

# Computing dominant poles of large second-order transfer functions

Joost Rommes\*<sup>1</sup> and Nelson Martins\*\*<sup>2</sup>

<sup>1</sup> NXP Semiconductors, Corp. I&T / DTF, High Tech Campus 37, Box WY4-01, 5656 AE, Eindhoven, The Netherlands

<sup>2</sup> CEPEL, Rio de Janeiro, Brazil

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## 1 Introduction

A transfer function of a large dynamical system often only has a small number of dominant poles compared to the number of state variables. Modal approximation techniques [1, 2, 3, 4] capture the dominant behavior of the system by projecting the state-space on the modes corresponding to the dominant poles. The computation of the dominant poles and modes requires specialized eigenvalue methods. In [5, 6, 7], algorithms were developed for the computation of dominant poles of single-input single-output (SISO) and multi-input multi-output (MIMO) transfer functions of large scale dynamical systems.

In this paper an efficient algorithm, the Quadratic Dominant Pole Algorithm (QDPA), for the computation of dominant poles of second-order transfer functions is described. Modal equivalents that are constructed by projecting the state-space matrices on the dominant left and right eigenspaces, preserve the structure of the original system. The dominant poles and modes can also be used to improve reduced-order models computed by rational Krylov based methods. For more details on the Quadratic Dominant Pole Algorithm, see [8, 9].

## 2 Second-order dynamical systems, transfer functions, and dominant poles

In this paper, the second-order dynamical systems  $(M, C, K, \mathbf{b}, \mathbf{c}, d)$  are of the form

$$\begin{cases} M\ddot{\mathbf{x}}(t) + C\dot{\mathbf{x}}(t) + K\mathbf{x}(t) &= \mathbf{b}u(t) \\ y(t) &= \mathbf{c}^*\mathbf{x}(t) + du(t), \end{cases} \quad (1)$$

where  $M, C, K \in \mathbb{R}^{n \times n}$  are the system matrices,  $\mathbf{b}, \mathbf{c}, \mathbf{x}(t) \in \mathbb{R}^n$ ,  $u(t), y(t), d \in \mathbb{R}$ . The vectors  $\mathbf{b}$  and  $\mathbf{c}$  are called the input and output vector, respectively. The transfer function  $H: \mathbb{C} \rightarrow \mathbb{C}$  of (1) is defined as  $H(s) = \mathbf{c}^*(s^2M + sC + K)^{-1}\mathbf{b} + d$ . The poles of  $H(s)$  are a subset of the eigenvalues  $\lambda_i \in \mathbb{C}$  of the quadratic eigenvalue problem (QEP)

$$(\lambda_i^2 M + \lambda_i C + K)\mathbf{x}_i = 0, \quad \mathbf{y}_i^*(\lambda_i^2 M + \lambda_i C + K) = 0, \quad \mathbf{x}_i \neq 0, \quad \mathbf{y}_i \neq 0, \quad (i = 1, \dots, 2n). \quad (2)$$

An eigentriplet  $(\lambda_i, \mathbf{x}_i, \mathbf{y}_i)$  is composed of an eigenvalue  $\lambda_i$  and corresponding right and left eigenvectors  $\mathbf{x}_i, \mathbf{y}_i \in \mathbb{C}^n$ .

By transforming [10, Section 3.5] QEP (2) and system (1) to linear equivalents, the partial fraction representation becomes  $H(s) = \mathbf{c}^*X(sI - \Lambda)^{-1}\Lambda Y^*\mathbf{b} + \sum_{i=1}^{2n} \frac{R_i}{s - \lambda_i}$ , where  $X = [\mathbf{x}_1, \dots, \mathbf{x}_{2n}]$ ,  $Y = [\mathbf{y}_1, \dots, \mathbf{y}_{2n}]$ , and  $R_i = (\mathbf{c}^*\mathbf{x}_i)(\mathbf{y}_i^*\mathbf{b})\lambda_i$ . The terms  $R_i$  are called the residues, and  $\mathbf{x}_i$  and  $\mathbf{y}_i$  are scaled so that  $-\mathbf{y}_i^*K\mathbf{x}_i + \lambda_i^2\mathbf{y}_i^*M\mathbf{x}_i = 1$ .

A pole  $\lambda_i$  of  $H(s)$  with corresponding right and left eigenvectors  $\mathbf{x}_i$  and  $\mathbf{y}_i$  is called the *dominant* pole if  $\widehat{R}_i = \frac{|R_i|}{\text{Re}(\lambda_i)} > \widehat{R}_j$ , for all  $j \neq i$ . More generally, a pole  $\lambda_i$  is called dominant if  $|\widehat{R}_i|$  is not very small compared to  $|\widehat{R}_j|$ , for all  $j \neq i$ . This can also be seen in the corresponding Bode-plot, which is a plot of  $|H(i\omega)|$  against  $\omega \in \mathbb{R}$ : peaks occur at frequencies  $\omega$  close to the imaginary parts of the dominant poles of  $H(s)$ .

