

GRAPH HOMOLOGY AND STABILITY OF COUPLED OSCILLATOR NETWORKS*

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Abstract. There are a number of models of coupled oscillator networks where the question of the stability of fixed points reduces to calculating the index of a graph Laplacian. Some examples of such models include the Kuramoto and Kuramoto–Sakaguchi equations as well as the swing equations, which govern the behavior of generators coupled in an electrical network. We show that the index calculation can be related to a dual calculation which is done on the first homology group of the graph, rather than the vertex space. We also show that this representation is computationally attractive for relatively sparse graphs, where the dimension of the first homology group is low, as is true in many applications. We also give explicit formulae for the dimension of the unstable manifold to a phase-locked solution for graphs containing one or two loops. As an application, we present some novel results for the Kuramoto model defined on a ring and compute the longest possible edge length for a stable solution.

Key words. spectral graph theory, dynamics on networks, graph Laplacian, coupled oscillators, Kuramoto model

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1. Introduction. Let $G = (V, E, \Gamma = \{\gamma_{v,w}\}_{v,w \in V})$ be a weighted graph. The Laplacian matrix¹ of G is the $|V| \times |V|$ matrix whose components are

$$(\mathcal{L}_G)_{vw} = \begin{cases} \gamma_{vw}, & v \neq w, \\ -\sum_u \gamma_{vu}, & v = w. \end{cases}$$

A classical result is that if $\gamma_{vw} \geq 0$, then \mathcal{L}_G is negative semidefinite, with the number of zero eigenvalues equal to the number of connected components of the graph G . If some of the γ_{vw} are negative, then \mathcal{L}_G can have positive eigenvalues as well. In [1, 2], the authors consider the problem where some subset of the γ_{vw} is negative, and show upper and lower bounds on the number of positive eigenvalues of \mathcal{L}_G that depend only on the “sign topology” of the graph, i.e., the pattern of which edges are positive and which are negative, regardless of their weights.

More specifically, the main result of [1] is as follows: assume that the graph is symmetric ($\gamma_{(v,w)} = \gamma_{(w,v)}$), and define G_+ (resp., G_-) to be the subgraphs of G containing only edges with positive (resp., negative) weights. Let $|G|$ be the number of components of a graph G . If $n_+(G)$ is the number of positive eigenvalues of \mathcal{L}_G , then

$$(1.1) \quad |G_+| - 1 \leq n_+(G) \leq |V| - |G_-|.$$

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¹Our Laplacian is the negative of the standard definition, but the reasons for this will be clear later.

Moreover, one can show that these bounds are saturated, i.e., for any integer in that range, there is some choice of weights that give exactly that many positive eigenvalues. (In fact, one can say much more about the genericity of the sets of weights that give these values; see [2] for more detail.)

It follows from the formula above that if G is a tree, then the number of positive eigenvalues is exactly the number of negatively weighted edges in the graph. To see this, notice that, for a tree, the number of connected components of G_+ is exactly one plus the number of negative edges: If we start with a tree with only positive edges, every flip of an edge from positive to negative will add another component to the positive subtree. Conversely, since the negative edges cannot form a cycle, the number of connected components of G_- is the number of vertices minus the number of negatively weighted edges. In short, if G is a tree, then $n_+(G)$ is independent of the weights of the edges, and very easy to compute by inspection.

Now, consider the case where the graph has one cycle and one negative edge. It is easy to see in this case that the upper bound in (1.1) is one, and the lower bound is either zero (if the negative edge lies in the cycle) or one (if the negative edge does not lie in the cycle). In the latter case, the number of positive eigenvalues is again independent of the weights. In the former case, the weight on the edge matters, and it is not hard to imagine that the relative magnitude of the weight on the negative edge compared to the other edges in the cycle determines the stability.

From these examples, it is at least plausible to conjecture that the stability computation has complexity that scales with the number of cycles. This is further supported by previous work of two of the authors [1], where the following result was established. Let $H_1(G)$ denote the additive group of cycles in the graph, defined as follows: For a graph G , choose and fix an orientation of the edges of G . A cycle in the graph is an assignment of integer weights to each edge in the graph such that, at every vertex, the sum of the incoming weights is equal to the sum of the outgoing weights. This defines a closed cycle on the graph in the natural manner, where the path traverses each edge a number of times equal to the weight on the edge. The addition of such paths is natural, and this defines an Abelian group. One of the main results of [1] is that if we fix the signs of the weights on each edge, but otherwise allow them to vary, then the difference between the largest possible value of $n_+(G)$ and the smallest possible value of $n_+(G)$ is equal to the dimension of the quotient of $H_1(G)$ by a certain subgroup. In particular, one quotients by all paths that travel only on positive or only on negative edges, so that the quotient space is spanned by cycles that have both signs. In short, counting the number of independent mixed cycles gives the difference between the upper and lower estimates in (1.1).

The result of the current paper is that the complexity in computing $n_+(G)$ for any particular choice of weights is exactly a function of the number of cycles in the graph. More specifically: If we denote the number of cycles by $C = |E| - |V| + 1$, then the main result of this paper is that the computation of $n_+(G)$ is equivalent to computing the index of a particular $C \times C$ matrix (what we call the “cycle intersection matrix”). Note that the naive approach to computing $n_+(G)$ is to compute the spectrum of a $|V| \times |V|$ matrix, so when $C \ll |V|$, this is clearly a much simpler problem.

1.1. Connection to nonlinearly coupled phase oscillators. As stated above, the problem we consider is a pure linear algebra problem and can be stated as such, but in fact it is inspired and informed by considering its “killer app,” i.e., understanding the dimension of the unstable manifold of a fixed point of a nonlinearly coupled network dynamical system.

Some examples of the kinds of models we are interested in include the Kuramoto problem on a general weighted connected graph $D = \{\delta_{ij}\}$, namely

$$(1.2) \quad \frac{d\theta_i}{dt} = \omega_i + \sum \delta_{ij} \sin(\theta_j - \theta_i).$$

More generally one can also consider the Kuramoto-Sakaguchi model [3, 4]

$$(1.3) \quad \frac{d\theta_i}{dt} = \omega_i + \sum \delta_{ij} \sin(\theta_j - \theta_i + \alpha),$$

as well as the swing equations, which govern the dynamics of a coupled system of generators and loads [5]

$$(1.4) \quad M \frac{d^2\theta_i}{dt^2} + D \frac{d\theta_i}{dt} = \omega_i + \sum \delta_{ij} \sin(\theta_j - \theta_i).$$

In each of these examples the dimension of the unstable manifold to a fixed point of the equations is given by the number of positive eigenvalues of a graph Laplacian.

The study of the fixed points of these systems, and their stability, has led to a huge number of papers in the literature [6], [7], [8], [9], [10], [11], [12], [13], [14], [15], [16], [17], [18], [19], [20], [21], [22], [23]. It has been widely observed that the behavior of solutions varies dramatically as the topology of the graph changes: These models are essentially trivial on a tree, but become more complex as the graph becomes denser. This again suggests that the number of loops in the graph should be important.

To date, most of the analytical work has focused on special graph configurations with a high degree of symmetry. The most-studied case, where the behavior of the model is very well understood, is the case of “all-to-all” coupling, where the underlying graph is the complete graph K_n , and all of the edge weights δ_{ij} are equal. A ground breaking paper [19] gave upper and lower bounds (which differed by 1) on the dimension of the unstable manifold to an arbitrary fixed point solution to this model. The authors of [24] gave a sufficient condition for the existence of a stable phase-locked solution, and two of the current authors [25] gave an explicit formula for the dimension of the unstable manifold. Similarly, [16] gives a proof of the existence of a critical coupling constant for the cases of the complete and bipartite graphs, and one of the current authors [26] gives some results on the minimizers and saddles for the Kuramoto on cyclic graphs. For general topologies, it was observed in [27] that if the graph is connected and all of the links are short (the angle differences are less than $\pi/2$), then the fixed point is stable. This was generalized in [28], where it was shown that there could be neutral links (angle differences of exactly $\pi/2$), as long as the induced subgraph given by the short links was connected, that also give a stable fixed point. This calculation corresponds to showing that a Laplacian with nonnegative links is stable (cf. (4.1), (4.2)), so in this language we are studying the stability of fixed points that have long links.

Finally, we want to point out papers on Laplacians directly related to the current work. In [29], the authors give a condition for the definiteness of a signed Laplacian (in our language, a condition that $n_+(G) = 0$) based on the effective resistance of the positive entries of the Laplacian of the graph. Those authors give a necessary and sufficient condition for definiteness in some cases, and we show below that at least in one example the two papers obtain the same bounds. However, the two papers ([29] and the current paper) generalize the problem in different ways, so that there are questions that each paper addressed that the other one does not. There is also

the work of [30] which studies the definiteness of a linear consensus problem, and the results again look similar, but we stress that here the framework is different. Here we choose the Laplacian to always have a zero row sum, but in [30] the diagonal term is chosen to be $-\sum |\gamma_{uv}|$ instead, more in line with the classical work on signed graphs done by [31], [32], and others.

