rangle + \sum_{\mathbf{k}} \gamma_{\mathbf{k}}(t) |b_1 b_2 \cdots b_N\rangle |1_{\mathbf{k}}\rangle.$$

Photon escape rates are a measure of localization and/or cooperative emission.

Escape rates are not a transport quantity.



More precisely : Photon escape rates

Evolution of the density matrix (Linblad form)

$$\frac{d\rho}{dt} = \frac{1}{i\hbar} \left(H_e \rho - \rho H_e^\dagger \right) + \Gamma_0 \sum_{i \neq j} \gamma_{ij} \Delta_i^+ \rho \Delta_j^-$$

$$H_e = \left(\hbar\omega_0 - i\frac{\hbar\Gamma_0}{2} \right) \sum_{i=1}^N |e_i\rangle\langle e_i| + \frac{\hbar\Gamma_0}{2} \sum_{i \neq j} V_{ij} \Delta_i^+ \Delta_j^-$$

$$V_{ij} = \beta_{ij} - i\gamma_{ij}$$

Photon escape rates from the atomic gas are
obtained from the eigenvalues of the
euclidean random matrix γ_{ij}

Eigenvalue density $P(\Gamma)$ of the $N \times N$ random matrix γ_{ij}

(Scalar case)

$$\gamma_{ij} = \frac{\sin k_0 r_{ij}}{k_0 r_{ij}} \quad x_{ij} \equiv k_0 r_{ij}$$

Defining dimensionless quantities $L^d = (\lambda a)^d$ $\lambda = 2\pi/k_0$

$$W = \frac{1}{k_0 l} = \frac{\pi}{2} \frac{\lambda}{L} \frac{N}{N_{\perp}}$$

Elastic mean free path

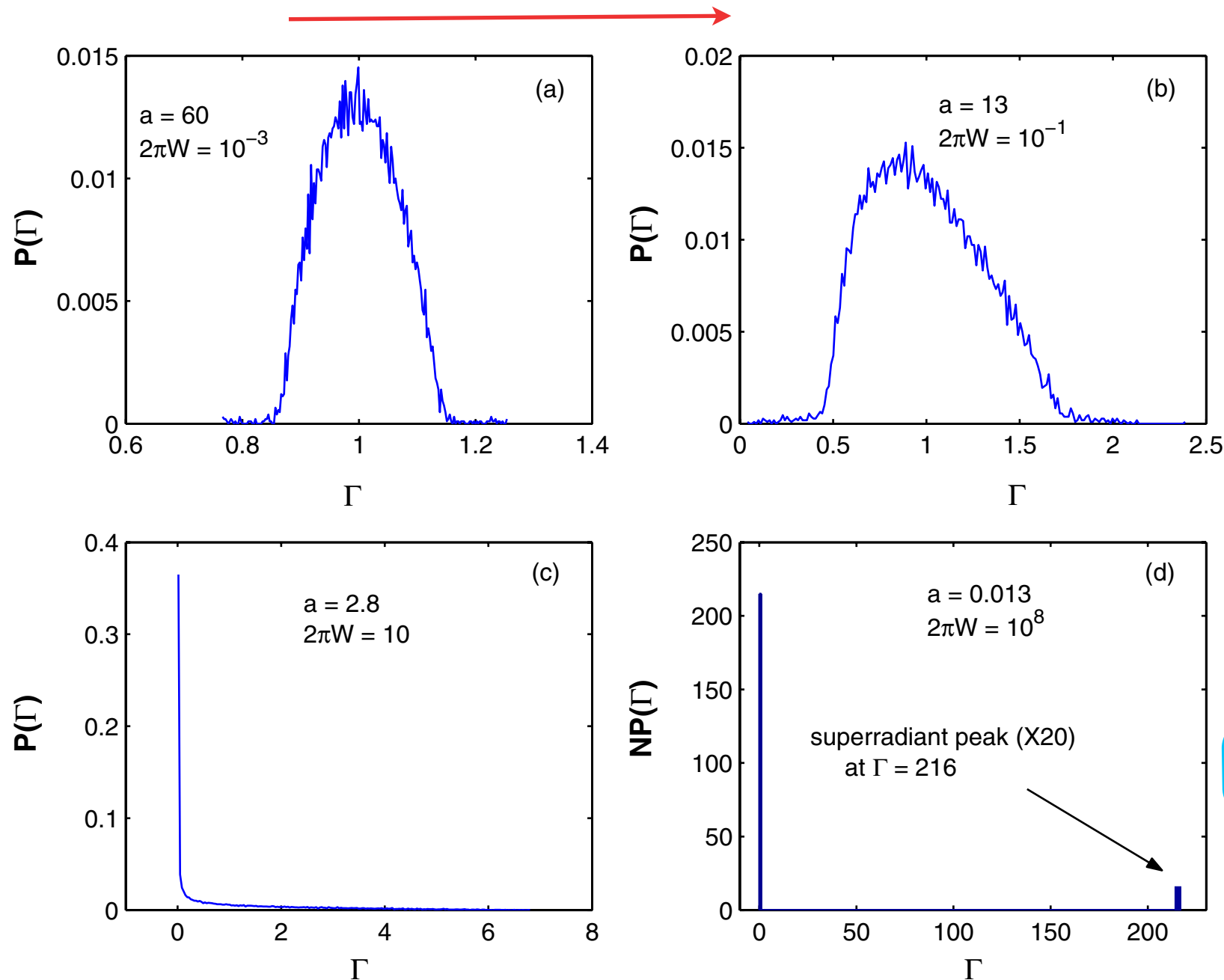
$$l = \frac{1}{n\sigma} = \frac{L^3}{N\lambda^2}$$

Number of transverse channels $(d = 3)$

$$N_{\perp} = (k_0 L)^2 / 4$$

Eigenvalue density $P(\Gamma)$

increasing disorder W



Dicke limit

localized photons

Numerical results $N=216$

Scaling ?

To characterize $P(\Gamma)$ we look for a scaling function $C(a,W)$

Relative number of localized states i.e. having a vanishing escape rate :

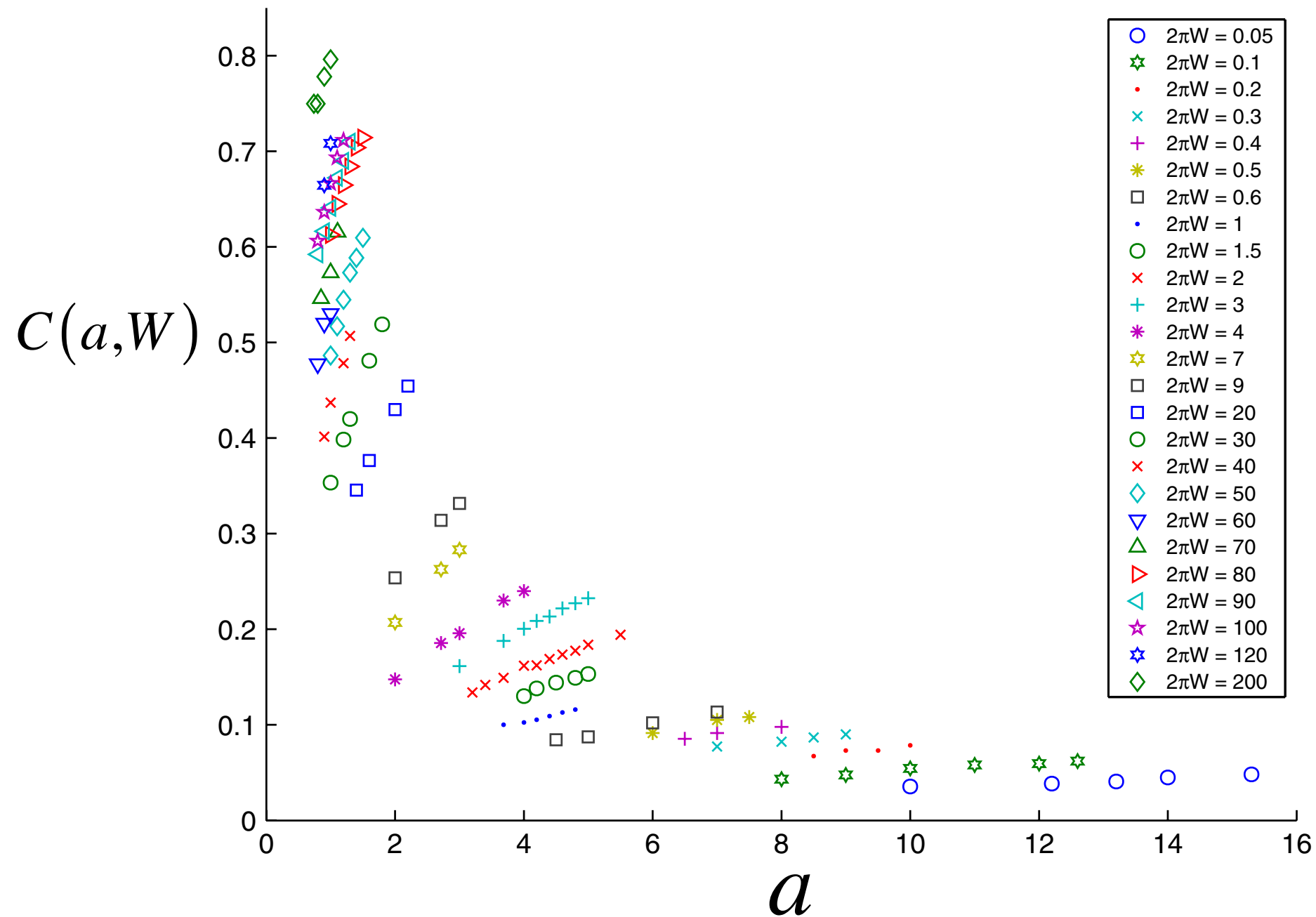
$$C(a,W) = 1 - 2 \int_1^{\infty} d\Gamma P(\Gamma)$$

$C(a,W)$ is defined between 0 and 1. At finite size, we expect the **scaling form**:

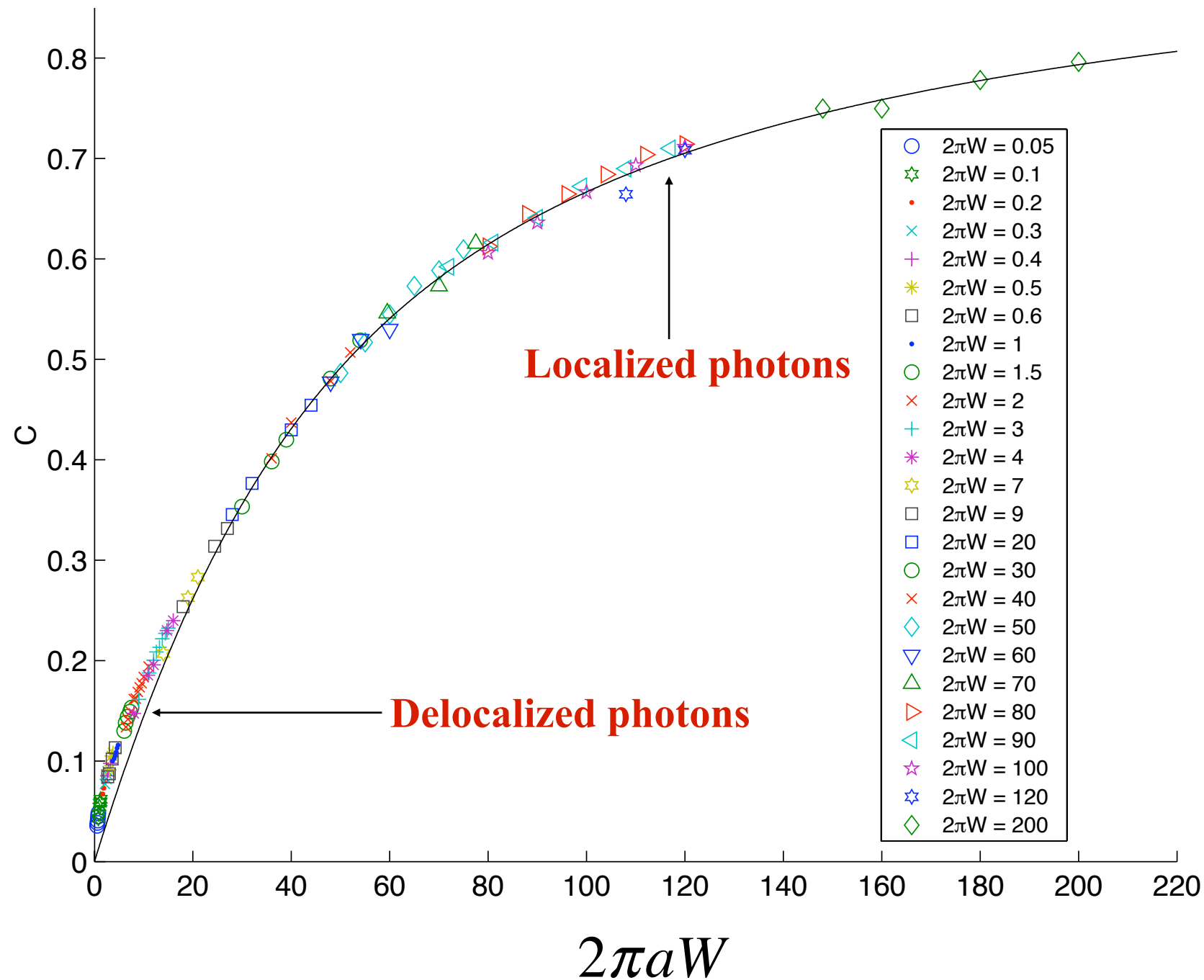
$$C(a,W) = f\left(a/\xi(W)\right)$$

Scaling behaviour

Large sample limit ($a \geq 1$)



Scaling behavior (large sample limit)



Is there a
localisation
phase
transition ?

$C(a, W)$ depends on $2\pi aW = \pi^2 N / N_{\perp}$

Is there a localisation phase transition ?

