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# **Developing an Open-Source Frequency Domain Modal Analysis Algorithm in Python**

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# 1 Introduction

## 1.1 Modal Analysis

In *linear* structural dynamics, *Modal Analysis* is universally regarded as the pre-eminent solution for the identification or characterization of a structure or system. Modal analysis provides a means to perform said identification by studying the structure's *modal properties or parameters*. Whether using mathematical modelling or experimental testing, modal analysis is a pillar for an engineer's understanding of a structure's response to various forms of excitation. This understanding is crucial for safety and complying with standards or for research areas such as structural health monitoring or system identification.

Fundamentally, modal analysis is the decomposition of the complex oscillatory behaviours of structures and systems into several smaller components called *modes*. Every mode is a numerical or mathematical representation of a specific vibration pattern corresponding to a *natural frequency*, the frequency, or set of, at which a system tends to oscillate at subject to an initial displacement or velocity, and an associated *mode shape*, a vector or function which describes the relative movement of the various degrees-of-freedom (DOFs) or axis points for each natural frequency, and *damping ratio*, a unitless parameter describing the energy dissipation properties of the system at the natural frequencies. The modal properties are defined by the interaction of the system's physical properties, its mass, stiffness, and damping. These inherent properties guide the system's vibration when subject to external forcing or initial displacements or velocities.

Based on the theory, the standard procedure for modal analysis begins with forming the equations of motion (EOMs) that represent a system. Typically, the EOMs are second order matrix differential equations. Considering a system with an arbitrary number  $N$ , degrees-of-freedom as shown in figure 1.1. Interpreting Newton's second law as:

$$\sum F_i = m_i \ddot{\mathbf{x}}_i$$

Where  $F_i$  is a force acting on the  $i$ -th DOF,  $m_i$  is the mass, and  $\mathbf{x}_i$  is the coordinate of said DOF. Or the Euler-Lagrange equation as such:

$$\frac{d}{dt} \left( \frac{\partial T}{\partial \dot{q}_i} \right) - \frac{\partial T}{\partial q_i} + \frac{\partial U}{\partial q_i} = Q_i$$

Where  $q_i$  is a generalized coordinate, for this system one can take  $q_i = \mathbf{x}_i$ ,  $T$  is the system's kinetic energy,  $U$  is the potential energy, and  $Q_i$  are the non-conservative forces which impart or dissipate energy from the system. One finds that the equations of motion for the system are defined as shown in Equation 1.1

$$\begin{aligned} m_1 \ddot{x}_1 + (c_1 + c_2) \dot{x}_1 - c_2 \dot{x}_2 + (k_1 + k_2) x_1 - k_2 x_2 &= F_1 \\ m_2 \ddot{x}_2 + (c_2 + c_3) \dot{x}_1 - c_2 \dot{x}_1 - c_3 \dot{x}_3 + (k_2 + k_3) x_2 - k_2 x_1 - k_3 x_3 &= F_2 \\ &\dots \\ m_N \ddot{x}_N + (c_N + c_{N+1}) \dot{x}_N - c_{N-1} \dot{x}_{N-1} + (k_N + k_{N+1}) x_N - k_{N-1} x_{N-1} &= F_N \end{aligned} \tag{1.1}$$

