

Not Yet

Introduction to Mesoscopics

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ORIGINS

E.P. Wigner, Conference on Neutron Physics by Time of Flight, November 1956

P.W. Anderson, “*Absence of Diffusion in Certain Random Lattices*”; Phys.Rev., **1958**, v.109, p.1492

L.D. Landau, ”*Fermi-Liquid Theory*” Zh. Exp. Teor. Fiz.,**1956**, v.30, p.1058

J. Bardeen, L.N. Cooper & J. Schriffer, “*Theory of Superconductivity*”; Phys.Rev., **1957**, v.108, p.1175.

Part 1 Without interactions

**Random Matrices, Anderson
Localization, and Quantum Chaos**

RANDOM MATRIX THEORY

$N \times N$

*ensemble of Hermitian matrices
with **random** matrix element*

$N \rightarrow \infty$

E_α

- spectrum (set of eigenvalues)

$$\delta_1 \equiv \langle E_{\alpha+1} - E_\alpha \rangle$$

- mean level spacing

$\langle \dots \rangle$

- ensemble averaging

$$s \equiv \frac{E_{\alpha+1} - E_\alpha}{\delta_1}$$

- spacing between nearest neighbors

$P(s)$

- distribution function of nearest neighbors spacing between

Spectral Rigidity

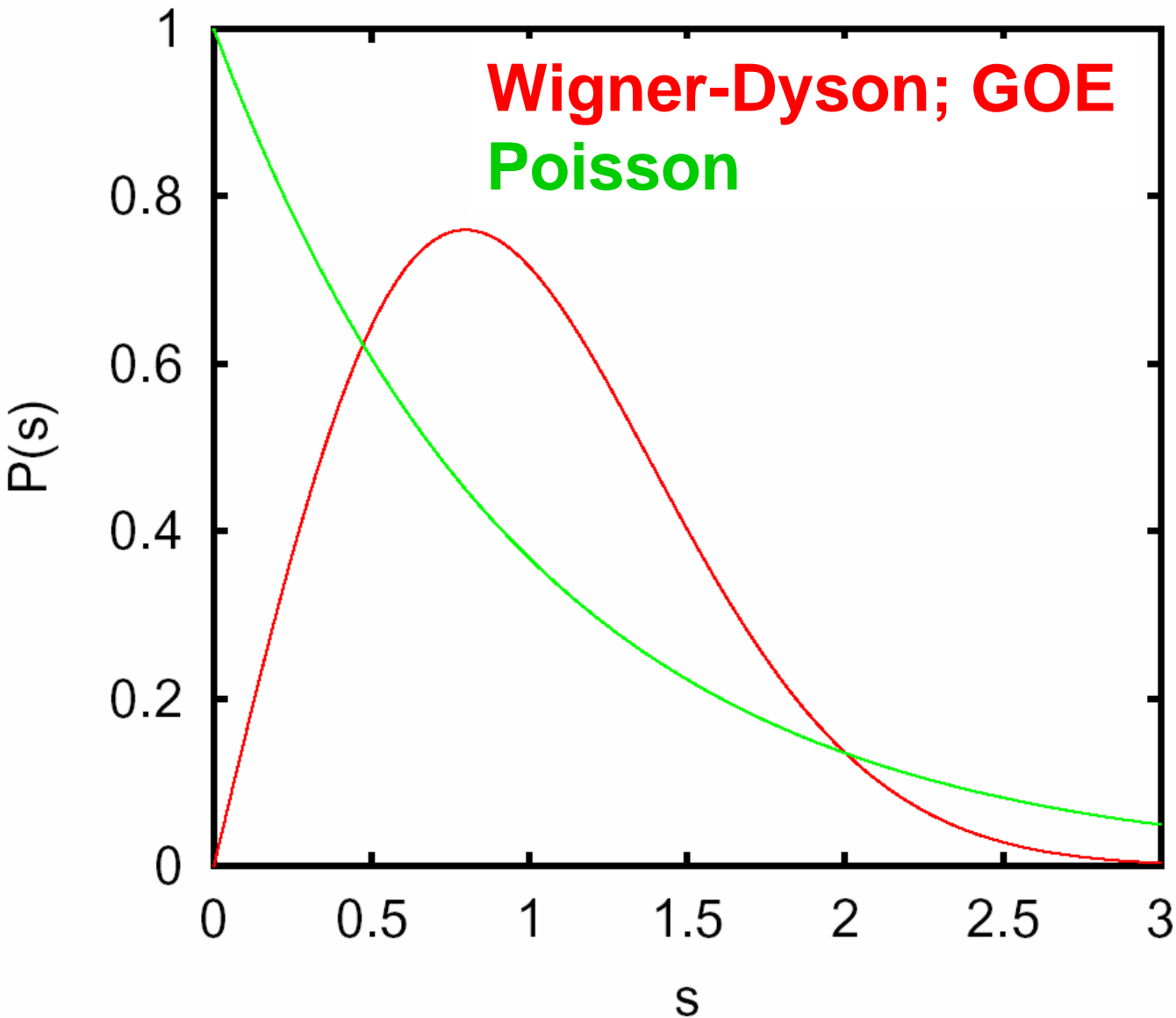
$$P(s = 0) = 0$$

Level repulsion

$$P(s \ll 1) \propto s^\beta \quad \beta=1,2,4$$

Dyson Ensembles and Hamiltonian systems

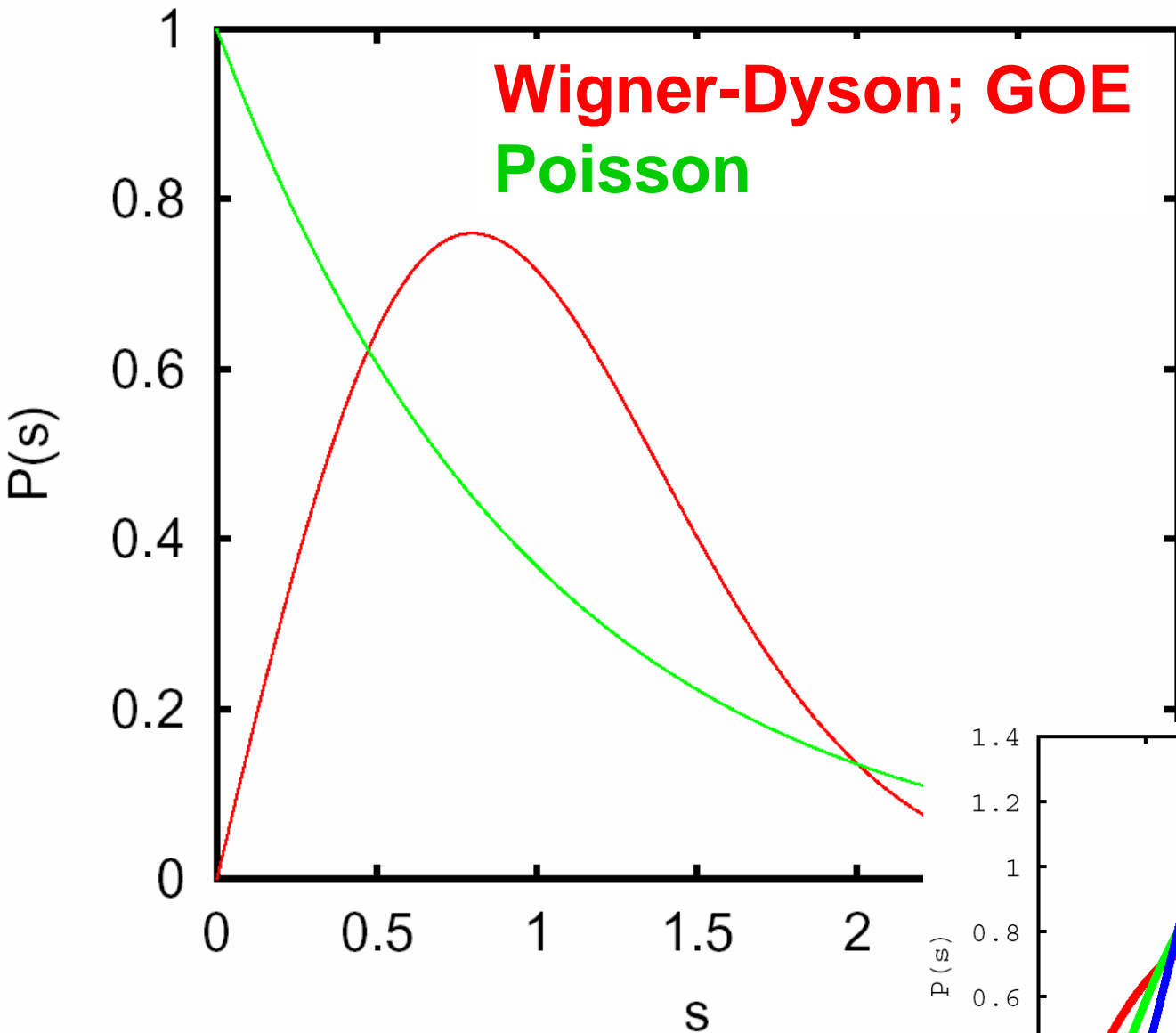
<u>Matrix elements</u>	<u>Ensemble</u>	β
real	orthogonal	1
complex	unitary	2
2×2 matrices	symplectic	4



Wigner-Dyson; GOE
Poisson

**Gaussian
Orthogonal
Ensemble**

Poisson – completely
uncorrelated
levels



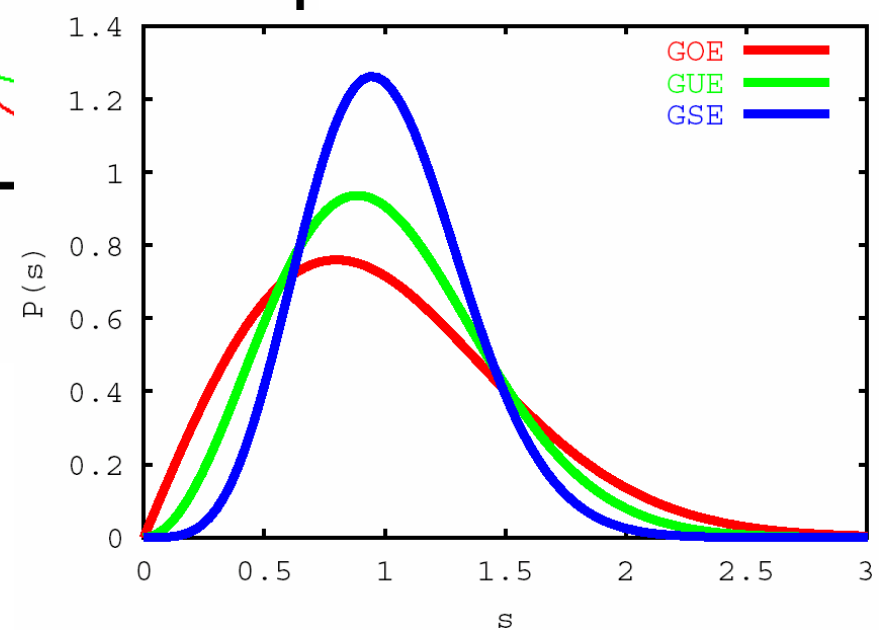
Gaussian
Orthogonal
Ensemble

Orthogonal
 $\beta=1$

Unitary
 $\beta=2$

Symplectic
 $\beta=4$

Poisson – completely uncorrelated levels



RANDOM MATRICES

$N \times N$ matrices with random matrix elements. $N \rightarrow \infty$

Dyson Ensembles

<u>Matrix elements</u>	<u>Ensemble</u>	<u>β</u>	<u>realization</u>
real	orthogonal	1	T-inv potential
complex	unitary	2	broken T-invariance (e.g., by magnetic field)
2×2 matrices	symplectic	4	T-inv, but with spin-orbital coupling

Reason for $P(s) \rightarrow 0$ when $s \rightarrow 0$:

$$\hat{H} = \begin{pmatrix} H_{11} & H_{12} \\ H_{12}^* & H_{22} \end{pmatrix}$$

$$E_2 - E_1 = \sqrt{(H_{22} - H_{11})^2 + |H_{12}|^2}$$

small

small

small

1. The assumption is that the matrix elements are statistically independent. Therefore probability of two levels to be degenerate vanishes.
2. If H_{12} is **real (orthogonal ensemble)**, then for s to be small **two statistically independent** variables ($(H_{22} - H_{11})$ and H_{12}) should be small and thus $P(s) \propto s$ $\beta = 1$
3. **Complex H_{12} (unitary ensemble)** \implies both $Re(H_{12})$ and $Im(H_{12})$ are statistically independent \implies **three** independent random variables should be small $\implies P(s) \propto s^2$ $\beta = 2$

Dyson Ensembles and Hamiltonian systems

<u>Matrix elements</u>	<u>Ensemble</u>	β	<u>realization</u>
real	orthogonal	1	T-inv potential
complex	unitary	2	broken T-invariance (e.g., by magnetic field)
2×2 matrices	symplectic	4	T-inv, but with spin-orbital coupling

ATOMS

Main goal is to classify the eigenstates in terms of the quantum numbers

NUCLEI

For the nuclear excitations this program does not work

E.P. Wigner:

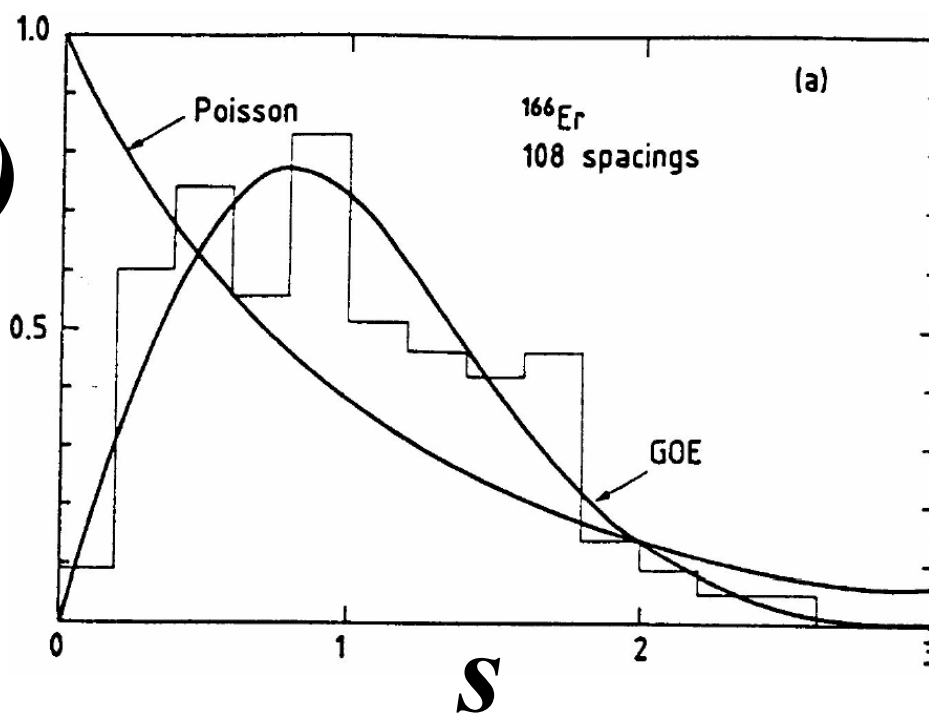
Study spectral **statistics** of a **particular** quantum system - a given nucleus

Random Matrices	Atomic Nuclei
<ul style="list-style-type: none">• <i>Ensemble</i>• <i>Ensemble averaging</i>	<ul style="list-style-type: none">• <i>Particular quantum system</i>• <i>Spectral averaging (over α)</i>

Nevertheless

Statistics of the nuclear spectra are almost exactly the same as the Random Matrix Statistics

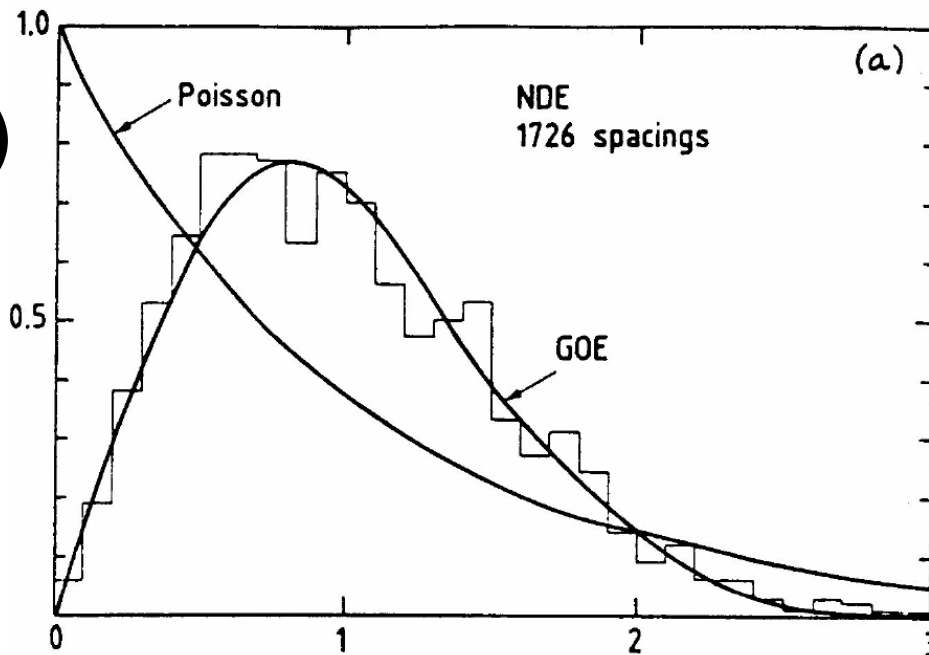
$P(s)$



Particular nucleus

^{166}Er

$P(s)$



Spectra of
several nuclei
combined (after
rescaling by the
mean level
spacing)

Q:

Why the random matrix theory (RMT) works so well for nuclear spectra

?

