

Integrable random matrix ensembles

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We propose new classes of random matrix ensembles whose statistical properties are intermediate between statistics of Wigner-Dyson random matrices and Poisson statistics. The construction is based on integrable N -body classical systems with a random distribution of momenta and coordinates of the particles. The Lax matrices of these systems yield random matrix ensembles whose joint distribution of eigenvalues can be calculated analytically thanks to integrability of the underlying system. Formulas for spacing distributions and level compressibility are obtained for various instances of such ensembles.

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I. INTRODUCTION

The theory of random matrices, introduced by Wigner in the 1950s, has proved to be a very useful tool in many fields of physics, from localisation theory to quantum transport (see e.g. [1] and references therein). In quantum chaos, a well accepted conjecture states that Wigner-Dyson random matrix ensembles describe statistical properties of spectra of quantum systems whose classical counterpart is chaotic [2], while statistics of integrable systems is best described by Poisson statistics of independent random variables [3]. The corresponding wave functions are extended in the chaotic case and localised in the integrable case. The choice of the random matrix ensemble suited to describe the statistical behaviour of a system depends on the symmetries of that system. In the usual setting [4], standard random matrix ensembles consist of matrices M with independent Gaussian random elements whose measure is invariant over conjugation

$$M \longrightarrow U^{-1} M U, \quad (1)$$

where U is an arbitrary matrix belonging to one of the three following groups of matrices: unitary, orthogonal, or symplectic. The unitary group defines the Gaussian Unitary Ensemble (GUE), which is supposed to describe statistical properties of energy levels of chaotic systems without time-reversal invariance. The orthogonal group corresponds to Gaussian Orthogonal Ensemble (GOE), used for time-reversal invariant chaotic systems. The symplectic group gives rise to Gaussian Symplectic Ensemble (GSE), applicable to time-reversal chaotic systems with half-integer spin without rotational symmetry.

Though many extensions and generalisations of random matrices have been proposed in order to best describe various models [5], the existence of a large invariance group as in (1) remains their characteristic feature. Without such invariance it is very difficult to connect analytically simple properties of matrix elements with complex properties of matrix eigenvalues. For all random matrix ensembles with invariance group it is possible to integrate over unnecessary variables in order to get explicitly the joint distribution of eigenvalues under the form

$$P(\lambda_1, \dots, \lambda_N) \sim \prod_{i < j} |\lambda_j - \lambda_i|^\beta e^{-\sum_k V(\lambda_k)}, \quad (2)$$

with $V(x)$ a system-dependent potential and β a parameter. For the Gaussian ensembles the potential is quadratic and the parameter β is equal to 1 for GOE, 2 for GUE, and 4 for GSE. All correlation functions for invariant ensembles can be calculated analytically [4]. However the resulting formulas are cumbersome. For the nearest-neighbour distribution $P(s)$, instead of the exact expression one often uses a simple surmise proposed by Wigner. This surmise has correct functional dependence at small and large argument and takes the form

$$P(s) = a s^\beta e^{-b s^2} \quad (3)$$

with constants a and b determined from the normalisation conditions

$$\int_0^\infty P(s) ds = \int_0^\infty s P(s) ds = 1. \quad (4)$$

The Wigner-type surmise for the probability $P(n, s)$ that between two eigenvalues separated by s there exist exactly $n - 1$ other levels (with $P(1, s) \equiv P(s)$) is [6]

$$P(n, s) = a_n s^{d_n} e^{-b_n s^2}, \quad d_n = n - 1 + \frac{1}{2} n(n + 1) \beta \quad (5)$$

and a_n, b_n are fixed by the normalisations

$$\int_0^\infty P(n, s) ds = 1, \quad \int_0^\infty s P(n, s) ds = n. \quad (6)$$

While for chaotic systems it is possible to argue that eigenstates may statistically be invariant under rotations, this is not the case for more general models. In order to describe statistical properties of such systems one has to consider non-invariant ensembles of random matrices. One of the most investigated examples is the three-dimensional Anderson model [7], with on-site disorder and nearest-neighbour coupling. Depending on the strength of the disorder, it can display metallic behaviour well described by standard random matrix ensembles, or insulator behaviour with Poisson-like spectrum. However, at the metal-insulator transition, spectral statistics are of an intermediate type and are not described by invariant ensembles [8]. Similar behaviours have been observed in pseudo-integrable billiards [9],

quantum maps corresponding to diffractive classical maps [10], or quantum Hall transitions [11]. Models have been proposed to describe such intermediate statistics [12], and random matrix ensembles which possess similar features have been constructed, e.g. power-law random banded matrix ensembles [13, 14].

The main purpose of this paper is to construct random matrix ensembles which are not invariant over rotations of eigenstates, but whose joint distributions of eigenvalues can nevertheless be calculated analytically. A short version of the paper has been published in [15]. All of these ensembles have intermediate statistics, and for certain of them spectral correlation functions, e.g. the nearest-neighbour distribution, are obtained explicitly. Eigenfunctions of these ensembles are neither localised (as for integrable systems) nor extended (as for chaotic models) but have fractal properties [16].

Random matrices of the proposed critical ensembles are constructed from the Lax matrices of classical integrable models. These models are systems of N classical particles labelled by an index i , $1 \leq i \leq N$ in a one-dimensional space. Each particle i is characterised by its position in space q_i and its momentum p_i . The dynamics of the particles is entirely described by the Hamiltonian $H(\mathbf{p}, \mathbf{q})$, where $\mathbf{p} = (p_1, \dots, p_N)$ and $\mathbf{q} = (q_1, \dots, q_N)$. The characteristic property of these models is the existence of a pair of $N \times N$ matrices L and M , called the Lax pair of the system [17], such that the equations of motion (the Hamilton equations, derived from the system Hamiltonian) are equivalent to

$$\frac{\partial L}{\partial t} = M L - L M. \quad (7)$$

The Lax matrix L is a matrix depending on momenta \mathbf{p} and coordinates \mathbf{q} . We propose to consider these Lax matrices as random matrices with a certain 'natural' measure of random variables p_j and q_j

$$dL = P(\mathbf{p}, \mathbf{q}) d^N \mathbf{p} d^N \mathbf{q}. \quad (8)$$

The explicit form of this measure depends on the system and will be discussed below. We do not impose any dynamics on variables \mathbf{p} and \mathbf{q} . The only information we use from the integrability of the underlying classical system is the existence and explicit form of action-angle variables $I_\alpha(\mathbf{p}, \mathbf{q})$ and $\phi_\alpha(\mathbf{p}, \mathbf{q})$. In particular, it is well known that the transformation from momenta and coordinates to action-angle variables is canonical, so that

$$\prod_j dp_j dq_j = \prod_\alpha dI_\alpha d\phi_\alpha. \quad (9)$$

Direct proof that the transformation is canonical is difficult in general, and implicit methods have been used to establish it for specific systems [18]-[20]. In the models we consider here, action variables turn out to be the eigenvalues λ_α of the Lax matrix, or a simple function of them. The canonical change of variables from momenta and coordinates to action-angle variables in (8) leads to a formal relation

$$dL = \mathcal{P}(\boldsymbol{\lambda}, \boldsymbol{\phi}) d^N \boldsymbol{\lambda} d^N \boldsymbol{\phi}, \quad (10)$$

where $\mathcal{P}(\boldsymbol{\lambda}, \boldsymbol{\phi}) \equiv P(\mathbf{p}(\boldsymbol{\lambda}, \boldsymbol{\phi}), \mathbf{q}(\boldsymbol{\lambda}, \boldsymbol{\phi}))$. The exact joint distribution of eigenvalues is then obtained by integration over angle variables, which can easily be performed in all cases considered, and yields

$$P(\boldsymbol{\lambda}) = \int \mathcal{P}(\boldsymbol{\lambda}, \boldsymbol{\phi}) d^N \boldsymbol{\phi}. \quad (11)$$

This scheme is general and can be adapted to many different models.

In this paper we consider in detail four typical models of N -particle classical integrable systems. The three first, labelled CM_r , CM_h , and CM_t , correspond to the rational, hyperbolic, and trigonometric Calogero-Moser models [21, 22]. The fourth model, labelled RS, is a trigonometric variant of the Ruijsenaars-Schneider model [23].

The Calogero-Moser models are defined by the Hamiltonian

$$H(\mathbf{p}, \mathbf{q}) = \frac{1}{2} \sum_j p_j^2 + g^2 \sum_{j < k} v(q_j - q_k), \quad (12)$$

where $v(\xi)$ is a potential depending on the distance between particles and g is a constant [24]. For the models considered here it has the form $v(\xi) = x^2(\xi)$, where

$$x(\xi) = \begin{cases} \frac{1}{\xi} & \text{model CM}_r \\ \frac{\mu/2}{\sinh(\mu\xi/2)} & \text{model CM}_h \\ \frac{\mu/2}{\sin(\mu\xi/2)} & \text{model CM}_t \end{cases}. \quad (13)$$

The Hamiltonian of our fourth model, the trigonometric Ruijsenaars-Schneider model, is [23]

$$H(\mathbf{p}, \mathbf{q}) = \sum_{j=1}^N \cos(p_j) \prod_{k \neq j} \left(1 - \frac{\sin^2 [\mu g/2]}{\sin^2 [\mu(q_j - q_k)/2]} \right)^{1/2} \quad \text{model RS .} \quad (14)$$

The plan of the paper is the following. Sections II, III, and IV are devoted to the construction of critical ensembles related respectively with the rational, hyperbolic, and trigonometric Calogero-Moser models. In each of these sections we briefly present the construction of the action-angle variables and choose a 'natural' measure of random momenta and coordinates which allows an easy change of variables as in (10). We then give explicit formulas for the joint distribution of eigenvalues for the resulting critical ensembles of Lax matrices. In section V this scheme is applied to the Ruijsenaars-Schneider model. For this model the joint distribution of eigenvalues takes a form which makes it suitable for the application of the transfer operator formalism. This approach is detailed in section VI, and in section VII it is applied to the analytic calculation of nearest-neighbour distributions for the RS model. The spectral compressibility for this model is obtained in section VIII.

For clarity we state below the principal results for the four models considered in this paper.

CM_r ensemble

The CM_r ensemble is defined as the ensemble of $N \times N$ Hermitian matrices of the form

$$L_{kr} = p_r \delta_{kr} + ig \frac{1 - \delta_{kr}}{q_k - q_r}, \quad (15)$$

with g a real constant. Positions \mathbf{q} and momenta \mathbf{p} are random variables distributed according to the density

$$P(\mathbf{p}, \mathbf{q}) \sim \exp \left[-A \left(\sum_j p_j^2 + g^2 \sum_{j \neq k} \frac{1}{(q_j - q_k)^2} \right) - B \sum_j q_j^2 \right], \quad (16)$$

with A and B arbitrary positive constants. The joint distribution of eigenvalues for this ensemble is then given by

$$P(\boldsymbol{\lambda}) \sim \exp \left[-A \sum_{\alpha} \lambda_{\alpha}^2 - g^2 B \sum_{\alpha \neq \beta} \frac{1}{(\lambda_{\alpha} - \lambda_{\beta})^2} \right]. \quad (17)$$

A characteristic property of this ensemble is the exponentially strong level repulsion: the nearest-neighbour spacing distribution $P(s)$ is characterised by

$$\ln P(s) \underset{s \rightarrow 0}{\sim} -\frac{b}{s^2} + \mathcal{O}(1). \quad (18)$$

We propose the following Wigner-type surmise for the next-to-nearest-neighbour spacing distributions $P(n, s)$, depending on four parameters:

$$P(n, s) = as^d \exp \left(-\frac{b}{s^2} - cs \right). \quad (19)$$

It contains two fitting constants depending on n . The other two are fixed by the normalisation (6).

CM_h ensemble

The CM_h ensemble is defined as the ensemble of $N \times N$ Hermitian matrices of the form

$$L_{kr} = p_r \delta_{kr} + ig \frac{\mu(1 - \delta_{kr})}{2 \sinh [\mu(q_k - q_r)/2]} \quad (20)$$

with g and μ real constants, and \mathbf{q} and \mathbf{p} distributed according to the density

$$P(\mathbf{p}, \mathbf{q}) \sim \exp \left[-A \left(\sum_j p_j^2 + g^2 \sum_{j \neq k} \frac{\mu^2}{4 \sinh^2 [\mu(q_j - q_k)/2]} \right) - B \sum_j \cosh \mu q_j \right]. \quad (21)$$

The exact joint distribution for this model is

$$P(\boldsymbol{\lambda}) \sim \exp \left(-A \sum_{\alpha} \lambda_{\alpha}^2 \right) \prod_{\alpha} K_0 \left(B \prod_{\beta \neq \alpha} \left| 1 + \frac{ig\mu}{\lambda_{\alpha} - \lambda_{\beta}} \right| \right) \quad (22)$$

where $K_0(x)$ is the modified Bessel function of the second kind. The nearest-neighbour spacing distribution has an exponential asymptotic similar to (18) but with $1/s$ leading term instead of $1/s^2$, namely

$$\ln P(s) \underset{s \rightarrow 0}{\sim} -\frac{b}{s} + \mathcal{O}(\ln s). \quad (23)$$

The Wigner-type surmise for CM_h is

$$P(n, s) = as^d \exp \left(-\frac{b}{s} - cs \right). \quad (24)$$

CM_t ensemble

Matrices from this ensemble correspond to a situation where μ in Eq. (20) is allowed to take pure imaginary values. They are of the form

$$L_{kr} = p_r \delta_{kr} + ig \frac{\mu(1 - \delta_{kr})}{2 \sin [\mu(q_k - q_r)/2]} \quad (25)$$

with g and μ real constants, and \mathbf{q} and \mathbf{p} distributed according to the density

$$P(\mathbf{p}, \mathbf{q}) \sim \exp \left[-A \left(\sum_j p_j^2 + g^2 \sum_{j \neq k} \frac{\mu^2}{4 \sin^2 [\mu(q_j - q_k)/2]} \right) \right] \quad (26)$$

with the restrictions that all q_j are between 0 and $2\pi/\mu$. The exact joint distribution of eigenvalues for this ensemble is

$$P(\boldsymbol{\lambda}) \sim \exp \left(-A \sum_{\alpha} \lambda_{\alpha}^2 \right) \chi(\boldsymbol{\lambda}), \quad (27)$$

where the function $\chi(\boldsymbol{\lambda})$ is equal to 1 if $\lambda_1 < \lambda_2 < \dots < \lambda_N$ and $\lambda_{\alpha+1} - \lambda_{\alpha} > \mu g$ for all α . The nearest-neighbour spacing distribution is given by a shifted Poisson distribution of the form

$$P(s) = \begin{cases} 0, & 0 < s < b \\ \frac{1}{1-b} \exp \left(-\frac{s-b}{1-b} \right), & s > b \end{cases} \quad (28)$$

with b some fitting constant.

