One Can Hear the Composition of a String: Experiments with an Inverse Eigenvalue Problem^{*}

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Abstract. To what extent do the vibrations of a mechanical system reveal its composition? Despite innumerable applications and mathematical elegance, this question often slips through those cracks that separate courses in mechanics, differential equations, and linear algebra. We address this omission by detailing a classical finite dimensional example: the use of frequencies of vibration to recover positions and masses of beads vibrating on a string. First we derive the equations of motion, then compare the eigenvalues of the resulting linearized model against vibration data measured from our laboratory's monochord. More challenging is the recovery of masses and positions of the beads from spectral data, a problem for which a variety of elegant algorithms exist. After presenting one such method based on orthogonal polynomials in a manner suitable for advanced undergraduates, we confirm its efficacy through physical experiment. We encourage readers to conduct their own explorations using the numerous data sets we provide.

Key words. beaded string, inverse eigenvalue problem, vibration, continued fractions

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1. Introduction. The 18th century witnessed revolutionary progress in the mathematical description of fundamental problems in mechanics, thanks to the collective efforts of natural philosophers such as Leonhard Euler, the Bernoulli family, d'Alembert, Lagrange, and others [5, 27]. These old masters developed predictive models: given the material properties of a system, along with its position and velocity at some initial time, determine the system's state at all future times. While such *forward models* give great insight, modern applications often present the problem backwards: We can measure how a system responds to some stimulus, and from such experiments seek to discover the system's composition. In many instances the response is an acoustic signature and so one is led to pose the backward, or *inverse*, problem in the form of the question, "Can one hear?" For example, Gopinath and Sondhi [15] ask, "Can one hear the shape of your throat?"; Kac [17] asks, "Can one hear the shape of a drum?"; Sekii and Shibahashi [22] ask, "Can one hear into the sun?"; Lin [20] asks, "Can one hear a crack in a beam?"; and Gutkin and Smilansky [16] ask, "Can one hear the

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shape of a graph?" Each of these investigations seeks to echo the success of Borg [2], Levinson [19], and Gelfand and Levitan [11] in their various proofs that one can hear a "potential," and Krein's demonstration that one can hear the mass distribution of a nonuniform string [18].

Krein's argument, as developed by Dym and McKean [9], proceeds from the beaded string (a massless thread supporting a finite number of point masses) to the general nonhomogeneous distribution of mass. With Gantmacher [10], Krein returned to the beaded case, resurrected the lovely work of Stieltjes [23] on continued fractions, and carefully developed the requisite matrix analysis and complex function theory. This finite-dimensional case has since been systematized, and numerous related algorithms now exist for its resolution [6]. Although these methods supply constructive means for determining masses and lengths from spectral data, this theory has, to our knowledge, remained untested. We here provide experimental confirmation on data taken from our own laboratory. Our larger aim, however, is to give an entry to inverse spectral theory, numerics and experiment that is accessible to students possessing a solid undergraduate background in linear algebra. Indeed, beaded string experiments form the culminating exercises in an optional one-credit physical laboratory that accompanies our junior-level Matrix Analysis course at Rice University.

The remainder of our tour is organized as follows. In section 2 we derive a system of differential equations for the displacement of a plucked beaded string and show how to express the solution in terms of eigenvalues and eigenvectors of underlying mass and stiffness matrices. In section 3 we introduce our experimental setup and explain how we measure the eigenvalues of a beaded string. In sections 4–8 we combine algorithms of de Boor and Golub [8] and Gladwell [12, section 4.4] to determine bead locations and masses from two sets of eigenvalues, and in section 9 we confirm, using real data, that one set of eigenvalues suffices to reveal the location and masses of symmetrically placed beads.

2. The Forward Problem. We thread a massless string through n beads, apply a known tension τ , and clamp its ends at a known distance, ℓ , apart. With reference to Figure 1 we denote the mass of the *j*th bead by m_j , and let ℓ_j denote the length between masses m_j and m_{j+1} (with ℓ_0 and ℓ_n denoting the length between the beads at each end and the clamped support). Following a small vertical pluck we presume that the *j*th bead suffers the planar displacement (x_j, y_j) and that the *j*th segment makes the angle ϕ_j with the horizontal. In this state the horizontal and vertical



Fig. I A string with n = 3 beads at rest (solid black) and in a deformed state (dotted).

ONE CAN HEAR THE COMPOSITION OF A STRING

components of the string restoring forces at the jth mass are

$$\tau(\cos\phi_j - \cos\phi_{j-1})$$
 and $\tau(\sin\phi_j - \sin\phi_{j-1})$,

respectively. The angles ϕ_{j-1} and ϕ_j can be determined from the horizontal and vertical displacements. As evident from Figure 1,

$$\cos \phi_j = \frac{\ell_j + (x_{j+1} - x_j)}{\sqrt{(\ell_j + (x_{j+1} - x_j))^2 + (y_{j+1} - y_j)^2}},\\ \sin \phi_j = \frac{y_{j+1} - y_j}{\sqrt{(\ell_j + (x_{j+1} - x_j))^2 + (y_{j+1} - y_j)^2}},$$

and similarly for ϕ_{j-1} . If the pluck is small and the string is taut, we may make the (customary) assumptions that $|y_{j+1} - y_j| \ll \ell_j$ and $|x_{j+1} - x_j| \ll \ell_j$ (for a careful alternative, see [1]), and so arrive at the approximations

$$\cos \phi_j \approx 1, \qquad \sin \phi_j \approx \frac{y_{j+1} - y_j}{\ell_j},$$
$$\cos \phi_{j-1} \approx 1, \qquad \sin \phi_{j-1} \approx \frac{y_j - y_{j-1}}{\ell_{j-1}}.$$

With these approximations the horizontal forces balance, so it remains only to balance the vertical restoring forces with their associated inertial terms ("ma = F"):

$$m_j y_j''(t) = \tau \left(\frac{y_{j+1}(t) - y_j(t)}{\ell_j} - \frac{y_j(t) - y_{j-1}(t)}{\ell_{j-1}} \right), \qquad j = 1, \dots, n.$$

For this equation to hold at the first and last masses (j = 1 and j = n), we define $y_0 = y_{n+1} = 0$, thus describing the fixed ends of the string. As a consequence of our first-order approximations to the sines and cosines, these *n* coupled equations are *linear*, and can be most conveniently organized into the matrix form

(2.1)
$$\mathbf{M}\mathbf{y}''(t) = -\mathbf{K}\mathbf{y}(t),$$

with state vector $\mathbf{y}(t)$, mass matrix \mathbf{M} ,

$$\mathbf{y}(t) = \begin{bmatrix} y_1(t) \\ y_2(t) \\ \vdots \\ y_n(t) \end{bmatrix}, \qquad \mathbf{M} = \begin{bmatrix} m_1 & & & \\ & m_2 & & \\ & & \ddots & \\ & & & m_n \end{bmatrix}$$

and stiffness matrix

(2.2)
$$\mathbf{K} = \tau \begin{bmatrix} \ell_0^{-1} + \ell_1^{-1} & -\ell_1^{-1} & & \\ -\ell_1^{-1} & \ell_1^{-1} + \ell_2^{-1} & \ddots & \\ & \ddots & \ddots & -\ell_{n-1}^{-1} \\ & & -\ell_{n-1}^{-1} & \ell_{n-1}^{-1} + \ell_n^{-1} \end{bmatrix},$$

with all unspecified entries equal to zero. The mass and stiffness matrices enjoy two lovely properties: they are symmetric, $\mathbf{M} = \mathbf{M}^T$ and $\mathbf{K} = \mathbf{K}^T$, and positive definite, meaning that

$$\mathbf{y}^T \mathbf{M} \mathbf{y} = \sum_{j=1}^n m_j y_j^2$$
 and $\mathbf{y}^T \mathbf{K} \mathbf{y} = \tau \left(\frac{y_1^2}{\ell_0} + \frac{y_n^2}{\ell_n} + \sum_{j=1}^{n-1} \frac{(y_j - y_{j+1})^2}{\ell_j} \right)$

are both positive for every nonzero vector $\mathbf{y} \in \mathbb{R}^{n,1}$

We are interested in the motion induced by an initial pluck of the string, whereby the masses are vertically displaced by the components of the vector \mathbf{y}_0 , then released. Thus we presume that

$$\mathbf{y}(0) = \mathbf{y}_0, \qquad \mathbf{y}'(0) = \mathbf{0}$$

Equation (2.1) is a typical second-order, constant-coefficient homogeneous system of equations, a problem routinely tackled with the help of some linear algebra.