Original answer:

*These are systems with a **large number of degrees of freedom**, and therefore the “complexity” is high*

Later it became clear that

there exist very “simple” systems with as many as 2 degrees of freedom ($d=2$), which demonstrate RMT - like spectral statistics

Classical ($\hbar = 0$) Dynamical Systems with d degrees of freedom

Integrable Systems

The variables can be separated and the problem reduces to d one-dimensional problems



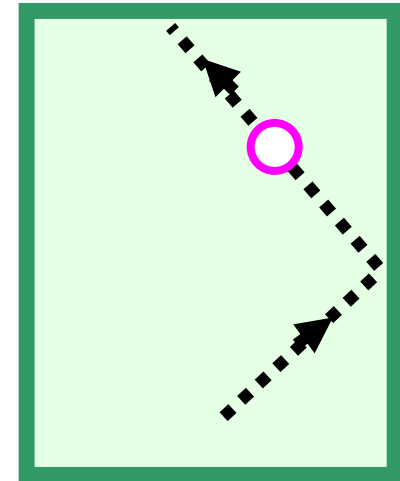
d integrals of motion

Examples

1. A ball inside rectangular billiard; $d=2$

- **Vertical** motion can be separated from the **horizontal** one

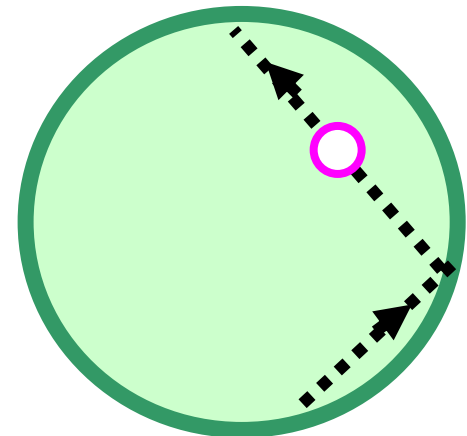
- **Vertical** and **horizontal** components of the momentum, are both integrals of motion



2. Circular billiard; $d=2$

- **Radial** motion can be separated from the **angular** one

- **Angular** momentum and **energy** are the integrals of motion



Classical Dynamical Systems with d degrees of freedom

Integrable Systems

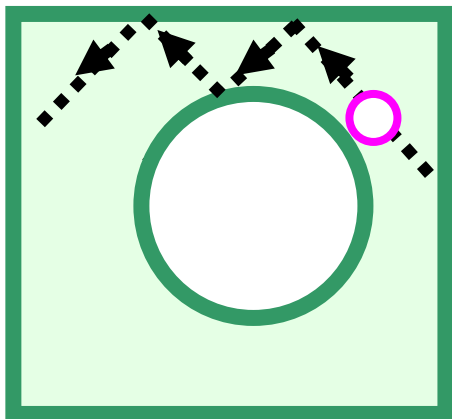
The variables can be separated \Rightarrow d one-dimensional problems \Rightarrow d integrals of motion

Rectangular and circular billiard, Kepler problem, . . . , 1d Hubbard model and other exactly solvable models, . .

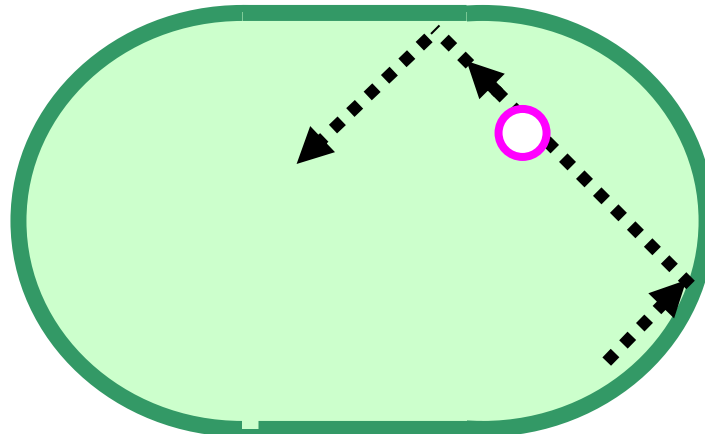
Chaotic Systems

The variables **can not** be separated \Rightarrow there is only one integral of motion - energy

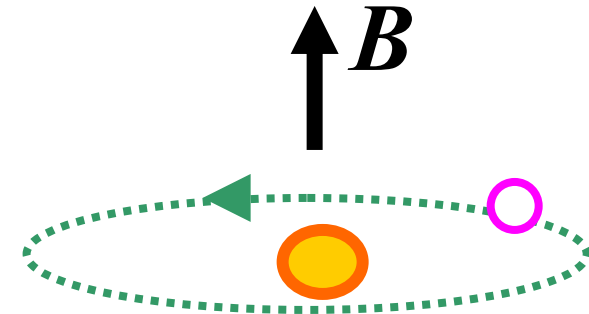
Examples



Sinai billiard



Stadium

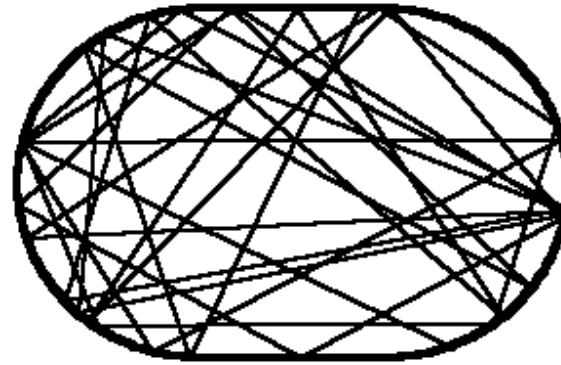
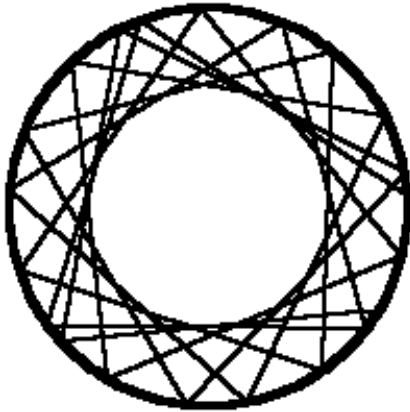


Kepler problem
in magnetic field

Classical Chaos

$$\hbar = 0$$

- *Nonlinearities*
- *Lyapunov exponents*
- *Exponential dependence on the original conditions*
- *Ergodicity*



Quantum description of any System with a finite number of the degrees of freedom is a linear problem - Shrodinger equation

Q: What does it mean Quantum Chaos ?

$\hbar \neq 0$

Bohigas – Giannoni – Schmit conjecture

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NUMBER 1

Characterization of Chaotic Quantum Spectra and Universality of Level Fluctuation Laws

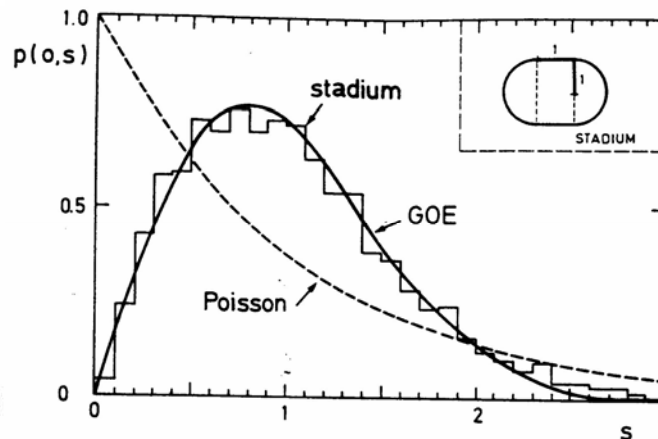
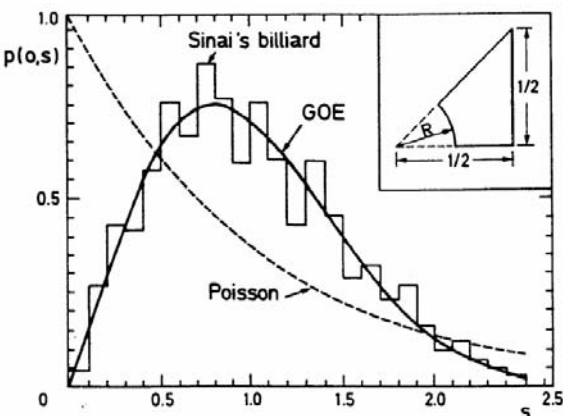
O. Bohigas, M. J. Giannoni, and C. Schmit

Division de Physique Théorique, Institut de Physique Nucléaire, F-91406 Orsay Cedex, France

(Received 2 August 1983)

It is found that the level fluctuations of the quantum Sinai's billiard are consistent with the predictions of the Gaussian orthogonal ensemble of random matrices. This reinforces the belief that level fluctuation laws are universal.

In summary, the question at issue is to prove or disprove the following conjecture: Spectra of time-reversal-invariant systems whose classical analogs are K systems show the same fluctuation properties as predicted by GOE



Chaotic classical analog



Wigner- Dyson spectral statistics



No quantum numbers except energy

Q: What does it mean **Quantum Chaos** ?

Two possible definitions

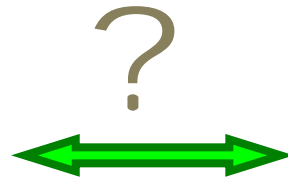
Chaotic
classical
analog

Wigner -
Dyson-like
spectrum

Classical

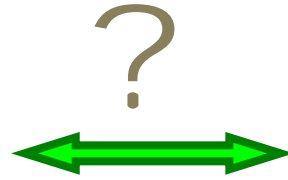
Quantum

Integrable

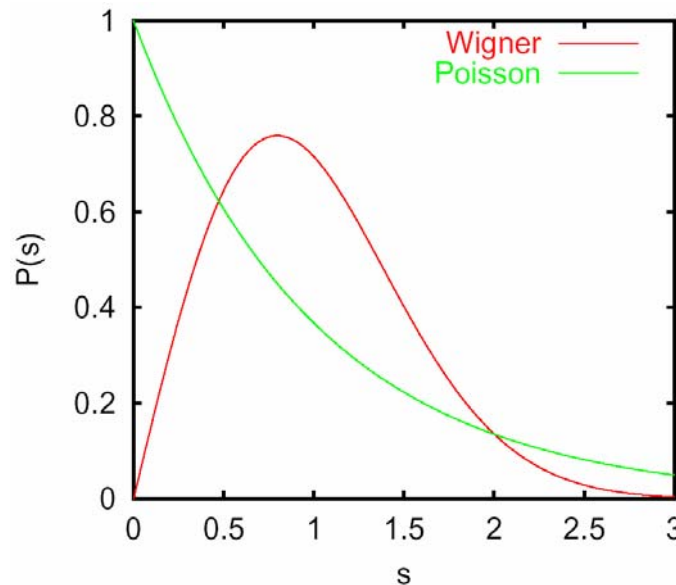


Poisson

Chaotic



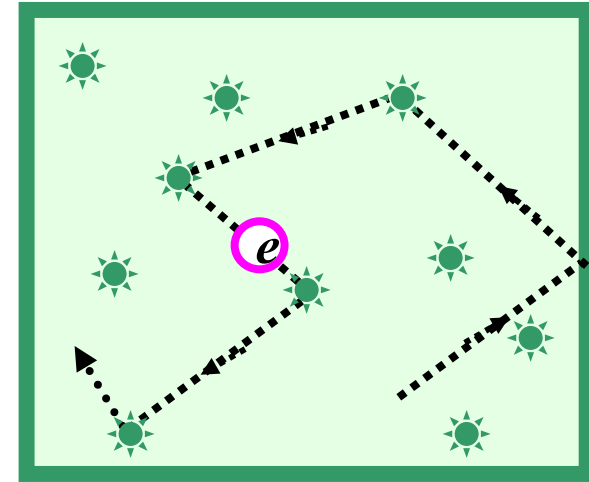
Wigner-Dyson



Poisson to Wigner-Dyson crossover

Important example: quantum particle subject to a random potential - disordered conductor

☼ *Scattering centers, e.g., impurities*



- As well as in the case of Random Matrices (RM) there is a luxury of ensemble averaging.
- The problem is much richer than RM theory
- There is still a lot of universality.

Anderson
localization (1958)

At strong enough disorder all eigenstates are **localized** in space

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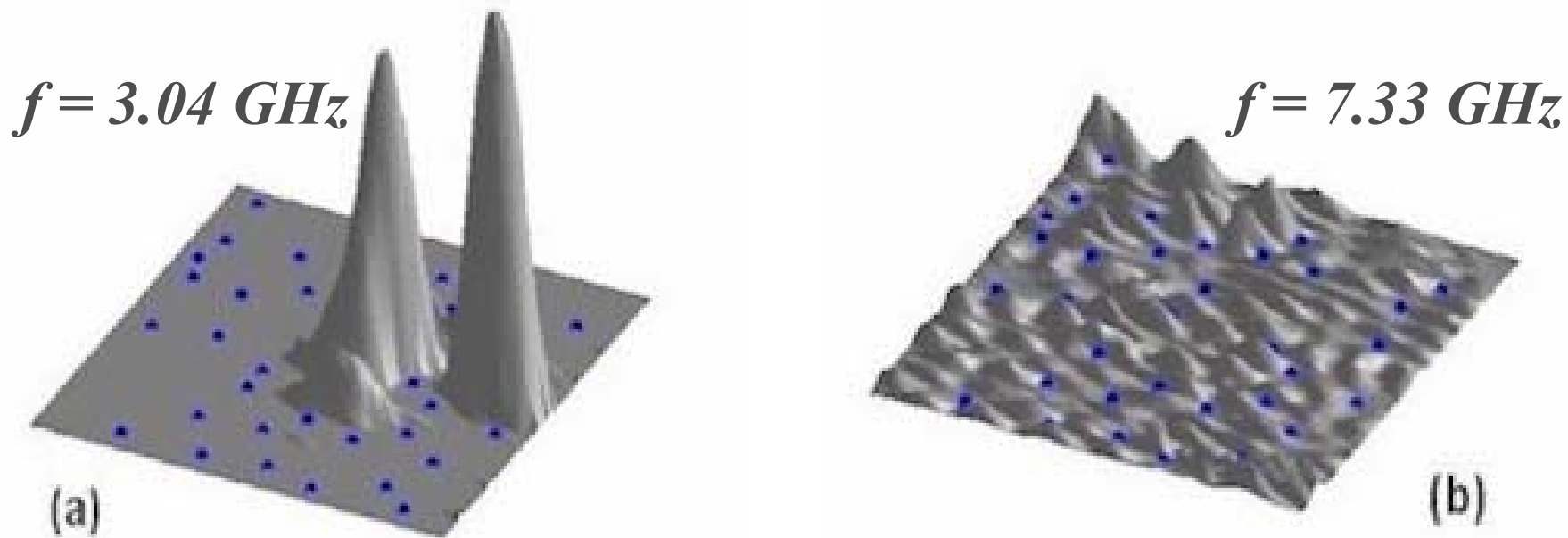
J. Bardeen, L.N. Cooper & J. Schriffer, “*Theory of Superconductivity*”; Phys.Rev., **1957**, v.108, p.1175.

Correlations due to Localization in Quantum Eigenfunctions of Disordered Microwave Cavities

Prabhakar Pradhan and S. Sridhar

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(Received 28 February 2000)



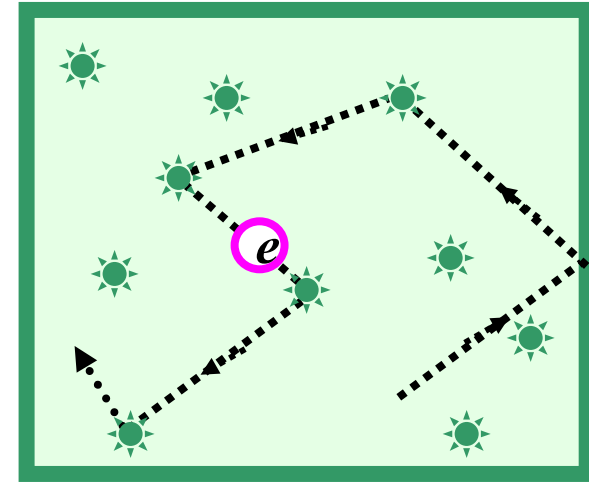
Anderson Insulator

Anderson Metal

Poisson to Wigner-Dyson crossover

Important example: quantum particle subject to a random potential - disordered conductor

☼ *Scattering centers, e.g., impurities*



Models of disorder:

Randomly located impurities

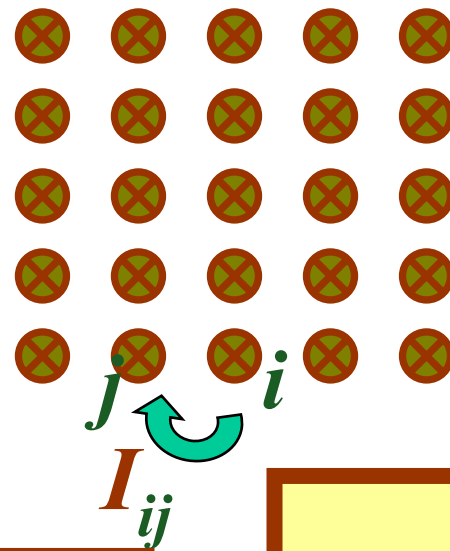
White noise potential

Lattice models

Anderson model

Lifshits model

Anderson Model



- *Lattice - tight binding model*
- *Onsite energies ϵ_i - **random***
- *Hopping matrix elements I_{ij}*

$$-W < \epsilon_i < W$$

uniformly distributed

$$I_{ij} = \begin{cases} I & \mathbf{i} \text{ and } \mathbf{j} \text{ are nearest neighbors} \\ 0 & \text{otherwise} \end{cases}$$

Anderson Transition

$$I < I_c$$

Insulator

*All eigenstates are **localized***

Localization length ξ

$$I > I_c$$

Metal

*There appear states **extended** all over the whole system*

Anderson Transition

$$I < I_c$$

Insulator

All eigenstates are localized
Localization length ξ

The eigenstates, which are localized at different places will not repel each other



