

A novel method to compute all eigenvalues of the polynomial eigenvalue problems in an open half plane

Xiaoping Chen^{1,2} · Xi Yang³ · Wei Wei³ · Jinde Cao⁴ · Shuai Tang²

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Abstract

In this paper, we introduce the linear fractional mapping and the contour integral method. Based on them, we develop a new numerical method to find all eigenvalues of the polynomial eigenvalue problems in an open half plane. Numerical examples are shown to illustrate the effectiveness of the proposed method.

Keywords Polynomial eigenvalue problem \cdot Linear fractional mapping \cdot Open half plane \cdot Contour integral method

Mathematics Subject Classification 15A18 · 65F15

1 Introduction

In this paper, we consider the polynomial eigenvalue problem (PEP)

$$P(\lambda)x = 0,\tag{1}$$

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 Xiaoping Chen xpchen@nuaa.edu.cn
 Xi Yang yangxi@nuaa.edu.cn; yangxi@lsec.cc.ac.cn
 Wei Wei shixueying26@163.com
 Jinde Cao jdcao@seu.edu.cn
 Shuai Tang tangshuaixs@163.com
 Department of Mathematics, Southeast University, Nanjing 210096, China
 Department of Mathematics, Taizhou University, Taizhou 225300, China
 Department of Mathematics, Nanjing University of Aeronautics and Astron

- ³ Department of Mathematics, Nanjing University of Aeronautics and Astronautics, Nanjing 210016, China
- ⁴ School of Mathematics and Research Center for Complex Systems and Network Sciences, Southeast University, Nanjing 210096, China



where $P(\lambda) = \lambda^d A_0 + \lambda^{d-1} A_1 + \dots + \lambda A_{d-1} + A_d$, $A_k \in C^{n \times n}$, $0 \le k \le m$. λ and x, known as the eigenvalue and the corresponding eigenvector, respectively, are the variables to be determined. The problem is very general and includes the standard eigenvalue problem $P(\lambda) = \lambda I - A$ (see, e.g., Wilkinson 1965), the generalized eigenvalue problem $P(\lambda) = \lambda A - B$ (see, e.g., Dai et al. 2015), the quadratic eigenvalue problem $P(\lambda) = \lambda^2 A + \lambda B + C$ (see, e.g., Tisseur and Meerbergen 2001) and the cubic eigenvalue problem $P(\lambda) = \lambda^3 A_0 + \lambda^2 A_1 + \lambda A_2 + A_3$ (see, e.g., Hwang et al. 2005).

The PEP arises in a number of various applications, and has received considerable attention in the last few decades. For example, it is ubiquitous in a wide range of problems, such as vibration analysis of viscoelastic systems (Adhikari and Pascual 2009), structural dynamic analysis (Gupta 1976), stability analysis of control systems (Higham and Tisseur 2002), numerical simulation of quantum dots (Hwang et al. 2004) and so on. Considerable efforts have been devoted to the polynomial eigenvalue problem in the literature. Gohberg et al. (1982) developed the mathematical theory concerning matrix polynomials. Gohberg et al. (1979), Dedieu and Tisseur (2003), Higham and Tisseur (2003) and Chu (2003) gave the perturbation theory for the polynomial eigenvalue problem. Higham et al. (2007), Tisseur (2000) and Lawrence and Corless (2015) analyzed backward error of the polynomial eigenvalue problem.

