Disordered Quantum Systems

Boris Altshuler

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Collaboration: Igor Aleiner, Columbia University

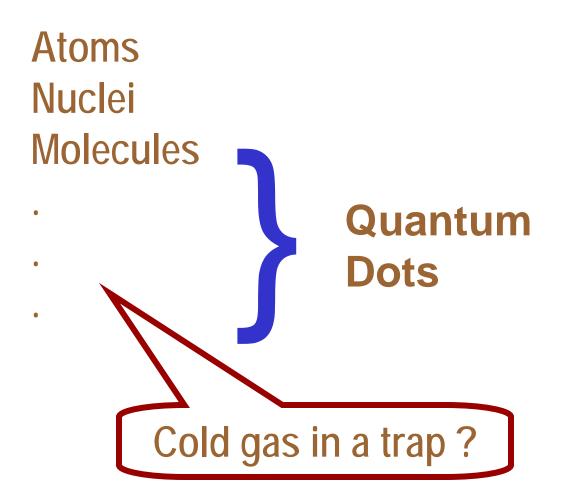
Part 1: Introduction

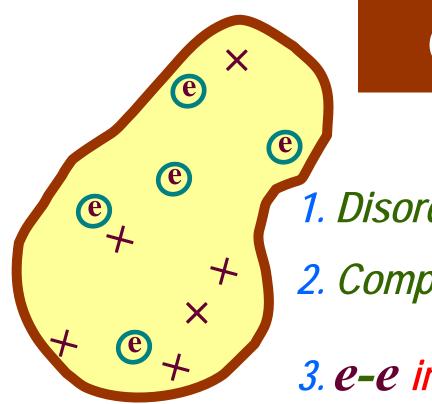
Part 2: BCS + disorder

INSTITUT HENRI POINCARE Centre Emile Borel Gaz quantiques

23 avril - 20 juillet 2007

Finite size quantum physical systems





Quantum Dot

Disorder (× – impurities)
 Complex geometry

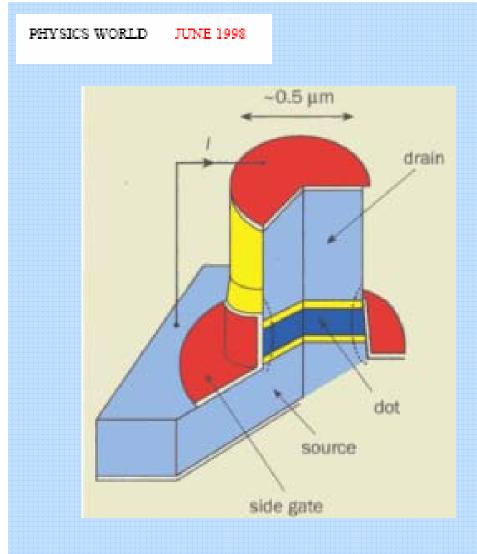
3. e-e interactions

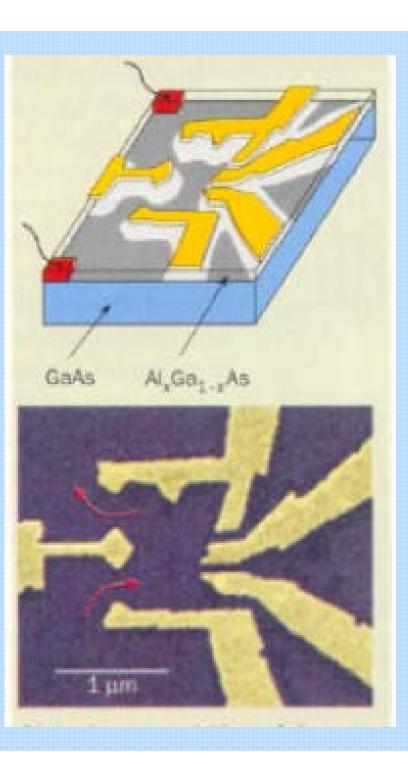
Realizations:

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

Quantum dots

Leo Kouwenhoven and Charles Marcus

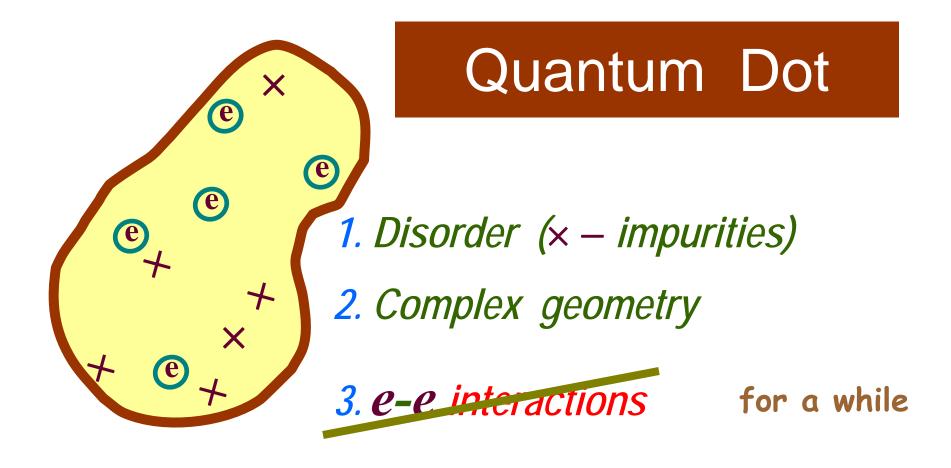




Finite number N of electrons: $\hat{H} \Psi_{\alpha} = E_{\alpha} \Psi_{\alpha}$

No interactions between electrons \rightarrow Shrodinger eqn in d dimensions

In the presence of the interactions between electrons \rightarrow Shrodinger equation in dN dimensions



Realizations:

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

I. Without interactions

Random Matrices, Anderson Localization Quantum Chaos

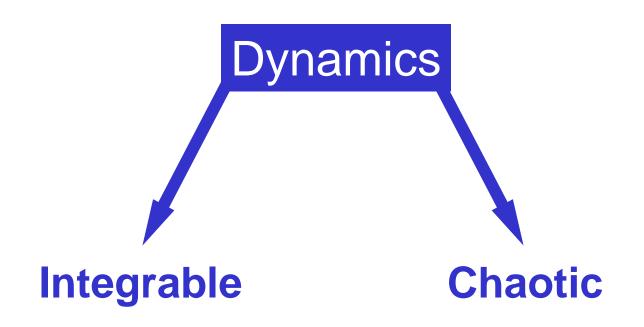
e 1. Disorder (× – impurities) e 2. Complex geometry e How to deal with disorder?

·Solve the Shrodinger equation exactly

•Start with plane waves, introduce the mean free path, and . . .

How to take quantum ? interference into account ?





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

Integrable Systems

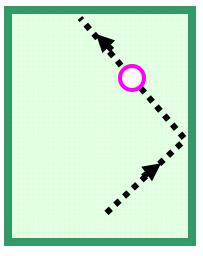
The variables can be separated and the problem reduces to *d* onedimensional problems



Examples

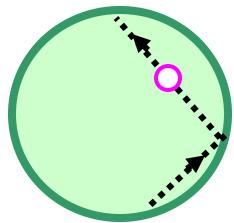
1. A ball inside rectangular billiard; d=2

- 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

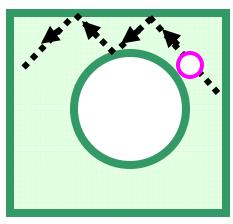
Integrable Systems

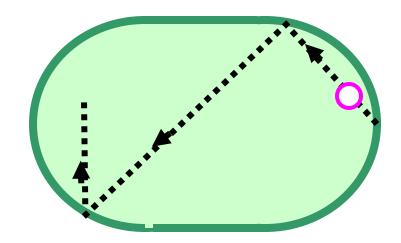
The variables can be separated $\Rightarrow d$ one-dimensional problems $\Rightarrow d$ integrals of motion

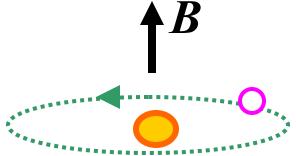
Rectangular and circular billiard, Kepler problem, ..., 1d Hubbard model and other exactly solvable models, ..

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

Examples







Kepler problem in magnetic field

Sinai billiard

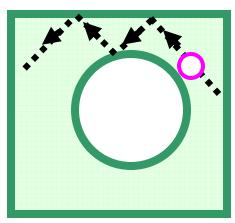
Stadium

The variables can not be separated ⇒ there is only one integral of motion - energy

Examples

Chaotic

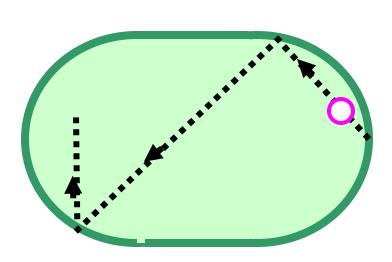
Systems



Sinai billiard



Yakov Sinai



Stadium



Leonid Bunimovich



Kepler problem in magnetic field

B

Johnnes Kepler

Integrable d-dimensional systems

d integrals of motion, *d* quantum numbers

$$I_k \qquad k = 1, 2, ..., d$$

Chaotic d-dimensional systems

The only conserved quantity is the energy Each eigenstate is characterized only by the eigenvalue of the Hamiltonian

Connection with the inverse problem:

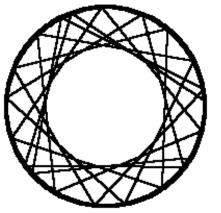
Q: Why original conditions can not be used as the integrals of motion ?