RS ensemble

The RS ensemble is defined as the ensemble of $N \times N$ matrices of the form [23]

$$L_{kr} = e^{i\sigma p_k/2} \tilde{W}_k^{1/2} \frac{\sin [\mu g \sigma / 2]}{\sin [\mu(q_k - q_r + g\sigma)/2]} \tilde{V}_r^{1/2} e^{i\sigma p_r/2} \quad (29)$$

with

$$\tilde{V}_k = \prod_{j \neq k} \frac{\sin [\mu(q_k - q_j - g\sigma)/2]}{\sin [\mu(q_k - q_j)/2]}, \quad \tilde{W}_k = \prod_{j \neq k} \frac{\sin [\mu(q_k - q_j + g\sigma)/2]}{\sin [\mu(q_k - q_j)/2]}. \quad (30)$$

Let $\tilde{\Omega}$ be the set of \mathbf{q} such that for all k the sign of both \tilde{V}_k and \tilde{W}_k is the same as the sign of $\sin(N\mu g\sigma/2)/\sin(\mu g\sigma/2)$. The matrix L is unitary if and only if $\mathbf{q} \in \tilde{\Omega}$. The variables \mathbf{q} and \mathbf{p} are chosen to be distributed according to the uniform density in the region where L is unitary. That is, we choose momentum variables p_j independent and uniformly distributed between 0 and $2\pi/\sigma$ and coordinate variables \mathbf{q} uniformly distributed over $\tilde{\Omega}$. In this case eigenvalues of the Lax matrices (29) are also uniformly distributed over $\tilde{\Omega}$. Choosing $\mu = 2\pi/N$, $\sigma = 1$ and $g = a$, we compute correlation functions of eigenvalues of matrix (29) for fixed a and $N \rightarrow \infty$. The results strongly depend on the integer part of a . For $0 < a < 1$ the nearest-neighbour spacing distribution is similar to (28) with constant b now equal to a . For $1 < a < 2$ the nearest-neighbour distribution takes the form

$$P(s) = \begin{cases} A^2 \sinh^2(\rho s) & \text{when } 1 < g < 4/3 \\ \frac{81}{64} s^2 & \text{when } g = 4/3 \\ A^2 \sin^2(\rho s) & \text{when } 4/3 < g < 2 \end{cases}. \quad (31)$$

Constants A and ρ are determined from the normalisation conditions (4). Other correlation functions are also obtained in section VII.

Numerical implementation

The results presented above are quite robust with respect to alterations in the distribution of \mathbf{q} and \mathbf{p} . In all models considered we chose (as explained above) a distribution of coordinates such that the q_j are confined to a finite interval while having a strong repulsion between each other. As may be expected physically (though we do not have a rigorous proof for this), numerical evidence shows that, if we keep these two characteristic features, spectral properties for $N \rightarrow \infty$ depend only weakly on the precise choice for the distribution of \mathbf{q} and \mathbf{p} .

From these considerations it is thus natural to use, rather than the exact complicated distribution of \mathbf{q} , the picket-fence configuration when all coordinates are just fixed and equally spaced. As all definitions of our ensembles involve only differences $q_j - q_k$ multiplied by a parameter (μ or g , depending on the model), we can without loss of generality choose to take $q_j = j$, $j = 1, \dots, N$.

For numerical implementation we chose $q_j = j$, and p_j as independent Gaussian variables with zero mean and with variance equal 1 (CM ensembles) or independent variables uniformly distributed between 0 and 2π (RS ensemble). For concreteness, we fixed $\mu = 4\pi/N$ for CM_h and CM_t and $\mu = 2\pi/N$ for RS. For such a choice, the $N \times N$ Lax matrices take the form

$$\begin{aligned} L_{kr} &= p_k \delta_{kr} + ig \frac{1 - \delta_{kr}}{k - r}, & \text{model CM}_r \\ L_{kr} &= p_k \delta_{kr} + ig \frac{2\pi(1 - \delta_{kr})}{N \sinh [2\pi(k - r)/N]}, & \text{model CM}_h \\ L_{kr} &= p_k \delta_{kr} + ig \frac{2\pi(1 - \delta_{kr})}{N \sin [2\pi(k - r)/N]}, & \text{model CM}_t \\ L_{kr} &= e^{ip_k} \frac{1 - e^{2i\pi g}}{N(1 - e^{2i\pi(k-r+g)/N})}, & \text{model RS} \end{aligned}. \quad (32)$$

For CM_t matrices with even N , to avoid the singularity we changed $N \rightarrow N + 1$ in the above formula. As the figures in the next sections show, despite this particular choice for the distribution of \mathbf{q} and \mathbf{p} , the agreement between the computed spectral statistics and analytical formulas is remarkable.

II. RATIONAL CALOGERO-MOSER MODEL

The first model we consider is the rational Calogero-Moser model CM_r [24], characterised by the Lax matrix

$$L_{kr} = p_r \delta_{kr} + ig \frac{1 - \delta_{kr}}{q_k - q_r}. \quad (33)$$

It depends on a real constant g and on a set of $2N$ random variables p_k and q_k whose distribution will be specified later on. We are interested in eigenvalues λ_α and eigenfunctions $u_k(\alpha)$ of this matrix (here and below we will use the Greek letters to label eigenvalues and corresponding eigenfunctions)

$$\sum_{r=1}^N L_{kr} u_r(\alpha) = \lambda_\alpha u_k(\alpha) . \quad (34)$$

To construct angle-action variables let us define the new quantities

$$Q_{\alpha\beta} = \sum_k u_k^*(\alpha) q_k u_k(\beta) . \quad (35)$$

From Eq. (33) one gets

$$L_{kr} q_r - q_k L_{kr} = -ig(1 - \delta_{kr}) . \quad (36)$$

Multiplying both sides by $u_k^*(\alpha) u_r(\beta)$ and summing over k and r one gets

$$Q_{\alpha\beta}(\lambda_\alpha - \lambda_\beta) = -ig(e_\alpha^* e_\beta - \delta_{\alpha\beta}) , \quad (37)$$

where

$$e_\alpha = \sum_k u_k(\alpha) . \quad (38)$$

For $\alpha = \beta$, Eq. (37) implies that $|e_\alpha|^2 = 1$, and one can choose the overall phase of the eigenvector $u_k(\alpha)$ in such a way that $e_\alpha = 1$. Let ϕ_α be new variables defined by

$$Q_{\alpha\alpha} = \phi_\alpha . \quad (39)$$

Then from Eq. (37) we have

$$Q_{\alpha\beta} = \phi_\alpha \delta_{\alpha\beta} - ig \frac{1 - \delta_{\alpha\beta}}{\lambda_\alpha - \lambda_\beta} . \quad (40)$$

The matrix Q can be seen as the dual matrix of L , with ϕ_α playing the role of momenta and λ_α the role of positions. In [18] it was proved that there is a canonical transformation from position and momentum variables (q_k, p_k) to action and angle variables $(\lambda_\alpha, \phi_\alpha)$. Showing that the transformation is canonical is a rather technical mathematical result. However one can easily check that the new variables λ_α and ϕ_α verify Hamilton-Jacobi equations (see Appendix A).

We now consider an ensemble of Hermitian matrices of the form (33) with random variables p_k and q_k drawn according to the measure

$$P(L)dL = \mathcal{N} \exp \left[-A \text{Tr} L^2 - B \sum_k q_k^2 \right] \prod_k dp_k dq_k , \quad (41)$$

where A and B are given constants and \mathcal{N} a normalisation factor. The first term in Eq. (41) is the analog of the usual Gaussian weight of RMT; the second term is a quadratic confinement potential. Since the action-angle transformation is canonical one has

$$\prod_k dp_k dq_k = \prod_\alpha d\lambda_\alpha d\phi_\alpha . \quad (42)$$

From Eq. (35), using orthogonality of eigenvectors one gets

$$\text{Tr} Q^2 = \sum_j q_j^2 . \quad (43)$$

Using these relations one can rewrite the distribution (41) in action-angle variables λ_α and ϕ_α as

$$P(L)dL = \mathcal{N} \exp \left[-A \sum_\alpha \lambda_\alpha^2 - B \left(\sum_\alpha \phi_\alpha^2 + g^2 \sum_{\alpha \neq \beta} \frac{1}{(\lambda_\alpha - \lambda_\beta)^2} \right) \right] \prod_\alpha d\lambda_\alpha d\phi_\alpha . \quad (44)$$

Integration over the ϕ_α gives a constant. We thus obtain the joint distribution of eigenvalues for the ensemble of random matrices L with the measure (41) as

$$P(\lambda_1, \dots, \lambda_N) \sim \exp \left[-A \sum_{\alpha} \lambda_{\alpha}^2 - Bg^2 \sum_{\alpha \neq \beta} \frac{1}{(\lambda_{\alpha} - \lambda_{\beta})^2} \right]. \quad (45)$$

Note that, similarly as in the standard RMT case, this joint eigenvalue distribution can be interpreted via the Coulomb gas model as the partition function of an ensemble of particles on a line, here with inverse square repulsion. After rescaling $x_k = \lambda_k (Bg^2/A)^{-1/4}$, equilibria positions of the particles at positions λ_α are given by

$$x_k = 2 \sum_{j \neq k} \frac{1}{(x_j - x_k)^3}, \quad 1 \leq k \leq N. \quad (46)$$

Such a relation characterises the zeros of Hermite polynomials of degree N (see also Eq. (10.3) of [24]). It is known from RMT [4] that the distribution of eigenvalues of Gaussian random ensembles has a similar property, which implies that the asymptotic density of eigenvalues is given by Wigner's semi-circle law.

An immediate consequence of the distribution (45) is the unusual very strong level repulsion at small distances. For all standard random matrix ensembles the nearest-neighbour distribution $P(s)$ behaves as s^β at small s . By contrast, in our case it follows from (45) that

$$P(s) \underset{s \rightarrow 0}{\sim} a e^{-b/s^2}. \quad (47)$$

As the potential between eigenvalues decreases as the inverse square of the distance between them, the probability of having a gap of size s for large s is exponentially small. We could not calculate exactly correlation functions for the distribution (45). However, combining the two asymptotics above, we build a Wigner-type surmise for the nearest-neighbour spacing distribution of the form

$$P(s) = a e^{-b/s^2 - cs}, \quad (48)$$

where b is a fitting constant, and constants a and c are determined from the normalisation conditions (4). For the n th nearest-neighbour spacing distributions $P(n, s)$, with $n \geq 2$, we conjecture, by analogy with the Wigner surmise (5) for standard random matrices, the form

$$P(n, s) = a s^d e^{-b/s^2 - cs} \quad (49)$$

with two fitting constants b and d .

To assess this conjecture we compare the analytical expressions (48)–(49) with numerical results, with the choice of parameters detailed in section I. Results displayed in Fig. 1 show that the agreement is remarkable.

III. HYPERBOLIC CALOGERO-MOSER MODEL

The Lax matrix for the hyperbolic Calogero–Moser model CM_h reads [24]

$$L_{kr} = p_r \delta_{kr} + ig(1 - \delta_{kr}) \frac{\mu}{2 \sinh(\mu(q_k - q_r)/2)}. \quad (50)$$

Let us define two matrices Q and R by

$$Q_{\alpha\beta} = \sum_k u_k^*(\alpha) e^{\mu q_k} u_k(\beta), \quad R_{\alpha\beta} = \sum_k u_k^*(\alpha) e^{-\mu q_k} u_k(\beta), \quad (51)$$

and two vectors

$$e_\alpha = \sum_k u_k(\alpha) e^{\mu q_k/2}, \quad f_\alpha = \sum_k u_k(\alpha) e^{-\mu q_k/2}. \quad (52)$$

From (50) one can get the two equivalent equations

$$e^{\mu q_k} L_{kr} - L_{kr} e^{\mu q_r} = ig\mu(1 - \delta_{kr}) e^{\mu(q_k + q_r)/2} \quad (53)$$

$$e^{-\mu q_r} L_{kr} - L_{kr} e^{-\mu q_k} = ig\mu(1 - \delta_{kr}) e^{-\mu(q_k + q_r)/2}. \quad (54)$$

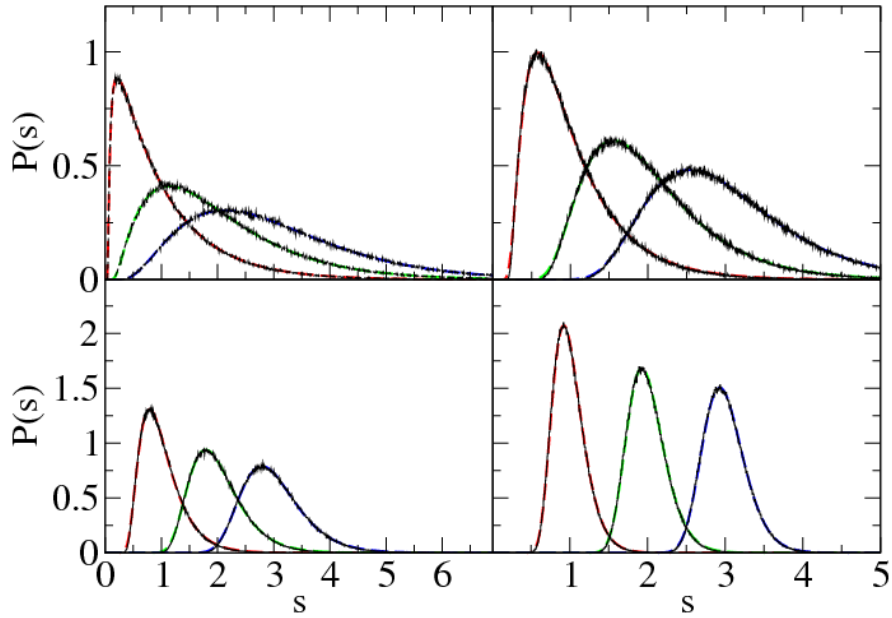


FIG. 1. (Color online). Nearest-neighbour distributions $P(n, s)$ for the random matrices of model CM_r , with $g = 0.05$ (top left), 0.25 (top right), 0.5 (bottom left), and 1 (bottom right), averaged over the central quarter of the unfolded spectrum for 8000 realisations of matrices of size $N = 512$. Solid lines are numerical results, dashed lines indicate the fits (48) for and (49) for $P(n, s)$ with (in each panel from left to right) $P(s) = P(1, s)$ (red), $P(2, s)$ (green) and $P(3, s)$ (blue).