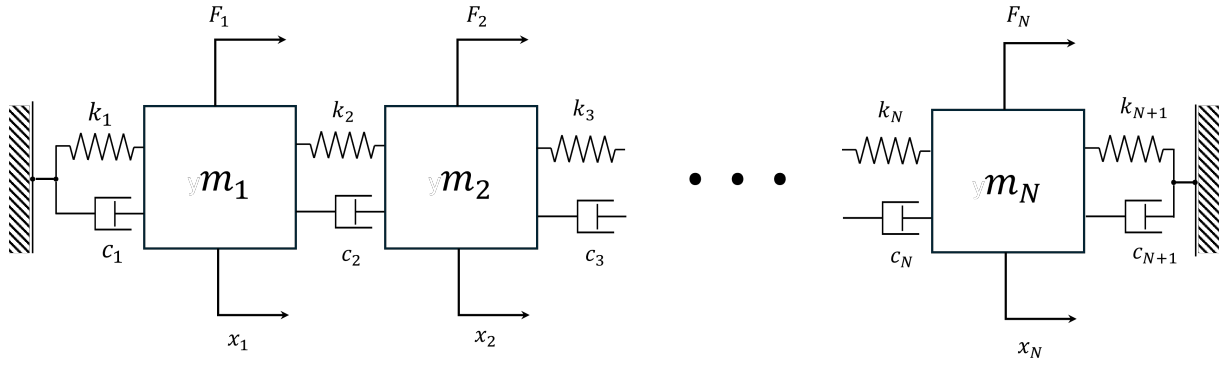


Figure 1.1: A "lumped-mass" system of  $N$  degrees-of-freedom

Assembling the physical parameters into respective matrices as highlighted in Equation 1.2, the characteristic differential equation of the system is found.

$$\begin{aligned}
 [\mathbf{K}] &= \begin{bmatrix} k_1 + k_2 & -k_2 & 0 & \dots & 0 \\ -k_2 & k_2 + k_3 & -k_3 & \dots & 0 \\ 0 & -k_3 & k_3 + k_4 & -k_4 & 0 \\ \dots & \dots & \dots & \ddots & \dots \\ 0 & 0 & 0 & -k_N & k_N + k_{N+1} \end{bmatrix} \\
 [\mathbf{C}] &= \begin{bmatrix} c_1 + c_2 & -c_2 & 0 & \dots & 0 \\ -c_2 & c_2 + c_3 & -c_3 & \dots & 0 \\ 0 & -c_3 & c_3 + c_4 & -c_4 & 0 \\ \dots & \dots & \dots & \ddots & \dots \\ 0 & 0 & 0 & -c_N & c_N + c_{N+1} \end{bmatrix} \\
 [\mathbf{M}] &= \text{diag}(m_i) \quad \forall i = 1, 2, \dots, N
 \end{aligned} \tag{1.2}$$

$$\therefore [\mathbf{M}]\ddot{\mathbf{x}} + [\mathbf{C}]\dot{\mathbf{x}} + [\mathbf{K}]\mathbf{x} = \mathbf{F} \tag{1.3}$$

For the sake of simplicity, assume that the vibrating system presented is left to vibrate freely, and is undamped, meaning  $\mathbf{F}$  and  $\mathbf{C}$  are zero. If the solution of the presented differential equation is harmonic such that  $\mathbf{x}(t) = \Psi e^{j\omega t}$ , one finds that equation 1.3 reduces to:

$$([\mathbf{K}] - \omega^2[\mathbf{M}])\Psi = 0 \tag{1.4}$$

By a rearrangement of the variables, one can manipulate the equation into becoming the general form for the eigenvalue problem:

$$\begin{aligned}
 [\mathbf{A}] &= [\mathbf{M}]^{-1}[\mathbf{K}] \\
 [\mathbf{A}]\Psi &= \omega^2\Psi
 \end{aligned} \tag{1.5}$$

Where  $\mathbf{A}$  is the so-called dynamical matrix,  $\omega^2$  is the eigenvalue — often referred to as an *eigenfrequency* — of the matrix, and  $\Psi$  is the corresponding eigenvector representing the modeshape vector.

In systems with a finite number of degrees of freedom—often referred to as discrete systems—the modal analysis is performed using the dynamical matrix  $[\mathbf{A}]$ , which encapsulates the system's inertial and elastic properties. Building upon the discrete case, in systems that are modelled as continuous functions, the physical parameters of the system are distributed spatially. This results in a partial differential equation that governs the motion, where finite-dimensional matrices are replaced by linear operators. Consequently, an eigenvalue problem is formed in

terms of a differential operator, with eigenvalues corresponding to the natural frequencies and the associated eigenfunctions representing the mode shapes in the spatial domain<sup>1</sup>. This principle, applied to transversal vibration of an Euler-Bernoulli beam results in the eigenvalue problem defined in equation 1.6:

$$\Delta^2 Y(x) = \beta^4 Y(x) \quad (1.6)$$

Where  $\Delta$  is the Laplacian operator (second partial derivative in Cartesian coordinate system), and  $\beta^4 = \frac{\omega^2}{c^2}$ , in which  $\omega$  is the natural frequency, and  $c^2$  is the ratio of the flexural rigidity of the beam, and its inertia.