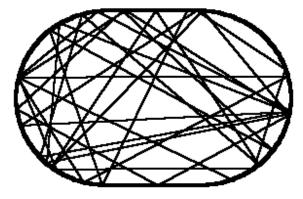
Classical Chaos $\hbar = 0$

•Nonlinearities

•Lyapunov exponents

Exponential dependence on the original conditions
Ergodicity

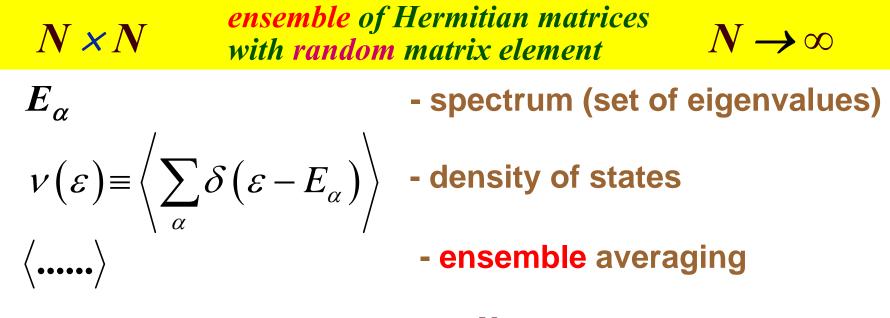


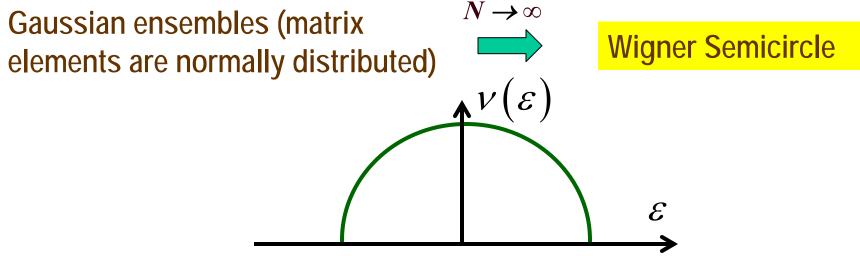


Quantum description of any System with a finite number of the degrees of freedom is a linear problem – Shrodinger equation

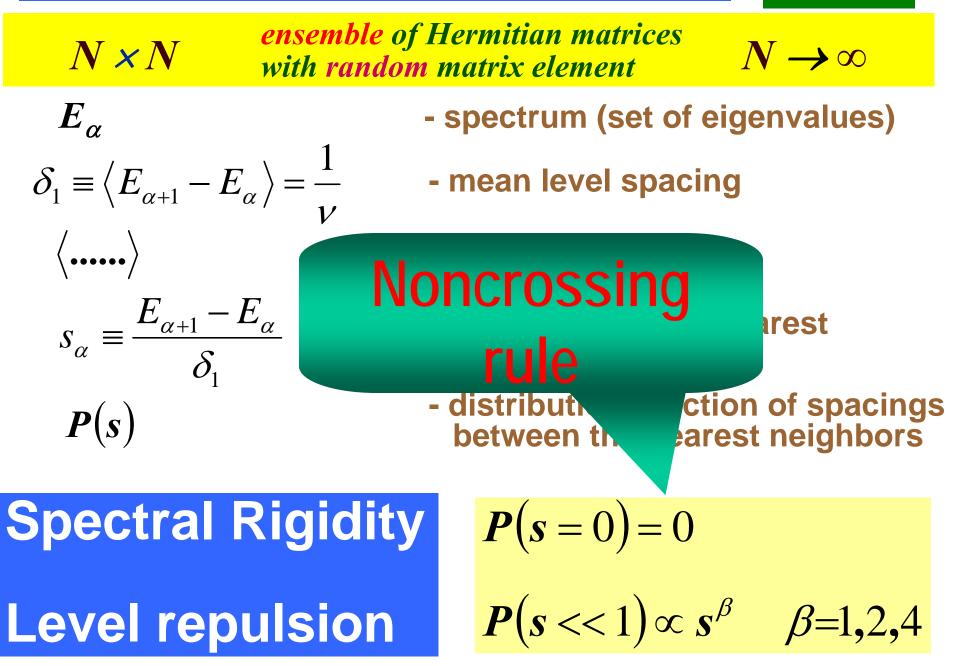
Q: What does it mean Quantum Chaos

RANDOM MATRICES





RANDOM MATRICES



Spectral

statistics

Noncrossing rule (theorem)

Suggested by Hund (Hund F. 1927 Phys. v.40, p.742)

Justified by von Neumann & Wigner (v. Neumann J. & Wigner E. 1929 Phys. Zeit. v.30, p.467)

Usually textbooks present a simplified version of the justification due to Teller (*Teller E., 1937 J. Phys. Chem 41 109*).

Arnold V. I., 1972 Funct. Anal. Appl.v. 6, p.94

Mathematical Methods of Classical Mechanics (Springer-Verlag: New York), Appendix 10, 1989

RANDOM MATRICES

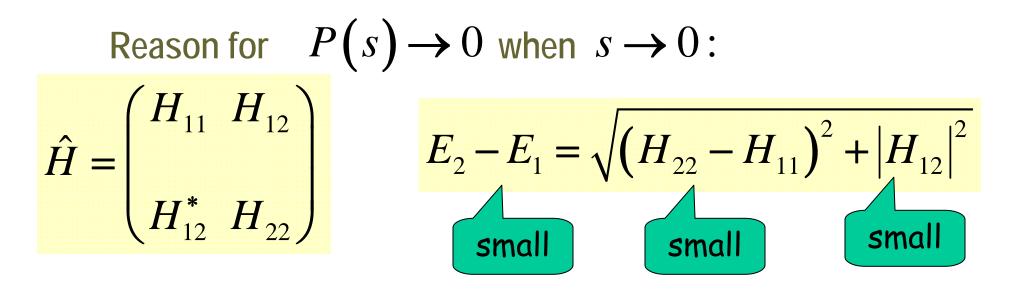


ensemble of Hermitian matrices with random matrix element

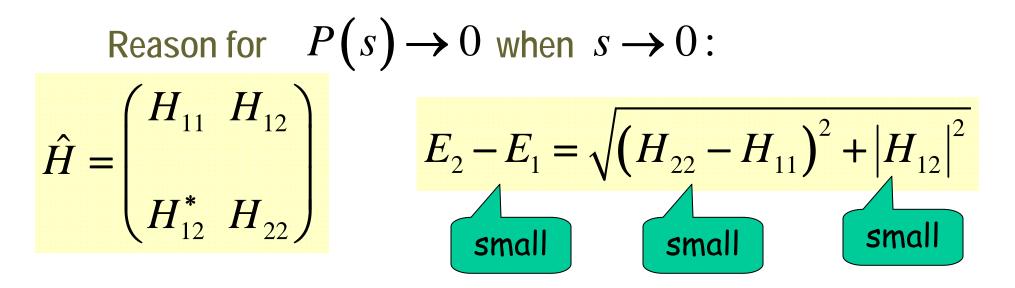
 $N \rightarrow \infty$

Dyson Ensembles

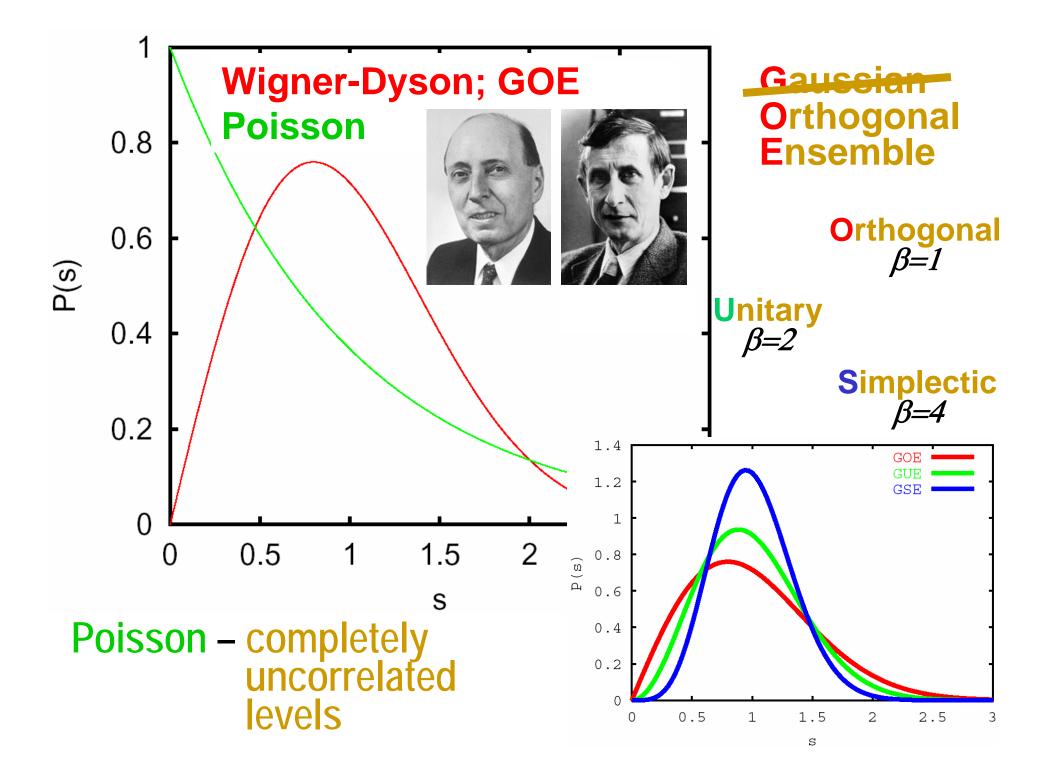
Matrix elements	Ensemble	ß
real	orthogonal	1
complex	unitary	2
2×2 matrices	simplectic	4