2.1. Solving the Differential Equation. By analogy with the harmonic motion experienced by a single tethered mass we put forward the educated guess that the solution of (2.1) takes the form

$$\mathbf{y}(t) = e^{i\omega t} \,\mathbf{v},$$

where the scalar ω is the frequency and the constant vector **v** somehow accounts for the interplay between the masses. On substituting our guess into (2.1) we find that ω and **v** must obey the *generalized eigenproblem*

(2.3)
$$\mathbf{K}\mathbf{v} = \omega^2 \mathbf{M}\mathbf{v},$$

that is, ω^2 is an eigenvalue with associated eigenvector **v** for the pair (**K**, **M**). We now argue that this pair has *n* positive distinct eigenvalues and *n* linearly independent real eigenvectors.

We begin by noting that $\mathbf{K}\mathbf{v} = \lambda \mathbf{M}\mathbf{v}$ can be transformed into the standard eigenproblem, $\mathbf{A}\mathbf{u} = \lambda \mathbf{u}$, via the substitutions

$$\mathbf{u} = \mathbf{M}^{1/2} \mathbf{v}$$
 and $\mathbf{A} = \mathbf{M}^{-1/2} \mathbf{K} \mathbf{M}^{-1/2}$

where $\mathbf{M}^{1/2}$ is the elementwise square root of \mathbf{M} (since \mathbf{M} is a diagonal matrix). We next note that, like \mathbf{M} and \mathbf{K} , the matrix \mathbf{A} is symmetric and positive definite. The spectral theorem (see, e.g., [21, Chap. 7] or [24, p. 61]) guarantees that \mathbf{A} has n positive real eigenvalues $\{\lambda_j\}_{j=1}^n$, and an orthonormal base of n real eigenvectors $\{\mathbf{u}_j\}_{j=1}^n$. It follows that $\{\mathbf{v}_j \equiv \mathbf{M}^{-1/2}\mathbf{u}_j\}_{j=1}^n$ is a basis of eigenvectors of the generalized problem (2.3), and we identify the frequencies as $\omega_j^2 = \lambda_j$. The orthogonality of the eigenvectors $\{\mathbf{u}_j\}$ of \mathbf{A} ensures the eigenvectors of (\mathbf{K}, \mathbf{M}) are \mathbf{M} -orthogonal:

(2.4)
$$\mathbf{v}_j^T \mathbf{M} \mathbf{v}_k = 0 \text{ if } j \neq k, \quad \mathbf{v}_j^T \mathbf{M} \mathbf{v}_j \neq 0.$$

It remains to demonstrate that the n eigenvalues are in fact distinct, or, in other words, that no eigenvalue may be associated with more than one eigendirection. If we

¹If the masses and lengths are uniform, say $m_j = 1/(n+1)$ and $\ell_j = 1/(n+1)$, and $\tau = 1$, then $\mathbf{M} = \mathbf{I}$ and \mathbf{K} becomes the familiar tridiagonal matrix that arises from the second-order finite difference discretization of the second spatial derivative; see, e.g., [25]. This reflects one way in which the wave equation $u_{tt} = u_{xx}$ can be derived as the limit of lumped masses.

express the *j*th row of $\mathbf{K}\mathbf{v} = \lambda \mathbf{M}\mathbf{v}$ in components with eigenvector $\mathbf{v} = [v_1, \dots, v_n]^T$, we find

(2.5)
$$\left(-\frac{\tau}{\ell_{j-1}}\right)v_{j-1} + \left(\frac{\tau}{\ell_{j-1}} + \frac{\tau}{\ell_j}\right)v_j + \left(-\frac{\tau}{\ell_j}\right)v_{j+1} = \lambda m_j v_j$$

where, by convention, $v_0 = v_{n+1} = 0$. Please note that if $v_1 = 0$, then the above implies that $v_2 = 0$ and so on. Hence, true eigenvectors obey $v_1 \neq 0$. Now if **w** also obeys $\mathbf{K}\mathbf{w} = \lambda \mathbf{M}\mathbf{w}$ for this same value of λ , then its components also satisfy (2.5), and so any linear combination of **v** and **w** will satisfy this equation. As $\mathbf{z} \equiv \mathbf{w} - (w_1/v_1)\mathbf{v}$ obeys $z_1 = 0$, it follows that $\mathbf{z} = \mathbf{0}$ and hence $\mathbf{w} = (w_1/v_1)\mathbf{v}$; i.e., our "new" eigenvector is simply a multiple of the original.

Since the eigenvectors $\{\mathbf{v}_k\}$ form a basis for *n*-dimensional space, we can write the solution to (2.1) as a time-varying linear combination:

$$\mathbf{y}(t) = \sum_{k=1}^{n} \gamma_k(t) \mathbf{v}_k$$

Substituting this expansion into the differential equation (2.1) gives

$$\sum_{k=1}^n \gamma_k''(t) \mathbf{M} \mathbf{v}_k = -\sum_{k=1}^n \gamma_k(t) \mathbf{K} \mathbf{v}_k = -\sum_{k=1}^n \omega_k^2 \gamma_k(t) \mathbf{M} \mathbf{v}_k.$$

To decompose into n independent equations, we premultiply this last equation by \mathbf{v}_j^T and appeal to (2.4), giving the familiar scalar equation

$$\gamma_j''(t) = -\omega_j^2 \gamma_j(t), \qquad j = 1, \dots, n$$

with solution

$$\gamma_j(t) = c_j \cos(\omega_j t) + s_j \sin(\omega_j t).$$

These c_j and s_j coefficients are determined by the pluck: $\mathbf{y}'(0) = \mathbf{0}$ implies each $s_j = 0$, while $\mathbf{y}(0) = \mathbf{y}_0$ implies that the c_j are the expansion coefficients of \mathbf{y}_0 in the eigenvector basis, which can be found by solving the linear system

$$\mathbf{y}_0 = \sum_{j=1}^n c_j \mathbf{v}_j.$$

3. Experimental Apparatus and Model Verification. How well does the model (2.1) just derived predict what really happens when a beaded string is plucked?

We investigate this question by conducting experiments on a high-precision monochord constructed by students in our laboratory at Rice University, shown in Figure 2. For the "massless string" we use a length of 0.015-inch-diameter nickel-plated steel musical wire donated by the Mapes Piano Wire Company. Tension is measured with a *force transducer* placed at the end of the string. The string then passes through a *collet*, which itself is mounted in a *collet vise*. The string proceeds through a photodetector that measures the vibrations at one point on the string. Brass beads are threaded onto the string, which continues through a second collet. (These beads have been carefully machined so as to snugly fit onto our wire.) Finally, the string is wound upon a spindle, which applies tension to the string. The experimenter winds the spin-



Fig. 2 The monochord loaded with beads.

dle until the string achieves a desired tension, then tightens the collet vises to fix the string at both ends (enforcing $y_0 = y_{n+1} = 0$).

A photodetector measures the displacement η_k at a single point along the string (not at a bead) at times $t_k = kh$ for some fixed time-step h. (The model only describes the motion of the beads, but the string itself must vibrate in concert: since we assume the string is perfectly elastic and the detector is placed between the fixed end and the first bead, these measurements are proportional to the first bead's displacement.) Consider a string loaded with five beads, as specified in Figure 3. We measure displacements for 10 sec. with h = 1/50000 sec., producing the samples $\{\eta_k\}$ shown on the left of Figure 4. (The magnitude of the displacements decay over the course of this ten second sample, reflecting some mild damping not captured by our simple physical model.) By analogy with the model (2.1), we expect that

(3.1)
$$\eta_k = \sum_{j=1}^n c_j \cos(t_k \sqrt{\lambda_j}) + \text{noise}$$

for some constants c_1, \ldots, c_n that depend on the initial pluck. The "noise" term captures errors both in our mathematical description of physical reality and in our ability to accurately measure that reality, as discussed in more detail in section 10.