In both cases, discrete and continuous, by mathematically manipulating the EOMs governing the system, differential equations are reduced to eigenvalue problems—a solution that is not only mathematically elegant and robust, as it is applicable to all linear systems, but also linguistically apt, given that "eigen" means "one's own". However, while this approach to modal analysis may be elegant, its practical applications rarely exist. Real world vibrating structures and systems rarely conform to the idealized assumptions of lumped-mass systems or Euler-Bernoulli beams with known boundary conditions, this limitation can be attributed to more complex geometries, or internal interactions between the components of a system. This limitation has incentivized academics and practitioners to approach modal analysis differently.

---

<sup>1</sup>An *eigenvector*, or an *eigenfunction*, of a matrix or linear operator defined on some vector/function space is any non-zero vector/function in said space that when multiplied or acted upon by the matrix/linear operator is equivalent to being multiplied by some scalar factor, said scalar factor is referred to as the *eigenvalue*.

Report Story:  
**Structural Dynamics in engineering**

- How studying structural mechanics has helped us have safer, lighter, greener structures.
- Modal Analysis, is regarded as **the** solution for linear structural dynamics.
- Maths of modal analysis:
  - EOM formulation.
  - Eigendecomposition of state matrices.
  - FRF in modal terms. (refer to plscf solution using that)
- This is in various industries, home appliances, aero, civil/structural, acoustics etc.
- Experimental Modal Analysis king.
- Curve-fitting methods for modal analysis.
- The need for algorithms in practice.

Onto software usage in engineering contexts

- engineers consistently rely on software tools, this is great as it streamlines the important processes.
- Important aspects of software dev. :
  - logic/control flow
  - complexities and big O notation
  - unit testing and integration testing.
  - version control.
  - foss vs. prop. discuss why it's cheaper in long run, and better for everyone involved.
- Revisit the aims and objectives.

## **2 Algorithms / development:**

### **maths-y bit**

- lsce (brief)
- lscf (brief)
- plscf (important bits and refer to appendix for full derivation, use own notation and wording).
- Why the companion matrix solution works.  $\text{eig}(C(p)) = \lambda_i \rightarrow p(\lambda) = 0$
- here one must explain what poles are for a system.
- lsfd for modeshapes.

## software-y bit

- time complexity optimization. discuss that O notation doesn't always equal less time.
- unit testing, explain pytest stuff and fixturing and bla bla bla.
- using numpy (BLAS routines/subroutines)
- user facing code.
- formatting guidelines.

## 3 Results

- plscf on simulated modal data.
  - clean
  - noisy
  - slightly nonlinear data (mimo where  $H_{ij} \approx H_{ji}$  but  $H_{ij} \neq H_{ji}$ )
  - high dofs.
- plscf on actual lab data.
- adam's plscf vs siemens lms polymax on same dataset.
- interpretation of stabilization diagram.

```
1 import numpy as np
2
3 def _make_polynomial_basis_fcn(
4     polynomial_order: int, frequency_vector: np.ndarray,
5     sampling_frequency: float
6 ):
7     """Function that creates a Polynomial Basis Function matrix
8
9     Args:
10         polynomial_order (int): Order of the polynomial created.
11             i.e. if 2 then  $P(x) = a_0 * x^0 + a_1 * x^1 + a_2 * x^2$ 
12             in the case of pLSCF,  $\Omega(w) = P(e^{\{jw\_delta\}t})$ 
13
14         frequency_vector (np.ndarray): vector of frequencies
15             measured or simulated, can be hz or rads-1.
16             MUST be either a row or column vector/1D Array
17
18     Returns:
19         Polynomial basis function matrix.
20     """
21     # the polynomial basis function is actually the vandermonde
22     # matrix for the polynomials, A and B.
23     dt = 1 / sampling_frequency # Sampling rate
```

```
20     s = np.exp(1.j*frequency_vector*dt) # this is the "x" in the
      polynomial
21     return np.vander(x=s,N=polynomial_order+1,increasing=True)
```