Multiplying both sides by $u_k(\alpha)^* u_r(\beta)$ and summing over all k and r one gets

$$Q_{\alpha\beta}(\lambda_\alpha - \lambda_\beta) = -ig\mu(e_\alpha^* e_\beta - Q_{\alpha\beta}) \quad (55)$$

$$R_{\alpha\beta}(\lambda_\alpha - \lambda_\beta) = ig\mu(f_\alpha^* f_\beta - R_{\alpha\beta}), \quad (56)$$

which implies that matrices Q and R take the form

$$Q_{\alpha\beta} = e_\alpha^* \frac{ig\mu}{\lambda_\beta - \lambda_\alpha + ig\mu} e_\beta, \quad R_{\alpha\beta} = f_\alpha^* \frac{-ig\mu}{\lambda_\beta - \lambda_\alpha - ig\mu} f_\beta. \quad (57)$$

By their definition (51), matrices Q and R are inverse of each other, so that $\sum_\gamma Q_{\alpha\gamma} R_{\gamma\beta} = \delta_{\alpha\beta}$ for all α, β . For $\alpha = \beta$ this condition implies that

$$-g^2 \mu^2 e_\alpha^* f_\alpha \sum_\gamma \frac{e_\gamma f_\gamma^*}{(\lambda_\gamma - \lambda_\alpha + ig\mu)^2} = 1 \quad (58)$$

(in particular, it follows that all e_α and f_α are nonzero). For $\alpha \neq \beta$ one obtains

$$\sum_\gamma \frac{e_\gamma f_\gamma^*}{(\lambda_\beta - \lambda_\alpha + ig\mu)(\lambda_\gamma - \lambda_\beta - ig\mu)} = 0. \quad (59)$$

Using the identity

$$\frac{1}{(\lambda_\beta - \lambda_\alpha + ig\mu)(\lambda_\gamma - \lambda_\beta - ig\mu)} = \left[\frac{1}{\lambda_\beta - \lambda_\gamma + ig\mu} - \frac{1}{\lambda_\beta - \lambda_\alpha + ig\mu} \right] \frac{1}{\lambda_\alpha - \lambda_\gamma}, \quad (60)$$

valid for $\alpha \neq \beta$, one concludes that

$$\sum_\gamma \frac{e_\gamma f_\gamma^*}{\lambda_\gamma - \lambda_\alpha + ig\mu} = c, \quad (61)$$

where c is a certain constant independent on α . According to this equation the quantities $b_\gamma = e_\gamma f_\gamma^*/c$ obey a system of linear equations of the form

$$\sum_\gamma \frac{b_\gamma}{x_\gamma - y_\alpha} = 1, \quad (62)$$

with $x_\gamma = \lambda_\gamma$ and $y_\alpha = \lambda_\alpha - ig\mu$. This equation coincides with Eq. (B1) in Appendix B. From (B2) it follows that

$$e_\alpha f_\alpha^* = ig\mu c V_\alpha, \quad (63)$$

where

$$V_\alpha = \prod_{\beta \neq \alpha} \left(1 + \frac{ig\mu}{\lambda_\alpha - \lambda_\beta} \right), \quad (64)$$

while (B4) implies that

$$\sum_\alpha e_\alpha f_\alpha^* = ig\mu c N. \quad (65)$$

It readily follows from the definition (52) of e_α and f_α that $\sum_\alpha e_\alpha f_\alpha^* = N$, thus the value of c is fixed by $ig\mu c = 1$. Equation (58) is then fulfilled as a direct consequence of (B3). Finally we have

$$e_\alpha f_\alpha^* = V_\alpha. \quad (66)$$

Let ϕ_α be new variables defined from diagonal elements of matrix Q by

$$Q_{\alpha\alpha} = |V_\alpha| e^{\mu\phi_\alpha}. \quad (67)$$

Then from (57) and (66) it follows that

$$R_{\alpha\alpha} = |V_\alpha| e^{-\mu\phi_\alpha}. \quad (68)$$

In the definition (52) of e_α it is convenient to choose the overall phase of the eigenvector $u_k(\alpha)$ in such a way that e_α be real. As $Q_{\alpha\alpha} = |e_\alpha|^2$ one has

$$e_\alpha = |V_\alpha|^{1/2} e^{\mu\phi_\alpha/2}. \quad (69)$$

Using Eq. (57), the matrix Q can now be expressed in terms of the new variables λ_α and ϕ_α , as

$$Q_{\alpha\beta} = |V_\alpha|^{1/2} e^{\mu\phi_\alpha/2} \frac{ig\mu}{\lambda_\beta - \lambda_\alpha + ig\mu} e^{\mu\phi_\beta/2} |V_\beta|^{1/2}. \quad (70)$$

As in the case of model CM_r , the matrix Q can be seen as the dual matrix of L . Indeed, Q coincides with the Lax matrix of the rational Ruijsenaars-Schneider model with coordinates λ_α and momenta ϕ_α [18]. In [18] it has been proved that the transformation from position and momentum variables (q_k, p_k) to action-angle variables $(\lambda_\alpha, \phi_\alpha)$ is canonical. Again one can check that the new variables λ_α and ϕ_α verify Hamilton-Jacobi equations.

We now consider an ensemble of Hermitian matrices of the form (50) with random variables p_k and q_k drawn according to the measure

$$P(L)dL = \mathcal{N} \exp \left[-A \text{Tr} L^2 - B \sum_k \cosh \mu q_k \right] \prod_k dp_k dq_k. \quad (71)$$

As in the case of model CM_r , Eq. (71) contains a standard RMT Gaussian weight and a confinement potential which can be rewritten as $\text{Tr} Q + \text{Tr} R$. Using (67), (68), and the fact that the transformation is canonical, we get the distribution in terms of the new variables λ_α and ϕ_α as

$$P(L)dL = \mathcal{N} \exp \left[-A \sum_\alpha \lambda_\alpha^2 - B \sum_\alpha |V_\alpha| \cosh \mu \phi_\alpha \right] \prod_\alpha d\lambda_\alpha d\phi_\alpha. \quad (72)$$

The joint distribution of eigenvalues is then obtained by integrating over the angle variables, using

$$\int_{-\infty}^{\infty} \exp[-B|V_\alpha| \cosh \mu\phi_\alpha] d\phi_\alpha = \frac{1}{\mu} K_0(B|V_\alpha|) \quad (73)$$

where K_0 is the modified Bessel function of the second kind. This yields the joint distribution of eigenvalues for model CM_h as

$$P(\lambda_1, \dots, \lambda_N) \sim \exp\left(-A \sum_{\alpha} \lambda_{\alpha}^2\right) \prod_{\alpha} K_0\left(B \prod_{\beta \neq \alpha} \left|1 + \frac{ig\mu}{\lambda_{\alpha} - \lambda_{\beta}}\right|\right). \quad (74)$$

This expression is exact but difficult to handle. In order to find a Wigner-type surmise for the nearest-neighbour distributions we consider the limiting behaviour $P(\boldsymbol{\lambda})$ when two nearby eigenvalues λ_1 and λ_2 get close to each other. Setting $s = \lambda_1 - \lambda_2$ we see that the factor $\exp(-b/s^2)$ in the case of model CM_r is replaced by a factor

$$K_0\left(B\sqrt{1 + \frac{g^2\mu^2}{s^2}}\right)^2 \underset{s \rightarrow 0}{\sim} s \exp\left(-\frac{2Bg\mu}{s}\right). \quad (75)$$

We therefore expect the nearest-neighbour spacing distribution to behave as

$$P(n, s) = as^d \exp(-b/s - cs). \quad (76)$$

In Fig. 2 we show the results of numerical computations of the nearest-neighbour spacing distributions for matrices of the form (50) with the choice of parameters and variables detailed in section I. The surmise (76) perfectly reproduces numerical results.

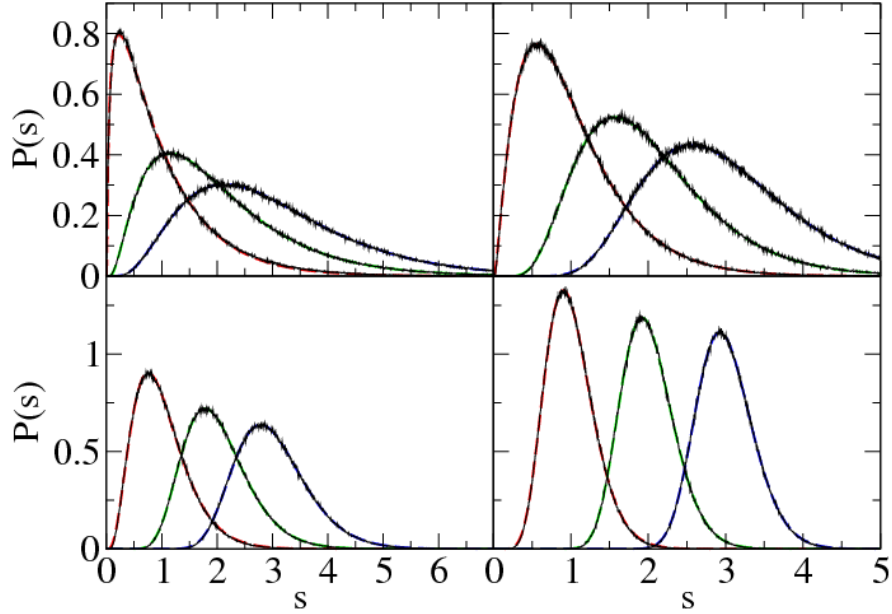


FIG. 2. (Color online). Nearest-neighbour spacing distributions $P(n, s)$ for the random matrices of model CM_h with $\mu = 4\pi/N$ and $g = 0.05$ (top left), 0.25 (top right), 0.5 (bottom left), and 1 (bottom right), averaged over the central quarter of the spectrum for 32000 realisations of matrices of size $N = 256$. Solid lines are numerical results, dashed lines indicate the fit (76) for $P(n, s)$ with (in each panel from left to right) $P(s) = P(1, s)$ (red), $P(2, s)$ (green) and $P(3, s)$ (blue).

In order to assess better the validity of the exponentially strong level repulsion for models CM_r and CM_h , we compare in Fig. 3 the beginning of the distributions $P(s)$ for these models. Clearly the $1/s^2$ repulsion for CM_r and the $1/s$ repulsion for CM_h fit numerical curves very well. However, the precision of our numerical results does not permit to confirm or reject the presence of the logarithmic term $d \ln s$ in $P(s)$ for CM_h model.

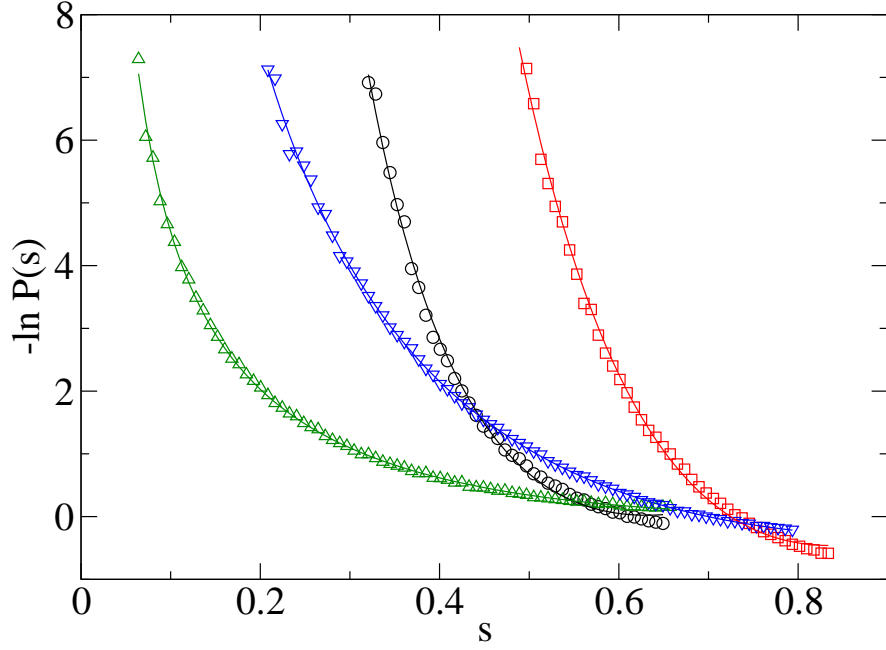


FIG. 3. (Color online). Nearest-neighbour spacing distributions $P(s)$ for the random matrices CM_r ($g = 0.5$, black circles) and $g = 1$, red squares) and CM_h ($g = .5$, green triangles up, and $g = 1$, blue triangles down), with $\mu = 4\pi/N$. Symbols are numerical results, solid lines indicate the fit $b/s^2 + cs - \ln a$ (model CM_r , Eq. (48)) and $b/s + cs - \ln a - d \ln s$ (model CM_h , Eq. (76)). Logarithm is natural.

IV. TRIGONOMETRIC CALOGERO-MOSER MODEL

For the trigonometric Calogero-Moser model CM_t the Lax matrix is [24]

$$L_{kr} = p_r \delta_{kr} + ig(1 - \delta_{kr}) \frac{\mu}{2 \sin(\mu(q_k - q_r)/2)}. \quad (77)$$

The only difference with the previous model, (50), is the sin function which replaces the sinh. The matrix (77) can be obtained from (50) by the substitution $\mu \rightarrow i\mu$. However the fact that positions of the particles are now defined on a circle (because of the sin function) makes the resulting spectral statistics entirely different from the previous models CM_r and CM_h .

To construct action-angle variable we introduce, as in the previous section, two matrices

$$Q_{\alpha\beta} = \sum_k u_k^*(\alpha) e^{i\mu q_k} u_k(\beta), \quad R_{\alpha\beta} = \sum_k u_k^*(\alpha) e^{-i\mu q_k} u_k(\beta), \quad (78)$$

and two vectors

$$e_\alpha = \sum_k u_k(\alpha) e^{i\mu q_k/2}, \quad f_\alpha = \sum_k u_k(\alpha) e^{-i\mu q_k/2}. \quad (79)$$

Following the same steps as above with μ replaced by $i\mu$, one gets

$$Q_{\alpha\beta} = f_\alpha^* \frac{g\mu}{\lambda_\alpha - \lambda_\beta + g\mu} e_\beta, \quad R_{\alpha\beta} = e_\alpha^* \frac{g\mu}{\lambda_\beta - \lambda_\alpha + g\mu} f_\beta. \quad (80)$$

Again, using the fact that Q is the inverse of R we obtain that

$$\sum_\gamma \frac{|e_\gamma|^2}{\lambda_\gamma - \lambda_\alpha - g\mu} = c_1, \quad \sum_\gamma \frac{|f_\gamma|^2}{\lambda_\gamma - \lambda_\alpha + g\mu} = c_2 \quad (81)$$

with certain constants c_1 and c_2 independent on α . Repeating the same arguments as in the previous section and using results of Appendix B one concludes that $c_2 = -c_1 = 1/(\mu g)$ and

$$|e_\alpha|^2 = V_\alpha, \quad |f_\alpha|^2 = W_\alpha, \quad (82)$$

with

$$V_\alpha = \prod_{\beta \neq \alpha} \left(1 - \frac{g\mu}{\lambda_\alpha - \lambda_\beta} \right), \quad W_\alpha = \prod_{\beta \neq \alpha} \left(1 + \frac{g\mu}{\lambda_\alpha - \lambda_\beta} \right). \quad (83)$$

The new variables ϕ_α are defined as above from the diagonal elements of matrix Q as follows

$$Q_{\alpha\alpha} = V_\alpha^{1/2} W_\alpha^{1/2} e^{i\mu\phi_\alpha}. \quad (84)$$

In the definition (79) of e_α one has a freedom to choose the overall phase of the eigenvector $u_k(\alpha)$. Since from (80) one must have $Q_{\alpha\alpha} = f_\alpha^* e_\alpha$, one can choose phases, for example, as follows

$$e_\alpha = V_\alpha^{1/2} e^{i\mu\phi_\alpha/2}, \quad f_\alpha = W_\alpha^{1/2} e^{-i\mu\phi_\alpha/2}. \quad (85)$$

Then the matrix Q can be expressed in terms of new variables λ_α and ϕ_α as

$$Q_{\alpha\beta} = e^{i\mu\phi_\alpha/2} W_\alpha^{1/2} \frac{g\mu}{\lambda_\alpha - \lambda_\beta + g\mu} V_\beta^{1/2} e^{i\mu\phi_\beta/2}. \quad (86)$$

The inverse matrix R plays a symmetric role, as it is obtained from Q by exchanging μ to $-\mu$. Again, there is a canonical transformation from position and momentum variables (q_k, p_k) to action and angle variables $(\lambda_\alpha, \phi_\alpha)$ [18].