The classical approach for solving the PEP is linearizing the problem (1) to produce an equivalent larger generalized eigenvalue problem (see, e.g., Gohberg et al. 1982; Higham et al. 2006; Mackey et al. 2006a, b), solved by any appropriate eigensolver. However, the linearization technique will enlarge the size of the original problem, and matrix structures and spectral properties of the problem (1) will not be preserved, and most of all, the linearized eigenvalue problem is more ill-conditioned (Tisseur 2000). In the last few years, lots of researchers have developed numerical methods that work directly with the original data of the polynomial eigenvalue problem for avoiding the above disadvantages. In these methods, the polynomial eigenvalue problem is projected onto a properly chosen low-dimensional subspace to obtain a new polynomial eigenvalue problem with matrix dimension of low order. Then, the reduced polynomial eigenvalue problem can be solved by the linearization method. These methods include Jacobi–Davidson method (Bai and Miao 2017a, b; Hwang et al. 2010; Sleijpen et al. 1996) and Krylov-type subspace method (Hoffnung et al. 2006). Jacobi-Davidson method targets at one eigenvalue at one time with local convergence. However, if the desired eigenvalues of the problem (1) form a cluster of nearby eigenvalues, then Jacobi-Davidson method has difficulties in detecting and resolving such a cluster.

Actually, recent focus in practical engineering, such as stability analysis of control systems, is how to compute all eigenvalues of the PEP in an open half plane, especially in the open right half plane. Besides, it is well known that a real symmetric or Hermitian matrix polynomial has a spectrum that is symmetric with respect to the real axis. Therefore, for the above matrix polynomials, we just need compute all eigenvalues in the open upper half plane instead of all eigenvalues of original matrix polynomial in the whole plane. Overall, a numerical method for computing all eigenvalues of the PEP (1) in an open half plane would be of particular interest.

The organization of the paper is as follows. In Sect. 2, we give a brief description of the linear fractional mapping. Through this mapping, the problem to find all eigenvalues of a PEP in the open half plane can be converted into a new PEP to find all eigenvalues in the open unit disk. Besides, we present the relation of the eigenvalues for the above eigenproblems. In Sect. 3, we introduce the contour integral method to compute all eigenvalues in the open unit disk. In Sect. 4, a novel algorithm to find all eigenvalues of the PEP in an open half plane is derived. Section 5 is devoted to some numerical experiments.

For convenience, we use the following notations: *I* denotes the identity matrix with suitable dimensions; tr(*A*) and det(*A*) denote the trace and the determinant of a matrix *A*, respectively; Re(λ) and Im(λ) denote the real part and imaginary part of a complex number λ , respectively; *i* denotes the imaginary unit.

2 Linear fractional transformation

To compute all eigenvalues of the PEP (1) in the open half plane, we first introduce the linear fractional transformation (Rudin 1987).

Definition 1 If a_1, a_2, a_3 and a_4 are complex numbers and $a_1a_4 - a_2a_3 \neq 0$ such that $\varphi(\lambda)$: $\lambda \to \omega$ where $\omega = \frac{a_1\lambda + a_2}{a_3\lambda + a_4}$, the mapping $\varphi(\lambda)$ is called a linear fractional transformation.

For the above transformation, every open half plane can be conformally mapped onto an open disk.

Next, we take the linear fractional mapping $\omega = \frac{\lambda - 1}{\lambda + 1}$, for example. The following result characterizes the relationship between all eigenvalues of some polynomial eigenproblem in one open half plane and all eigenvalues of another corresponding eigenproblem in the open unit disk.

Theorem 2 For a linear fractional transformation $\omega = \frac{\lambda-1}{\lambda+1}$, the problem to find all eigenvalues of some polynomial eigenproblem $P(\lambda)x = 0$ (1) in the open right half plane $(Re(\lambda) > 0)$ can be converted into a new problem to find all eigenvalues of another eigenproblem $B(\omega)x = 0$ in the open unit disk, where $B(w) = (w+1)^d A_d + (w+1)^{d-1}(1-w)A_{d-1}+\dots+(w+1)(1-w)^{d-1}A_1+(1-w)^d A_0$. Besides, the corresponding eigenvalues of the above two eigenproblems follow the equality $\lambda = \frac{\omega+1}{1-\omega}$.

Proof Substitute $\lambda = \frac{\omega+1}{1-\omega}$ into $P(\lambda)x = 0$, then the conclusion holds clearly since ω is a conformal one-to-one mapping.

Motivated by the above idea, we can also find all eigenvalues of a PEP in the open upper half plane, in the open left half plane or in the open lower half plane. For simplicity, the corresponding result of finding all eigenvalues of a PEP $P(\lambda)x = 0$ in the open upper half plane is shown as follows.