Consider the Jacobian matrix, for a system with  $n_{outputs} = 3$ :

$$J = \begin{pmatrix} X_1 & 0 & 0 & Y_1 \\ 0 & X_2 & 0 & Y_2 \\ 0 & 0 & X_3 & Y_3 \end{pmatrix} \quad (3.1)$$

$$J^H J = \begin{pmatrix} X_1^H X_1 & 0 & 0 & X_1^H Y_1 \\ 0 & X_2^H X_2 & 0 & X_1^H Y_2 \\ 0 & 0 & X_2^H X_2 & X_1^H Y_3 \\ X_1^* Y_1^T & X_2^* Y_2^T & X_3^* Y_3^T & Y_1^H Y_1 + Y_2^H Y_2 + Y_3^H Y_3 \end{pmatrix} \quad (3.2)$$

if:

$$R_o = Re(X_o^H X_o)$$

$$S_o = Re(X_o^H Y_o)$$

$$T_o = Re(Y_o^H Y_o)$$

$$2Re(J^H J)\theta = 2Re \begin{pmatrix} X_1^H X_1 & 0 & 0 & X_1^H Y_1 \\ 0 & X_2^H X_2 & 0 & X_1^H Y_2 \\ 0 & 0 & X_2^H X_2 & X_1^H Y_3 \\ X_1^* Y_1^T & X_2^* Y_2^T & X_3^* Y_3^T & Y_1^H Y_1 + Y_2^H Y_2 + Y_3^H Y_3 \end{pmatrix} \begin{pmatrix} \beta_1 \\ \beta_2 \\ \beta_3 \\ \alpha \end{pmatrix} \quad (3.3)$$

$$= 2 \begin{pmatrix} R_1 \beta_1 + S_1 \alpha \\ R_2 \beta_2 + S_2 \alpha \\ R_3 \beta_3 + S_3 \alpha \\ S_1^T \beta_1 + S_2^T \beta_2 + S_3^T \beta_3 + (T_1 + T_2 + T_3) \alpha \end{pmatrix} \quad (3.4)$$

which corresponds to the solutions in the normal equations in terms of alpha and beta.

$$l^{LS}(\theta) = tr\{\theta^T Re(J^H J)\theta\} \quad (3.5)$$

from vector calculus, if

$$\mathbf{f} = \mathbf{x}^T \mathbf{B} \mathbf{x} \quad (3.6)$$

then,

$$\frac{\partial \mathbf{f}}{\partial \mathbf{x}} = 2\mathbf{B}\mathbf{x} \quad (3.7)$$

then the normal equation for polymax should be:

$$\frac{\partial l^{LS}}{\partial \theta} = 2Re(J^H J)\theta \quad (3.8)$$

## A Derivation of p-LSCF modal parameter estimation method

### A.1 The right-matrix rational fractional model

The polyreference least-squares complex frequency domain method employs a right matrix fractional model to fit MIMO Frequency Response Function measurements into a set of rational polynomial transfer functions:

$$[H(\omega)] = [N(\omega)][D(\omega)]^{-1} \quad (\text{A.1})$$

Such that  $H(\omega) \in \mathbb{C}^{N_{\text{outputs}} \times N_{\text{inputs}}}$  is the FRF matrix, where  $D(\omega) \in \mathbb{C}^{N_{\text{inputs}} \times N_{\text{inputs}}}$ , is the denominator matrix polynomial, and  $N(\omega) \in \mathbb{C}^{N_{\text{outputs}} \times N_{\text{inputs}}}$ , is the numerator matrix polynomial. The rows corresponding to each output  $o$  in the FRF matrix can be represented as such:

$$\langle H_o(\omega) \rangle = \langle N_o(\omega) \rangle [D(\omega)]^{-1} \quad (\text{A.2})$$