To assess the accuracy of the model, we shall investigate whether the series of measurements $\{\eta_k\}$ for the five-beaded string in Figure 3 indeed oscillate at the frequencies predicted by the analysis in section 2. To do so, we compute the discrete Fourier transform (DFT) of the data. A detailed discussion of the DFT is beyond our scope, but excellent expositions can be found in [4, 24], and the operation can be implemented in just a few lines of MATLAB:

<pre>freq = 2*pi*[0:N-1]/N*sample_rate;</pre>	% set up vector of frequencies
<pre>semilogy(freq,abs(fft(eta)))</pre>	% plot magnitude of Fourier coefs
xlim([0 700])	% set axis to relevant frequencies

These operations produce a plot that shows the component of the signal over a range of frequencies as shown on the right in Figure 4. A signal behaving like $t \mapsto \cos(\omega t)$ should produce a peak in the DFT at $\omega \sec^{-1}$. By (3.1), we expect our signal to be dominated by combinations of $\cos(t\sqrt{\lambda_j})$ terms, and so we should find peaks precisely at $\sqrt{\lambda_j}$, where λ_j is an eigenvalue of (**K**, **M**). As the beads are not point masses, their finite diameters restrict the string's ability to vibrate freely; this could effectively shorten the total length of the string. In Figure 4 we predict a range for each eigenvalue, with the lower end determined by the actual length of the string, and upper end derived from the shorter string with the bead diameters removed.

Gómez et al. [14] provide complementary experimental work for continuous strings with one or two beads, including mode visualizations.



Fig. 3 Configuration of the five-bead experiment described in section 3. The string has total length of 112.4 cm and is drawn to a tension of 1.706×10^7 dyn. The beads have diameter of 5/8 in. (beads 1 and 4), 3/4 in. (beads 2 and 5), and 1/2 in. (bead 3). (Bead widths are exaggerated relative to the string length in our plots.)



Fig. 4 Displacement of the five-bead string in Figure 3 in the time domain (left) and frequency domain (right). Notice the five prominent peaks on the right, each corresponding to an eigenvalue of the pair (\mathbf{K}, \mathbf{M}) : The gray shaded regions denote the predicted location of the peaks from the mathematical model, $\omega_i = \sqrt{\lambda_i}$, based on a string of full length (left boundaries) and shortened by the bead diameters (right boundaries).

4. Inner Products and Jacobi Matrices. Having seen the predictive ability of the forward model (2.1), we now address a more interesting—and challenging—problem: Given knowledge of eigenvalues (e.g., as discerned from the peaks in Figure 4), can we "hear the beads on the string"? Can we determine the bead positions and masses? Numerous elegant algorithms solve this problem. Here we present a general approach that first recovers the symmetric tridiagonal matrix $M^{-1/2}KM^{-1/2}$ using an orthogonal polynomial algorithm of de Boor and Golub [8], followed by extraction of bead lengths and masses using a technique of Gladwell [12, section 4.4]. Some students may prefer Krein's continued fraction approach (see Supplement II of [10]), which is more elementary but specialized. We provide a detailed guide to this technique online [7], which can readily substitute for sections 4–7.

4.1. Orthogonal Polynomials with Discrete Inner Products. We begin the inversion process with an excursion into the beautiful, important subject of orthogonal polynomials (see [13] for theory and applications). Let \mathcal{P}_k denote the set of polynomials of degree k or less (with real coefficients). A k-degree polynomial p is monic if its leading coefficient is one, i.e., $p(z) = z^k + r(z)$ for some $r \in \mathcal{P}_{k-1}$. The set \mathcal{P}_{n-1} forms an n-dimensional vector space, and just as we have the dot product on n-dimensional Euclidean space, we can also specify an *inner product* on \mathcal{P}_n . Toward this end, choose real numbers for

nodes:
$$\xi_1 < \xi_2 < \cdots < \xi_n$$

and weights:
$$w_1, w_2, \ldots, w_n > 0$$
,

and define the *inner product* of polynomials p and q to be

(4.1)
$$\langle p,q\rangle = \sum_{j=1}^{n} w_j p(\xi_j) q(\xi_j).$$

One can readily verify that this function obeys the axioms required for an inner product on \mathcal{P}_{n-1} ; see, e.g., [21, p. 286]. We say p and q are orthogonal when $\langle p, q \rangle = 0$, and use the inner product to define a norm: $||p||^2 = \langle p, p \rangle \ge 0$.

The inner product is central to our development, which closely follows the important early work of de Boor and Golub [8]. The first step involves the construction of a sequence of monic polynomials of increasing degree that are all orthogonal to one another. The monic degree-0 polynomial must be

$$p_0(z) = 1.$$

To build the monic degree-1 polynomial, multiply $p_0(z)$ by z (to get a monic polynomial of degree 1), then subtract out the projection of zp_0 onto the span of p_0 . This amounts to a single step of the Gram–Schmidt orthogonalization process [21, p. 307]:

$$p_1(z) = zp_0(z) - \frac{\langle zp_0, p_0 \rangle}{\langle p_0, p_0 \rangle} p_0(z) =: zp_0(z) - a_1 p_0(z).$$

Step k + 1 follows similarly, via Gram–Schmidt:

$$p_{k+1}(z) = zp_k(z) - \sum_{j=1}^k \frac{\langle zp_k, p_j \rangle}{\langle p_j, p_j \rangle} p_j(z).$$

Now we make an essential simplifying observation: p_k must be orthogonal to any lower degree polynomial (which must be a linear combination of p_0, \ldots, p_{k-1} , as these orthogonal polynomials form a basis for \mathcal{P}_{k-1}). Since $zp_j \in \mathcal{P}_{j+1}$, notice that

$$\langle zp_k, p_j \rangle = \langle p_k, zp_j \rangle = 0, \qquad j < k-1.$$

The formula for p_{k+1} thus reduces to a *three-term recurrence relation*:

$$p_{k+1}(z) = zp_k(z) - \frac{\langle zp_k, p_k \rangle}{\langle p_k, p_k \rangle} p_k(z) - \frac{\langle zp_k, p_{k-1} \rangle}{\langle p_{k-1}, p_{k-1} \rangle} p_{k-1}(z)$$

The coefficient of p_{k-1} is always negative: since $zp_{k-1} = p_k + r$ for some $r \in \mathcal{P}_{k-1}$,

$$\frac{\langle zp_k, p_{k-1} \rangle}{\langle p_{k-1}, p_{k-1} \rangle} = \frac{\langle p_k, zp_{k-1} \rangle}{\langle p_{k-1}, p_{k-1} \rangle} = \frac{\langle p_k, p_k + r \rangle}{\langle p_{k-1}, p_{k-1} \rangle} = \frac{\langle p_k, p_k \rangle}{\langle p_{k-1}, p_{k-1} \rangle} = \frac{\|p_k\|^2}{\|p_{k-1}\|^2} \ge 0.$$

Thus we write

$$p_{k+1}(z) = zp_k(z) - \frac{\langle zp_k, p_k \rangle}{\langle p_k, p_k \rangle} p_k(z) - \frac{\|p_k\|^2}{\|p_{k-1}\|^2} p_{k-1}(z)$$

(4.2)
$$=: zp_k(z) - a_{k+1}p_k(z) - b_k^2 p_{k-1}(z);$$

the label " b_k^2 " emphasizes that the constant $b_k^2 := \|p_k\|^2 / \|p_{k-1}\|^2$ is positive. We see that orthogonal polynomials intimately link to the constants a_1, \ldots, a_n and b_1, \ldots, b_{n-1} . File this fact away for the moment, and return to the matrices at hand.

