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Quantum chaos on discrete graphs

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Abstract

Adapting a method developed for the study of quantum chaos on quantum (metric) graphs (Kottos and Smilansky 1997 Phys. Rev. Lett. 79 4794, Kottos and Smilansky 1999 Ann. Phys., NY 274 76), spectral $\zeta$ functions and trace formulae for discrete Laplacians on graphs are derived. This is achieved by expressing the spectral secular equation in terms of the periodic orbits of the graph and obtaining functions which belong to the class of $\zeta$ functions proposed originally by Ihara (1966 J. Mat. Soc. Japan 18 219) and expanded by subsequent authors (Stark and Terras 1996 Adv. Math. 121 124, Kotani and Sunada 2000 J. Math. Sci. Univ. Tokyo 7 7). Finally, a model of ‘classical dynamics’ on the discrete graph is proposed. It is analogous to the corresponding classical dynamics derived for quantum graphs (Kottos and Smilansky 1997 Phys. Rev. Lett. 79 4794, Kottos and Smilansky 1999 Ann. Phys., NY 274 76).

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1. Introduction and preliminaries

Some ten years ago quantum (metric) graphs were proposed as a convenient paradigm for the study of quantum chaos in compact [1] and scattering [5] systems. The crucial point which highlighted the close similarity between metric graphs—whose only claim to complexity is their topology—and chaotic Hamiltonian flows is the formal similarity between the trace formulae [6, 7] which express the spectral densities as sums over periodic orbits. Requiring additionally that the lengths of the bonds are rationally independent, and that the graph is well connected, render the spectrum of the Schrödinger operator on graphs sufficiently disordered to display spectral statistics which are consistent with the predictions of random matrix theory. Another important contact point was the identification of the classical dynamics on the graph, which is derived from the quantum evolution, and which describes, under certain conditions,
an exponential approach to equilibrium, in analogy with mixing Hamiltonian flows. The \( \zeta \) function for the Peron–Frobenius operator on the graph can be written in terms of the same periodic orbits which are used for the quantum spectral \( \zeta \) function—in close similarity with the expansion of the Ruelle \( \zeta \) function as a sum over periodic orbits for chaotic Hamiltonian flows. The derivation of the trace formula for quantum graphs which was presented in [1] differs from the original method [6]. It uses another approach (called sometimes the ‘scattering approach’), which reveals in a natural way the underlying classical dynamics.

Discrete graphs, where only the graph topology and not its metric plays a role, are mostly studied in number theory, combinatorics, etc. There is abundant literature relating to various aspects of graphs. Much of the relevant material to the present discussion can be found in [8, 9]. Audrey Terras’ review [10] surveys the field and its relation to quantum chaos. The present work attempts to highlight further this quantum chaos connection, by proposing trace formulæ and spectral \( \zeta \) functions, and linking them with the Ihara \( \zeta \) function [2] and some of its recent generalizations. To introduce these concepts, a few preliminaries and definitions are necessary, and they are provided below.

A graph \( G \) consists of \( V \) vertices connected by \( B \) bonds. The \( V \times V \) connectivity (or adjacency) matrix \( C \) is defined such that \( C_{i,j} = 1(0) \) if the vertices \( i, j \) are connected (disconnected). Graphs with parallel bonds or loops are excluded. The valency (some times referred to as the degree) of a vertex is the number of bonds which emanate from a vertex. It is denoted by \( v_i = \sum_{j=1}^{V} C_{i,j} \). To any bond \( b = (i, j) \) one can assign an arbitrary direction, resulting in two directed bonds, \( d = (i, j) \) and \( \hat{d} = (j, i) \). Thus, the graph can be viewed as \( V \) vertices connected by bonds \( b = 1, \ldots, B \) or by \( 2B \) directed bonds \( d = 1, \ldots, 2B \). (The notation \( b \) for bonds and \( d \) for directed bonds will be used throughout.) It is convenient to associate with each directed bond \( d = (j, i) \) its origin \( o(d) = i \) and terminus \( t(d) = j \) so that \( d \) points from the vertex \( i \) to the vertex \( j \). The bond \( d' \) follows \( d \) if \( t(d') = o(d') \). A periodic orbit (cycle) of length \( n \) is a sequence of \( n \) successively following directed bonds \( d_1, \ldots, d_n \) and \( d_1 \) follows \( d_n \). Cyclic permutations of the bonds generate the same periodic orbit. A primitive periodic orbit is an orbit which cannot be written as a repetition of a shorter periodic orbit. The set of primitive \( n \)-periodic orbits will be denoted by \( P(n) \), and \( P = \bigcup_{n=2}^{\infty} P(n) \). An important subset of \( P(n) \) is the set of \( n \) primitive periodic orbits without backscatter, namely, periodic orbits where \( d_{i+1} \neq d_i \). The corresponding sets will be denoted by \( C(n) \) and \( C = \bigcup_{n=2}^{\infty} C(n) \).