- Microscopic QED approach

Large disorder limit $N \gg N_{\perp}$

- Phenomenological Markov process
(Small world networks)

For the whole range of disorder

Microscopic QED approach

Large disorder limit $N \gg N_{\perp}$

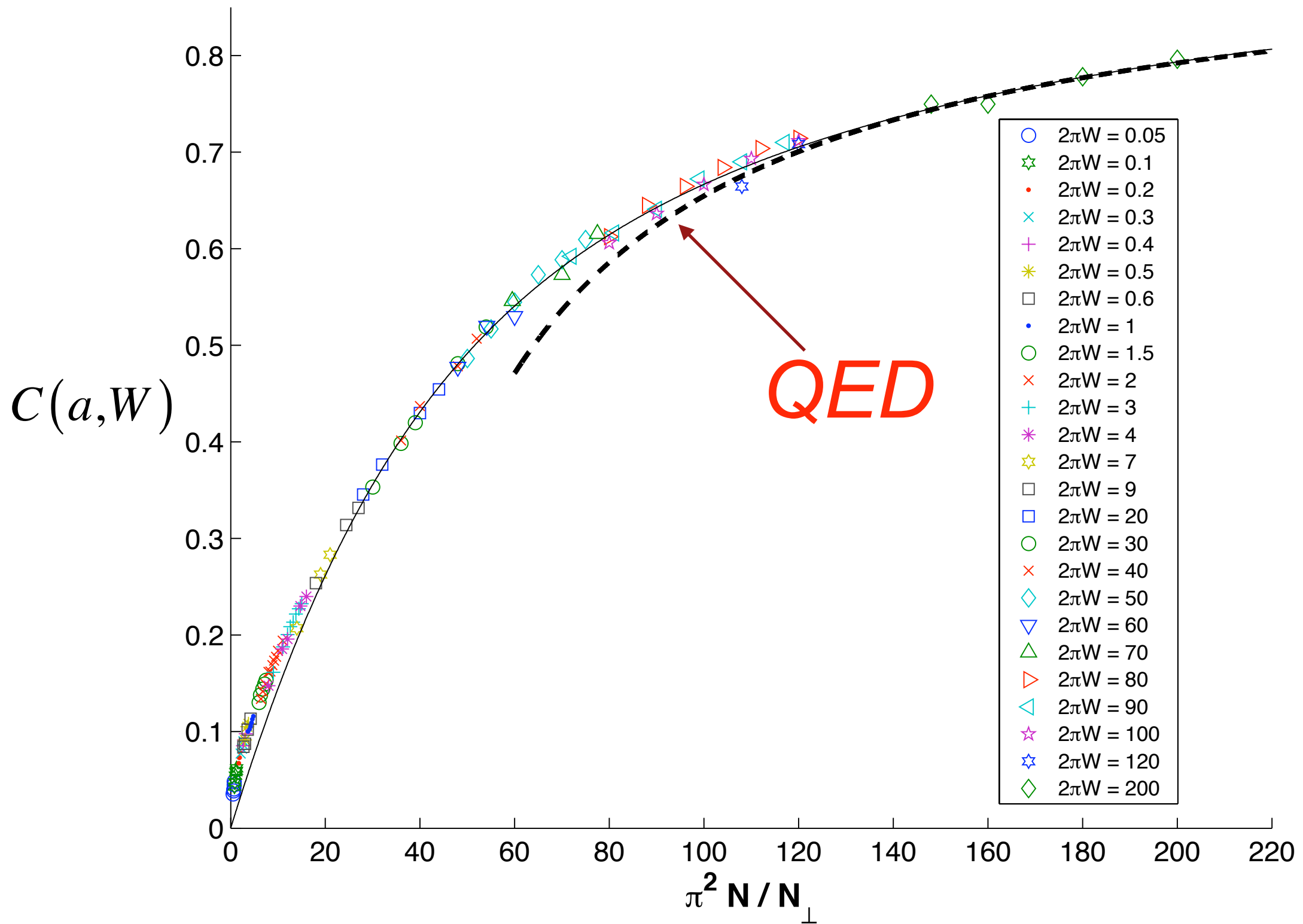
Resummation of the cumulants of $P(\Gamma)$ leads to the asymptotic behavior

$$P(\Gamma) = \left(1 - \frac{3N_{\perp}}{2N}\right) \delta(\Gamma) + 3\Gamma \left(\frac{N_{\perp}}{N}\right)^3 \quad \text{for } \Gamma \leq N/N_{\perp}$$

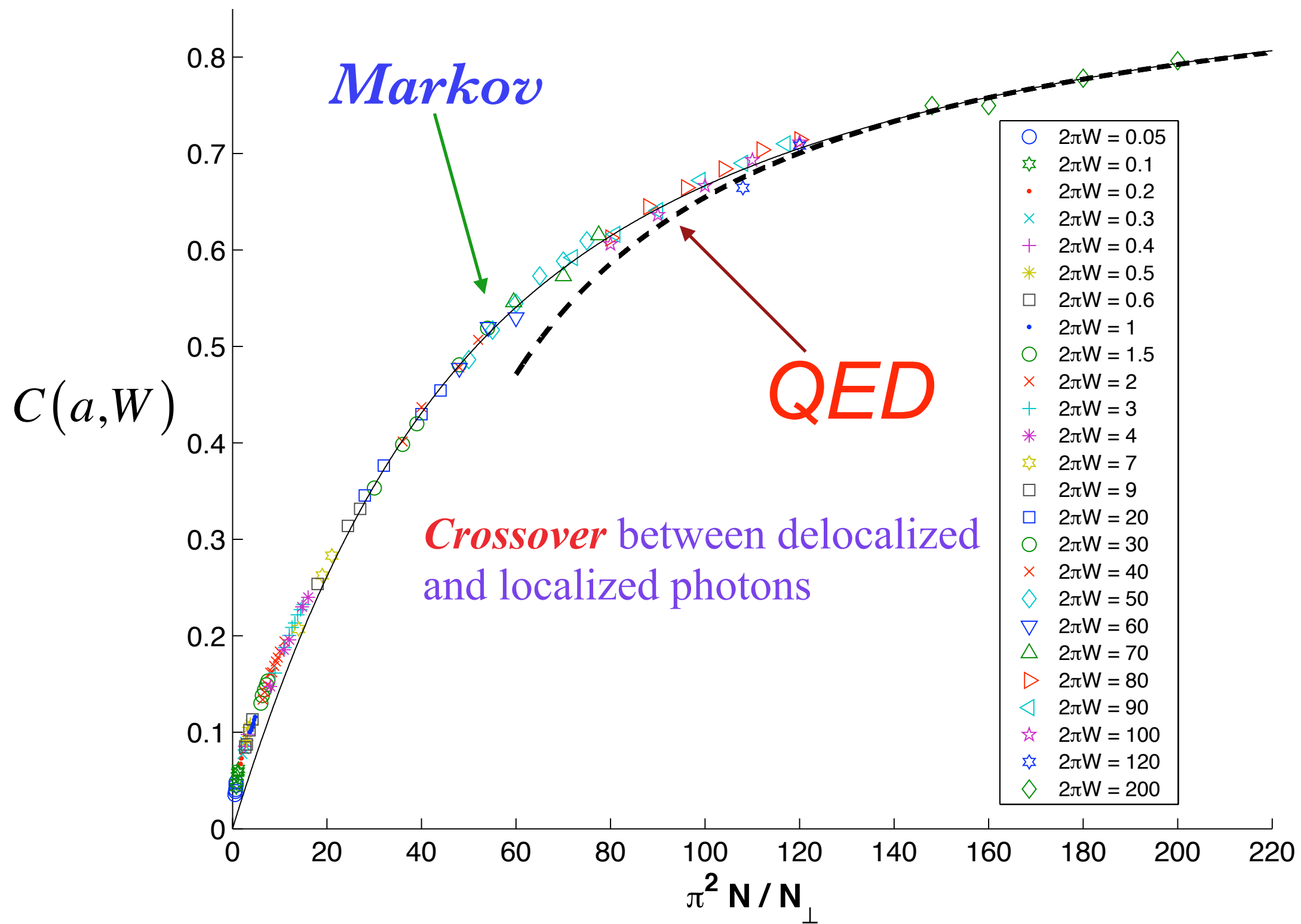
$$P(\Gamma) = 0 \quad \text{otherwise}$$

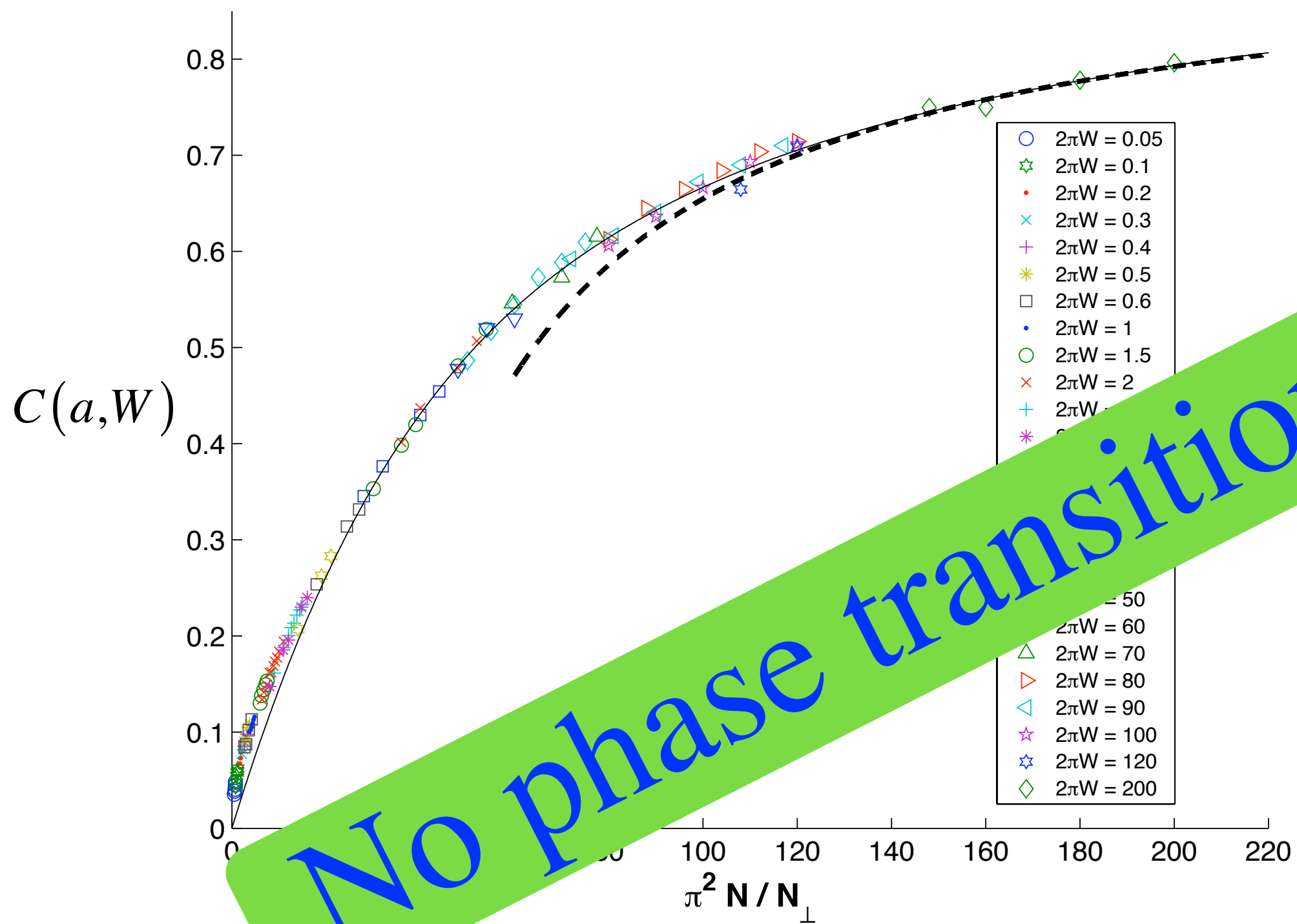
so that

$$C\left(\frac{N}{N_{\perp}}\right) = 1 - 3\frac{N_{\perp}}{N}$$



Phenomenological Markov process (Small world networks)





Dependence upon the space dimension ?

disorder driven localisation transition
(Anderson)

One-dimensional random atomic gas : Absence of single atom limit (Wigner-Weisskopf)

$d = 1$: Same expression of the effective atomic Hamiltonian H_e ,

$$H_e = \left(\hbar\omega_0 - i\frac{\hbar\Gamma_0}{2} \right) \sum_{i=1}^N |e_i\rangle\langle e_i| + \frac{\hbar\Gamma_0}{2} \sum_{i \neq j} V_{ij} \Delta_i^+ \Delta_j^-$$

with $V_{ij} = \beta_{ij} - i\gamma_{ij}$ but $\gamma_{ij} = \cos k_0 r_{ij}$ instead of $\gamma_{ij} = \frac{\sin k_0 r_{ij}}{k_0 r_{ij}}$

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Two limits :

$a = L/\lambda \gg 1$ dilute large sample limit (Wigner-Weisskopf + disorder effects)

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Two limits :

$a = L/\lambda \gg 1$ dilute large sample limit (Wigner-Weisskopf + disorder effects)

$a \ll 1$ Dicke limit (cooperative effects are expected) $\gamma_{ij} = \cos k_0 r_{ij} = \begin{pmatrix} 1 & 1 & \dots & 1 \\ 1 & 1 & \dots & 1 \\ \vdots & \vdots & & \vdots \\ 1 & 1 & \dots & 1 \end{pmatrix}$

so that $P(\Gamma) = \frac{1}{N} [(N-1)\delta(\Gamma) + \delta(\Gamma - N)]$

Method : Decomposition into a product of matrices

$N \times N$ matrix $U_{ij} = \cos k_0 r_{ij}$ can be written $U = \frac{1}{2} A^\dagger A$

with A is the $2 \times N$ matrix defined by $A_{0j} = e^{ik_0 r_j}$ and $A_{1j} = e^{-ik_0 r_j}$

