# On the Pseudospectra of Matrix Polynomials

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http://www.maths.man.ac.uk/~nareports/narep445.pdf

## 0. Prologue

Consider the differential equation

 $A_m u^{(m)}(t) + A_{m-1} u^{(m-1)} + \dots +$ 

 $+\cdots + A_1 u'(t) + A_0 u(t) = f(t)$  (1)

and the difference equation

 $A_m u_{j+m} + A_{m-1} u_{j+m-1} + \dots +$ 

 $+\cdots + A_1 u_{j+1} + A_0 u_j = f_j,$  (2)

where  $A_j \in \mathbb{C}^{n \times n}$ ,  $u(t), u_j \in \mathbb{C}^n$  and  $\underline{\det A_m \neq 0}$ .

Applying the Laplace transformation to (1) or the Z-transformation to (2) yields the *matrix polynomial* 

 $P(\lambda) = A_m \lambda^m + A_{m-1} \lambda^{m-1} + \dots + A_1 \lambda + A_0.$ 

A  $0 \neq x_0 \in \mathbb{C}^n$  is an *eigenvector* of  $P(\lambda)$  corresponding to the *eigenvalue*  $\lambda_0 \in \mathbb{C}$  if

 $P(\lambda_0)x_0 = 0$  (eigenproblem).

If, in addition,  $x_1, x_2, \ldots, x_k \in \mathbb{C}^n$  satisfy

$$\sum_{j=1}^{\xi} \frac{1}{j!} P^{(j)}(\lambda_0) x_{\xi-j} = 0 \quad ; \quad \xi = 1, 2, \dots, k,$$

then  $x_0, x_1, \ldots, x_k$  is a *Jordan chain* of  $P(\lambda)$ . The solution of (1) is of the form

$$u(t) = X_P e^{t J_P} c + \int_{t_0}^t X_P e^{(t-s) J_P} Y_P f(s) ds$$

and the solution of (2) is of the form

$$u_j = X_P J_P^j c + \sum_{i=0}^{\nu-1} X_P J_P^{\nu-i-1} Y_P f_i$$

#### **<u>1. Introduction</u>**

The *spectrum* of  $P(\lambda) = A_m \lambda^m + A_{m-1} \lambda^{m-1} + \dots + A_1 \lambda + A_0$ is the set of all eigenvalues of  $P(\lambda)$ ,

 $\sigma(P) = \{\lambda \in \mathbb{C} : \det P(\lambda) = 0\},\$ 

where det  $P(\lambda)$  is a scalar polynomial of degree n m, with leading coefficient det  $A_m \neq 0$ .

We are interested in the spectra of perturbations of  $P(\lambda)$  of the form

 $P_{\Delta}(\lambda) = (A_m + \Delta_m)\lambda^m + (A_{m-1} + \Delta_{m-1})\lambda^{m-1}$ 

 $+\cdots+(A_1+\Delta_1)\lambda+A_0+\Delta_0,$ 

where  $\Delta_0, \Delta_1, \ldots, \Delta_m \in \mathbb{C}^{n \times n}$  are arbitrary.

For a given  $\varepsilon > 0$  and a given set of nonnegative weights  $\mathbf{w} = \{w_0, w_1, \dots, w_m\}$ , the  $\varepsilon$ *pseudospectrum* of  $P(\lambda)$  with respect to  $\mathbf{w}$ (introduced by Tisseur and Higham, 2001) is

 $\sigma_{\varepsilon,\mathbf{w}}(P) = \{\lambda \in \mathbb{C} : \det P_{\Delta}(\lambda) = 0, \}$ 

 $\|\Delta_j\| \leq \varepsilon \, w_j, \ j = 0, 1, \dots, m\}.$ 

 $w_0, w_1, \ldots, w_m \ge 0$  allow freedom in how perturbations are measured; for example, in an absolute sense when  $w_0 = w_1 = \cdots = w_m = 1$ , or in a relative sense when  $w_j = ||A_j||$ . Different values for  $w_j$  admit different levels of confidence in  $A_j$ .