Theorem 3 For a linear fractional transformation $\omega = \frac{\lambda - i}{\lambda + i}$, the problem to find all eigenvalues of $P(\lambda)x = 0$ (1) in the open upper half plane $(Im(\lambda) > 0)$ can be converted into a new problem to find all eigenvalues of $B(\omega)x = 0$ in the open unit disk, and the corresponding eigenvalues follow the equality $\lambda = \frac{\omega + 1}{1 - \omega}i$.

Proof The proof is similar to that of Theorem 2. Hence, it is omitted.

3 Computing all eigenvalues of a PEP in the open unit disk

In this section, we will briefly present the algorithm to compute all eigenvalues of a PEP (1) in the open unit disk, see Asakura et al. (2010) for details.

For a PEP (1), we define a function $f(\lambda) = \det(P(\lambda))$, then $f(\lambda)$ is an analytical function in λ . As we know, λ is an eigenvalue of a PEP (1) if and only if $f(\lambda) = 0$.

Let Γ be the boundary of an open unit disk for the PEP (1). The center and radius of the open unit disk are denoted as γ and ρ , respectively. Let m ($m \le n$) be a positive integer and the complex moment u_p be

$$u_p = \frac{1}{2\pi i} \oint_{\Gamma} (\lambda - \gamma)^p \frac{f'(\lambda)}{f(\lambda)} dz,$$

where p = 0, 1, ..., 2m - 1.

Using these complex moments u_p , we define the following two $m \times m$ Hankel matrices

$$H_m = \begin{pmatrix} u_0 & u_1 & \cdots & u_{m-1} \\ u_1 & u_2 & \cdots & u_m \\ \vdots & \vdots & & \vdots \\ u_{m-1} & u_m & \cdots & u_{2m-2} \end{pmatrix}$$

and

$$H_{m}^{<} = \begin{pmatrix} u_{1} & u_{2} & \cdots & u_{m} \\ u_{2} & u_{3} & \cdots & u_{m+1} \\ \vdots & \vdots & & \vdots \\ u_{m} & u_{m+1} & \cdots & u_{2m-1} \end{pmatrix}.$$

Now, we consider that the integration is evaluated via a trapezoidal rule on the circle Γ . Let N be the number of sample points on the circle Γ and $v_j = \gamma + \rho \exp(2\pi ji/N)(j = 0, 1, ..., N-1)$. According to the trapezoidal rule, the complex moments u_p can be approximately obtained by the following formula

$$u_p \approx \widehat{u}_p = \frac{1}{N} \sum_{j=0}^{N-1} (v_j - \gamma)^{p+1} \frac{f'(v_j)}{f(v_j)}.$$

Based on the Trace-Theorem of Devidenko (see Davidenko 1960) $\frac{f'(\lambda)}{f(\lambda)} = \text{tr}(T^{-1}(\lambda)T'(\lambda)),$ \hat{u}_p can be rewritten as

$$\widehat{u}_p = \frac{1}{N} \sum_{j=0}^{N-1} (\nu_j - \gamma)^{p+1} \operatorname{tr} \left(T^{-1}(\nu_j) T'(\nu_j) \right).$$
(2)

Using \hat{u}_p (p = 0, 1, ..., 2m - 1), we form the following two Hankel matrices $\hat{H}_m^<$ and \hat{H}_m

$$H_m \approx \widehat{H}_m = \begin{pmatrix} \widehat{u}_0 & \widehat{u}_1 & \cdots & \widehat{u}_{m-1} \\ \widehat{u}_1 & \widehat{u}_2 & \cdots & \widehat{u}_m \\ \vdots & \vdots & & \vdots \\ \widehat{u}_{m-1} & \widehat{u}_m & \cdots & \widehat{u}_{2m-2} \end{pmatrix}$$
(3)

and

$$H_m^{<} \approx \widehat{H}_m^{<} = \begin{pmatrix} \widehat{u}_1 & \widehat{u}_2 & \cdots & \widehat{u}_m \\ \widehat{u}_2 & \widehat{u}_3 & \cdots & \widehat{u}_{m+1} \\ \vdots & \vdots & & \vdots \\ \widehat{u}_m & \widehat{u}_{m+1} & \cdots & \widehat{u}_{2m-1} \end{pmatrix}.$$
 (4)

The numerical algorithm to compute all eigenvalues of the PEP (1) in the open unit disk can be described as follows.