The discrete Laplacian on a graph is defined as

\[
L \equiv -C + D, \tag{1}
\]

where \( C \) is the connectivity matrix and \( D \) is a diagonal matrix with \( D_{i,i} \equiv cdv_i \). It is a self-adjoint operator whose spectrum consists of \( V \) non-negative real numbers. The spectrum is determined as the zeros of the secular function (characteristic polynomial):

\[
Z_L(\lambda) \equiv \det(\lambda I^V - L). \tag{2}
\]

Here, \( \lambda \) is the spectral parameter and \( I^V \) is the unit matrix in \( V \) dimensions. The lowest eigenvalue is 0, and it is simple if and only if the graph is connected.

It is sometimes convenient to generalize the Laplacian (1) by replacing the matrix \( C \) by a matrix \( \tilde{C} \) whose zero entries coincide with those of \( C \), but arbitrary, strictly positive weights \( w_{i,j} (= w_{j,i}) \) replace the values 1 where \( C_{i,j} = 1 \). One then defines \( \tilde{D}_{i,j} \equiv u_i = \sum_j \tilde{C}_{i,j} \) and the generalized Laplacian is

\[
\tilde{L} \equiv -\tilde{C} + \tilde{D}. \tag{3}
\]
The spectrum of $\tilde{L}$ consists of the zeros of the secular equation (characteristic polynomial) $Z_{\tilde{L}}(\lambda) \equiv \det(\lambda I(V) - \tilde{L})$. The spectrum is non-negative, $0$ is in the spectrum and it is a simple eigenvalue if and only if the graph is connected.

The focus of the present work is on $\zeta$ functions and trace formulae for discrete graphs.

This research subject was initiated by Ihara [2] who defined a $\zeta$ function for a graph as

$$\zeta(u)^{-1} = \prod_n (1 - u^n)^{|C(n)|},$$

where $|C(n)|$ is the cardinality of the set $C(n)$, and $u \in \mathbb{C}$ with $|u|$ sufficiently small to ensure the convergence of the infinite product. Following Ihara’s original work, several authors (see, e.g., [11] for a survey of the methods) have proved that

$$\zeta(u)^{-1} = (1 - u^2)^{r-1} \det(I(V) - uC + u^2Q).$$

(5)

Here, $r \equiv B - V + 1$ is the rank of the graph (the number of independent cycles on the graph or, equivalently, the rank of its fundamental group). $I(V)$ is the unit matrix in $V$ dimensions, $C$ is the connectivity matrix and the diagonal matrix $Q \equiv D - I(V)$. If the graph is $v$-regular, that is $v_i = v \forall i$, the non-trivial poles of the Ihara $\zeta$ (the trivial poles are at $u = \pm 1$) can be easily computed from the eigenvalues of the graph Laplacian (1).

The following $\zeta$ function defined by Stark [12] will serve as an example of the more recent developments in the field. Consider a matrix $Y$ in the space of directed bonds

$$Y_{d', d} \equiv \eta_{d', d} \delta_{d, d'} \delta_{d, 0}(1 - \delta_{d', d}),$$

(6)

where $\eta_{d', d}$ are arbitrary. Note that matrix elements between reversed bonds are excluded. Associate with any primitive periodic orbit $c \in C$ the amplitude

$$f_c \equiv \eta_{d_c, d_{c-1}} \eta_{d_{c-1}, d_{c-2}} \cdots \eta_{d_2, d_1} \eta_{d_1, d_c}.$$  

(7)

Then,

$$\zeta_E(Y)^{-1} = \prod_{c \in C} (1 - f_c) = \det(I^{(2B)} - Y),$$

(8)

where $I^{(2B)}$ is the unit matrix in $2B$ dimensions. This result will be used in the last section.