Note that for  $\varepsilon = 0$ ,  $\sigma_{0,\mathbf{w}}(P) = \sigma(P)$ .

For  $P(\lambda) = I\lambda - A$   $(A \in \mathbb{C}^{n \times n})$ ,  $\sigma(P)$  coincides with the spectrum of A,  $\sigma(A)$ . If, in addition,  $\mathbf{w} = \{w_0, w_1\} = \{1, 0\}$ , then  $\sigma_{\varepsilon, \mathbf{w}}(P)$  coincides with the  $\varepsilon$ -pseudospectrum of A,

 $\sigma_{\varepsilon}(A) = \{\lambda \in \mathbb{C} : \lambda \in \sigma(A + \Delta_0), \|\Delta_0\| \le \varepsilon\}.$ 

For the spectral norm, defining the scalar  $q_{\mathbf{w}}(\lambda) = w_m \lambda^m + w_{m-1} \lambda^{m-1} + \dots + w_1 \lambda + w_0$ , one of the main tools is the formula (Tisseur-Higham, 2001) ( $s_{\min}$ : min. singular value)

 $\sigma_{\varepsilon,\mathbf{w}}(P) = \{\lambda \in \mathbb{C} : s_{\min}(P(\lambda)) \le \varepsilon q_{\mathbf{w}}(|\lambda|) \}.$ 

As the eigenvalues of  $P_{\Delta}(\lambda)$  are continuous,

 $\partial \sigma_{\varepsilon, \mathbf{w}}(P) = \{\lambda \in \mathbb{C} : s_{\min}(P(\lambda)) = \varepsilon q_{\mathbf{w}}(|\lambda|) \}.$ 

### 2. Examples (using the spectral norm)

**Example 1** (*A wing problem*) The eigenproblem of the matrix polynomial

$$Q(\lambda) = \begin{bmatrix} 17.6 & 1.28 & 2.89 \\ 1.28 & 0.824 & 0.413 \\ 2.89 & 0.413 & 0.725 \end{bmatrix} \lambda^2 + \\ + \begin{bmatrix} 7.66 & 2.45 & 2.1 \\ 0.23 & 1.04 & 0.223 \\ 0.6 & 0.756 & 0.658 \end{bmatrix} \lambda + \begin{bmatrix} 121 & 18.9 & 15.9 \\ 0 & 2.7 & 0.145 \\ 11.9 & 3.64 & 15.5 \end{bmatrix}.$$

arose from a study of the oscillations of a wing in an airstream. The eigenvalues of  $Q(\lambda)$  are

 $-0.88 \pm i 8.44, \ 0.09 \pm i 2.52, \ -0.92 \pm i 1.76.$ 

Perturbations are measured in the absolute sense, i.e.,  $w_0 = w_1 = w_2 = 1$ .



 $\partial \sigma_{\varepsilon,\mathbf{w}}(Q)$  for  $\varepsilon = 0.01, 0.1, 0.15 < s_{\min}(A_2)$ .



The less sensitive eigenvalues.



 $\partial \sigma_{\varepsilon,\mathbf{w}}(Q)$  for  $\varepsilon = 0.18 > s_{\min}(A_2)$ .



The less sensitive eigenvalues.

**Example 2** (A vibrating system) The  $3 \times 3$  selfadjoint matrix polynomial

$$P(\lambda) = \begin{bmatrix} 1 & 0 & 0 \\ 0 & 2 & 0 \\ 0 & 0 & 5 \end{bmatrix} \lambda^2 + \begin{bmatrix} 0 & 0 & 0 \\ 0 & 3 & -1 \\ 0 & -1 & 6 \end{bmatrix} \lambda + \begin{bmatrix} 2 & -1 & 0 \\ -1 & 3 & 0 \\ 0 & 0 & 10 \end{bmatrix}$$

corresponds to a mass-spring model described by Falk (1960). The eigenvalues of  $P(\lambda)$  are

 $-0.08 \pm i 1.45, -0.75 \pm i 0.86, -0.51 \pm i 1.25.$ 

Perturbations are measured in a relative sense, i.e.,  $w_0 = ||A_0|| = 10$ ,  $w_1 = ||A_1|| = 6.3$  and  $w_2 = ||A_2|| = 5$ .