4.2. Eigenvalues of Symmetric Tridiagonal Matrices. The matrices modeling beaded string vibrations can be written in the form

$$\mathbf{A}_{n} = \mathbf{M}^{-1/2} \mathbf{K} \mathbf{M}^{-1/2} = \begin{bmatrix} a_{1} & b_{1} & & \\ b_{1} & a_{2} & \ddots & \\ & \ddots & \ddots & b_{n-1} \\ & & b_{n-1} & a_{n} \end{bmatrix},$$

where $a_k = \tau(1/\ell_{k-1} + 1/\ell_k)/m_k$ and $b_k = -\tau/(\ell_k \sqrt{m_k m_{k+1}}) < 0$; see (2.5). The matrix \mathbf{A}_n (more precisely, $-\mathbf{A}_n$) is an example of a *Jacobi matrix*, a widely studied family having many fascinating properties. To analyze the eigenvalues, consider the characteristic polynomial $q_n(z) := \det(z\mathbf{I} - \mathbf{A}_n)$. Expand this determinant along the last row and column of $z\mathbf{I} - \mathbf{A}_n$ to arrive at the formula

$$\det(z\mathbf{I} - \mathbf{A}_n) = (z - a_n)\det(z\mathbf{I} - \mathbf{A}_{n-1}) - b_{n-1}^2\det(z\mathbf{I} - \mathbf{A}_{n-2}).$$

In other words,

(4.3)
$$q_n(z) = (z - a_n)q_{n-1}(z) - b_{n-1}^2q_{n-2}(z)$$

which is exactly the recurrence (4.2) for the orthogonal polynomial p_n . Jacobi matrices and orthogonal polynomials thus enjoy a rich intertwined history, with applications spanning from numerical analysis to mathematical physics.

5. The Inverse Eigenvalue Problem for Jacobi Matrices. We now wish to use the connection between Jacobi matrices and orthogonal polynomials to build a unique Jacobi matrix using only knowledge of eigenvalues. In section 6 we shall adjust our approach to suit the data typically available from our laboratory, and then show how to extract bead positions and masses from Jacobi matrix entries in section 7.

Suppose we do not know the entries that make up the matrix \mathbf{A}_n , but we know its eigenvalues, along with those of its $(n-1) \times (n-1)$ block, \mathbf{A}_{n-1} :

eigenvalues of
$$\mathbf{A}_n$$
: $\lambda_1 < \lambda_2 < \cdots < \lambda_n$;
eigenvalues of \mathbf{A}_{n-1} : $\mu_1 < \mu_2 < \cdots < \mu_{n-1}$.

We can use special orthogonal polynomials to recover the entries of \mathbf{A}_n . First note that the eigenvalues of \mathbf{A}_n and \mathbf{A}_{n-1} never coincide, and in fact weave between each other,

(5.1)
$$\lambda_1 < \mu_1 < \lambda_2 < \dots < \lambda_{n-1} < \mu_{n-1} < \lambda_n$$

an instance of Cauchy's interlacing theorem; see, e.g., [21, p. 552].

5.1. Custom-Tailoring an Inner Product. Given the parallel between the entries in \mathbf{A}_n and the coefficients of the orthogonal polynomial recurrence relation (4.2), one might wonder: Does there exist some inner product (defined by nodes $\xi_1 < \cdots < \xi_n$ and (positive) weights w_1, \ldots, w_n) in which the characteristic polynomials

(5.2)
$$q_n(z) = \det(z\mathbf{I} - \mathbf{A}_n) = \prod_{j=1}^n (z - \lambda_j),$$

(5.3)
$$q_{n-1}(z) = \det(z\mathbf{I} - \mathbf{A}_{n-1}) = \prod_{j=1}^{n-1} (z - \mu_j)$$

are monic orthogonal polynomials? Given such an inner product, we could identify

(5.4)
$$a_n = \frac{\langle zq_{n-1}, q_{n-1} \rangle}{\langle q_{n-1}, q_{n-1} \rangle},$$

then run recurrence (4.2) backward to compute

(5.5)
$$b_{n-1}^2 q_{n-2}(z) = (z - a_n)q_{n-1}(z) - q_n(z).$$

To split b_{n-1}^2 from q_{n-2} , use the fact that q_{n-2} must be monic. Repeat this process with q_{n-2} in (5.4) to recover a_{n-1} , then use (5.5) for b_{n-2} , and so on, to build \mathbf{A}_n .

To discover such an inner product (i.e., suitable nodes and weights), we will explore some properties this inner product would need to obey. For one thing, q_n should be orthogonal to any lower degree polynomial in \mathcal{P}_{n-1} . We can always build a polynomial $r \in \mathcal{P}_{n-1}$ that interpolates $q_n(z)$ at the *n* nodes $z = \xi_1, \ldots, \xi_n$:

$$r(\xi_j) = q_n(\xi_j), \quad j = 1, \dots, n.$$

(This is an extension of the fact that one can always construct a linear polynomial through any two points, a quadratic through any three points, etc.) Since $r \in \mathcal{P}_{n-1}$,

$$0 = \langle q_n, r \rangle = \sum_{j=1}^n w_j q_n(\xi_j) r(\xi_j) = \sum_{j=1}^n w_j q_n(\xi_j)^2.$$

As the weights must be positive, we conclude that

$$q_n(\xi_j) = 0, \qquad j = 1, \dots, n.$$

That is, the *n* nodes ξ_1, \ldots, ξ_n must coincide with the *n* roots of q_n :

(5.6)
$$\xi_j = \lambda_j, \qquad j = 1, \dots, n.$$

To finish determining the inner product, we must specify the weights w_1, \ldots, w_n . In this endeavor we are aided by the Lagrange interpolating polynomials, objects that arise in a basic numerical analysis class (see, e.g., [26, section 6.2]). Define $\delta_k \in \mathcal{P}_{n-1}$ by the following *n* conditions: δ_k passes through zero at the nodes ξ_j for $j \neq k$, while taking the value one at ξ_k . You can quickly verify that such polynomials have the form

(5.7)
$$\delta_k(z) = \prod_{\substack{j=1\\j \neq k}}^n \frac{z - \xi_j}{\xi_k - \xi_j} \in \mathcal{P}_{n-1}.$$

Since $q_n(\xi_k) = 0$ for all k,

(5.8)
$$q'_n(\xi_k) = \lim_{z \to \xi_k} \frac{q_n(z) - q_n(\xi_k)}{z - \xi_k} = \lim_{z \to \xi_k} \frac{(z - \xi_k) \prod_{\substack{j=1; j \neq k \\ z - \xi_k}}^n (z - \xi_j)}{z - \xi_k} = \prod_{\substack{j=1 \\ j \neq k}}^n (\xi_k - \xi_j),$$

ONE CAN HEAR THE COMPOSITION OF A STRING

the product of denominators in the formula (5.7) for δ_k . Since $q_{n-1} \in \mathcal{P}_{n-1}$ is monic, the leading term of δ_k matches that of $q_{n-1}/q'_n(\xi_k)$. Hence for some $r_k \in \mathcal{P}_{n-2}$,

(5.9)
$$\delta_k(z) = \frac{q_{n-1}(z)}{q'_n(\xi_k)} + r_k(z)$$

By the definition of the inner product and the nature of the Lagrange polynomials,

$$\langle q_{n-1}, \delta_k \rangle = \sum_{j=1}^n w_j q_{n-1}(\xi_j) \delta_k(\xi_j) = w_k q_{n-1}(\xi_k).$$

Solve this equation for the weight w_k , using the form (5.9) and the fact that $r_k \in \mathcal{P}_{n-2}$ to find

$$w_{k} = \frac{\langle q_{n-1}, \delta_{k} \rangle}{q_{n-1}(\xi_{k})} = \frac{\langle q_{n-1}, q_{n-1} \rangle}{q'_{n}(\xi_{k})q_{n-1}(\xi_{k})} = \frac{\|q_{n-1}\|^{2}}{q'_{n}(\xi_{k})q_{n-1}(\xi_{k})}, \qquad k = 1, \dots, n.$$

Do you notice a subtle drawback of this formula? The expression for w_k involves $||q_{n-1}||$, which we can only compute once all the weights are known. There is an easy dodge: the term $||q_{n-1}||^2$ is independent of k, so it affects all the weights the same way. Orthogonality—the property we care most about—is independent of the collective scaling of the weights, so we can select any convenient scaling, such as

(5.10)
$$w_k = \frac{1}{q'_n(\xi_k)q_{n-1}(\xi_k)}, \qquad k = 1, \dots, n.$$

There can be no division by zero in this formula, since the nodes $\xi_k = \lambda_k$ never coincide with the roots μ_1, \ldots, μ_{n-1} of q_{n-1} by the interlacing property (5.1); however, computational subtleties can emerge when computing these weights, as they can vary substantially in magnitude with the index k, especially when n is large.