In the next section, other $\zeta$ functions are defined, discussed and expressed as rational functions which are reminiscent of (5) and (8), but are different in many respects. Trace formulae for the spectra of the Laplacians (1), (3) will also be derived. In the last section, the approach developed here will be compared with its analogues in the theory of quantum graphs, and the ‘classical dynamics’ on the discrete graph will be proposed.

2. Secular functions, $\zeta$ functions and trace formulae

To start, an alternative form of the secular equations for the Laplacians (1), (3) will be derived. It is convenient to begin with a detailed derivation for the traditional Laplacian (1). The necessary modifications for the generalized form will be indicated later. For both Laplacians, the secular function will be shown to take the form

$$Z_S(\lambda) = \frac{1}{2^B} \left( \det U(\lambda) \right)^{-\frac{1}{2}} \det(I^{(2B)} - U(\lambda))$$

(9)

where $U(\lambda)$ is a unitary matrix of dimension $2B$ which depends on the spectral parameter $\lambda$. By construction, $Z_S(\lambda)$ is real for $\lambda \in \mathbb{R}$, and its zeros will be shown to coincide (with their multiplicity) with the spectrum of the Laplacian. Thus, $Z_S(\lambda)$ and $Z_L(\lambda)$ can differ at most by a multiplicative function of $\lambda$ which does not vanish for real $\lambda$. This construction of the secular function paraphrases the ‘scattering approach’ introduced in [1] for quantum graphs.
To compute an eigenvector \( \psi = (\psi_1, \ldots, \psi_V) \) of \( L \), corresponding to an eigenvalue \( \lambda \), the following steps are taken. To each bond \( b = (i, j) \) one associates a bond wavefunction

\[
\psi_b(x) = a_b e^{i \frac{\pi}{4} x} + a_{\hat{b}} e^{-i \frac{\pi}{4} x}, \quad x \in \{\pm 1\}
\]  

subject to the condition

\[
\psi_b(1) = \psi_i, \quad \psi_b(-1) = \psi_j. \tag{11}
\]

Consider any vertex indexed by \( i \) and the bonds \((b_1, b_2, \ldots, b_{v_i})\) which emanate from \( i \). The corresponding bond wavefunctions have to satisfy three requirements in order to form a proper eigenvector of \( L \).

(I) **Uniqueness.** The value of the eigenvector at the vertex \( i, \psi_i \), computed in terms of the bond wavefunctions is the same for all the bonds emanating from \( i \). The following \( v_i - 1 \) independent equalities express this requirement:

\[
a_{b_1} e^{i \frac{\pi}{4}} + a_{\hat{b}_1} e^{-i \frac{\pi}{4}} = a_{b_2} e^{i \frac{\pi}{4}} + a_{\hat{b}_2} e^{-i \frac{\pi}{4}} = \cdots = a_{b_{v_i}} e^{i \frac{\pi}{4}} + a_{\hat{b}_{v_i}} e^{-i \frac{\pi}{4}}. \tag{12}
\]

(II) **\( \psi \) is an eigenvector of \( L \).** At the vertex \( i, \sum_{j=1}^{v_i} L_{i,j} \psi_j = \lambda \psi_i \). In terms of the bond wavefunctions this reads

\[
- \sum_{l=1}^{v_i} [a_{b_l} e^{-i \frac{\pi}{4}} + a_{\hat{b}_l} e^{i \frac{\pi}{4}}] = (\lambda - v_i) \frac{1}{v_j} \sum_{m=1}^{v_j} [a_{b_m} e^{i \frac{\pi}{4}} + a_{\hat{b}_m} e^{-i \frac{\pi}{4}}]. \tag{13}
\]

To get the equation above, \( \psi_i \) was presented as

\[
\psi_i = \frac{1}{v_i} \sum_{j=1}^{v_i} (a_{b_j} e^{i \frac{\pi}{4}} + a_{\hat{b}_j} e^{-i \frac{\pi}{4}}). \tag{14}
\]

