Defect Analysis of 1D Spring-Mass Systems via Laplace Transform and Asymptotics Larry Guan¹; Neil Jerome Egarguin, M.S.¹; Daniel Onofrei, Ph.D.¹ ¹Department of Mathematics, University of Houston, Houston, USA

INTRODUCTION

Spring-mass systems have utility in modeling multiple physical phenomena involving elastic deformation and wave propagation. In the scope of this project: we focus on linear spring-mass systems and use experimental data to analytically locate and enumerate defects along the chain. A particular application of this project concerns limited access pipelines, where we can determine the number of blockages present and their severity.

Consider a chain of n blocks linked in series with springs within two rigid, immobile walls on a flat surface. All springs are massless, identical, and uniformly stiff (k = 1 N/m). Frictional effects dampen each block in the chain uniformly (d = 0.1). All blocks are assigned an integer indicative of their location along the chain (1, 2, ..., n) and are identical in size and shape. Most blocks have a fixed mass of 1 kg, but some of them have unknown mass of non-unity (the "defects"). The locations of each defect are also unknown.

If we apply an impulse to the first block and observe ONLY the vibrations of the first block, is it possible to:

- count the number of defects in the chain?
- recover each defects' masses and locations?





(In the case of one defect: the defective block has mass m_i and is situated at location *j*.)

THEORY

• Establish the system of differential equations describing the motions of each block.

$$\begin{cases} \delta(t) = x_1'' + dx_1' + 2x_1 - x_2 \\ 0 = x_2'' + dx_2' + 2x_2 - x_1 - x_3 \\ \vdots \\ 0 = m_j x_j'' + dx_j' + 2x_j - x_{j-1} - x_{j+1} \\ \vdots \\ 0 = x_{n-1}'' + dx_{n-1}' + 2x_{n-1} - x_{n-2} - x_{n+1} \\ 0 = x_n'' + dx_n' + 2x_n - x_{n-1} \end{cases}$$

Transform the system into the Laplace domain and rewrite as a matrix equation.

$$(A + E)\tilde{x} = \mathbf{b}$$

$$A = \begin{pmatrix} h & 1 & 0 & \cdots & 0 \\ 1 & h & 1 & \ddots & \vdots \\ 0 & \ddots & \ddots & \ddots & 0 \\ \vdots & \ddots & 1 & h & 1 \\ 0 & \cdots & 0 & 1 & h \end{pmatrix} E = \begin{pmatrix} 0 & \cdots & 0 & \cdots & 0 \\ \vdots & \ddots & \vdots & \ddots & \vdots \\ 0 & \cdots & 0 & \cdots & 0 \end{pmatrix} \mathbf{b} = \begin{pmatrix} -1 \\ 0 \\ \vdots \\ 0 \end{pmatrix}$$

$$h = -(s^2 + ds + 2)$$

Explicitly invert A with a method from "Analytical Inversion of Symmetric Tridiagonal Matrices" by Hu and O'Connell. $(A^{-1} = R)$

$$R_{i,j} = \frac{\cosh[\lambda(n+1-i-j)] - \cosh[\lambda(n+1-|j-i|)]}{2\sinh(\lambda)\sinh[\lambda(n+1)]}$$

$$\lambda = \operatorname{arccosh}(-h/2)$$

• Compute an alternate form of $\widetilde{x_1}(s)$ with the matrix equation.

$$\widetilde{x_1}(s) = -R_{1,1} + (1 - m_j)s^2 \frac{|R_{1,j}|^2}{1 + (1 - m_j)s^2R_{j,j}}$$

• This equation is for one defect. When an arbitrary number of defects are present, there are additional additive terms of similar structure representing each defect (one term per defect).

METHOD

Algebraically manipulate the alternate form of $\widetilde{x_1}(s)$ under full expansion such that all terms containing a^{j} remain on one side of the equation.

$$\frac{\mathrm{D}}{s^2}(\widetilde{x_1}(s) + R_{1,1}) = a^{-2j} \frac{(1 - m_j)|r_{1,j}|^2}{\mathrm{D} + (1 - m_j)s^2 r_{j,j}}$$

where:
$$a = \frac{1}{2}(-h + \sqrt{h^2 - 4})$$

$$D = \stackrel{2}{1} + a^{-2n-4} - a^{-2} - a^{-2n-2}$$
$$r_{1,j} = a^{-2} + a^{2j-2n-2} - 1 - a^{2j-2n-4}$$

Insert a new parameter j_{test} and define the left side as a function.

$$g(s, j_{\text{test}}) = \frac{D}{s^2} \left(\widetilde{x_1}(s) + R_{1,1} \right) a^{2j_{\text{test}}}$$
$$g(s, j_{\text{test}}) \approx \frac{1 - m_j}{m_1} a^{2j_{\text{test}} - 2j} \quad \text{(for large values of)}$$