The row vector numerator polynomial for the  $o^{th}$  output, and the denominator matrix polynomial are defined in terms of a polynomial basis function,  $\Omega(\omega)$ , and their respective polynomial coefficients,  $\beta$  and  $\alpha$  as such:

$$\langle N_o(\omega) \rangle = \sum_{r=1}^p \Omega_r(\omega) \langle \beta_{or}(\omega) \rangle \quad (\text{A.3})$$

$$[D(\omega)] = \sum_{r=1}^p \Omega_r(\omega) [\alpha_r] \quad (\text{A.4})$$

With the polynomial basis function  $\Omega_r(\omega) = e^{j\omega\Delta t r}$ . Although not initially obvious as polynomials with conventional form  $p(x) = c_0 + c_1x + c_2x^2 + \dots + c_nx^n$ , the basis functions are expressed in the  $s$ -domain where  $s = e^{j\omega\Delta t}$ . The polynomial coefficients,  $\alpha_r \in \mathbb{R}^{N_{\text{inputs}} \times N_{\text{inputs}}}$  and  $\beta_{or} \in \mathbb{R}^{1 \times N_{\text{inputs}}}$ , are assembled into matrix form:

$$\beta_o = \begin{pmatrix} \beta_{o0} \\ \beta_{o1} \\ \beta_{o2} \\ \dots \\ \beta_{op} \end{pmatrix} \in \mathbb{R}^{(p+1) \times N_{\text{inputs}}} \quad (\text{A.5})$$

$$\alpha = \begin{pmatrix} \alpha_0 \\ \alpha_1 \\ \alpha_2 \\ \dots \\ \alpha_p \end{pmatrix} \in \mathbb{R}^{N_{\text{inputs}} * (p+1) \times N_{\text{inputs}}} \quad (\text{A.6})$$

$$\theta = \begin{pmatrix} \beta_0 \\ \beta_1 \\ \beta_2 \\ \dots \\ \beta_{N_o} \\ \alpha \end{pmatrix} \in \mathbb{R}^{(N_{\text{outputs}} + N_{\text{inputs}})(p+1) \times N_{\text{inputs}}} \quad (\text{A.7})$$

## A.2 Minimizing the sum of the squared residuals

The collection of both sets of coefficients into one variable  $\theta$ , makes performing the least squares problem simpler, in a sense, as it becomes the one unknown in this least squares model. As typical in any fitting method, one must minimize the error between the model and the real or measured value. The nonlinear least-squares error for:

Measured FRF:  $\hat{H}_o(\omega_k)$

Model FRF:  $H_o(\omega_k)$

is weighted such that:

$$\epsilon_o^{NLS}(\theta, \omega_k) = w_o(\omega_k)(H_o(\omega_k) - \hat{H}_o(\omega_k)) \quad (\text{A.8})$$

Where  $\epsilon_o^{NLS} \in \mathbb{C}^{1 \times N_{inputs}}$ ,  $w_o(\omega_k)$  is a scalar weighing function which captures the variation and deviation between multiple inputs on the same measurement point, and  $\forall k = 0, 1, 2, \dots, N_{frequency}$ . Said weighing function is typically denoted by

$$w_o(\omega_k) = \frac{1}{\sqrt{\text{var}[H_o(\omega_k)]}} \quad (\text{A.9})$$

(See [reference for weighted linear regressions] for more information on weighted least squares.) One can then define the nonlinear cost function as the sum of the error "squared", (hermitian inner product), over the data points, in this case, spectral lines and outputs;

$$l^{NLS}(\theta) = \sum_{o=1}^{N_{out}} \sum_{k=1}^{N_f} \text{tr}\{(\epsilon_o^{NLS}(\theta, \omega_k))^H \epsilon_o^{NLS}(\theta, \omega_k)\} \quad (\text{A.10})$$

In this equation,  $\text{tr}\{\bullet\}$  denotes the trace of a matrix, also known as the sum of diagonal elements, and  $\bullet^H$  denotes the Hermitian (conjugate) transpose. The trace operator is used as the trace of a product of 2 matrices  $\mathbf{A} \in \mathbb{C}^{m \times n}$  and  $\mathbf{B} \in \mathbb{C}^{n \times m}$  will equal the sum of each individual element in  $\mathbf{A}$  with the individual elements of  $\mathbf{B}$ . This provides a sum of all square residuals/errors in the cost function.