Algorithm 1. An algorithm for computing all eigenvalues of the PEP (1) in the open unit disk

- 1. Choose the number N of sample points on the circle Γ , the center γ and the radius ρ of the circle Γ .
- 2. Compute $v_j = \gamma + \rho \exp(2\pi i(j)/N)$, where j = 0, 1, ..., N 1.
- 3. Compute \hat{u}_p (p = 0, 1, ..., 2m 1) with (2).
- 4. Form two $m \times m$ Hankel matrices \widehat{H}_m and $\widehat{H}_m^<$ with (3) and (4), respectively.
- 5. Compute all eigenvalues λ_k (k = 1, 2, ..., m) of $\widehat{H}_m^< \lambda \widehat{H}_m$.

Remark 3.1 The low order $\widehat{H}_m^< - \lambda \widehat{H}_m$ may be solved by the QZ method (Moler and Stewart 1973).

According to the Theorem 3.5 in Asakura et al. (2010), the eigenvalues λ_k (k = 1, 2, ..., m) of $\hat{H}_m^< x = \lambda \hat{H}_m x$ are the eigenvalues of the PEP (1) in the circle Γ .

4 Computing all eigenvalues of a PEP in an open half plane

For the linear fractional transformation

$$\omega = \frac{\lambda - 1}{\lambda + 1},\tag{5}$$

substitute (5) into $P(\lambda)x = 0$, then it follows that

$$B(w) = (w+1)^d A_d + (w+1)^{d-1} (1-w) A_{d-1} + \dots + (w+1)(1-w)^{d-1} A_1 + (1-w)^d A_0.$$

Next use Algorithm 1 to compute all eigenvalues ω_k of $B(\omega)x = 0$ in the open unit disk, and then substitute ω_k into $\lambda = \frac{\omega+1}{1-\omega}$. Thus, we compute all eigenvalues λ_k of original eigenvalue problem in the open right half plane according to Theorem 1.

From the above results and notations, we can derive the following algorithm to find all eigenvalues of the PEP in the open right half plane.

Algorithm 2. The algorithm to compute all eigenvalues of a polynomial eigenvalue problem in the open right half plane

- 1. Form a new polynomial eigenvalue problem $B(\omega)x = 0$, where $B(\omega) = (\omega+1)^d A_d + (\omega+1)^{d-1}(1-\omega)A_{d-1} + \dots + (\omega+1)(1-\omega)^{d-1}A_1 + (1-\omega)^d A_0$.
- 2. Compute all eigenvalues ω_k of $B(\omega)x = 0$ in the open unit disk by Algorithm 1.
- 3. Compute $\lambda_k = \frac{\omega_k + 1}{1 \omega_k}$.

Similarly, we can also get the algorithm to find all eigenvalues of a polynomial eigenvalue problem in the open upper half plane.



Algorithm 3. The algorithm to compute all eigenvalues of a polynomial eigenvalue problem in the open upper half plane

- 1. Form a new polynomial eigenvalue problem $B(\omega)x = 0$, where $B(\omega) = (\omega + 1)^d i^d A_d + (\omega + 1)^{d-1} i^{d-1} (1 - \omega) A_{d-1} + \dots + (\omega + 1) i (1 - \omega)^{d-1} A_1 + (1 - \omega)^d A_0$.
- 2. Compute all eigenvalues ω_k of $B(\omega)x = 0$ in the open unit disk by Algorithm 1.
- 3. Compute $\lambda_k = \frac{\omega_k + 1}{1 \omega_k} i$.