Together (12) and (13) provide \( v_i \) homogeneous linear relations between the \( 2v_i \) coefficients \( a_d \), where \( d \) stand for directed bonds which are either incoming to \((i(d) = i)\) or outgoing from \((o(d) = i)\) the vertex \( i \). Using these equations, the outgoing coefficients are expressed in terms of the incoming ones,

\[
a_d = \sum_{d' : o(d') = i} \sigma_{d,d'}^{(i)}(\lambda) a_{d'}, \quad \forall \ d : o(d) = i, \tag{15}
\]

where,

\[
\sigma_{d,d'}^{(i)}(\lambda) = i \left( \delta_{d,d'} - \frac{2}{v_i} \frac{1}{1 - i \left( 1 - \frac{1}{v_i} \right)} \right) = i \left( \delta_{d,d'} - \frac{1}{v_i} \left( 1 + e^{i \nu_i(\lambda)} \right) \right)
\]

\[
e^{i \nu_i(\lambda)} = \frac{1 + i \left( 1 - \frac{1}{v_i} \right)}{1 - i \left( 1 - \frac{1}{v_i} \right)}. \tag{16}
\]

The vertex scattering matrices \( \sigma^{(i)}(\lambda) \) are the main building blocks of the present approach. They distinguish clearly between backscatter transitions \((\hat{d} = d')\) and the transitions to other bonds, for which the same strength is given, independently of the original and the final bonds. A straightforward computation shows that for real \( \lambda \) the vertex scattering matrices are unitary matrices. They are the discrete analogues of the vertex scattering matrices derived for the Schrödinger equation on graphs [1].

(III) **Consistency.** The linear relation between the incoming and the outgoing coefficients (15) must be satisfied simultaneously at all the vertices. However, a directed bond \((i, j)\) when
observed from the vertex $j$ is outgoing, while when observed from $i$ it is incoming. This consistency requirement is implemented by introducing the evolution operator $U_{d',d}(\lambda)$ in the $2B$-dimensional space of directed bonds,

$$U_{d',d}(\lambda) = \delta_{d'(1),d(1)}\sigma_{d',d}(\lambda).$$

($U$ is also referred to in the literature as the bond scattering matrix [1]). The evolution operator is unitary $UU^\dagger = I(2B)$ for $\lambda \in \mathbb{R}$ due to the unitarity of its constituents $\sigma(i)$. Denoting by $a$ the $2B$-dimensional vector of the directed bonds coefficients $a_d$ defined above, the consistency requirement reduces to

$$U(\lambda)a = a.$$  \hspace{1cm} (18)

This can be satisfied only for those values of $\lambda$ for which

$$\xi(\lambda) \equiv \det(I(2B) - U(\lambda)) = 0.$$  \hspace{1cm} (19)

For real $\lambda$ the spectrum of $U(\lambda)$ is restricted to the unit circle. Therefore $|\xi(\lambda)|$ is finite for all $\lambda \in \mathbb{R}$. Due to (16) the matrix elements of $U(\lambda)$ are the ratios of monomials in $\lambda$. These two properties imply that $\xi(\lambda) = p(\lambda)/q(\lambda)$ where $p$ and $q$ are polynomials of the same degree in $\lambda$, and their degree is at most $2B$. The zeros of $q(\lambda)$ coincide with the poles of $\det U(\lambda)$. They are complex because $|\det U(\lambda)| = 1$ for $\lambda \in \mathbb{R}$. A straightforward computation yields

$$\det U(\lambda) = \prod_{j=1}^V \frac{1 + i(1 - \frac{\lambda}{v_j})}{1 - i(1 - \frac{\lambda}{v_j})}, \quad \Rightarrow \quad q(\lambda) = \text{const} \prod_{j=1}^V \left(1 - i \left(1 - \frac{\lambda}{v_j}\right)\right).$$  \hspace{1cm} (20)