Equipped with these nodes and weights, we can repeatedly use (5.4)–(5.5) to reconstruct \mathbf{A}_n . Before so doing, we must make some accommodations for the data that are readily accessible in our laboratory.

6. Adapting the Inverse Jacobi Algorithm for Experimental Data. In our string laboratory we experimentally determine the length of the string, $\ell = \sum_{k=0}^{n} \ell_k$, the tension, τ , and the eigenvalues $\lambda_1 < \cdots < \lambda_n$ of the matrix $\mathbf{A}_n = \mathbf{M}^{-1/2} \mathbf{K} \mathbf{M}^{-1/2}$ having the form

$$\begin{bmatrix} \frac{\tau}{\ell_0 m_1} + \frac{\tau}{\ell_1 m_1} & -\frac{\tau}{\ell_1 \sqrt{m_1 m_2}} \\ -\frac{\tau}{\ell_1 \sqrt{m_1 m_2}} & \frac{\tau}{\ell_1 m_2} + \frac{\tau}{\ell_2 m_2} & \ddots \\ & \ddots & \ddots & -\frac{\tau}{\ell_{n-1} \sqrt{m_n m_{n-1}}} \\ & -\frac{\tau}{\ell_{n-1} \sqrt{m_n m_{n-1}}} & \frac{\tau}{\ell_{n-1} m_n} + \frac{\tau}{\ell_n m_n} \end{bmatrix}$$

These eigenvalues give the nodes of our inner product, $\xi_k := \lambda_k, k = 1, ..., n$. To compute the weights for the inner product, we need the second set of eigenvalues

STEVEN J. COX, MARK EMBREE, AND JEFFREY M. HOKANSON

 $\mu_1 < \cdots < \mu_{n-1}$ coming from \mathbf{A}_{n-1} . How can we experimentally measure these values? This would amount to fixing the *n*th mass, thus making the string one bead shorter. One could imagine building an apparatus to implement this condition, but it demands knowledge of the internal composition of the string: namely, the *location* of the *n*th mass, hence the quantity ℓ_n . We prefer a different strategy that will avoid such an intrusion into the interior of the string. We have laboratory data for the string fixed at both ends (the "fixed-fixed" string), giving measured eigenvalues $\lambda_1 < \cdots < \lambda_n$. Now suppose we leave the string configuration the same, except now the right end is attached by a ring to a frictionless vertical pole, so that end remains level with the *n*th bead (the "fixed-flat" string): measure the eigenvalues $\hat{\lambda}_1 < \cdots < \hat{\lambda}_n$ of this modified string. (In section 8 we will find a more palatable route to these data using only our regular fixed-fixed experiment.)

The beads of the fixed-flat string obey the same equations of motion (2.1) derived for the fixed-fixed string save for the last bead, which is governed by

$$m_n y_n''(t) = -\tau \Big(\frac{y_n(t) - y_{n-1}(t)}{\ell_{j-1}} \Big);$$

that is, $y_{n+1} = y_n$ replaces the fixed condition $y_{n+1} = 0$. This change in the equation for the *n*th bead gives a fixed-flat matrix $\widehat{\mathbf{A}}_n$ that differs from the fixed-fixed matrix only in the bottom-right entry:

(6.1)
$$\widehat{\mathbf{A}}_n = \mathbf{A}_n - \frac{\tau}{\ell_n m_n} \mathbf{e}_n \mathbf{e}_n^{\mathrm{T}}.$$

Here \mathbf{e}_k denotes the *k*th column of the identity matrix, whose dimension is clear from the context. Label the (n, n) entry of $\widehat{\mathbf{A}}_n$ as \widehat{a}_n (to distinguish it from the (n, n) entry a_n of \mathbf{A}_n). Now partition the two Jacobi matrices as

$$\mathbf{A}_{n} = \begin{bmatrix} \mathbf{A}_{n-1} & b_{n-1}\mathbf{e}_{n-1} \\ b_{n-1}\mathbf{e}_{n-1}^{\mathrm{T}} & a_{n} \end{bmatrix} = \begin{bmatrix} \mathbf{A}_{n-2} & b_{n-2}\mathbf{e}_{n-2} & \mathbf{0} \\ b_{n-2}\mathbf{e}_{n-2}^{\mathrm{T}} & a_{n-1} & b_{n-1} \\ \mathbf{0} & b_{n-1} & a_{n} \end{bmatrix}$$

and

$$\widehat{\mathbf{A}}_{n} = \begin{bmatrix} \mathbf{A}_{n-1} & b_{n-1}\mathbf{e}_{n-1} \\ b_{n-1}\mathbf{e}_{n-1}^{\mathrm{T}} & \widehat{a}_{n} \end{bmatrix} = \begin{bmatrix} \mathbf{A}_{n-2} & b_{n-2}\mathbf{e}_{n-2} & \mathbf{0} \\ b_{n-2}\mathbf{e}_{n-2}^{\mathrm{T}} & a_{n-1} & b_{n-1} \\ \mathbf{0} & b_{n-1} & \widehat{a}_{n} \end{bmatrix}.$$

With these forms, the characteristic polynomials for \mathbf{A}_n and $\widehat{\mathbf{A}}_n$ are

$$q_n(z) = \det(z\mathbf{I} - \mathbf{A}_n) = (z - a_n)\det(z\mathbf{I} - \mathbf{A}_{n-1}) + b_{n-1}\det\left(\begin{bmatrix}z\mathbf{I} - \mathbf{A}_{n-2} & \mathbf{0}\\-b_{n-2}\mathbf{e}_{n-2}^{\mathrm{T}} & -b_{n-1}\end{bmatrix}\right),$$
$$\widehat{q}_n(z) = \det(z\mathbf{I} - \widehat{\mathbf{A}}_n) = (z - \widehat{a}_n)\det(z\mathbf{I} - \mathbf{A}_{n-1}) + b_{n-1}\det\left(\begin{bmatrix}z\mathbf{I} - \mathbf{A}_{n-2} & \mathbf{0}\\-b_{n-2}\mathbf{e}_{n-2}^{\mathrm{T}} & -b_{n-1}\end{bmatrix}\right),$$

whose difference is

(6.2)
$$\widehat{q}_n(z) - q_n(z) = (a_n - \widehat{a}_n) \det(z\mathbf{I} - \mathbf{A}_{n-1})$$
$$= (a_n - \widehat{a}_n)q_{n-1}(z).$$

168

ONE CAN HEAR THE COMPOSITION OF A STRING

This suggests a way to access the eigenvalues $\mu_1 < \cdots < \mu_{n-1}$ of \mathbf{A}_{n-1} needed for the weights: from the measured eigenvalues $\lambda_1, \ldots, \lambda_n$ and $\hat{\lambda}_1, \ldots, \hat{\lambda}_n$, build the characteristic polynomials q_n and \hat{q}_n ; by (6.2), μ_1, \ldots, μ_{n-1} are roots of their difference. In practice this is a bad idea, for polynomial roots often vary greatly when coefficients are perturbed (inevitable with experimental data). De Boor and Golub [8] suggest a better route to the weights that avoids the eigenvalues of \mathbf{A}_{n-1} . The formula (5.10) for the weights only requires the values $q_{n-1}(\lambda_k)$. We can access these quantities by recalling that the eigenvalues $\lambda_1, \ldots, \lambda_n$ are the roots of q_n , hence from (6.2),