- 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 $P(s) \propto s$ $\beta = 1$



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- 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 $\implies P(s) \propto s^2 \qquad \beta = 2$



RANDOM MATRICES

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

$N \rightarrow \infty$

No conservation laws ⇒ no quantum numbers except the energy •

N imes N matrices with random matrix elements. $N o \infty$

Spectral Rigidity
Level repulsion
$$P(s \ll 1) \propto s^{\beta}$$
 $\beta = 1, 2, 4$

2

4

Dyson Ensembles

- Matrix elements Ensemble β
- real orthogonal
- complex unitary

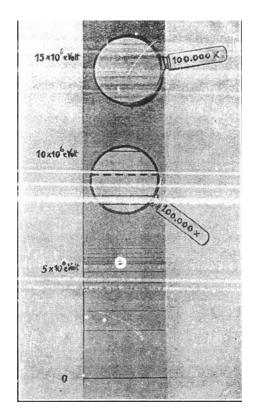
 2×2 matrices simplectic

Realizations

- **T-inv potential**
- broken T-invariance (e.g., by magnetic field)
- T-inv, but with spinorbital coupling

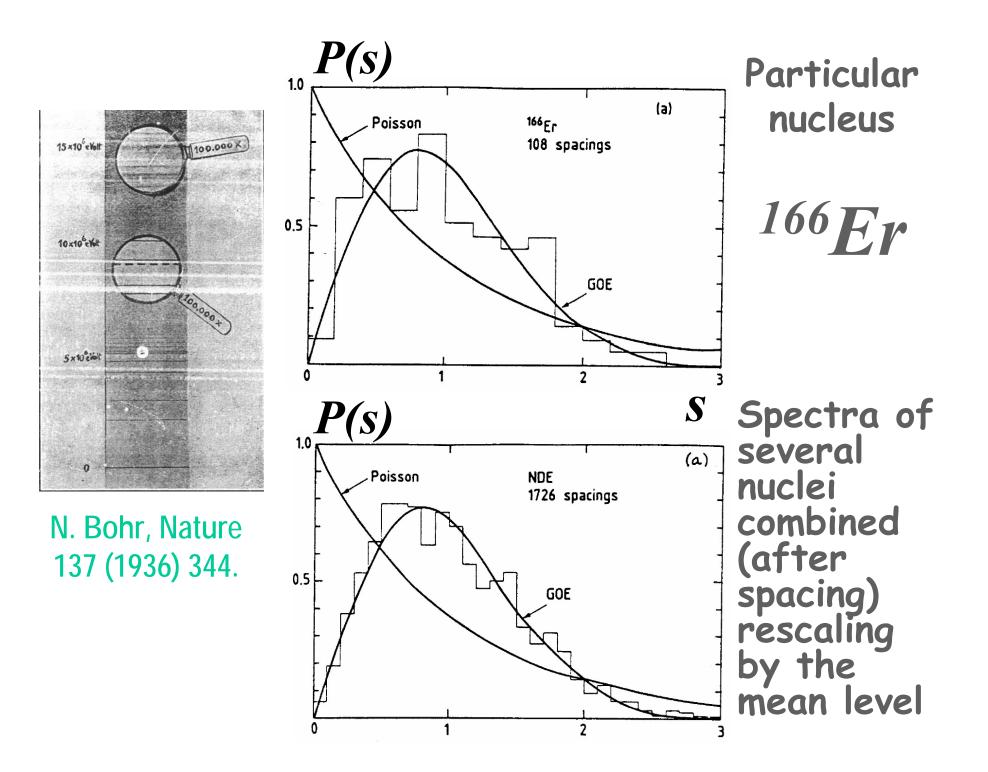
ATOMS Main goal is to classify the eigenstates in terms of the quantum numbers

NUCLEI For the nuclear excitations this program does not work



N. Bohr, Nature 137 (1936) 344.

ATOMSMain goal is to classify the eigenstates
in terms of the quantum numbersNUCLEIFor the nuclear excitations this
program does not workE.P. Wigner
(Ann.Math, v.62, 1955)Study spectral statistics of
a particular quantum system
- a given nucleus



ATOMS Main goal in terms o	is to classify the eigenstates f the quantum numbers	
NUCLEI For the nuclear excitations this program does not work		
E.P. Wigner (Ann.Math, v.62, 1955)	Study spectral statistics of a particular quantum system - a given nucleus	

Random Matrices	Atomic Nuclei
• Ensemble	• Particular quantum system
• Ensemble averaging	• Spectral averaging (over $lpha$)

Nevertheless are almost exactly the same as the Random Matrix Statistics

Why the random matrixtheory (RMT) works so wellfor nuclear spectra

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

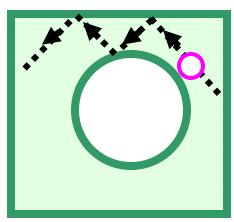
there exist very "simple" systems with as many as 2 degrees of freedom (d=2), which demonstrate RMT - like spectral statistics

The variables can not be separated ⇒ there is only one integral of motion - energy

Examples

Chaotic

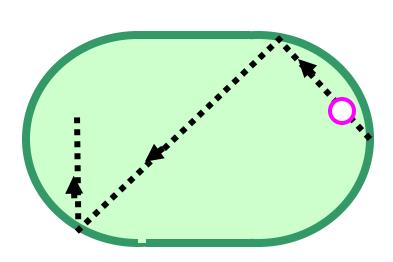
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$\hbar \neq 0$ Bohigas – Giannoni – Schmit conjecture

VOLUME 52

2 JANUARY 1984

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 timereversal-invariant systems whose classical analogs are K systems show the same fluctuation properties as predicted by GOE

$\hbar \neq 0$ Bohigas – Giannoni – Schmit conjecture

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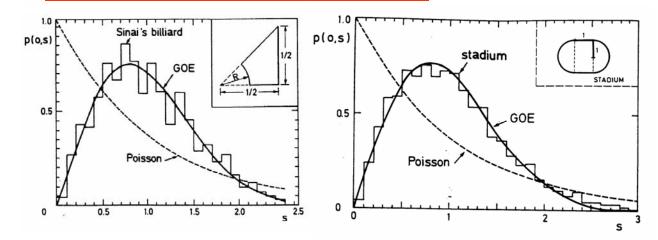
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In