An important consequence of (82) is that for all α we should have

$$V_\alpha > 0, \quad W_\alpha > 0. \quad (87)$$

These inequalities impose non-trivial restrictions on eigenvalues λ_α , as we will see now. We label eigenvalues so that $\lambda_1 < \lambda_2 < \dots < \lambda_N$, and we consider the function

$$h(x) = \sum_\gamma \frac{V_\alpha}{\lambda_\gamma - x} + \frac{1}{\mu g}. \quad (88)$$

It has N poles at $x = \lambda_\gamma$, and according to Eqs. (81) and (85) it has N zeros at $x = \lambda_\alpha + \mu g$. If all numerators V_α are positive then the derivative of h is positive, and it is easy to check from the graph of the function that between two consecutive poles there is one and only one zero. Suppose $\mu g > 0$. Then for $x \rightarrow -\infty$ the function $h(x)$ has a strictly positive limit. The lowest zero $\lambda_1 + \mu g$ must thus lie in the interval $]\lambda_1, \lambda_2[$. More generally one must have $\lambda_\alpha + \mu g \in]\lambda_\alpha, \lambda_{\alpha+1}[$ for $1 \leq \alpha \leq N-1$, while the largest zero $\lambda_N + \mu g$ lies in the interval $]\lambda_N, \infty[$. Thus eigenvalues fulfill the inequalities

$$\lambda_{\alpha+1} - \lambda_\alpha > g\mu. \quad (89)$$

Conversely, if these inequalities are fulfilled then trivially all V_α are positive. Therefore, (89) are the necessary and sufficient conditions for the positivity of all V_α . In particular, eigenvalues of the Lax matrix (77) obey inequalities (89) for any choice of the \mathbf{q} . These results adapt straightforwardly to the case where μg is negative.

Since in (77) the q_k only appear as an argument in the sin function, there is no need to choose a confining potential for the particle distribution as in models CM_r and CM_h . We consider the probability distribution of \mathbf{p} and \mathbf{q} in the form

$$P(\mathbf{p}, \mathbf{q}) \sim \exp \left[-A \left(\sum_j p_j^2 + g^2 \sum_{j \neq k} \frac{\mu^2}{4 \sin^2 [\mu(q_k - q_r)/2]} \right) \right] \quad (90)$$

with the restrictions that all q_j are between 0 and $2\pi/\mu$. Since the change of variables from \mathbf{p} and \mathbf{q} to λ_α and ϕ_α is canonical and the restrictions (89) do not depend on phase variables the joint distribution of eigenvalues is

$$P(\lambda_1, \dots, \lambda_N) \sim \exp \left(-A \sum_\alpha \lambda_\alpha^2 \right) \chi(\lambda_1, \dots, \lambda_N), \quad (91)$$

where the function $\chi(\boldsymbol{\lambda})$ is equal to 1 if (89) is fulfilled for all α , and 0 otherwise.

It turns out that model CM_t is very similar to a fourth model, the Ruijsenaars-Schneider model, that we will consider in the next section. Therefore we postpone analytical calculations of the nearest-neighbour spacing distributions to section VI. The nearest-neighbour spacing distributions $P(n, s)$ are shifted Poisson distributions of the form

$$P(n, s) = \begin{cases} 0, & 0 < s < nb \\ \frac{(s - nb)^{n-1}}{(n-1)!(1-b)^n} e^{-(s-nb)/(1-b)}, & s > nb \end{cases} \quad (92)$$

with some numerical constant b . In Fig. 4 we show the results of numerical computations for matrices of the form (77) with the choice of parameters and variables detailed in section I.

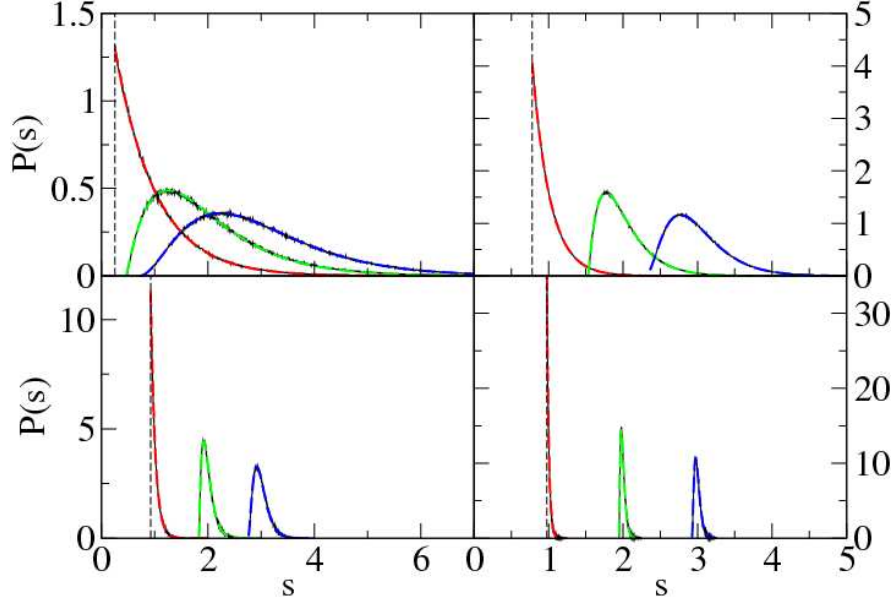


FIG. 4. (Color online). Same as Fig. 2 for model CM_t with $P(n, s)$ the shifted Poisson distribution (92).

V. RUIJSENAARS-SCHNEIDER MODEL

The Calogero-Moser models considered in the previous sections are such that there exists a matrix Q which is, in a certain sense, dual to the Lax matrix L . Namely, the canonical transformation from variables (p_k, q_k) to action-angle variables $(\lambda_\alpha, \phi_\alpha)$ is such that the action variables λ_α are eigenvalues of L and angle variables ϕ_α are related to $Q_{\alpha\alpha}$ in a simple way. In fact, the matrix Q is also a Lax matrix, corresponding to a possibly different Hamiltonian, and L plays the role of a matrix dual to Q for the inverse of the canonical transformation [18]. The rational Calogero-Moser system is self-dual since matrices L and Q , given by (33) and (40), are equal up to labelling of the variables.

The model we consider in this section is related to the above models in that its Lax matrix L and the dual matrix Q are a kind of generalisation of those of model CM_t (86). The treatment of this model closely follows the previous section.

The Lax matrix for Ruijsenaars-Schneider model is the $N \times N$ unitary matrix given by [20]

$$L_{kr} = e^{i\sigma p_k/2} \tilde{W}_k^{1/2} \frac{\sin[\mu g \sigma/2]}{\sin[\mu(q_k - q_r + g\sigma)/2]} \tilde{V}_r^{1/2} e^{i\sigma p_r/2}. \quad (93)$$

with

$$\tilde{V}_k = \prod_{j \neq k} \frac{\sin[\mu(q_k - q_j - g\sigma)/2]}{\sin[\mu(q_k - q_j)/2]}, \quad \tilde{W}_k = \prod_{j \neq k} \frac{\sin[\mu(q_k - q_j + g\sigma)/2]}{\sin[\mu(q_k - q_j)/2]}. \quad (94)$$

This Lax matrix is related with the Hamiltonian (14) by

$$H(\mathbf{p}, \mathbf{q}) = \frac{1}{2} \text{Tr}(L + L^\dagger) . \quad (95)$$

It is convenient to introduce the vectors

$$\tilde{e}_k = \tilde{V}_k^{1/2} e^{i\sigma p_k/2} e^{i\mu q_k/2} , \quad \tilde{f}_k = \tilde{W}_k^{1/2} e^{i\sigma p_k/2} e^{-i\mu q_k/2} , \quad (96)$$

so that matrix L can be rewritten as

$$L_{kr} = \tilde{f}_k^* e^{-i\mu q_k/2} \frac{\sin [\mu g \sigma / 2]}{\sin [\mu (q_k - q_r + g \sigma) / 2]} e^{-i\mu q_r/2} \tilde{e}_r . \quad (97)$$

As in the previous section, the condition that the matrix (93) is unitary imposes certain restrictions on the coordinates \mathbf{q} , which we will discuss later. Assuming that the Lax matrix is unitary, we choose to denote its eigenvalues by $e^{i\sigma \lambda_\alpha}$. The dual matrices for the Ruijsenaars-Schneider model are defined by [20]

$$Q_{\alpha\beta} = \sum_k u_k^*(\alpha) e^{i\mu q_k} u_k(\beta) , \quad R_{\alpha\beta} = Q_{\beta\alpha}^* = \sum_k u_k^*(\alpha) e^{-i\mu q_k} u_k(\beta) , \quad (98)$$

and the vectors e_α and f_α by

$$e_\alpha = \sum_k u_k(\alpha) \tilde{e}_k , \quad f_\alpha = \sum_k u_k(\alpha) \tilde{f}_k . \quad (99)$$

From Eq. (97) one has

$$e^{i\mu(q_k+q_r)/2} L_{kr} \sin [\mu(q_k - q_r + g \sigma) / 2] = \tilde{f}_k^* \sin [\mu g \sigma / 2] \tilde{e}_r . \quad (100)$$

Multiplying this expression by $u_k^*(\alpha) u_r(\beta)$ and summing both sides over k and r leads to

$$\frac{1}{2i} Q_{\alpha\beta} \left(e^{i\sigma \lambda_\beta + i\sigma \mu g / 2} - e^{i\sigma \lambda_\alpha - i\sigma \mu g / 2} \right) = f_\alpha^* \sin [\mu g \sigma / 2] e_\beta , \quad (101)$$

which yields the analogue of (80),

$$Q_{\alpha\beta} = f_\alpha^* e^{-i\sigma \lambda_\alpha / 2} \frac{\sin [\sigma g \mu / 2]}{\sin [\sigma (\lambda_\beta - \lambda_\alpha + g \mu) / 2]} e^{-i\sigma \lambda_\beta / 2} e_\beta . \quad (102)$$

Let us rewrite

$$Q_{\alpha\beta} = f_\alpha^* \frac{e^{-i\tau} - 1}{e^{i(\sigma \lambda_\alpha - \tau)} - e^{i\sigma \lambda_\beta}} e_\beta , \quad (103)$$

where we have set $\tau = \sigma \mu g$. From its definition (98) it is clear that Q has to be an unitary matrix, i.e. $\sum_\gamma Q_{\alpha\gamma} Q_{\beta\gamma}^* = \delta_{\alpha\beta}$. Selecting terms with $\alpha = \beta$ and $\alpha \neq \beta$ yields the two equations

$$|\rho|^2 |f_\alpha|^2 \sum_\gamma \frac{|e_\gamma|^2}{|e^{i(\sigma \lambda_\alpha - \tau)} - e^{i\sigma \lambda_\gamma}|^2} = 1 , \quad (104)$$

$$\sum_\gamma \frac{|e_\gamma|^2}{(e^{i(\sigma \lambda_\alpha - \tau)} - e^{i\sigma \lambda_\gamma})(e^{-i(\sigma \lambda_\beta - \tau)} - e^{-i\sigma \lambda_\gamma})} = 0 , \quad (105)$$

where we have set $\rho = e^{-i\tau} - 1$. Using the identity

$$\frac{1}{(e^{i(\sigma \lambda_\alpha - \tau)} - e^{i\sigma \lambda_\gamma})(e^{-i(\sigma \lambda_\beta - \tau)} - e^{-i\sigma \lambda_\gamma})} = \frac{e^{i(\sigma \lambda_\beta - \tau)}}{e^{i(\sigma \lambda_\alpha - \tau)} - e^{i\sigma \lambda_\beta - \tau}} \left(\frac{e^{i\sigma \lambda_\gamma}}{e^{i(\sigma \lambda_\alpha - \tau)} - e^{i\sigma \lambda_\gamma}} + \frac{e^{i\sigma \lambda_\gamma}}{e^{i\sigma \lambda_\gamma} - e^{i(\sigma \lambda_\beta - \tau)}} \right) , \quad (106)$$

it follows from (105) that there exists a constant c such that

$$\sum_{\gamma} \frac{|e_{\gamma}|^2 e^{i\sigma\lambda_{\gamma}}}{e^{i\sigma\lambda_{\gamma}} - e^{i(\sigma\lambda_{\alpha} - \tau)}} = c. \quad (107)$$

Equations (B1)–(B2) allow to obtain

$$|e_{\alpha}|^2 = -c\rho V_{\alpha} e^{-i\tau(N-1)/2}, \quad (108)$$

while from Eq. (104) one obtains, using Eq. (B3)

$$|f_{\alpha}|^2 = -\frac{1}{c\rho} W_{\alpha} e^{i\tau(N-1)/2}, \quad (109)$$

with new vectors V_{α} and W_{α} defined by

$$V_{\alpha} = \prod_{\beta \neq \alpha} \frac{\sin[\sigma(\lambda_{\alpha} - \lambda_{\beta} + g\mu)/2]}{\sin[\sigma(\lambda_{\alpha} - \lambda_{\beta})/2]}, \quad W_{\alpha} = \prod_{\beta \neq \alpha} \frac{\sin[\sigma(\lambda_{\alpha} - \lambda_{\beta} - g\mu)/2]}{\sin[\sigma(\lambda_{\alpha} - \lambda_{\beta})/2]}. \quad (110)$$

Using (B5) we get

$$\sum_{\alpha} |e_{\alpha}|^2 = c(1 - e^{-i\tau N}). \quad (111)$$

There is some overall freedom in the definition of vectors \tilde{e}_k and \tilde{f}_k in Eq. (96), as one could multiply \tilde{e}_{α} by some constant factor and divide \tilde{f}_{α} by the same factor. This in turn entails the same freedom for vectors e_{α} and f_{α} in (99). If one chooses for instance $\sum_j |\tilde{e}_j|^2 = t$ then from unitarity of the transformation in (99) one has $\sum_{\alpha} |e_{\alpha}|^2 = t$, which fixes the value

$$c = \frac{t}{1 - e^{-i\tau N}} \quad (112)$$

and thus

$$|e_{\alpha}|^2 = \frac{t \sin(\tau/2)}{\sin(N\tau/2)} V_{\alpha}, \quad |f_{\alpha}|^2 = \frac{\sin(N\tau/2)}{t \sin(\tau/2)} W_{\alpha}. \quad (113)$$

By definition $t = \sum_j |\tilde{e}_j|^2$ is positive. A convenient choice is to take

$$t = \left| \frac{\sin(N\tau/2)}{\sin(\tau/2)} \right|. \quad (114)$$

Since $|e_{\alpha}|^2$ and $|f_{\alpha}|^2$ are non-negative one concludes that V_{α} and W_{α} have the same sign as $\sin(N\tau/2)/\sin(\tau/2)$ for all α . The choice (114) implies that

$$|e_{\alpha}|^2 = |V_{\alpha}|, \quad |f_{\alpha}|^2 = |W_{\alpha}|. \quad (115)$$