U real symmetric matrix, its non vanishing eigenvalues are obtained from those of the 2×2 matrix U^\dagger

$$U^\dagger = \frac{1}{2} \begin{pmatrix} N & M \\ M^* & N \end{pmatrix}.$$

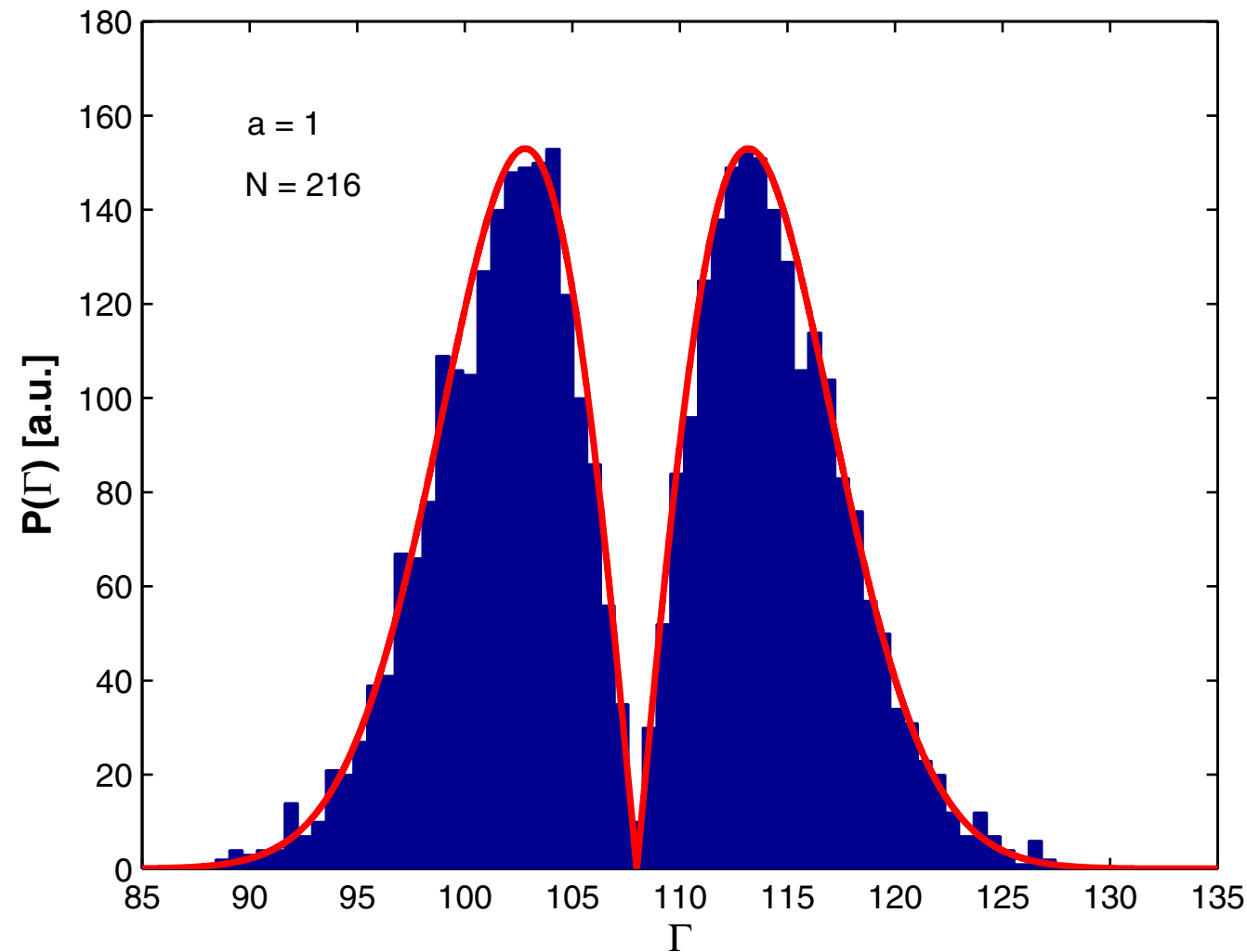
where $M = \sum_{k=1}^N e^{2ik_0 r_k}$ is a random variable.

The two eigenvalues of U^\dagger are $\lambda_{\pm} = \frac{N \pm |M|}{2},$

and the spectrum of U is

$$P(\Gamma) = \frac{1}{N} [(N-2)\delta(\Gamma) + \delta(\Gamma - \lambda_+) + \delta(\Gamma - \lambda_-)].$$

One-dimensional random atomic gas : Absence of single atom limit



Subradiant mode is
not represented

$$|M|^2 = N + \sum_{p \neq q} e^{2ik_0(r_p - r_q)}$$

Rayleigh distribution $P(|M|) = \frac{2|M|}{N} e^{-\frac{|M|^2}{N}}$

One-dimensional random atomic gas

- $d=1$: no crossover between localised and delocalised photons.
- Single atom (Wigner-Weisskopf) limit is never reached.
- Results in $d=1$ are valid for both **ordered** and **disordered** media (M is not a random variable)
- Cooperative effects (not disorder) is the mechanism underlying photon localisation in $d=1$.

Two-dimensional random atomic gas : Marchenko-Pastur distribution

In $d = 2$ the same expression of the effective atomic Hamiltonian H_e holds,

with $V_{ij} = \beta_{ij} - i\gamma_{ij}$ but $\gamma_{ij} = J_0(k_0 r_{ij})$ instead of $\gamma_{ij} = \frac{\sin k_0 r_{ij}}{k_0 r_{ij}}$

Two-dimensional random atomic gas : Marchenko-Pastur distribution

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The $d = 1$ trick does not work.

Instead we use the general decomposition: $U = H T H^\dagger$,

where $T(M \times M)$, $H(N \times M)$

(S. Skipetrov and also “Free probability theory” (Voiculescu), “Wireless communications” (Debbah, Tulino, Verdu), “spin glasses” I. Kanter et al.

Two-dimensional random atomic gas : Marchenko-Pastur distribution

Spectrum of $\gamma_{ij} = J_0(k_0 r_{ij})$

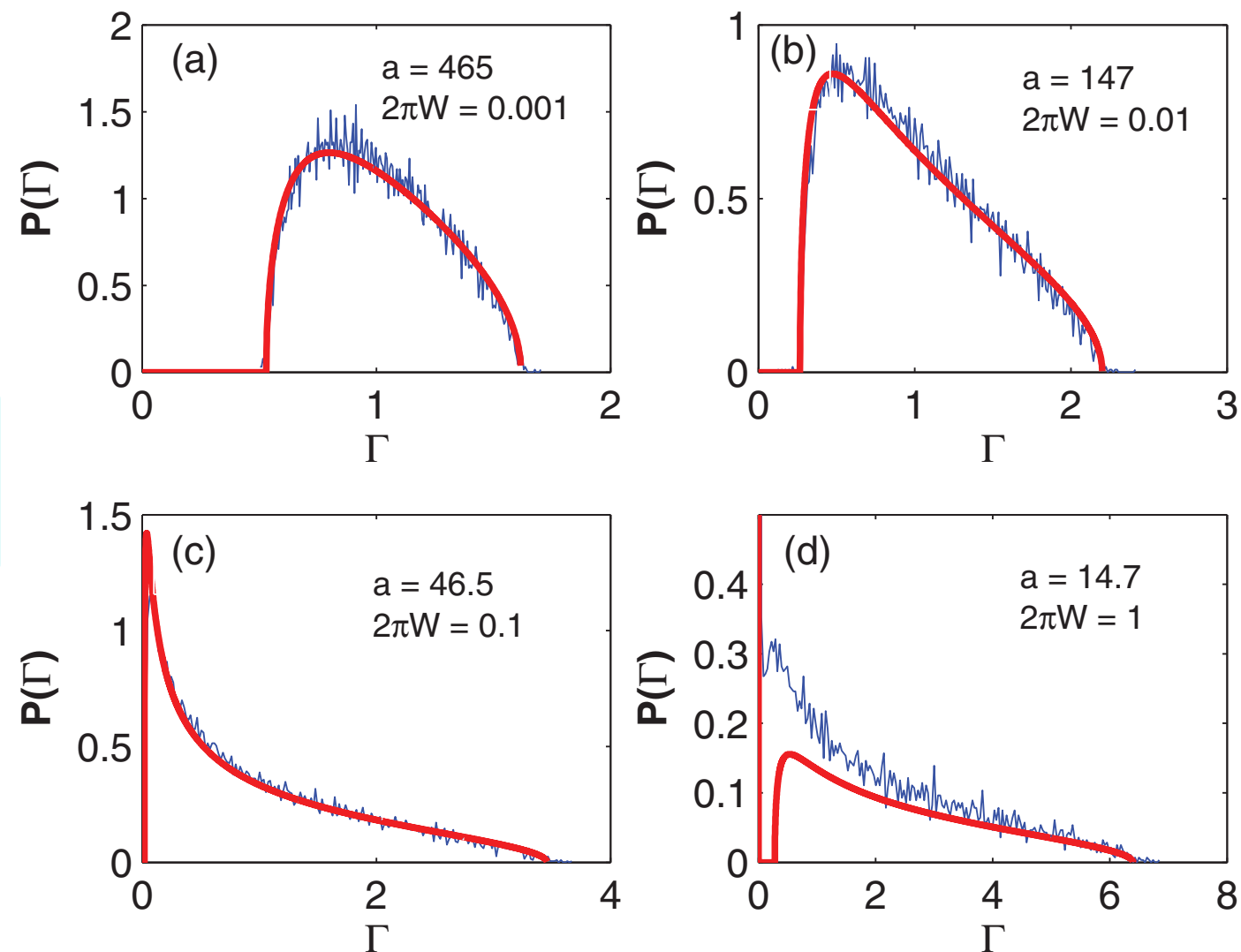
For $a \gg 1$, dilute limit,

$$P(\Gamma) \simeq \left(1 - \frac{M}{N}\right)^+ \delta(\Gamma) + \frac{\sqrt{(\Gamma - \Gamma_-)^+(\Gamma_+ - \Gamma)^+}}{2\pi \frac{N}{M} \Gamma},$$

$$x^+ = \max(0, x)$$

$$\Gamma_{\pm} = (1 \pm \sqrt{N/M})^2$$

$$M = 2\pi a$$



Disorder parameter

$$L^d = (\lambda a)^d$$

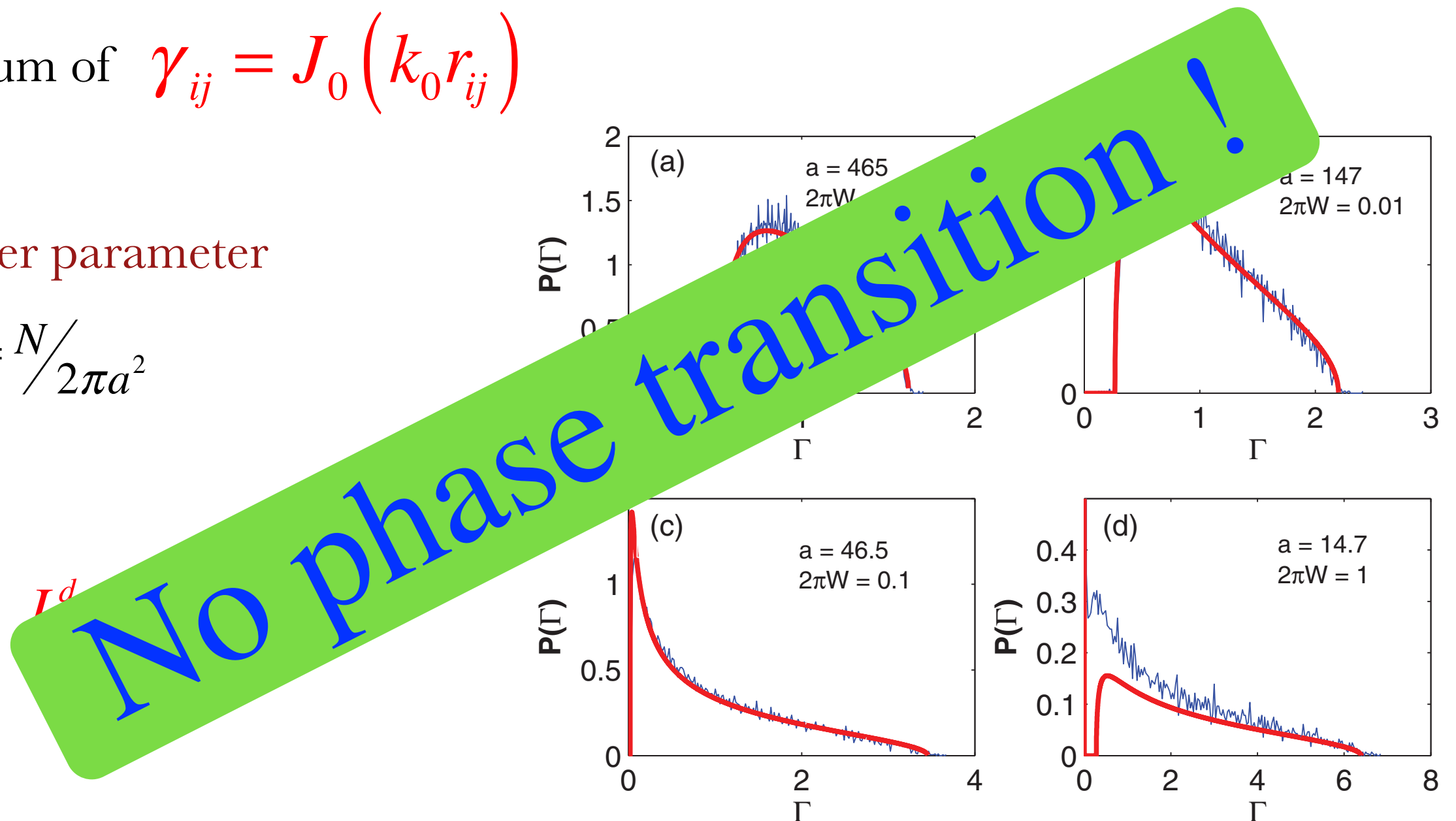
$$W = N / 2\pi a^2$$

Two-dimensional random atomic gas : Marchenko-Pastur distribution

Spectrum of $\gamma_{ij} = J_0(k_0 r_{ij})$

Disorder parameter

$$W = N / 2\pi a^2$$

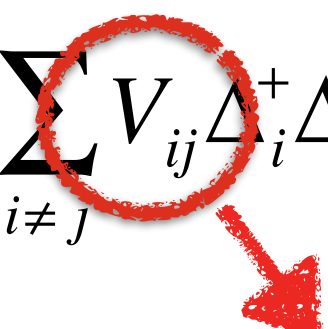


Numerical results $N=216$

Eigenvalues of the non Hermitian random Hamiltonian

Time evolution of the ground state population is driven by the eigenvalues of the random matrix γ_{ij}

while the effective Hamiltonian is

$$H_e = \left(\hbar\omega_0 - i\frac{\hbar\Gamma_0}{2} \right) \sum_{i=1}^N |e_i\rangle\langle e_i| + \frac{\hbar\Gamma_0}{2} \sum_{i \neq j} V_{ij} \Delta_i^+ \Delta_j^-$$


