

# On the Eigenpair Problem

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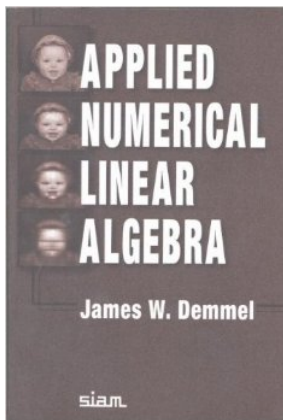
Joint work with D. Armentano, C. Beltrán, P. Bürgisser, and  
M. Shub

Let  $A \in \mathbb{C}^{n \times n}$

We say that  $\lambda \in \mathbb{C}$  is an **eigenvalue** of  $A$  when there exists  $v \in \mathbb{C}^{n \times n} \setminus \{0\}$  such that

$$Av = \lambda v.$$

In this case we say that  $v$  is an **eigenvector**.



So the problem of devising an algorithm [for the eigenvalue problem] that is numerically stable and globally (and quickly!) convergent remains open. [p. 139]

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  - ▶ The QR algorithm with Rayleigh Quotient shift fails for open sets of real input matrices.
  - ▶ It is unknown whether the Francis (double) shift algorithm converges generally on real or complex matrices.

[\*\*\* No theoretical understanding \*\*\*]

**Theorem** We exhibit an algorithm which on input a matrix  $A$  with complex Gaussian entries generates (with probability 1) an “excellent” approximation to all the (eigenvalue, eigenvector) pairs of  $A$ . Moreover, the running time of this algorithm is polynomial in  $n$  on the average.

This algorithm is numerically stable.

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- ▶ *Approximation à la Smale.* An appropriate version of Newton's iteration, starting at  $\tilde{\zeta}$ , converges immediately, quadratically fast, to  $\zeta$ .

# Newton's method

We define a **Newton map** associated to  $A$

$$N_A : \mathbb{C} \times (\mathbb{C}^n \setminus \{0\}) \rightarrow \mathbb{C} \times (\mathbb{C}^n \setminus \{0\}).$$

Let  $\mathcal{V} := \{(A, \lambda, v) \mid (A - \lambda \text{Id})v = 0\}$ .



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Given  $(A, \lambda, v) \in \mathcal{V}$  we say that  $(\zeta, w) \in \mathbb{C} \times (\mathbb{C}^n \setminus \{0\})$  is an **approximate eigenpair** of  $A$  with associated eigenpair  $(\lambda, v)$  when for all  $k \geq 1$  the  $k$ th iterate  $N_A^k(\zeta, w)$  of the Newton map at  $(\zeta, w)$  is well defined and satisfies

$$\text{dist}((N_A^k(\zeta, w)), (\lambda, v)) \leq \left(\frac{1}{2}\right)^{2^k - 1} \text{dist}((\zeta, w), (\lambda, v)).$$

# Well-posedness

A triple  $(A, \lambda, v)$  is **well-posed** when  $\lambda$  is a simple eigenvalue of  $A$ . We write  $(A, \lambda, v) \in \mathcal{W}$ . Otherwise, it is said to be **ill-posed**.

**Proposition** Ill-posed eigenpairs have no approximate eigenpairs à la Smale.

**Proposition** The set  $\Sigma$  of matrices  $A$  with multiple eigenvalues has (real) codimension 2 in  $\mathbb{C}^{n \times n}$ .

Given  $(A, \lambda, \nu) \in \mathbb{C}^{n \times n} \times \mathbb{C} \times \mathbb{P}(\mathbb{C}^n)$ , we let  $A_{\lambda, \nu} : T_\nu \rightarrow T_\nu$  be

$$A_{\lambda, \nu} := P_{\nu^\perp} \circ (A - \lambda \text{Id})|_{T_\nu}.$$

We define the **condition number** of the triple  $(A, \lambda, \nu)$  by

$$\mu(A, \lambda, \nu) := \|A\|_F \|A_{\lambda, \nu}^{-1}\|,$$

**Theorem** Let  $A \in \mathbb{C}^{n \times n}$  with  $\|A\|_F = 1$  and  $(\lambda, \nu), (\lambda_0, \nu_0) \in \mathbb{C} \times (\mathbb{C}^n \setminus \{0\})$ . If  $(\lambda, \nu)$  is a well-posed eigenpair of  $A$  and

$$\text{dist}((\lambda, \nu), (\lambda_0, \nu_0)) < \frac{c_0}{\mu(A, \lambda, \nu)}$$

then  $(\lambda_0, \nu_0)$  is an approximate eigenpair of  $A$  with associated eigenpair  $(\lambda, \nu)$ . One may choose  $c_0 = 0.2881$ .

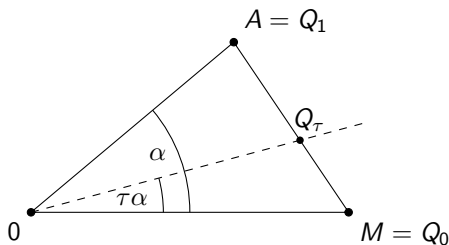
# The Algorithm

We are given  $A \in \mathbb{C}^{n \times n}$  and an initial triple  $(M, \lambda_0, \nu_0)$  in  $\mathcal{W}$ .