Thus, $\det U$ has exactly $V$ complex poles, implying that the degree of $p(\lambda)$ which equals the degree of $q(\lambda)$ is also $V$. Note finally that the zeros of $p(\lambda)$ coincide with the zeros of the secular function $Z_L(\lambda) = \det(\lambda I(V) - L)$ which is also a polynomial of degree $V$. Hence, $p(\lambda)$ and $Z_L(\lambda)$ are identical up to a constant factor. It is convenient to define the secular equation so that it is real on the real axis. This can be achieved by multiplying $\xi(\lambda)$ by $(\det U(\lambda))^{-\frac{1}{2}}$. A further factor of $2^{-B}$ normalizes the resulting function to approach 1 as $|\lambda| \to \infty$. The resulting secular equation reads

$$Z_S(\lambda) = \frac{1}{2^B} (\det U(\lambda))^{-\frac{1}{2}} \det(I(2B) - U(\lambda))$$

$$= \frac{1}{2^B} \prod_{j=1}^V \left(1 + i\left(1 - \frac{\lambda}{v_j}\right)\right)^{-\frac{1}{2}} p(\lambda) q(\lambda) = \frac{\det(\lambda I(V) - L)}{\prod_{j=1}^V \left(v_j^2 + (v_j - \lambda)^2\right)^{\frac{B}{2}}}.$$  \hspace{1cm} (21)

This expression for the secular equation is the basis for the further results of the present work. To begin, use is made of the fact that the spectrum of $U(\lambda)$ for $\text{Im}(\lambda) < 0$ is confined to the interior of the unit circle. Thus, for any $\lambda$ with an arbitrarily small (but finite) negative imaginary part we expand

$$\log \det(I(2B) - U(\lambda)) = -\sum_{n=1}^\infty \frac{1}{n} \text{tr} U^n(\lambda),$$  \hspace{1cm} (22)

and

$$\text{tr} U^n(\lambda) = \sum_{m|m|n} \sum_{p \in P(m)} a_p(\lambda).$$  \hspace{1cm} (23)

The sum above is over all the primitive periodic orbits $p$ with period $m$ which is a divisor of $n$, $p = d_1, \ldots, d_m$ and

$$a_p(\lambda) = \sigma_{d_1,d_m}(\lambda) \cdots \sigma_{d_2,d_1}(\lambda).$$  \hspace{1cm} (24)
The explicit dependence of $a_p(\lambda)$ on $\lambda$ is obtained from the following expressions for the vertex scattering matrix elements:

$$
\sigma_{d',d} = \begin{cases}
\frac{4}{v_j^2 + (v_j - \lambda)^2} & \text{for } d' \neq \hat{d}, \\
1 - \frac{4(v_j - 1)}{v_j^2 + (v_j - \lambda)^2} & \text{for } d' = \hat{d},
\end{cases}
$$

$$
\left[ \frac{4}{v_j^2 + (v_j - \lambda)^2} \right]^{1/2} e^{i \arctan(1 - \frac{z}{\gamma})/2} \text{ for } d' \neq \hat{d},
$$

$$
\left[ 1 - \frac{4(v_j - 1)}{v_j^2 + (v_j - \lambda)^2} \right]^{1/2} e^{-i \arctan \frac{2v_j - \lambda}{v_j^2 - v_j - \lambda^2}} \text{ for } d' = \hat{d},
$$

where $j = t(d) = o(d')$. The explicit expressions above were written so that for real $\lambda$, the absolute square of $a_p$ is a product of ‘transition probabilities’, while the phase of $a_p$ is a sum which plays the role of the ‘action’ or ‘length’ associated with the periodic orbit. Substituting (23) into (22) and summing over the repetition numbers $\frac{n_m}{m}$ one gets

$$
\det(I(V) - U(\lambda)) = \prod_{p \in P} (1 - a_p(\lambda)).
$$

The $\zeta$ function which is introduced in the present work is defined as

$$
\zeta_S(\lambda)^{-1} = \prod_{p \in P} (1 - a_p(\lambda)).
$$

Combining (22) and (26) with the definition of $\zeta_S(\lambda)$ gives

$$
\zeta_S(\lambda)^{-1} = \frac{\det(\lambda I(V) - L)}{\prod_{\text{min}(v_j + i(v_j - \lambda))}}.
$$