The new action variables are the λ_{α} , and angle variables ϕ_{α} are defined by

$$Q_{\alpha\alpha} = (V_{\alpha} W_{\alpha})^{1/2} e^{i\mu\phi_{\alpha}}. \quad (116)$$

Equation (102) implies that $Q_{\alpha\alpha} = f_{\alpha}^* e_{\alpha} e^{-i\sigma\lambda_{\alpha}}$. In view of (115) and (116) one can choose the phases of the eigenvectors $u_k(\alpha)$ such that e_{α} and f_{α} defined by (99) can be expressed as

$$e_{\alpha} = V_{\alpha}^{1/2} e^{i\mu\phi_{\alpha}/2} e^{i\sigma\lambda_{\alpha}/2}, \quad f_{\alpha} = W_{\alpha}^{1/2} e^{-i\mu\phi_{\alpha}/2} e^{-i\sigma\lambda_{\alpha}/2}. \quad (117)$$

In terms of the new variables, Q thus reads

$$Q_{\alpha\beta} = e^{i\mu\phi_{\alpha}/2} W_{\alpha}^{1/2} \frac{\sin[\sigma g\mu/2]}{\sin[\sigma(\lambda_{\beta} - \lambda_{\alpha} + g\mu)/2]} V_{\beta}^{1/2} e^{i\mu\phi_{\beta}/2}. \quad (118)$$

Comparing (93) and (118) one sees that for model RS the dual matrix matrix Q is obtained from L by changing $\sigma \leftrightarrow \mu$, $g \rightarrow -g$, $\mathbf{p} \rightarrow \phi$, and $\mathbf{q} \rightarrow \lambda$. It means that this model is self-dual (the matrices L and Q are the same up to a change of notation). One can show that the transformation from (q_k, p_k) to action and angle variables $(\lambda_\alpha, \phi_\alpha)$ is canonical [20]. As mentioned, a consequence of the unitarity of the matrix Q is that V_α and W_α have the same sign as $\sin(N\tau/2)/\sin(\tau/2)$ for all α . This implies that certain inequalities have to be verified by the λ_α , which we now derive in a way similar as in the previous section.

Let us define the function

$$h(x) = \sum_{\gamma} V_{\gamma} \cot [(x - \sigma\lambda_{\gamma})/2] - t \cot(N\tau/2), \quad (119)$$

which is periodic with period 2π and can be considered as a function on the unit circle. It has N poles at $x = \sigma\lambda_{\gamma}$. Taking the imaginary part of Eq. (107), with c given by (112) and e_{α} given by (117), one obtains that $h(x)$ has N zeros at $x = \sigma\lambda_{\gamma} - \tau$. When all V_{γ} are positive, the same arguments as in the previous section imply that between two consecutive poles of $h(x)$ there must be exactly one zero. It means that between two nearby eigenvalues $\sigma\lambda_{\alpha}$ there is one and only one number of the form $\sigma\lambda_{\gamma} - \tau$. A similar reasoning starting from matrix $R = Q^{\dagger}$ leads to the conclusion that between two consecutive $\sigma\lambda_{\gamma}$ there must also be one and only one number of the form $\sigma\lambda_{\alpha} + \tau$. These two conditions (shift by $+\tau$ or $-\tau$) are equivalent, thus we can restrict ourselves to a shift by $+\tau$, i.e. the condition that the sets $\{\sigma\lambda_{\gamma}, 1 \leq \gamma \leq N\}$ and $\{\sigma\lambda_{\gamma} + \tau, 1 \leq \gamma \leq N\}$ intertwine on the unit circle. It is a necessary and sufficient condition for the matrix (102) to be an unitary matrix. A similar conclusion is readily obtained when all V_{γ} are negative.

The conditions implied by this type of intertwining have been discussed in [27, 28]. There is an fundamental difference between these eigenvalue conditions in the RS model and in the CM_t model discussed in the previous section. In the CM_t case the Lax matrix is Hermitian, and the λ_{α} can take values on the whole real axis. In the RS case, as the Lax matrix is unitary, the $\sigma\lambda_{\alpha}$ lie on the unit circle. Therefore poles and zeros cannot be ordered in a simple way as in the previous case, and the analysis of the previous section does not apply.

For completeness we shortly repeat the arguments of the papers [27, 28]. Let us put all eigenvalues $\sigma\lambda_{\alpha}$ of the Lax matrix (93) on the unit circle and divide the circle into sectors with angle τ . Denote the (positive) angular distance from the boundaries of the k th sector in counter clockwise direction by x_k and in clockwise direction by y_k , as in Fig. 5. After a shift by τ , the intertwining relations imply that only one of the two points corresponding to x_k and y_k

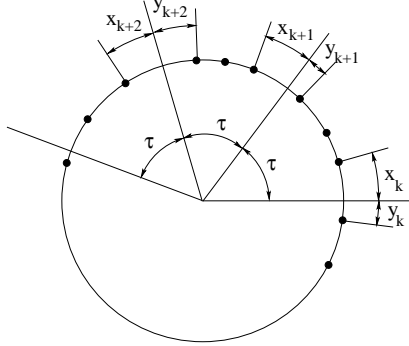


FIG. 5. Division of the circle into sectors of angle τ . Only sectors numbers k , $k+1$, and $k+2$ are indicated. Black circles indicate the positions of the $\sigma\lambda_{\alpha}$, where $\exp(i\sigma\lambda_{\alpha})$ are the eigenvalues of the RS Lax matrix.

will fall in-between points corresponding to x_{k+1} and y_{k+1} . The first case corresponds to $x_k > x_{k+1}$ and $y_{k+1} > y_k$. In the second case the inequalities are reversed and $x_k < x_{k+1}$ and $y_{k+1} < y_k$. In both cases the inequality

$$(y_{k+1} - y_k)(x_{k+1} - x_k) < 0 \quad (120)$$

is fulfilled.

Let us consider consecutive sectors of angle τ as in Fig. 5. Denote the number of eigenvalues in each sector by n_k . After a shift by τ , eigenphases from the k th sector will move into the $(k+1)$ th sector. The n_k shifted points divide this sector into $n_k + 1$ intervals. As was proved above, eigenphases in the $(k+1)$ th sector have to intertwine with these shifted eigenphases. Therefore all $n_k + 1$ intervals except the first and the last will be occupied. The first will be occupied if $x_{k+1} < x_k$ and the last interval will be occupied provided $y_{k+1} > y_{k+2}$. These statements can be rewritten in the form of the recurrence relation

$$n_{k+1} = n_k - 1 + \Theta(x_k - x_{k+1}) + \Theta(y_{k+1} - y_{k+2}), \quad (121)$$

where $\Theta(t)$ is the Heaviside step function, $\Theta(t) = 1$ when $t > 0$ and $\Theta(t) = 0$ for $t < 0$. From (120) it follows that this relation can be rewritten in the form

$$n_{k+1} = n_k - 1 + \Theta(x_k - x_{k+1}) + \Theta(x_{k+2} - x_{k+1}) . \quad (122)$$

We now specialise to the case where τ depends on the size N of the Lax matrix. We set

$$\tau = \frac{2\pi}{N}a \quad (123)$$

with fixed a . The case of integer a is trivial and we thus assume that a is not an integer. The total number of sectors of angle τ in the unit circle is

$$K = \left[\frac{N}{a} \right] \quad (124)$$

where $[t]$ denotes the integer part of t . Suppose the beginning of the first sector lies at position $\sigma\lambda_1$. We choose to consider that this eigenvalue does not belong to the first sector, i.e. $y_1 = 0$. Applying Eq. (120) for $k = 0$ implies that $x_1 > x_0$. Thus necessarily $n_2 = n_1 + 1$ and for $k \geq 3$ one easily gets from Eq. (122) that

$$n_k = n_1 + \Theta(x_{k+1} - x_k) . \quad (125)$$

The total number of eigenvalues lying into all K sectors obeys the inequalities

$$N - n_1 - 1 \leq \sum_{k=1}^K n_k \leq N - 1 . \quad (126)$$

The right-hand side inequality comes from the fact that when a is not an integer the union of all K intervals does not overlap the whole circle: in particular, it does not contain the first eigenvalue $\sigma\lambda_1$ from which we start our sectors. The left-hand side inequality is a consequence of the fact that if eigenvalues were shifted by $-\tau$, the first sector in the opposite direction would have exactly the same number of eigenvalues as the second sector, i.e. $n_1 + 1$ eigenvalues: as all K sectors does not cover the whole circle and the uncovering region is smaller than the sector of angle τ , it follows that the number of eigenvalues in all K sectors is larger than $N - (n_1 + 1)$.

From (125) one easily obtains a second inequality

$$Kn_1 + 1 \leq \sum_{k=1}^K n_k \leq K(n_1 + 1) . \quad (127)$$

From these two inequalities it follows that

$$\frac{N}{K+1} - 1 \leq n_1 \leq \frac{N-2}{K} . \quad (128)$$

By definition of K we have

$$\frac{N}{a} - 1 < K < \frac{N}{a} . \quad (129)$$

Substituting in the right-hand side of (128) the minimum of K and in the left-hand side the maximum value of K one gets

$$a - 1 - \frac{a^2}{N+a} < n_1 < a + \frac{a(a-2)}{N-a} , \quad (130)$$

which entails that for N large enough $a - 1 < n_1 < a$. Since n_1 has to be an integer, $n_1 = [a]$. This means that for sufficiently large N , within an interval of length τ from any eigenvalue there are always exactly $[a]$ eigenvalues.

The inverse statement is also true. If within an interval of length τ from any eigenvalue there exist exactly $[a]$ other eigenvalues then all V_α and W_α have the sign of $\sin(\tau/2)/\sin(N\tau/2)$. To see this let us consider e.g. V_α given by Eq. (110). It is a product of terms $\sin x/\sin(x-\tau)$ where $x = \sigma(\lambda_\alpha - \lambda_\beta)/2$. As the distance between two eigenvalues may be restricted, $0 < \sigma\lambda_\alpha - \sigma\lambda_\beta < 2\pi$, x obeys inequality $0 < x < \pi$. Therefore $\sin(x) > 0$, and $\sin(x-\tau)$ is negative when $0 < x < \tau$ and positive when $\tau < x < \pi$. If within an interval of length τ from $\sigma\lambda_\alpha$ there are exactly $[a]$

eigenvalues, then in the product formula for V_α there are exactly $[a]$ negative terms, so that its total sign is $(-1)^{[a]}$. For N large enough the sign of $\sin(\tau/2)/\sin(N\tau/2)$ with $\tau = 2\pi a/N$ is the sign of $\sin \pi a$, which is precisely $(-1)^{[a]}$.

The above arguments prove that for sufficiently large N (whose value depends only on a) the necessary and sufficient condition for the unitarity of matrix Q is that at distance $|\tau|$ from any eigenvalue there exist $[a]$ other eigenvalues.

As matrices L and Q are dual, the unitarity condition for L can be readily deduced: it is that at distance $|\tau|$ from one coordinate μq_k there are exactly $[a]$ other coordinates. Note that in [20] only the case $0 < a < 1$ had been considered.

These restrictions determine the allowed region in coordinate space. We choose as the 'natural' measure of momenta and coordinates the uniform distribution for momenta (between 0 and $2\pi/\sigma$) and coordinates uniformly distributed in the allowed region as explained above. After the change of variables from coordinates and momentum to action-angle variables it follows that the resulting distribution of eigenvalues will be also uniform but in the allowed region of eigenvalues with the only restriction that any interval $]\sigma\lambda_\alpha, \sigma\lambda_\alpha + 2\pi a/N[$ contains exactly $[a]$ eigenvalues. In next section we will show that it is possible to calculate asymptotic expressions for the joint distribution of eigenvalues by using a transfer operator technique.

For numerical investigations, we consider an ensemble of unitary matrices of the form (97), with p_k chosen as independent random variables uniformly distributed between 0 and 2π and the picket fence distribution of coordinates $q_j = j$, $1 \leq j \leq N$ (see section I). Choosing constants such that $\mu = 2\pi/N$, $\sigma = 1$, and $g = a$, so that $\tau = 2\pi a/N$, direct calculations yield

$$\tilde{V}_k = \tilde{W}_k = \left| \frac{\sin N\tau/2}{N \sin \tau/2} \right|. \quad (131)$$

With a slightly different choice of phases in (93), matrix L simplifies to

$$L_{kr} = \frac{e^{i\Phi_k}}{N} \frac{1 - e^{2i\pi a}}{1 - e^{2i\pi((k-r+a)/N)}}, \quad (132)$$

where we denote $p_k = \Phi_k$. This is a particular specialisation of the Lax matrix for the Ruijsenaars-Schneider model. In the form (132) but with $a = bN$ with fixed b it first appeared in [27] as a result of the quantisation of a classical parabolic map on the torus proposed in [10]. When b is a rational number the map considered in [10] corresponds to a pseudo-integrable map of exchange of two intervals. For this particular case where a depends on N , the spectral statistics of the unitary matrix (132) has been obtained analytically in [27] and [28] without knowledge of the relation with the Lax matrix of the Ruijsenaars-Schneider model. In the present case, where a is a fixed parameter independent on N , results are quite different. Analytical calculations of spectral correlation functions for this model are performed in the next sections.

To illustrate the accuracy of the analytical results that we derive in the next sections, we show in Fig. 6 results of numerical computations for matrices of the form (132) for different values of the parameter a . Agreement is remarkable for all parameter values.

VI. JOINT DISTRIBUTION OF EIGENVALUE SPACINGS FOR MODEL RS

We now calculate asymptotic expressions for the joint distribution of eigenvalue spacings for the Lax matrix ensemble corresponding to the RS model. As discussed in the previous section, eigenvalues are such that within an interval of length a from any eigenvalue there are exactly $[a]$ other eigenvalues. We introduce the rescaled nearest-neighbour spacings

$$\xi_k \equiv \frac{N}{2\pi} (\lambda_{\alpha+1} - \lambda_\alpha). \quad (133)$$

The answer strongly depends on the integer part of a . Therefore we consider different cases separately.