Poisson spectral statistics

$$I > I_c$$

Metal

There appear states extended all over the whole system

Any two extended eigenstates repel each other

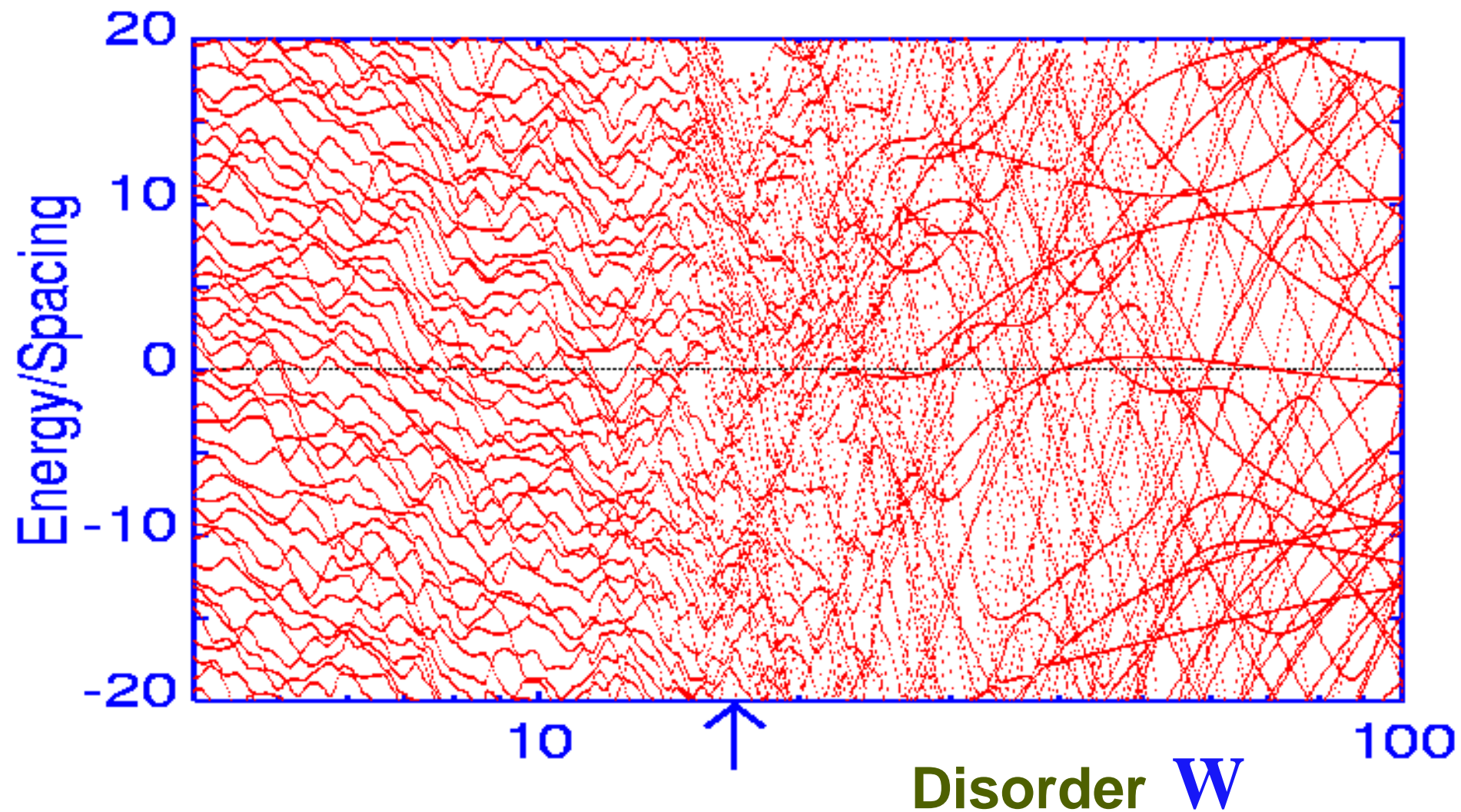


Wigner – Dyson spectral statistics

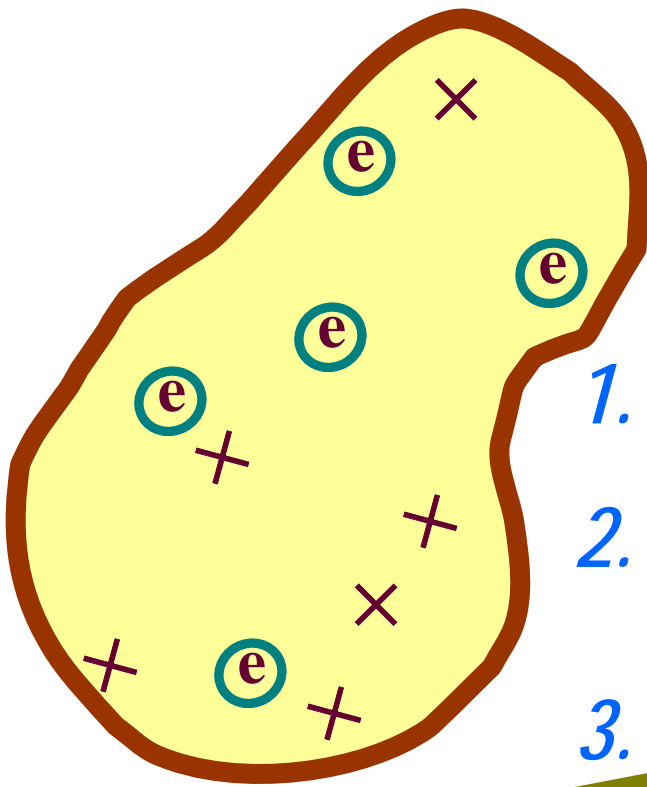
Zharekeshev & Kramer.

Exact diagonalization of the Anderson model

3D cube of volume 20x20x20



Quantum Dot



1. Disorder (x – impurities)

2. Complex geometry

3. ~~*e-e interactions*~~

for a while

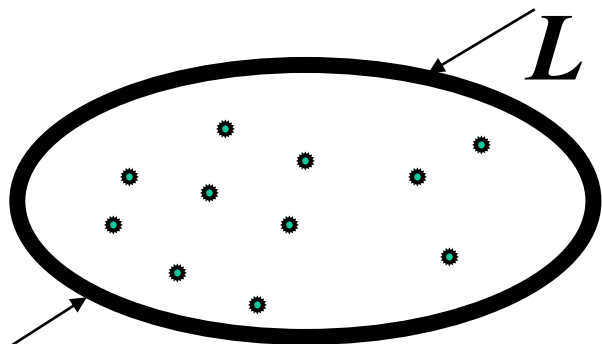
Realizations:

- Metallic clusters
- Gate determined confinement in 2D gases (e.g. *GaAs/AlGaAs*)
- Carbon nanotubes

One-particle problem (*Thouless, 1972*)

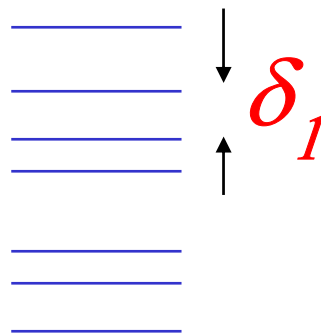
Energy scales

1. Mean level spacing



$$\delta_1 = 1/v \times L^d$$

energy



L is the system size;

d is the number of dimensions

2. Thouless energy

$$E_T = hD/L^2$$

D is the diffusion const

E_T has a meaning of the *inverse diffusion time* of the traveling through the system or the *escape rate* (for open systems)

$$g = E_T / \delta_1$$

dimensionless
Thouless
conductance

$$g = Gh/e^2$$

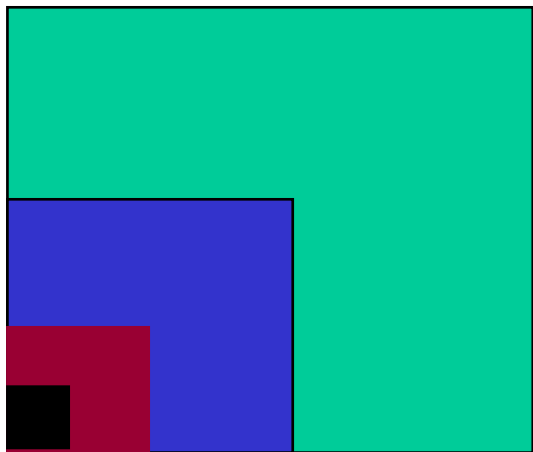
Scaling theory of Localization

(Abrahams, Anderson, Licciardello and Ramakrishnan 1979)

$$g = E_T / \delta_1$$

Dimensionless *Thouless*
conductance

$$g = Gh/e^2$$



$$L = 2L = 4L = 8L \dots$$

without quantum corrections

$$E_T \propto L^{-2} \quad \delta_1 \propto L^{-d}$$

$$E_T \longrightarrow E_T \longrightarrow E_T \longrightarrow E_T$$

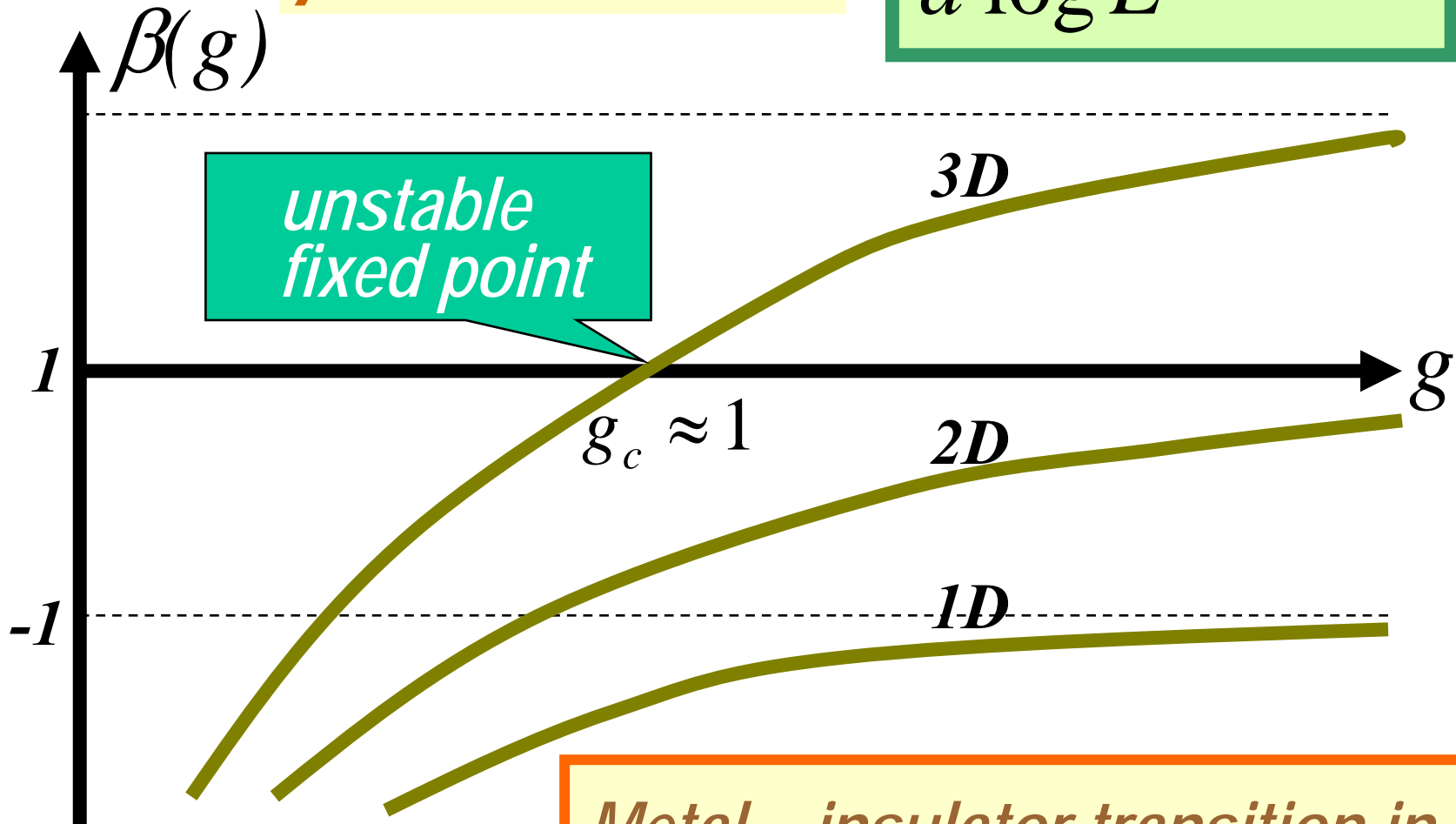
$$\delta_1 \longrightarrow \delta_1 \longrightarrow \delta_1 \longrightarrow \delta_1$$

$$g \longrightarrow g \longrightarrow g \longrightarrow g$$

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

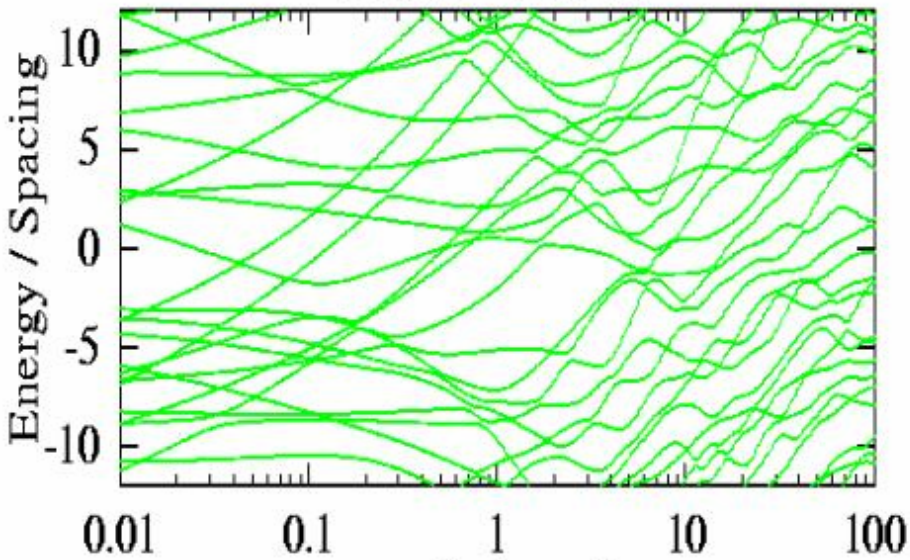
β - function

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

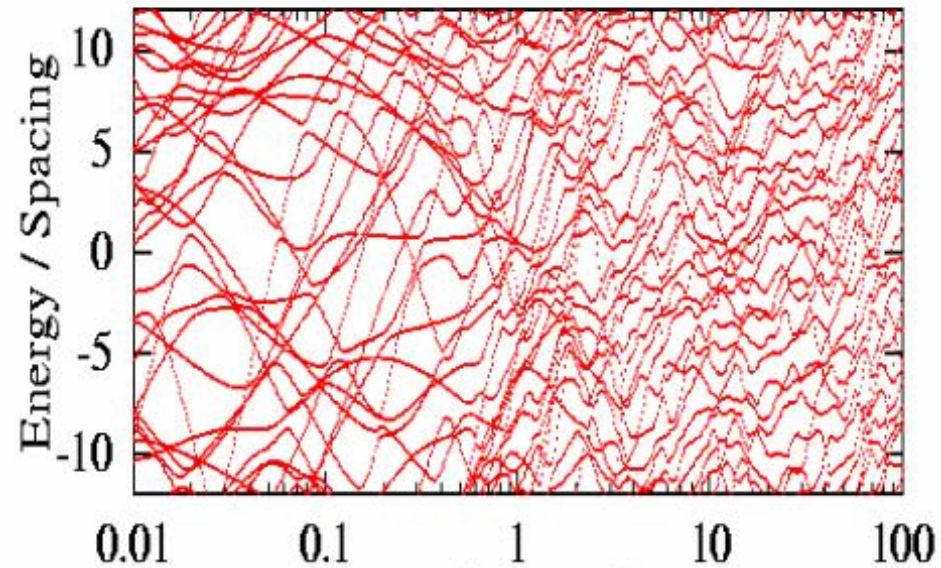


Metal – insulator transition in 3D
All states are localized for $d=1,2$

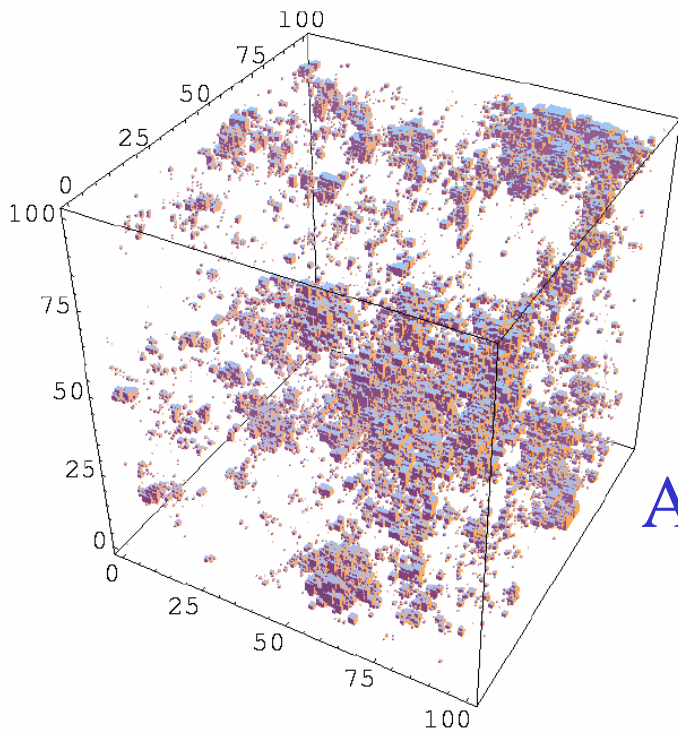
volume = $8 \times 8 \times 8$



volume = $20 \times 20 \times 20$



Critical electron eigenstate at the Anderson transition

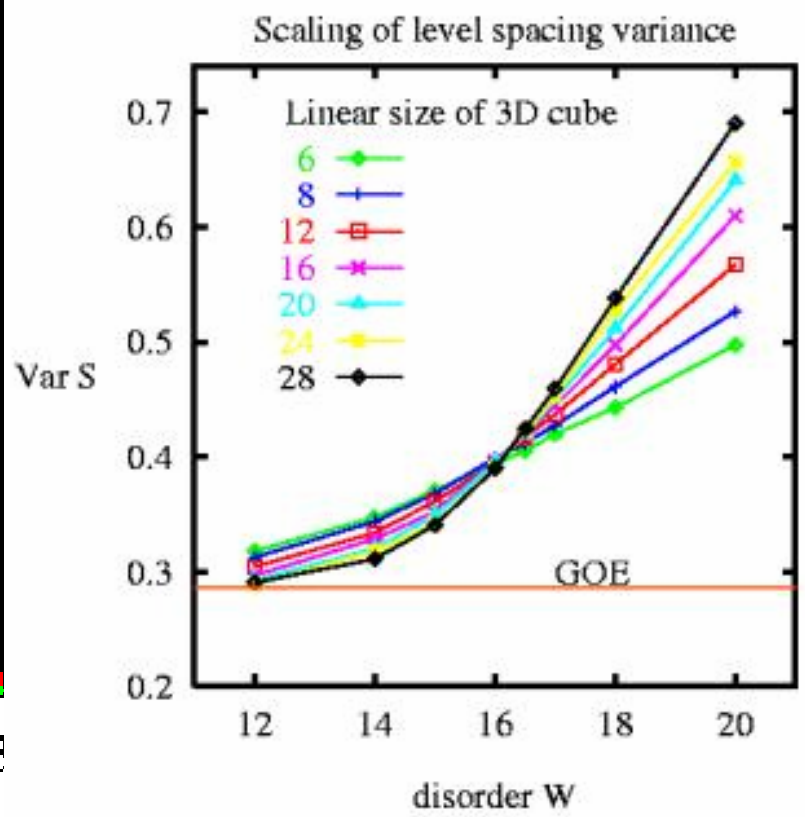
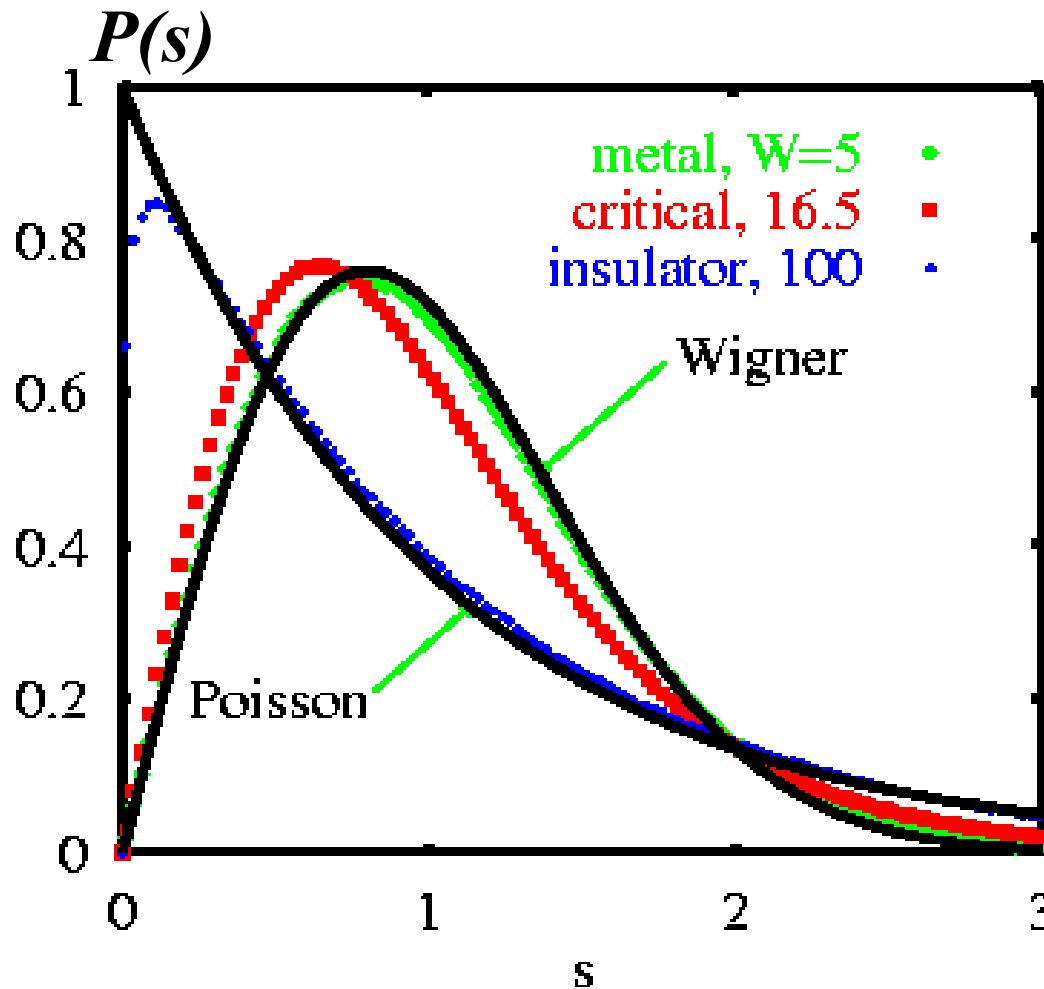