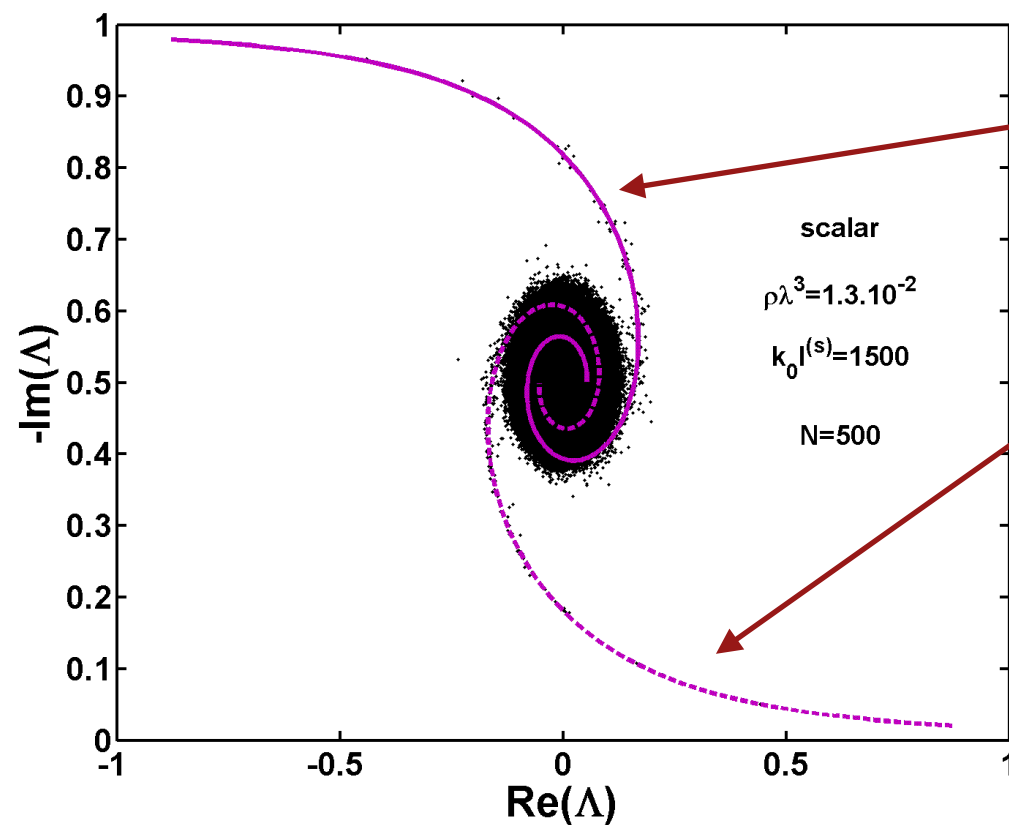
$$V_{ij} = \beta_{ij} - i\gamma_{ij}$$

Study the complex eigenvalues of H_e

$$E_n - i\hbar \frac{\Gamma_n}{2} \equiv \hbar\omega_0 + \hbar\Gamma_0 \Lambda_n$$

Eigenvalues of the non Hermitian random Hamiltonian

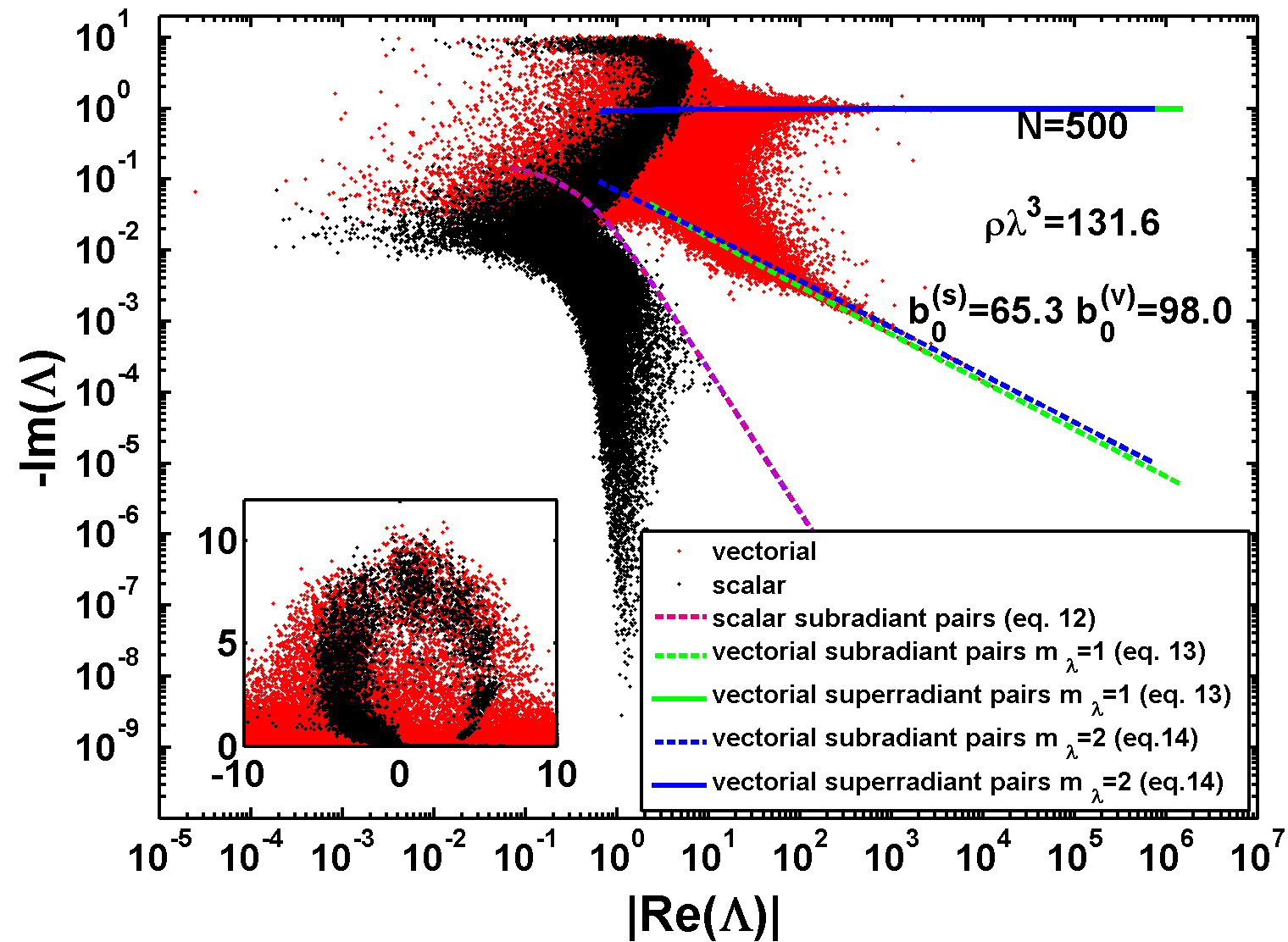
N=2 atoms case : The spectrum of H_e can be obtained explicitly (for both scalar and vectorial case)



Cooperative pairs

$$E_n - i\hbar \frac{\Gamma_n}{2} \equiv \hbar\omega_0 + \hbar\Gamma_0 \Lambda_n$$

For N large and in the dense limit



Thouless parameter : localisation phase transition

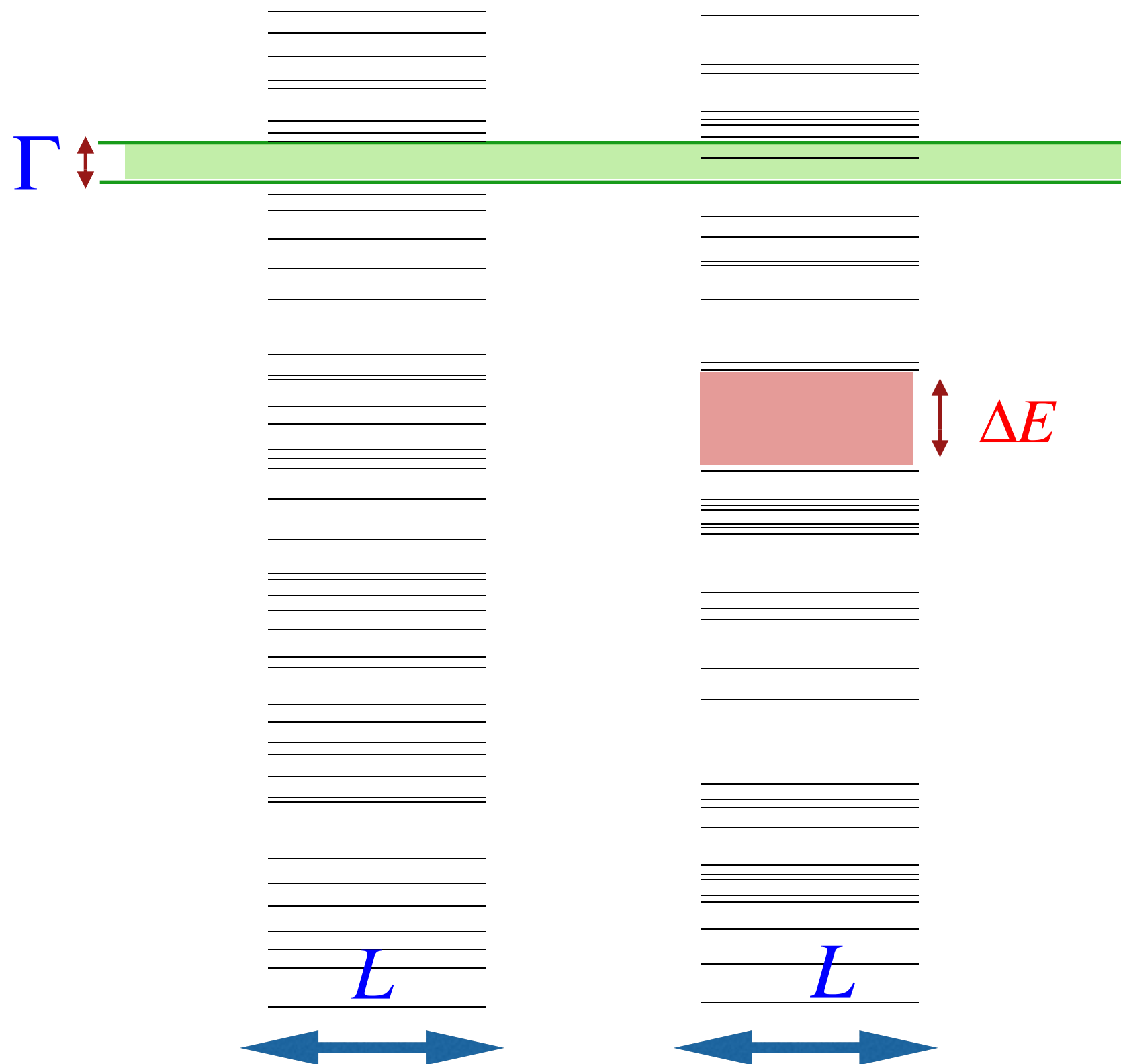
Edwards & Thouless ('72), Thouless ('77)

Coupling between open quantum systems
Transport (conductance)

Using Random matrix theory :

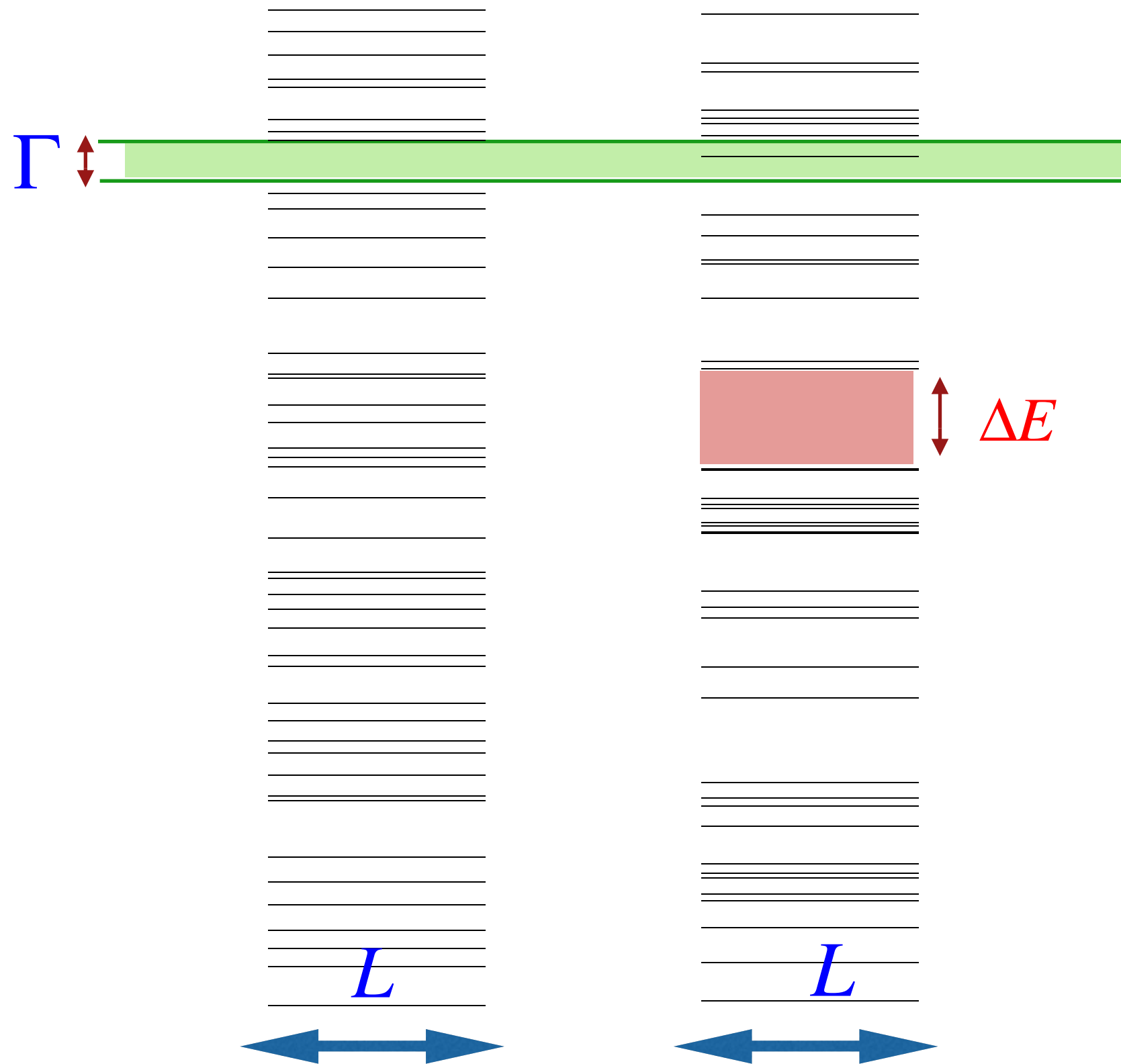
G. Montambaux, E.A., (1992), I. Guarneri et al. (1994)