 $\partial \sigma_{\varepsilon,\mathbf{w}}(P)$  for  $\varepsilon = 0.02, 0.05, 0.1.$ 

#### **Example 3** (A gyroscopic system)

Let *B* be the  $10 \times 10$  nilpotent matrix with ones on the subdiagonal and zeros elsewhere. Define  $\hat{M} = (4I_{10} + B + B^T)/6$ ,  $\hat{G} = B - B^T$  $\hat{K} = B + B^T - 2I_{10}$ , and set

- $M = I_{10} \otimes \hat{M} + 1.30 \hat{M} \otimes I_{10},$
- $G = 1.35I_{10} \otimes \hat{G} + 1.10\hat{G} \otimes I_{10},$
- $K = I_{10} \otimes \hat{K} + 1.20 \hat{K} \otimes I_{10}.$

The  $100 \times 100$  matrix polynomial  $M\lambda^2 + G\lambda + K$ corresponds to a gyroscopic system (Mehrmann-Watkins, 2001). Adding the damping matrix  $D = \text{tridiag}\{-0.1, 0.3, -0.1\}$  to G yields

#### $R(\lambda) = M\lambda^2 + (G+D)\lambda + K.$

Perturbations are measured in the absolute sense.



 $\partial \sigma_{\varepsilon,\mathbf{w}}(R)$  for  $\varepsilon = 0.004, 0.02, 0.1.$ 

### 3. General Properties

Consider an  $n \times n$  matrix polynomial  $P(\lambda) = A_m \lambda^m + A_{m-1} \lambda^{m-1} + \dots + A_1 \lambda + A_0.$ 

**Proposition 1** If the coefficients of  $P(\lambda)$  are all real or all hermitian, then for any  $\varepsilon > 0$  and  $\mathbf{w} = \{w_0, w_1, \dots, w_m\}, \sigma_{\varepsilon, \mathbf{w}}(P)$  is symmetric with respect to the real axis.

**Proof** Based on the observation

 $\|\Delta_j\| = \|\overline{\Delta}_j\| = \|\Delta_j^*\|. \quad \Box$ 

**Theorem 2** The pseudospectrum  $\sigma_{\varepsilon,\mathbf{w}}(P)$  is bounded if and only if  $0 \notin \sigma_{\varepsilon w_m}(A_m)$ , i.e., if and only if

 $s_{\min}(A_m) \geq \varepsilon w_m.$ 

**Proof** Suppose  $0 \notin \sigma_{\varepsilon w_m}(A_m)$  and define

 $\zeta_{\varepsilon} = \min\{|\det(A_m + \Delta_m)| : ||\Delta_m|| \le \varepsilon w_m\} > 0.$ 

Then there is an  $M_{\varepsilon} > 0$  such that for any associated perturbation

 $P_{\Delta}(\lambda) = (A_m + \Delta_m)\lambda^m + \dots + (A_1 + \Delta_1)\lambda + A_0 + \Delta_0$ 

and for any  $\lambda \in \mathbb{C}$  with  $|\lambda| > M_{\varepsilon}$ ,

 $|\det P_{\Delta}(\lambda) - \det(A_m + \Delta_m)\lambda^{nm}| < \zeta_{\varepsilon}|\lambda^{mn}|$ 

 $\leq |\det(A_m + \Delta_m)\lambda^{mn}|,$ 

i.e., det  $P_{\Delta}(\lambda) \neq 0$ . Hence,  $\sigma_{\varepsilon, \mathbf{w}}(P)$  is bounded.