We show in this paper that the intuition put forward in the preceding paragraphs is correct. We show that, rather than computing the number of positive eigenvalues of the Laplacian, an object that acts on the vertex space of the graph, there is a dual calculation that can be posed on the cycle space. This dual problem gives an alternative method for computing the dimension of the unstable manifold to a particular fixed point. This approach is particularly attractive for problems in which the graph is relatively sparse, and the dimension of the cycle space is much smaller than the dimension of the vertex space. We note that there are some very important applications for which this condition is satisfied. One example is the graphs associated with electrical power networks. Such networks are governed by the swing equations (1.4) and these tend to be reasonably sparse. The following data is from Cotilla-Sanchez et al. [33], who calculated the number of edges and vertices for the IEEE 300 model (a standard test power system) as well as the Eastern (EI), Western (WI), and Texas (TI) interconnects, the largest components of the North American electrical network:²

TABLE 1
Sparsities of various electrical networks.

Network	Edges E	Vertices N	Number of cycles $C = E - N + 1$	Sparsity $\frac{C}{N}$
IEEE 300	409	300	110	.37
EI	52075	41228	10848	.26
WI	13734	11432	2303	.2
TI	5532	4513	1020	.23

As can be seen from Table 1, electrical networks tend to be quite sparse, with the dimension of the cycle space being 20–40% of dimension of the vertex space. For instance, if one wanted to understand the stability of a fixed point of the Texas Interconnect, the dual problem presented in this paper would reduce the calculation to the eigenvalues of a 1020×1020 matrix, rather than the 4513×4513 eigenvalue problem determined by a naive stability calculation.

2. Summary of results. In this section, we define our notation and state the two main theorems of the paper, and give a sense of the main ideas of the proofs. The complete proof of the main results will be deferred to section 5.

2.1. Definitions.

DEFINITION 2.1. A weighted graph G is the triple $G = (V, E, \Gamma)$, where V is a set of vertices, $E \subseteq V \times V$ is a set of edges, and $\Gamma = \{\gamma_e\}_{e \in E}$ is a set of edge weights. By this definition, the graphs in this paper are always simple (at most one edge between vertices) and undirected. We use the convention throughout that $\gamma_e \neq 0$ for all $e \in E$ (an edge weight of zero corresponds to the edge not being in E , in other words).

²A small amount of processing has been done on this data, including combining parallel edges and removing disconnected buses. Other sources give slightly different numbers without changing the general picture.

DEFINITION 2.2. Let $G = (V, E, \Gamma)$ be a weighted graph. The Laplacian matrix of G is the $|V| \times |V|$ matrix whose entries are given by

$$(\mathcal{L}_G)_{v,w} = \begin{cases} \gamma_{(v,w)}, & v \neq w, \\ -\sum_u \gamma_{(v,u)}, & v = w. \end{cases}$$

Since the graph is undirected, the Laplacian matrix is always symmetric and has real spectrum. We denote by $n_+(G), n_0(G), n_-(G)$ the number of positive, zero, and negative eigenvalues of \mathcal{L}_G , and refer to this triple as the spectral index of G . Since, for any G , $\mathcal{L}_G \mathbf{1} = \mathbf{0}$, \mathcal{L}_G has a zero eigenvalue and thus a zero determinant. We will colloquially refer to a Laplacian as stable if $n_+(\mathcal{L}_G) = 0$ and unstable otherwise.

DEFINITION 2.3. As mentioned in the previous definition, $\det(\mathcal{L}_G) = 0$ for any G . Let us write the eigenvalues of \mathcal{L}_G as $\lambda_1 = 0, \lambda_2, \dots, \lambda_n$. We define the reduced determinant of \mathcal{L}_G , $\det_{\text{red}}(\mathcal{L}_G)$, to be the product $\lambda_2 \times \dots \times \lambda_n$. Alternatively, we define $\det_{\text{red}}(\mathcal{L}_G)$ as the product of the nonzero eigenvalues of \mathcal{L}_G if \mathcal{L}_G has a simple zero eigenvalue, and zero otherwise.

Remark 2.4. The following are all equivalent definitions of $\det_{\text{red}}(\mathcal{L}_G)$. We will use many of these results in the body of this paper.

- Since \mathcal{L}_G is symmetric and $\mathbf{1} \in \ker(\mathcal{L}_G)$, we can consider \mathcal{L}_G as a linear map from $\mathbf{1}^\perp$ to $\mathbf{1}^\perp$. Then $\det_{\text{red}}(\mathcal{L}_G)$ is the determinant of this restricted linear map.
- $\det_{\text{red}}(\mathcal{L}_G)$ is the linear term in the characteristic polynomial.
- $\det_{\text{red}}(\mathcal{L}_G) = \det(\mathcal{L}_G + \frac{1}{|V|} \mathbf{1} \otimes \mathbf{1})$.

The proofs of all of these are reasonably straightforward exercises in linear algebra.

We also use the following main result.

THEOREM 2.5 (Kirchhoff's matrix tree theorem [34]). Let G be a graph, and let $ST(G)$ be the set of spanning trees (connected acyclic subgraphs) of G . Then we have the formula

$$\det_{\text{red}}(\mathcal{L}_G) = |V| \sum_{T \in ST(G)} \prod_{e \in T} \gamma_e.$$

In other words we take the sum over all spanning trees of the product over all edges in the spanning tree of the edge weights. Alternatively, if $\det_{ij}(\cdot)$ denotes the $|V| - 1 \times |V| - 1$ minor determinant computed by removing the i th row and j th column, then

$$\det_{\text{red}}(\mathcal{L}_G) = |V|(-1)^{i+j} \det_{ij}(\mathcal{L}_G),$$

so any $|V| - 1 \times |V| - 1$ minor (up to a sign) enumerates all of the spanning trees of the graph.

The proof of this theorem is in many graph theory textbooks; see, for instance, the texts of Tutte [35] or Godsil and Royle [36].

DEFINITION 2.6. Let $G = G(V, E, \Gamma)$ be a connected weighted graph. Choose and fix and orient for each edge (so that each edge has both a head and tail vertex). Define the signed incidence matrix of G as the $|V| \times |E|$ matrix whose entries are given by

$$(B_G)_{v,e} = \begin{cases} 1, & e = (v, *), \\ -1, & e = (*, v), \\ 0 & \text{else.} \end{cases}$$

The nullspace of the matrix is called the cycle space. It is a standard result [36] that if G is connected, then the dimension of this nullspace is $C := |E(G)| - |V(G)| + 1$. We call this integer the cycle rank of the graph. Note also that while the definition of B depends on the choice of orientation, the number C does not.

Using the Rank–Nullity theorem, the above is equivalent to saying that B_G^\top has nullity one, a fact we also use below. It is also not hard to see that $B_G^\top \mathbf{1} = \mathbf{0}$, so in fact $\ker(B_G^\top) = \text{span}(\mathbf{1})$ whenever G is connected.

We define the tree set $\text{Tree}(G)$ to be the set of edges of G that are contained in every spanning tree of G . The complement of $\text{Tree}(G)$ is the cycle set $\text{Cycle}(G)$, consisting of every edge which is included in some cycle in the graph. In terms of the incidence matrix, e is in the cycle set if and only if there is a vector $x \in N(B_G)$ with $x_e \neq 0$.

DEFINITION 2.7. Let y_1, y_2, \dots, y_C be a basis for the cycle space. Define the $|E| \times C$ matrix Y_G whose columns are given by the y_i , and let D_G be the $|E| \times |E|$ diagonal matrix where the e th entry is γ_e . Finally, we define the cycle intersection matrix, or the cycle form

$$(2.1) \quad \mathcal{Z}_G := -Y_G^\top D_G^{-1} Y_G,$$

i.e., the matrix with components

$$(\mathcal{Z}_G)_{ij} = - \sum_{e \in E} \gamma_e^{-1} y_{i,e} y_{j,e}.$$

Note that \mathcal{Z}_G is always a $C \times C$ matrix. The exact form of \mathcal{Z}_G will depend on the choice of basis, and we will specify the exact choice of basis in (5.3). For now, we simply note that any property that we need for \mathcal{Z}_G will be independent of this choice.

LEMMA 2.8. Let $G = (V, E, \Gamma)$ be a connected weighted graph. Define D as in Definition 2.7 and B as in Definition 2.6. Then

$$\mathcal{L}_G = -B_G D_G B_G^\top.$$

Proof. Let us first assume that $v \neq w$. Then

$$(B_G D_G B_G^\top)_{v,w} = \sum_{e,e'} B_{ve} D_{ee'} B_{e'w}^\top = \sum_e B_{ve} B_{we} \gamma_e.$$

If v, w are adjacent, then either $e = (v, w)$ or $e = (w, v)$ is an oriented edge. In either case, $B_{ve} B_{we} = -1$. Moreover, if e is an edge not incident on both v and w , then $B_{ve} B_{we} = 0$, and thus the sum has a single term, and $(B_G D_G B_G^\top)_{vw} = -\gamma_e$. Therefore, all of the off-diagonal terms are the same as $-\mathcal{L}_G$.

Finally, notice that since $B_G^\top \mathbf{1} = \mathbf{0}$, we have $B_G D_G B_G^\top \mathbf{1} = \mathbf{0}$, and from this it follows that the diagonal terms must be correct as well. \square

2.2. Theorems. The following are the main results of this paper. The first result gives a formula for the number of positive eigenvalues of the Laplacian in terms of the number of edges with negative edge weights together with the number of positive eigenvalues of the cycle intersection matrix.