- Systematically determine *j* by plotting $g(s, j_{test})$ for j_{test} from 2 to n. If...
 - $j_{\text{test}} < j: \lim_{s \to \infty} [g(s, j_{\text{test}})] = 0$; Block is not a defect.
 - $j_{\text{test}} > j: \lim_{s \to \infty} [g(s, j_{\text{test}})] = \pm \infty$; Block is not a defect.
 - $j_{\text{test}} = j: \lim_{s \to \infty} [g(s, j_{\text{test}})] = \frac{1 m_j}{m_i}$; Block is a defect.
- m_i can be recovered from the finite limit.

$$m_j \approx \lim_{s \to \infty} \left[\frac{1}{1 + g(s, j_{\text{test}})} \right]$$

- Repeat iteratively. Loop back to the algebraic isolation of the next successive defect's factor each time a defect is discovered, and terminate at the end of the chain. The iteration count of this algorithm gives the number of defective blocks.
- The lack of access to a laboratory setting compelled numerical synthesis of the vibrational data for $\widetilde{x_1}(s)$ from the matrix equation $(A + E)\widetilde{x} = \mathbf{b}$. Such a system is impossible to solve exactly, so $\widetilde{x_1}(s)$ is numerically
 - approximate (though, to a very high degree of precision).
- The alternate form of $\widetilde{x_1}(s)$, being analytically exact, acted as smoothed noise-free data and was considered "perfect."

RESULTS

One Defect

• Data for $\widetilde{x_1}(s)$ was synthesized with the following parameters: • $n = 10; m_i = 2; j = 4$

Plots of g using the synthetic data of $\widetilde{x_1}(s)$:





- Algorithm detected one defect of correct mass at the right location. Convergence to zero for $j_{\text{test}} < j$ and divergence for $j_{\text{test}} > j$.
- As *s* grows large, numerical instabilities overtake smoothness of g.
- Extremely small numerical imprecisions in the simulation of $\widetilde{x_1}(s)$ (~10⁻¹⁵⁰) due to software limitations are blown up as *s* grows.
 - Analogous to measurement errors and cannot be remedied.

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RESULTS (cont.)





Plots of g using the "perfect" data of $\widetilde{x_1}(s)$:

 $g(s, j_{test})$ for $j_{test} \leq j$



- Expected behavior is observed.
- g is smooth everywhere.
- Noise-free data is impossible to obtain in reality.

Two Defects

$$\frac{\widetilde{x_1}(s) + R_{1,1}}{s^2} a^{2j_{\text{test}}} \approx \left(\frac{1 - m_{j_1}}{m_{j_1}} a^{-2j_1} + \frac{1 - m_{j_2}}{m_{j_2}} a^{-2j_2}\right) a^2$$

- Data for $\widetilde{x_1}(s)$ was synthesized with the following parameters:
 - $n = 20; m_{j_1} = 3; j_1 = 5; m_{j_2} = 8; j_2 = 16$
- Algorithm found the first defect with correct mass at the right location.
- Defect #2 is irrecoverable for the following reasons:
 - Multiplicative coupling of m_{j_1} and m_{j_2} . The alternate form of $\widetilde{x_1}(s)$ requires knowledge of m_{j_2} when trialing for j_2 .
 - Exacerbated numerical imprecision. Finite convergence is never observed if m_{i_2} is approximated out during the second iteration. Exponentiation rules. When trialing for j_2 , no such value of j_{test} can
 - force convergence to a finite value.

CONCLUSION & FUTURE WORK

- Though theoretically sound, this algorithm managed to detect only the first defect in the chain. Furthermore, the numerical instabilities in plotting $g(s, j_{test})$ irreconcilably conflicted with the expectation of smooth, gradual convergence to a finite value due to the imprecision of $\widetilde{x_1}(s)$ – fundamentally similar to measurement noise. Using exact data removed the instability, but that is impossible to procure in reality.
- In the future: we'll revisit the problem in the time domain, where the first block vibration curve can be found by rewriting the system of differential equations into a first order system and solving with matrix exponentials.

$$\mathbf{u}' = \mathbf{f} + A\mathbf{u}$$

$$A = \begin{pmatrix} 0 & 1 & 0 & \cdots & \cdots & \cdots \\ -2k & -d & k & \ddots & \ddots & \ddots \\ 0 & 0 & 0 & 1 & \ddots & \ddots \\ k & 0 & -2k & -d & k & \ddots \\ \ddots & \ddots & \ddots & \ddots & \ddots & \ddots \end{pmatrix} \quad \mathbf{f} = \begin{pmatrix} 0 \\ \delta(t) \\ 0 \\ \vdots \\ 0 \end{pmatrix} \quad \mathbf{u} =$$

REFERENCES

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u_2 u_{2n}

2j_{test}