This is one of the main results of the present work. It provides a ‘Ihara’-like identity which expresses an infinite product over primitive periodic orbits on the graph in terms of the characteristic polynomial of the graph discrete Laplacian. The main difference is that here all the periodic orbits, including orbits with backscattering, contribute to the product and that the amplitudes $a_p$ depend on the spectral parameter in a more complicated way. To get a closer look at the $\zeta_S$ function and its relation to the Ihara $\zeta$ function, it is instructive to write $\zeta_S$ for a general $v$-regular graph. For this purpose, it is convenient to define a new complex variable

$$
z = \frac{1 + i(1 - \frac{z}{\gamma})}{1 - i(1 - \frac{z}{\gamma})}
$$

which is a $1 \leftrightarrow 1$ map of $\mathbb{R}$ to the unit circle in $\mathbb{C}$. With these simplifications, $\zeta_S(z)$ (28) reduces to

$$
\zeta_S(z)^{-1} = \left( \frac{2z}{z + 1} \right)^v \det \left( C + iv \frac{z - 1}{z + 1} I(V) \right).
$$

It is convenient to define $\gamma_S(z) = z^\frac{v}{2} \zeta_S(z)$, in terms of which a functional equation for $\zeta_S$ can be written as

$$
\gamma_S(z^{-1}) = (\gamma_S(z^*)^*)^*,
$$

where $(\cdot)^*$ stands for complex conjugation. Functional equations of similar type are satisfied also by the Ihara $\zeta$ function (for $v$-regular graphs) as well as by most other functions of this genre. Typically, functional equations enable the analytical continuation of $\zeta$ functions which are defined by infinite products beyond their radius of convergence. Here, it also provides the analytic continuation of $\zeta_S(z)$ to the exterior of the unit disc.
The periodic amplitudes $a_p(z)$ simplify considerably for $v$-regular graphs. Denote by $n_p$ the period of the primitive periodic orbit $p$ and by $\beta_p$ the number of vertices in $p$ where backscattering occurs: $\beta_p = \sharp \{ i : d_i = d_{i+1}, d_i \in p, d_{n_p+1} = d_1 \}$. Then,

$$a_p(z) = e^{-i\frac{\pi}{2}n_p} \left( 1 + \frac{z}{v} \right)^{n_p-\beta_p} \left( -1 \right)^{\beta_p} \left( 1 - \frac{1 + z}{v} \right)^{\beta_p}.$$(32)

The results above pave the way to the derivation of trace formulae for the discrete Laplacians. Trace formulae provide a powerful tool in spectral theory. They express the spectral density (written formally as a sum of Dirac $\delta$ functions located at the spectral set) in terms of information derived from the manifold metric. The spectral density is written as a sum of two contributions—both of which have a geometric origin. The first is a smooth function of $\lambda$, whose asymptotic limit at $\lambda \to \infty$ was first studied by Weyl. The second contribution is an infinite sum over periodic geodesics on the manifold. The equality between the spectral density and its geometric representation should be understood only in the sense of distributions. An analogous trace formula will be derived now for the discrete Laplacian. Making use of Cauchy theorem and the fact that $Z_S(\lambda)$ is analytic in the vicinity of the real $\lambda$ axis, and real on it, one can write

$$d(\lambda) = \sum_{j=1}^{V} \delta(\lambda - \lambda_j) \quad (33)$$

Using

$$Z_S(\lambda) = \frac{1}{Z^B} (\det U(\lambda))^{-\frac{1}{2}} \det(I^{(2B)} - U(\lambda)), \quad (35)$$

the explicit form of $\det U(\lambda) (20)$ and the periodic orbit expansion (23), one gets

$$d(\lambda) = \frac{1}{\pi} \sum_{j=1}^{V} \frac{1}{\sqrt{1 + (1 - \frac{1}{n_p})^2}} \frac{1}{\sqrt{1 - i (1 - \frac{1}{n_p})}} \sum_{r=1}^{\infty} \sum_{p \in P} n(p) |a_p(\lambda)|^r e^{i\phi_p(\lambda)}. \quad (36)$$

The first term is the ‘smooth’ (Weyl) contribution to the spectral density. It consists of a sum of Lorenzians with poles at $\lambda_j = v_j (1 \pm i)$. The second term is an infinite sum over periodic orbits analogous in structure to the ‘fluctuating’ contributions which appear in any trace formula of this kind. It can be written explicitly using (25). Noting that the $a_p(\lambda)$ are complex numbers with $\lambda$ dependent phases $\phi_p(\lambda)$, the periodic orbit sum in the trace formula is a fluctuating function of $\lambda$. It is the term which turns the rhs of (36) to a distribution when $\epsilon \to 0$.