Conductance g

$100 \times 100 \times 100$

Anderson model cube

Anderson transition in terms of pure level statistics



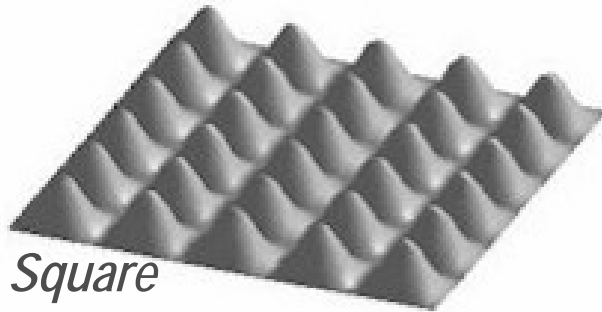
Correlations due to Localization in Quantum Eigenfunctions of Disordered Microwave Cavities

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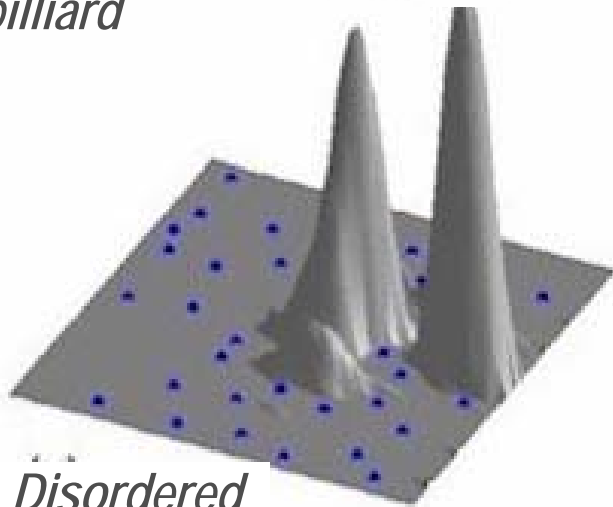
Department of Physics, Northeastern University, Boston, Massachusetts 02115

(Received 28 February 2000)

Integrable



*Square
billiard*

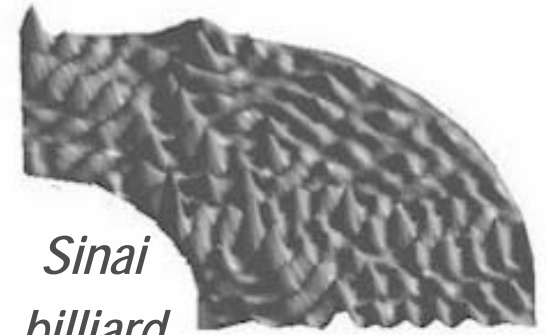


*Disordered
localized*

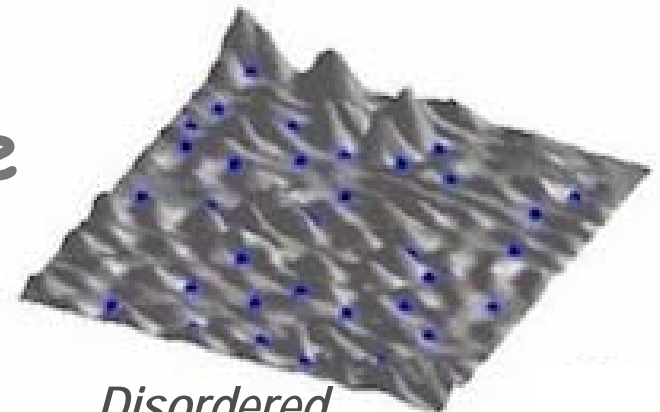
All chaotic systems resemble each other.

All integrable systems are integrable in their own way

Chaotic



*Sinai
billiard*



*Disordered
extended*

Disordered Systems:

$$E_T > \delta_1; \quad g > 1$$

*Anderson metal;
Wigner-Dyson spectral statistics*

$$E_T < \delta_1; \quad g < 1$$

*Anderson insulator;
Poisson spectral statistics*

Q: *Is it a generic scenario for the Wigner-Dyson to Poisson crossover ?*

Speculations

Consider an *integrable* system. Each state is characterized by a set of quantum numbers.

It can be viewed as a point in the space of quantum numbers. The whole set of the states forms a *lattice* in this space.

A *perturbation* that violates the integrability provides matrix elements of the *hopping* between different sites (*Anderson model* !?)

Q: *Does Anderson localization provide a generic scenario for the Wigner-Dyson to Poisson crossover ?*

Consider an *integrable* system. Each state is characterized by a *set of quantum numbers*.

It can be viewed as a point in the *space of quantum numbers*. The whole set of the states forms a *lattice* in this space.

A *perturbation* that violates the integrability provides matrix elements of the *hopping* between different sites (*Anderson model !?*)

Weak enough hopping - Localization - Poisson
Strong hopping - transition to Wigner-Dyson

The very definition of the localization is **not invariant** - one should specify in which space the eigenstates are localized.

Level statistics **is invariant**:

Poissonian
statistics

\exists basis where the
eigenfunctions are localized

Wigner -Dyson
statistics

\forall basis the eigenfunctions
are extended

Example 1

Doped semiconductor

Low concentration of donors

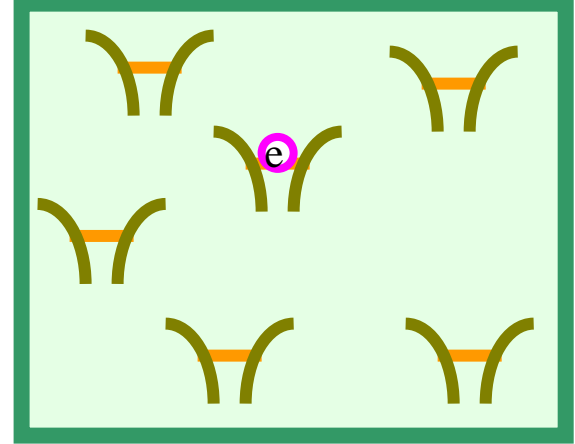


Electrons are localized on donors \Rightarrow **Poisson**

Higher donor concentration



Electronic states are extended \Rightarrow **Wigner-Dyson**

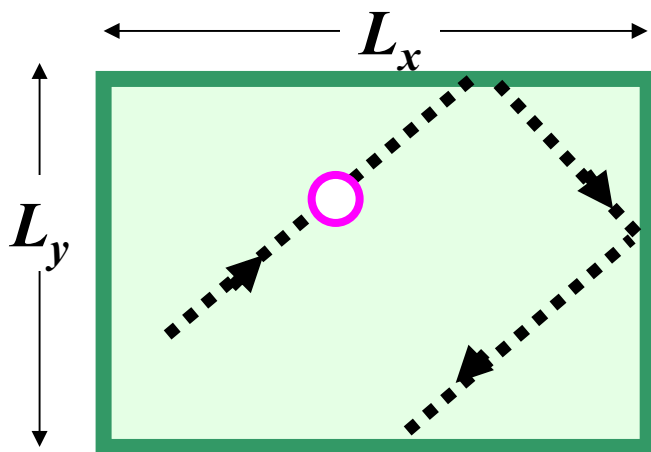


Example 2

Rectangular billiard

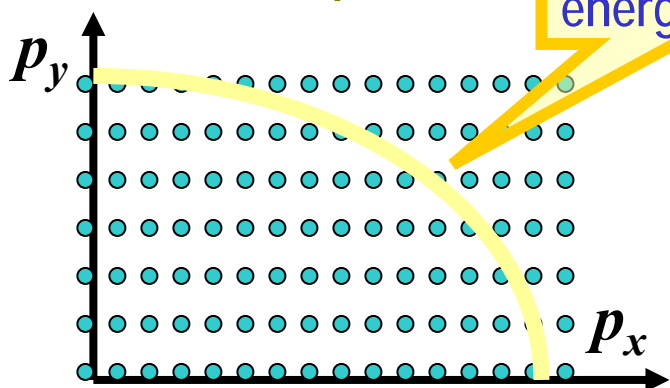
Two integrals of motion

$$p_x = \frac{\pi n}{L_x}; \quad p_y = \frac{\pi m}{L_x}$$



Lattice in the momentum space

Line (surface) of constant energy



Ideal billiard

- localization in the momentum space \Rightarrow **Poisson**

Deformation or smooth random potential

- delocalization in the momentum space \Rightarrow **Wigner-Dyson**

Localization and diffusion in the angular momentum space

Diffusion and Localization in Chaotic Billiards

Fausto Borgonovi,^{1,3,4} Giulio Casati,^{2,3,5} and Baowen Li^{6,7}

¹Dipartimento di Matematica, Università Cattolica, via Trieste 17, 25121 Brescia, Italy

²Università di Milano, sede di Como, Via Lucini 3, Como, Italy

³Istituto Nazionale di Fisica della Materia, Unità di Milano, via Celoria 16, 22100, Milano, Italy

⁴Istituto Nazionale di Fisica Nucleare, Sezione di Pavia, Pavia, Italy

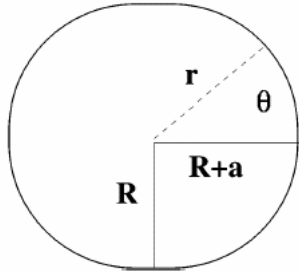
⁵Istituto Nazionale di Fisica Nucleare, Sezione di Milano, Milano, Italy

⁶Department of Physics and Centre for Nonlinear and Complex Systems, Hong Kong Baptist University, Hong Kong

⁷Center for Applied Mathematics and Theoretical Physics, University of Maribor, Krekova 2, 2000 Maribor, Slovenia

(Received 29 July 1996)

$$\varepsilon \equiv \frac{a}{R}$$



$\varepsilon > 0$ Chaotic stadium

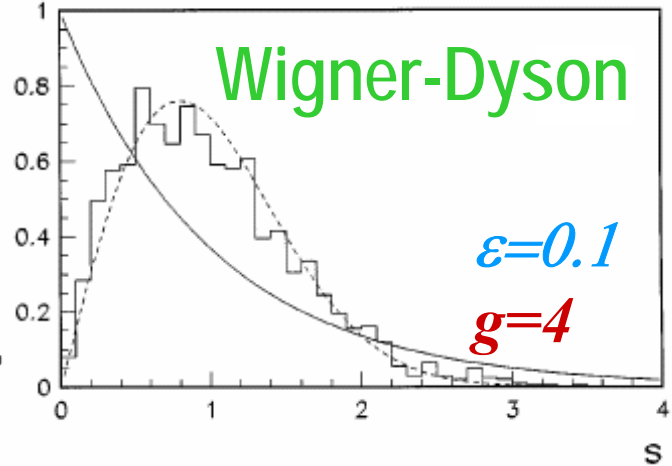
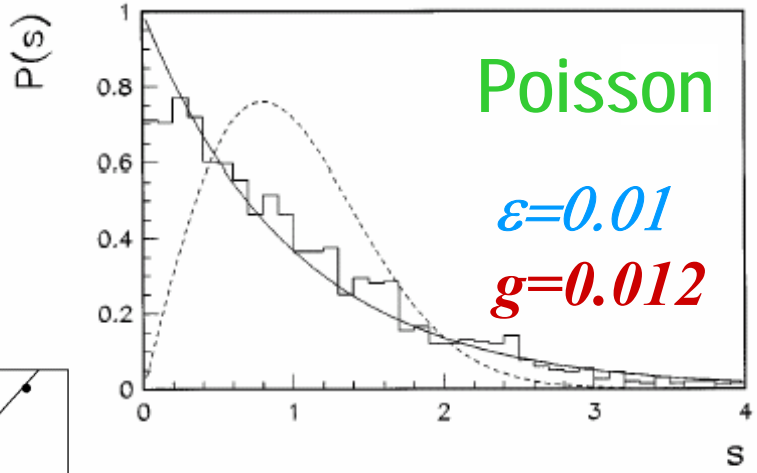
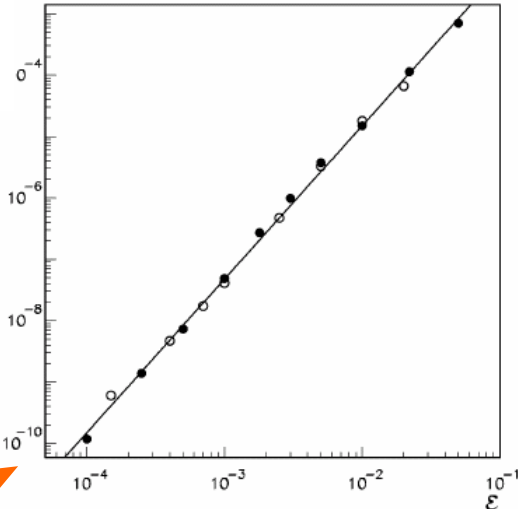
$\varepsilon \rightarrow 0$ Integrable circular billiard

Angular momentum is the integral of motion

$$\hbar = 0; \quad \varepsilon \ll 1$$

Diffusion in the angular momentum space

$$D \propto \varepsilon^{5/2}$$



1D Hubbard Model on a periodic chain

$$H = t \sum_{i,\sigma} \left(c_{i,\sigma}^+ c_{i+1,\sigma} + c_{i+1,\sigma}^+ c_{i,\sigma} \right) + U \sum_{i,\sigma} n_{i,\sigma} n_{i,-\sigma} + V \sum_{i,\sigma,\sigma'} n_{i,\sigma} n_{i+1,\sigma'}$$

$V = 0$ Hubbard model

integrable

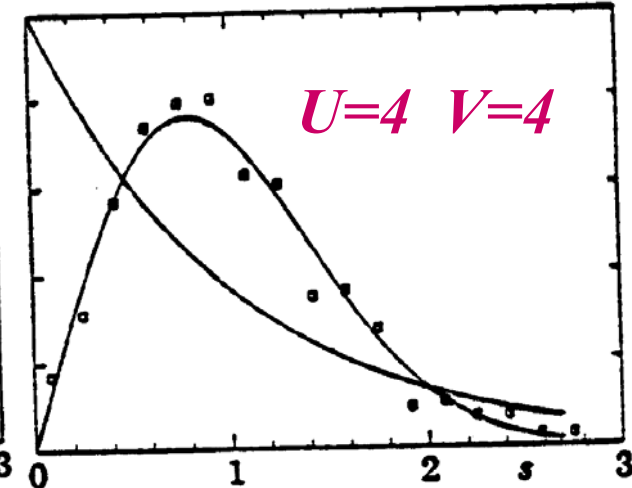
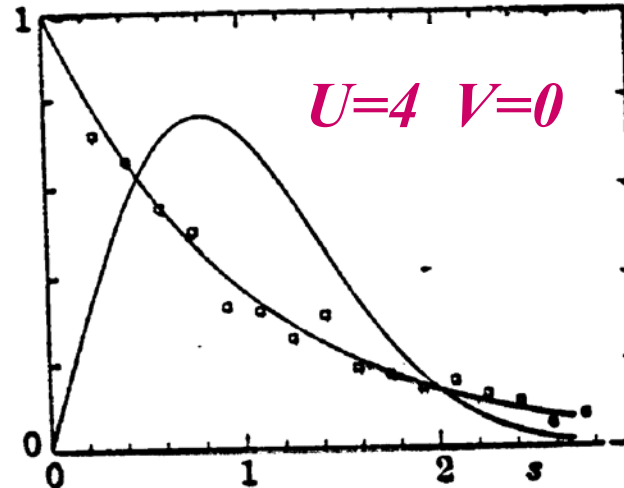
Onsite interaction