5 Numerical examples

In this section, we report some numerical examples to show the effectiveness of the proposed algorithms. All computations are carried out in Matlab (2012b). In our examples, λ_k^* and λ_k are the *k*th exact eigenvalue and the *k*th approximate eigenvalue of the PEP (1), respectively. All eigenvalues λ_k are rounded to 10 decimal places in Tables 1, 2, 3 and 4.

Example 1 Consider the generalized eigenvalue problem with

$$P(\lambda) x = (\lambda A_0 + A_1) x = 0,$$

where $A_1 = \begin{pmatrix} 1 & 1 & 0 & 0 & 0 & \cdots & 0 \\ 0 & 3 & 1 & 0 & 0 & \cdots & 0 \\ 0 & 0 & -1 & 1 & 0 & \cdots & 0 \\ 0 & 0 & 0 & -2 & \ddots & \ddots & \vdots \\ \vdots & \vdots & \ddots & \ddots & \ddots & \ddots & \ddots & \vdots \\ 0 & \cdots & \cdots & \ddots & 0 & -97 & 1 \\ 0 & \cdots & \cdots & 0 & 0 & 4 \end{pmatrix}$ and $A_0 = I \in C^{100 \times 100}.$

It is easy to verify that all eigenvalues λ_k^* of this question in the right half plane are 1, 3 and 4.

The numerical results are obtained with the parameters N = 2000, $\gamma = 0$ and $\rho = 1$. The eigenvalues λ_k computed by Algorithm 2 are given in Table 1. It shows that the numerical results are very close to the exact eigenvalues.

Table 1 Numerical results in Example 1	k	λ_k	$ \lambda_k - \lambda_k^* $
	1	1.0000000000	1.9e-12
	2	3.000000000	1.8e-10
	3	4.0000000000	1.8e-10
Table 2 Numerical results in Example 2	k	λ_k	$ \lambda_k - \lambda_k^* $
	1	1.000000000	5.8e-15
	2	2.0000000000	1.0e-14



8.3e-13

1.6e - 12

1.8e-13

Table 3Numerical results inExample 3	k	λ_k	$ \lambda_k - \lambda_k^* $
	1	0.0257024260 + 0.4701394322i	6.5e-16
	2	0.4373052984 + 0.9271387105i	9.6e-15
	3	2.0226888997	3.2e-14
	4	0.4373052984 - 0.9271387105i	7.2e-15
	5	0.0257024260 - 0.4701394322 <i>i</i>	2.4e-16
Table 4 Numerical results in Example 4	k	λ_k	$ \lambda_k - \lambda_k^* $
	1	1.4928844738 + 1.2068208717i	1.3e-13
	2	0.7671255871 + 0.7267920145i	1.4e-12
	3	0.5715381735 + 0.6134359109i	8.1e-13

-0.5715381735 + 0.6134359109i

-0.7671255871 + 0.7267920145i

-1.4928844738 + 1.2068208717i

Example 2 In this example we consider the quadratic eigenvalue problem with

$$P(\lambda)x = \left(\begin{pmatrix} 0 & 1 \\ -2 & 3 \end{pmatrix} + \lambda \begin{pmatrix} 7 & -5 \\ 10 & -8 \end{pmatrix} + \lambda^2 \begin{pmatrix} 1 & 0 \\ 0 & 1 \end{pmatrix} \right) x = 0.$$

We can verify that all eigenvalues λ_k^* of $P(\lambda)x = 0$ in the open right half plane are 1 and 2.

The numerical results for Algorithm 2 with the parameters N = 20, $\gamma = 0$ and $\rho = 1$ are shown in Table 2. From Table 2, the numerical results are in good agreement with the exact eigenvalues.

