

## Lecture 9: Examples for the use of Integral Eqs.

### 1 Vibration of an inhomogeneous clamped string[1]

Consider a stretched metallic string clamped between two horizontal points. The distance between the fixed points is  $L$ , and the mass per unit length of the string is not a constant, and hence the speed of propagation of the waves depends on the location along the string. When a disturbance is imparted to the string, the particles in the string vibrate in the vertical direction with a distribution of frequencies that is to be determined.

We denote by  $y(x, t)$  the small displacement of a mass-point on the string in the vertical direction away from the equilibrium position  $y = 0$  at a horizontal distance  $x$  of the point from the left end at time  $t$ . The wave equation is

$$\frac{\partial^2 y}{\partial x^2} - \frac{\rho}{T} \frac{\partial^2 y}{\partial t^2} = 0 \quad (1)$$

where  $\rho(x)$  is the density of the string material and  $T$  is the tension along the string. We define the dimensionless function  $\mathcal{R}(x)$  which describes the variation of the density with  $x$  as

$$\rho(x) = \rho_0 \mathcal{R}(x) \quad (2)$$

where  $\rho_0$  is some fixed value of the density. If we define the reference speed  $c$  by

$$\frac{\rho_0}{T} = \frac{1}{c^2} \quad (3)$$

then Eq. (1) becomes

$$\frac{\partial^2 y}{\partial x^2} - \frac{1}{c^2} \mathcal{R}(x) \frac{\partial^2 y}{\partial t^2} = 0 \quad (4)$$

A separation of variables  $y(x, t) = \psi(x) A(t)$  yields the two equations

$$\frac{d^2 \psi(x)}{dx^2} + \lambda \mathcal{R}(x) \psi(x) = 0 \quad (5)$$

and

$$\frac{d^2 A}{dt^2} = -\lambda c^2 A. \quad (6)$$

If  $\lambda$  is known then the solution of Eq. (6) is

$$A(t) = a \cos(wt) + b \sin(wt); \quad w = c\sqrt{\lambda}. \quad (7)$$

The values of  $\lambda$  are obtained from the solution of Eq. (5). This is a Sturm Liouville eigenvalue equation with an infinite number of solutions determined by

the boundary conditions. The solutions  $\psi_n(x)$  with eigenvalues  $\lambda_n$  form a complete set with  $n = 1, 2, \dots$  in terms of which a general solution of Eq. (4) is given by

$$\psi(x, t) = \sum_{n=1} [a_n \cos(w_n t) + b_n \sin(w_n t)] \psi_n(x) \quad (8)$$

The focus of the present example is how to obtain the eigensolutions of Eq. (5) under the boundary conditions

$$\psi_n(0) = \psi_n(L) = 0. \quad (9)$$

and for  $t = 0$

$$y(x, 0) = f(x), \text{ and } dy(x, 0)/dt = g(x). \quad (10)$$

Here  $f(x)$  describes the "shape" of the string at  $t = 0$ , and  $g(x)$  describes the initial velocity of the mass-pieces of the string. One can visualize this as a hammer striking the string at a certain point, imparting to the string at this point a "dent", and a velocity to the particles in the region of the dent. Once  $f$  and  $g$  are given, then the coefficients in Eq. (8) are determined by the integrals

$$a_n = \int_0^L f(x) \psi_n(x) dx \text{ and } b_n = \frac{1}{w_n} \int_0^L g(x) \psi_n(x) dx \quad (11)$$

### 1.1 Fourier solution of Eq. (9.6)

The fourier functions  $\phi_\ell(x)$  that have the same boundary conditions (9) as  $\psi_n(x)$  are

$$\phi_\ell(x) = \sqrt{\frac{2}{L}} \sin(k_\ell x), \quad k_\ell = \ell\pi/L, \quad \ell = 1, 2, \dots \quad (12)$$

These functions form an orthogonal basis in terms of which the functions  $\psi_n$  can be represented

$$\psi^{(N)}(x) = \sum_{\ell=1}^N d_\ell \phi_\ell(x) \quad (13)$$

The subscript  $n$  has been dropped at this stage, but the truncation parameter  $N$  has been included. Inserting the expansion (13) into Eq. (5), multiplying on the left with  $\phi_{\ell'}(x)$ , and integrating over  $x$  from 0 to  $L$ , one obtains

$$\sum_{\ell=1}^N R_{\ell'\ell} d_\ell = \frac{1}{\lambda} k_{\ell'}^2 d_{\ell'} \quad (14)$$

where

$$\mathbf{R}_{\ell'\ell} = R_{\ell'\ell} = \int_0^L \phi_{\ell'}(x) \mathcal{R}(x) \phi_\ell(x) dx. \quad (15)$$

and where  $\mathbf{R}$  is a symmetric  $N \times N$  matrix.