To prove the converse, assume that  $\sigma_{\varepsilon,\mathbf{w}}(P)$  is bounded but there is an associated

 $P_{\widehat{\Delta}}(\lambda) = (A_m + \widehat{\Delta}_m)\lambda^m + \dots + (A_1 + \widehat{\Delta}_1)\lambda + A_0 + \widehat{\Delta}_0,$ 

with  $\det(A_m + \hat{\Delta}_m) = 0$ . One of the coefficients of  $\det P_{\hat{\Delta}}(\lambda)$ , let of  $\lambda^{\tau}$ , is  $\beta_{\tau} \neq 0$ . Construct a sequence  $\{\hat{\Delta}_{m,k}\}_{k\in\mathbb{N}} \subset \mathbb{C}^{n\times n}$  such that  $\lim_{k\to\infty} \hat{\Delta}_{m,k} = \hat{\Delta}_m$ , and for every  $k \in \mathbb{N}$ ,

 $\det(A_m + \hat{\Delta}_{m,k}) \neq 0$  and  $\|\hat{\Delta}_{m,k}\| \leq \varepsilon w_m$ .

Since  $\sigma_{\varepsilon, \mathbf{w}}(P)$  is bounded, the  $(nm - \tau)$ th elementary symmetric function of the zeros of

 $\det[(A_m + \hat{\Delta}_{m,k})\lambda^m + \dots + (A_1 + \hat{\Delta}_1)\lambda + A_0 + \hat{\Delta}_0],$ 

which is equal to  $\pm \beta_{\tau} / \det(A_m + \hat{\Delta}_{m,k})$ , is bounded for all k; this is a contradiction.

**Theorem 3** If  $\sigma_{\varepsilon,\mathbf{w}}(P)$  is bounded, then it has no more than nm con. components, and any associated  $P_{\Delta}(\lambda)$  has the same number ( $\geq 1$ ) of eigenvalues with  $P(\lambda)$  in each one of these components, counting multiplicities.

**Proof** By Theorem 2, for any associated

 $P_{\Delta}(\lambda) = (A_m + \Delta_m)\lambda^m + \dots + (A_1 + \Delta_1)\lambda + A_0 + \Delta_0,$ 

det $(A_m + \Delta_m) \neq 0$ . Thus,  $P_{\Delta}(\lambda)$  has exactly nm eigenvalues, counting multiplicities, as does every

 $P_{\Delta,t}(\lambda) = (1-t)P(\lambda) + tP_{\Delta}(\lambda); \quad t \in [0,1].$ 

As t varies from 0 to 1, the eigenvalues of  $P_{\Delta,t}(\lambda)$  trace continuous paths from the eigenvalues of  $P(\lambda)$  to the eigenvalues of  $P_{\Delta}(\lambda)$ .

### 4. A Curve-Tracing Algorithm

Recall that for the spectral norm,

 $\partial \sigma_{\varepsilon,\mathbf{w}}(P) = \{\lambda \in \mathbb{C} : s_{\min}(P(\lambda)) = \varepsilon q_{\mathbf{w}}(|\lambda|) \}.$ 

For convenience, define

$$g_P(x,y) = s_{\min}(P(x+iy)) ; x,y \in \mathbb{R}$$

and

$$g_P(\lambda) = s_{\min}(P(\lambda)) ; \quad \lambda \in \mathbb{C}.$$

**Theorem 4** (Sun, 1988) Let  $\lambda_0 = x_0 + iy_0 \in \mathbb{C} \setminus \sigma(P)$ . If  $s_{\min}(P(\lambda_0))$  is a simple singular value of  $P(\lambda_0)$ , and  $u_0$ ,  $v_0$  are associated left and right singular vectors, respectively, then

$$\nabla g_P(\lambda_0) = \left( \mathsf{Re}\left( u_0^* \frac{\partial P(\lambda_0)}{\partial x} v_0 \right), \mathsf{Re}\left( u_0^* \frac{\partial P(\lambda_0)}{\partial y} v_0 \right) \right).$$