Example 3 We consider the polynomial eigenvalue problem with

4

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$$P(\lambda) x = (\lambda^3 A_0 + \lambda^2 A_1 + \lambda A_2 + A_3) x = 0,$$

where $A_3 = \begin{pmatrix} -16 & -4 & 7 \\ -14 & 7 & 13 \\ 6 & 8 & 7 \end{pmatrix}, A_2 = \begin{pmatrix} 0 & 0 & 0 \\ 0 & 0 & 0 \\ 0 & 0 & 0 \end{pmatrix}, A_1 = \begin{pmatrix} 2 & -6 & 1 \\ -2 & 22 & 11 \\ 7 & -1 & 1 \end{pmatrix}$ and
 $A_0 = \begin{pmatrix} -4 & 3 & 12 \\ -17 & -11 & 0 \\ 1 & -1 & 3 \end{pmatrix}.$

It is easy to verify that all eigenvalues λ_k^* of this question in the right half plane are 2.02268890, 0.43730530 + 0.92713871*i*, 0.43730530 - 0.92713871*i*, 0.02570243 + 0.47013943*i*, 0.02570243 - 0.47013943*i*.

The numerical results are obtained with the parameters N = 1000, $\gamma = 0$ and $\rho = 1$. The computed eigenvalues λ_k are given in Table 3. It is clearly seen from Table 3 that the numerical results obtained by Algorithm 2 are very close to the exact eigenvalues.

Example 4 We consider another polynomial eigenvalue problem with

$$P(\lambda)x = (\lambda^{4}A_{0} + \lambda^{3}A_{1} + \lambda^{2}A_{2} + \lambda A_{3} + A_{4})x = 0$$

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where
$$A_4 = \begin{pmatrix} -2 & 4 & 11 \\ -5 & 7 & 8 \\ 6 & 8 & 9 \end{pmatrix}$$
, $A_3 = \begin{pmatrix} 0 & 0 & 0 \\ 0 & 0 & 0 \\ 0 & 0 & 0 \end{pmatrix}$, $A_2 = \begin{pmatrix} 0 & 5 & 0 \\ 0 & 4 & 0 \\ 0 & 3 & 0 \end{pmatrix}$, $A_1 = \begin{pmatrix} 0 & 0 & 0 \\ 0 & 0 & 0 \\ 0 & 0 & 0 \end{pmatrix}$ and $A_0 = \begin{pmatrix} 6 & -6 & 1 \\ -2 & 22 & 10 \\ 7 & -1 & 3 \end{pmatrix}$.

It is not difficult to verify that all eigenvalues λ_k^* in the open upper half plane are 1.49288447 + 1.20682087i, 0.76712559 + 0.72679201i, 0.57153817 + 0.61343591i, -0.57153817 + 0.61343591i, -0.76712559 + 0.72679201i, -1.49288447 + 1.20682087i.

The numerical results for Algorithm 3 are given in Table 4 using the parameters N = 1000, $\gamma = 0$ and $\rho = 1$. From Table 4, we observe that the numerical results coincide with the exact eigenvalues.

6 Conclusion

In this paper, we have presented a novel method to compute all eigenvalues of the PEP in an open half plane. This approach first transforms the original PEP into a new PEP via a linear fractional transformation, then employs the contour integral method to compute all eigenvalues of the new PEP in the open unit disk, and substitute the above eigenvalues into the linear fractional transformation to find all eigenvalues of original eigenproblem lastly. Our numerical results show that the proposed method is very useful and effective for computing all eigenvalues of the PEP in an open half plane.

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References

Adhikari S, Pascual B (2009) Eigenvalues of linear viscoelastic systems. J Sound Vib 325:1000-1011

- Asakura J, Sakurai T, Tadano H, Kimura K (2010) A numerical method for polynomial eigenvalue problems using contour integral. Jpn J Ind Appl Math 27:73–90
- Bai Z-Z, Miao C-Q (2017a) On local quadratic convergence of inexact simplified Jacobi–Davidson method. Linear Algebra Appl 520:215–241
- Bai Z-Z, Miao C-Q (2017b) On local quadratic convergence of inexact simplified Jacobi–Davidson method for interior eigenpairs of Hermitian eigenproblems. Appl Math Lett 72:23–28
- Chu EK-W (2003) Perturbation of eigenvalues for matrix polynomials via the Bauer–Fike theorems. SIAM J Matrix Anal Appl 25:551–573
- Dai H, Bai Z-Z, Wei Y (2015) On the solvability condition and numerical algorithm for the parameterized generalized inverse eigenvalue problems. SIAM J Matrix Anal Appl 36:707–726