Wigner-Dyson spectral statistics

No quantum

numbers except

energy

Chaotic

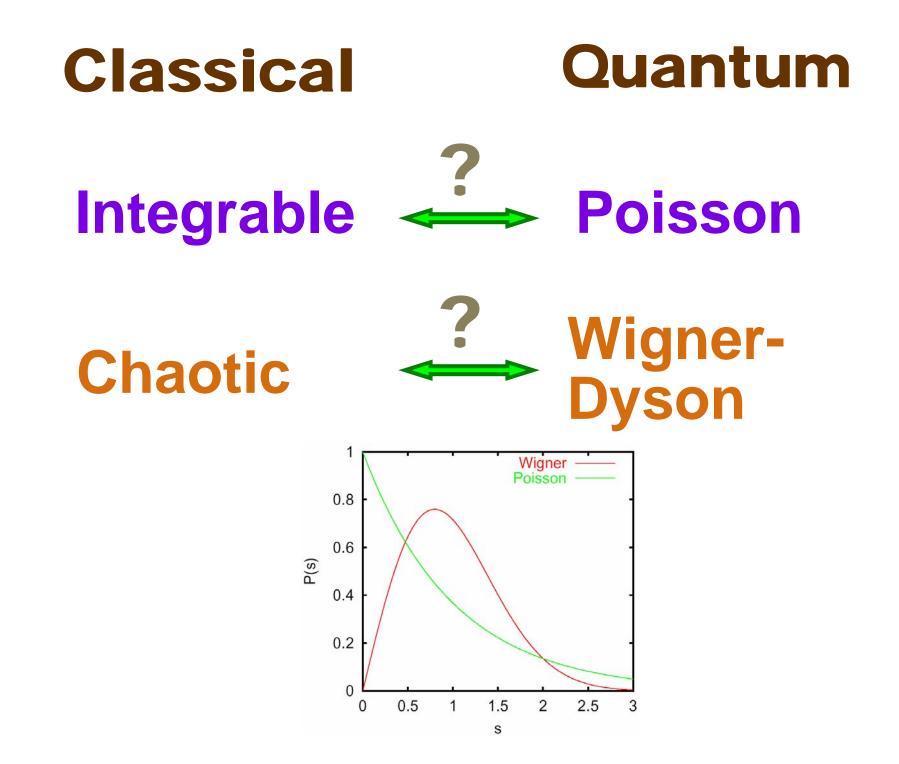
classical analog

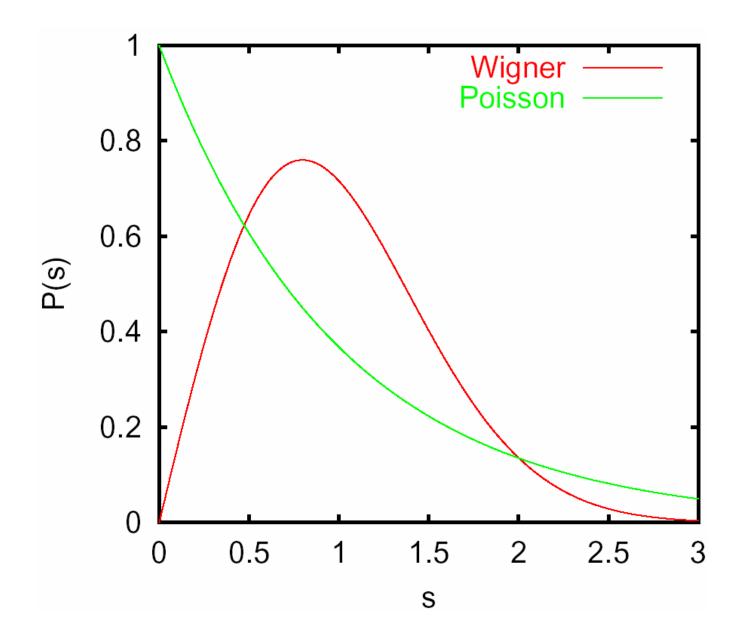
Q: What does it mean Quantum Chaos **?**

Two possible definitions

Chaotic classical analog

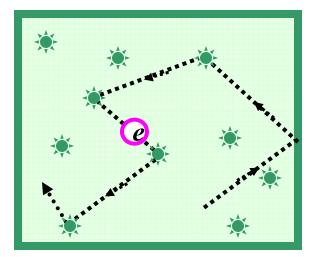
Wigner -Dyson-like spectrum





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

* Scattering centers, e.g., impurities



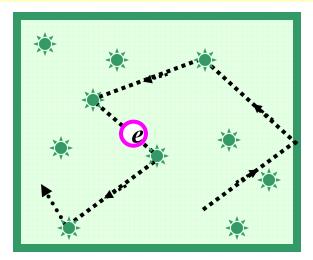
Important example: quantum particle subject to a random potential – disordered conductor

* 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.



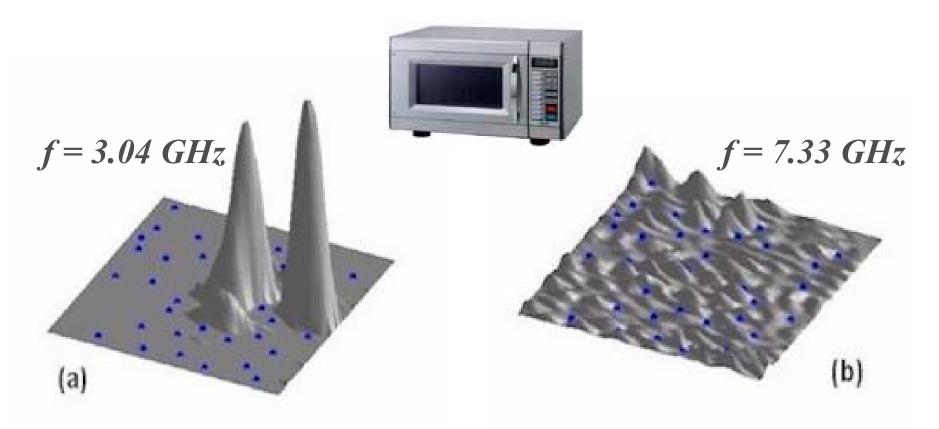


Anderson localization (1956)

At strong enough disorder all eigenstates are localized in space

Correlations due to Localization in Quantum Eigenfunctions of Disordered Microwave Cavities

Prabhakar Pradhan and S. Sridhar Department of Physics, Northeastern University, Boston, Massachusetts 02115 (Received 28 February 2000)



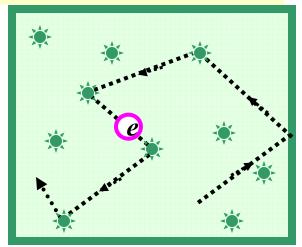
Anderson Insulator

Anderson Metal

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

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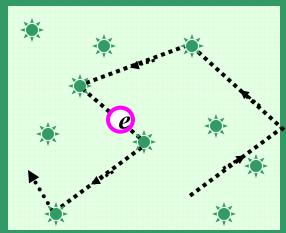
Models of disorder: Randomly located impurities



$$U(\vec{r}) = \sum_{i} u\left(\vec{r} - \vec{r}_{i}\right)$$

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

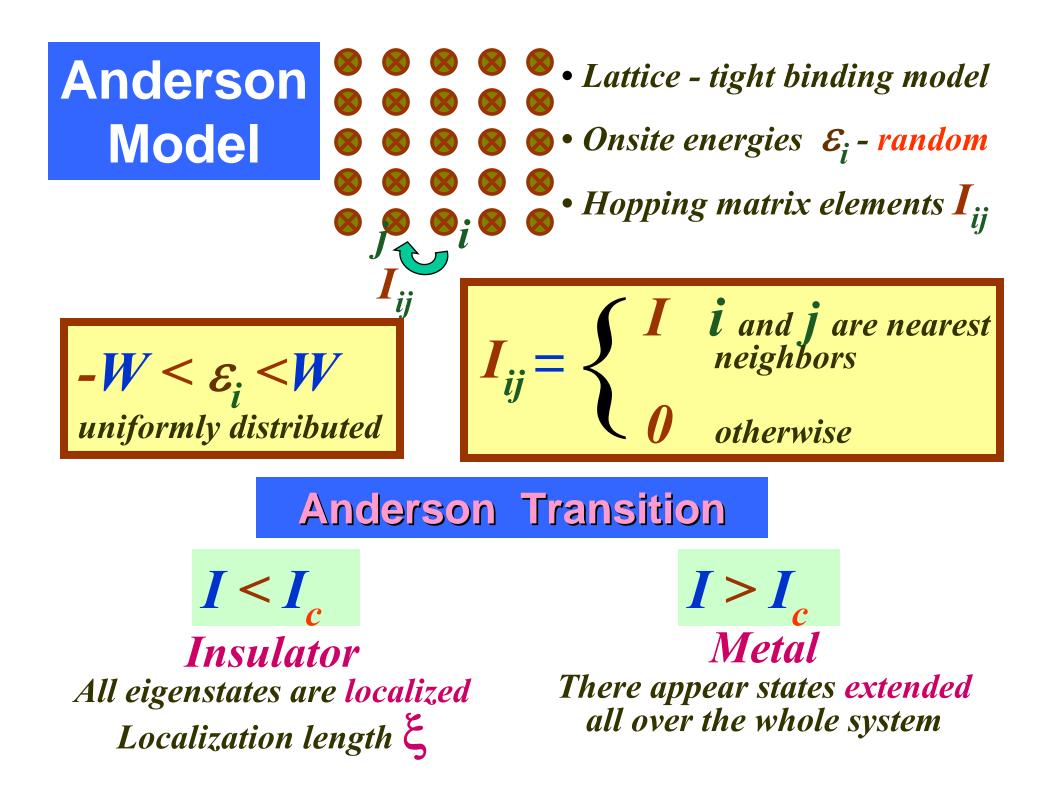
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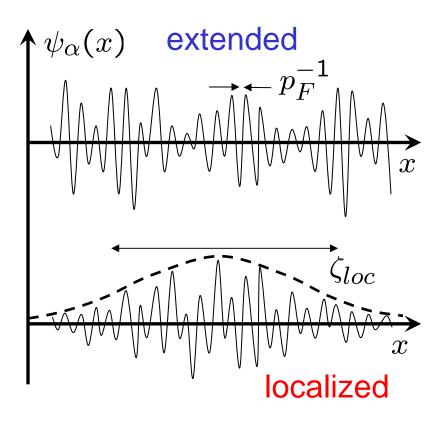
Models of disorder:
Randomly located impurities $U(\vec{r}) = \sum_{i} u(\vec{r} - \vec{r_i})$ White noise potential $u(\vec{r}) \rightarrow \lambda \delta(\vec{r})$ $\lambda \rightarrow 0$ $c_{im} \rightarrow \infty$ Anderson model – tight-binding model with onsite disorder

Lifshits model - tight-binding model with offdiagonal disorder

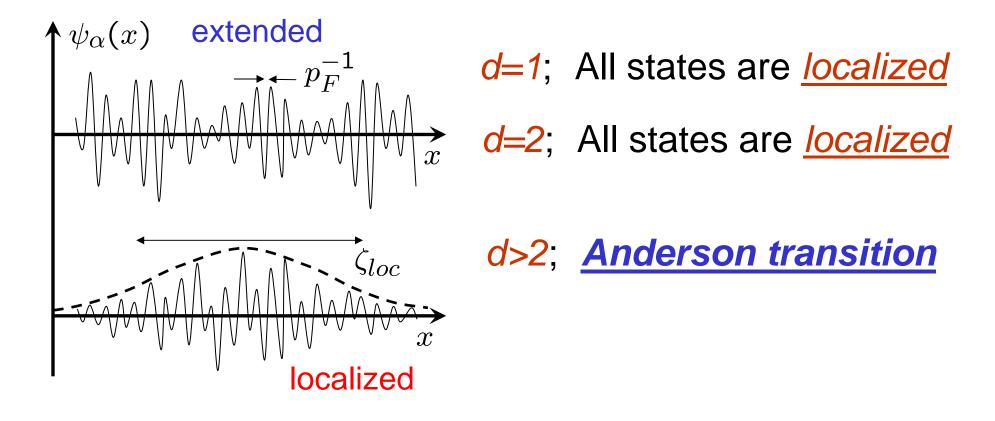
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- .
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Localization of single-electron wave-functions: $\left[-\frac{\nabla^2}{2m} + U(\boldsymbol{r}) - \epsilon_F\right]\psi_{\alpha}(\boldsymbol{r}) = \xi_{\alpha}\psi_{\alpha}(\boldsymbol{r})$



