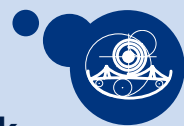


Frequency-weighted damping via nonsmooth optimization and fast computation of QEPs with low-rank updates



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DUBROVNIK]

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Introduction and motivation

We consider damped linear vibration system

$$M\ddot{q}(t) + C(v)\dot{q}(t) + Kq(t) = f(t),$$
$$q(0) = q_0, \quad \text{and} \quad \dot{q}(0) = \dot{q}_0.$$

Where $M, C(v), K \in \mathbb{R}^{n \times n}$ system matrices, $v \in \mathbb{R}^s$ parameter vector.

- M, K are positive definite Hermitian matrices
- $C(v) = C_{int} + C_{ext}(v)$, $C_{int} > 0$ internal damping, $C_{ext}(v) \geq 0$ external damping.
- $C_{ext}(v) = \sum_{i=1}^s v_i g_i g_i^T$
- $C_{int} = \alpha_c C_{crit}$, where $C_{crit} = 2M^{1/2} \sqrt{M^{-1/2} K M^{-1/2}} M^{1/2}$.



Linearization

- Let Φ simultaneously diagonalize pair M and K

$$\Phi^T K \Phi = \Omega^2 = \text{diag}(\omega_1^2, \dots, \omega_n^2) \quad \text{and} \quad \Phi^T M \Phi = I.$$

Note that internal damping is s.t. $\Phi^T C_{int} \Phi = \alpha \Omega$.

With $q = \Phi q_\Phi$ and $y_1 = \Omega q_\Phi$, $y_2 = \dot{q}_\Phi$ we have

$$\frac{d}{dt} \begin{bmatrix} y_1 \\ y_2 \end{bmatrix} = \underbrace{\begin{bmatrix} 0 & \Omega \\ -\Omega & -\Phi^T C(v) \Phi \end{bmatrix}}_{A(v)} \begin{bmatrix} y_1 \\ y_2 \end{bmatrix}.$$

We obtain first order differential equation:

$$\dot{y} = Ay, \quad \text{with solution} \quad y = e^{At} y_0, \quad \text{where } y_0 \text{ is initial data.}$$



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We consider parameter dependant QEP:

$$(\lambda^2(v)M + \lambda(v)C(v) + K) x(v) = 0.$$

With $w_1(v) = \Omega\Phi^{-1}x(v)$ and $w_2(v) = \lambda(v)\Omega^{-1}w_1(v)$ we have

$$\lambda(v) \begin{bmatrix} w_1(v) \\ w_2(v) \end{bmatrix} = \underbrace{\begin{bmatrix} 0 & \Omega \\ -\Omega & -\Phi^T C(v) \Phi \end{bmatrix}}_{A(v)} \begin{bmatrix} w_1(v) \\ w_2(v) \end{bmatrix}.$$



Very important question arises in considering such systems:

For the given mass (M) and stiffness (K) determine the best (optimal) damping which will insure optimal evanescence.

(M, K) , A not stable

(M, K, C) , A stable



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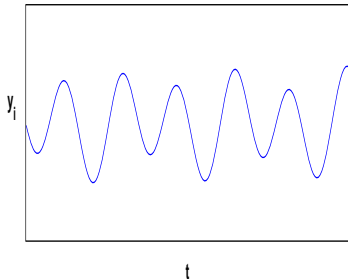
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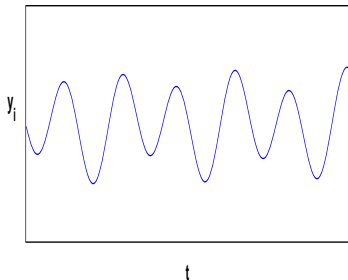
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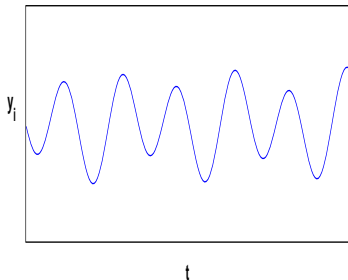
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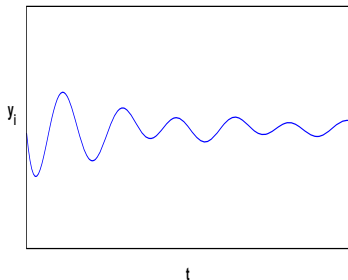
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Very important question arises in considering such systems:

How to re-design a given damped mechanical system, such that a new system does not have eigenvalues in some "dangerous" part of a complex plane, typically called the resonance band.