A. $0 < a < 1$

The simplest case corresponds to $0 < a < 1$. In this case the only restriction is that the distance between the nearest eigenvalues is larger than a , namely $\xi_k \geq a$. It is convenient to rewrite this restriction as follows. The joint probability density of having $N + 1$ eigenvalues inside an interval of length L is given by

$$p(\xi_1, \xi_2, \dots, \xi_N) = \frac{1}{Z_N(L)} \prod_{j=1}^N g(\xi_j) \delta \left(L - \sum_{k=1}^N \xi_k \right) \quad (134)$$

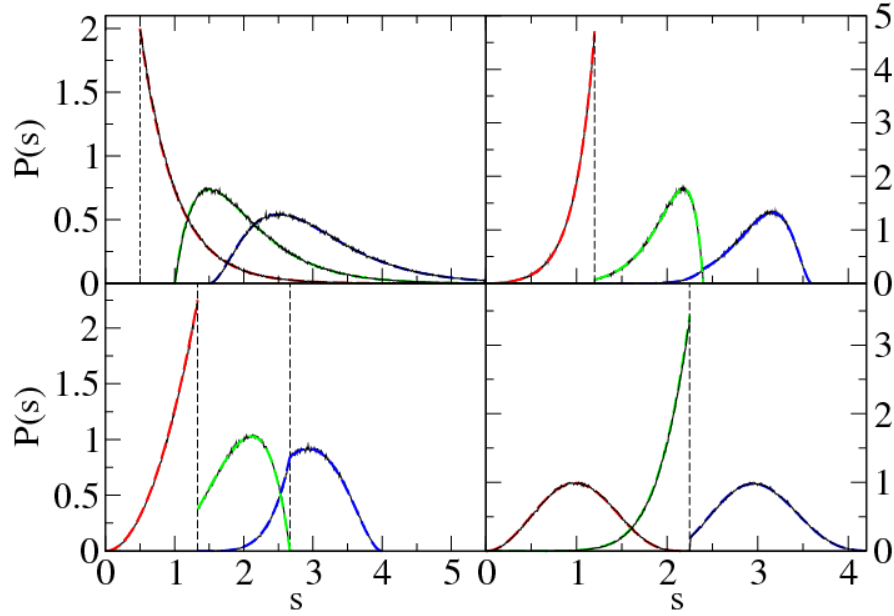


FIG. 6. Nearest-neighbour distributions $P(n, s)$ for the random matrices of model RS with $\mu = 2\pi/N$ and $g = 1/2$ (top left), $6/5$ (top right), $4/3$ (bottom left), and $9/4$ (bottom right), averaged over 1000 realisations of matrices of size $N = 701$. Solid lines are numerical results, dashed lines indicate the analytical expressions (170) for $g = 1/2$, (176)–(180) for $g = 1.2$ and $g = 4/3$, and (192)–(194) for $g = 9/4$. In each panel (from left to right) $P(s) = P(1, s)$ (red), $P(2, s)$ (green) and $P(3, s)$ (blue).

where

$$g(x) = \begin{cases} 0 & \text{when } x < a \\ 1 & \text{otherwise} \end{cases}, \quad (135)$$

and $Z_N(L)$ is the normalisation constant

$$Z_N(L) = \int_0^\infty d\xi_1 \dots \int_0^\infty d\xi_N \prod_{j=1}^N g(\xi_j) \delta\left(L - \sum_{k=1}^N \xi_k\right). \quad (136)$$

We are interested in the joint probability distribution of n consecutive spacings

$$p(\xi_1, \dots, \xi_n) = \int_0^\infty d\xi_{n+1} \dots \int_0^\infty d\xi_N p(\xi_1, \xi_2, \dots, \xi_N) \quad (137)$$

when $L, N \rightarrow \infty$ with mean level spacing $\Delta = L/N$ remaining constant. In the following we set $\Delta = 1$. We shall proceed as it was done in [29]. The multiple integrals in (137) are easily calculated by introducing the function

$$h_{n,N}(L) = \int_0^\infty d\xi_{n+1} \dots \int_0^\infty d\xi_N \prod_{j=1}^N g(\xi_j) \delta\left(L - \sum_{k=1}^N \xi_k\right), \quad (138)$$

whose Laplace transform reads

$$g_{n,N}(t) = \lambda(t)^{N-n} \prod_{k=1}^n g(\xi_k) e^{-t\xi_k}. \quad (139)$$

Here

$$\lambda(t) = \int_0^\infty g(x) e^{-tx} dx = \frac{e^{-ta}}{t} \quad (140)$$

is the Laplace transform of $g(x)$. The inverse Laplace transform of $\lambda(t)^{N-n}$ is then

$$\frac{1}{2i\pi} \int_{c-i\infty}^{c+i\infty} \lambda(t)^{N-n} \exp(Lt) dt = \frac{1}{2i\pi} \int_{c-i\infty}^{c+i\infty} \frac{1}{\lambda(t)^n} \exp(N(\ln \lambda(t) + \Delta t)) dt. \quad (141)$$

The large- N behaviour of (141) is obtained by saddle-point approximation. The value c corresponds to the solution of the saddle-point equation

$$1 + \frac{\lambda'(c)}{\lambda(c)} = 0 \quad (142)$$

(recall that $\Delta = 1$). The solution reads

$$c = \frac{1}{1-a}. \quad (143)$$

Similarly, the inverse Laplace transform of $g_{n,N}(t)$ can be calculated by saddle-point approximation. Rather than calculating explicitly all prefactors coming from the integration, it is easier to observe that the large- N behaviour of the normalisation factor $Z_N = h_{0,N}(L)$ is obtained similarly. One finally gets

$$p(\xi_1, \dots, \xi_n) = \frac{e^{na/(1-a)}}{(1-a)^n} \prod_{j=1}^n g(\xi_j) e^{-\xi_j/(1-a)}. \quad (144)$$

B. $1 < a < 2$

We now consider the case $1 < a < 2$. According to previous section for sufficiently large matrix size, any eigenvalue x_k is such that there exists exactly 1 eigenvalue in the interval $]x_k, x_k + 2\pi a/N[$. In other words, the constraints on x_k are that $x_{k+1} \in]x_k, x_k + 2\pi a/N[$ and $x_{k+2} > x_k + 2\pi a/N$. In terms of the differences (133) between consecutive eigenvalues, the above restriction are equivalent to

$$0 < \xi_{k+1} < a, \quad a < \xi_{k+1} + \xi_k. \quad (145)$$

Introduce the function $f(x)$ by

$$f(x) = \begin{cases} 1 & \text{when } 0 < x < a \\ 0 & \text{otherwise} \end{cases} \quad (146)$$

and $g(x)$ as in (135). Then the joint probability density of $N + 1$ eigenvalues inside an interval of length L is given by the following expression

$$p(\xi_1, \xi_2, \dots, \xi_N) = \frac{1}{Z_N(L)} \prod_{j=1}^N f(\xi_j) g(\xi_j + \xi_{j+1}) \delta \left(L - \sum_{k=1}^N \xi_k \right) \quad (147)$$

where $Z_N(L)$ is the normalisation constant

$$Z_N(L) = \int_0^\infty d\xi_1 \dots \int_0^\infty d\xi_N \prod_{j=1}^N f(\xi_j) g(\xi_j + \xi_{j+1}) \delta \left(L - \sum_{k=1}^N \xi_k \right). \quad (148)$$

The large- N behaviour of the joint probability distribution of n consecutive spacings (137) can be then obtained as above, following [29]. We review the main steps of the procedure. Introducing the function

$$h_{n,N}(L) = \int_0^\infty d\xi_{n+1} \dots \int_0^\infty d\xi_N \prod_{j=1}^N f(\xi_j) g(\xi_j + \xi_{j+1}) \delta \left(L - \sum_{k=1}^N \xi_k \right), \quad (149)$$

its Laplace transform reads

$$g_{n,N}(t) = \int_0^\infty d\xi_{n+1} \dots \int_0^\infty d\xi_N \prod_{j=1}^N e^{-t\xi_j} f(\xi_j) g(\xi_j + \xi_{j+1}). \quad (150)$$

This quantity can be seen as a product of transfer operators

$$g_{n,N}(t) = K_t(\xi_1, \xi_2)K_t(\xi_2, \xi_3) \dots K_t(\xi_{n-1}, \xi_n) \quad (151)$$

$$\times \int_0^\infty d\xi_{n+1} \dots \int_0^\infty d\xi_N K_t(\xi_n, \xi_{n+1}) \dots K_t(\xi_{N-1}, \xi_N)K_t(\xi_N, \xi_1), \quad (152)$$

where the transfer operator, $K_t(\xi, \xi')$ is defined as

$$K_t(\xi, \xi') = f(\xi)g(\xi + \xi')f(\xi') e^{-t(\xi+\xi')/2}. \quad (153)$$

This is a real symmetric operator. Its real eigenvalues, $\lambda_j(t)$, and eigenfunctions, $\phi_j(t; \xi)$, verify

$$\int_0^\infty K_t(\xi, \xi')\phi_j(t; \xi')d\xi' = \lambda_j(t)\phi_j(t; \xi). \quad (154)$$

As this operator is real symmetric its eigenfunctions can be chosen to be orthonormal

$$\int_0^\infty \phi_j(t; \xi)\phi_k(t; \xi)d\xi = \delta_{jk}. \quad (155)$$

The transfer operator can be expanded over the basis of eigenfunctions as

$$K_t(\xi, \xi') = \sum_j \lambda_j(t)\phi_j(t; \xi)\phi_j(t; \xi'). \quad (156)$$

In the large- N limit the dominant contribution to (151) is given by the largest eigenvalue $\lambda_0(t)$. Then using the orthogonality property (155) of the ϕ_j we get

$$g_{n,N}(t) \underset{N \rightarrow \infty}{\sim} \prod_{k=1}^{n-1} K_t(\xi_k, \xi_{k+1})\lambda_0(t)^{N-n+1}\phi_0(t; \xi_n)\phi_0(t; \xi_1). \quad (157)$$

The large- N behaviour of $h_{n,N}(L)$ is obtained by performing the inverse Laplace transform of $\lambda_0(t)^{N-n+1}$. As in (141) the leading term is obtained by saddle-point approximation; here the saddle-point is such that

$$1 + \frac{\lambda_0'(c)}{\lambda_0(c)} = 0 \quad (158)$$

(again we take $\Delta = 1$). Calculating in a similar way the large- N behaviour of the normalisation factor $Z_N = h_{0,N}(L)$, one finally gets

$$p(\xi_1, \xi_2, \dots, \xi_n) = \frac{1}{\lambda_0^{n-1}(c)}\phi_0(c; \xi_1)K_c(\xi_1, \xi_2)K_c(\xi_2, \xi_3) \dots K_c(\xi_{n-1}, \xi_n)\phi_0(c; \xi_n). \quad (159)$$

C. $m < a < m + 1$

The general case $m < a < m + 1$ can be treated in a similar way. In this case each interval $]x_k, x_k + 2\pi a/N[$ contains exactly m eigenvalues x_{k+1}, \dots, x_{k+m} . In terms of the rescaled differences (133) between consecutive eigenvalues this condition is equivalent to the following two inequalities

$$0 < \xi_k + \xi_{k+1} + \dots + \xi_{k+m-1} < a \quad (160)$$

$$a < \xi_k + \xi_{k+1} + \dots + \xi_{k+m}. \quad (161)$$

In analogy with Eq. (147), the joint probability of eigenvalues spacings thus reads

$$p(\xi_1, \xi_2, \dots, \xi_N) = \frac{1}{Z_N(L)} \prod_{j=1}^N f(\xi_j + \dots + \xi_{j+m-1})g(\xi_j + \dots + \xi_{j+m})\delta\left(L - \sum_{k=1}^N \xi_k\right) \quad (162)$$

with f and g defined by (146) and (135). The large- N behaviour is calculated as above by introducing a transfer operator, which in this case depends on two sets of variables $\boldsymbol{\xi} = (\xi_1, \dots, \xi_m)$ and $\boldsymbol{\xi}' = (\xi'_1, \dots, \xi'_m)$ shifted by one unit i.e. $\xi_2 = \xi'_1$, $\xi_3 = \xi'_2$, \dots , $\xi_m = \xi'_{m-1}$ (see e.g. [29]). The explicit form of the transfer operator is the following

$$K(\boldsymbol{\xi}, \boldsymbol{\xi}') = \delta(\xi_2 - \xi'_1) \dots \delta(\xi_m - \xi'_{m-1}) \\ \times e^{-t\xi_1/2} f(\xi_1 + \dots + \xi_m) g(\xi_1 + \dots + \xi_m + \xi'_m) f(\xi'_1 + \dots + \xi'_m) e^{-t\xi'_m/2} . \quad (163)$$

The eigenvalue equation

$$\int K(\boldsymbol{\xi}, \boldsymbol{\xi}') \phi(\boldsymbol{\xi}') d\boldsymbol{\xi}' = \lambda \phi(\boldsymbol{\xi}) \quad (164)$$

reduces to a one-dimensional equation because of the δ -functions appearing in the definition of the transfer operator. This equation can be written in the form

$$e^{-t\xi_1/2} \int_0^\infty e^{-tz/2} g(\xi_1 + \xi_2 + \dots + \xi_m + z) \phi(t; \xi_2, \dots, \xi_m, z) dz = \lambda(t) \phi(t; \xi_1, \dots, \xi_m) . \quad (165)$$

Here it is implicitly assumed that all variables $\xi_j > 0$ and

$$\phi(t; \xi_1, \dots, \xi_m) = 0 \text{ when } \xi_1 + \dots + \xi_m > a . \quad (166)$$

As above the largest eigenvalue, $\lambda_0(t)$ as well as the corresponding eigenfunction $\phi_0(t; \boldsymbol{\xi})$, calculated at point $t = c$ obeying the same saddle-point condition Eq. (158), determine all correlation functions in the limit of large N . The joint probability of n consecutive spacings takes a different form for $n \leq m$ and $n > m$. For $n \leq m$

$$p(\xi_1, \dots, \xi_n) = \int_0^a d\xi_{n+1} \dots \int_0^a d\xi_m \phi_0(c; \xi_1, \dots, \xi_m) \phi_0(c; \xi_m, \dots, \xi_1), \quad (167)$$

while for $n > m$

$$p(\xi_1, \dots, \xi_n) = \lambda_0(c)^{-n+m} \phi_0(c; \xi_n, \dots, \xi_{n-m}) \phi_0(c; \xi_1, \dots, \xi_m) e^{-c \sum_{s=1}^n \xi_s} \\ \times \prod_{j=1}^{n-m+1} f(\xi_j + \dots + \xi_{j+m-1}) \prod_{j=1}^{n-m} g(\xi_j + \dots + \xi_{j+m}). \quad (168)$$

VII. NEAREST-NEIGHBOUR SPACING DISTRIBUTIONS FOR MODEL RS

In the previous section we have derived expressions for the joint distribution of eigenvalue spacings $p(\xi_1, \dots, \xi_n)$. From these expressions the n th nearest-neighbour spacing distribution can be calculated as

$$P(n, s) = \int_0^\infty d\xi_1 \dots \int_0^\infty d\xi_n p(\xi_1, \dots, \xi_n) \delta\left(s - \sum_{i=1}^n \xi_i\right) . \quad (169)$$

A. $0 < a < 1$

For $0 < a < 1$ these integrals are easily calculable (e.g. by Laplace transform) and from the joint distribution Eq. (144) we obtain

$$P(n, s) = \begin{cases} \frac{e^{(na-s)/(1-s)} (s-na)^{n-1}}{(1-a)^n (n-1)!}, & s \geq na \\ 0, & 0 < s < na \end{cases} . \quad (170)$$

Comparison with numerical simulations is displayed at Fig. 6.