Davidenko DF (1960) Algorithms for λ -matrices. Sov Math 1:316–319

- Dedieu J, Tisseur F (2003) Perturbation theory for homogeneous polynomial eigenvalue problems. Linear Algebra Appl 358:71–94
- Gohberg I, Lancaster P, Rodman L (1979) Perturbation theory for divisors of operator polynomials. SIAM J Math Anal 10:1161–1183

Gohberg I, Lancaster P, Rodman L (1982) Matrix polynomials. Academic Press, New York

Gupta KK (1976) On a finite dynamic element method for free vibration analysis of structures. Comput Methods Appl Mech Eng 9:105–120



Higham NJ, Tisseur F (2002) More on pseudo spectra for polynomial eigenvalue problems and applications in control theory. Linear Algebra Appl 351:435–453

Higham NJ, Tisseur F (2003) Bounds for eigenvalues of matrix polynomials. Linear Algebra Appl 358:5–22

Higham NJ, Mackey DS, Tisseur F (2006) The conditioning of linearizations of matrix polynomials. SIAM J Matrix Anal Appl 28:1005–1028

- Higham NJ, Li R-C, Tisseur F (2007) Backward error of polynomial eigenproblems solved by linearization. SIAM J Matrix Anal Appl 29:1218–1241
- Hoffnung L, Li R-C, Ye Q (2006) Krylov type subspace methods for matrix polynomials. Linear Algebra Appl 415:52–81
- Huang W-Q, Li T-X, Li Y-T, Lin W-W (2013) A semiorthogonal generalized Arnoldi method and its variations for quadratic eigenvalue problems. Numer Linear Algebra Appl 20:259–280
- Hwang T-M, Lin W-W, Wang W-C, Wang W (2004) Numerical simulation of three dimensional pyramid quantum dot. J Comput Phys 196:208–232
- Hwang T-M, Lin W-W, Liu J-L, Wang W (2005) Jacobi–Davidson methods for cubic eigenvalue problems. Numer Linear Algebra Appl 12:605–624
- Hwang F-N, Wei Z-H, Hwang T-M, Wang W (2010) A parallel additive Schwarz preconditioned Jacobi– Davidson algorithm for polynomial eigenvalue problems in quantum dot simulation. J Comput Phys 229:2932–2947
- Lawrence PW, Corless RM (2015) Backward error of polynomial eigenvalue problems solved by linearization of Lagrange interpolants. SIAM J Matrix Anal Appl 36:1425–1442
- Mackey DS, Mackey N, Mehl C, Mehrmann V (2006a) Vector spaces of linearizations for matrix polynomials. SIAM J Matrix Anal Appl 28:971–1004
- Mackey DS, Mackey N, Mehl C, Mehrmann V (2006b) Structured polynomial eigenvalue problems: good vibrations from good linearization. SIAM J Matrix Anal Appl 28:1029–1051
- Moler CB, Stewart GW (1973) An algorithm for generalized matrix eigenvalue problems. SIAM J Numer Anal 10:241–256
- Rudin W (1987) Real and complex analysis. McGraw-Hill, New York
- Sleijpen GLG, Booten AGL, Fokkema DR, van der Vorst HA (1996) Jacobi–Davidson type methods for generalized eigenproblems and polynomial eigenproblems. BIT 36:595–633
- Tisseur F (2000) Backward error and condition of polynomial eigenvalue problems. Linear Algebra Appl 309:339–361
- Tisseur F, Meerbergen K (2001) The quadratic eigenvalue problem. SIAM Rev 43:235-286
- Wilkinson JH (1965) The algebraic eigenvalue problem. Clarendon Press, Oxford

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