THEOREM 2.9. Let $G = (V, E, \Gamma)$ be a connected, weighted graph, and let \mathcal{L}_G be its Laplacian matrix. Then the number of positive eigenvalues of the Laplacian is the

number of negative edges in the graph minus the number of positive eigenvalues of the cycle intersection matrix, i.e.,

$$(2.2) \quad n_+(\mathcal{L}_G) = \#\{e \in E \mid \gamma_e < 0\} - n_+(\mathcal{Z}_G).$$

Recall in the definition of \mathcal{Z}_G that one makes a choice of basis; however, it follows from Lemma A.3 that the index of \mathcal{Z}_G is independent of this choice of basis.

In particular, we have the inequality

$$n_+(\mathcal{L}_G) \leq \#\{e \in \text{Tree}(G) \mid \gamma_e < 0\}.$$

From (2.2), the maximal number of negative edges a graph can have and still have a stable Laplacian is the cycle number C . In the case where the Laplacian is stable and there are exactly C negative edges, then there is choice of basis cycles so that there is exactly one negative edge in each cycle in the basis. In other words if there are exactly C negative edges, then the graph with the negative edges removed must be a tree.

Remark 2.10. Intuitively this theorem makes precise the intuition that having redundancy in a network increases robustness and should tend to stabilize solutions. Redundancy in a network means having multiple paths between vertices, which is exactly the information that is encoded in the cycle space. The first term on the right-hand side of the formula in Theorem 2.9 is the number of instabilities in the absence of redundancy (i.e., for a tree) while the second term counts the number of these instabilities that are suppressed by the redundancy in the network (cf. [29] for a connection in terms of redundancies and effective resistance of alternate paths in the network).

Remark 2.11. The basis for the cycle space can be chosen in the following way: Let G_+ be the subgraph containing only the edges of G with positive weight. The cycle space of G_+ forms a subspace S_+ of the full cycle space of G . Similarly, let G_- be the subgraph containing only negatively weighted edges, and let S_- be the corresponding cycle space. Note that S_+ and S_- are necessarily orthogonal. The basis for the cycle space of the full graph G can be chosen as the union of bases for S_+ and S_- together with a set of “mixed” cycles that cannot be decomposed into a sum of cycles containing only edges of one sign. These mixed cycles are a basis for the quotient of the full group of cycles by the subgroups S_{\pm} . In this basis the cycle intersection form can be written in the following partitioned form:

$$\mathcal{Z}_G = \left(\begin{array}{c|cc} A_M & B_{M,+} & B_{M,-} \\ \hline B_{M,+}^t & \mathcal{Z}_{G_+} & 0 \\ B_{M,-}^t & 0 & \mathcal{Z}_{G_-} \end{array} \right),$$

where \mathcal{Z}_{G_+} and \mathcal{Z}_{G_-} are the cycle intersection forms for the subgraphs G_+ and G_- , respectively.

First, we note that \mathcal{Z}_{G_+} is negative definite and \mathcal{Z}_{G_-} is positive definite, so that $n_+(\mathcal{Z}_G)$ is at least as large as the dimension of the cycle space of G_- and no larger than the dimension of the full cycle space minus the dimension of the cycle space of G_+ . These inequalities are equivalent to the inequalities derived previously by the authors [1].

Second, we note that the calculation of the index of \mathcal{Z}_G can be further reduced to a calculation on the space of mixed cycles. In particular, by applying the Haynsworth theorem to (2.11) we have that

$$n_+(\mathcal{Z}_G) = n_+(A_M - B_{M,+}\mathcal{Z}_{G_+}^{-1}B_{M,+}^t - B_{M,-}\mathcal{Z}_{G_-}^{-1}B_{M,-}^t),$$

and thus the calculation can be reduced to finding the index of a matrix whose size is the number of mixed cycles. However, since the above formula requires inversion of the matrices $\mathcal{Z}_{G_{\pm}}$, it is not obvious that this is more attractive than working on the full cycle space.

The second theorem is a similar result at the level of determinants rather than the spectral index. This result is not really needed for the applications, but it is proven with similar ideas and provides a nice dual form of the famous matrix tree theorem. This is essentially a weighted version of a theorem proven by Sjogren [37], though the proof given here is quite different.

THEOREM 2.12. *Let $G = (V, E, \Gamma)$ be connected, and define \mathcal{L}_G as above. Let $N_G = |V(G)|$, and then*

$$(2.3) \quad \frac{1}{N_G} \det_{\text{red}}(\mathcal{L}_G) = \det(\mathcal{Z}(G)) \times \prod_{e \in E} \gamma_e.$$

The quantity $\det(\mathcal{Z}(G))$ is a homogeneous Laurent polynomial in the edge weights which generates labeled tree complements: each term in the Laurent polynomial represents a set of edges that, when removed from the graph, yield a tree.

Remark 2.13. There is an obvious connection between the two theorems, but they are not quite equivalent since they address different quantities. For example, Theorem 2.12 seems stronger since it gives an actual number, but it only implies Theorem 2.9 modulo 2.

We defer the proof of these results until a later section, but we will briefly summarize the two main ideas here. The first is that for any G , there is a covering tree T (which is a weighted tree), and a graph map $\varphi: T \rightarrow G$, which is surjective on vertices and bijective on edges. In other words if one identifies certain vertices of the tree T , one obtains G . The identification of vertices in the graph corresponds at the level of linear algebra to restriction of the Laplacian to a certain subspace, so the Laplacian on any graph can be realized as the restriction of the Laplacian on a tree to an appropriate subspace.

The second idea is to use the Haynsworth theorem [38], which is stated in Lemma A.1. The Haynsworth theorem is a kind of duality result relating the number of positive eigenvalues of a matrix to the number of positive eigenvalues of a $k \times k$ principle minor matrix and the number of positive eigenvalues of the complementary $(n-k) \times (n-k)$ minor of the inverse matrix. In other words, modulo some invertibility assumptions, if M is self-adjoint and S is a subspace of \mathbb{R}^N , then

$$n_+(M) = n_+(M|_S) + n_+(M^{-1}|_{S^\perp}).$$

Brushing aside some technical points associated with invertibility which will be addressed below, the basic observation is that if we take M to be the Laplacian on the covering tree and $M|_S$ to be the Laplacian on the graph G , then one can compute that $M^{-1}|_{S^\perp}$ is exactly the cycle intersection form \mathcal{Z}_G . For a tree it is easy to see that the number of positive eigenvalues is exactly the number of edges with negative edge weights, and formula (2.2) follows. There is an analogous identity at the level of determinants, giving (2.3).

3. Linear examples.

3.1. Diamond graph. We consider a simple graph with cycle rank $C = 2$, the diamond graph shown in the left frame of Figure 2, with five edges and four vertices. The Laplace matrix is given by

$$\mathcal{L}_G = \begin{pmatrix} -(\gamma_a + \gamma_d + \gamma_e) & \gamma_a & \gamma_e & \gamma_d \\ \gamma_a & -(\gamma_a + \gamma_b) & \gamma_b & 0 \\ \gamma_e & \gamma_b & -(\gamma_b + \gamma_c + \gamma_e) & \gamma_c \\ \gamma_d & 0 & \gamma_c & -(\gamma_c + \gamma_d) \end{pmatrix}.$$

By the matrix tree theorem, Theorem 2.5, we can obtain the eight labeled spanning trees by considering any cofactor (here we choose the 4, 4 cofactor):

$$(3.1) \quad \det \mathcal{L}_{44}(G) = \begin{vmatrix} -(\gamma_a + \gamma_d + \gamma_e) & \gamma_a & \gamma_e \\ \gamma_a & -(\gamma_a + \gamma_b) & \gamma_b \\ \gamma_e & \gamma_b & -(\gamma_b + \gamma_c + \gamma_e) \end{vmatrix} \\ = \gamma_a \gamma_b \gamma_c + \gamma_a \gamma_b \gamma_d + \gamma_a \gamma_c \gamma_d + \gamma_a \gamma_c \gamma_e + \gamma_a \gamma_d \gamma_e + \gamma_b \gamma_c \gamma_d + \gamma_b \gamma_c \gamma_e + \gamma_b \gamma_d \gamma_e.$$

Let us orient this graph clockwise around the outside, i.e., $1 \rightarrow 4 \rightarrow 3 \rightarrow 2 \rightarrow 1$, and then choose the directed edge $1 \rightarrow 3$. With this orientation, this gives the incidence matrix

$$B_G = \begin{pmatrix} -1 & 0 & 0 & 1 & 1 \\ 1 & -1 & 0 & 0 & 0 \\ 0 & 1 & -1 & 0 & -1 \\ 0 & 0 & 1 & -1 & 0 \end{pmatrix},$$

and a basis for the cycle space is given by

$$y_1 = (1, 1, 0, 0, 1)^t, \\ y_2 = (0, 0, -1, -1, 1)^t.$$

For ease of notation, we write $\rho_i = 1/\gamma_i$ (we can think of ρ_i as resistances and γ_i as capacitances): The cycle intersection matrix \mathcal{Z}_G is given by

$$\mathcal{Z}_G = - \begin{pmatrix} \rho_a + \rho_b + \rho_e & \rho_e \\ \rho_e & \rho_c + \rho_d + \rho_e \end{pmatrix}.$$

The determinant of this matrix is

$$(3.2) \quad \det \mathcal{Z}_G = \rho_a \rho_c + \rho_a \rho_d + \rho_a \rho_e + \rho_b \rho_c + \rho_b \rho_d + \rho_b \rho_e + \rho_c \rho_e + \rho_d \rho_e.$$

It is not hard to see that if we multiply (3.2) by $\gamma_a \gamma_b \gamma_c \gamma_d \gamma_e$, then we obtain the expression in (3.1), verifying Theorem 2.12. The $\det(\mathcal{Z}_G)$ generates labeled tree complements: Each term in this polynomial represents a pair of edges that, when removed from the graph, yield a tree.