Optimization criteria

- Minimization of spectral abscissa

$$\alpha_{MCK}(v) \rightarrow \min_v,$$

where $\alpha_{MCK}(v) = \max_k \operatorname{Re} \lambda_k(v)$ and $\lambda_k(v)$ are the eigenvalues of

$$(\lambda^2(v)M + \lambda(v)C(v) + K) x(v) = 0.$$



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Optimization criteria-idea

The undesirable frequency band: $[\omega - b, \omega + b]$, for some given $b > 0$, where $\omega \in \mathbb{R}$ is an undesirable frequency.

$$\begin{aligned} \min_{v \in \mathbb{R}^s} \quad & \max\{\operatorname{Re} \lambda(v) : \lambda(v) \in \Lambda(v) \text{ and } \operatorname{Im} \lambda(v) \in [\omega - b, \omega + b]\} \\ \text{s.t.} \quad & \alpha_{\text{MCK}}(v) \leq \text{tol}_{\text{sa}} \text{ for some } \text{tol}_{\text{sa}} < 0, \\ & v_j \geq 0 \text{ for } j = 1, \dots, s \end{aligned}$$



Using ellipse instead of band

Let $E = (a, b, c)$ denote the axis-aligned ellipse

$$\frac{(x - \operatorname{Re} c)^2}{a^2} + \frac{(y - \operatorname{Im} c)^2}{b^2} = 1,$$

where $a, b > 0$ respectively denote the semi-major and -minor axes and $c \in \mathbb{C}$ is the center of the ellipse. Identifying \mathbb{R}^2 with \mathbb{C} , consider the following algebraic distance $d : \mathbb{C} \mapsto [0, \infty)$ of a point $z \in \mathbb{C}$ to this ellipse, i.e.,

$$d(z; E) := \frac{(\operatorname{Re}(z - c))^2}{a^2} + \frac{(\operatorname{Im}(z - c))^2}{b^2}.$$



Using ellipse instead of band

Measure of the distance of the spectrum to the undesirable frequency $[\omega - b, \omega + b]$, we define

$$d_{\Lambda, E}(v) := \min\{d(\lambda(v); E) : \lambda(v) \in \Lambda(v)\},$$

where $E = (a, b, \mathbf{i}\omega)$. Similarly,

$$d_{\Lambda, \mathcal{E}}(v) := \min\{d_{\Lambda, E_j}(v) : E_j \in \mathcal{E}\},$$

where $E_j := (a_j, b_j, \mathbf{i}\omega_j)$ is defining the j th ellipse for the j th undesirable frequency band $[\omega_j - b_j, \omega_j + b_j]$ with relative importance $a_j > 0$ and $\mathcal{E} := \{E_1, \dots, E_k\}$ is the set of k corresponding ellipses.



New optimization criteria

- Frequency isolation while minimizing spectral abscissa (FI1)

FI1:
$$\begin{aligned} \min_{v \in \mathbb{R}^s} \quad & \alpha_{\text{MCK}}(v) \\ \text{s.t.} \quad & d_{\Lambda, \mathcal{E}}(v) \geq 1, \\ & \alpha_{\text{MCK}}(v) \leq \text{tol}_{\text{sa}} \text{ for some } \text{tol}_{\text{sa}} < 0, \\ & v_j \geq 0 \text{ for } j = 1, \dots, s. \end{aligned}$$