(6.3)
$$\widehat{q}_n(\lambda_k) = (a_n - \widehat{a}_n)q_{n-1}(\lambda_k), \qquad k = 1, \dots, n.$$

This equation specifies $q_{n-1}(\lambda_k) = q_{n-1}(\xi_k)$ up to the term $\hat{a}_n - a_n$, which is constant for all k values, and hence can be omitted from all weights (like $||q_{n-1}||^2$ earlier), as we are unconcerned about scalings of the inner product. Thus use the scaled weights

(6.4)
$$w_k = \frac{1}{q'_n(\xi_k)\hat{q}_n(\xi_k)}, \qquad k = 1, \dots, n.$$

To obtain q_{n-1} , use (6.2) along with the fact that q_{n-1} must be monic, i.e., $a_n - \hat{a}_n$ is the leading coefficient of the degree n-1 polynomial $\hat{q}_n - q_n$. Since

$$q_n(z) = \prod_{j=1}^n (z - \lambda_j) = z^n - \left(\sum_{j=1}^n \lambda_j\right) z^{n-1} + \cdots,$$
$$\widehat{q}_n(z) = \prod_{j=1}^n (z - \widehat{\lambda}_j) = z^n - \left(\sum_{j=1}^n \widehat{\lambda}_j\right) z^{n-1} + \cdots,$$

we can compute

$$a_n - \widehat{a}_n = \sum_{j=1}^n (\lambda_j - \widehat{\lambda}_j)$$

and thus

(6.5)
$$q_{n-1}(z) = \frac{\widehat{q}_n(z) - q_n(z)}{\sum_{j=1}^n (\lambda_j - \widehat{\lambda}_j)}.$$

6.1. An Algorithm for Recovering the Jacobi Matrix. We are ready to build \mathbf{A}_n from the measured eigenvalues $\lambda_1 < \cdots < \lambda_n$ and $\hat{\lambda}_1 < \cdots < \hat{\lambda}_n$ from the fixed-fixed and fixed-flat strings. We will represent the orthogonal polynomials q_k by their values at the nodes ξ_1, \ldots, ξ_n , rather than by their coefficients. As such, (5.5) will not suffice to reveal b_k^2 . Instead, take the inner product of (5.5) (with k replacing n-1) with $zq_k(z)$ to obtain

(6.6)
$$b_k^2 = \frac{\langle zq_k, zq_k \rangle - a_{k+1} \langle q_k, zq_k \rangle - \langle q_{k+1}, zq_k \rangle}{\langle q_{k-1}, zq_k \rangle}$$

The denominator appears troubling, as q_{k-1} will not have been determined at this stage. There is a slick work-around:

$$\langle q_{k-1}, zq_k \rangle = \langle zq_{k-1}, q_k \rangle = \langle q_k + r, q_k \rangle = \langle q_k, q_k \rangle$$

for some $r \in \mathcal{P}_{k-1}$. Hence we can use the formula for b_k^2 given in step 5(a) in Figure 5. Since $b_k = -\tau/(\ell_k \sqrt{m_k m_{k+1}}) < 0$, we define $b_k := -\sqrt{b_k^2}$. 1. Use (5.6) to determine the nodes of the inner product: $\xi_k := \lambda_k.$ 2. Use (6.4) to construct the weights, with $q'_n(\xi_k)$ given by (5.8): $w_{k} := \frac{1}{q_{n}'(\xi_{k})\widehat{q}_{n}(\xi_{k})} = \frac{1}{\left(\prod_{j=1}^{n}(\xi_{k} - \xi_{j})\right) \left(\prod_{j=1}^{n}(\xi_{k} - \widehat{\lambda}_{j})\right)}.$ 3. Determine the values of q_n and q_{n-1} at the nodes; see (6.5): $q_n(\xi_k) := 0,$ $q_{n-1}(\xi_k) := \frac{\widehat{q}_n(\xi_k) - q_n(\xi_k)}{\sum_{j=1}^n (\lambda_j - \widehat{\lambda}_j)} = \frac{\prod_{j=1}^n (\xi_k - \lambda_j)}{\sum_{j=1}^n (\lambda_j - \widehat{\lambda}_j)}.$ 4. Compute, via the inner product (4.1). $a_n := \frac{\langle zq_{n-1}, q_{n-1} \rangle}{\langle q_{n-1}, q_{n-1} \rangle} = \frac{\sum_{j=1}^n w_j \xi_j q_{n-1}(\xi_j)^2}{\sum_{j=1}^n w_j q_{n-1}(\xi_j)^2}.$ 5. For $k = n - 1, n - 2, \dots, 1$ (a) Compute $b_k = -\sqrt{b_k^2}$ via the approach described after (6.6): $b_k^2 := \frac{\langle zq_k, zq_k \rangle - a_{k+1} \langle q_k, zq_k \rangle - \langle q_{k+1}, zq_k \rangle}{\langle q_k, q_k \rangle}.$ (b) Compute q_{k-1} at the nodes: $q_{k-1}(\xi_j) := \frac{(\xi_j - a_{k+1})q_k(\xi_j) - q_{k+1}(\xi_j)}{b_i^2}.$ (c) Define $a_k := \frac{\langle zq_{k-1}, q_{k-1} \rangle}{\langle q_{k-1}, q_{k-1} \rangle}.$

Fig. 5 Algorithm for recovering the main diagonal (a_1, \ldots, a_n) and off-diagonal (b_1, \ldots, b_{n-1}) of the Jacobi matrix \mathbf{A}_n from the fixed-fixed and fixed-flat eigenvalues $\lambda_1 < \cdots < \lambda_n$ and $\widehat{\lambda}_1 < \cdots < \widehat{\lambda}_n$. Adapted from de Boor and Golub [8].

7. Relating the Jacobi Matrix Approach to Beaded Strings. We can use the procedure detailed in Figure 5 to recover the entries of A_n , but how then do we access the masses and lengths from which these entries derive? Gladwell describes a procedure for mass-spring networks that we can adapt for our setting [12, p. 77ff].

Let **e** denote the vector with one in each entry, $\mathbf{e} = [1, 1, ..., 1]^{\mathrm{T}}$, while \mathbf{e}_1 and \mathbf{e}_n denote the first and last columns of the identity matrix. Observe from (2.2) that

(7.1)
$$\mathbf{K}\mathbf{e} = \frac{\tau}{\ell_0}\mathbf{e}_1 + \frac{\tau}{\ell_n}\mathbf{e}_n,$$

so knowledge of **K** would reveal ℓ_0 and ℓ_n via (7.1), from which we could determine the other lengths. However, at this stage we only know \mathbf{A}_n , not **K**. Using **K** = $\mathbf{M}^{1/2}\mathbf{A}_{n}\mathbf{M}^{1/2}$, write

$$\mathbf{M}^{1/2}\mathbf{A}_{n}\mathbf{M}^{1/2}\mathbf{e} = \frac{\tau}{\ell_{0}}\mathbf{e}_{1} + \frac{\tau}{\ell_{n}}\mathbf{e}_{n}.$$

Premultiplying by $\mathbf{M}^{-1/2}$ and labeling $\mathbf{d} := \mathbf{M}^{1/2} \mathbf{e} = [\sqrt{m_1}, \dots, \sqrt{m_n}]^T$ gives

(7.2)
$$\mathbf{A}_n \mathbf{d} = \frac{\tau}{\ell_0 \sqrt{m_1}} \mathbf{e}_1 + \frac{\tau}{\ell_n \sqrt{m_n}} \mathbf{e}_n.$$