Moreover, we can use Theorem 2.9 to compute bounds on the edge weights to determine stability. Writing

$$w_1 = \rho_a + \rho_b, \quad w_2 = \rho_c + \rho_d, \quad w_{12} = \rho_e$$

gives

$$(3.3) \quad \mathcal{Z}_G = - \begin{pmatrix} w_1 + w_{12} & w_{12} \\ w_{12} & w_2 + w_{12} \end{pmatrix}.$$

3.1.1. One negative edge. Let us first consider the case where there is one negative edge. In this case, \mathcal{Z}_G has either one or zero positive eigenvalues. If there is one, then \mathcal{L}_G is stable, and if zero, it is unstable.

Note, however, that $\det \mathcal{Z}_G < 0$ if and only if there is one positive and one negative eigenvalue, which gives the condition

$$(w_1 + w_{12})(w_2 + w_{12}) - w_{12}^2 < 0$$

or

$$w_1 w_2 + w_1 w_{12} + w_2 w_{12} < 0.$$

If we assume that $\gamma_e < 0$, then, solving for w_{12} , we obtain the inequality

$$w_{12} < -\frac{w_1 w_2}{w_1 + w_2}$$

or

$$\rho_e < -\frac{\rho_a + \rho_b + \rho_c + \rho_d}{(\rho_a + \rho_b)(\rho_c + \rho_d)}.$$

Notice, for example, that if we assume that $\gamma_a \approx 0$, then the ρ_a terms dominate, and the condition we obtain for stability is $\gamma_e > -(\rho_c + \rho_d)$. Here γ_b is not relevant, and this makes sense: As $\gamma_a \rightarrow 0$, the a, b path is almost broken as a loop and thus cannot significantly affect stability, so the condition for stability will be a balance between e and c, d .

On the other hand, if one of the edges on the outside is negative (say without loss of generality γ_a), we obtain the inequality

$$w_1 < \frac{w_2 w_{12}}{w_2 + w_{12}}$$

or

$$(3.4) \quad \rho_a < -\frac{(\rho_c + \rho_d) \rho_e}{\rho_c + \rho_d + \rho_e} - \rho_b.$$

Note that every ρ on the right-hand side is positive. Note the particular role played by the ρ_b term: If we make γ_b larger, this makes ρ_b smaller (in absolute value), making it more difficult for ρ_a to cause the instability.

We note here that the inequality (3.4) is the same as would be obtained by Theorem III.3 of [29]. Stated concisely, that theorem says that as long as the weight on the single negative edge is less in absolute value than the reciprocal of the effective resistance between the two nodes (considering only the positive edges), then the Laplacian is stable. But it is not hard to see that the effective resistance between nodes 1, 2 is the negative of the right-hand side of (3.4), thus giving the same condition for stability.

3.1.2. Two negative edges. Now, assume the graph has two negative edges. Note that we cannot have both a, b or both c, d negative, since this automatically gives an unstable Laplacian, as follows from Theorem 2.10 of [1]: If edges a, b both have negative weights, then the negative edges disconnect the graph (more specifically, the subgraph containing only positive edges has more than one connected component), which implies that the graph Laplacian is unstable no matter how small the weights are in absolute value. Therefore, for stability, we need that at least two of w_1, w_2, w_{12} are negative.

Using Theorem 2.9, we have that \mathcal{L}_G is stable if and only if \mathcal{Z}_G has two positive eigenvalues or $-\mathcal{Z}_G$ has two negative eigenvalues. Looking at (3.3) (and noting the minus out front), this is the same as saying $-\mathcal{Z}_G$ has a positive determinant and a negative trace, so

$$w_1 w_2 + w_1 w_{12} + w_2 w_{12} > 0, \quad w_1 + w_2 + 2w_{12} < 0.$$

By symmetry, we need only consider the case where $w_1, w_2 < 0$ and where $w_{12} < 0$. In the former case, there is a negative edge on each of the “outsides” of the graph, and in the latter there is one in the middle.

3.2. One-cycle graphs. Let us now consider the general one-cycle graph and consider the condition for stability. Note that such a graph can have at most one negative edge and still be stable, and, moreover, it cannot be contained in the tree part of the graph, but must be part of the cycle. Therefore, we can simplify and consider R_N , the ring graph on N vertices: This graph has N vertices and N edges, and let us label them so that edge i goes from $i \rightarrow i+1$, modulo N .

The cycle can be represented by the vector $\mathbf{1}$, and thus we have that \mathcal{Z}_G is the 1×1 matrix (i.e., the scalar)

$$\mathcal{Z}_G = - \sum_{i=1}^N \frac{1}{\gamma_i}.$$

Thus to obtain a stable Laplacian, we need that $\mathcal{Z}_G > 0$ or

$$\sum_{i=1}^N \frac{1}{\gamma_i} < 0.$$

Assuming that $\gamma_1 < 0$ and the remainder are positive, the condition for stability is

$$\gamma_1 > \frac{-1}{\sum_{i=2}^n \frac{1}{\gamma_i}}.$$

If all of the positive edges have an equal weight, $\gamma_i \equiv \gamma$, then we obtain

$$\gamma_1 > -\gamma/(n-1).$$

In this case, we see that increasing γ helps with stability, whereas increasing the number of edges, all things being equal, hurts stability.

3.3. Two-cycle graphs. We now can consider the general graph with cycle number $C = 2$; in fact, we will reuse many of the computations of section 3.1.

As in the previous section, we simplify by assuming that there are no edges in $\text{Tree}(G)$, i.e., each edge is in one or more cycles. Let us split the edges into three sets S_1, S_2, S_{12} ; the edges in S_1 are those only in cycle one, those in S_2 are those only in cycle two, and finally S_{12} are those edges both in cycles one and two. It is clear that we can choose S_1, S_2 to be nonempty, but for some graphs, it may then be the case that S_{12} is empty (e.g., a graph with a figure-eight structure). Let us reorder the edges so that when ordered, the sets S_1, S_2, S_{12} come in order, and we write $k_* = |S_*|$. From this we see that we can choose the cycle basis as

$$y_1 = (1, 1, \dots, 1, 0, \dots, 0, -1, \dots, -1), \quad y_2 = (0, 0, \dots, 0, 1, \dots, 1, 1, \dots, 1),$$

where the three groups of edges are of lengths k_1, k_2, k_{12} . If we then define

$$w_* = \sum_{e \in S_*} \frac{1}{\gamma_e},$$

then we see that the cycle form \mathcal{Z}_G has the same structure as in (3.3), and the subsequent analysis holds over for this case.

The simplest case is the one where $S_{12} = \emptyset$ and $w_{12} = 0$. In this case \mathcal{Z}_G is diagonal, and the number of positive eigenvalues is exactly the number of negative diagonal entries. For example, let us say that the edge e for which $\gamma_e < 0$ is in cycle 1, and thus we need $w_1 < 0$ to have a stable Laplacian. This gives

$$\frac{1}{\gamma_e} + \sum_{f \neq e, f \in S_1} \frac{1}{\gamma_f} < 0$$

or

$$\gamma_e > \frac{-1}{\sum_{f \neq e, f \in S_1} 1/\gamma_f}.$$

This is exactly the condition as in the one-cycle case. Similarly, if the graph has two negative edges, we need both $w_1, w_2 < 0$ and the conditions on each cycle are decoupled.

In the case where $w_{12} > 0$, this gives an analysis similar to that given above. Let us consider the case where there is one negative edge, and it is in both cycles one and two. We again have the condition

$$w_{12} < -\frac{w_1 w_2}{w_1 + w_2}.$$

If, for example, we consider the case where all of the positive edges have the same weight γ , and the one negative edge has weight $\gamma_e < 0$, then we have

$$w_1 = k_1/\gamma, \quad w_2 = k_2/\gamma, \quad w_{12} = \frac{k_{12} - 1}{\gamma} + \frac{1}{\gamma_e}.$$

This simplifies to

$$(3.5) \quad \frac{1}{\gamma_e} < -\frac{1}{\gamma} \left(\frac{k_1 k_2}{k_1 + k_2} + (k_{12} - 1) \right).$$

Again, increasing γ helps for stability since it makes the condition easier to achieve, but increasing any of k_1, k_2, k_{12} makes the condition harder to achieve.

The interpretation of these facts is as follows: Clearly, increasing the strength of the positive couplings will make it harder for the negative coupling to destabilize the entire system, and in fact from (3.5) we see that the two weights are effectively linearly related. Moreover, we see that increasing the number of links in either of the cycles makes it easier to destabilize the system, similarly to the case of a single cycle.