Thouless parameter - Resonance overlap



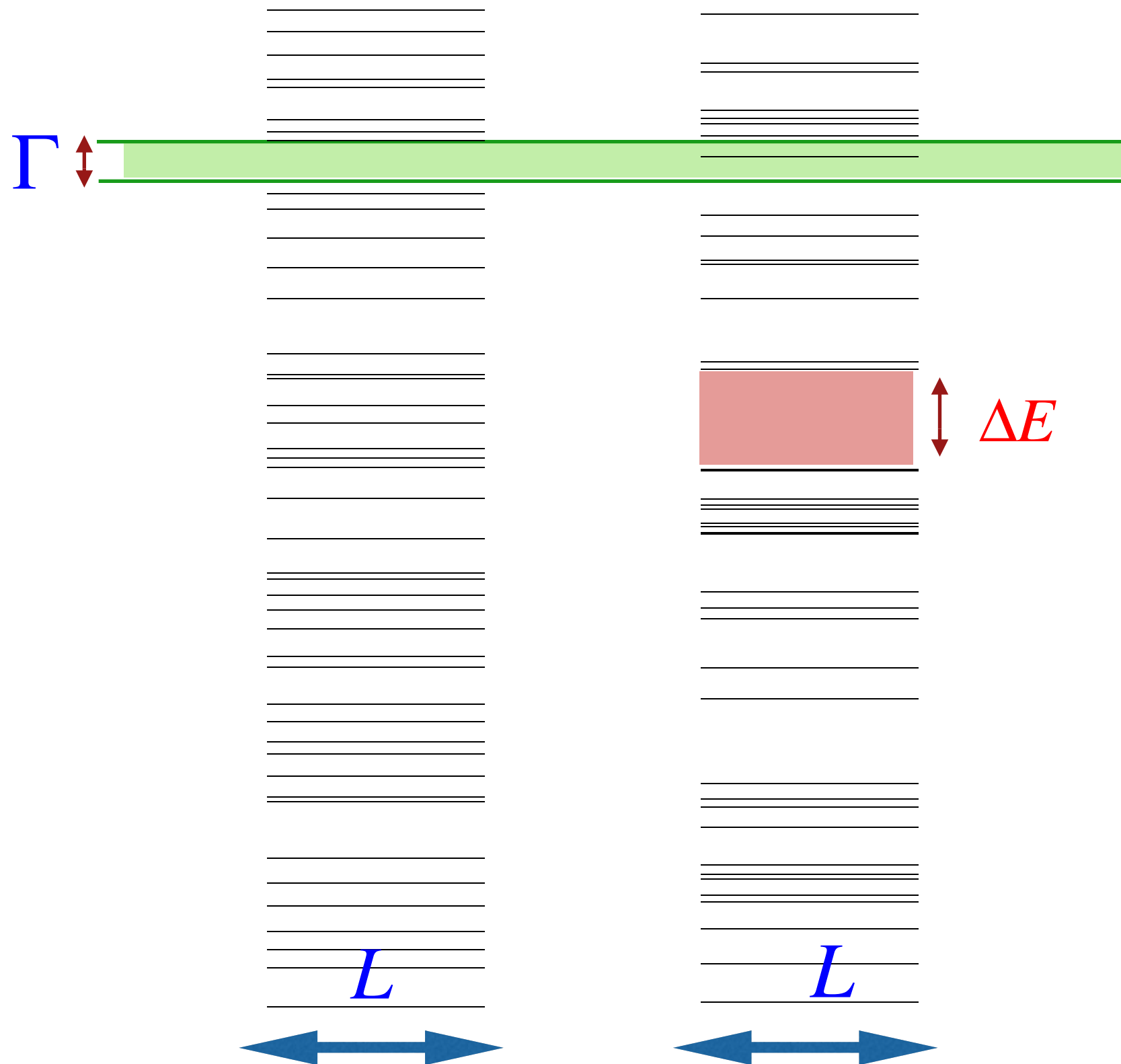
Energy spectrum
of two random
quantum systems

Thouless parameter - Resonance overlap



$$g(L) = \left\langle \frac{\langle \Gamma \rangle_i}{\langle \Delta E \rangle_i} \right\rangle$$

Thouless parameter - Resonance overlap



$$g(L) = \left\langle \frac{\langle \Gamma \rangle_i}{\langle \Delta E \rangle_i} \right\rangle$$

$$g(L \rightarrow \infty) \gg 1$$

large overlap : delocalised
states - conductor

$$g \ll 1$$

Small overlap : localised
states - insulator

Thouless parameter - localisation phase transition

Scaling and its meaning : (P.W. Anderson *et al.*, 1979)

If we know $g(L)$, we know it at any scale :

$$g(L(1 + \varepsilon)) = f(g(L), \varepsilon)$$

Scaling behavior :

$$g(L, W) = f\left(\frac{L}{\xi(W)}\right)$$

$\xi(W)$ is the localization length

Thouless parameter - localisation phase transition

Scaling and its meaning : (P.W. Anderson *et al.*, 1979)

If we know $g(L)$, we know it at any scale :

$$g(L(1 + \varepsilon)) = f(g(L), \varepsilon)$$

$$\beta(g) = \frac{d \ln g}{d \ln L}$$

is a function of g only.

Scaling behavior :

$$g(L, W) = f\left(\frac{L}{\xi(W)}\right)$$

$\xi(W)$ is the localization length

Thouless parameter - localisation phase transition

$d = 2$

$d = 3$

$g(L, W)$

Anderson phase transition

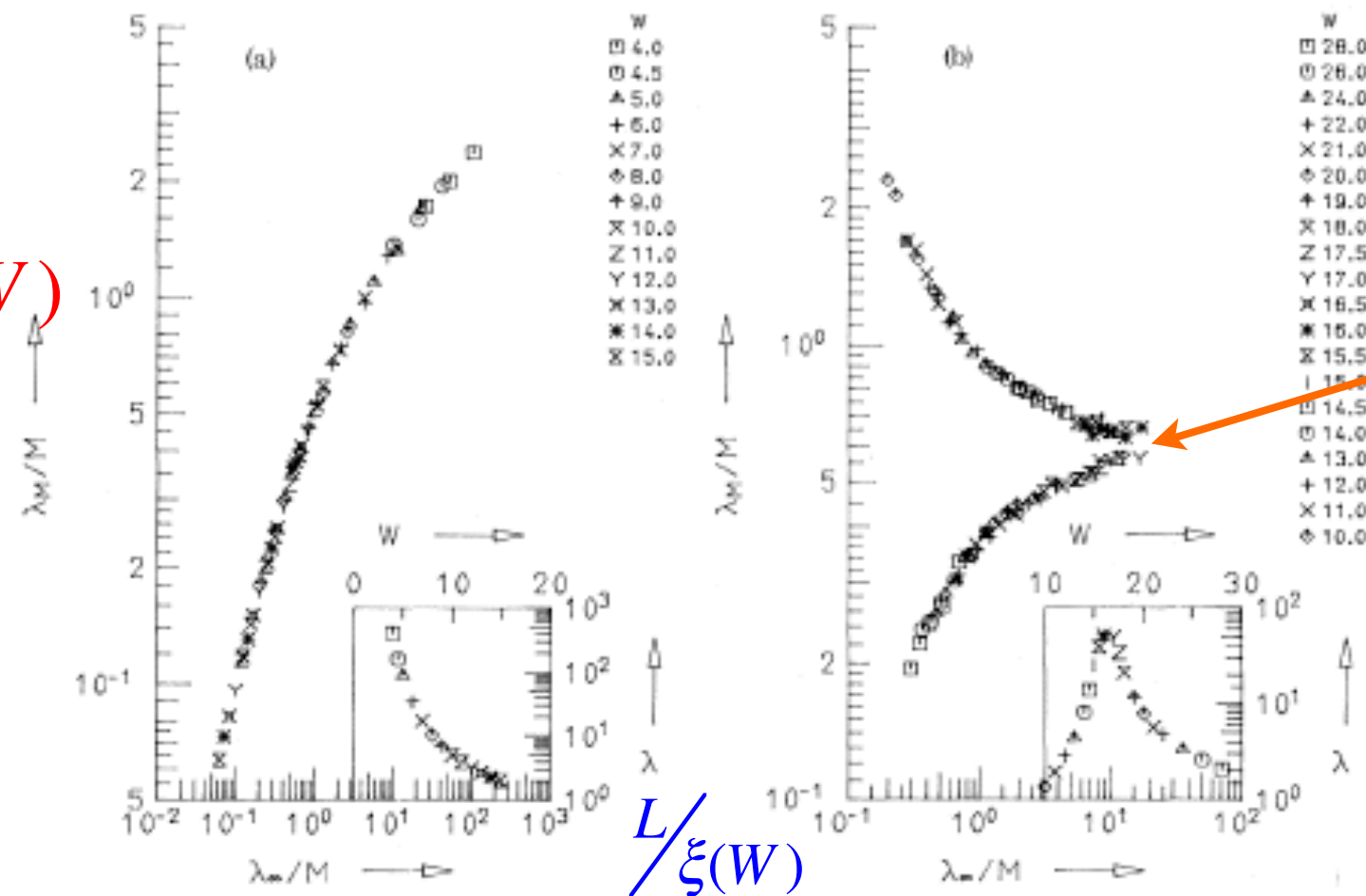


FIG. 1. Scaling function λ_M/M vs λ_m/M for the localization length λ_M of a system of thickness M for (a) $d=2$ ($M \geq 4$) and (b) $d=3$ ($M \geq 3$). Insets show the scaling parameter λ_m as a function of the disorder W .

Scaling behaviour - phase transition

Thouless scaling parameter (conductance)

$$g(L) = \left\langle \frac{\langle \Gamma \rangle_i}{\langle \Delta E \rangle_i} \right\rangle$$

$$E_n - i\hbar \Gamma_n / 2 \equiv \hbar \omega_0 + \hbar \Gamma_0 \Lambda_n$$

Scaling behaviour - phase transition

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because of the constraint : $\langle \Gamma \rangle_i = -2\text{Tr}(\Lambda)/N = 1$

Scaling behaviour - phase transition

Thouless scaling parameter (conductance)

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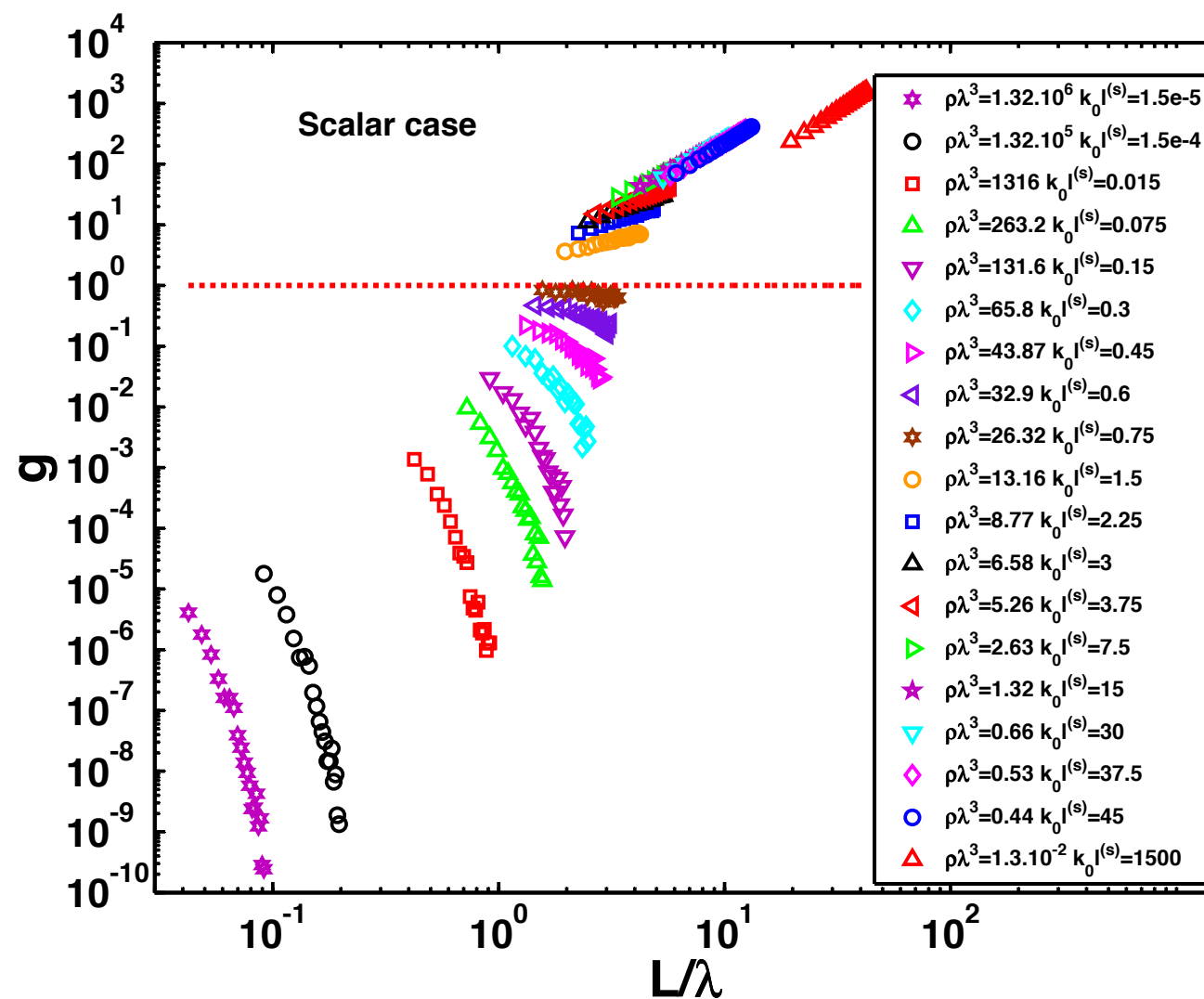
because of the constraint : $\langle \Gamma \rangle_i = -2\text{Tr}(\Lambda)/N = 1$