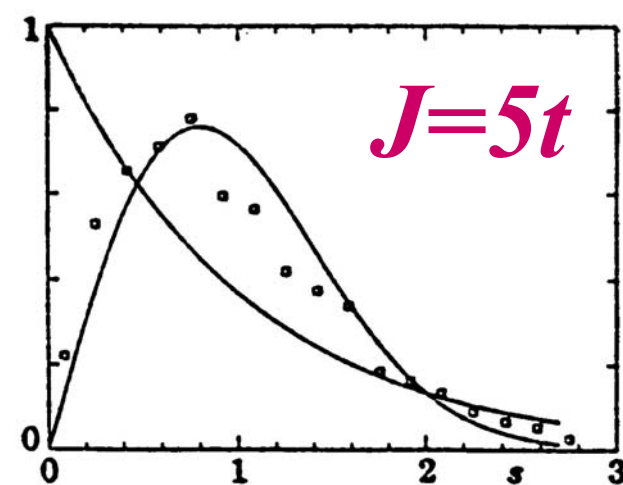
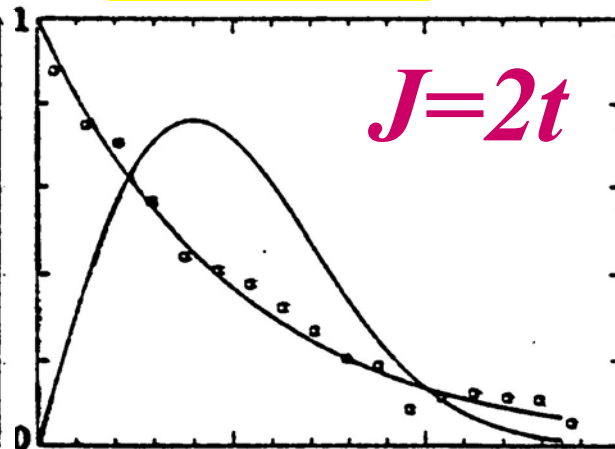
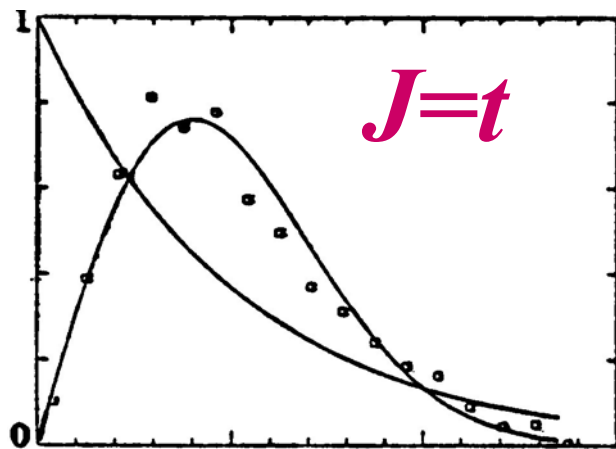
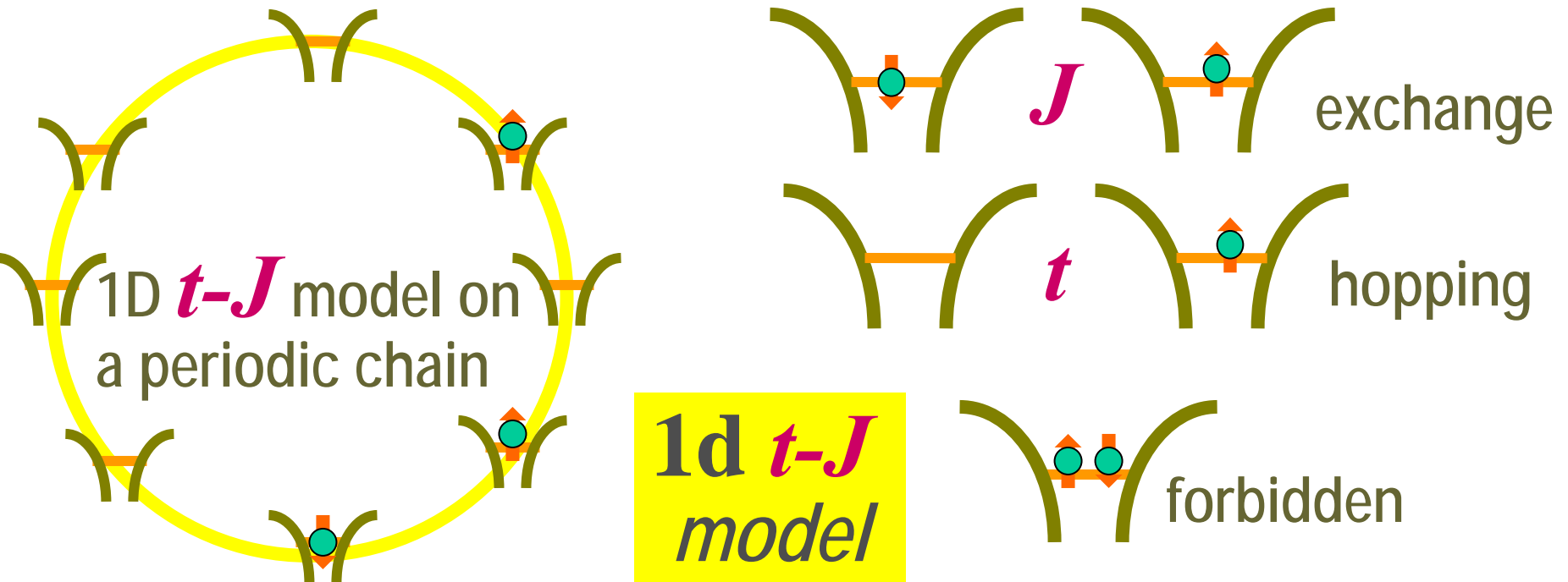
n. neighbors interaction

$V \neq 0$ extended Hubbard model

nonintegrable

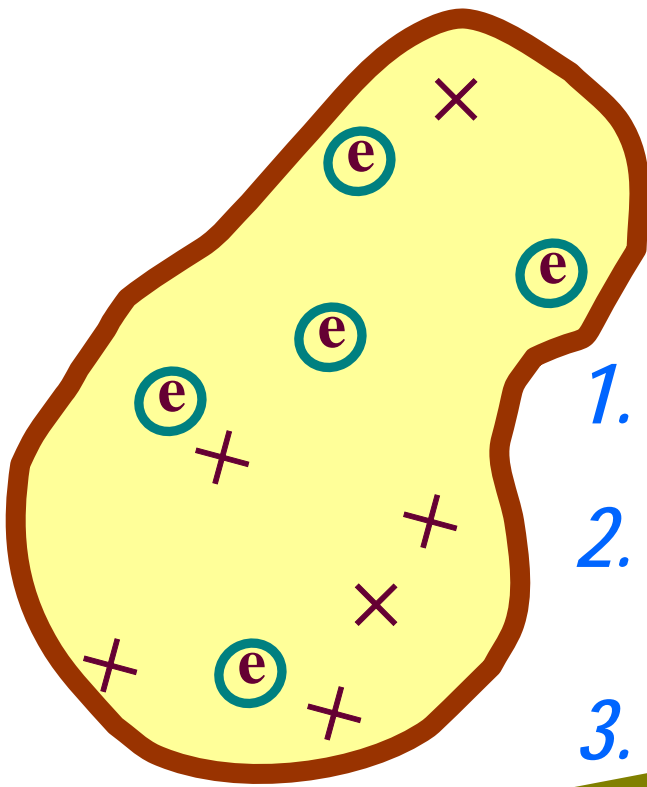
12 sites
 3 particles
 Zero total spin
 Total momentum $\pi/6$





$N=16$; one hole

Quantum Dot



1. Disorder (x – impurities)

2. Complex geometry

3. ~~*e-e interactions*~~

for a while

Realizations:

- Metallic clusters
- Gate determined confinement in 2D gases (e.g. *GaAs/AlGaAs*)
- Carbon nanotubes
-
-

Part 2 Disorder/Chaos + Interactions

Zero-Dimensional Fermi-liquid

ORIGINS

E.P. Wigner, Conference on Neutron Physics by Time of Flight, November **1956**

P.W. Anderson, “*Absence of Diffusion in Certain Random Lattices*”; Phys.Rev., **1958**, v.109, p.1492

L.D. Landau, “*Fermi-Liquid Theory*” Zh. Exp. Teor. Fiz., **1956**, v.30, p.1058

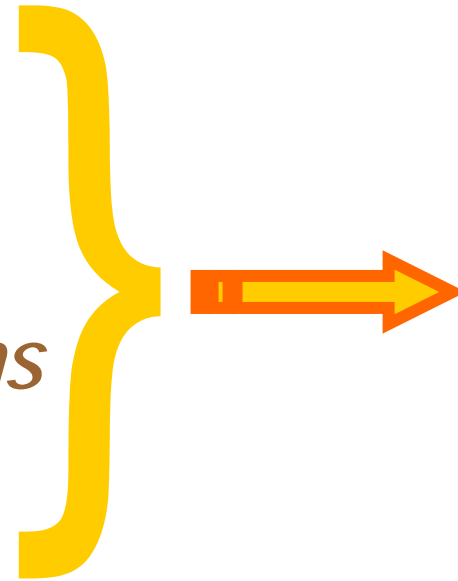
J. Bardeen, L.N. Cooper & J. Schriffer, “*Theory of Superconductivity*”; Phys.Rev., **1957**, v.108, p.1175.

What does it mean - **non-Fermi liquid** ?

What does it mean **Fermi liquid** ?

Fermi Liquid

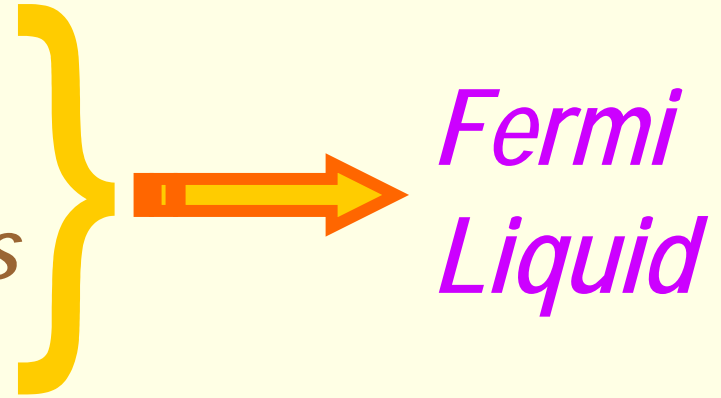
- *Fermi statistics*
- *Low temperatures*
- *Not too strong interactions*
- *Translation invariance*



*Fermi
Liquid*

What does it mean?

- *Fermi statistics*
- *Low temperatures*
- *Not too strong interactions*
- *Translation invariance*



It means that

1. *Excitations are similar to the excitations in a Fermi-gas:*
 - a) *the same quantum numbers – momentum, spin $\frac{1}{2}$, charge e*
 - b) *decay rate is small as compared with the excitation energy*
2. *Substantial renormalizations. For example, in a Fermi gas*

$$\partial n / \partial \mu, \quad \gamma = c / T, \quad \chi / g \mu_B$$

are all equal to the one-particle density of states. These quantities are different in a Fermi liquid

Signatures of the Fermi - Liquid state ?!

1. Resistivity is proportional to T^2 :

L.D. Landau & I.Ya. Pomeranchuk “*To the properties of metals at very low temperatures*”; Zh.Exp.Teor.Fiz., 1936, v.10, p.649

...The increase of the resistance caused by the interaction between the electrons is proportional to T^2 and at low temperatures exceeds the usual resistance, which is proportional to T^5 .

... the sum of the momenta of the interaction electrons can change by an integer number of the periods of the reciprocal lattice. Therefore the momentum increase caused by the electric field can be destroyed by the interaction between the electrons, not only by the thermal oscillations of the lattice.

Signatures of the Fermi - Liquid state ?!

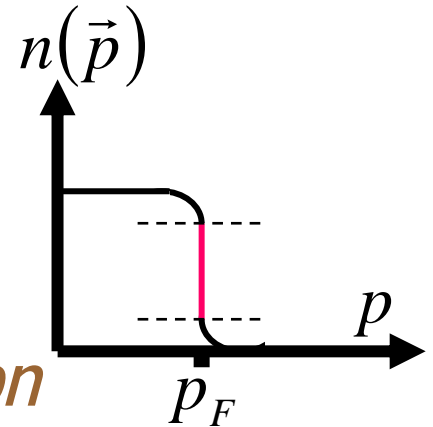
1. Resistivity is proportional to T^2 :

L.D. Landau & I.Ya. Pomeranchuk “To the properties of metals at very low temperatures”; Zh.Exp.Teor.Fiz., 1936, v.10, p.649

Umklapp electron – electron scattering dominates the charge transport (?!)

2. Jump in the momentum distribution function at $T=0$.

2a. Pole in the one-particle Green function



$$G(\varepsilon, \vec{p}) = \frac{Z}{i\varepsilon_n - \xi(\vec{p})}$$

Fermi liquid = $0 < Z < 1$ (?!)

Landau Fermi - Liquid theory

Momentum

$$\vec{p}$$

Momentum distribution

$$n(\vec{p})$$

Total energy

$$E\{n(\vec{p})\}$$

Quasiparticle energy

$$\xi(\vec{p}) \equiv \delta E / \delta n(\vec{p})$$

Landau f-function

$$f(\vec{p}, \vec{p}') \equiv \delta \xi(\vec{p}) / \delta n(\vec{p}')$$

Q:

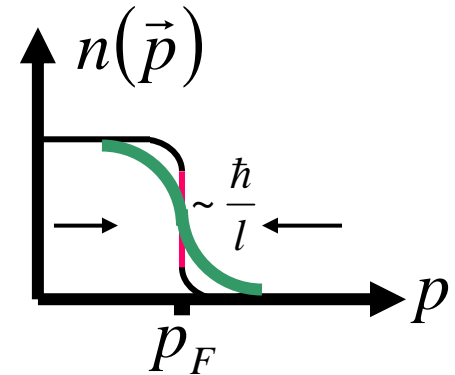
*Can Fermi – liquid survive without the **momenta***

*Does it make sense to speak about the **Fermi – liquid state** in the presence of a **quenched disorder***

?

Q: Does it make sense to speak about the *Fermi-liquid* state in the presence of a *quenched disorder* ?

1. Momentum *is not* a good quantum number – the momentum uncertainty is inverse proportional to the *elastic mean free path*, l . The step in the momentum distribution function is broadened by this uncertainty



2. Neither resistivity nor its temperature dependence is determined by the *umklapp processes* and thus does not behave as T^2

3. Sometimes (e.g., for random quenched magnetic field) the disorder averaged one-particle Green function even without interactions *does not have a pole* as a function of the energy, ϵ . The residue, Z , makes no sense.

Nevertheless even in the presence of the disorder

I. Excitations are *similar* to the excitations in a disordered *Fermi-gas*.

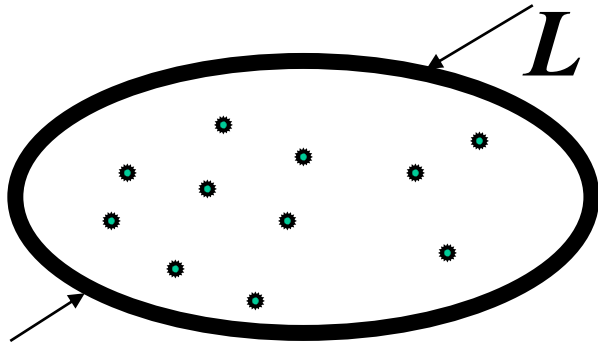
II. Small decay rate

III. Substantial renormalizations

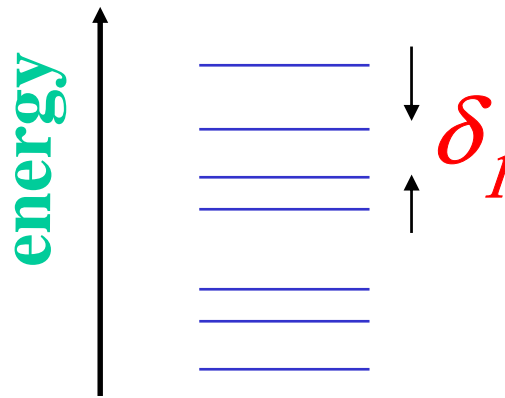
One-particle problem (*Thouless, 1972*)

Energy scales

1. Mean level spacing



$$\delta_1 = 1/v \times L^d$$



L is the system size;

d is the number of dimensions

2. Thouless energy

$$E_T = hD/L^2$$

D is the diffusion const

E_T has a meaning of the *inverse diffusion time* of the traveling through the system or the *escape rate* (for open systems)

$$g = E_T / \delta_1$$

dimensionless
Thouless
conductance

$$g = Gh/e^2$$

Zero Dimensional Fermi Liquid

Finite
System



Thouless
energy E_T

$$\varepsilon \ll E_T \xrightarrow{\text{def}} \text{0D}$$

At the same time, we want the typical energies, ε , to exceed the mean level spacing, δ_1 :

$$\delta_1 \ll \varepsilon \ll E_T$$

$$g \equiv \frac{E_T}{\delta_1} \gg 1$$

$N \times N$
Random Matrices

*Quantum Dots with
dimensionless
conductance g*

$N \rightarrow \infty$

*The same statistics of the
random spectra and one-
particle wave functions
(eigenvectors)*

$g \rightarrow \infty$

Two-Body Interactions

$$|\alpha, \sigma\rangle$$

Set of one particle states. σ and α label correspondingly *spin* and *orbit*.

$$\hat{H}_0 = \sum_{\alpha} \epsilon_{\alpha} a_{\alpha, \sigma}^{\dagger} a_{\alpha, \sigma}$$

$$\hat{H}_{\text{int}} = \sum_{\substack{\alpha, \beta, \gamma, \delta \\ \sigma, \sigma'}} M_{\alpha\beta\gamma\delta} a_{\alpha, \sigma}^{\dagger} a_{\beta, \sigma'}^{\dagger} a_{\gamma, \sigma} a_{\delta, \sigma'}$$

ϵ_{α} -one-particle orbital energies

$M_{\alpha\beta\gamma\delta}$ -interaction matrix elements

Nuclear
Physics

$$\epsilon_{\alpha}$$

are taken from the *shell model*

$$M_{\alpha\beta\gamma\delta}$$

are assumed to be *random*

Quantum
Dots

$$\epsilon_{\alpha}$$

RANDOM; Wigner-Dyson statistics

$$M_{\alpha\beta\gamma\delta}$$

?????????