## 3 Quadratic Dominant Pole Algorithm

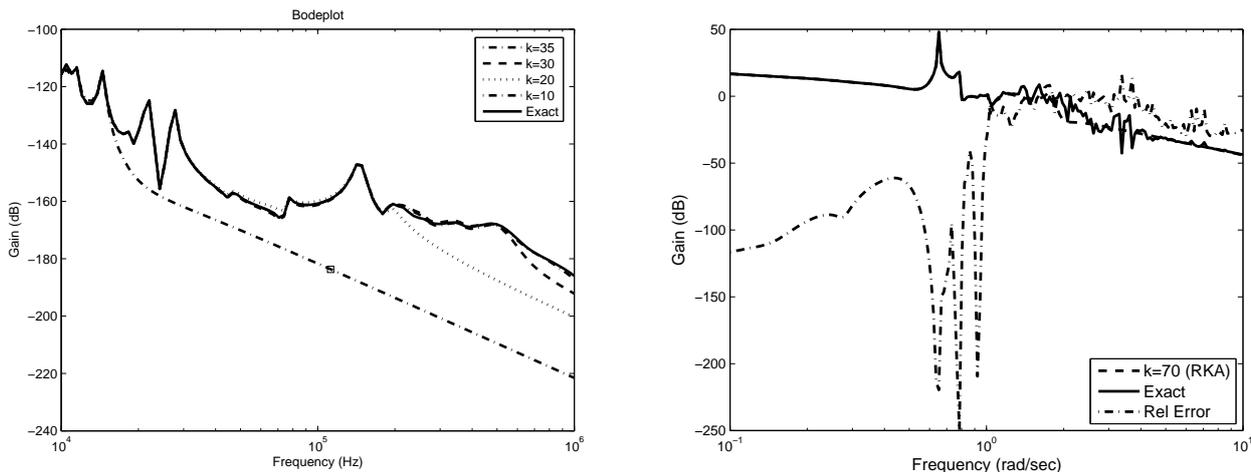
The poles of  $H(s)$  are the  $\lambda \in \mathbb{C}$  for which  $\lim_{s \rightarrow \lambda} |H(s)| = \infty$  and hence  $\lim_{s \rightarrow \lambda} 1/H(s) = 0$ . In other words, the poles are the roots of  $1/H(s)$  and a good candidate to find these roots is Newton's method: noting that  $H'(s) = -\mathbf{c}^*(s^2M + sC + K)^{-1}(2sM + C)(s^2M + sC + K)^{-1}\mathbf{b}$  and starting with initial pole estimate  $s_0$  gives the following scheme:

$$s_{k+1} = s_k + \frac{1}{H(s_k)} \frac{H^2(s_k)}{H'(s_k)} = s_k - \frac{\mathbf{c}^*\mathbf{v}}{\mathbf{w}^*(2s_k M + C)\mathbf{v}},$$

where  $\mathbf{v} = (s_k^2 M + s_k C + K)^{-1}\mathbf{b}$  and  $\mathbf{w} = (s_k^2 M + s_k C + K)^{-*}\mathbf{c}$ .

\* Corresponding author: e-mail: joost.rommes@nxp.com, Phone: +00 31 (0)40 27 49798, Fax: +00 31 (0)40 27 46276

\*\* e-mail: nelson@cepel.br



**Fig. 1** Exact and reduced system transfer functions for the gyro (left) and vibrating body (right).

QDPA can be extended with subspace acceleration (keep search spaces for right and left eigenvectors) and a selection strategy (select the most dominant approximation every iteration) to improve global convergence to the most dominant poles, and deflation to avoid recomputation of already found poles. See [8] for more details.

#### 4 Numerical results and conclusions

The left figure in Fig. 1 shows frequency response Bode plots of reduced order models based on 10, 20, 30, and 35 poles and corresponding eigenvectors, for a micro-mechanical gyro with  $n = 17361$  degrees of freedom [11]. As can be seen from the matching of the peaks, QDPA finds the dominant poles.

The figure at the right in Fig. 1 shows the frequency response of a 70th order Second-Order Arnoldi [12] reduced model of vibrating body from sound radiation analysis ( $n = 17611$  degrees of freedom), that was computed using the complex part  $i\beta$  of dominant poles  $\lambda = \alpha + i\beta$  (computed by QDPA) as interpolation points. This model is more accurate than reduced order models based on standard Krylov methods and matches the peaks up to  $\omega = 1$  rad/s, because of use of shifts near the resonance frequencies.

The Quadratic Dominant Pole Algorithm (QDPA) is an efficient and effective method for the computation of dominant poles of second-order transfer functions. The dominant poles and corresponding left and right eigenvectors can be used to construct structure-preserving modal equivalents and to determine interpolation points for rational Krylov based model order reduction methods. QDPA can be generalized to MIMO systems and higher-order systems, and can be used for the computation of dominant zeros as well. For more details and results, see [8].

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