So far, the discussion was restricted to the Laplacians (1). The extension to the generalized Laplacians starts by modifying the definition of the bond wavefunctions (10) to read

$$\psi_b = \sqrt{w_b} (a_b e^{\frac{i}{2} \lambda} + a^* b e^{-\frac{i}{2} \lambda}). \quad (37)$$

Then, following the same steps as above, the vertex scattering matrices are derived, and they take the form

$$\tilde{\sigma}^{(i)}_{d,d'}(\lambda) = i \left( \delta_{d,d'} \right) \left( 1 + e^{i\omega(\lambda)} \right) \sqrt{w_d w_{d'}}; \quad e^{i\omega(\lambda)} = \frac{1 + i(1 - \frac{1}{n_p})}{1 - i(1 - \frac{1}{n_p})}, \quad (38)$$
where \( u_j = \sum_{j} w_{i,j} \) as defined previously. The subsequent derivation follows the same steps, resulting in the generalized \( \xi_{S} \) function:

\[
\xi_{S}(\lambda) \equiv \prod_{p \in P} \left( \frac{1}{1 - a_p(\lambda)} \right) = \det(\lambda \mathcal{I}(V) - \tilde{\mathcal{L}}) \prod_{V} \frac{1}{(u_j + i(u_j - \lambda))}. 
\] (39)

A trace formula is also derived in the same way:

\[
d(\lambda) = \sum_{j=1}^{V} \delta(\lambda - \tilde{\lambda}_j) = \frac{1}{\pi} \lim_{\epsilon \to 0^+} \Im \frac{d}{d\lambda} \log Z_{S}(\lambda - i\epsilon) = \frac{1}{\pi} \sum_{j=1}^{V} \frac{1}{1 + \left(1 - \frac{\lambda}{u_j}\right)^2} - \frac{1}{\pi} \sum_{r=1}^{\infty} \sum_{p \in P} \frac{1}{n(p)} |a_p|^2 e^{i\phi_p(\lambda)}. 
\] (40)

The expressions for \( a_p(\lambda) \) can be derived by a simple modification of (25) and therefore they will not be written here.

3. Classical dynamics

The present approach emerges from the alternative secular function for the spectrum of Laplacians, based on the quantum evolution operator \( U(\lambda) \) in the space of directed bond amplitudes \( \mathbf{a} \in l^2(\mathbb{C}^2)^B \). Consider \( U(\lambda) \) as a quantum map which maps this \( 2B \)-dimensional space onto itself. \( U \) is unitary and hence the map conserves the \( l^2 \) norm—the quantum probability. The condition \( U(\lambda) \mathbf{a} = \mathbf{a} \) can be interpreted as a requirement that \( \lambda \) is an eigenvalue if there exists a non-trivial vector \( \mathbf{a} \) which is stationary under the action of the quantum map [17]. The requirement of stationarity is naturally associated with the eigenvalue being in the spectrum of the underlying Hamiltonian.

The building blocks for the theory are the vertex scattering matrices. Similar matrices appear in the theory of quantum graphs. There they emerge when the Schrödinger equation on the graph is augmented by vertex boundary conditions which render the resulting operator self-adjoint. The self-adjoint extension is not unique and depends on the parameter \( k \) and on an arbitrary parameter \( \kappa \) which interpolates between the ‘Dirichlet’ (\( \kappa = 0 \)) and the ‘Neumann’ (\( \kappa = \infty \)) boundary conditions [1, 18]. The scattering matrices for discrete graphs are obtained from their quantum graph analogues by replacing \( \kappa/k \) by \( \lambda \).

The unitary quantum evolution operator is the starting point for the construction of a classical evolution on the discrete graph. The classical ‘phase space’ in this case is the probability vectors \( \rho \in l^2(\mathbb{R}^2)^B \) where the components of \( \rho \) are interpreted as the probabilities to find the classical system on the corresponding directed bonds. The classical transition matrix is constructed from the quantum probability to make a transition from \( d \) to \( d' \):

\[
M_{d,d'} = |U_{d,d'}|^2. 
\] (41)

The unitarity of \( U \) implies that \( M \) is bi-stochastic, namely, \( \sum_{d} M_{d,d} = \sum_{d'} M_{d,d} = 1 \). This transition matrix induces a discrete, random walk dynamics in phase space. If \( n \) denotes the discrete ‘time’,

\[
\rho(n+1) = M \rho(n). 
\] (42)