The vector \mathbf{d} that solves this linear system will thus reveal the bead masses. We could solve the system (7.2) for \mathbf{d} using Gaussian elimination, except we lack a formula for the right-hand side. We can obtain the right-hand side up to a scaling factor through the following strategy. Solve the two linear systems

$$\mathbf{A}_n \mathbf{x} = \mathbf{e}_1, \qquad \mathbf{A}_n \mathbf{y} = \mathbf{e}_n$$

for the vectors \mathbf{x} and \mathbf{y}^2 . Now

$$\mathbf{A}_n \Big(\frac{\tau}{\ell_0 \sqrt{m_1}} \mathbf{x} + \frac{\tau}{\ell_n \sqrt{m_n}} \mathbf{y} \Big) = \frac{\tau}{\ell_0 \sqrt{m_1}} \mathbf{e}_1 + \frac{\tau}{\ell_n \sqrt{m_n}} \mathbf{e}_n = \mathbf{A}_n \mathbf{d}.$$

Since \mathbf{A}_n is nonsingular (it is positive definite, as described in section 2), we can access \mathbf{d} as

(7.3)
$$\mathbf{d} = \frac{\tau}{\ell_0 \sqrt{m_1}} \mathbf{x} + \frac{\tau}{\ell_n \sqrt{m_n}} \mathbf{y}.$$

One obstacle remains: What are the coefficients $\tau/(\ell_0\sqrt{m_1})$ and $\tau/(\ell_n\sqrt{m_n})$? We can resolve this issue using our experimental data. Since the trace of a matrix (sum of the diagonal elements) is the sum of the eigenvalues [21, p. 494], the experimentally measured eigenvalues reveal, by way of (6.1),

(7.4)
$$\frac{\tau}{\ell_n m_n} = \operatorname{tr}(\mathbf{A}_n) - \operatorname{tr}(\widehat{\mathbf{A}}_n) = \sum_{k=1}^n (\lambda_k - \widehat{\lambda}_k).$$

Thus the second coefficient in (7.3) is given by

(7.5)
$$\frac{\tau}{\ell_n \sqrt{m_n}} = \frac{\tau \sqrt{m_n}}{\ell_n m_n} = \sqrt{m_n} \sum_{k=1}^n (\lambda_k - \widehat{\lambda}_k).$$

To obtain the first coefficient in (7.3), note that the *n*th row of (7.3) gives

$$\sqrt{m_n} = \frac{\tau}{\ell_0 \sqrt{m_1}} x_n + \frac{\tau}{\ell_n \sqrt{m_n}} y_n,$$

which can be rearranged as

(7.6)
$$\frac{\tau}{\ell_0\sqrt{m_1}} = \frac{1}{x_n} \left(\sqrt{m_n} - \frac{\tau}{\ell_n\sqrt{m_n}}y_n\right) = \frac{\sqrt{m_n}}{x_n} \left(1 - y_n \sum_{k=1}^n (\lambda_k - \widehat{\lambda}_k)\right).$$

²Alternatively, the entries in \mathbf{x} and \mathbf{y} can be obtained using explicit formulas involving minors of \mathbf{A}_n ; see equation (4.4.12) in Gladwell [12].

1. Solve
$$\mathbf{A}_n \mathbf{x} = \mathbf{e}_1$$
 and $\mathbf{A}_n \mathbf{y} = \mathbf{e}_n$ for \mathbf{x} and \mathbf{y} .
2. $\gamma_2 := \sum_{j=1}^n (\lambda_j - \hat{\lambda}_j)$ and $\gamma_1 := (1 - y_n \gamma_2)/x_n$.
3. $\widetilde{\mathbf{d}} := \mathbf{d}/\sqrt{m_n} = \gamma_1 \mathbf{x} + \gamma_2 \mathbf{y}$.
4. $\ell_j m_n := -\tau/(b_j \widetilde{d}_j \widetilde{d}_{j+1})$ for $j = 1, \dots, n-1$.
 $\ell_0 m_n := \tau/(a_1 \widetilde{d}_1^2 - \tau/(\ell_1 m_n))$.
 $\ell_n m_n := \tau/(a_n \widetilde{d}_n^2 - \tau/(\ell_{n-1} m_n))$.
5. $m_n := \left(\sum_{j=0}^n \ell_j m_n\right)/\ell$.

Fig. 6 Algorithm for extracting bead masses and locations from a Jacobi matrix \mathbf{A}_n (main diagonal entries a_1, \ldots, a_n and off-diagonal entries b_1, \ldots, b_{n-1}), given knowledge of the tension τ , total length of the string, ℓ , and the fixed-fixed and fixed-flat eigenvalues. (For the symmetric strings in section 8, be sure to use $\ell = L/2$.)

Given coefficients (7.5) and (7.6), (7.3) gives **d** up to a factor of $\sqrt{m_n}$:

(7.7)
$$\widetilde{\mathbf{d}} := \frac{\mathbf{d}}{\sqrt{m_n}} = \frac{1}{x_n} \Big(1 - y_n \sum_{k=1}^n (\lambda_k - \widehat{\lambda}_k) \Big) \mathbf{x} + \Big(\sum_{k=1}^n (\lambda_k - \widehat{\lambda}_k) \Big) \mathbf{y},$$

and so

(7.8)
$$\mathbf{K}/m_n = (\mathbf{M}^{1/2} \mathbf{A}_n \mathbf{M}^{1/2})/m_n = \operatorname{diag}(\widetilde{\mathbf{d}}) \mathbf{A}_n \operatorname{diag}(\widetilde{\mathbf{d}}).$$

Recalling (2.2), the (j, j+1) entries of (7.8) give $-\tau/(\ell_j m_n) = b_j \tilde{d}_j \tilde{d}_{j+1}$, i.e.,

(7.9)
$$\ell_j m_n = -\frac{\tau}{b_j \tilde{d}_j \tilde{d}_{j+1}}, \qquad j = 1, \dots, n-1.$$

We can then find the end lengths from the (1,1) and (n,n) entries of (7.8):

(7.10)
$$\ell_0 m_n = \frac{\tau}{a_1 \tilde{d}_1^2 + \tau/(\ell_1 m_n)}, \qquad \ell_n m_n = \frac{\tau}{a_n \tilde{d}_n^2 + \tau/(\ell_{n-1} m_n)}.$$

We now have expressions for all the masses via (7.7) and lengths via (7.9)–(7.10), up to the scaling factor m_n . We do not know m_n (as it is an interior property of the string), but we do know the total string length, ℓ , and from this we can compute $m_n = \left(\sum_{j=0}^n \ell_j m_n\right)/\ell$. This overall recovery process is summarized in Figure 6.

7.1. The Special Case of n = 1 Bead. The case of n = 1 bead, in which case $\mathbf{A}_n = a_1$ is a 1×1 matrix, requires special attention. Recovery of \mathbf{A}_n from λ_1 and $\hat{\lambda}_1$ becomes trivial: $\mathbf{A}_n = \lambda_1$. The reader can then verify, starting from (7.4), that

$$\ell_0 m_1 = \frac{\tau}{\widehat{\lambda}_1}, \qquad \ell_1 m_1 = \frac{\tau}{\lambda_1 - \widehat{\lambda}_1}, \qquad m_1 = \frac{\tau \lambda_1}{\ell \widehat{\lambda}_1 (\lambda_1 - \widehat{\lambda}_1)}$$



Fig. 7 A string with six beads with lengths and masses arranged symmetrically about the middle of the string, denoted by the vertical gray line.

8. Symmetrically Loaded Strings. Having seen how to determine the bead locations and masses from two sets of eigenvalues, we must now address the practical issue of experimentally measuring these data. When the string is fixed at both ends, we can approximate the eigenvalues using the reliable experiments performed in section 3. The fixed–flat string poses a greater challenge. Fortunately, in one interesting special case we can determine these eigenvalues without any modification to the experimental apparatus. (Boyko and Pivovarchik recently proposed an alternative: clamping the string at an interior point and measuring the spectra on each substring [3].)