Localization of single-electron wave-functions: $\left[-\frac{\nabla^2}{2m} + U(\boldsymbol{r}) - \epsilon_F\right]\psi_{\alpha}(\boldsymbol{r}) = \xi_{\alpha}\psi_{\alpha}(\boldsymbol{r})$



Anderson Transition

 $I < I_c$

Insulator All eigenstates are localized Localization length ξ *Metal There appear states extended all over the whole system*

I > I

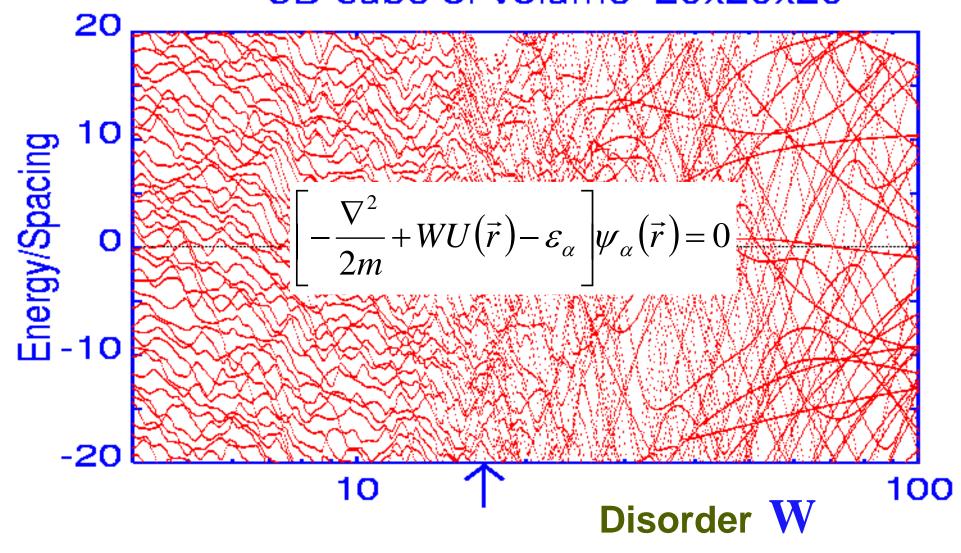
The eigenstates, which are localized at different places will not repel each other Any two extended eigenstates repel each other

Poisson spectral statistics

Wigner – Dyson spectral statistics

Zharekeschev & Kramer.

Exact diagonalization of the Anderson model 3D cube of volume 20x20x20



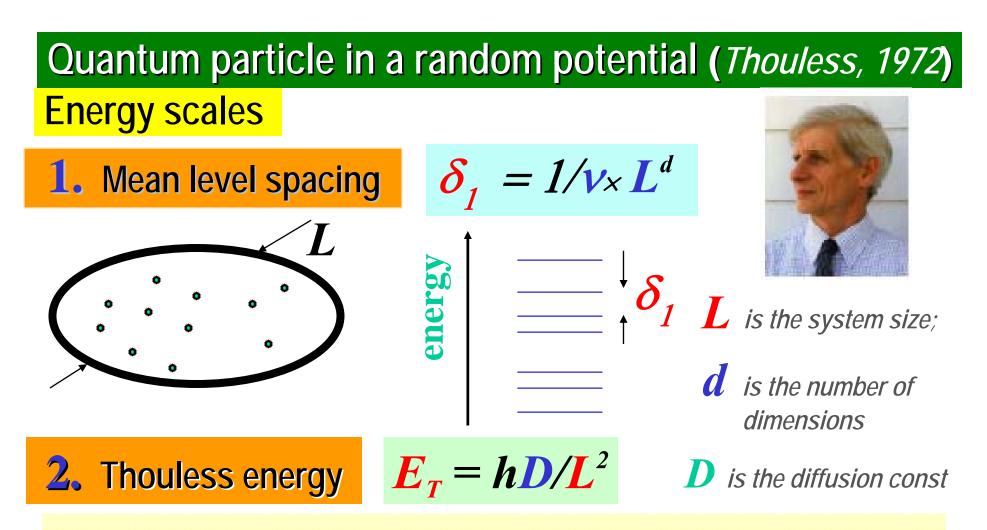
Q: What does it mean Quantum Chaos **?**

Two possible definitions

Chaotic ? Wigner classical Dyson-like analog spectrum

Are the two definitions equivalent?

Maybe not because of the localization!

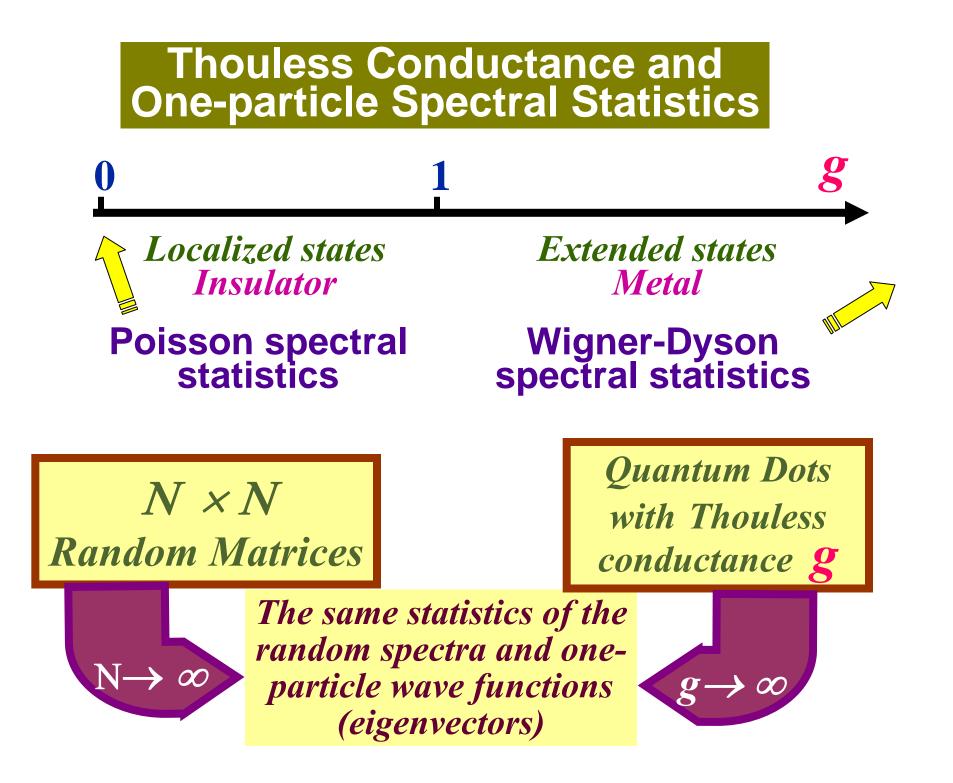


 E_T has a meaning of the inverse diffusion time of the traveling through the system or the escape rate (for open systems)

 $\mathbf{g} = \mathbf{E}_T / \delta_1$

dimensionless Thouless conductance





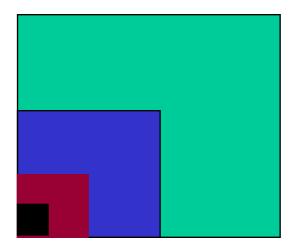
Scaling theory of Localization

(Abrahams, Anderson, Licciardello and Ramakrishnan 1979)