Eq. (14) can be transformed into a simple eigenvalue equation by defining the diagonal matrix

$$\mathbf{k}^{-1} = \begin{pmatrix} k_1^{-1} & & & & \\ & k_2^{-1} & & & \\ & & k_3^{-1} & & \\ & & & \ddots & \\ & & & & k_N^{-1} \end{pmatrix} \quad (16)$$

in terms of which the matrix  $\mathbf{M}_{fourier}$  can be defined as

$$\mathbf{M}_{fourier} = \mathbf{k}^{-1} \mathbf{R} \mathbf{k}^{-1}. \quad (17)$$

The column vectors  $(u) = (u_1, u_2, ..u_N)^T$  and  $(d) = (d_1, d_2, ..d_N)^T$  related by

$$(u) = \mathbf{k}(d) \quad (18)$$

can also be defined, where  $\mathbf{k}$  is the diagonal matrix of the values  $k_i$ ,  $i = 1, 2, ..N$ . With these transformations the generalized eigenvalue Eq. (14) can be transformed into an ordinary eigenvalue Eq.

$$\mathbf{M}_{Fourier}(u)_n = \frac{1}{\lambda_n}(u_n), \quad n = 1, 2, ..N \quad (19)$$

where  $\mathbf{M}_{Fourier}$  is a symmetric matrix and  $n$  represents each different eigenvalue and eigenfunction. The eigenvalues are displayed in Fig.2 and 4 of the .pdf page 4 of Ref. [1], corresponding to an inhomogeneity

$$\mathcal{R}(x) = 1 + F_0 x^2 \quad (20)$$

with  $F_0 = 2$  or  $4$  in units of inverse length, and  $x$  is the distance along the string. For the present calculation  $L = 1$ . The inhomogeneity (20) corresponds to a 20% or 40% increase in density along the string. Nevertheless, the frequencies are still harmonic to a surprisingly good approximation, as shown in Fig. 4 of Ref. [1].

The main "computational" observation made in this sub-section is that the eigenvalues of a matrix, even though symmetric, loose accuracy towards the high eigenvalues. An iterative way of calculating eigenvalues will be described in the last section, that serve as a benchmark result because of its high accuracy ( $10^{-11}$ ), and is not obtained as an eigenvalue of a matrix.

## 1.2 Eigenvalues in terms of an integral equation.

In the present section we transform Eq. (5) into an integral equation by "inverting" the operator  $d^2/dx^2$  by means of a special energy independent Greens function

$$\frac{1}{\lambda_n} \psi_n = - \int_0^L \mathcal{G}_0(r, r') \mathcal{R}(r') \psi_n(r') dr', \quad (21)$$

with  $n = 1, 2, \dots, N$ , where  $N$  is the number of Chebyshev polynomials used in the expansion of all the functions involved. Here

$$\begin{aligned}\mathcal{G}_0(r, r') &= -F(r)G(r')/L \text{ for } r \leq r' \\ \mathcal{G}_0(r, r') &= -F(r')G(r)/L \text{ for } r' \leq r\end{aligned}\quad (22)$$

where

$$F(r) = r \text{ and } G(r) = L - r, \quad (23)$$

and the factor  $1/L$  is the inverse of the Wronskian between  $F$  and  $G$ . Both  $F$  and  $G$  obey  $d^2/dr^2(F \text{ and } G) = 0$ , they are linearly independent of each other, and provide the appropriate boundary conditions for  $\psi$ . The Greens function above is the limiting result for  $k \rightarrow 0$  of the Greens function  $\mathcal{G}_0$  described by Eq. (6) of Lecture 6. An additional difference to Eq. (6) of Lecture 6 is that Eq. (21) lacks a driving term  $F$ , i.e., Eq. (21) is now an eigenvalue equation.

Similarly to Eq. (13) of Lecture 6, the spectral matrix that represents Eq. (21) is

$$\mathbf{M}_{IEM} = -C^{-1}(1/L)(L/2) [F_D C S_R C^{-1} G_D + G_D C S_L C^{-1} F_D] \mathcal{R}_D C \quad (24)$$

and the corresponding eigenvalue equation is

$$\mathbf{M}_{IEM}(a_n) = -\frac{1}{\lambda_n}(a_n). \quad (25)$$

It is remarkable that the two totally different matrices  $\mathbf{M}_{Fourier}$  and  $\mathbf{M}_{IEM}$  give the same eigenvalues  $\lambda_n$ , albeit with different accuracies. The accuracies are illustrated in Fig. 10 of page 9 of Ref. [1]. The iterative method used to provide an accuracy comparison standard are also described in Ref. [1], but will not be elaborated here.

The basic conclusion of the present lecture is that there are always several different computational methods to solve a problem, and comparison between such methods is desirable in order to gain insight into the accuracy of each of these methods.

## References

- [1] G. Rawitscher and J. Liss, "*The vibrating inhomogeneous string*", Am. J. Phys. 79, 417-429, (2011)