Suppose that the number of beads, N, is even, and the masses of the beads and lengths between them are symmetric about the midpoint of the string. To distinguish this case from the general scenario considered earlier, we shall denote these masses by M_1, \ldots, M_N , and the lengths by L_0, \ldots, L_N , with the total length $L = \sum_{j=0}^N L_j$. Thus, the symmetric arrangement requires that

$$M_{1} = M_{N} \qquad L_{0} = L_{N}$$

$$M_{2} = M_{N-1} \qquad L_{1} = L_{N-1}$$

$$\vdots \qquad \vdots$$

$$M_{N/2} = M_{N/2+1}, \qquad L_{N/2-1} = L_{N/2+1}$$

as illustrated for N = 6 in Figure 7. We can install such a symmetric arrangement on our laboratory's monochord (as pictured in Figure 2), then experimentally measure the eigenvalues when both ends of the string are fixed; let us label these eigenvalues $\Lambda_1 < \Lambda_2 < \cdots < \Lambda_N$. In Figure 8 we show the eigenvectors for the configuration shown in Figure 7. This illustration reveals a remarkable property: eigenvectors corresponding to the odd eigenvalues Λ_1 , Λ_3 , and Λ_5 are all symmetric about the middle of the string: If we cut the string in half, these eigenvectors would be *fixed* on the left end and be *flat* at the right. On the other hand, eigenvectors corresponding to the even eigenvalues Λ_2 , Λ_4 , and Λ_6 are all antisymmetric about the midpoint: If we cut the string in half, these eigenvectors would be *fixed at zero at both ends*. These observations hold for all symmetric string configurations, and they hint at a key fact:

The N eigenvalues of a symmetric beaded string fixed at both ends exactly match the N/2 fixed-fixed and N/2 fixed-flat eigenvalues associated with half of the string.

More precisely, consider a length $\ell = L/2$ string having n = N/2 beads with masses

$$m_j = M_j, \quad j = 1, \dots, n,$$

separated by lengths

$$\ell_j = L_j, \quad j = 0, \dots, n-1,$$

$$\ell_n = L_n/2.$$



Fig. 8 Eigenvectors for the string shown in Figure 7. In each plot, the vertical displacement of the kth mass indicates the kth entry of the corresponding eigenvector \mathbf{v}_j of $\mathbf{M}^{-1}\mathbf{K}$.



Fig. 9 Eigenvectors for the half-string corresponding to the symmetric configuration in Figure 7. The eigenvectors on the left satisfy the fixed-flat conditions; those on the right satisfy fixed-fixed conditions. Compare to the eigenvectors for the symmetric string in Figure 8.

The odd eigenvalues for the symmetric string match the eigenvalues for the fixed–flat half-string,

$$\widehat{\lambda}_j = \Lambda_{2j-1}, \quad j = 1, \dots, n,$$

while the even eigenvalues for the symmetric string match the eigenvalues for the fixed-fixed half string,

$$\lambda_j = \Lambda_{2j}, \quad j = 1, \dots, n.$$

Figure 9 shows eigenvectors for the half-string corresponding to the symmetric string

174

in Figure 7. The fixed-flat eigenvectors, shown in the left column, correspond to the left halves of the eigenvectors shown on the left of Figure 8, and the eigenvalues match Λ_1 , Λ_3 , and Λ_5 exactly. Similarly, the fixed-fixed eigenvectors on the half-string, shown on the right of Figure 9, equal the left halves of the eigenvectors on the right of Figure 8, and the eigenvalues perfectly match Λ_2 , Λ_4 , and Λ_6 .

We thus have a method for obtaining the data required for the inversion procedure described in the last section, *provided our string has symmetrically arranged beads*.

- 1. Pluck the symmetric string and record the N = 2n eigenvalues $\Lambda_1, \ldots, \Lambda_{2n}$.
- 2. Relabel these eigenvalues for the corresponding half-string: $\lambda_j = \Lambda_{2j}$ and $\widehat{\lambda}_j = \Lambda_{2j-1}$ for j = 1, ..., n.
- 3. Apply the inversion algorithms (Figures 5 and 6) to the half-string eigenvalues to obtain ℓ_0, \ldots, ℓ_n and m_1, \ldots, m_n .
- 4. Recover the parameters for the original symmetric string:

$$L_0 = L_N = \ell_0, \quad \dots, \quad L_{n-1} = L_{n+1} = \ell_{n-1}, \quad L_n = 2\ell_n;$$

 $M_1 = M_N = m_1, \quad \dots, \quad M_n = M_{n+1} = m_n.$

9. Experimental Results for the Inverse Problem. How well does this algorithm perform when applied to real data collected from a string with symmetrically configured beads? Figures 10 and 11 provide data for strings with four and six beads. For each string we collect displacement data for ten seconds at 50,000 samples per second. As described in section 3, the discrete Fourier transform (DFT) of these data reveals peaks that should correspond to $\sqrt{\Lambda_j}$ for $j = 1, \ldots, N$. Deriving from these peaks estimates of the eigenvalues Λ_j , we sort the eigenvalues into fixed-flat and fixed-fixed eigenvalues as described in the last section, then feed them to the inversion algorithms in Figures 5 and 6. The results are illustrated in Figures 10 and 11. We see quite satisfactory agreement for the four bead case, with all quantities recovered to a relative error less than 3.5%. Some challenges begin to emerge with six beads. For one, the frequency plot in Figure 11 reveals a number of secondary peaks. Moreover, the forward experiment illustrated in Figure 4 hints that the finite width of the beads may introduce some uncertainty. For these data, the recovered length L_0 suffers from an 18% relative error; the other lengths and masses are a bit more accurate.

10. Further Explorations. We encourage readers to conduct explorations of both the forward and inverse problem using data sets we provide on the website

http://www.caam.rice.edu/~beads

which includes time series data and peak locations for numerous two-, four-, and six-bead systems.

As you will observe in your own experiments, the inverse procedure we describe is subjected to a variety of errors. To begin with, our measurements introduce imprecision: we measure lengths accurate to $\pm 2 \text{ mm}$, tension to $\pm 10^4 \text{ dyn}$, masses to $\pm 0.01 \text{ g}$, and frequencies to $\pm 0.7 \text{ sec}^{-1}$. Our mathematical model is only an approximation of the true system; we account for neither nonlinear effects nor the damping that causes our string to eventually stop vibrating. The string itself is neither perfectly flexible nor massless, as the model assumes, nor are the beads point-masses. (Figure 4 hints at one possible effect of the bead widths on the eigenvalues.) The model describes asymptotically small vibrations, but the amplitude of our vibrations must be large enough to be measured by the photodetector. Given this litany of errors, you might be surprised at the accuracy achieved in the experiments described in section 9!



Fig. 10 Experimental results illustrating the recovery of positions and masses for a string with four symmetrically-placed beads. The top plot shows the DFT of the displacement data; the four peaks give values for $\sqrt{\Lambda_j}$, shown as gray vertical lines. The middle illustrations compare the positions and masses recovered from these experimental eigenvalues to the "true" positions and masses measured directly, with the corresponding data presented in the table.

There is one further source of error that merits consideration, particularly when we consider applying this experiment to a string with many beads. The numerical algorithm detailed in section 5 will incur rounding errors when implemented in floating point computer arithmetic. Roughly speaking, the algorithm in Figure 5 can exhibit significant errors when n > 50, in MATLAB's default double-precision arithmetic. (One should also scale the units appropriately, as physical values such as $\tau = 10^7$ dyn can lead to overflow as n gets large.) Readers can explore this instability by computing eigenvalues for hypothetical symmetric strings using the **K** and **M** matrices from section 2 (use eig(K,M) in MATLAB), then feeding these "exact" data to the inversion algorithm. How accurately do you recover the lengths and masses that you started with? How does this accuracy depend on the number of beads, as n gets very large? Interested readers can study alternative algorithms such as the continued fractions approach (see the extended manuscript available at [7]) and the Lanczos method (see [6, section 4.2]).



Fig. 11 Repetition of Figure 10, but now for a string with six beads.

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