

Gergen Lectures      Duke University

October 22-24, 2001

Random Matrix Theory

Lecture 1 It is a great pleasure and honor for me to give the year 2001 Gergen Lectures. I am also very pleased to be here with so many of my friends. My [topic] is random matrix theory with emphasis on the relationship to integrable systems. In this set [first] of 6 lectures, I am going to begin with some

- general remarks about integrable systems. Then I will
- make some general and historical remarks about random matrix theory (RMT)

- universality conjecture in RMT

In the  $\boxed{2^{\text{nd}}}$  lecture, I will

- introduce a variety of objects and techniques from modern theory of integrable systems

In the  $\boxed{3^{\text{rd}}}$  lecture, I will show how

- to use these techniques to solve the universality conjecture following the approach of T. Kriecherbauer, K. McLaughlin, S. Venakides, X. Zhou and P. D.

In the  $\boxed{4^{\text{th}}}$  and final lecture, I will show how

- to use these techniques to analyze the so-called Ulam problem in the theory of random permutations.

### Integrable systems

The modern Theory of integrable systems began in with the discovery of Gardner, Greene, Kruskal and Miura of a method for integrating the Korteweg de Vries (KdV) equation

$$u_t + uu_x + u_{xxx} = 0 \quad , \quad x \in \mathbb{R}, \quad t \geq 0$$

$$u(x, t=0) \rightarrow 0 \quad \text{as} \quad |x| \rightarrow \infty$$

This equation arises in the Theory of water waves

(3)

Initially the discovery of Gardner et al was regarded as providing a method of solution for a rather thin set of evolutionary equations, but by the early 1980's it started to become clear that the discovery of Gardner et al was just the first glimpse of a far more general integrable method that would eventually have applications across the broad spectrum of problems in pure and applied mathematics. In the narrowest sense, an integrable system is a

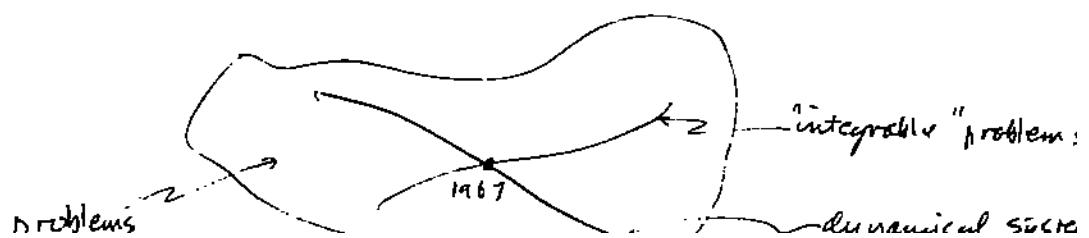
Hamiltonian dynamical system — finite or infinite dimensional — with "enough" integrals of the motion all of whose Poisson brackets are zero, to solve the system in some "explicit" form. Now it has been a rather extraordinary experience in the field over the last 30 years, that many systems which are of great mathematical and physical interest, which may be Hamiltonian, and may not even be dynamical, can be solved "explicitly" using techniques that have a direct link back to the method of solution for the KdV equation discovered by Gardner et al. The kind of developments that I have in mind are for example

- The solution of the classical Schottky problem in algebraic geometry in terms of the solution of

- The introduction of quantum groups
- integrable statistical models — connection to Jones polynomials
- 2D quantum gravity and the work of Witten and Kondratenko on the I<sub>1</sub>A<sub>1</sub>V hierarchy
- nonlinear special function theory (Painlevé theory)
- conformal field theory — work of Knizhnik, Zuber
- \* • random matrix theory
- combinatorial problems of Robinson-Schensted-Knuth

The thrust of these lectures is to illustrate how the integrable methods interact with one of these top — random matrix theory.