$$\text{tr}\{\mathbf{A}^H \mathbf{B}\} = \text{tr}\{\mathbf{A} \mathbf{B}^H\} = \text{tr}\{\mathbf{B}^H \mathbf{A}\} = \text{tr}\{\mathbf{B} \mathbf{A}^H\} = \sum_{i=1}^m \sum_{j=1}^n a_{ij} b_{ij} \quad (\text{A.11})$$

## A.3 Linearizing the error

One can then obtain the polynomial coefficients through minimizing the cost function in A.10, by setting the derivative  $\frac{\partial l^{NLS}}{\partial \theta}$  equal to zero, however a nonlinear cost function will yield nonlinear derivative equations, (typically called normal equations in linear regression). A subsequent linearization of the cost function can approximate (suboptimally) the least squares problem, this is achieved through right multiplying the cost function with denominator polynomial  $\mathbf{D}$ . This gives a linear error:

$$\begin{aligned} \epsilon_o^{LS}(\omega_k, \theta) &= w_o(\omega_k)(N_o(\omega_k, \beta_o) - \hat{H}_o(\omega_k)D(\omega_k, \alpha)) \\ &= w_o(\omega_k) \sum_{r=0}^p (\Omega_r(\omega_k)\beta_{or} - \Omega_r(\omega_k)\hat{H}_o(\omega_k)\alpha_r) \end{aligned} \quad (\text{A.12})$$

Stacking the error in terms for all spectral lines in one matrix  $E_o^{LS}(\theta) \in \mathbb{C}^{N_f \times N_{in}}$ :

$$E_o^{LS}(\theta) = \begin{pmatrix} \epsilon_o^{LS}(\omega_1, \theta) \\ \epsilon_o^{LS}(\omega_2, \theta) \\ \epsilon_o^{LS}(\omega_3, \theta) \\ \vdots \\ \epsilon_o^{LS}(\omega_{N_f}, \theta) \end{pmatrix} = \begin{pmatrix} X_o & Y_o \end{pmatrix} \begin{pmatrix} \beta_o \\ \alpha \end{pmatrix} \quad (\text{A.13})$$

Here, new variables  $\mathbf{X}$  and  $\mathbf{Y}$  are introduced:

$$X_o = \begin{pmatrix} w_o(\omega_1) \left( \Omega_0(\omega_1) + \Omega_1(\omega_1) \dots \Omega_p(\omega_1) \right) \\ w_o(\omega_2) \left( \Omega_0(\omega_2) + \Omega_1(\omega_2) \dots \Omega_p(\omega_2) \right) \\ \vdots \\ w_o(\omega_{N_f}) \left( \Omega_0(\omega_{N_f}) + \Omega_1(\omega_{N_f}) \dots \Omega_p(\omega_{N_f}) \right) \end{pmatrix} \in \mathbb{C}^{N_f \times (p+1)} \quad (\text{A.14})$$

$$Y_o = \begin{pmatrix} -w_o(\omega_1) \left( \Omega_0(\omega_1) + \Omega_1(\omega_1) \dots \Omega_p(\omega_1) \right) \otimes \hat{H}_o(\omega_1) \\ -w_o(\omega_2) \left( \Omega_0(\omega_2) + \Omega_1(\omega_2) \dots \Omega_p(\omega_2) \right) \otimes \hat{H}_o(\omega_2) \\ \vdots \\ -w_o(\omega_{N_f}) \left( \Omega_0(\omega_{N_f}) + \Omega_1(\omega_{N_f}) \dots \Omega_p(\omega_{N_f}) \right) \otimes \hat{H}_o(\omega_{N_f}) \end{pmatrix} \in \mathbb{C}^{N_f \times N_{in}(p+1)} \quad (\text{A.15})$$