Consider the line segment  $[M, A]$  in  $\mathbb{C}^{n \times n}$  with endpoints  $M$  and  $A$

$$[M, A] = \{Q_\tau \in \mathbb{C}^{n \times n} \mid \tau \in [0, 1]\}$$

with  $Q_\tau$  being the only point in  $[M, A]$  such that  $d_{\mathbb{S}}(M, Q_\tau) = \tau\alpha$ .



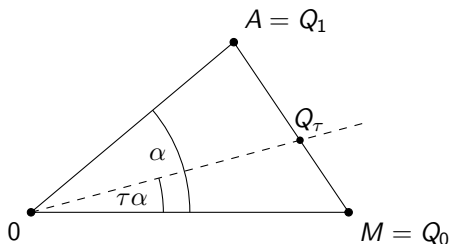
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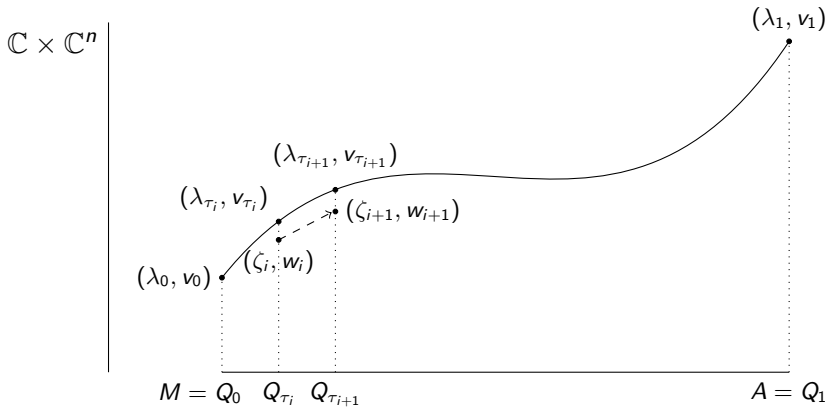
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When  $\tau$  moves from 0 to 1 the eigenpair  $(\lambda_\tau, v_\tau)$  of  $Q_\tau$  moves from  $(\lambda_0, v_0)$  to an eigenpair  $(\lambda_1, v_1)$  of  $A = Q_1$ .

We want to “follow” the curve  $(\lambda_\tau, v_\tau)$ .



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## Path-follow

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**Input:**  $A \in \mathbb{C}^{n \times n}$  and  $(M, \lambda_0, v_0) \in \mathcal{W}$

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$\alpha := d_{\mathbb{S}}(M, A)$ ,  $r := \|A\|_F$ ,  $s := \|M\|_F$

$\tau := 0$ ,  $Q := M$ ,  $(\zeta, w) := (\lambda_0, v_0)$

repeat

$$\Delta\tau := \frac{0.001461}{\alpha \mu^2(Q, \zeta, w)}$$

$$\tau := \min\{1, \tau + \Delta\tau\}$$

$$t := \frac{s}{s + r(\sin \alpha \cot(\tau\alpha) - \cos \alpha)}$$

$$Q := tA + (1 - t)M$$

$$(\zeta, w) := N_Q(\zeta, w)$$

until  $\tau = 1$

return  $(\zeta, w)$

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**Output:**  $(\zeta, w) \in \mathbb{C} \times \mathbb{C}^n$ , approximate eigenpair of  $A$ .

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# Complexity

Cost of Path-follow on input  $(A, M, \lambda_0, v_0)$ :

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$$\begin{array}{ccc} & \nearrow & \nwarrow \\ & K(A, M, \lambda_0, v_0) & \mathcal{O}(n^3) \end{array}$$

Path-follow terminates (i.e.,  $K(A, M, \lambda_0, v_0) < \infty$ ) iff  $(Q_\tau, \lambda_\tau, v_\tau) \in \mathcal{W}$  for all  $\tau \in [0, 1]$ .

## Theorem

$$K(A, M, \lambda_0, \nu_0) \leq C \int_0^1 \mu^2(Q_T, \lambda_T, \nu_T) dT.$$

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**Theorem** We have

$$\mathbb{E}_{A \sim N(0, \text{Id})} \mathbb{E}_{i \sim \{1, \dots, n\}} K(A, M, \lambda_0, v_0) = \mathcal{O}(n^5)$$

and, consequently,

$$\mathbb{E}_{A \sim N(0, \text{Id})} \mathbb{E}_{i \sim \{1, \dots, n\}} \text{cost}(A, M, \lambda_0, v_0) = \mathcal{O}(n^8).$$

One can also (deterministically) compute **all** the eigenpairs of  $A$ .  
The average total cost is  $\mathcal{O}(n^9)$ .

## A finer randomization?

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To do so, it is enough to draw a diagonal matrix whose diagonal elements  $(\lambda_1, \dots, \lambda_n)$  follow the spectrum law of the Ginibre ensemble.

Can we do this in  $\mathcal{O}(n^6)$  operations?