Allow me a very schematic moment. Consider the following picture

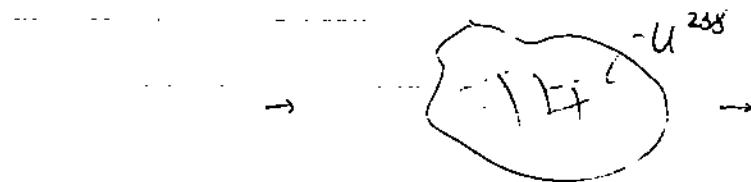


In 1967, these two "manifolds" had a "transverse intersection". The initial thought was that to develop the ideas of Gardner et al. one should move along the "pde" manifold. But this turned out to be too limiting: we now know that in order to reach the full development of the method one should move in the "transverse" direction. And this is the direction we will move in the next 2+ lectures.

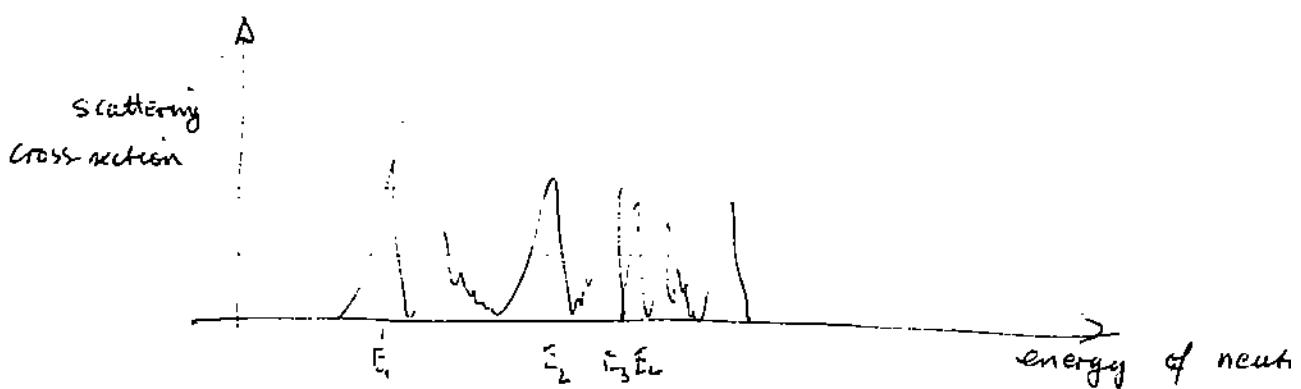
### RMT

Now what is random matrix theory? RMT was first studied in the 1930's in the context of mathematical statistics by Hsu, Wishart and others but it was Wigner in the early 1950's who introduced RMT to mathematical physics. Wigner was concerned with the scattering of neutrons off (large)

nuclei



Schematically one sees the following:



The energies  $E_i$  for which one obtains a large scattering cross section are called resonances, and for a heavy nucleus like  $U^{238}$ , say, there can be hundreds of them. In theory, one could write down many Schrödinger equations for 16 neutrons + nucleus  $U$  and try to solve it numerically to compute these  $E_i$ . But in Wigner's time and still in our time,

also into the foreseeable future, this was not a real approach, and so people began to think that it was more appropriate to give the resonances a statistical meaning. But what should the statistical model be? At this point Wigner put forward the remarkable hypothesis that the (high) resonances  $E_i$  behave like the  ~~$\lambda$~~  eigenvalues of a (large) matrix. It is hard to overemphasize what a radical and revolutionary theory this was: we all recall that when we first were learning some physics, we understood that the detail of the model were paramount: if you changed the force law in the equations of motion, the behavior of the system would change. But now all of that was out the window. The precise mechanism was

no longer important. All that Wigner (<sup>, and later Dyson,</sup> required was that

or real symmetric

- The matrices be Hermitian (so that the eigenvalues were real)

- That the ensemble behaves "naturally" under certain  $\begin{matrix} N \times N \\ \text{Hermitian} \end{matrix}$   $\begin{matrix} N \times N \\ \text{real symmetric} \end{matrix}$   $\begin{matrix} 2N \times 2N \\ \text{Hermitian}, s \\ H = J H^T \end{matrix}$  physical symmetry groups (GUE, GOE, GSE ensemble — more later)