4. Nonlinear example—Kuramoto model on a ring. The examples in the previous section studied the stability of linear Laplacians, but we can also use the current results to obtain novel results for nonlinearly coupled oscillators as well—in particular, we consider the Kuramoto model.

It has been known for a long time that if we have a fixed point for the Kuramoto model where all of the links are “short,” i.e., the angle difference across a link is less than $\pi/2$ in absolute value, then this gives a stable fixed point; see [28]. We could then pose the question of whether there exist solutions with “long” links that are still stable (this would correspond to having negative off-diagonal terms in the Jacobian).

Intuitively, it makes sense that if we make the weight δ_{ij} on an edge arbitrarily small, then this is possible, but we could then pose this question: Can we have a stable long-link solution if the edge weights are all equal? We will show that this answer is affirmative on the ring by a judicious choice of frequencies. In particular, we will answer the following question for the ring topology: Given a specific graph topology for the Kuramoto system with equal edge weights, what is the largest possible link for a stable fixed point of the system?

We consider the Kuramoto model as in (1.2), where we assume that the graph is a ring with n vertices and that the edge weights are all equal, and by rescaling we can assume that they are all one, so that $\delta_{i,i+1} = 1$ (here and below this sum on the indices is understood modulo n). Let θ^* be a fixed point of (1.2). From this we have

$$(4.1) \quad 0 = \omega_i + \sin(\theta_{i+1}^* - \theta_i^*) + \sin(\theta_{i-1}^* - \theta_i^*),$$

and the Jacobian at this point is given by J , where

$$(4.2) \quad J_{i,i\pm 1} = \cos(\theta_i^* - \theta_{i\pm 1}^*),$$

and $J_{i,i} = -J_{i,i+1} - J_{i,i-1}$. Note that the matrix has zero row sum and is thus a signed Laplacian. Then we see from the previous section that it is a stable fixed point if one of the following two conditions is met:

1. For all $i = 1, \dots, n$, $\gamma_{i,i+1} := \cos(\theta_i^* - \theta_{i+1}^*) > 0$;
2. one of the $\gamma_{i,i+1}$ is negative, and

$$(4.3) \quad \sum_{i=0}^{n-1} \frac{1}{\gamma_{i,i+1}} = \sum_{i=0}^{n-1} \frac{1}{\cos(\theta_i^* - \theta_{i+1}^*)} < 0.$$

Note that $\cos(\theta) > 0$ if and only if $|\theta| < \pi/2$, and thus the sign of the cosine of the angle difference tells us whether a link is long or short. These conditions can be reinterpreted as saying the following: The fixed point is stable if all of the links are short, or if there is one long link, then we need the sum of the reciprocals of the edge weights to be negative. The question then remains: Is there a configuration θ^* that is a fixed point of (1.2), has one long link, and satisfies (4.3)?

Let us write $\zeta_i = \theta_{i+1} - \theta_i$. Note that every term in the right-hand side of (4.1) will be a ζ_i , except for the term $\theta_{n-1} - \theta_0$, which we can write as $\sum_{i=0}^{n-2} \zeta_i$. Then (4.3)

becomes

$$(4.4) \quad \sum_{i=0}^{n-1} \frac{1}{\cos(\zeta_i)} + \frac{1}{\cos\left(\sum_{i=0}^{n-1} \zeta_i\right)} < 0.$$

Without loss of generality, let us assume that the longest link is given by the gap $\theta_{n-1} - \theta_0$, so that if there is one long link, it is represented by the last term in (4.4).

We can then pose this problem in an alternate form as follows: For any $\eta \in (0, 2\pi)$, let us find the maximum of

$$f(\zeta) = \sum_{i=0}^{n-2} \frac{1}{\cos(\zeta_i)} \quad \text{subject to } g(\zeta) = \sum_{i=0}^{n-2} \zeta_i = \eta.$$

Then it is possible to realize a fixed point with largest gap η if and only if

$$\min_{\zeta} f(\zeta) + \frac{1}{\cos \eta} < 0.$$

The nice thing we see from the constrained optimization problem is that since the constraint is symmetric up to any permutation of the variables, any internal optimizer must be as well. (We could also use Lagrange multipliers to obtain the same result.) Thus the maximum of f is achieved only when all of the ζ_i are equal, and (4.4) becomes

$$(4.5) \quad \frac{n-1}{\cos(\zeta)} + \frac{1}{\cos((n-1)\zeta)} < 0.$$

Summarizing our conditions, if we assume that $\cos(\zeta)$ is positive, then if and only if the conditions

$$\cos((n-1)\zeta) > 0 \quad \text{or} \quad \left(\cos((n-1)\zeta) < 0 \quad \text{and} \quad \frac{n-1}{\cos(\zeta)} + \frac{1}{\cos((n-1)\zeta)} < 0 \right)$$

hold, then we have a stable fixed point for the solution $\theta_k^* = k\zeta$. Note that both of these conditions can be simplified to require that their product be positive, i.e.,

$$(4.6) \quad h_n(\zeta) := (n-1) \frac{\cos((n-1)\zeta)}{\cos(\zeta)} + 1 > 0.$$

To exhibit a “long link” solution, we need to show that there is ζ^* with roots of $h_n(\zeta^*) = 0$ and $|n\zeta^* \pmod{2\pi}| \in (\pi/2, 3\pi/2)$. We will show numerically that this is possible for every n ; see Table 2.

Here we have found the first root of $h_n(\zeta)$ and compute $n\zeta_n$: This is the distance between θ_0 and θ_{n-1} if we choose the solution $\theta_k = k\zeta^*$. Note that in each case, the length of the link is greater than a quarter turn (note that we are presenting the length divided by 2π , so that a link is “long” if and only if it is greater than $1/4$), but this distance decreases monotonically to $1/4$ as $n \rightarrow \infty$.

Note that in general, $h_n(\zeta)$ has multiple roots, so that the first root may not always give the largest link, but we have found empirically that this is so. As such, the numbers in Table 2 represent the longest link possible for a stable fixed point of the Kuramoto system on a ring of size n .

TABLE 2
Length of longest possible link for a stable solution on the ring with n vertices.

n	$n\zeta^*/2\pi$
3	0.447
4	0.392
5	0.358
10	0.297
20	0.272
30	0.264
40	0.261
50	0.258

We can also compute the values of ω_k that give this solution. For all $k = 1, \dots, n-2$, we have $\theta_{k+1} - \theta_k = \theta_{k-1} - \theta_k = \zeta$, so that $\omega_k = 0$. However, we also have

$$\omega_{n-1}(\zeta) := -\sin(\theta_0 - \theta_{n-1}) - \sin(\theta_{n-2} - \theta_{n-1}) = \sin((n-1)\zeta) + \sin(\zeta),$$

and $\omega_0 = -\omega_{n-1}$.

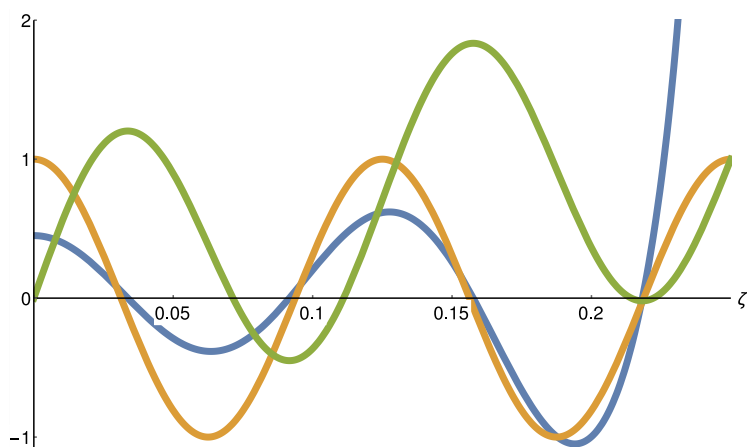


FIG. 1. We plot several of the functions referenced in the text for $n = 9$ vertices in the ring: $0.05 * h_9(\zeta)$ (blue), $\omega_8(\zeta)$ (orange), $\cos(8\zeta)$ (purple). (Figure appears in color online only.)

We plot several of the curves for the case of $n = 9$ in Figure 1. We have scaled $h_9(\zeta)$ so that all the curves are of the same order of magnitude; note that only its sign matters. As shown above, where the blue curve is positive, the solution is stable. The purple curve is the cosine of the length of the $(n-1, 0)$ link: since we see that this becomes negative before $h_9(\zeta)$ does, there is a stable long link configuration. Finally, note that the zeros of $h_9(\zeta)$ correspond to the critical points of $\omega_8(\zeta)$ (this can be checked), and thus the $\omega_8(\zeta)$ curve can be interpreted as a series of stable and unstable branches: Whenever $\omega_8(\zeta)$ is increasing in ζ , this corresponds to a stable solution, and when it is decreasing, this corresponds to a 1-saddle. As we slide up and down in this graph, this leads to saddle-node bifurcations in the obvious manner. Moreover, note that the $\omega = 0$ slice is the one considered in [26]—it was shown there that for $n = 9$ there are two nontrivial sinks and two corresponding saddles, and this is recovered here. By decreasing or increasing ω_7 , we can cause these saddles to collide with either of two sinks.