Our continuation method for drawing

 $\partial \sigma_{\varepsilon,\mathbf{w}}(P) = \{\lambda \in \mathbb{C} : g_P(\lambda) - \varepsilon q_{\mathbf{w}}(|\lambda|) = 0\}$ is an extension of (Brühl, 1996), and consists of an initial step to find a starting point on  $\partial \sigma_{\varepsilon,\mathbf{w}}(P)$  followed by a sequence of "predictor" steps tangential to  $\partial \sigma_{\varepsilon,\mathbf{w}}(P)$  and "corrector" steps to go back to  $\partial \sigma_{\varepsilon,\mathbf{w}}(P)$ .

**Initial Step:** For calculation of a first point on  $\partial \sigma_{\varepsilon, \mathbf{w}}(P)$ , let  $\lambda_0 \in \sigma_{\varepsilon, \mathbf{w}}(P) \setminus \sigma(P)$  and  $d_0 \in \mathbb{C}$  be nonzero. Then use Newton's method to solve

#### $g_P(\lambda_0 + t d_0) - \varepsilon q_{\mathbf{w}}(|\lambda_0 + t d_0|) = 0$

along the straight line  $\{\lambda_0 + t d_0 : t \in \mathbb{R}\}$ . Set  $t_0 = 0$ , and assume that  $g_P$  is differentiable at  $\lambda_0$  and  $\nabla g_P(\lambda_0)$  is given by Theorem 4.

The first Newton iterate gives

$$t_1 = -\frac{g_P(\lambda_0) - \varepsilon q_{\mathbf{w}}(|\lambda_0|)}{(g_P(\lambda_0 + t d_0) - \varepsilon q_{\mathbf{w}}(|\lambda_0 + t d_0|))'}$$

and the point

$$z_{1} = \lambda_{0} - \frac{g_{P}(\lambda_{0}) - \varepsilon q_{\mathbf{w}}(\lambda_{0})}{(\operatorname{Re} d_{0}, \operatorname{Im} d_{0}) \cdot \nabla [g_{P}(\lambda_{0}) - \varepsilon q_{\mathbf{w}}(|\lambda_{0}|)]} d_{0}.$$
(3)

Since  $\lambda_0 \in \sigma_{\varepsilon, \mathbf{w}}(P)$ , for suitable direction  $d_0$ , we estimate a point of  $\partial \sigma_{\varepsilon, \mathbf{w}}(P)$  by repeating (3) until  $|s_{\min}(P(z)) - \varepsilon q_{\mathbf{w}}(|z|)|$  is small enough. In our examples, only a few iterations are required.

For 
$$d_0 = \nabla [g_P(\lambda_0) - \varepsilon q_{\mathbf{w}}(|\lambda_0|)]$$
, (3) implies  
 $z_1 = \lambda_0 - (g_P(\lambda_0) - \varepsilon q_{\mathbf{w}}(|\lambda_0|)) \overline{d_0}^{-1}$ . (4)

**Prediction:** Assuming that  $z_{k-1} \in \partial \sigma_{\varepsilon, \mathbf{w}}(P)$ has been computed and  $\tau_k$  is the corresponding step-length, the (tangential) prediction for the *k*th boundary point of  $\sigma_{\varepsilon, \mathbf{w}}(P)$ ,  $z_k$ , is

$$\widehat{z}_{k} = z_{k-1} + \tau_{k} \left( i \frac{\nabla \left[ g_{P}(z_{k-1}) - \varepsilon \, q_{\mathbf{w}}(|z_{k-1}|) \right]}{\left| \nabla \left[ g_{P}(z_{k-1}) - \varepsilon \, q_{\mathbf{w}}(|z_{k-1}|) \right] \right|} \right).$$