During the late 50's, 60's and early 70's, various research (particularly, <sup>both sides</sup>) began to test Wigner's hypothesis against real experimental

data in a variety of physical situations and the

results were pretty impressive. A classic reference for

RMT is

- Random matrices — M.-L. Mehta, 2nd Ed

and I invite you to look at p17 for a comp

of 1726 nuclear level spacings against predictions of

the GOE ensemble.

procedure: first we must rescale the resonances so

the eigenvalues of the random matrices so that

number of resonances / unit interval

=

expected number of eigenvalues / unit interval

then we compare the statistics.

More concretely, we typically have a situation where

we are interested in resonances in an interval  $(E, E + \Delta E)$

where  $\Delta E \ll E$  but  $\Delta E$  contains many resonances

We then rescale the resonances in  $\Delta E$  so that the

average density is one. We then do a similar

scaling for the eigenvalues in some ensemble, and

only at that point, after we have adjusted

our "microscopes" do we compare statistics. A scientist

approaches such problems with two instruments in

— a microscope and a list of ensembles

account for

- The dials on the microscope are adjusted to the macroscopic situation, and vary from system to system,
- but once the "slide" is in focus, one sees universal behavior described by one of the entries in the list of ensembles.

Having given me a schematic moment, I now ask you for a small philosophical moment.

Now it is always something of a mysterious process when begin to think of something that is quite deterministic in a statistical way. Nevertheless, it

for example

take a die, which surely obeys the laws of solid mechanics, but we readily and intuitively understand it as a statistical object. Moreover we "know" what

the stochastic model should be: all 6 sides have equal probability.

Looked at from this point of view, what Wigner had to do was to see the neutron + nucle-

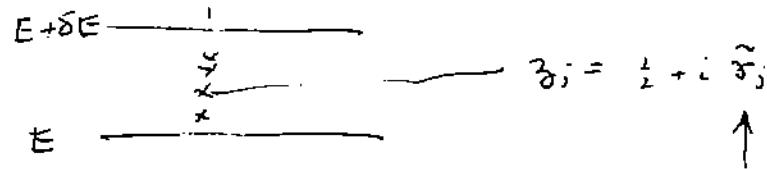
for its stochasticity. In all such problems there is an  
some singular / asymptotic process } involved and when we cross some "Bayesian  
point, phase space opens up, and all bets are off.  
We are standing, as it were, on 42<sup>nd</sup> corner of 8<sup>th</sup>  
Av 42<sup>nd</sup> Street and we are watching - this little kid  
play 3 card Monte: if we are fast enough, we  
can follow all his moves but then "poof!" the card  
are out there, and all 3 <sup>cards</sup> are equally likely

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Now in the early 1970's a very remarkable  
thing happened. Montgomery, quite independently of the  
other goings on, began thinking that the zeros of the  
Riemann zeta function on the critical line  $\Re s = \frac{1}{2}$ ,  
should also be viewed statistically. And so, assuming  
and rescaling the imaginary parts of the zeros

$$\left( = \frac{1}{N} \# \underbrace{\left\{ \text{pairs } (i, j) , \quad i \leq i+j \leq N, \quad \text{st } a \leq \tilde{x}_i - \tilde{x}_j \leq b \right\}} \right)$$

for the (rescaled) zeros in an interval  $[(E, E + \delta E) \subset E >> 1,$



$$\text{rescaled: } \tilde{\delta_j} = : \frac{\delta_j \log \delta_j}{2\pi} :$$

$$\#\{x_i \in T\} \approx |T| \text{ as}$$

and he found a limiting formula  $R = R(a, s)$

$$\left( \int_a^b \left[ 1 - \left( \frac{\sin \pi u}{2\pi u} \right)^2 \right] du \right)$$

for the 2 pt. function as  $E \rightarrow \infty$ ,  $\#\tilde{\delta}_i + \delta E \rightarrow \infty$ . (6)