 $g = E_T / \delta_1$

Dimensionless Thouless conductance

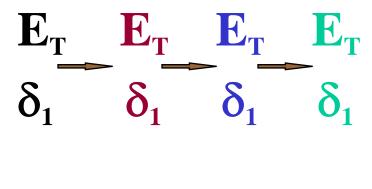




$$\mathbf{L} = 2\mathbf{L} = 4\mathbf{L} = 8\mathbf{L} \dots$$

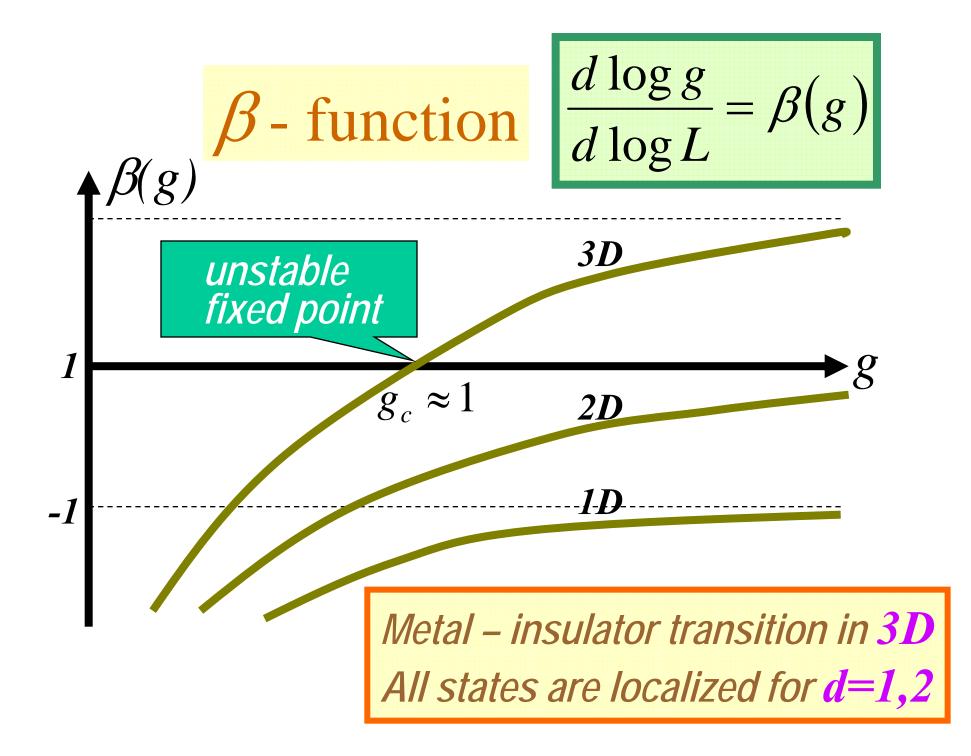
without quantum corrections

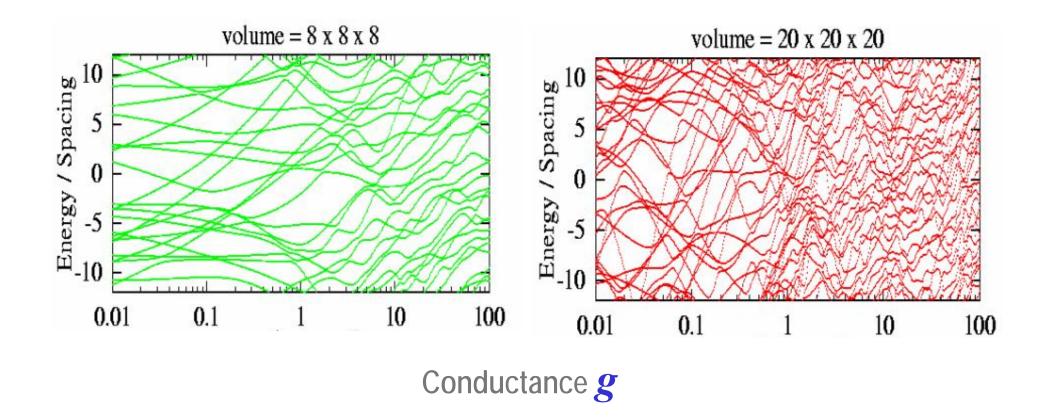
 $E_T \propto L^{-2} \quad \delta_1 \propto L^{-d}$



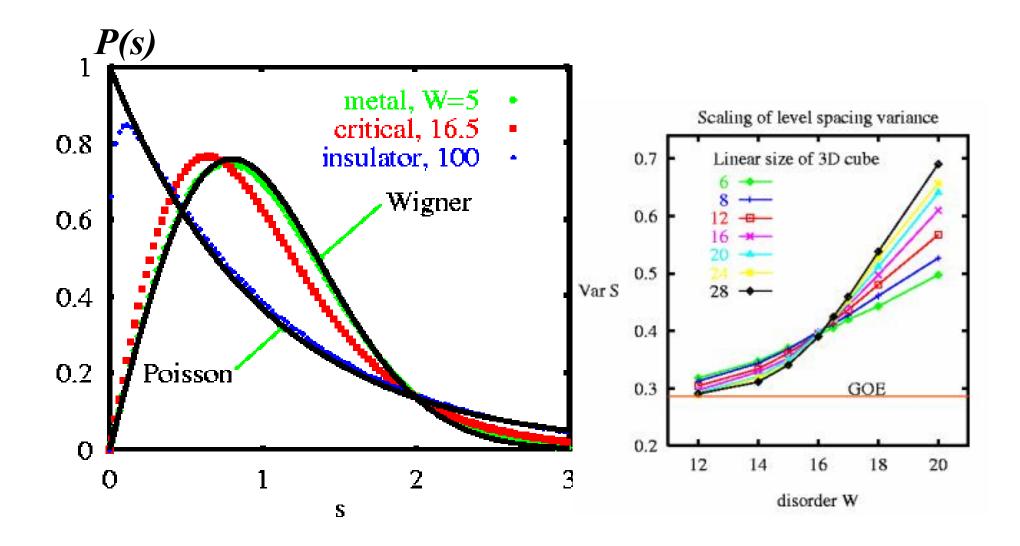
 $\mathbf{g} \longrightarrow \mathbf{g} \longrightarrow \mathbf{g} \longrightarrow \mathbf{g}$

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





Anderson transition in terms of pure level statistics

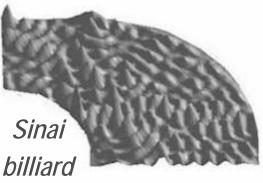


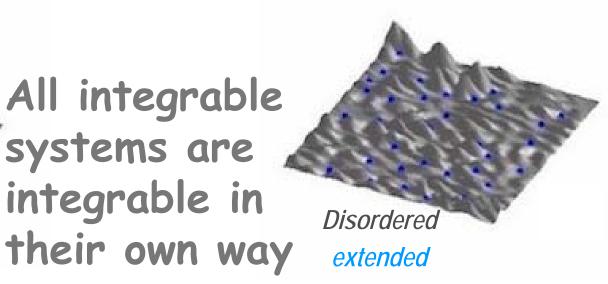
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Integrable

All chaotic systems resemble each other. Chaotic





billiard

Square

Disordered localized





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

Anderson metal; statistics

Anderson insulator; **Poisson** spectral statistics

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.

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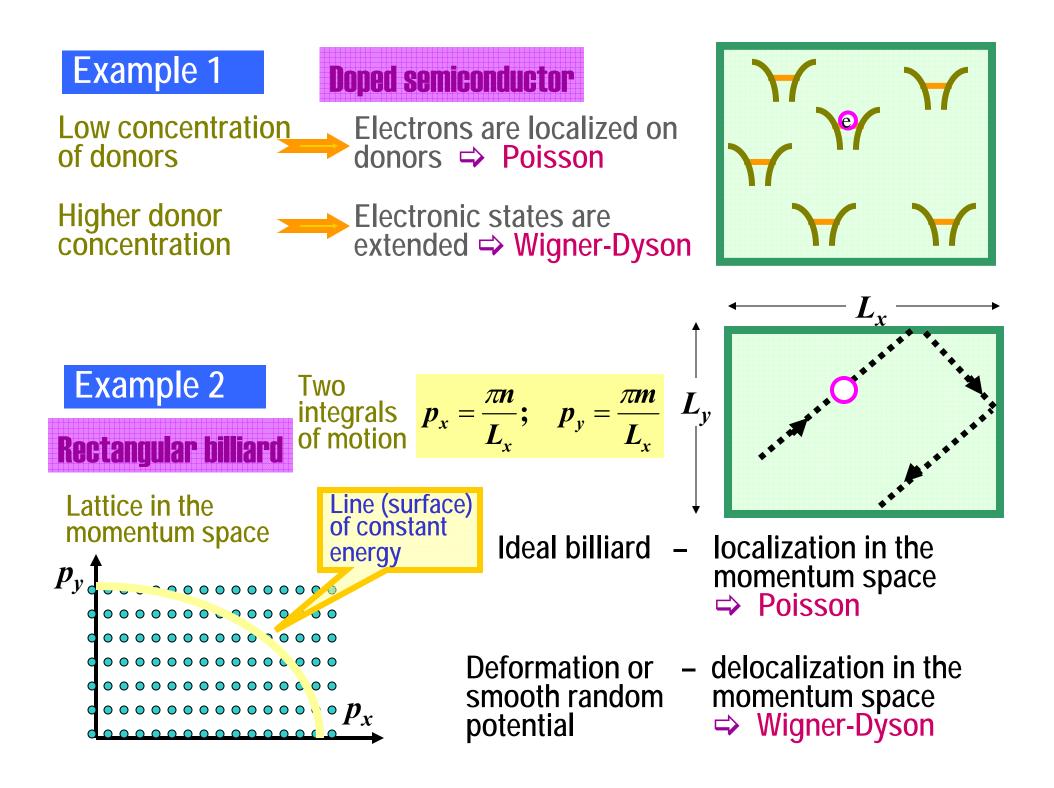
Weak enough hopping - Localization - Poisson Strong hopping - transition to Wigner-Dyson The very definition of the localization is not invariant - one should specify in which space the eigenstates are localized.