5. Proofs of theorems.

5.1. Proof of Theorem 2.9.

5.1.1. Definition of covering tree.

DEFINITION 5.1. Let $G = (V, E, \Gamma)$ be a connected graph. We define the universal cover of G , denoted by $U(G)$. If G is a tree, then $U(G) = G$. If G is not a tree, then $U(G)$ is a countably infinite tree defined as follows:

- Choose a root vertex $r \in V(G)$.
- The vertices of $U(G)$ are nonbacktracking rooted paths in G , i.e., sequences of the form $w = (r, v_2, v_3, \dots, v_n)$, where $(v_i, v_{i+1}) \in E$, and $v_{i-1} \neq v_{i+1}$ for all i .
- We say that two vertices $w_1, w_2 \in V(U(G))$ are adjacent if one is an extension of the other, i.e.,

$$w_1 = (r, v_2, v_3, \dots, v_n), \quad w_2 = (r, v_2, v_3, \dots, v_n, v_{n+1}),$$

or vice versa.

- Finally, we define the weight of the edge (w_1, w_2) in the example above to be the same as the weight on the edge (v_{n-1}, v_n) .

DEFINITION 5.2 (fundamental domain). For any graph $G = (V, E, \Gamma)$, we define T , a fundamental domain of the cover of G , as follows. First, construct the universal cover $U(G)$. Choose T to be any finite connected subtree of $U(G)$ such that each edge of G appears once. Notice that by definition, G and T have the same edge sets, so we will write E for either set. Since there are $|E|$ edges in T and it is a tree, then $|V(T)| = |E| + 1$. We will also use the notation $N_G = |V(G)|$ and $N_T = |V(T)|$ throughout.

Remark 5.3. The universal cover has certain properties:

1. If G is not a tree, then $U(G)$ has countably many vertices. The degree of the node $w = (r, v_2, \dots, v_n)$ in $U(G)$ is the same as the degree of v_n in G . In particular, $U(G)$ is locally finite.
2. The map

$$\begin{aligned} \varphi: U(G) &\rightarrow G, \\ (r, v_2, \dots, v_n) &\mapsto v_n \end{aligned}$$

is a covering map, i.e., it is a map of graphs, it is surjective, and it is an isomorphism in the neighborhood of any vertex.

3. The construction of T is independent (up to isomorphism) of the choice of the root vertex r .
4. Finally, it should be noted that we can construct T directly without using the universal cover; see Example 5.5 for an example.

PROPOSITION 5.4. Let $G = (V, E, \Gamma)$ be a connected weighted graph, and define T as above. Then the graph map $\varphi: T \rightarrow G$ given by

$$\varphi((r, v_2, \dots, v_n)) = v_n$$

is a surjective map. Note that $|V(T)| - |V(G)| = C$, i.e., the difference in the number of vertices of a fundamental domain of the covering tree and the original graph itself is the cycle rank of the graph.

Proof. Assume that $w_1, w_2 \in V(T)$ and $(w_1, w_2) \in E(T)$. Since w_1, w_2 are adjacent in $U(G)$, we have that

$$w_1 = (r, v_1, \dots, v_k), \quad w_2 = (r, v_1, \dots, v_k, v_{k+1}),$$

and $(v_k, v_{k+1}) \in E(G)$. Then $\varphi(w_1) = v_k$ and $\varphi(w_2) = v_{k+1}$ are adjacent in G whenever w_1, w_2 are adjacent in T . Therefore, φ is a graph homomorphism. To see that it is surjective, note that every edge of G appears in T , and therefore, its incident vertices must as well. Finally, note that

$$|V(T)| - |V(G)| = |E| + 1 - |V(G)| = C.$$

This completes the proof. \square

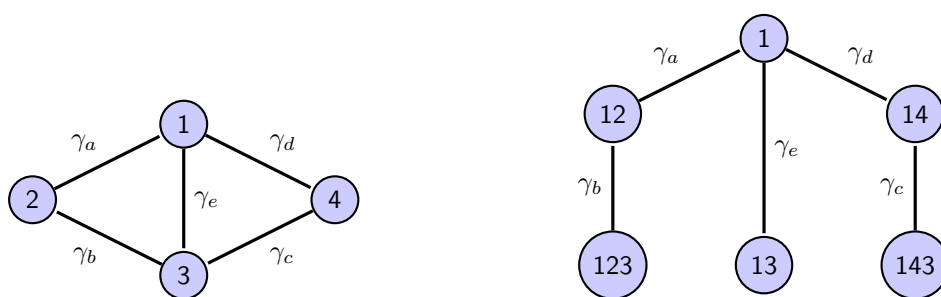


FIG. 2. The diamond graph G , and a choice of fundamental domain of its universal cover T . Note that G and T have the same edge sets with five edges, and T necessarily has six vertices.

Example 5.5. Let us consider the diamond graph in Figure 2. If we choose the root vertex as 1, then 2, 3, 4 are neighbors and so we add 12, 13, 14 to T . Since γ_b, γ_c have not yet appeared, we add them as in the tree. Of course, this is not the only covering tree we can obtain, even assuming that we choose 1 as the root; for example, if we remove vertex 143 and add a vertex 1234, this would also be a covering tree.

5.1.2. The projections.

DEFINITION 5.6. Let G be a connected graph, and let $T, \varphi: T \rightarrow G$ be as defined above. We define X to be the $N_T \times N_G$ matrix whose elements are

$$X_{a,v} = \begin{cases} 1, & \varphi(a) = v, \\ 0, & \varphi(a) \neq v. \end{cases}$$

We will also refer to the columns by the vectors $x^v, v \in V(G)$.

LEMMA 5.7. Let $G = (V, E, \Gamma)$ be a connected weighted graph, and T, X be as in Definitions 5.2 and 5.6. Then

$$(5.1) \quad \mathcal{L}_G = X^\top \mathcal{L}_T X$$

or

$$(5.2) \quad (\mathcal{L}_G)_{v,w} = \langle x^v, \mathcal{L}_T x^w \rangle.$$

Proof. Let us use formula (5.2). Writing this out, we have

$$\langle x^v, \mathcal{L}_T x^w \rangle = \sum_{i,j} x_i^v (\mathcal{L}_T)_{ij} x_j^w = \sum_{i,j} (\mathcal{L}_T)_{ij} 1_{\varphi(i)=v} 1_{\varphi(j)=w}.$$

If we can show that there exists a unique pair i, j such that $\varphi(i) = v$, $\varphi(j) = w$, and $(\mathcal{L}_T)_{ij} = \gamma_{(v,w)}$, then we are done. But notice that since $\varphi(i) = v$, and $\varphi(j) = w$, if $(i, j) \in E(T)$, then the weight on the edge (i, j) must be $\gamma_{(v,w)}$, since φ is a graph homomorphism. Moreover, since each edge of G can appear at most once in T , this means that there is only one such pair. \square

DEFINITION 5.8. *The map $\varphi: T \rightarrow G$ is surjective but in general not injective. For each $v \in V(G)$ with $|\varphi^{-1}(v)| > 1$, choose one representative of $\varphi^{-1}(v)$ as the primary representative of this set. Now, for each $w \in \varphi^{-1}(v)$ with $w \neq \text{prim}(v)$, define the vector $q^w \in \mathbb{R}^{N_T}$ by*

$$q_i^w = \begin{cases} -1, & i = \text{prim}(v), \\ 1, & i = w, \\ 0 & \text{else.} \end{cases}$$

We define Q as the matrix whose columns are the q^w .

LEMMA 5.9. *Let $S \subseteq \mathbb{R}^{N_T}$ be the span of the x^v defined in Definition 5.6. Then the vectors q^w form a basis for S^\perp . In particular, $\dim(\text{span}(q^w)) = \text{rank}(Q) = c$, the cycle rank of G .*

Proof. First, we see that each q^w is orthogonal to S . Consider x^v where $v = \varphi(w)$. This vector is equal to 1 in both the $\text{prim}(v)$ and the w index, and therefore, $\langle q^w, x^v \rangle = 0$. Moreover, if we consider x^v with $v \neq \varphi(w)$, then this vector is equal to zero on these indices, and again $\langle q^w, x^v \rangle = 0$. Since $\dim(\text{span}(S)) = \text{rank}(X) = |V(G)|$, then $\text{rank}(Q) = |E| + 1 - \dim S = |E| - |V(G)| + 1 = c$. Finally, note that the q^w are clearly independent by construction, so by dimension counting they form a basis for S^\perp . \square