happened next is very well-known (I even asked Ayson

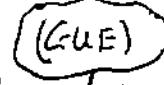
to authenticate this version). Id outcomes met Ayson at 1

at the Institute in Princeton, and when he told him

about his calculations, Dyson immediately wrote down

a formula and asked Montgomery, "Did you get this

(1)

had he known the answer, Dyson said "Well, if the zeros of the zeta function behaved like the eigenvalues of a random GUE matrix, this would have to be the answer!" Indeed what Montgomery had obtained for  $R$  was precisely the 2-pt function for the eigenvalues of a (large) random  matrix.

At this point, the cat was out of the bag. People began asking whether their favorite list of #'s behaved like the eigenvalues of a random matrix. Through the 80's an extraordinary variety of systems were investigated from this point of view with astounding results; for example, if we take a region in the plane



and look at the eigenvalues  $0 = \lambda_1 \leq \lambda_2 \leq \dots$  of

the Dirichlet Laplacian in this region they they too  
(after the standard scaling) behave like the eigenvalues  
of a random GUE matrix (see Mehta p13)

In the late 1980's, Montgomery's work was followed  
up numerically by Odlyzko, who then confirmed  
Montgomery's work to high accuracy & also investigated  
other statistics such as the nearest neighbor spacing  
again there was incredible agreement with random  
matrix theory. In recent years, the work of Montgomery  
Odlyzko has been a wonderful springboard for  
Sarnak - Rudnick and then Sarnak - Katz, to prove all  
kinds of GUE (& GSE) random matrix properties  
for the zeros of all kinds of automorphic L functions

Now up until very recently, the physical and mathematical phenomena which were investigated, concerned the eigenvalues of a random matrix in the bulk of the spectrum.

But in the last 2 yrs or so a very interesting class of problems have started to appear in combinatorics and also in statistical particle models, which concern the eigenvalues at the top of the spectrum. For example, consider the following version of solitaire called "patience sorting". Suppose we have N cards numbered linearly  $1, 2, \dots, N$  for convenience.

Shuffle the cards

1 2 3 - -  $\infty$

Now take the top card (1c)

Eventually you end up with  $P_N(\pi)$ . files Question

Putting uniform distribution on a set of

permutation shuffles  $(\pi^3 - S_N)$ , how does  $P_N(\pi)$  behave as  $N \rightarrow \infty$ ? differently Or said, how

big a table do I need typically to play the game with  $N$  cards?

To state the answer to this question we need to introduce some notation

(which we will discuss in the 4<sup>th</sup> lecture,  
the Theorem is the following (Baik, Johansson, P.D.)

As  $N \rightarrow \infty$ ,  $P_N$ , suitably centred and scaled behaves statistically like the largest eigenvalue of a random GUE matrix

More precisely suppose  $\lambda_1^{(N)}(u) \geq \lambda_2^{(N)}(u) \geq \dots \geq \lambda_N^{(N)}(u)$

(1)

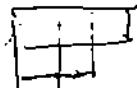
Are the eigenvalues of a (large)  $N \times N$  GUE matrix

then

$$\lim_{N \rightarrow \infty} \text{Prob} \left( \frac{p_N - 2\sqrt{N}}{N^{1/6}} \leq t \right) = \lim_{N \rightarrow \infty} \text{Prob} \left( \frac{\sum_{i=1}^{N^2} a_i - \sqrt{N}}{2^{-1/2} N^{1/6}} \leq t \right) \\ = F(t)$$

$F(t)$  is called the Tracy-Widom distribution after it was discovered and can be expressed in terms of a solution of the Painlevé-II equation.

There are many equivalent formulations of the patience sorting problem — longest increasing subsequence of a random permutation, the height of a nucleate droplet in a supersaturated medium in the <sup>no-called</sup> Polyakov growth model, the number of boxes in the first row of a Young diagram under Plancherel measure



out that the # of boxes in the 2<sup>nd</sup> row behaves statistically like the 2<sup>nd</sup> largest eigenvalue  $\lambda_2^{(n)}$  of a random GUE matrix, and so on.