Level statistics is invariant:

Poissonian statistics

basis where the eigenfunctions are localized

Wigner -Dyson statistics basis the eigenfunctions

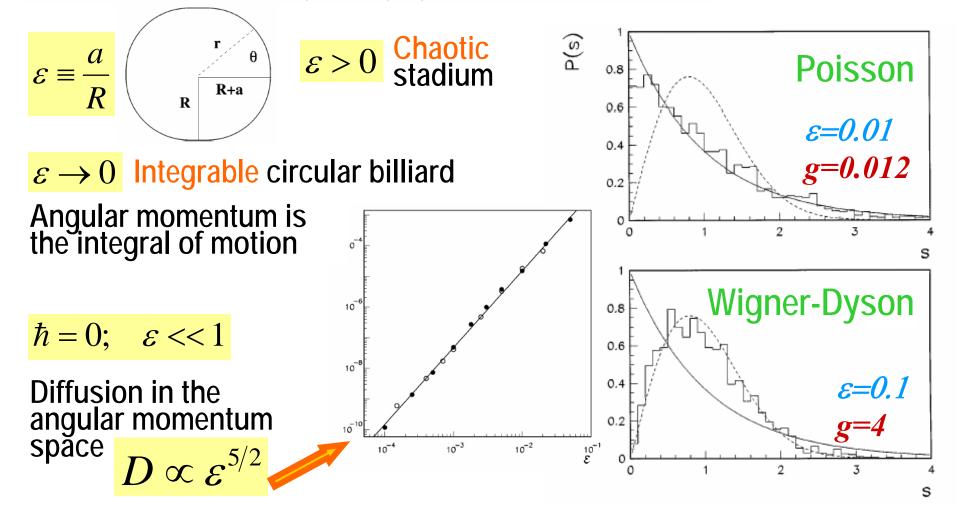


2 December 1996

Diffusion and Localization in Chaotic Billiards

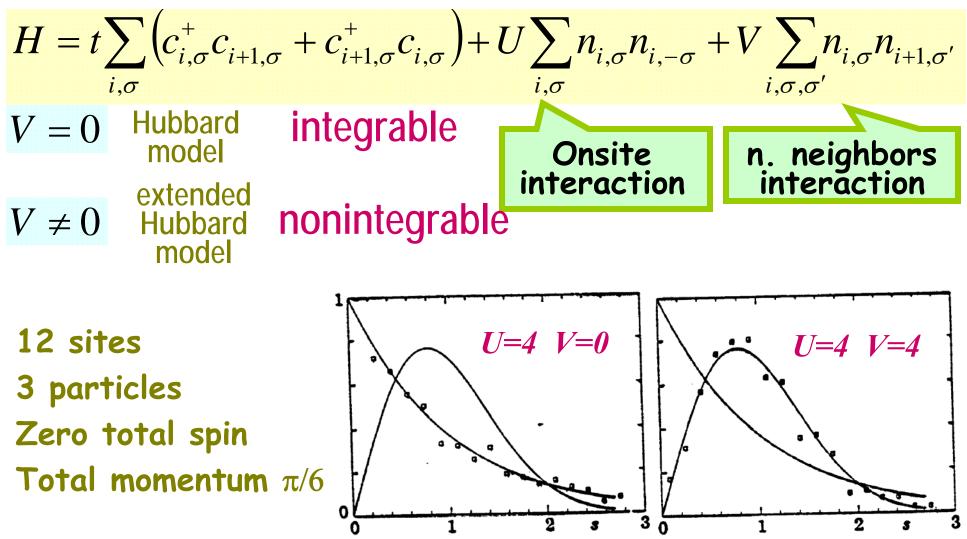
Fausto Borgonovi,^{1,3,4} Giulio Casati,^{2,3,5} and Baowen Li^{6,7} ¹Dipartimento di Matematica, Università Cattolica, via Trieste 17, 25121 Brescia, Italy ²Università di Milano, sede di Como, Via Lucini 3, Como, Italy ³Istituto Nazionale di Fisica della Materia, Unità di Milano, via Celoria 16, 22100, Milano, Italy ⁴Instituto Nazionale di Fisica Nucleare, Sezione di Pavia, Pavia, Italy ⁵Instituto 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 ^{a7}Center for Applied Mathematics and Theoretical Physics, University of Maribor, Krekova 2, 2000 Maribor, Slovenia (Received 29 July 1996)

Localization and diffusion in the angular momentum space



D.Poilblanc, T.Ziman, J.Bellisard, F.Mila & G.Montambaux *Europhysics Letters*, v.22, p.537, 1993

1D Hubbard Model on a periodic chain



Finite number N of electrons:

$$\hat{H}\Psi_{\alpha} = E_{\alpha}\Psi_{\alpha}$$

No interactions between electrons → Shrodinger eqn in d dimensions

Integrable system – each energy is conserved Poissonian many-body spectrum

In the presence of the interactions between electrons \rightarrow Shrodinger eqn in dN dimensions

Finite number N of electrons:

$$\hat{H}\Psi_{\alpha} = E_{\alpha}\Psi_{\alpha}$$

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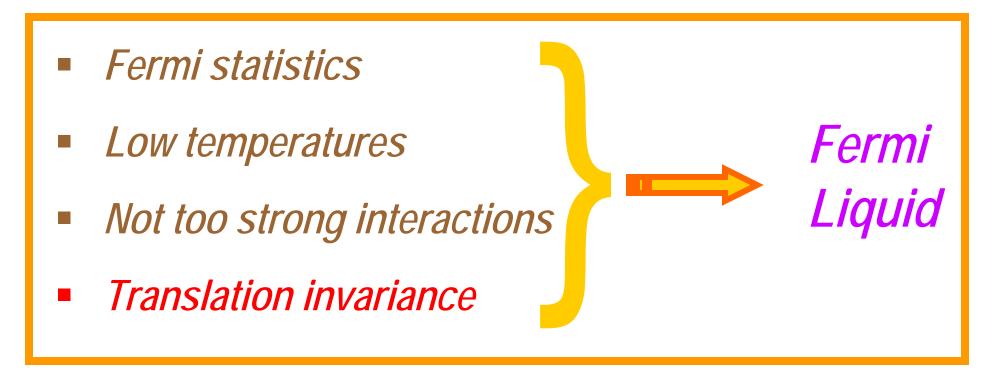
Q. Can interaction between the particles drive ? this system into chaos and make it ergodic?

Random Matrics statistics of nuclear spectra

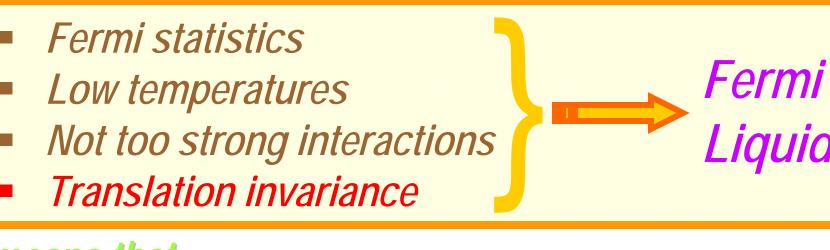
II. With interactions

Fermi Liquid and Disorder Zero Dimensional Fermi Liquid





What does it mean?



It means that

- Excitations are similar to the excitations in a Fermi-gas:

 a) the same quantum numbers momentum, spin ½, charge e
 b) decay rate is small as compared with the excitation energy
- 2. Substantial renormalizations. For example, in a Fermi gas

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

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 moments 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 (?!) $n(\vec{p})$

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



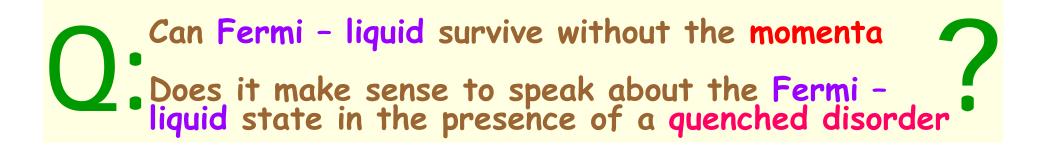
$$G(\varepsilon, \vec{p}) = \frac{Z}{i\varepsilon_n - \xi(\vec{p})}$$

 p_F

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

Landau Fermi - Liquid theory

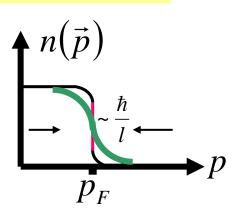
Momentum	\vec{p}
Momentum distribution	$n(\vec{p})$
Total energy	$E\{n(\vec{p})\}$
Quasiparticle energy	$\xi(\vec{p}) \equiv \delta E / \delta n(\vec{p})$
Landau f-function	$f(\vec{p},\vec{p}') \equiv \delta\xi(\vec{p})/\delta n(\vec{p}')$



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

 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 Does it make sense to speak about the Fermi –
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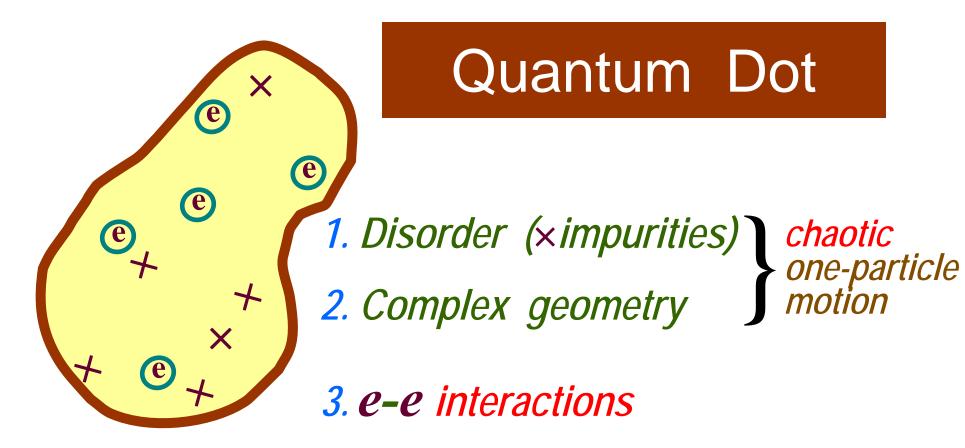
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- 2. Neither resistivity nor its temperature dependence is determined by the umklapp processes and thus does not behave as T^2
- *3.* Sometimes (e.g., for random quenched magnetic field) the disorder averaged oneparticle Green function even without interactions does not have a pole as a function of the energy, *c*. The residue , *Z*, makes no sense.