Where  $\otimes$  is the Kronecker product. In these equations,  $\mathbf{X}$  is used to capture the frequency content of the least squares problem, and  $\mathbf{Y}$  is used to capture both the frequency content and the measured response data. Using these matrices, one can reconstruct the nonlinear cost function into one that is linear:

$$\begin{aligned} l^{LS}(\theta) &= \sum_{o=1}^{N_{out}} \sum_{k=1}^{N_f} \text{tr} \{ (\epsilon_o^{LS}(\omega_k, \theta))^H \epsilon_o^{LS}(\omega_k, \theta) \} \\ &= \sum_{o=1}^{N_{out}} \text{tr} \left\{ (E_o^{LS}(\theta))^H E_o^{LS}(\theta) \right\} \\ &= \sum_{o=1}^{N_{out}} \text{tr} \left\{ \begin{pmatrix} \beta_o^T & \alpha^T \end{pmatrix} \begin{pmatrix} X_o^H \\ Y_o^H \end{pmatrix} \begin{pmatrix} X_o & Y_o \end{pmatrix} \begin{pmatrix} \beta_o \\ \alpha \end{pmatrix} \right\} \end{aligned} \quad (\text{A.16})$$

If one defines a *Jacobian* matrix  $\mathbf{J} \in \mathbb{C}^{N_f N_{out} \times (N_{in} + N_{out})(p+1)}$  for the problem as such:

$$\mathbf{J} = \begin{pmatrix} X_1 & 0 & \dots & 0 & Y_1 \\ 0 & X_2 & \dots & 0 & Y_2 \\ \dots & \dots & \dots & \dots & \dots \\ 0 & 0 & 0 & X_{N_{out}} & Y_{N_{out}} \end{pmatrix} \quad (\text{A.17})$$

The cost function can be represented as:

$$l^{LS}(\theta) = \text{tr} \{ \theta^T \mathbf{J}^H \mathbf{J} \theta \} \quad (\text{A.18})$$

To obtain real values of  $\theta$ , one must place a constraint on the cost function such that:

$$l^{LS}(\theta) = \text{tr} \{ \theta^T \text{Re}(\mathbf{J}^H \mathbf{J}) \theta \} \quad (\text{A.19})$$

Where the Gramian matrix of  $\mathbf{J}$  can be represented in terms a set of variables,  $\mathbf{R}$ ,  $\mathbf{S}$  and  $\mathbf{T}$ :

$$\text{Re}(\mathbf{J}^H \mathbf{J}) = \begin{pmatrix} R_1 & 0 & \dots & 0 & S_1 \\ 0 & 0 & \dots & 0 & S_2 \\ \dots & \dots & \dots & \dots & \dots \\ 0 & 0 & \dots & R_{N_{out}} & S_{N_{out}} \\ S_1^T & S_2^T & \dots & S_{N_{out}}^T & \sum_{o=1}^{N_{out}} T_o \end{pmatrix} \quad (\text{A.20})$$

in which:

$$R_o = Re(X_o^H X_o) \quad (\text{A.21})$$

$$S_o = Re(X_o^H Y_o) \quad (\text{A.22})$$

$$T_o = Re(Y_o^H Y_o) \quad (\text{A.23})$$

#### A.4 The Normal Equations, and extracting the modal parameters

The cost function can then be minimized in terms of  $\alpha$  and  $\beta$  to find the best least-squares fit:

$$\frac{\partial l^{LS}(\theta)}{\partial \beta_o} = 2(R_o \beta_o + S_o \alpha) = 0 \quad (\text{A.24})$$

$$\forall O = 1, 2, \dots, N_{out}$$

$$\frac{\partial l^{LS}(\theta)}{\partial \alpha} = 2 \sum_{o=1}^{N_{out}} (S_o^T \beta_o + T_o \alpha) \quad (\text{A.25})$$

Giving normal equations of this least squares problem in terms of the wanted polynomial coefficients, one can also assemble those normal equations into 1 equation:

$$\frac{\partial l^{LS}(\theta)}{\partial \theta} = 2Re(\mathbf{J}^H \mathbf{J})\theta = 0 \quad (\text{A.26})$$