LEMMA 5.10. *The Laplacian $\mathcal{L}_T: \mathbb{R}^{N_T} \rightarrow \mathbb{R}^{N_T}$ is not invertible. Recall that by definition all of the edge weights are chosen to be nonzero and T is a tree; it follows from this that \mathcal{L}_T has a one-dimensional nullspace given by $\text{span}(\mathbf{1})$. In this case, the canonical map*

$$\tilde{\mathcal{L}}_T: \mathbb{R}^{N_T} / \{\mathbf{1}\} \rightarrow \mathbb{R}^{N_T} / \{\mathbf{1}\}$$

is invertible. (We can define $\tilde{\mathcal{L}}_T$ explicitly as follows: Let P be the orthogonal projection onto $\mathbf{1}^\perp$, and then $P\mathcal{L}_TP$ maps $\mathbb{R}^{N_T} / \{\mathbf{1}\}$ to itself; we denote this map by $\tilde{\mathcal{L}}_T$.) We then have that

$$Q^\top (\tilde{\mathcal{L}}_T)^{-1} Q = \mathcal{Z}_G.$$

Proof. In general, for any graph H , the incidence matrix B_H is a map from $\mathbb{R}^{|E(H)|}$ to $\mathbb{R}^{|V(H)|}$. In particular, since $|E(T)| = |E(G)|$ and $|V(T)| = |E(G)| + 1$, this means that

$$B_T: \mathbb{R}^{|E(T)|} \rightarrow \mathbb{R}^{|V(T)|}$$

is a map of co-rank one, and as such B_T^\top is a map with a one-dimensional nullspace,

which is $\text{span}(\mathbf{1})$. Therefore we can define \tilde{B} to be the canonical map so that

$$\tilde{B}: \mathbb{R}^{|E(T)|} \rightarrow \mathbb{R}^{|E(T)|+1}/\{\mathbf{1}\}$$

is invertible: To define this, we first define \tilde{B}^\top to be the induced map on the quotient and then $\tilde{B} = (\tilde{B}^\top)^\top$. In particular, note that $\tilde{B} = B$ on $\mathbf{1}^\perp$.

For each w such that $q^w \neq \mathbf{0}$, consider the unique path in T from w to $\text{prim}(\varphi(w))$. Let us write this path as

$$w = v_0 \rightarrow v_1 \rightarrow v_2 \rightarrow \cdots \rightarrow v_{n-1} \rightarrow v_n = \text{prim}(\varphi(w)).$$

Define a vector $r^w \in \mathbb{R}^{|E(T)|}$ as follows:

$$r_e^w = \begin{cases} 1, & e = (v_i, v_{i+1}), \exists i, \\ -1, & e = (v_{i+1}, v_i), \exists i, \\ 0 & \text{else.} \end{cases}$$

We first compute that

$$B_T r^w = q^w.$$

To see this, note that $B_T r^w = 0$ on any vertex not in the path, and, moreover, will be zero on any vertex on the interior of the path due to cancellation. We need only consider the endpoints.

First, assume that $w \rightarrow v_1$ is the chosen orientation for the edge. Then $(B_T r^w)_w = ((B_T)_{(w, v_1)})r_{(w, v_1)} = 1 \cdot 1 = 1$. If, on the other hand, $v_1 \rightarrow w$ is the orientation, then we have $(B_T r^w)_w = ((B_T)_{(w, v_1)})r_{(w, v_1)} = -1 \cdot -1 = 1$ as well. Similarly, $(B_T r^w)_{\text{prim}(\varphi(w))} = -1$. If we define

$$(5.3) \quad Y = \tilde{B}_T^{-1} Q,$$

then the columns of Y form a basis for the cycle space of G .

By Lemma 2.8, $\mathcal{L}_T = -B_T D B_T^\top$ (note that $D_G = D_T$ so that we can drop the subscript), and it is easy to see that similarly, $\tilde{\mathcal{L}}_T = -\tilde{B}_T D \tilde{B}_T^\top$. From this we have

$$Q^\top (\tilde{\mathcal{L}}_T)^{-1} Q = -Q^\top \tilde{B}_T^{-\top} D^{-1} \tilde{B}_T^{-1} Q = -Y^\top D^{-1} Y,$$

giving the definition (2.1). \square

Proof of Theorem 2.9. Recall the definition of $M := \tilde{\mathcal{L}}_T$. From Lemma 5.7 we have that

$$\tilde{\mathcal{L}}_G = X^\top \tilde{\mathcal{L}}_T X,$$

and Lemma A.3 gives that

$$n_+(\mathcal{L}_G) = n_+(M|_S),$$

since the columns of X form a basis for S . From Lemma 5.10 we have that

$$\mathcal{Z}_G = Q^\top (\tilde{\mathcal{L}}_T)^{-1} Q,$$

and Lemma A.3 again gives that

$$n_+(\mathcal{Z}_G) = n_+((M)^{-1}|_{S^\perp}),$$

since the columns of Q form a basis for S^\perp . Using Lemma A.1 gives

$$n_+(\mathcal{L}_G) = n_+(\mathcal{L}_T) - n_+(\mathcal{Z}_G).$$

Recall that T is a tree. Using Theorem 2.10 of [1], this implies that the number of positive eigenvalues of \mathcal{L}_T are given exactly by the number of negative edges in $E(T)$. Recalling that the edge sets of G and T are the same, we are done.

Finally, to prove the last sentence of the theorem, we again appeal to Theorem 2.10 of [1]. One consequence of this theorem is that for \mathcal{L}_G to be stable, G_+ must be connected. If the number of negative edges in G is the same as the number of cycles, this means that G_+ is a tree, and thus is a spanning tree of G . For each negative edge in G , consider the two incident vertices and the unique path in G_+ connecting them. This plus the addition of the negative edge make a cycle in G , and it is clear that each negative edge will be in a different cycle. Using this choice of cycles, we now have exactly one negative edge in each cycle. \square

Remark 5.11. Instead of defining $\tilde{\mathcal{L}}_T$ as an invertible operator, we could have equivalently chosen \mathcal{L}_T^\dagger as the Moore–Penrose pseudoinverse of \mathcal{L}_T , and the formula $Q^\top \mathcal{L}_T^\dagger Q = \mathcal{Z}_G$ would still hold. To see this, write $\mathcal{L}_T^\dagger = -(B_T D_T B_T^\top)^\dagger = -B_T^\top D_T^{-1} B_T^\dagger$, and note that the computation in the proof of Lemma 5.10 shows that $Y = B_T^\dagger Q$, since $B_T Y = B_T B_T^\dagger Q = \text{Proj}_{1^\perp} Q = Q$. From this we obtain

$$Q^\top \mathcal{L}_T^\dagger Q = -Q^\top B_T^\top D_T^{-1} B_T^\dagger Q = -Y^\top D_T^{-1} Y = \mathcal{Z}_G$$

as before. However, the approach we use above meshes better with Lemma A.1.

Example 5.12. There are many matrices defined along the path to proving Theorem 2.9, so we work out all of them for the graph considered in section 3.1 and shown in Figure 2.

We choose the convention that the vertices of G are ordered as 1, 2, 3, 4 and the vertices of T are ordered 1, 12, 14, 123, 13, 143. There is only one vertex in $V(G)$ that has multiple preimages, and we decide that $13 = \text{prim}(3)$. Finally, we take care to choose the same orientation on T that corresponds to the orientation we took on the graph G in section 3.1. Then we have

$$X = \begin{pmatrix} 1 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 \\ 0 & 0 & 0 & 1 \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 1 & 0 \end{pmatrix}, \quad Q = \begin{pmatrix} 0 & 0 \\ 0 & 0 \\ 0 & 0 \\ 1 & 0 \\ -1 & -1 \\ 0 & 1 \end{pmatrix}, \quad B_T = \begin{pmatrix} -1 & 0 & 0 & 1 & 1 \\ 1 & -1 & 0 & 0 & 0 \\ 0 & 0 & 1 & -1 & 0 \\ 0 & 1 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & -1 \\ 0 & 0 & -1 & 0 & 0 \end{pmatrix}.$$

These are the patterns we can see in the matrix X : Each row has a single 1, and the sum of any column is the size of its preimage under φ .

We first compute

$$B_T^\dagger = \frac{1}{6} \begin{pmatrix} -2 & 4 & -2 & 4 & -2 & -2 \\ -1 & -1 & -1 & 5 & -1 & -1 \\ 1 & 1 & 1 & 1 & 1 & -5 \\ 2 & 2 & -4 & 2 & 2 & -4 \\ 1 & 1 & 1 & 1 & -5 & 1 \end{pmatrix}, \quad Y = B_T^\dagger Q = \begin{pmatrix} 1 & 0 \\ 1 & 0 \\ 0 & -1 \\ 0 & -1 \\ 1 & 1 \end{pmatrix},$$

the columns of which are the cycles of the graph G .

5.2. Proof of Theorem 2.12. We again start with several lemmas.

DEFINITION 5.13. Recall $G, T, \varphi: T \rightarrow G$. For each $v \in V(G)$, if $|\varphi^{-1}(v)| = 1$, let $\psi(v) = \varphi^{-1}(v)$; else, $\psi(v) = \text{prim}(v)$ as defined in Definition 5.8. Then order the vertices of $V(T)$ so that we list all of the vertices in the range of ψ first and then all of the other vertices in some order. It is easy to see the range of ψ as N_G elements.

We then define the first N_G columns of U in the same way as we did in the definition of X , i.e., for $v \in V(G)$, define the v th column of U , u^v by

$$(u^v)_a = \begin{cases} 1, & \varphi(a) = v, \\ 0 & \text{else,} \end{cases}$$

and for columns $i = N_G + 1, \dots, N_T$, define u^i to be the standard basis function with a 1 in the i th slot. By construction, U is a lower triangular matrix with 1s on the diagonal, so $\det(U) = 1$.