Now what are these three basic distributions mentioned above that were singled out by Wigner and Dyson (on the basis of the behavior of the system under time reversal and change of ~~reference frame~~ reference frame)?

\* i) Gaussian Unitary Ensemble (GUE) consisting of

(a)  $N \times N$  Hermitian matrices  $M = (M_{ij})$

(b) with probability distribution

$$P(M)dM = P(M) \prod_{i=1}^N dM_{ii} \prod_{i>j} d\text{Im} M_{ij} \prod_{i>j} d\text{Im} M_{ji}$$

↑  
normalization density

which is invariant under unitary conjugation  $U$

$$M \rightarrow U^\dagger M U = M'$$

$$\text{i.e. } P(M')dM' = P(M)dM$$

(c)  $(M_{11}, \text{Re } M_{12}, \text{Im } M_{12})$  are independent so

$$P(M) = \prod_{i=1}^N \varphi^{(1)}(M_{ii}) \prod_{i>j} \varphi^{(2)}(M_{ij}) \prod_{i>j} \varphi^{(3)}(M_{ji})$$

(2c)

2) Gaussian Orthogonal ensemble (GOE)

(a)  $N \times N$  real symmetric

3) Gaussian Symplectic ensemble (GSE)

(a)  $2N \times 2N$  Hermitian self dual  $M = M^*$

$$M = J M^T J^{-1}$$

$$J = \begin{pmatrix} 0 & I \\ -I & 0 \end{pmatrix}$$

(b)  $\dots$

involved unit  $M \mapsto U^T M U$ ,  $U$  unit  
symplectic

$$U^T J U = J$$

(c)  $\dots$

The analysis of GUE is the simplest and I am going to  
restrict myself to this case throughout these lectures

So focusing on GUE, the first Theorem in the business  
is that if  $P(M)$  satisfies (a)-(b)-(c) then necessarily

$$P(M) = \text{const } e^{-\frac{1}{2}(\alpha M^2 + \beta M + \delta)}$$

where  $\alpha > 0, \beta, \delta \in \mathbb{R}$ . Centering and rescaling we have the

GUE distribution  
↑ Gaussian

$$(2) P(M) = \frac{1}{\sqrt{\pi}} e^{-\frac{1}{2}M^2} dM$$

Now here comes the problem: whereas conditions (a) & (b) are physical, (c) is just a device and has no physical basis.

If we just assume (a) and (b) we find that

$$(3) \quad P(m) dm = \frac{1}{Z_n} e^{-\text{Tr } V(m)} dm$$

for some real valued function  $V$ ,  $V(x) \rightarrow \infty$  as  $|x| \rightarrow \infty$   
(are called Unitary Ensembles)

Ensembles of matrices of type (3). Now physicists turn

This all on its head and say "as there is no physical

way to distinguish between different choices of  $V$ , then

whatever answers we compute, the answer must be independent  
of  $V$ "

This is a rough form of what is meant by

universality in random matrix theory. In the

coming lectures we will focus on a particular <sup>basic</sup> statistic

(

The probability  $\overset{(v)}{P}(a,b)$  that a (large) matrix has no eigenvalues in an interval  $(a,b)$ , and we will show that  $\overset{(v)}{P}(a,b)$  indeed has a universal form independent of  $V$ . And in proving this, we will employ techniques that have arisen in integrable Theory over the years and can be traced back in some form to the ~~original~~ <sup>original</sup> method of solution for the kdv equation discovered in 1967 by Gardner, Greene, Kruskal and Miura.

To summarize, what has emerged, somewhat mysteriously and to everyone's surprise in a very powerful heuristic: all kinds of things physical and mathematical, behave, in some limit, like random matrices. Next time you have a system that you want to ~~approximate~~ model with a large matrix, go for it. You have a good chance of being in such a way