Matrix Elements

$$\hat{H}_{\text{int}} = \sum_{\substack{\alpha, \beta, \gamma, \delta \\ \sigma, \sigma'}} M_{\alpha\beta\gamma\delta} a_{\alpha, \sigma}^+ a_{\beta, \sigma'}^+ a_{\gamma, \sigma} a_{\delta, \sigma'}$$

Matrix
Elements $M_{\alpha\beta\gamma\delta}$

Diagonal - $\alpha, \beta, \gamma, \delta$ are equal *pairwise*

$\alpha = \gamma$ and $\beta = \delta$ or $\alpha = \delta$ and $\beta = \gamma$ or $\alpha = \beta$ and $\gamma = \delta$

Offdiagonal - *otherwise*

It turns
out that

in the limit $g \rightarrow \infty$

- *Diagonal* matrix elements are *much bigger* than the *offdiagonal* ones

$$M_{\text{diagonal}} \gg M_{\text{offdiagonal}}$$

- *Diagonal* matrix elements in a particular sample do not fluctuate - *selfaveraging*

Toy model:

Short range **e-e** interactions

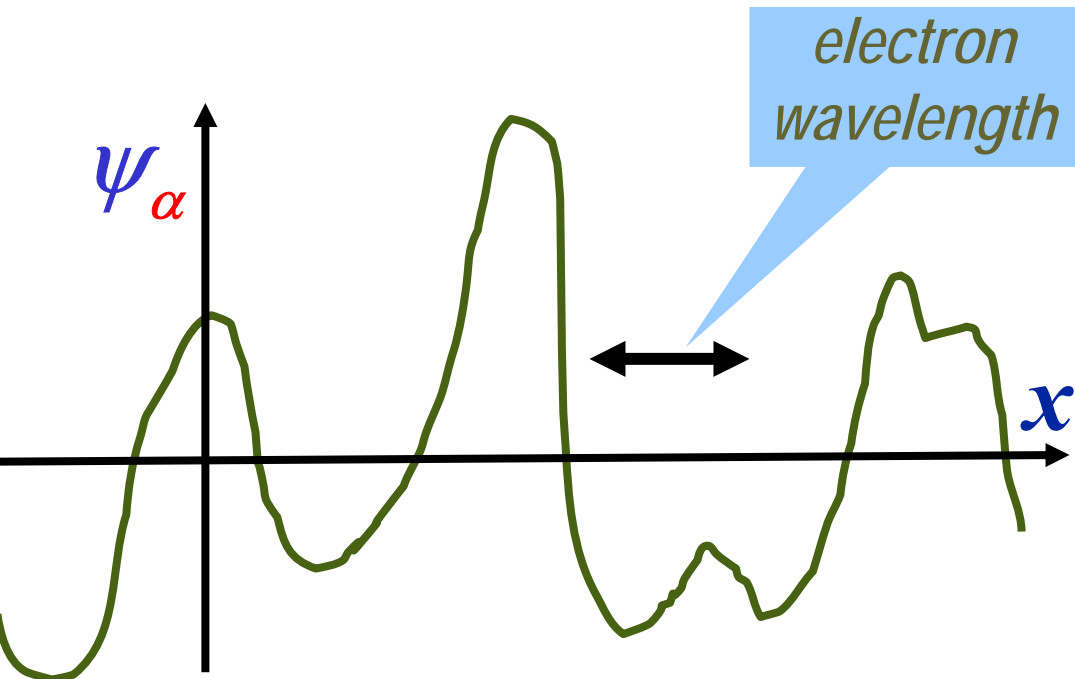
$$U(\vec{r}) = \frac{\lambda}{\nu} \delta(\vec{r})$$

λ is dimensionless coupling constant
 ν is the electron density of states

$$M_{\alpha\beta\gamma\delta} = \frac{\lambda}{\nu} \int d\vec{r} \psi_{\alpha}^*(\vec{r}) \psi_{\beta}^*(\vec{r}) \psi_{\gamma}(\vec{r}) \psi_{\delta}(\vec{r})$$

$$\psi_{\alpha}(\vec{r})$$

one-particle
eigenfunctions



$\Psi_{\alpha}(\mathbf{x})$ is a random
function that
rapidly oscillates

$$|\psi_{\alpha}(\mathbf{x})|^2 \geq 0$$

$\psi_{\alpha}(\mathbf{x})^2 \geq 0$ as long as
T-invariance
is preserved

In the limit

$$g \rightarrow \infty$$

- *Diagonal matrix elements are much bigger than the offdiagonal ones*

$$M_{\text{diagonal}} \gg M_{\text{offdiagonal}}$$

- *Diagonal matrix elements in a particular sample do not fluctuate - selfaveraging*

$$M_{\alpha\beta\alpha\beta} = \frac{\lambda}{v} \int d\vec{r} |\psi_{\alpha}(\vec{r})|^2 |\psi_{\beta}(\vec{r})|^2$$

$$|\psi_{\alpha}(\vec{r})|^2 \Rightarrow \frac{1}{\text{volume}}$$

$$M_{\alpha\beta\alpha\beta} = \lambda \delta_1$$

More general: *finite range interaction potential* $U(\vec{r})$

$$M_{\alpha\beta\alpha\beta} = \frac{\lambda}{v} \int |\psi_{\alpha}(\vec{r}_1)|^2 |\psi_{\beta}(\vec{r}_2)|^2 U(\vec{r}_1 - \vec{r}_2) d\vec{r}_1 d\vec{r}_2$$

The same conclusion

Universal (Random Matrix) limit - Random Matrix **symmetry** of the correlation functions:

All correlation functions are invariant under arbitrary orthogonal transformation:

$$\tilde{\psi}_\mu(\vec{r}) = \sum_\nu \int d\vec{r}_1 O_\mu^\nu(\vec{r}, \vec{r}_1) \psi_\nu(\vec{r}_1)$$

$$\int d\vec{r}_1 O_\mu^\nu(\vec{r}, \vec{r}_1) O_\nu^\eta(\vec{r}_1, \vec{r}') = \delta_{\mu\eta} \delta(\vec{r} - \vec{r}')$$

There are **only** three operators, which are quadratic in the fermion operators a^+ , a , and invariant under **RM** transformations:

$$\hat{n} = \sum_{\alpha, \sigma} a_{\alpha, \sigma}^+ a_{\alpha, \sigma}$$

total number of particles

$$\hat{S} = \sum_{\alpha, \sigma_1, \sigma_2} a_{\alpha, \sigma_1}^+ \vec{\sigma}_{\sigma_1, \sigma_2} a_{\alpha, \sigma_2}$$

total spin

$$\hat{T}^+ = \sum_{\alpha} a_{\alpha, \uparrow}^+ a_{\alpha, \downarrow}^+$$

????

Charge conservation
(gauge invariance) -no \hat{T} or \hat{T}^+ only $\hat{T} \hat{T}^+$

Invariance under
rotations in spin space -no \hat{S} only \hat{S}^2

Therefore, in a very general case

$$\hat{H}_{\text{int}} = eV\hat{n} + E_c\hat{n}^2 + J\hat{S}^2 + \lambda_{BCS}\hat{T}^+\hat{T}.$$

Only three coupling constants describe **all** of the effects of e-e interactions

In a very general case **only three** coupling constants describe **all** effects of electron-electron interactions:

$$\hat{H} = \sum_{\alpha} \varepsilon_{\alpha} n_{\alpha} + \hat{H}_{\text{int}}$$

$$\hat{H}_{\text{int}} = eV\hat{n} + E_c\hat{n}^2 + J\hat{S}^2 + \lambda_{\text{BCS}}\hat{T}^+\hat{T}.$$

I.L. Kurland, I.L. Aleiner & B.A., 2000

See also

P.W. Brouwer, Y. Oreg & B.I. Halperin, 1999

H. Baranger & L.I. Glazman, 1999

H-Y Kee, I.L. Aleiner & B.A., 1998

In a very general case **only three** coupling constants describe **all** effects of electron-electron interactions:

$$\hat{H} = \sum_{\alpha} \varepsilon_{\alpha} n_{\alpha} + \hat{H}_{\text{int}}$$

$$\hat{H}_{\text{int}} = eV\hat{n} + E_c\hat{n}^2 + J\hat{S}^2 + \lambda_{\text{BCS}}\hat{T}^+\hat{T}.$$

For a short range interaction with a coupling constant λ

$$E_c = \frac{\lambda\delta_1}{2} \quad J = -2\lambda\delta_1 \quad \lambda_{\text{BCS}} = \lambda\delta_1(2 - \beta)$$

where δ_1 is the one-particle mean level spacing

$$\hat{H} = \hat{H}_0 + \hat{H}_{\text{int}}$$

$$\hat{H}_0 = \sum_{\alpha} \varepsilon_{\alpha} n_{\alpha}$$

$$\hat{H}_{\text{int}} = eV\hat{n} + E_c\hat{n}^2 + J\hat{S}^2 + \lambda_{\text{BCS}}\hat{T}^+\hat{T}.$$

Only one-particle part of the Hamiltonian, \hat{H}_0 , contains randomness



$$\hat{H} = \hat{H}_0 + \hat{H}_{\text{int}}$$

$$\hat{H}_0 = \sum_{\alpha} \varepsilon_{\alpha} n_{\alpha}$$

$$\hat{H}_{\text{int}} = eV\hat{n} + E_c\hat{n}^2 + J\hat{S}^2 + \lambda_{BCS}\hat{T}^+\hat{T}.$$

E_c determines the charging energy
(Coulomb blockade)

J describes the spin exchange interaction

λ_{BCS} determines effect of superconducting-like pairing

ORIGINS

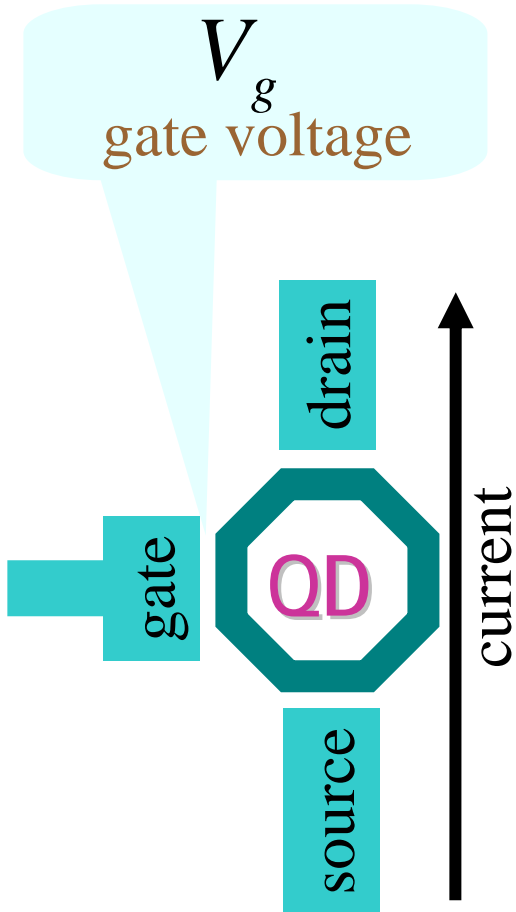
E.P. Wigner, Conference on Neutron Physics by Time of Flight, November **1956**

P.W. Anderson, “*Absence of Diffusion in Certain Random Lattices*”; Phys.Rev., **1958**, v.109, p.1492

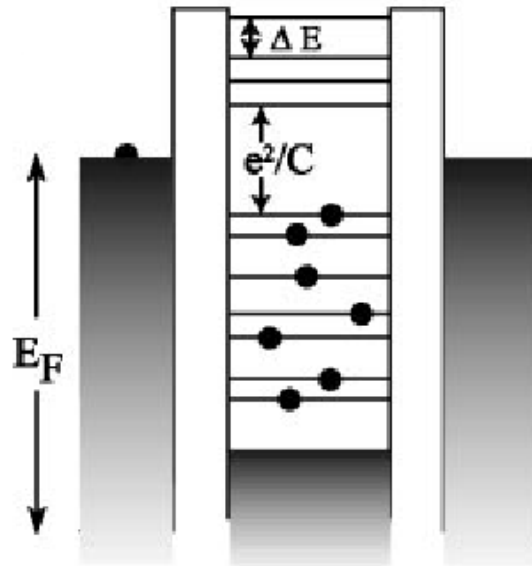
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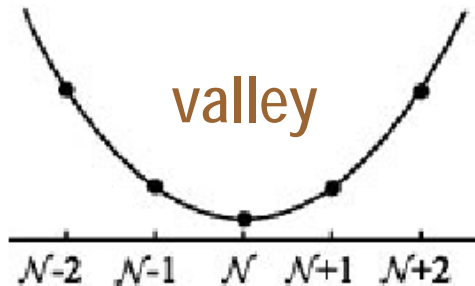
Example 1: Coulomb Blockade



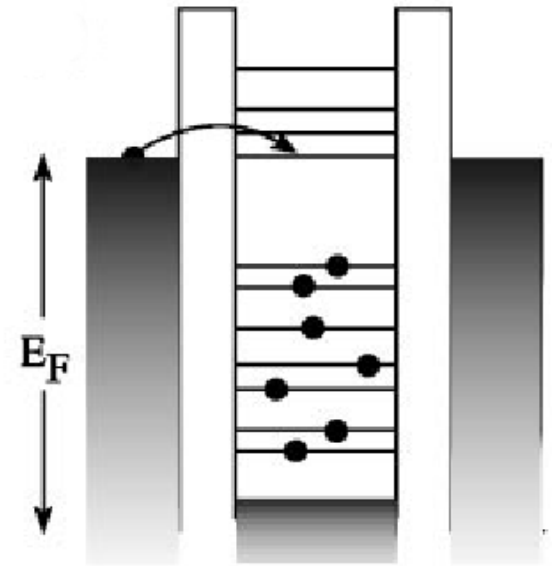
$$e \alpha V_g = N \frac{e^2}{C}$$



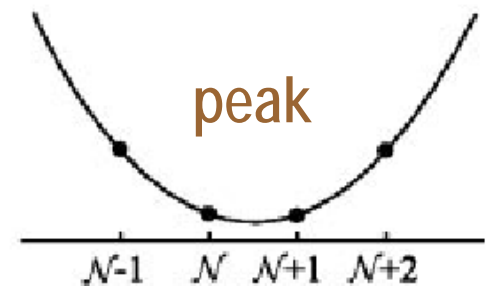
Coulomb Blockade

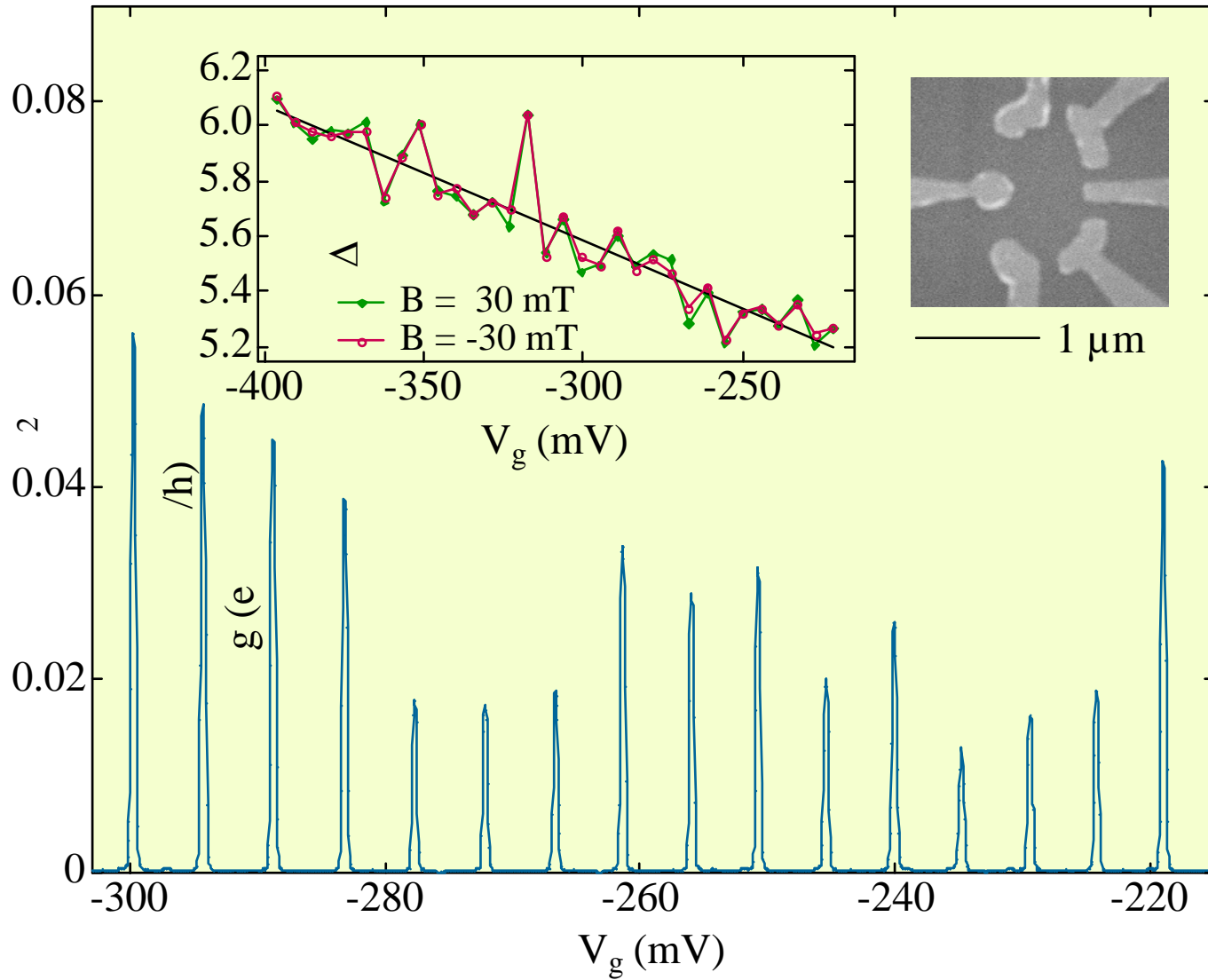


$$e \alpha V_g = (N+1/2) \frac{e^2}{C}$$



$N \rightarrow N+1$ transition





Coulomb Blockade Peak Spacing
Patel, et al. PRL 80 4522 (1998)
(Marcus Lab)