LEMMA 5.14. The $N_G \times N_G$ upper-left block of $U^\top \mathcal{L}_T U$ is \mathcal{L}_G .

Proof. The proof is more or less the same as Lemma 5.7. Notice that once we have chosen this ordering, the first N_G columns of U is the matrix X as defined in Definition 5.6 with the corresponding ordering of vertices. \square

LEMMA 5.15. Let H be a symmetric matrix with a one-dimensional nullspace, and let x span this nullspace. Then for any y, z , we have

$$\det(H + y \otimes z) = \frac{\langle x, y \rangle \langle x, z \rangle}{\langle x, x \rangle} \det_{\text{red}} H.$$

Proof. Since H is self-adjoint, it has real eigenvalues and an orthonormal eigenbasis. Let us write $\lambda_1 = 0$ with $\lambda_2, \dots, \lambda_n$ not zero, and let w_1, \dots, w_n be the corresponding eigenvectors. Let W be the matrix whose columns are the w_i ; since the eigenvectors form an orthonormal basis, $W^{-1} = W^\top$.

Let M_1 be the diagonal matrix whose diagonal entries are the λ_i , and define $M_2 = (W^\top y) \otimes (W^\top z)$, i.e.,

$$(M_2)_{ij} = \langle w_i, y \rangle \langle w_j, z \rangle.$$

Notice that $H = WM_1W^\top$, since this is just its diagonalization, and

$$WM_2W^\top = W((W^\top y) \otimes (W^\top z))W^\top = WW^\top(y \otimes z)WW^\top = y \otimes z,$$

where we have used the identity $(Ax_1) \otimes (Bx_2) = A(x_1 \otimes x_2)B^\top$. Thus we have

$$H + y \otimes z = W(M_1 + M_2)W^\top,$$

and since W is unitary this gives

$$\det(H + y \otimes z) = \det(M_1 + M_2).$$

To compute the last term, notice that we can use the multilinearity of the determinant by rows and break this up into 2^n terms. However, note two things: the matrix M_2 is rank one and thus has only one linearly independent row, so if we choose more than one row from M_2 , the determinant is zero. Also, the first row of M_1 is zero, so to obtain a nonzero determinant, we must choose the first row of M_2 . Therefore, the only term which gives a nonzero determinant is the one where we choose the first row from M_2 and all of the other rows from M_1 . Expanding by minors in the first column, the determinant of this matrix is

$$\langle w_1, y \rangle \langle w_1, z \rangle \lambda_2 \times \cdots \times \lambda_n.$$

Since we chose $\lambda_1 = 0$, this means that w_1 is the unit vector that spans $\ker(H)$, and also notice that $\det_{\text{red}} H$ is by definition $\lambda_2 \times \cdots \times \lambda_n$. Thus we have

$$\det(H + y \otimes z) = \langle w_1, y \rangle \langle w_1, z \rangle \det_{\text{red}} H.$$

If x is a general vector in $\ker(H)$, then x is a scalar multiple of w_1 and we have

$$\det(H + y \otimes z) = \frac{\langle x, y \rangle \langle x, z \rangle}{\langle x, x \rangle} \det_{\text{red}} H.$$

This completes the proof. \square

Proof of Theorem 2.12. Recall the definition of U in Definition 5.13. Define

$$A = U^\top (\mathcal{L}_T + \mathbf{1} \otimes \mathbf{1}) U = U^\top \mathcal{L}_T U + U^\top \mathbf{1} \otimes U^\top \mathbf{1}.$$

If we can establish the following three identities:

$$(5.4) \quad \det(A) = N_T^2 \prod_{e \in E(G)} \gamma_e,$$

$$(5.5) \quad \det(A|_S) = \frac{N_T^2}{N_G} \det_{\text{red}} \mathcal{L}_G,$$

$$(5.6) \quad \det((A^{-1})|_{S^\perp}) = \det(\mathcal{Z}_G),$$

then, using Lemma A.2, we obtain

$$N_T^2 \prod_{e \in E(G)} \gamma_e = \frac{\frac{N_T^2}{N_G} \det_{\text{red}} \mathcal{L}_G}{\det(\mathcal{Z}_G)},$$

and this proves the theorem. So it remains to show these three identities.

Identity (5.6) follows directly from Lemma 5.10. To establish (5.4), we note that

$$\det(A) = \det(U^\top (\mathcal{L}_T + \mathbf{1} \otimes \mathbf{1}) U) = \det(\mathcal{L}_T + \mathbf{1} \otimes \mathbf{1}),$$

since $\det(U) = 1$. By Lemma 5.15, we have

$$\det(\mathcal{L}_T + \mathbf{1} \otimes \mathbf{1}) = \frac{\langle \mathbf{1}, \mathbf{1} \rangle \langle \mathbf{1}, \mathbf{1} \rangle}{\langle \mathbf{1}, \mathbf{1} \rangle} \det_{\text{red}} \mathcal{L}_T = N_T \det_{\text{red}} \mathcal{L}_T,$$

and since there is only one spanning tree of T , by the matrix tree theorem,

$$\det_{\text{red}} \mathcal{L}_T = N_T \prod_{e \in E(G)} \gamma_e.$$

To establish (5.5), note that if we use the basis for \mathbb{R}^{N_T} made from the columns of U , then $A|_S$ is just the upper left $N_G \times N_G$ block of A , which is itself the upper-left block of $U^\top \mathcal{L}_T U + U^\top \mathbf{1} \otimes U^\top \mathbf{1}$. By Lemma 5.14, the upper-left block of the first term is just \mathcal{L}_G , and the upper-left block of the second term can be written as

$$\tilde{U}^\top \mathbf{1} \otimes \tilde{U}^\top \mathbf{1},$$

where \tilde{U} is just the first N_G columns of U . In particular, the v th row of $\tilde{U}^\top \mathbf{1}$ is a zero-one vector with as many ones as vertices in $\varphi^{-1}(v)$, and therefore, $u := \tilde{U}^\top \mathbf{1}$ is a column vector of height N_G whose v th entry is $|\varphi^{-1}(v)|$.

Putting this together, and again using Lemma 5.15, gives

$$\begin{aligned} \det(A|_S) &= \det(\mathcal{L}_G + u \otimes u) = \frac{\langle u, \mathbf{1} \rangle \langle u, \mathbf{1} \rangle}{\langle \mathbf{1}, \mathbf{1} \rangle} \det_{\text{red}} \mathcal{L}_G \\ &= \frac{\left(\sum_{v \in V(G)} \varphi^{-1}(v) \right)^2}{N_G} \det_{\text{red}} \mathcal{L}_G = \frac{N_T^2}{N_G} \det_{\text{red}} \mathcal{L}_G. \end{aligned}$$

This completes the proof. \square

Appendix A. Classical lemmas. This follows from the following duality formula for the number of negative eigenvalues (or more generally the inertia) of a matrix. This result seems to have been frequently rediscovered but the first instance we have found in the literature is due to Haynsworth [38]. This result has been used frequently in the nonlinear waves literature, in the context of the stability of traveling wave solutions.

LEMMA A.1 (Haynsworth). *Suppose that M is a nonsingular $N \times N$ Hermitian matrix. Let S be a subspace of \mathbb{R}^N , let S^\perp be the orthogonal subspace, let P_S be the orthogonal projection onto S , and let $M|_S = P_S M P_S$ be the restriction of M to S . If $M|_S$ is nonsingular, then*

$$n_+(M) = n_+(M|_S) + n_+((M^{-1})|_{S^\perp}).$$

Note that, in her original paper, Haynsworth states this theorem slightly differently in terms of the Schur complement of the matrix, but it is obviously equivalent.

LEMMA A.2. *Suppose that M is a nonsingular $N \times N$ Hermitian matrix. Let S be a subspace of \mathbb{R}^N , let S^\perp be the orthogonal subspace, let P_S be the orthogonal projection onto S , and let $M|_S = P_S M P_S$ be the restriction of M to S . If $M|_S$ is nonsingular, then*

$$\det(M) = \frac{\det(M|_S)}{\det(M^{-1}|_{S^\perp})}.$$

This is more or less the determinantal version of Lemma A.1 and could be proved in a similar fashion. It can also be found explicitly in [39].

LEMMA A.3 (Sylvester's law of inertia). *Suppose M is Hermitian and U is nonsingular. Then*

$$(A.1) \quad n_-(M) = n_-(U^\top MU).$$

In particular, if A is an $n \times n$ matrix, $S \subset \mathbb{R}^n$ is a k -dimensional subspace, P_S is the orthogonal projection onto S , v_1, \dots, v_k is a basis for S , and $V = (v_1, \dots, v_k)$ is the $k \times k$ matrix whose i th column is v_i , then

$$n_+(P_S A P_S) = n_+(V^\top A V).$$

Proof. The first statement is just the standard formulation of Sylvester's theorem [40], so we will prove the second part. Notice that if we choose w^1, \dots, w^k to be an orthonormal basis for S , and let W be the $k \times k$ matrix whose columns are the w^k , then $P_S A P_S = W^\top A W$. Choosing $Z = W^{-1}V$, we have

$$Z^\top (W^\top A W) Z = V^\top A V,$$

and by the first statement this implies that these two matrices have the same signature. \square

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