New optimization criteria

- Frequency isolation while maximizing the major axis of the ellipses (FI2)

$$\mathbf{FI2:} \quad \max_{v \in \mathbb{R}^s} \sum_{j=1}^k \phi_j a_{\Lambda, E_j}(v)$$

$$\begin{aligned} \text{s.t.} \quad & \alpha_{\text{MCK}}(v) \leq \text{tol}_{\text{sa}} \text{ for some } \text{tol}_{\text{sa}} < 0, \\ & v_j \geq 0 \text{ for } j = 1, \dots, s. \end{aligned}$$

where $a_{\Lambda, E}(v) := \min\{a(\lambda(v); E) : \lambda(v) \in \Lambda(v)\}$,

$$a(z; E) := \begin{cases} \frac{b|\operatorname{Re} z|}{\sqrt{b^2 - (\operatorname{Im} z - \omega)^2}}, & \text{if } \operatorname{Im} z \in (\omega - b, \omega + b), \\ \infty & \text{otherwise.} \end{cases}$$



GRANSO: GRAdient-based Algorithm for Non-Smooth Optimization

Requires gradients:

$$\begin{aligned}\frac{\partial \lambda(v)}{\partial v_j} \Big|_{v=\hat{v}} &= -\frac{\hat{x}^* (\lambda(\hat{v}) g_j g_j^T) \hat{x}}{\hat{x}^* (2\lambda(\hat{v}) M + C(\hat{v})) \hat{x}}, \\ \frac{\partial \alpha_{\text{MCK}}(v)}{\partial v_j} \Big|_{v=\hat{v}} &= -\text{Re} \frac{\partial \lambda(v)}{\partial v_j} \Big|_{v=\hat{v}}, \\ d'(z(t); E) &= 2 \left(\frac{\text{Re}(z(t) - c) \cdot \text{Re} z'(t)}{a^2} + \frac{\text{Im}(z(t) - c) \cdot \text{Im} z'(t)}{b^2} \right), \\ a'(z(t); E) &= \frac{b \text{sgn}(\text{Re } z(t)) \cdot \text{Re } z'(t)}{(b^2 - (\text{Im } z(t) - \omega)^2)^{1/2}} + \frac{b |\text{Re } z(t)| (\text{Im } z(t) - \omega) \cdot \text{Im } z'(t)}{(b^2 - (\text{Im } z(t) - \omega)^2)^{3/2}}.\end{aligned}$$



Apply s times efficient algorithm for computing eigenvalues of (CSymDPR1) matrix $A = D + \rho uu^T$.

The eigenvalues of A are the zeros of the secular function:

$$f(\lambda) = 1 + \rho \sum_{i=1}^{2n} \frac{u_i^2}{d_i - \lambda} = 1 + \rho u^T (D - \lambda I)^{-1} u,$$

and the corresponding eigenvectors are given by

$$w_i = \frac{y_i}{\|y_i\|_2}, \text{ where } y_i = (D - \lambda_i I)^{-1} u, \quad i = 1, \dots, 2n.$$



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initial vector $y_0 \neq 0$

Rayleigh quotient iteration

$$\lambda = 0$$

until convergence

1. $y^1 := (A - \lambda I)^{-1}y^0$
2. $\mu = \frac{y^{1T}y^0}{y^{0T}y^0}$
3. $\lambda = \lambda + \frac{1}{\mu}$
4. $y^0 = y^1$

Modified Rayleigh quotient iteration

$$\lambda = 0$$

until convergence

1. $\mu = \eta \frac{y^{0T}(A - \lambda I)y^0}{y^{0T}y^0}$
2. $\lambda = \lambda + \mu$
3. $y^0 := (D - \lambda I)^{-1}u$

compute eigenvectors

N. Jakovčević Stor, I. Slapničar, and Z. Tomljanović. *Fast computation of optimal damping parameters for linear vibrational systems*. arXiv e-prints, 2020



$$C_{ext}(v) = \sum_{i=1}^s v_i g_i g_i^T = G \text{diag}(v) G^T$$

How to get from $Ay = \lambda y$, where

$$A = \begin{bmatrix} 0 & \Omega \\ -\Omega & -\alpha\Omega - \Phi^T G \text{diag}(v) G^T \Phi \end{bmatrix},$$

to multiple CSymDPR1 eigenvalue problem?