Instead we define :

$$g \equiv \left\langle \frac{1}{\langle \frac{1}{\Gamma} \rangle_i \langle \Delta E \rangle_i} \right\rangle$$

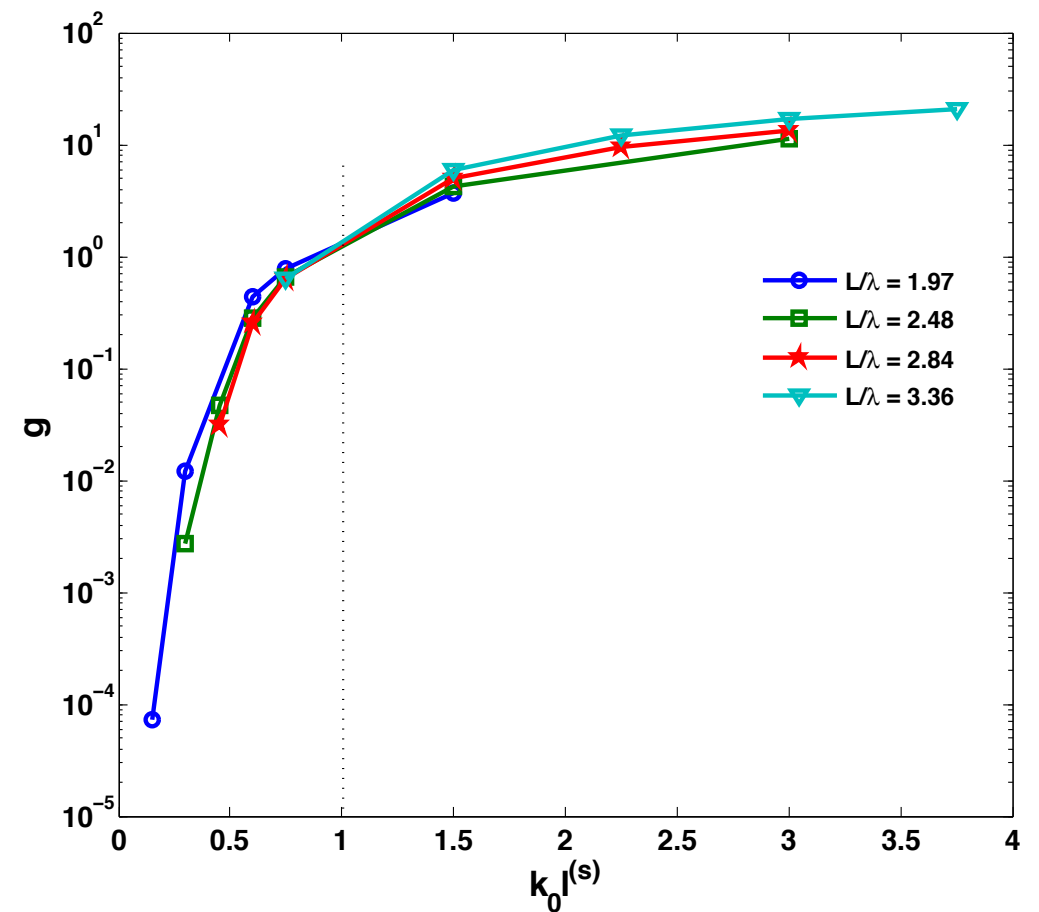
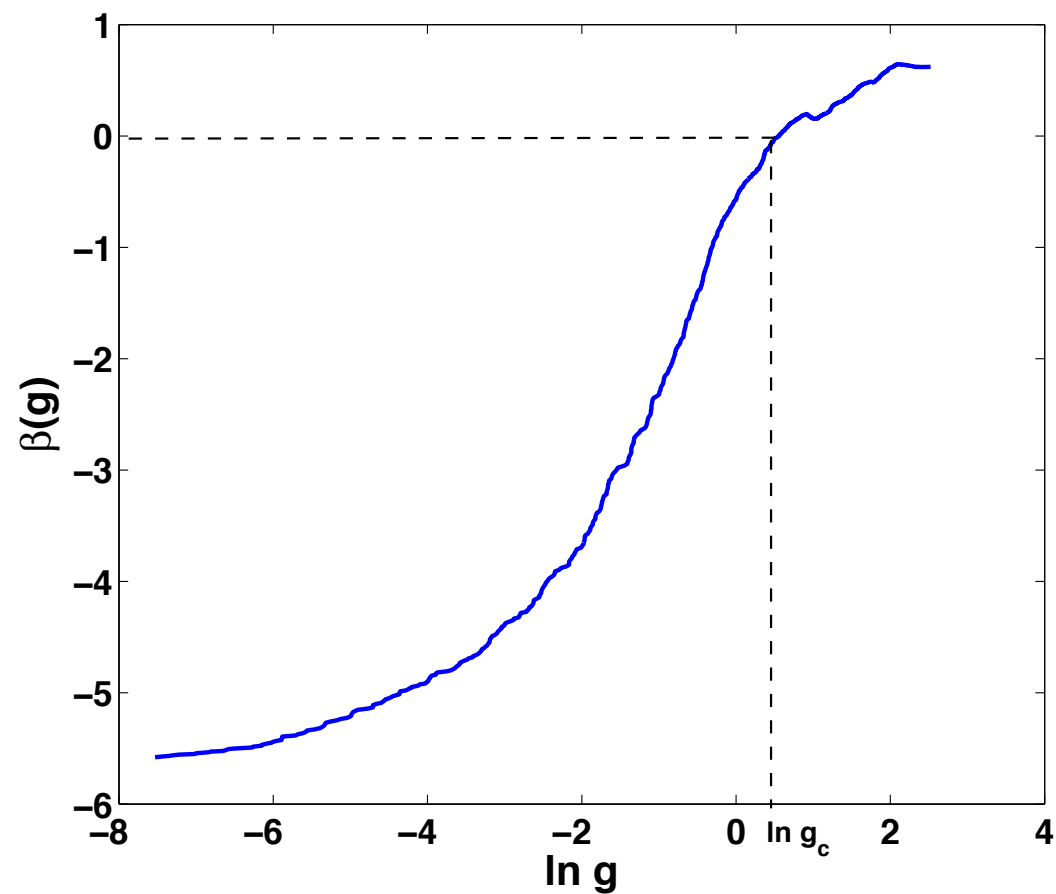
Scaling behaviour - phase transition

$$g \equiv \left\langle \frac{1}{\left\langle \frac{1}{\Gamma} \right\rangle_i \langle \Delta E \rangle_i} \right\rangle$$



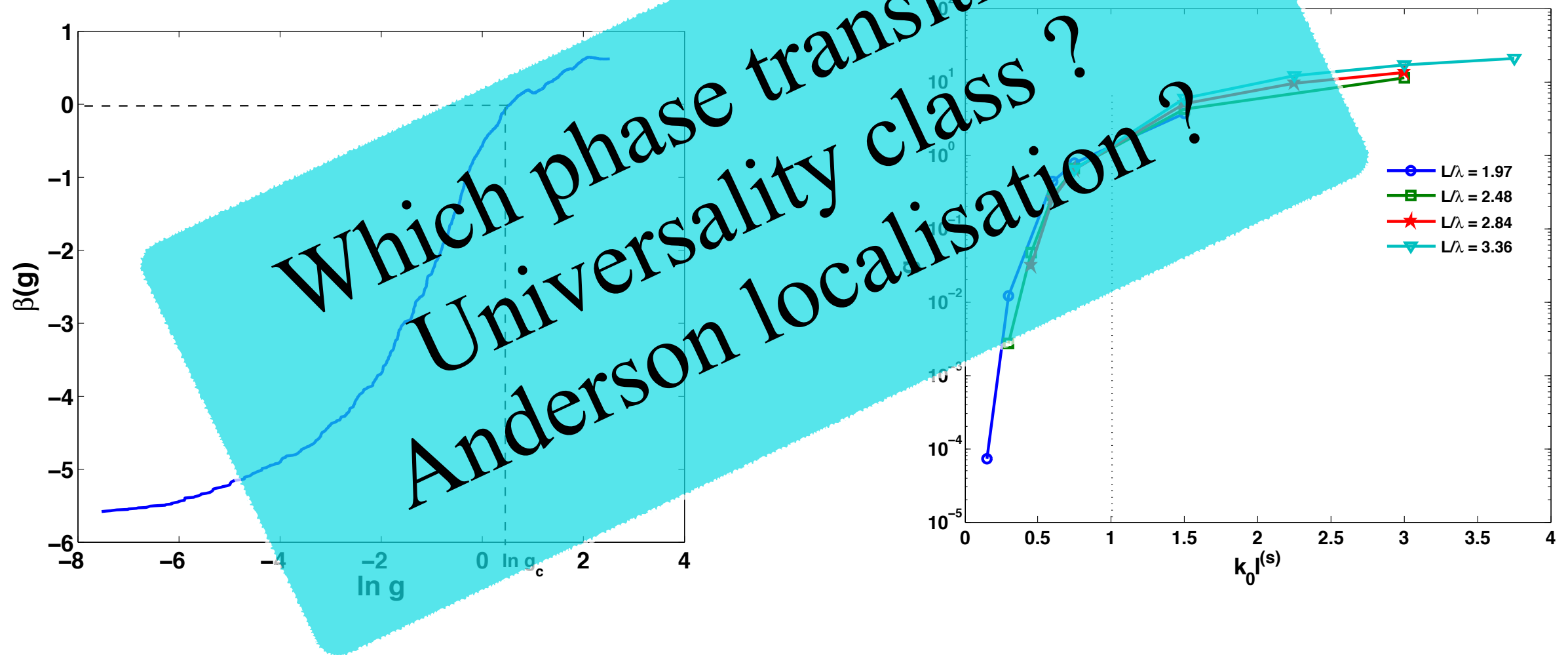
Critical point for a phase transition :