**Correction:** For small  $\tau_k$ , the correction step is a single Newton iterate for the equation  $g_P(\hat{z}_k + t d_k) - \varepsilon q_W(|\hat{z}_k + t d_k|) = 0$ , with an appropriate direction  $d_k$  and initial  $t_0 = 0$ . In our examples, has been found that one Newton step gives satisfactory performance, although the effect of taking more steps could be a subject for further investigation. A natural choice for  $d_k \in \mathbb{C}$  is

$$\widehat{d}_k = \nabla \left[ g_P(\widehat{z}_k) - \varepsilon \, q_{\mathbf{w}}(|\widehat{z}_k|) \right].$$

In this case, the step (4) is written

$$z_k = \hat{z}_k - \left(g_P(z_{k-1}) - \varepsilon \, q_{\mathbf{w}}(|z_{k-1}|)\right) \, \hat{d}_k^{-1}$$

and the estimation of  $z_k$  requires the computation of  $s_{\min}(P(z_{k-1}))$ ,  $s_{\min}(P(\hat{z}_k))$  and their associated left and right singular vectors.



Choosing the direction  $\hat{d}_k$  in the correction step.

The computation of  $s_{\min}(P(\hat{z}_k))$  and the corresponding singular vectors can be avoided (and the computational cost of the algorithm reduced by about a half) if the correction step is taken in the direction of

$$d_k = \nabla \left[ g_P(z_{k-1}) - \varepsilon \, q_{\mathbf{w}}(|z_{k-1}|) \right]$$

and (4) is written in the form

 $z_k = \hat{z}_k - (g_P(z_{k-1}) - \varepsilon q_{\mathbf{w}}(|z_{k-1}|)) \overline{d_k}^{-1}.$ 



Choosing the direction  $d_k$  in the correction step.

## 5. Comments on the Algorithm

(a) It tracks the boundary of that con. component of  $\sigma_{\varepsilon,\mathbf{w}}(P)$  containing  $\lambda_0$ . For a complete picture, it may be necessary to repeat the procedure for several values of  $\lambda_0$ .

(b) It does not require <u>a priori</u> knowledge of the size of  $\sigma_{\varepsilon,\mathbf{w}}(P)$ , since it sketches the con. components of  $\sigma_{\varepsilon,\mathbf{w}}(P)$  one after the other by using starting points close to eigenvalues.

(c) The size of the step-lengths,  $\tau_k$ , in the prediction step affects the accuracy and the computational cost of the algorithm, and it is important to obtain criteria for their selection (Bekas-Gallopoulos, 2001).

(d) The algorithm may lose its path near boundary points where  $\nabla [g_P(\lambda) - \varepsilon q_w(|\lambda|)]$  does not exist or it is zero, and near points where the distance between con. components of  $\sigma_{\varepsilon,w}(P)$ becomes small. Some of these difficulties can be solved by choosing a smaller step-length (increasing the cost). See below  $\sigma_{\varepsilon,w}(P)$  of Example 2, for  $\varepsilon = 0.06$ .



Constant step-lengths  $\tau = 0.03$  and  $\tau = 0.003$ .

## 6. Some Open Questions

(a) What else can we say about the topological and geometrical properties of  $\sigma_{\varepsilon,\mathbf{w}}(P)$  and its con. components?

(b) When  $\sigma_{\varepsilon,\mathbf{w}}(P)$  is unbounded, how many con. components it may have?

(c) Is it true that a bounded con. component  $\mathcal{G}$  of  $\sigma_{\varepsilon,\mathbf{w}}(P)$  contains 2 eigenvalues of  $P(\lambda)$  if and only if an associated perturbation  $P_{\Delta}(\lambda)$  has a multiple eigenvalue in  $\mathcal{G}$ ?

(d) How can  $\sigma_{\varepsilon,\mathbf{w}}(P)$  be used in studying the stability of the spectral factorization of  $P(\lambda)$ ?

(e) Can the path-tracing algorithm be a part of a parallel algorithm?

#### 7. References

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