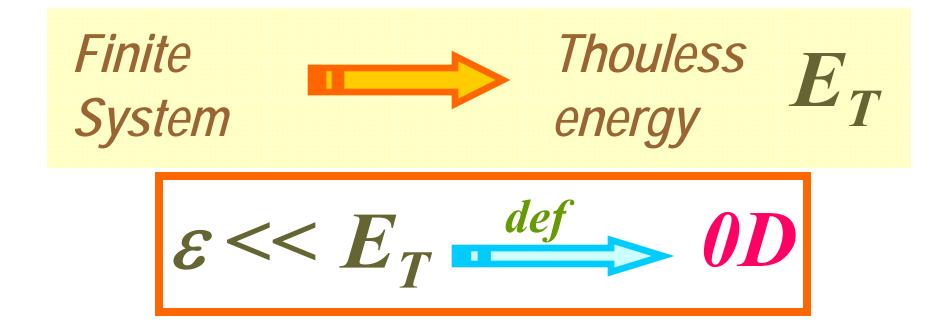
Nevertheless even in the presence of the disorder

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



Realizations:

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



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

$$\delta_1 << \varepsilon << E_T$$

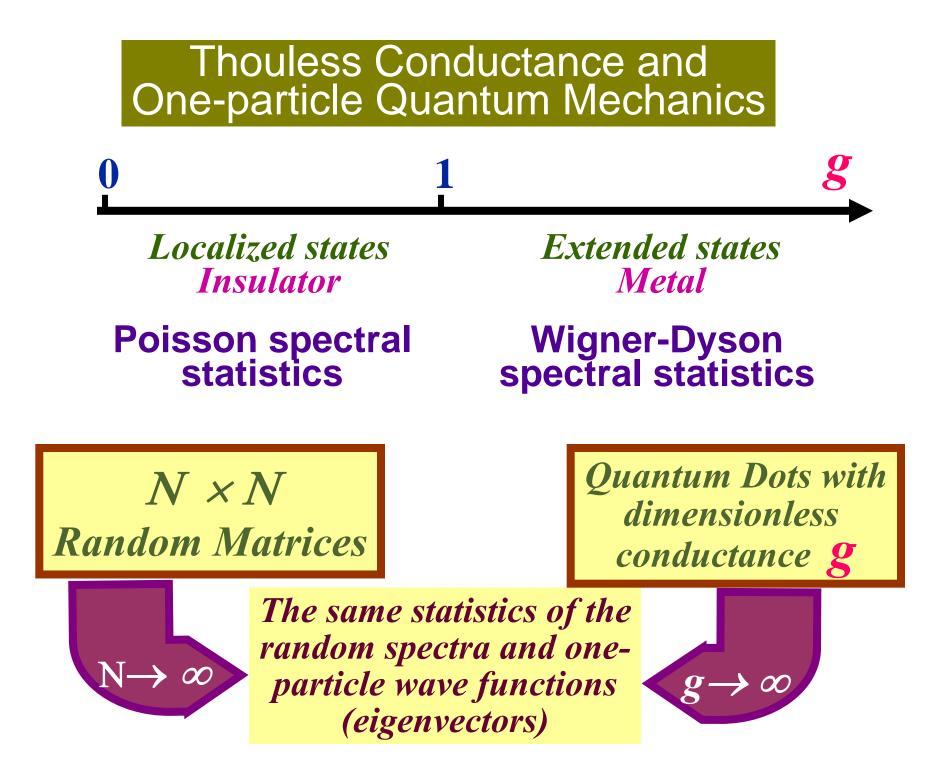
$$g \equiv \frac{E_T}{\delta_1} >> 1$$

Two-Body
InteractionsSet of one particle states. σ
and α label correspondingly
spin and orbit. $\hat{H}_0 = \sum_{\alpha} \varepsilon_{\alpha} a^+_{\alpha,\sigma} a_{\alpha,\sigma}$ $\hat{H}_{int} = \sum_{\substack{\alpha,\beta,\gamma,\delta\\\sigma,\sigma'}} M_{\alpha\beta\gamma\delta} a^+_{\alpha,\sigma} a^+_{\beta,\sigma'} a_{\gamma,\sigma} a_{\delta,\sigma'}$

 \mathcal{E}_{α} -one-particle orbital energies

 $M_{lphaeta\gamma\delta}$ -interaction matrix elements

<i>Nuclear Physics</i>	\mathcal{E}_{α}	are taken from the shell model
	Μ _{αβγδ}	are assumed to be random
Quantum Dots	\mathcal{E}_{α}	RANDOM; Wigner-Dyson statistics
	$M_{_{lphaeta\gamma\delta}}$???????



Matrix Elements

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

Μαβγδ

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 $g \rightarrow \infty$

Matrix Elements

• Diagonal matrix elements are much bigger than the offdiagonal ones

 $M_{\rm diagonal} >> M_{\rm offdiagonal}$

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

Toy model: Short range *e-e* interactions $U(\vec{r}) = \frac{\lambda}{-\delta} \delta(\vec{r}) \qquad \frac{\lambda}{V} \text{ is dimensionless coupling constant}$ \boldsymbol{v} is the electron density of states

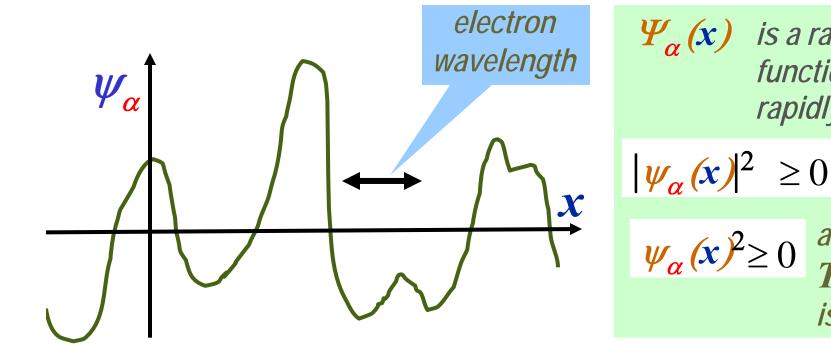
$$M_{\alpha\beta\gamma\delta} = \frac{\lambda}{\nu} \int d\vec{r} \,\psi *_{\alpha} (\vec{r}) \psi *_{\beta} (\vec{r}) \psi_{\gamma}(\vec{r}) \psi_{\delta}(\vec{r})$$

$$\psi_{\alpha}(\vec{r})$$
one-particle
eigenfunctions

is a random

function that

rapidly oscillates



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

In the limit



 Diagonal matrix elements are much bigger than the offdiagonal ones

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

$$M_{\alpha\beta\alpha\beta} = \frac{\lambda}{\nu} \int d\vec{r} |\psi_{\alpha}(\vec{r})|^{2} |\psi_{\beta}(\vec{r})|^{2}$$
$$\implies M_{\alpha\beta\alpha\beta} = \lambda \delta_{1}$$
$$|\psi_{\alpha}(\vec{r})|^{2} \Rightarrow \frac{1}{\text{volume}}$$

<u>More general</u>: finite range interaction potential $U(\vec{r})$

$$M_{\alpha\beta\alpha\beta} = \frac{\lambda}{\nu} \int \left| \psi_{\alpha}(\vec{r}_{1}) \right|^{2} \left| \psi_{\beta}(\vec{r}_{2}) \right|^{2} U(\vec{r}_{1} - \vec{r}_{2}) d\vec{r}_{1} d\vec{r}_{2}$$

The same conclusion

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

All correlation functions are invariant under arbitrary orthogonal transformation:

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

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

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

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

$$\hat{S} = \sum_{\alpha,\sigma_{1},\sigma_{2}} a_{\alpha,\sigma_{1}}^{+} \vec{\sigma}_{\sigma_{1},\sigma_{2}} a_{\alpha,\sigma_{2}}$$

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

$$\tilde{K}^{+} = \sum_{\alpha} a_{\alpha,\uparrow}^{+} a_{\alpha,\downarrow}^{+}$$

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

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

Therefore, in a very general case

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

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:

$$\begin{split} \hat{H} &= \sum_{\alpha} \varepsilon_{\alpha} n_{\alpha} + \hat{H}_{int} \\ \hat{H}_{int} &= eV\hat{n} + E_c\hat{n}^2 + J\hat{S}^2 + \lambda_{BCS}\hat{K}^+\hat{K}. \end{split}$$

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

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For a short range interaction with a coupling constant ${\cal X}$

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

where δ_1 is the one-particle mean level spacing

$$\hat{H} = \hat{H}_0 + \hat{H}_{int} \qquad \qquad \hat{H}_0 = \sum_{\alpha} \mathcal{E}_{\alpha} n_{\alpha}$$

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

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

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

- *E_c* determines the charging energy (Coulomb blockade)
- J describes the spin exchange interaction



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

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Isn't it a Fermi liquid ?

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