B. $1 < a < 2$

When $1 < a < 2$ the joint distribution is given by Eq. (159). What remains is to calculate the largest eigenvalue of the transfer operator K_t , as well as its associated eigenfunction. As K_t is a positive operator, the analog of the Perron-Frobenius theorem states that the eigenvector corresponding to the largest eigenvalue is positive. Orthogonality of the eigenfunctions implies that the converse is also true. Thus if one finds a positive eigenfunction then the corresponding eigenvalue is the largest one. The eigenvalue equation (154) is equivalent to

$$e^{-t\xi/2} \int_{a-\xi}^a e^{-t\xi'/2} \phi(\xi') d\xi' = \lambda \phi(\xi). \quad (171)$$

Let us look for solutions of Eq. (171) positive on $[0, a]$ under the form $\phi(\xi) = \sinh \rho \xi$, with ρ some unknown complex parameter. Since Eq. (171) should hold for all $\xi \in [0, a]$ we get the necessary condition

$$t = -2\rho \coth(\rho a). \quad (172)$$

When $t < -2/a$, Eq. (172) admits two real solutions $\rho = \pm \rho_0$. Thus $\phi(\xi) = \sinh \rho_0 \xi$ with $\rho_0 > 0$ solution of Eq. (172) is a positive solution of Eq. (171). If $t > -2/a$, Eq. (172) admits two pure imaginary solutions $\rho = \pm i\rho_0$, thus $\phi(\xi) = \sin \rho_0 \xi$ with $\rho_0 > 0$ and $i\rho_0$ solution of Eq. (172) is a positive solution of Eq. (171). Finally if $t = -2/a$, $\rho = 0$ is the unique solution to Eq. (172). In that case $\phi(\xi) = \xi$ is a solution of Eq. (171) which is positive on $[0, a]$. Thus for all t we have a positive solution to Eq. (171). Properly normalised, this solution gives the eigenvector $\phi_0(t; \xi)$. The corresponding eigenvalue is given by

$$\lambda_0(t) = \frac{e^{(\rho-t/2)a}}{\rho - t/2}, \quad (173)$$

with ρ an implicit function of t .

The saddle-point c is a solution of Eq. (158). For λ_0 given by Eq. (173) the condition becomes

$$1 + 2a(2-a)\rho^2 - \cosh 2\rho a + 2\rho(a-1) \sinh 2\rho a = 0 \quad (174)$$

and the saddle-point c is obtained from ρ through Eq. (172). Equivalently, this condition can be expressed as

$$a = \frac{2z^2 - z \sinh 2z}{z^2 + \sinh^2 z - z \sinh 2z}, \quad z = \rho a. \quad (175)$$

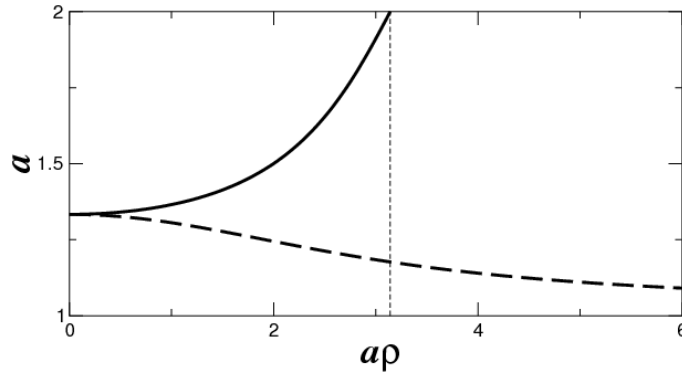


FIG. 7. (Color online). Graph of Eq. (175) for real (dashed) or pure imaginary (solid) $z = \rho a$. Dashed vertical line indicates the abscissa equal π .

In Fig. 7 we plot a as a function of $z = \rho a$. For $1 < a < 4/3$ Eq. (175) has a unique real solution $\rho_0 > 0$, and $\phi_0(\xi) = \sinh \rho_0 \xi$ is a positive eigenfunction of the transfer operator. For $4/3 < a < 2$ Eq. (175) has a unique pure imaginary solution $i\rho_0$ with $\rho_0 > 0$. Furthermore in that latter case $\rho_0 a \in]0, \pi[$, so that $\phi_0(\xi) = \sin \rho_0 \xi$ is an eigenfunction of the transfer operator which is positive on $[0, a]$. At $a = 4/3$ the unique solution is $\rho = 0$ and $\phi_0(\xi) = \xi$ is a positive eigenfunction of the transfer operator. The n th nearest-neighbour spacing distribution can now be calculated from Eq. (159).

In the case $n = 1$ it directly gives us the nearest-neighbour spacing distribution $P(s) = A^2 \phi_0(s)^2$, where A is the normalisation constant. It is nonzero only for $s \in [0, a]$, where it takes the following form

$$P(s) = \begin{cases} A^2 \sinh^2(\rho s) & \text{when } 1 < a < 4/3 \\ \frac{81}{64} s^2 & \text{when } a = 4/3 \\ A^2 \sin^2(\rho s) & \text{when } 4/3 < a < 2 \end{cases}. \quad (176)$$

Constants A and ρ can be determined either by solving Eq. (175) and normalizing the eigenfunction $\phi_0(\xi)$, or equivalently by imposing the normalisation conditions (4). The next-to-nearest distribution, $P(2, s)$ is non-zero only when $a < s < 2a$ and within this interval it is given by

$$P(2, s) = \frac{A^2}{\lambda_0(c)} e^{-cs/2} \int_{s-a}^a \phi_0(\xi) \phi_0(s - \xi) d\xi. \quad (177)$$

In particular for $a = 4/3$ all integrals can be calculated analytically and $P(2, s)$ has the form

$$P(2, s) = \left(-\frac{3}{2} + \frac{27}{16}s - \frac{81}{512}s^3\right) e^{3s/4-1}. \quad (178)$$

In a similar manner one can obtain the higher nearest-neighbour functions. For example, $P(3, s)$ is given by the formula

$$P(3, s) = \frac{A^2 e^{-cs/2}}{\lambda_0(c)^2} \int_0^a d\xi_1 \phi_0(\xi_1) \int_{a-\xi_1}^a d\xi_2 e^{-c\xi_2/2} \int_{a-\xi_2}^a d\xi_3 \phi_0(\xi_3) \delta\left(s - \sum_{i=1}^n \xi_i\right). \quad (179)$$

It is non-zero only when $a < s < 3a$. In particular for $a = 4/3$ we obtain

$$p(3, s) = \begin{cases} \left(\frac{3}{4} - \frac{81}{32}s + \frac{81}{512}s^3\right) e^{3s/4-1} + \frac{81}{64}s^2 & \text{when } 4/3 < s < 8/3 \\ \left(-\frac{9}{4} + \frac{27}{32}s - \frac{81}{512}s^3\right) e^{3s/4-1} + 9e^{3s/2-4} & \text{when } 8/3 < s < 4 \end{cases}. \quad (180)$$

In Fig. 6 these formulas are compared with numerical simulations and show a remarkable agreement.

C. $2 < a < 3$

For $2 < a < 3$ the joint distribution is given by (168). The largest eigenvalue and corresponding eigenfunction of the transfer operator (163) are solution of the eigenvalue equation (165). For $m = 2$ it takes the form

$$e^{-t\xi_1/2} \int_{a-\xi_1-\xi_2}^{a-\xi_2} e^{-t\xi_3/2} \phi(\xi_2, \xi_3) d\xi_3 = \lambda \phi(\xi_1, \xi_2). \quad (181)$$

Let us look for solutions of the form similar to Bethe Ansatz

$$\begin{aligned} \phi(\xi_1, \xi_2) = & e^{\alpha\xi_1 + \beta\xi_2} + e^{-\beta\xi_1 + (\alpha - \beta - \mu)\xi_2} + e^{(-\alpha + \beta + \mu)\xi_1 - \alpha\xi_2} \\ & - e^{-\beta\xi_1 - \alpha\xi_2} - e^{\alpha\xi_1 + (\alpha - \beta - \mu)\xi_2} - e^{(-\alpha + \beta + \mu)\xi_1 + \beta\xi_2}, \end{aligned} \quad (182)$$

where we have set $\mu = -t/2$. As Eq. (181) has to be fulfilled for all ξ_1, ξ_2 , this function is a solution of (181) if and only if the following conditions are valid

$$\frac{e^{a(\mu + \beta)}}{\mu + \beta} = \frac{e^{a(\mu - \alpha)}}{\mu - \alpha} = \frac{e^{a(\alpha - \beta)}}{\alpha - \beta} = -\frac{\lambda}{a}. \quad (183)$$

From the first equality in Eq. (183) one can express μ as a function of α and β . After inspection we found that the solutions of the above equations have the following form

$$\alpha a = \frac{1}{2}x_1 + ix_2, \quad \beta a = -\frac{1}{2}x_1 + ix_2, \quad \mu a = \frac{1}{2}x_1 + x_3 \quad (184)$$

with real parameters x_1, x_2 , and x_3 . Under this substitution the eigenfunction (182) is transformed to

$$\phi(\xi_1, \xi_2) = e^{x_1(\xi_1 - \xi_2)/2a} \left(\sin \frac{x_2(\xi_1 + \xi_2)}{a} - e^{(x_1 - x_3)\xi_2/a} \sin \frac{x_2\xi_1}{a} - e^{-(x_1 - x_3)\xi_1/a} \sin \frac{x_2\xi_2}{a} \right) \quad (185)$$

where from (183) x_1 , x_2 , and x_3 must fulfill the following equalities

$$\frac{e^{x_1}}{x_1} = \frac{e^{x_3+ix_2}}{x_3+ix_2} = \frac{e^{x_3-ix_2}}{x_3-ix_2} = -\frac{\lambda}{a} \quad (186)$$

and depend on time t through the relation

$$-\frac{ta}{2} = \frac{x_1}{2} + x_3. \quad (187)$$

This implies that

$$x_3 = \frac{x_2}{\tan x_2} \quad (188)$$

and, consequently, x_1 is related with x_2 as follows

$$\frac{e^{x_1}}{x_1} = \frac{\sin x_2}{x_2} e^{x_2/\tan x_2}. \quad (189)$$

The eigenfunction corresponding to the largest eigenvalue of the transfer operator is thus given by (185) with x_1 , x_2 , and x_3 real parameters depending on t , which must verify (188) and (189) and be such that $\phi(\xi_1, \xi_2)$ is a positive function over $[0, a]^2$.

The saddle-point condition is again given by Eq. (158). Using (186)–(189) we get a second relation between x_1 and x_2 , namely

$$a = \frac{1}{1-1/x_1} + \frac{2 - \sin(2x_2)/x_2}{1 + \sin^2(x_2)/x_2^2 - \sin(2x_2)/x_2}. \quad (190)$$

Equations (189) and (190) determine parameters x_1 and x_2 at a given a . To get a positive eigenfunction ϕ it is necessary to get the solutions in the intervals

$$x_1 < 0, \quad \pi < x_2 < 2\pi. \quad (191)$$

The knowledge of these parameters allows us to calculate the eigenfunction (185), from which the nearest-neighbour distributions can be deduced through Eq. (168). The first distributions read

$$P(s) = A \int_0^{a-s} \phi(s, y) \phi(y, s) dy, \quad (192)$$

$$P(2, s) = A \int_0^s \phi(s-y, y) \phi(y, s-y) dy, \quad (193)$$

and

$$P(3, s) = \frac{A}{\lambda} \int_{s-a}^a dx e^{\mu x} \int_{s-a}^{s-x} dy e^{\mu y} \phi(s-x-y, x) \phi(s-x-y, y) dy, \quad (194)$$

with A the normalisation constant

$$A = \left(\int_0^a dx \left(\int_0^{a-x} dy \phi(x, y) \phi(y, x) \right) \right)^{-1}. \quad (195)$$

These analytical expressions perfectly agree with numerical simulations, as shown in Fig. 6.

VIII. LEVEL COMPRESSIBILITY FOR MODEL RS

The expressions for the joint distribution of eigenvalue spacings $p(\xi_1, \dots, \xi_n)$ obtained in section VI allow to derive formulas for the level compressibility χ , which characterises the asymptotic behaviour of the number variance.

The number variance $\Sigma^2(L)$ is the average variance of the number of energy levels in an interval of length L . It is defined from the two-point correlation function $R_2(s) = \sum_{n=1}^{\infty} P(n, s)$ as

$$\Sigma^2(L) = L - 2 \int_0^L ds (L-s)(1-R_2(s)). \quad (196)$$

For systems with intermediate spectral statistics, $\Sigma^2(L) \sim \chi L$ for large L . In order to obtain the large- N behaviour of the level compressibility we calculate the Laplace transform of the two-point correlation function. It has a series expansion of the form

$$g_2(t) = \int_0^L ds R_2(s) e^{-ts} = \frac{1}{t} + \frac{\chi - 1}{2} + O(t) \quad (197)$$

which allows us to obtain χ (see [29] for more detail).

A. Case $0 < a < 1$

The n th nearest-neighbour spacing distributions are given by Eq. (170). Summation over n gives the two-point correlation function, and its Laplace transform is readily obtained, yielding

$$g_2(t) = \frac{1}{e^{at}(1+t-at) - 1}. \quad (198)$$

Small- t expansion of $g_2(t)$ gives

$$\chi = (1-a)^2. \quad (199)$$

B. Case $1 < a < 2$

The functions $P(n, s)$ are given by Eqs. (159) and (169). Their Laplace transform reads

$$g(n, t) = \frac{1}{\lambda_0^{n-1}(c)} \int_0^\infty d\xi_1 \dots \int_0^\infty d\xi_n \times \phi_0(c; \xi_1) K_c(\xi_1, \xi_2) K_c(\xi_2, \xi_3) \dots K_c(\xi_{n-1}, \xi_n) \phi_0(c; \xi_n) e^{-t(\xi_1 + \dots + \xi_n)}, \quad (200)$$

where as in the previous sections $\lambda_0(c)$ is the largest eigenvalue of the transfer operator (153) and $\phi_0(c; \xi)$ its associated eigenfunction, both taken at the saddle-point c . Using the definition (153) of the transfer operator, we see that $g(n, t)$ can be rewritten

$$g(n, t) = \frac{1}{\lambda_0^{n-1}(c)} \int_0^\infty d\xi \int_0^\infty d\xi' \phi_0(c; \xi) e^{-t\xi/2} K_{c+t}(\xi, \xi')^{n-1} \phi_0(c; \xi') e^{-t\xi'/2}. \quad (201)$$

Replacing the transfer operator by its expansion (156) and summing over n we get

$$g_2(t) = \sum_j \frac{\lambda_0(c)}{\lambda_0(c) - \lambda_j(c+t)} \left(\int_0^\infty d\xi \phi_0(c; \xi) \phi_j(c+t; \xi) e^{-t\xi/2} \right)^2. \quad (202)$$

One can check, using normalisation (155) of the eigenfunctions and the saddle-point condition (158), that the leading-order term is given by $g_2(t) \sim 1/t$. The next-order term can be simplified using the normalisation of ϕ_0 . It yields

$$g_2(t) = \frac{1}{t} - \frac{\lambda_0''(c)}{2\lambda_0'(c)} - 1 + o(t^2), \quad (203)$$

from which one gets

$$\chi = -1 - \frac{\lambda_0''(c)}{\lambda_0'(c)}. \quad (204)$$

Here $\lambda_0(t)$ is given by (173) (with ρ depending on t through (172)), and c is given by condition (158). After calculation, χ can be expressed as a function of ρ at the saddle-point. We get

$$\chi = \left(\frac{a^2}{4} - \frac{4a(1-a)z^2 + a^2 \sinh^2 z}{(2z - \sinh 2z)^2} \sinh^2 z \right) \frac{\sinh^2 z}{z^2}, \quad z = \rho a, \quad (205)$$

with ρ the real positive solution ρ_0 of (175) for $1 < a < 4/3$ or the pure imaginary solution $i\rho_0$ of (175) for $4/3 < a < 2$. For $a = 4/3$, the limit $\rho \rightarrow 0$ in (205) gives $\chi = 4/9$.