This Markovian evolution preserves the \( l^1 \) norm—the classical probability. The spectrum of \( M \) is confined to the interior of the unit circle. 1 is always an eigenvalue corresponding to an eigenvector with equal components which describes the system in an equilibrated state. When
the eigenvalue 1 is the only eigenvalue on the unit circle, the classical dynamics drives the system to equilibrium at a rate which depends on the distance of the next highest eigenvalue to the unit circle. This classical dynamics is identical to that which was introduced in the study of quantum graphs [1]. It plays an important role in the theory of spectral statistics on quantum graphs [1, 19–21]. Finally, the analogue of the Ruelle $\zeta$ function for the evolution induced by $P$ can be easily written starting with the secular function

$$Z_M(\mu) \equiv \text{det}(I^{(2B)} - \mu M). \quad (43)$$

The periodic orbit sum is identical to the trace formula (36) in which the amplitudes $a_p$ are replaced by their absolute squares.

To emphasize the intricate connections between the concepts developed here and their predecessors [2–4], consider a $v$-regular graph ($v > 2$), and the classical evolution operator obtained for the spectral parameter $\lambda = v + i(v - 2)$, corresponding to $z = v - 1$ in (29). At this value, $\sigma_d, d = 0$ and $\sigma_{d', d} = 1$ for $d' \neq d$. The resulting classical evolution matrix $M^2$ needs to be multiplied by $(v - 1)^{-1}$ to make it a legitimate (probability conserving) evolution operator. The resulting evolution does not permit backscatter and, therefore, the secular equation (43) can be computed using Stark’s $\zeta$ function (8), with $Y = \frac{\mu}{\mu^2 + 1}M^2$. The product over the set of non-backscattering primitive periodic orbits becomes identical to that appearing in the Ihara zeta function (4). Using (5), one finally gets

$$Z_{M^2}(\mu) = \prod_n \left(1 - \left(\frac{\mu}{v - 1}\right)^n\right)^{C(n)}$$

$$= \left(1 - \left(\frac{\mu}{\mu^2 + 1}\right)^{v-1}\right) \text{det} \left(I^{(\Psi)} \left(1 + \frac{\mu^2}{\mu^2 + 1}\right) - \frac{\mu}{v - 1}C\right). \quad (44)$$

Thus, the spectrum $m_j$ of $M^2$ consists of $(r - 1)$-fold degenerate eigenvalues at $m_j^{(\pm)} = \pm \frac{1}{v - 1}$ and the rest which can be computed from the spectrum of the discrete Laplacian $\lambda_j$:

$$m_j^{\pm} = \frac{(v - \lambda_j) \pm \sqrt{(v - \lambda_j)^2 - 4(v - 1)}}{2(v - 1)}. \quad (45)$$

The eigenvalue 0 of the Laplacian corresponds to the eigenvalues 1 and $\frac{1}{v - 1}$ of $M^2$. The gap in the classical evolution spectrum is determined by the first non-zero eigenvalue of $L$.

The comment above may have interesting and novel consequences going beyond its anecdotal appearance. In quantum graphs, one can choose vertex scattering matrices from a much larger variety than offered by the vertex scattering matrices (16). Thus, it is possible to construct vertex scattering matrices which do not scatter backwards, but with equal scattering probability to the other vertices. The unitarity is maintained by a proper choice of the phases of the scattering amplitudes [23]. In such cases, and for $v$-regular graphs, the classical analogues are identical with $M^2$ and (44) is applicable. Working with such systems is particularly interesting because in quantum chaos the gap between the eigenvalue 1 and the rest of the spectrum determines whether the spectrum of the $U$ matrix (and hence of the Schrödinger operator) displays the statistics predicted by random matrix theory, in the limit of large graphs. For non-backscattering dynamics (44) reduces the problem to the study of the spectrum of the Laplacian. The behaviour of the gap in the Laplacian spectrum of large graphs is an important subject in the theory of discrete graphs and number theory, related amongst others to the Ramanujan conjecture [22]. A detailed discussion of this connection will take the present manuscript far afield, and it is deferred to a future publication.
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