$$A = \begin{bmatrix} 0 & \Omega \\ -\Omega & -\alpha\Omega \end{bmatrix} + \begin{bmatrix} 0 & 0 \\ 0 & -\Phi^T G \text{diag}(v) G^T \Phi \end{bmatrix}$$





$$P^T \cdot \setminus A = \begin{bmatrix} 0 & \Omega \\ -\Omega & -\alpha\Omega \end{bmatrix} + \begin{bmatrix} 0 & 0 \\ 0 & -\Phi^T G \text{diag}(v) G^T \Phi \end{bmatrix} / \cdot P$$





$$P^T A P = \left(\begin{bmatrix} D_1 & & \\ & \ddots & \\ & & D_n \end{bmatrix} - \hat{G} \begin{bmatrix} v_1 & & \\ & \ddots & \\ & & v_s \end{bmatrix} \hat{G} \right), \text{ where}$$

$$D_i = \begin{bmatrix} 0 & \omega_i \\ -\omega_i & -\alpha\omega_i \end{bmatrix} \text{ and } \hat{G} = P^T \begin{bmatrix} 0 \\ \Phi^T G \end{bmatrix}.$$



$$\Psi^{-1} \backslash P^T A P = \left(\begin{bmatrix} D_1 & & \\ & \ddots & \\ & & D_n \end{bmatrix} - \hat{G} \begin{bmatrix} v_1 & & \\ & \ddots & \\ & & v_s \end{bmatrix} \hat{G} \right), \quad / \cdot \Psi$$

where $\Psi = \begin{bmatrix} \Psi_1 & & \\ & \ddots & \\ & & \Psi_n \end{bmatrix}$



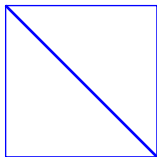


$$\tilde{A} = D - U \begin{bmatrix} v_1 & & \\ & \ddots & \\ & & v_s \end{bmatrix} Z^T = D - \sum_{j=1}^s v_j u_j z_j^T,$$

$$D = \Psi^{-1} \begin{bmatrix} D_1 & & \\ & \ddots & \\ & & D_n \end{bmatrix} \Psi,$$

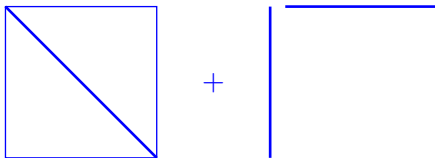
$$U = \Psi^{-1} \hat{G}, \quad \text{and} \quad Z = \Psi^T \hat{G},$$





+ ... +







Algorithm 1 Frequency-weighted damping optimization algorithm

Input: M and K , $\alpha \geq 0$ for C_{int} and G , set of k ellipses \mathcal{E} , weights $[\phi_1, \dots, \phi_k]$ with each $\phi_j \in (0, 1]$ for ellipse $E_j \in \mathcal{E}$, $\text{tol}_{\text{sa}} < 0$, initial viscosity $v_{\text{init}} \in \mathbb{R}_+^s$, and $\text{approach} \in \{1, 2\}$.

Output: Computed for optimized viscosities $v_{\text{opt}} \in \mathbb{R}_+^s$ for either FI1 or FI2

- 1: $[\Phi, \Omega]$ matrices from linearization
 - 2: $[\Psi, D, U, Z]$ matrices that construct low-rank structure
 - 3: **if** $\text{approach} = 1$ **then**
 - 4: $v_{\text{opt}} \leftarrow$ solution returned by GRANSO for FI1 initialized at v_{init}
 - 5: **else**
 - 6: $v_{\text{opt}} \leftarrow$ solution returned by GRANSO for FI2 initialized at v_{init}
 - 7: **end if**
-

