

9. EIGENVALUES: important in engineering applications such as structural engineering (a te example of bridge).
Associated with resonances of system, dynamical systems, stability...

§9.1 Linear Algebra preliminaries

Def (eigenvalue, eigenvector) Let $A \in \mathbb{C}^{n \times n}$. A non-zero vector $x \in \mathbb{C}^n$ is an eigenvector of A and λ its corresponding eigenvalue if

$$Ax = \lambda x$$

Def (Linear independence) Let $\{v_1, v_2, \dots, v_n\}$ be a set of vectors. The set is linearly independent if:

$$\sum_{i=1}^k \alpha_i v_i = 0 \Leftrightarrow \alpha_i = 0, i = 1 \dots k.$$

Theorem (Basis): Let $\{v_1, \dots, v_n\}$ be a set of n lin. indep vectors of \mathbb{R}^n . Then:

$$\forall x \in \mathbb{R}^n \exists \beta_i \text{ s.t. } x = \sum_{i=1}^n \beta_i v_i$$

Proof: $A = [v_1 | v_2 | \dots | v_n] \in \mathbb{R}^{n \times n}$.

$$\{v_1, v_2, \dots, v_n\} \text{ lin indep} \Leftrightarrow \left(A \underline{\alpha} = \sum_{i=1}^n \alpha_i v_i = 0 \Leftrightarrow \underline{\alpha} = 0 \right)$$

$\Leftrightarrow A$ is invertible (nonsingular)

$$\text{Then } \underline{x} = A \underline{\beta}, \text{ where } \underline{\beta} = A^{-1} \underline{x}$$

Def (Dimension) The dimension of a vector subspace is the max. number of linearly independent vectors spanning the set.

Eigenvalue decomposition

For a square matrix $A \in \mathbb{C}^{n \times n}$ the eigenvalue decomposition is a factorization:

$$A = X \Lambda X^{-1}$$

(note: is not always possible!)

$$\Leftrightarrow AX = X \Lambda$$

$$\Leftrightarrow A [x_1 | x_2 | \dots | x_n] = [x_1 | x_2 | \dots | x_n] \begin{bmatrix} \lambda_1 & & \\ & \dots & \\ & & \lambda_n \end{bmatrix}$$

Def (Characteristic polynomial) The characteristic poly of

some $A \in \mathbb{C}^{m \times m}$ is

$$P_A(\lambda) = \det(\lambda I - A) = m \text{ degree poly.}$$

Theorem: λ is an eigenvalue of $A \Leftrightarrow P_A(\lambda) = 0$

proof: λ eigenvalue $\Leftrightarrow \exists x \neq 0$ s.t. $\lambda x - Ax = 0$

$\Leftrightarrow \lambda I - A$ is singular

$$\Leftrightarrow \det(\lambda I - A) = 0$$

This theorem means that even if $A \in \mathbb{R}^n$, spectrum may be complex.
(physically complex eigenvalues give oscillatory behaviour).

Def (Algebraic Multiplicity)

Fundamental theorem of algebra $\Rightarrow P_A(\lambda) = (\lambda - \lambda_1)(\lambda - \lambda_2) \dots (\lambda - \lambda_m)$

The algebraic multiplicity is the multiplicity of λ as a root of p . (ie how many times $\lambda - \lambda$ appears in $P_A(\lambda)$).

Def (Geometric Multiplicity)

Let λ be an eigenvalue of A .

$$E_\lambda = \{x \mid Ax = \lambda x\} = \text{vector space}$$

= eigenspace or invariant subspace of A . $[AE_\lambda \subseteq E_\lambda]$

geometric multiplicity of $\lambda = \dim E_\lambda = \dim \text{null}(\lambda I - A)$.

Theorem: Let $A \in \mathbb{C}^{m \times m}$, A has m eigenvalues counted with algebraic multiplicity. (easy corollary of fundamental theorem of algebra)

Similarity transformations If $X \in \mathbb{C}^m$ is non-singular, then the map $A \rightarrow X^{-1}AX$ is a similarity transformation.