$g(L)$ becomes L - independent i.e. $\beta(g) = \frac{d \ln g}{d \ln L} = 0$



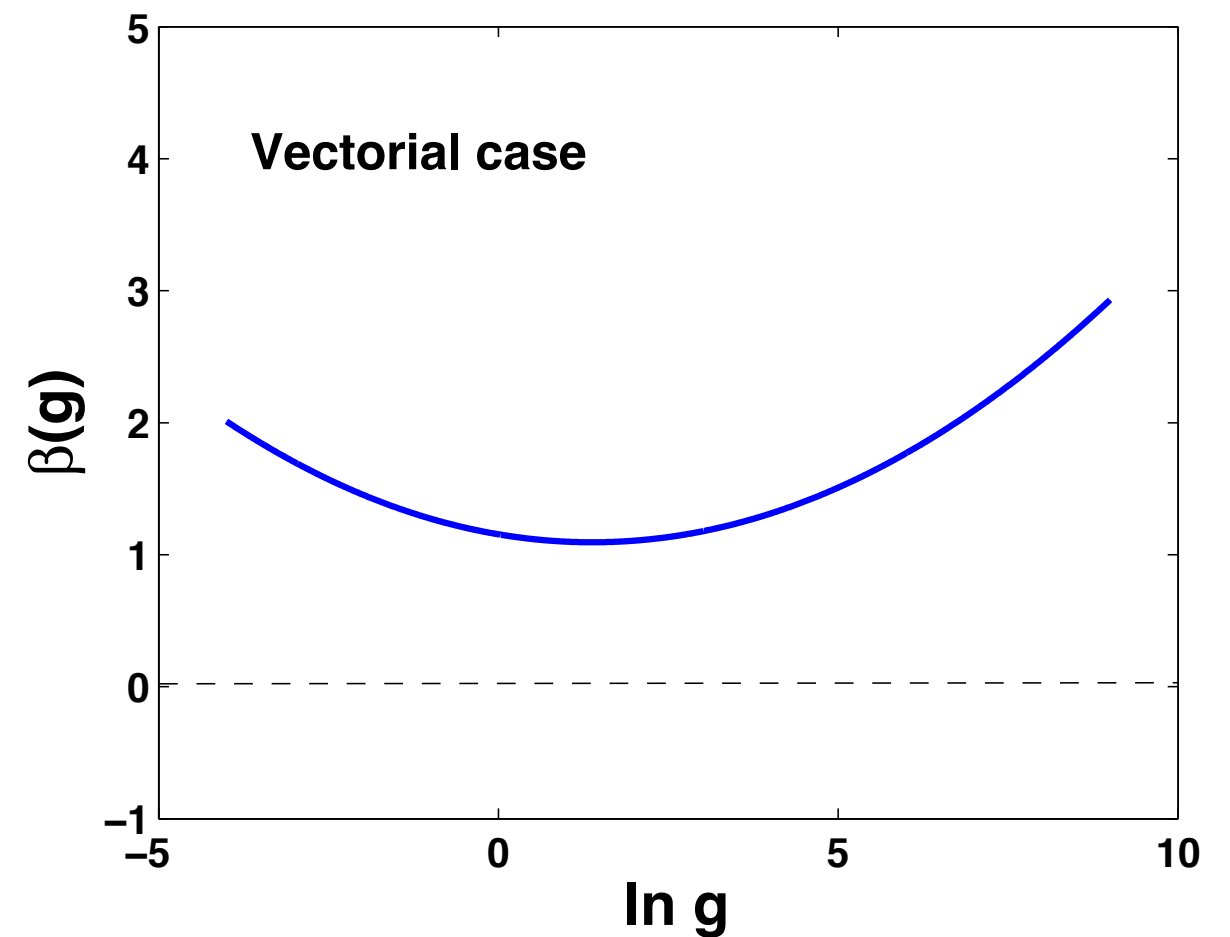
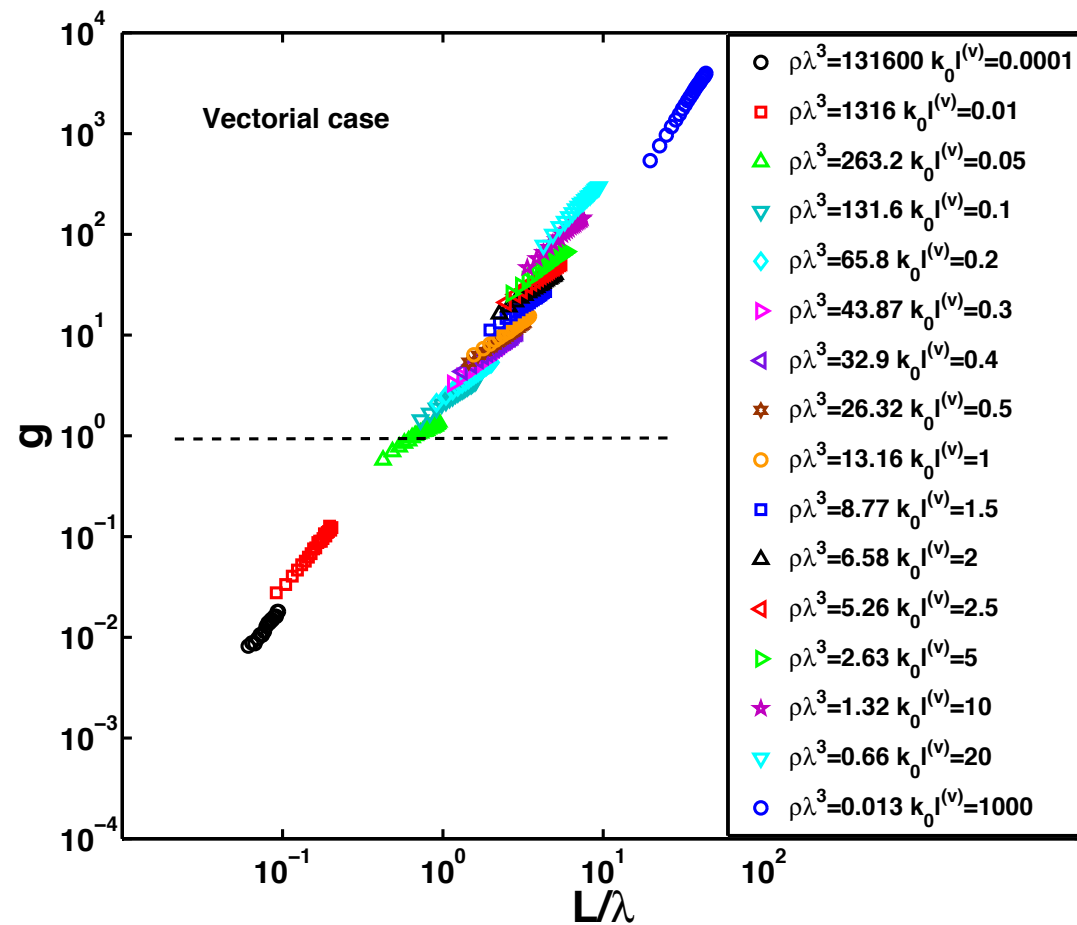
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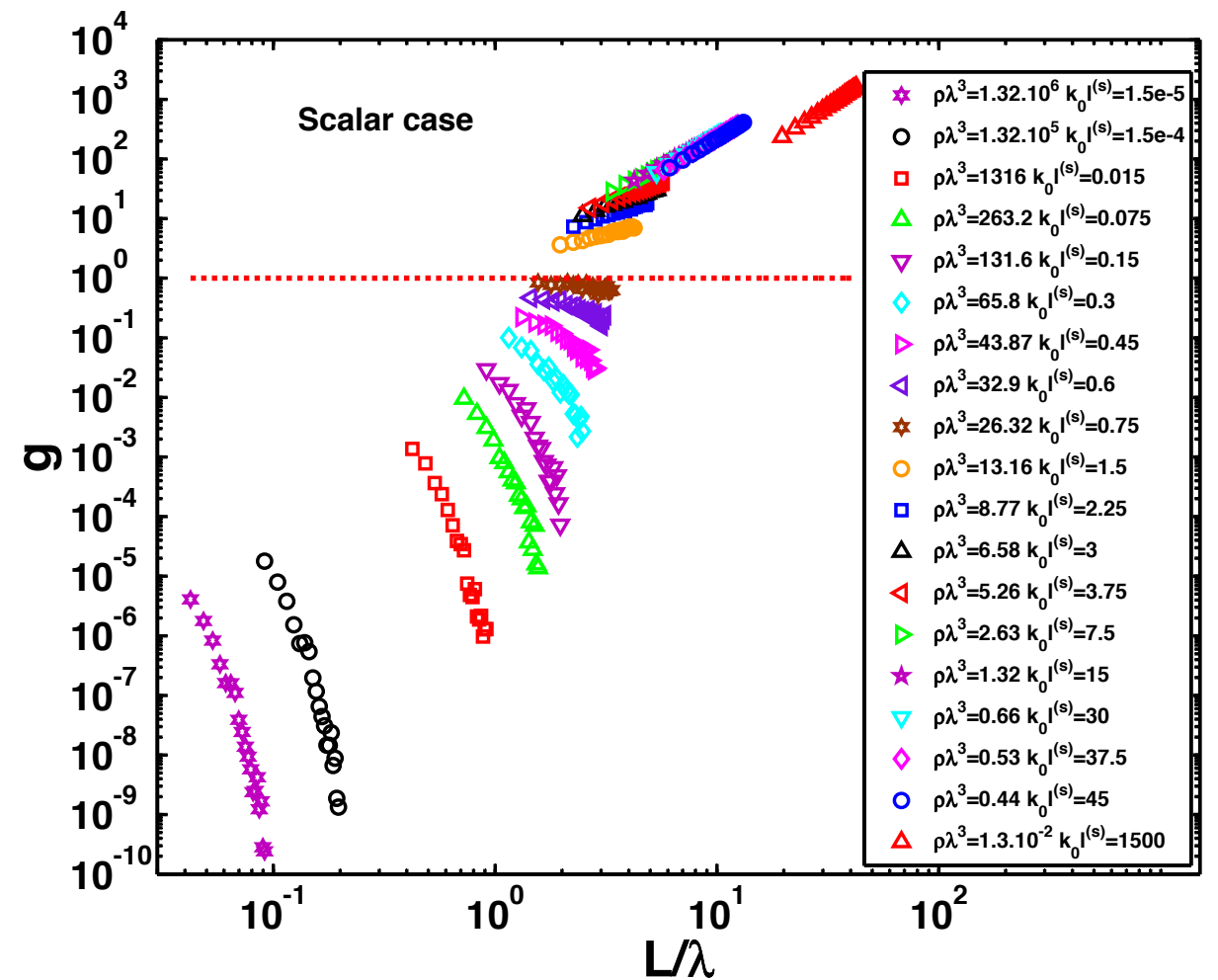
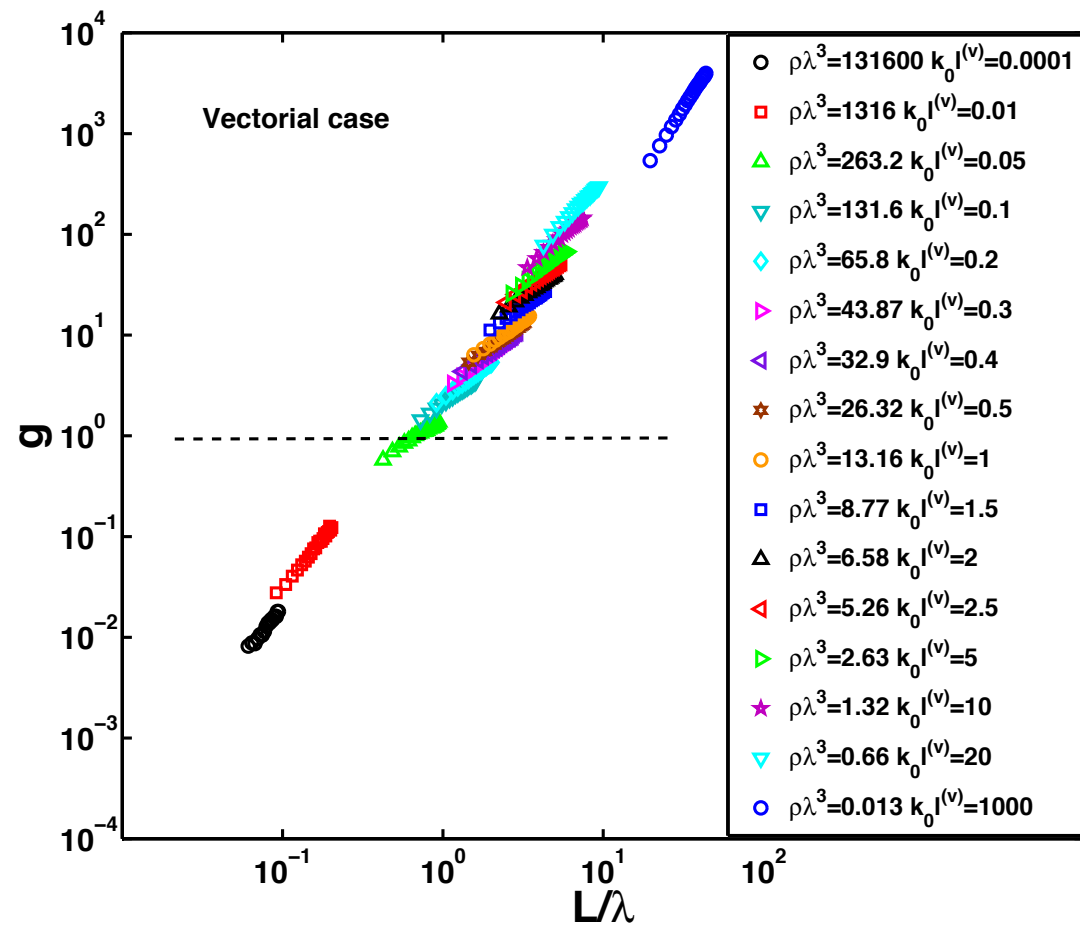
Scaling behaviour - phase transition to make things more complicated

Vector case - polarised waves



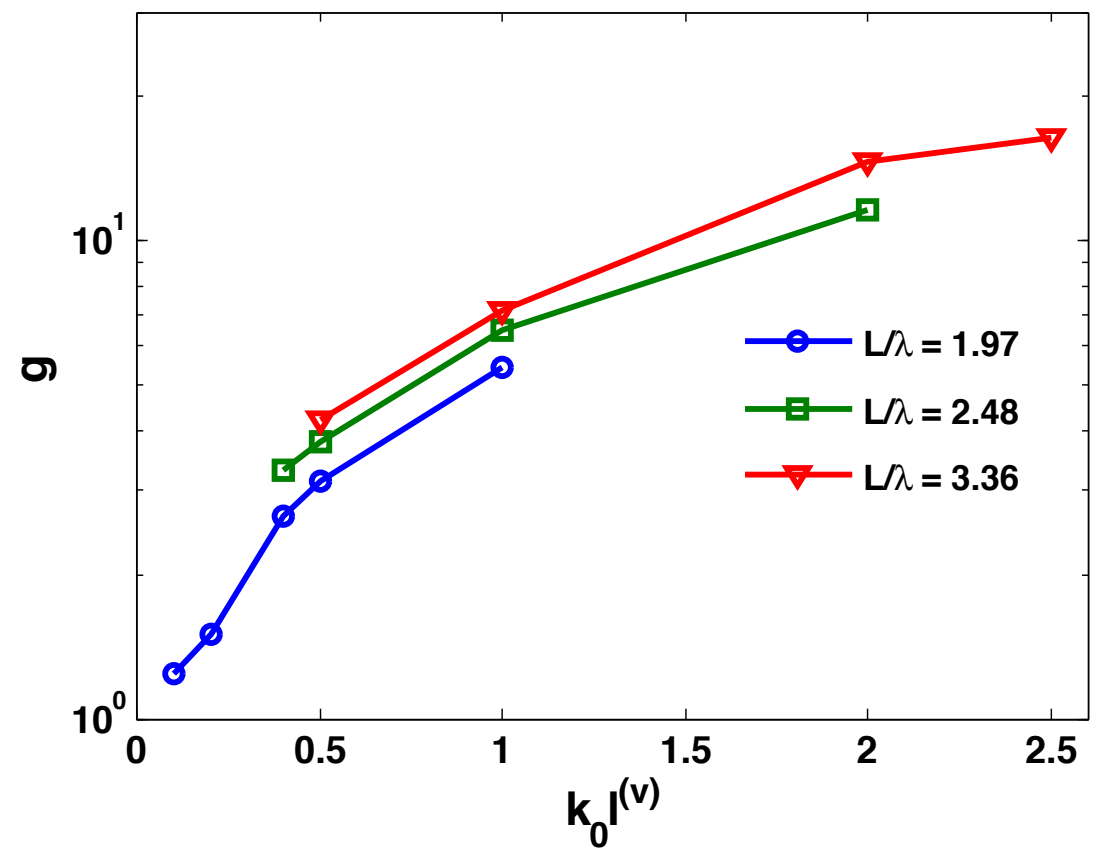
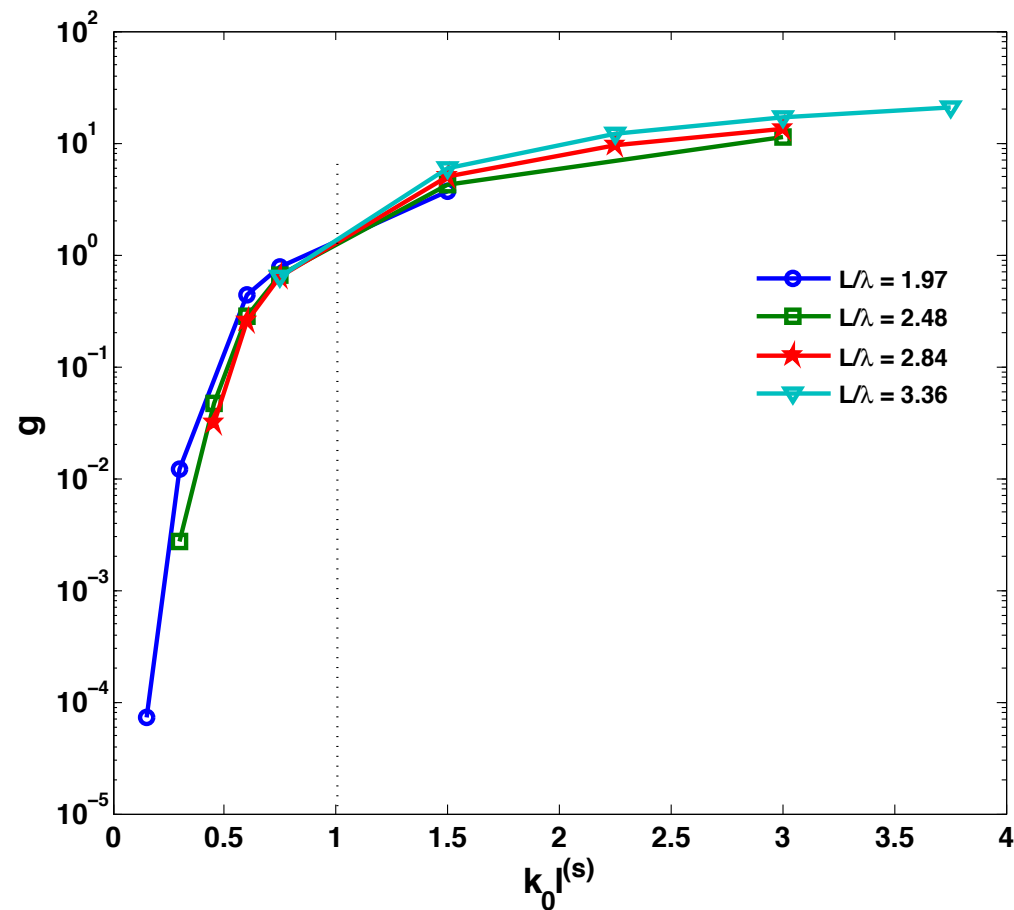
Scaling behaviour - phase transition to make things more complicated