C. Case $2 < a < 3$

As in the previous case χ is given by (204) with $\lambda_0(t)$ given by (186), with x_1, x_2, x_3 and t related through (187)–(189). From (187)–(188), parameter x_1 can be expressed as

$$x_1 = -2\frac{x_2}{\tan x_2} - at. \quad (206)$$

Differentiating both (189) and (206) with respect to time we obtain dx_1/dt and dx_2/dt as a function of x_1 and x_2 , and then similarly d^2x_1/dt^2 and d^2x_2/dt^2 . Using (186), the saddle-point condition (158) can be rewritten

$$1 + \frac{dx_1}{dt} \left(1 - \frac{1}{x_1}\right) = 0, \quad (207)$$

and from (204) χ can then be expressed as

$$\chi = \frac{1}{(1-x_1)^2} + \frac{d^2x_1}{dt^2} \left(1 - \frac{1}{x_1}\right). \quad (208)$$

Using the expression obtained d^2x_1/dt^2 we finally obtain χ as a function of x_1 and x_2 , with x_1, x_2 obtained as solution of (189)–(190). Inverting (190) we get

$$x_1 = \frac{a \sin^2 x_2 + (a-2)x_2^2 + (1-a)x_2 \sin 2x_2}{(a-1) \sin^2 x_2 + (a-3)x_2^2 + (2-a)x_2 \sin 2x_2}. \quad (209)$$

After some manipulation χ simplifies to

$$\begin{aligned} \chi = & \frac{1}{a(\sin^2 x_2 + x_2^2 - x_2 \sin 2x_2)^2} \left[(a-3)^2(a-2)x_2^4 \right. \\ & - (a-3)(a-1)(2a-5)x_2^3 \sin 2x_2 + 2(a-2)((\cos 2x_2 + 2)(a-1)(a-2) - 3)x_2^2 \sin^2 x_2 \\ & \left. - 2a(a-2)(2a-3)x_2 \cos x_2 \sin^3 x_2 + a(a-1)^2 \sin^4 x_2 \right]. \end{aligned} \quad (210)$$

Figure 8 is a plot of the level compressibility χ . The theoretical prediction obtained from (199), (205) and (210) agrees with numerical data.

D. Asymptotics in the vicinity of integer a

For integer a the spectrum is rigid and thus the level compressibility is expected to take the value 0. Here we consider the first-order expansion of χ in the vicinity of integer a . We will show that at lowest order the expansion of χ around $a = n$ is given by $\chi \simeq (1-a)^2/n^2$.

We first consider the expansion around $a = 1$. Let $a = 1 + \epsilon$. For $a < 1$ we have $\chi = (1-a)^2 = \epsilon^2$, thus expansion is trivial. For $a > 1$ χ is given by (205) with a and z related by (175). At $a = 1$ the solution of (175) is $z = \infty$. An asymptotic expansion of (205) and (175) yields

$$a = \frac{2z}{2z-1} + F(z)e^{-2z} + o(e^{-2z}) \quad (211)$$

where $F(z)$ is some rational fraction in z . Thus at first order

$$z = \frac{1}{2} \left(1 + \frac{1}{\epsilon}\right). \quad (212)$$

Expanding χ for large z to lowest order gives

$$\chi = a(a-1) + a^2 \left(\frac{1}{4z^2} - \frac{1}{2z}\right) + G(z)e^{-2z} + o(e^{-2z}) \quad (213)$$

with $G(z)$ is some rational fraction in z . Using (212) one gets $\chi \simeq \epsilon^2 = (1-a)^2$.

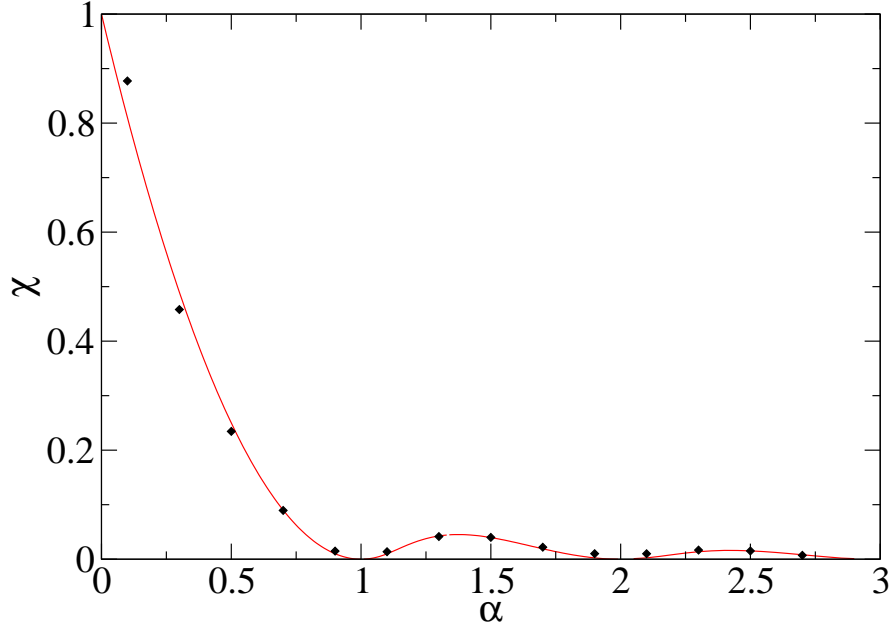


FIG. 8. Level compressibility χ . Black diamonds are the numerical values extracted from a cubic fit of $\Sigma^2(L)$ over the range $L \in [0, 80]$, with $\Sigma^2(L)$ the variance of the number of levels in an interval of length L averaged over 50 windows of length L , calculated from the unfolded spectrum with mean level spacing $\Delta = 1$, for 10000 realisations of the random matrix with matrix size $N = 256$. Red curve is the theoretical prediction (199), (205) and (210).

Suppose now that $a = 2 + \epsilon$. For $a < 2$ χ is given by (205), with $\rho = i\rho_0$ solution of (175). Equivalently, χ is given by

$$\chi = \left(\frac{a^2}{4} + \frac{4a(1-a)z^2 + a^2 \sin^2 z}{(2z - \sin 2z)^2} \sin^2 z \right) \frac{\sin^2 z}{z^2}, \quad (214)$$

with z the real positive solution of

$$a = \frac{2z^2 - z \sin 2z}{z^2 - 2z \sin 2z + \sin^2 z}. \quad (215)$$

At point $a = 2$ the solution is $z = \pi$. Expanding both sides of (215) at lowest order in ϵ with $z = \pi + z_1\epsilon$ we get $z_1 = \pi/2$. Inserting this expansion for z in (214) we obtain $\chi = \epsilon^2/4 + o(\epsilon^3)$.

For $a > 2$ χ is given by (210), with x_1 and x_2 specified by (189)–(190). At $a = 2$, we have

$$\chi = \frac{2 \sin^4 x_2 - x_2^3 \sin 2x_2}{2(x_2^2 + \sin^2 x_2 - x_2 \sin 2x_2)} \quad (216)$$

which vanishes for $x_2 = \pi$. For $x_2 = \pi + t_1\epsilon + t_2\epsilon^2$ we have the expansion

$$\chi = \left(\frac{1}{2} - \frac{t_1}{\pi} \right) \epsilon + \left(-\frac{5}{4} + \frac{9t_1}{2\pi} - \frac{3t_1^2}{\pi^2} - \frac{t_2}{\pi} \right) \epsilon^2 + o(\epsilon^3). \quad (217)$$

Equation (189) is equivalent to

$$\exp \left(x_1 - \frac{x_2}{\tan x_2} \right) = \frac{x_1 \sin x_2}{x_2} \quad (218)$$

and the small- ϵ expansion of both members of this equation reads

$$x_1 - \frac{x_2}{\tan x_2} = -\frac{\pi}{\epsilon t_1} + \frac{\pi t_2}{t_1} - 1 + o(1), \quad (219)$$

$$\frac{x_1 \sin x_2}{x_2} = -\frac{\pi t_1 - 2t_1^2}{\pi^2} \epsilon^2 + \frac{\pi^2 t_1 - 5\pi t_1^2 + 6t_1^3 + \pi^2 t_2 - 4\pi t_1 t_2}{\pi^3} \epsilon^3 + o(\epsilon^4) \quad (220)$$

(we use (209) to obtain the expansion of x_1). This implies that the two first terms in the expansion (220) must vanish, thus $t_1 = \pi/2$ and $t_2 = 0$. Putting these values into (217) gives $\chi = \epsilon^2/4 + o(\epsilon^3)$.

The same result can be obtained from (210) and (189)–(190) for $a < 3$. At $a = 3$ again x_2 takes the value π , and an expansion $x_2 = \pi + t_1\epsilon$ gives $t_1 = \pi/3$, whence $\chi = \epsilon^2/9 + o(\epsilon^3)$.

IX. CONCLUSION

In this paper we construct new random matrix ensembles with unusual properties. Random matrices from these ensembles are Lax matrices of N -body integrable classical systems with a certain measure of momenta and coordinates. Though such matrices are not invariant over rotation of the basis (as usual random matrix ensembles) the joint distribution of their eigenvalues can be calculated analytically. Four different models are considered in detail. Three of them correspond to rational, hyperbolic, and trigonometric Calogero-Moser models. The fourth is related to the trigonometric Ruijsenaars-Schneider model. For the trigonometric Calogero-Moser model and the Ruijsenaars-Schneider model spectral correlation functions are calculated explicitly. For rational and hyperbolic Calogero-Moser models Wigner-type surmises are proposed. Our formulas are in a good agreement with results of direct numerical calculations.

Appendix A: Hamilton-Jacobi equations

In this appendix we check that the action-angle variables λ_α and ϕ_α for model CM_r verify Hamilton-Jacobi equations by calculating their time derivative. We use the fact that for the Lax pair (L, M) the matrix M can be seen as a time derivative operator for the eigenfunctions of the matrix L . Namely, if $(u_k)_{1 \leq k \leq N}$ is a normalised eigenvector of L , then

$$\dot{u}_k = \sum_r M_{kr} u_r \quad \text{and} \quad \dot{u}_k^* = - \sum_r u_r^* M_{rk} \quad (\text{A1})$$

(here $*$ denotes complex conjugation). For model CM_r , the Lax matrix M is given by

$$M_{kr} = -ig \delta_{kr} \sum_{j \neq k} \frac{1}{(q_k - q_j)^2} + ig(1 - \delta_{kr}) \frac{1}{(q_k - q_r)^2}. \quad (\text{A2})$$

One can easily check that for $1 \leq k, r \leq N$ one has

$$p_r \delta_{kr} + M_{kr}(q_k - q_r) = L_{kr}. \quad (\text{A3})$$

Deriving (39) with respect to time, using the definition of Q , yields

$$\dot{\phi}_\alpha = \sum_k [\dot{u}_k^*(\alpha) q_k u_k(\alpha) + u_k^*(\alpha) \dot{q}_k u_k(\alpha) + u_k^*(\alpha) q_k \dot{u}_k(\alpha)]. \quad (\text{A4})$$

From Hamilton-Jacobi equations $\dot{q}_k = p_k$. Using (A1), we get that the time derivative of ϕ_α is given by

$$\dot{\phi}_\alpha = \sum_{k,r} u_k^*(\alpha) [p_k \delta_{kr} + M_{kr}(q_k - q_r)] u_r(\alpha) = \sum_{k,r} u_k^*(\alpha) L_{kr} u_r(\alpha) = \lambda_\alpha. \quad (\text{A5})$$

The time derivative of λ_α is easily obtained from (7), (A1) and (34), yielding $\dot{\lambda}_\alpha = 0$. This shows that the λ_α and ϕ_α verify Hamilton-Jacobi equations.

Appendix B: Identities

The purpose of the Appendix is to give, for completeness, the proofs of certain often used formulas.

Let coefficients b_m obey the following system of linear equations for all $n = 1, \dots, N$ with known x_m and y_m

$$\sum_{m=1}^N \frac{b_m}{x_m - y_n} = 1. \quad (\text{B1})$$

Then b_m for all $m = 1, \dots, N$ are expressed through x_m and y_m by using e.g. the Cauchy determinants

$$b_m = \frac{\prod_n (x_m - y_n)}{\prod_{s \neq m} (x_m - x_s)}. \quad (\text{B2})$$

The following identities are also useful. For all $l = 1, \dots, N$ one has

$$\sum_{m=1}^N \frac{b_m}{(x_m - y_l)^2} = -\frac{\prod_{n \neq l} (y_l - y_n)}{\prod_s (y_l - x_s)}, \quad (\text{B3})$$

$$\sum_{m=1}^N b_m = \sum_{m=1}^N (x_m - y_m), \quad (\text{B4})$$

and

$$\sum_{m=1}^N \frac{b_m}{x_m} = 1 - \prod_n \frac{y_n}{x_n}. \quad (\text{B5})$$

A simple way to check (B2) is to consider the function

$$f_n(x) = \frac{\prod_{r \neq n} (x - y_r)}{\prod_s (x - x_s)} = \frac{\prod_r (x - y_r)}{(x - y_n) \prod_s (x - x_s)}. \quad (\text{B6})$$

Asymptotically this function decreases as $1/x$ when $x \rightarrow \infty$, so that the integral over a large contour encircling all poles equals 1. Rewriting this integral as the sum over all finite poles gives

$$1 = \sum_m \frac{\prod_r (x_m - y_r)}{(x_m - y_n) \prod_{s \neq m} (x_m - x_s)} \quad (\text{B7})$$

which proves (B2).

The equality (B3) can be obtained by the integration of the function

$$\bar{f}_l(x) = \frac{\prod_n (x - y_n)}{(x - y_l)^2 \prod_n (x - x_n)} \quad (\text{B8})$$

over a contour which includes all poles. As this function decreases as $1/x^2$ when $x \rightarrow \infty$ the integral equals zero. Taking the sum over poles at $x = x_n$ with all $n = 1, \dots, N$ and at $x = y_l$ one verifies (B3).

To get (B4) one has to integrate the function

$$\tilde{f}(x) = \prod_{n=1}^N \frac{x - y_n}{x - x_n} \quad (\text{B9})$$

over the large contour and compare the residues at infinity and at $x = x_n$.

Let us now consider the function

$$\hat{f}(x) = \frac{\prod_n (x - y_n)}{x \prod_n (x - x_n)}. \quad (\text{B10})$$

It decreases as $1/x$ when $x \rightarrow \infty$ and has poles at $x = 0$ and $x = x_m$ with $m = 1, \dots, N$. Integrating it over a contour encircling all poles one obtains (B5).

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