The denominator coefficients  $\alpha$  are used to obtain the poles and the modal participation factors, which are sufficient information for the constructing a stabilization diagram. Hence, one can further reduce the normal equations by setting:

$$\beta_o = R_o^{-1} S_o \alpha \quad (\text{A.27})$$

This yields the reduced normal equation:

$$\left\{ 2 \sum_{o=1}^{N_{out}} (T_o - S_o^T R_o^{-1} S_o) \right\} \alpha = 0 \quad (\text{A.28})$$

$$\mathbf{M} \alpha = 0$$

For a non-trivial solution to the normal equation, a constraint is set on  $\alpha$ , where:

$$\alpha_p = \mathbf{I}_{N_{in}} \quad (\text{A.29})$$

The rest of the denominator coefficients are then found using:

$$\mathbf{M}(1 : N_{in} * p, 1 : N_{in} * p) \begin{pmatrix} \alpha_0 \\ \alpha_1 \\ \alpha_2 \\ \dots \\ \alpha_{p-1} \end{pmatrix} = \mathbf{M}(1 : N_{in} * p, N_{in} * p + 1 : N_{in} * (p + 1)) \quad (\text{A.30})$$

The least-squares estimate for  $\alpha$  is then:

$$\hat{\alpha}_{LS} = \begin{Bmatrix} \alpha \\ \mathbf{I}_{N_{in}} \end{Bmatrix}$$

This makes the denominator polynomial  $\mathbf{D}$  a *monic* polynomial. Based on the fundamental definition of system poles, which is the points at which the system's response is "infinite", one can say that for an arbitrary rational polynomial  $p(x)$ :

$$p(x) = \frac{a_0 + a_1x + a_2x^2 + \dots + a_nx^n}{b_0 + b_1x + b_2x^2 + \dots + b_nx^n} \quad (\text{A.31})$$

To achieve said "infinite" value, one must find values of  $x$  that represent the roots of the denominator polynomial. In the pLSCF model, one can exploit the monic property of the denominator polynomial. Frobenius companion matrices are square matrices that represent monic polynomials, given one has a monic polynomial

$$p(x) = x^n + a_{n-1}x^{n-1} + a_{n-2}x^{n-2} + \dots + a_1x + a_0$$

The companion matrix of said polynomial is defined as:

$$C(p) = \begin{bmatrix} 0 & 0 & \dots & 0 & -a_0 \\ 1 & 0 & \dots & 0 & -a_1 \\ 0 & 1 & \dots & 0 & -a_2 \\ \vdots & \vdots & \ddots & \vdots & \vdots \\ 0 & 0 & \dots & 1 & -a_{n-1} \end{bmatrix} \quad (\text{A.32})$$

A property of the companion matrix  $C(p) \in \mathbb{R}^{n \times n}$  is that its eigenvalues  $\lambda_1, \lambda_2, \dots, \lambda_n$  are the roots of  $p(x)$ , where  $p(\lambda_i) = 0, \forall i = 1, 2, \dots, n$ . This property aids in finding the system poles, after constructing the companion matrix for the denominator polynomial, an Eigendecomposition of the matrix yields the discrete time poles as the eigenvalues, and the corresponding eigenvectors are the modal participation factors:

$$C(\mathbf{D}) = \begin{pmatrix} 0 & I & \dots & 0 & 0 \\ 0 & 0 & I & \dots & 0 \\ \dots & \dots & \dots & \dots & \dots \\ 0 & 0 & \dots & 0 & I \\ -\alpha_0^T & -\alpha_1^T & -\alpha_2^T & \dots - \alpha_{p-2}^T & -\alpha_{p-1}^T \end{pmatrix} \quad (\text{A.33})$$

$$C(\mathbf{D})[\mathbf{L}] = \Lambda[\mathbf{L}] \quad (\text{A.34})$$

Where the matrix  $[\mathbf{L}]$  is the Eigenmatrix (matrix with columns as eigenvectors), representing the modal participation factors of each mode, and the matrix  $\Lambda$  contains the discrete time poles on its diagonal elements. The transpose of a companion matrix, this however does not affect the numerical value of the poles or participation factors [reference proof]. A  $p$ -order right matrix-fraction polynomial estimation should yield  $pN_{in}$  number of poles.

## A.5 Finding the modeshapes using the Least-Squares Frequency Domain (LSFD) method

This part should be in actual report: When fitting theoretical models to experimental data, the difficulty does not typically lie in the mathematical framework enabling the modelling process. Instead, the challenge is in constructing a model that provides physically meaningful insights. Given that the goal of many modal parameter estimation methods is to find a set of poles which As apparent, it is straightforward to find the polynomial coefficients using measured data.