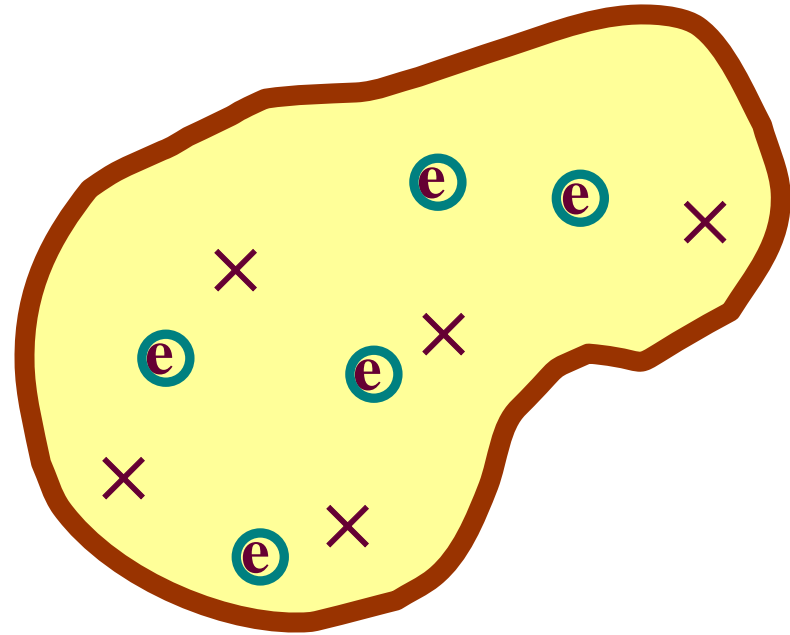
Example 2: Spontaneous Magnetization

1. Disorder
(\times impurities)

2. Complex
geometry

3. e - e interactions

chaotic
one-particle
motion

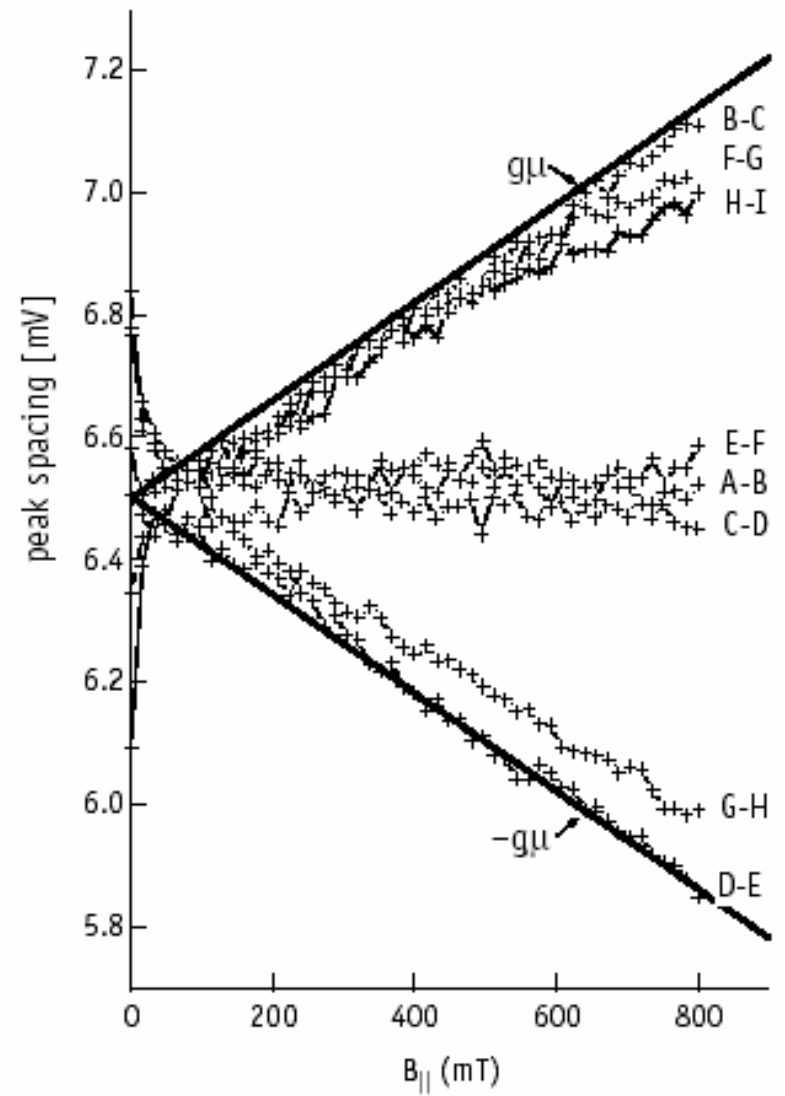
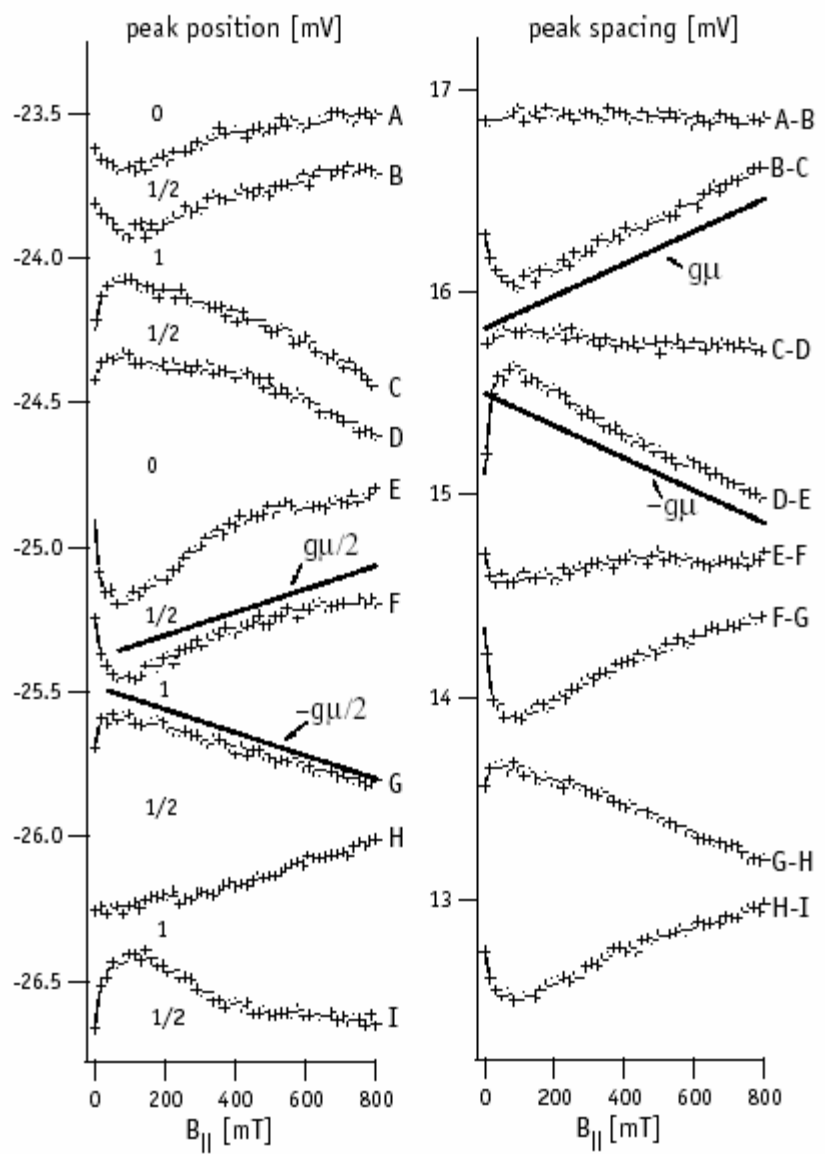


Q

- *What is the **spin** of the Quantum?*
- *Dot in the ground state*

?

How to measure the Magnetization – **motion** of the Coulomb blockade peaks in the **parallel** magnetic field



In the presence of magnetic field

$$\hat{H}_{\text{int}} = \sum_{\alpha} \varepsilon_{\alpha} n_{\alpha} + J \hat{S}^2 + \vec{B} \bullet \hat{S}$$

Scaling:

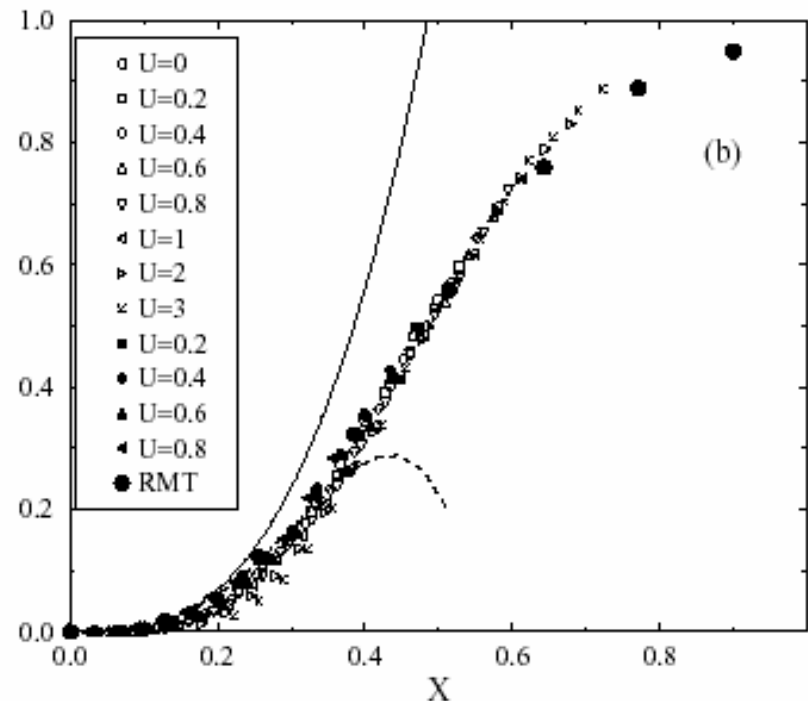
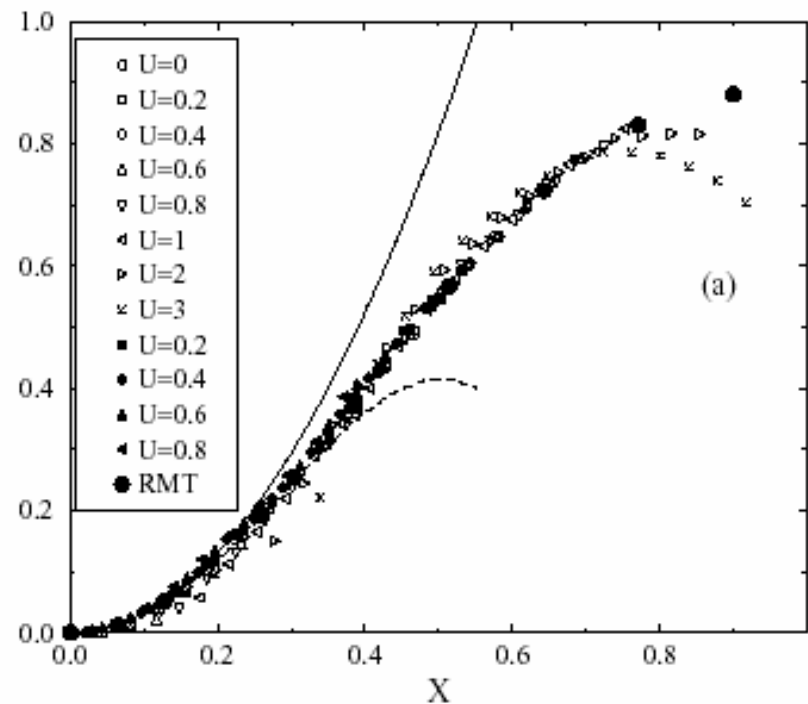
the probability to find a ground state at a given magnetic field, B , with a given spin, S , depends on the combination rather than on B and J separately

$$X = J + g\mu_B \frac{B}{2S}$$

Probability to observe a triplet state as a function of the parameter X

● - *results of the calculation based on the universal Hamiltonian with the RM one-particle states*

The rest – exact diagonalization for Hubbard clusters with disorder. No adjustable parameters



$$\hat{H} = \hat{H}_0 + \hat{H}_{\text{int}}$$

$$\hat{H}_0 = \sum_{\alpha} \varepsilon_{\alpha} n_{\alpha}$$

$$\hat{H}_{\text{int}} = eV\hat{n} + E_c\hat{n}^2 + J\hat{S}^2 + \lambda_{\text{BCS}}\hat{T}^+\hat{T}.$$

- I. Excitations are *similar* to the excitations in a disordered Fermi-gas.
- II. Small decay rate
- III. Substantial renormalizations

Isn't it a Fermi liquid ?

Fermi liquid behavior follows from the fact that different wave functions are almost uncorrelated

CONCLUSIONS

Anderson localization provides a generic scenario for the transition between chaotic and integrable behavior.

One-particle chaos + moderate interaction of the electrons \mapsto to a rather simple Hamiltonian of the system, which can be called Zero-dimensional Fermi liquid.

The main parameter that justifies this description is the Thouless conductance, which is supposed to be large

Excitations are characterized by their one-particle energy, charge and spin, **but not by their momentum.**

These excitations have the lifetime, which is proportional to the Thouless conductance, i.e., is long.

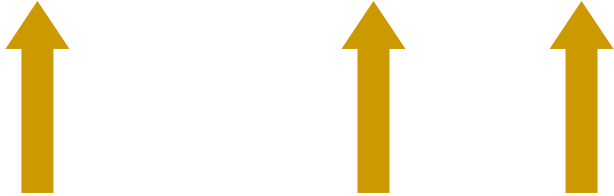
This approach allows to describe Coulomb blockade (renormalization of the compressibility), as well as the substantial renormalization of the magnetic susceptibility and effects of superconducting pairing

BCS Hamiltonian

Finite systems

$$\hat{H} = \sum_{\alpha} \varepsilon_{\alpha} n_{\alpha} + E_c \hat{n}^2 + J \hat{S}^2 + \lambda_{BCS} \hat{T}^+ \hat{T}.$$

$$\hat{T}^+ = \sum_{\alpha} a_{\alpha, \uparrow}^+ a_{\alpha, \downarrow}^+$$



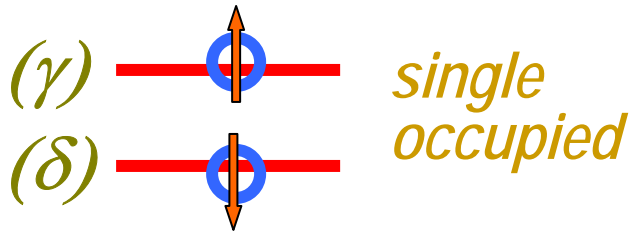
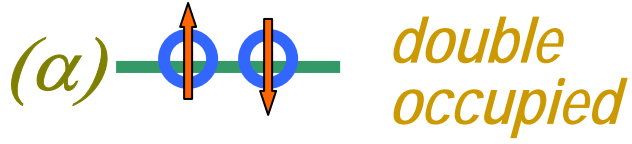
commute with
each other



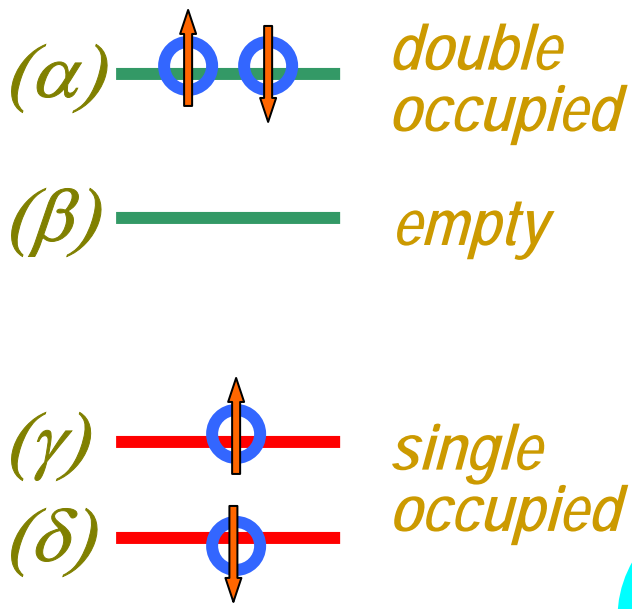
does not commute
with K.E.

$$\hat{H} = \sum_{\alpha} \varepsilon_{\alpha} n_{\alpha} + \lambda_{BCS} \hat{T}^+ \hat{T}.$$

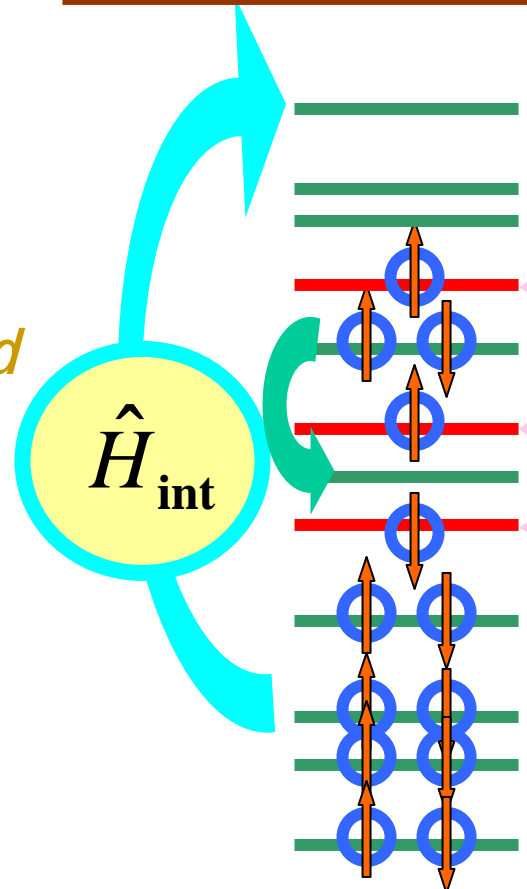
$$\hat{H} = \sum_{\alpha} \varepsilon_{\alpha} n_{\alpha} + \lambda_{BCS} \hat{T}^+ \hat{T}.$$



$$\hat{H} = \sum_{\alpha} \varepsilon_{\alpha} n_{\alpha} + \lambda_{BCS} \hat{T}^+ \hat{T}.$$



$a_{\alpha\uparrow}^+ a_{\alpha\downarrow}^+ a_{\beta\uparrow} a_{\beta\downarrow}$ mixes (α) and (β)
 at the same time $a_{\alpha\uparrow}^+ a_{\alpha\downarrow}^+ a_{\beta\uparrow} a_{\beta\downarrow} (\gamma, \delta) = 0$



This **single-occupied** states are not effected by the interaction.
 They are **blocked**
 The Hilbert space is separated into **two independent Hilbert subspaces**

Blocking effect.
(V.G.Soloviev, 1961)

Theory of Superconductivity*

J. BARDEEN, L. N. COOPER,[†] AND J. R. SCHRIEFFER[‡]
Department of Physics, University of Illinois, Urbana, Illinois

(Received July 8, 1957)

$$H_{\text{red}} = 2 \sum_{k > k_F} \epsilon_k b_k^* b_k + 2 \sum_{k < k_F} |\epsilon_k| b_k b_k^* - \sum_{\mathbf{k}\mathbf{k}'} V_{\mathbf{k}\mathbf{k}'} b_{\mathbf{k}'}^* b_{\mathbf{k}}. \quad (2.14)$$

$$b_{\mathbf{k}} = c_{-\mathbf{k}\downarrow} c_{\mathbf{k}\uparrow},$$

$$b_{\mathbf{k}}^* = c_{\mathbf{k}\uparrow}^* c_{-\mathbf{k}\downarrow}^*$$

One-particle
energy

BCS
interaction

The energy is
measured relative
to the Fermi level

BCS Hamiltonian; no single-occupied states

n one-particle levels; one-particle energies ε_α

$$\hat{H}_{BCS} = \sum_{\substack{0 \leq \alpha \leq n-1 \\ \sigma = \uparrow, \downarrow}} \varepsilon_\alpha a_{\alpha, \sigma}^+ a_{\alpha, \sigma} - \lambda_{BCS} \sum_{\substack{0 \leq \alpha, \beta \leq n-1 \\ \alpha \neq \beta}} a_{\alpha, \uparrow}^+ a_{\alpha, \downarrow}^+ a_{\beta, \uparrow} a_{\beta, \downarrow}$$

Anderson spin chain

$$\hat{T}_\alpha^z = \frac{1}{2} \left(-1 + \sum_{\sigma = \uparrow, \downarrow} a_{\alpha, \sigma}^+ a_{\alpha, \sigma} \right) \quad \hat{T}_\alpha^+ = a_{\alpha, \uparrow}^+ a_{\alpha, \downarrow}^+ \quad \hat{T}_\alpha^- = a_{\alpha, \uparrow} a_{\alpha, \downarrow}$$