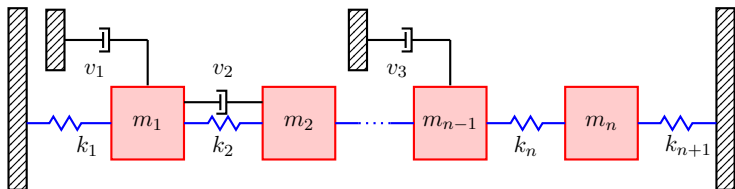


Figure: n -mass oscillator

$$\begin{aligned}
 n &= 200 \cdot i, \quad i = 1, \dots, 10, \\
 M &= \text{diag}(m_1, m_2, \dots, m_n), \\
 m_i &= 10 + \frac{990}{n-1} \cdot (i-1), \quad i = 1, \dots, n
 \end{aligned}$$

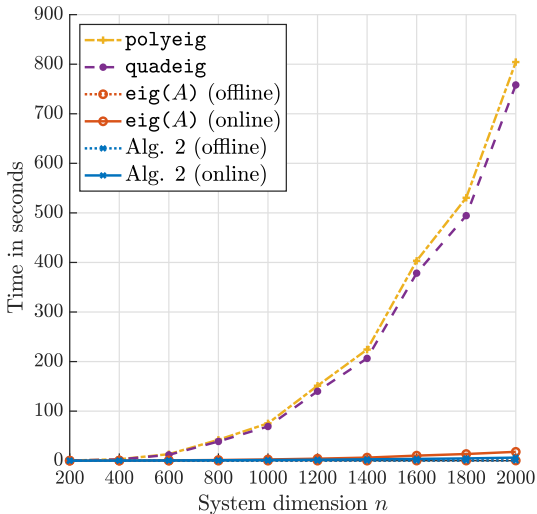


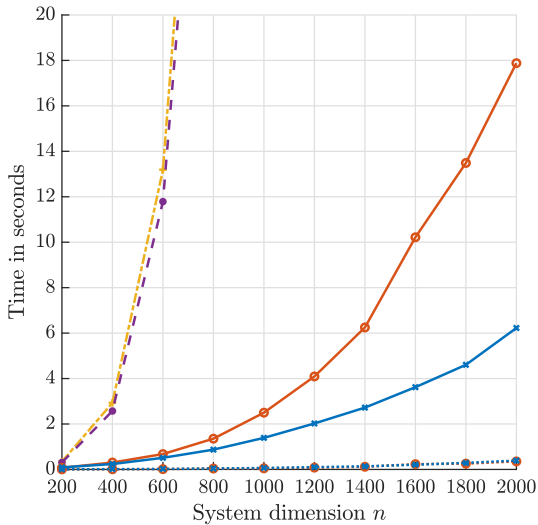
Example

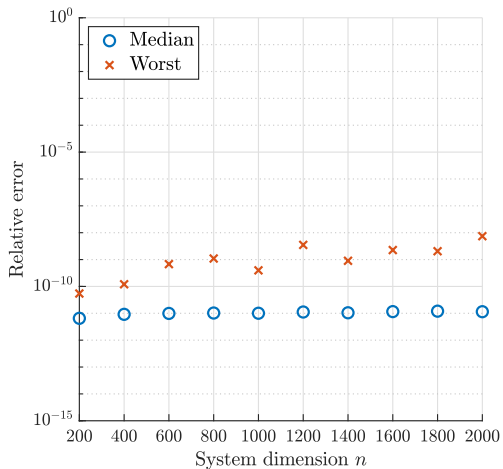
$$K = \begin{bmatrix} k_1+k_2 & -k_2 & & & \\ -k_2 & \ddots & \ddots & & \\ & \ddots & \ddots & -k_n & \\ & & -k_n & k_n+k_{n+1} & \end{bmatrix}, \quad k_i = 5, i = 1, \dots, n+1$$

$$C_{\text{ext}} = v_1 e_k e_k^T + v_2 (e_j - e_{j+1})(e_j - e_{j+1})^T + v_3 e_l e_l^T, \quad \alpha = 0.004$$
$$v = [v_1, v_2, v_3]^T, \quad v_1, v_2, v_3 \in [0.1, 1.1]$$