Vector case - polarised waves



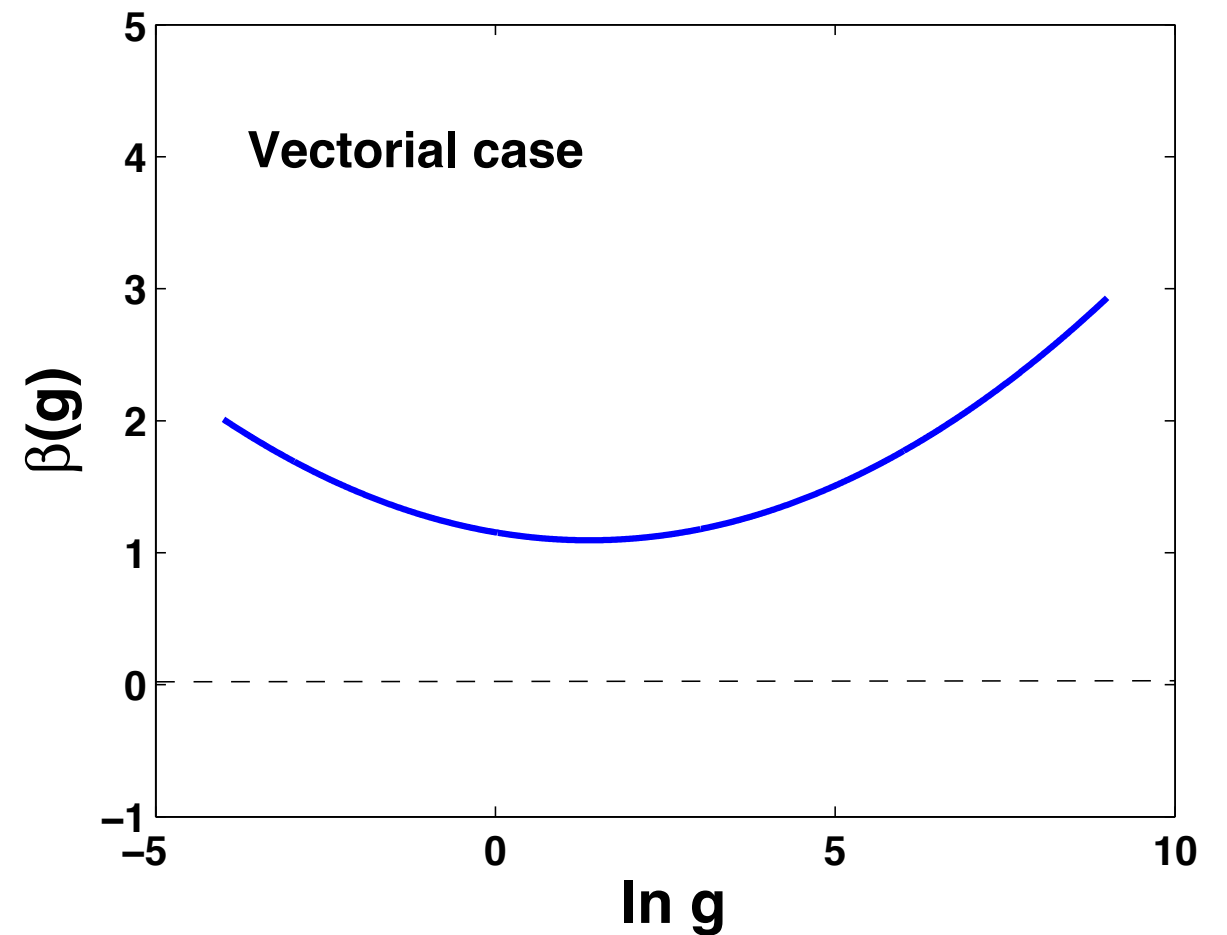
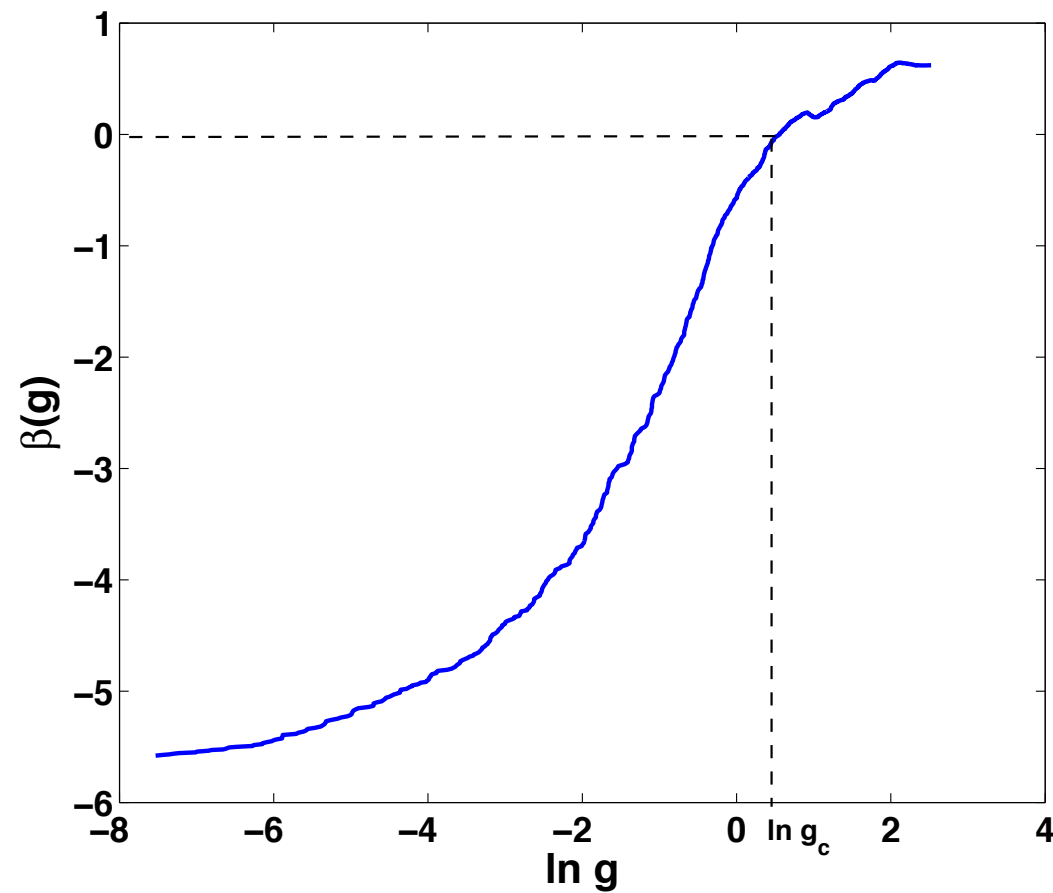
Scaling behaviour - phase transition to make things more complicated

Vector case - polarised waves



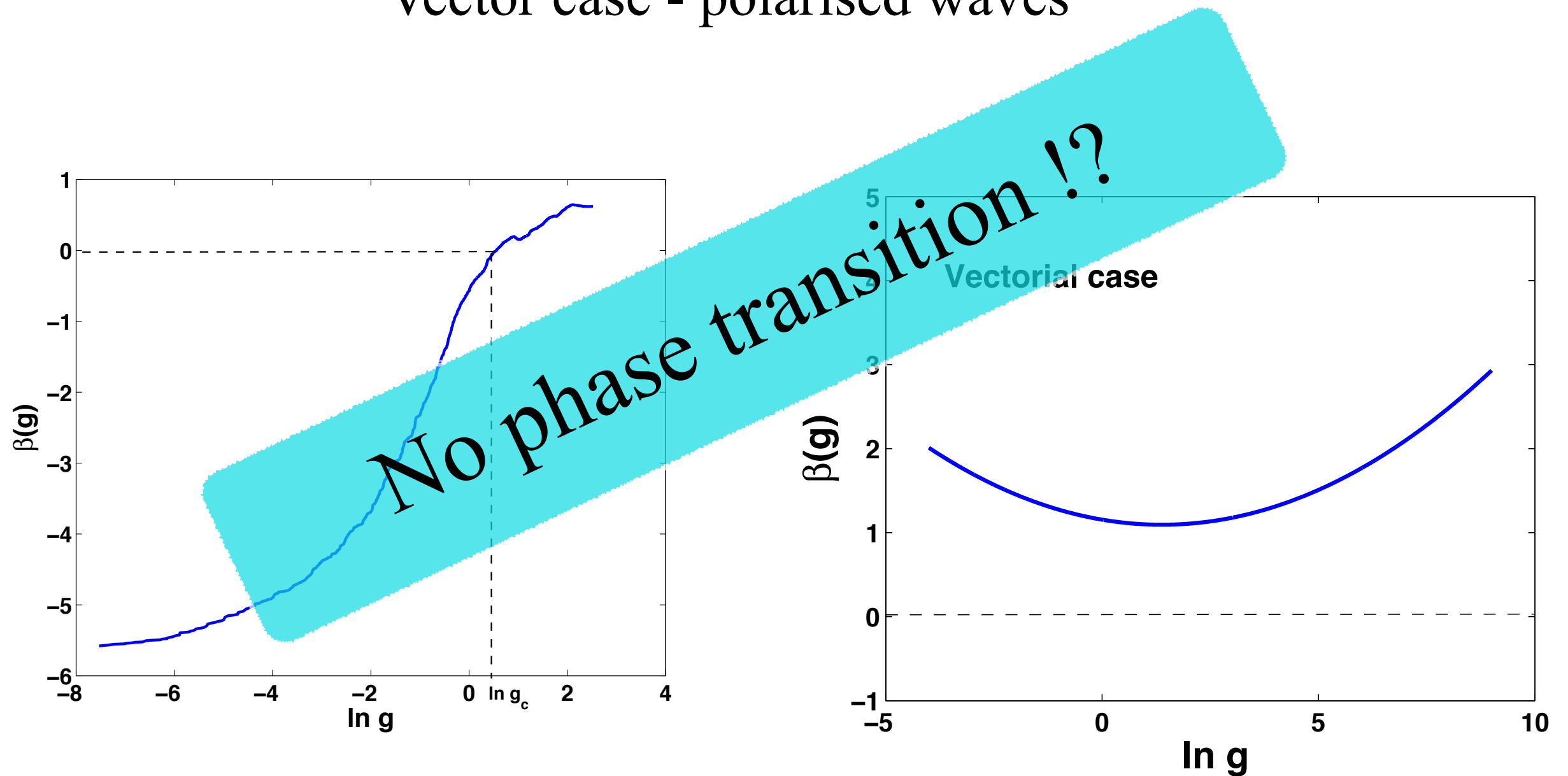
Scaling behaviour - phase transition to make things more complicated

Vector case - polarised waves



Scaling behaviour - phase transition to make things more complicated

Vector case - polarised waves



Conclusion - Summary

- Study of the scaling properties of the Non Hermitian Euclidean random Hamiltonian

$$H_e = \left(\hbar\omega_0 - i\frac{\hbar\Gamma_0}{2} \right) \sum_{i=1}^N |e_i\rangle\langle e_i| + \frac{\hbar\Gamma_0}{2} \sum_{i \neq j} V_{ij} \Delta_i^+ \Delta_j^-$$

$$\text{with } V_{ij} = \beta_{ij} - i\gamma_{ij}$$

- H_e accounts for **cooperative** properties of the atomic gas (Super- and Sub-radiance). It also depends on the disorder.
- The radiation pattern is well accounted by the part γ_{ij} of the interaction.
- The distribution of eigenvalues of γ_{ij} exhibits scaling properties but there is *no indication of the existence of a phase transition* driven either by disorder or interactions.

- The interplay between disorder and cooperative effects depend upon the space dimensionality.
- For $d = 2, 3$, there is a *crossover* between a delocalised (Wigner-Weisskopf) regime and a behaviour driven by cooperative effects (eventually Dicke regime)
- For $d = 1$, there is no single atom limit.
- The eigenvalue distribution of the whole Hamiltonian H_e exhibits also scaling properties. *A critical behaviour is obtained for scalar waves* using a conveniently defined Thouless conductance for that problem.
- *The critical behaviour disappears for vector waves.*
- The nature and universality of this transition is still unclear.
- Set of new experimental efforts to probe the interplay of disorder and cooperative effects (R. Kaiser, A. Browaeys, M. Havey,...)

Thank you for your attention.