SU_2
algebra

$$\hat{H}_{BCS} = \sum_{0 \leq \alpha \leq n-1} \varepsilon_\alpha \hat{T}_\alpha^z - \lambda_{BCS} \sum_{0 \leq \alpha \neq \beta \leq n-1} \hat{T}_\alpha^+ \hat{T}_\alpha^- = \sum_{0 \leq \alpha \leq n-1} \varepsilon_p \hat{T}_p^z - \lambda_{BCS} \hat{L}_+ \hat{L}_-$$

$$\hat{L}_\pm \equiv \sum_\alpha \hat{T}_\alpha^\pm$$

Anderson spin chain

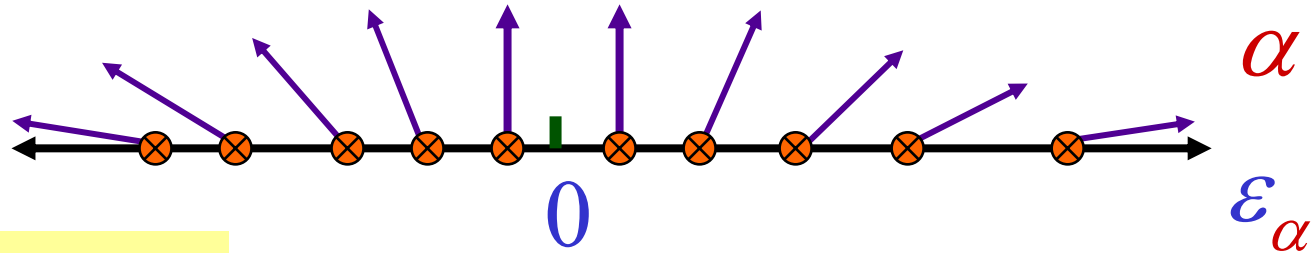
$$\hat{T}_\alpha^z = \frac{1}{2} \left(-1 + \sum_{\sigma=\uparrow,\downarrow} a_{\alpha,\sigma}^+ a_{\alpha,\sigma} \right)$$

$$\hat{T}_\alpha^+ = a_{\alpha,\uparrow}^+ a_{\alpha,\downarrow}^+$$

$$\hat{T}_\alpha^- = a_{\alpha,\uparrow} a_{\alpha,\downarrow}$$

$$\vec{\hat{L}} \equiv \sum_{\alpha} \vec{\hat{T}}_{\alpha}$$

$$\hat{H}_{BCS} = \sum_{0 \leq \alpha \leq n-1} \varepsilon_{\alpha} \hat{T}_{\alpha}^z - \lambda_{BCS} \sum_{0 \leq \alpha \neq \beta \leq n-1} \hat{T}_{\alpha}^+ \hat{T}_{\alpha}^- = \sum_{0 \leq \alpha \leq n-1} \varepsilon_p \hat{T}_p^z - \lambda_{BCS} \hat{L}_+ \hat{L}_-$$



$$\sum \hat{T}_{\alpha}^+ \equiv \frac{1}{\lambda_{BCS}} \hat{\Delta}$$

Superconducting order parameter

$$[\hat{N}, \hat{\Delta}] \neq 0$$

$$\sum \hat{T}_{\alpha}^z \equiv \hat{N}$$

Total number of the particles

Anderson spin chain

$$\hat{T}_\alpha^z = \frac{1}{2} \left(-1 + \sum_{\sigma=\uparrow,\downarrow} a_{\beta,\sigma}^+ a_{\beta,\sigma} \right) \quad \hat{T}_\alpha^+ = a_{\alpha,\uparrow}^+ a_{\alpha,\downarrow}^+ \quad \hat{T}_\alpha^- = a_{\alpha,\uparrow} a_{\alpha,\downarrow}$$

$$\hat{H}_{BCS} = \sum_{0 \leq \alpha \leq n-1} \varepsilon_\alpha \hat{T}_\alpha^z - \lambda_{BCS} \sum_{0 \leq \alpha \neq \beta \leq n-1} \hat{T}_\alpha^+ \hat{T}_\alpha^- = \sum_{0 \leq \alpha \leq n-1} \varepsilon_p \hat{T}_p^z - \lambda_{BCS} \hat{L}_+ \hat{L}_-$$

$$\sum \hat{T}_\alpha^z \equiv \hat{N}$$

Total number of the particles

For a fixed number of the particles (closed system) we can add the term $-gN^2 = \text{const}$ to the hamiltonian:

$$\hat{H}_{BCS} \sum_{0 \leq \alpha \leq n-1} \varepsilon_\alpha \hat{T}_\alpha^z - \lambda_{BCS} \left(\hat{L} \right)^2$$

$$\hat{L} \equiv \sum_{\alpha} \hat{T}_\alpha^z$$

$$\hat{H}_{BCS} = \sum_{0 \leq \alpha \leq n-1} \varepsilon_{\alpha} \hat{T}_{\alpha}^z - \lambda_{BCS} \left(\hat{L} \right)^2$$

$$\hat{L} \equiv \sum_{\alpha} \hat{T}_{\alpha}$$

Integrable model

Richardson solution - Bethe Ansatz

$$\frac{-1}{\lambda_{BCS}} + \sum_{0 \leq \beta \leq n-1} \frac{2}{E_{\alpha} - E_{\beta}} = \sum_{0 \leq \beta \leq n-1} \frac{1}{E_{\alpha} - 2\varepsilon_{\beta}}$$

$$E = \sum_{0 \leq \alpha \leq n-1} E_{\alpha}$$

n equations for the parameters E_p

How to describe dynamics in the time domain?

Gaudin Magnets

$$\hat{H}_\alpha = 2 \sum_{\substack{0 \leq \beta \leq n-1 \\ \beta \neq \alpha}} \frac{\hat{T}_\alpha \hat{T}_\beta}{\varepsilon_\alpha - \varepsilon_\beta} - A \hat{T}_\alpha^z \quad \alpha = 0, 1, 2, \dots, n-1$$

$$[\hat{H}_\alpha, \hat{H}_\beta] = 0$$

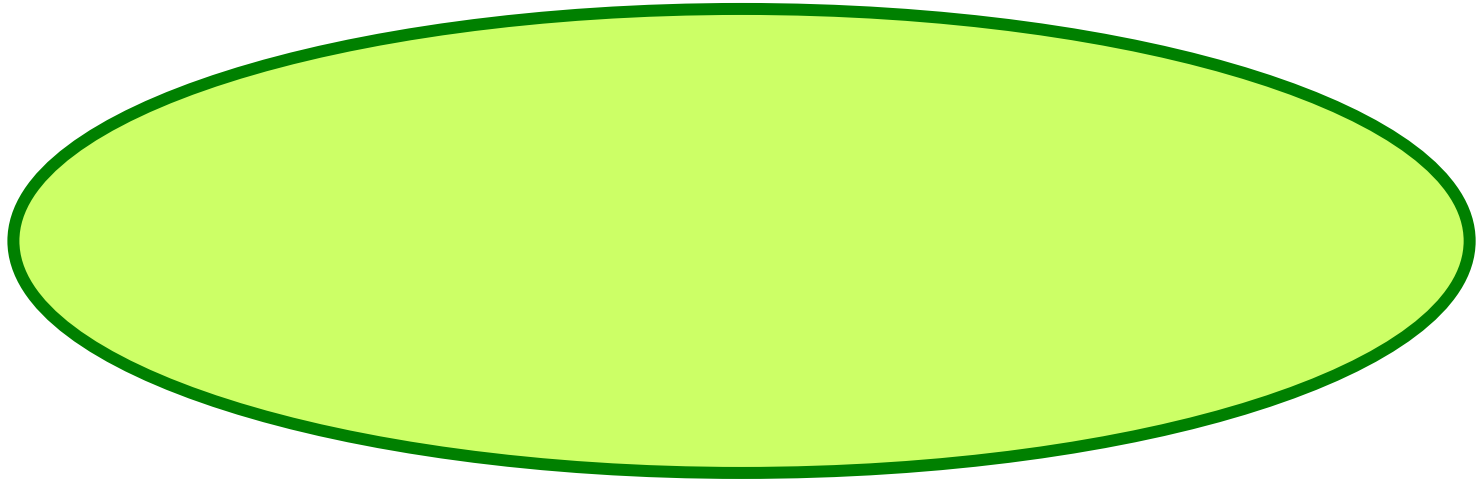
\hat{H}_0 – Hamiltonian

$\hat{H}_{\beta>0}$ – integrals of motion

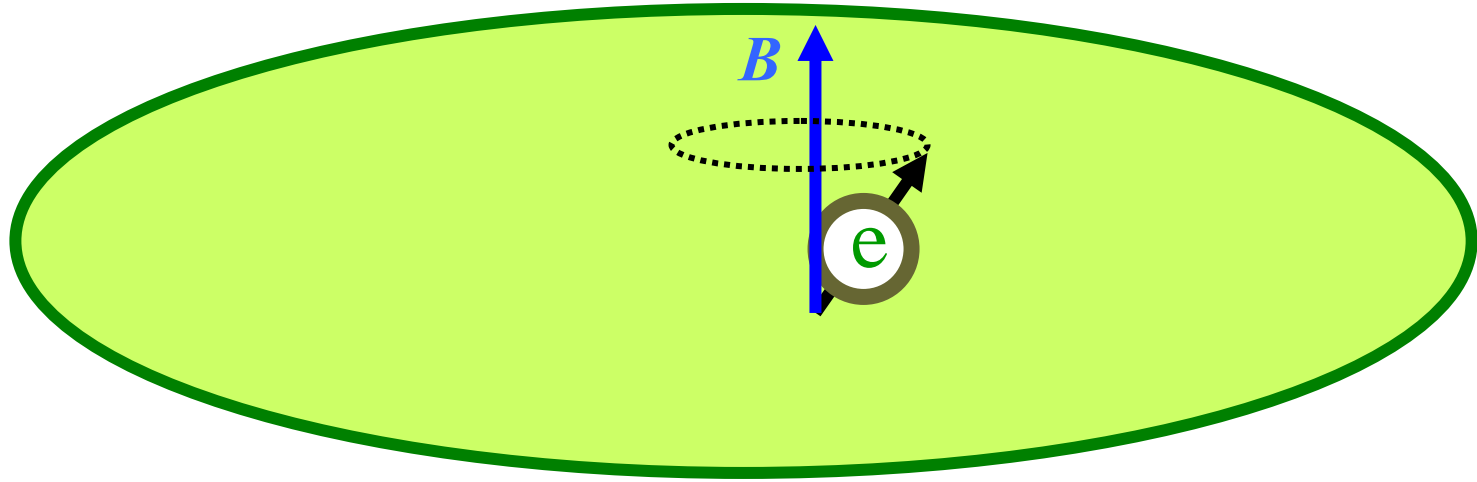
BCS Hamiltonian

$$\sum_{0 \leq \alpha \leq n-1} \varepsilon_\alpha \hat{H}_\alpha = \hat{H}_{BCS} + const \quad \lambda_{BCS} = \frac{2}{A}$$

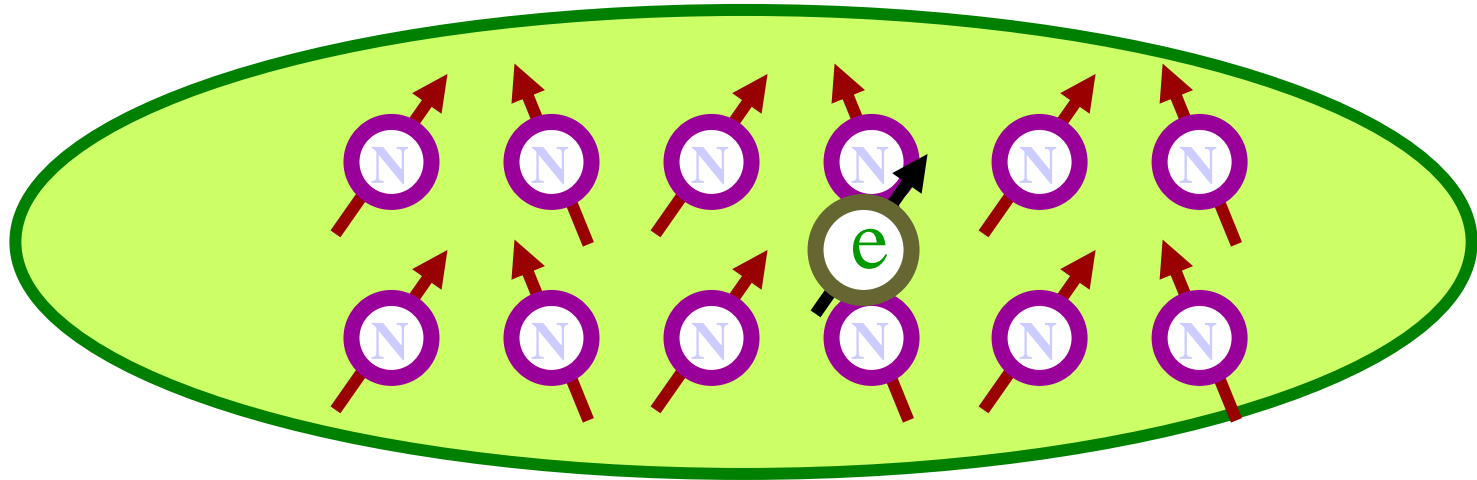
Overhauser interaction of electronic spins in quantum dots



Overhauser interaction of electronic spins in quantum dots



Overhauser interaction of electronic spins in quantum dots



- More than 10^6 of nuclear spins per electron
- Exchange interaction of the electronic spin with the nuclear ones.
- Collective effect of the nuclei on the electron is pretty strong.
- No interaction between the nuclear spins

$$\hat{H}_{eN} = \sum_{1 \leq \alpha \leq n-1} \gamma_{\alpha} \hat{S}_0 \hat{S}_{\alpha} - B \hat{S}_0^z$$

$$\gamma_{\alpha} \propto |\psi(\vec{r}_{\alpha})|^2$$

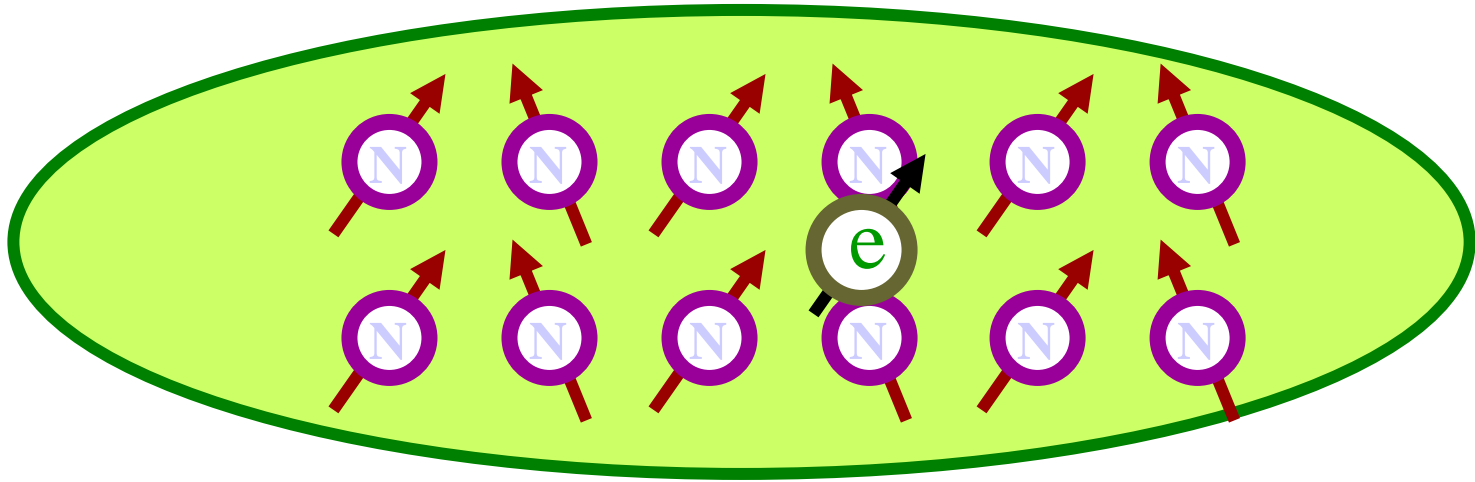
\hat{S}_0 – *spin of the electron*

$\hat{S}_{\alpha > 0}$ – *nuclear spins*

$\psi_{\alpha}(\vec{r})$ – *the electron w.f.*

B – *magnetic field || z*

Overhauser interaction of electronic spins in quantum dots



$$\hat{H}_{eN} = \sum_{1 \leq \alpha \leq n-1} \gamma_{\alpha} \hat{S}_0^{\hat{z}} \hat{S}_{\alpha}^{\hat{z}} - B \hat{S}_0^{\hat{z}}$$

Central spin problem

$$\hat{H}_{\alpha} = \sum_{0 \leq \beta \leq n-1, \beta \neq \alpha} \frac{2 \hat{T}_{\alpha}^{\hat{z}} \hat{T}_{\beta}^{\hat{z}}}{\epsilon_{\alpha} - \epsilon_{\beta}} - A \hat{T}_{\alpha}^{\hat{z}}$$

Gaudin problem

$$\hat{H}_{eN} = \hat{H}_0$$

$$\hat{S}_{\alpha}^{\hat{z}} \Rightarrow \hat{T}_{\alpha}^{\hat{z}}$$

$$\gamma_{\alpha} \Rightarrow \frac{2}{\epsilon_0 - \epsilon_{\alpha}}$$

$$B \Rightarrow A$$

$$\hat{H}_{eN} = \sum_{1 \leq \alpha \leq n-1} \gamma_{\alpha} \hat{S}_0^z \hat{S}_{\alpha}^z - B \hat{S}_0^z$$

Central spin problem

$$\hat{H}_{\alpha} = \sum_{0 \leq \beta \leq n-1, \beta \neq \alpha} \frac{2 \hat{T}_{\alpha}^z \hat{T}_{\beta}^z}{\varepsilon_{\alpha} - \varepsilon_{\beta}} - A \hat{T}_{\alpha}^z$$

Gaudin problem

$$\begin{aligned} \hat{S}_{\alpha} &\Rightarrow \hat{T}_{\alpha} \\ \hat{H}_{eN} = \hat{H}_0 &\quad \gamma_{\alpha} \Rightarrow \frac{2}{\varepsilon_0 - \varepsilon_{\alpha}} \\ B &\Rightarrow A \end{aligned}$$

Idea:

- Can we consider the classical dynamics of these Hamiltonians?
- Can it be described explicitly?
- What are connections between the classical and quantum dynamics?

Substitute quantum spin operators by classical vectors

$$\hat{T}_{\alpha}^z \Leftarrow \vec{s}_{\alpha}$$

$$\hat{L}^z \Leftarrow \vec{J}$$

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