$$(k, j, l) = \left(\frac{n}{10}, \frac{3n}{10}, \frac{5n}{10} \right)$$









Example

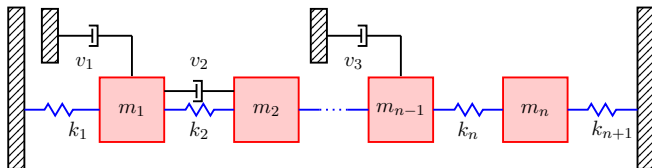


Figure: n -mass oscillator

$$n = 1000,$$

$$m_i = m_{n+1-i} = \frac{2n-i}{200} \quad \text{for } i = 1, \dots, \frac{n}{2},$$

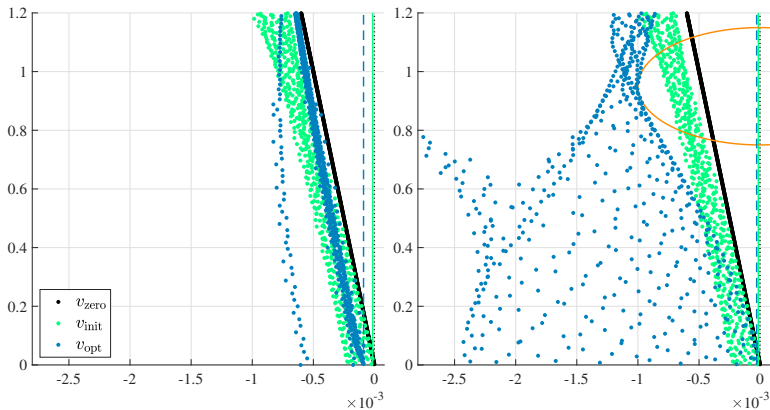
$$k_i = 5 \quad \text{for } i = 1, \dots, n+1,$$

$$(j, k, l) = (100, 400, 900),$$

$$v_{\text{init}} = \text{ones}(3, 1).$$



$$E = (0.001, 0.2, 0.95i)$$



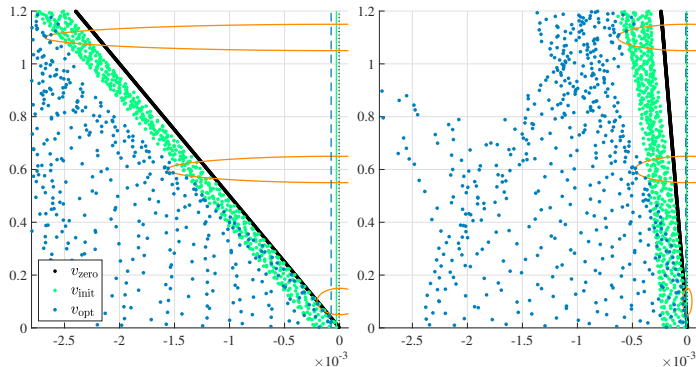
(a) Spectral abscissa minimization only

$$v_{\text{opt}} = \begin{bmatrix} 314.7 \\ 94.1 \\ 124.5 \end{bmatrix}$$

(b) Approach 1 (Fixed ellipses)

$$v_{\text{opt}} = \begin{bmatrix} 8.032 \\ 8.184 \\ 9.970 \end{bmatrix}$$

$$E_1 = (\sim, 0.05, 0.1i), \phi_1 = 1, \quad E_2 = (\sim, 0.05, 0.6i), \phi_2 = 0.2, \quad E_3 = (\sim, 0.05, 1.1i), \phi_3 = 0.1$$



(c) $\alpha = 0.004,$

$$v_{\text{opt}} = \begin{bmatrix} 8.137 \\ 7.147 \\ 1.789 \end{bmatrix}$$

(d) $\alpha = 0.0004,$

$$v_{\text{opt}} = \begin{bmatrix} 8.294 \\ 7.767 \\ 1.673 \end{bmatrix}$$



Introduction and motivation
Solving optimization problem
Numerical experiments



Example - eigenvalue approximation
Example - optimization criteria



Thank you for your attention!