A and B are similar if there is a similarity transf s.t.

$$A = X^{-1}BX. \quad (\sim \text{change of basis})$$

Theorem (similar matrices) if X is nonsingular then A and $X^{-1}AX$ have the same $p_A(z)$, eigenvalues and algebraic (geom.) multiplicities.

Proof:
$$p_{X^{-1}AX}(z) = \det(zI - X^{-1}AX) = \det(X^{-1}(zI - A)X)$$
$$= \det(zI - A) = p_A(z)$$

proves same eigenvalues and algebraic multiplicity

Also if E_λ is an eigenspace for A then $X^{-1}E_\lambda$ is an eigenspace for $X^{-1}AX$ as well.

Theorem. algebraic multiplicity \geq geom. multiplicity. (59)

Matrices for which geom. multiplicity $<$ algebraic mult are called defective.

Example:

$$A = \begin{bmatrix} 2 & \\ & 2 \end{bmatrix} \quad B = \begin{bmatrix} 2 & 1 & \\ & 2 & 1 \\ & & 2 \end{bmatrix} \quad p_A(z) = (z-2)^3$$

$$p_B(z) = (z-2)^3$$

\Rightarrow same eigenvalues and algebraic mult.
($\lambda=2$) (alg. mult of A is 3)

Three lin indep eigenvectors of A are: \Rightarrow geom mult is 3
 $\underline{e}_1, \underline{e}_2, \underline{e}_3$.

For B the only eigenvector possible is \underline{e}_1 . \Rightarrow geom mult is 1.

Theorem A non defective $\Leftrightarrow A = X \Lambda X^{-1}$ (diagonalizable)

Def A matrix Q is unitary if $Q^* Q = I$.

The columns of a unitary matrix are orthogonal.

$$Q^{-1} = Q^*$$

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

$$A = A^* \Rightarrow A = Q \Lambda Q^*, \quad \begin{array}{l} Q \text{ unitary} \\ \Lambda \text{ diagonal} \end{array}$$

(Hermitian) and eigenvalues are real.

Theorem: A is unitarily diagonalizable iff it is normal,

i.e. $A^* A = A A^*$.

Schur Factorization (behind famous QR algo)

$$A = QTQ^* \quad T = (\nabla) \quad Q = \text{unitary}$$

This is an eigenvalue revealing factorization because A and T are similar.
 \Rightarrow eigenvalues of A appear on diagonal of T .

Theorem: $\forall A \in \mathbb{C}^{n \times n}$ admits a Schur factorization

proof: By induction on size of A .

$n=1$ trivial

Assume any matrix of size n has a Schur factorization.

Consider $A \in \mathbb{C}^{(n+1) \times (n+1)}$, let x be an eigenvector of A w/ eigenvalue λ .

$$U = \begin{bmatrix} \frac{x}{\|x\|} & \dots \end{bmatrix} = \text{unitary.}$$

complete basis

Then: $U^* A U = \begin{bmatrix} \lambda & B \\ 0 & C \end{bmatrix}$ By induction hyp: $C = V T V^*$, $T = (\nabla)$
 $V = \text{unitary}$

$$\text{let } Q = U \begin{bmatrix} 1 & 0 \\ 0 & V \end{bmatrix}$$

$$\begin{aligned} \text{Then } Q^* A Q &= \begin{bmatrix} 1 & 0 \\ 0 & V^* \end{bmatrix} U^* A U \begin{bmatrix} 1 & 0 \\ 0 & V \end{bmatrix} \\ &= \begin{bmatrix} 1 & 0 \\ 0 & V^* \end{bmatrix} \begin{bmatrix} \lambda & B \\ 0 & C \end{bmatrix} \begin{bmatrix} 1 & 0 \\ 0 & V \end{bmatrix} \\ &= \begin{bmatrix} \lambda & B \\ 0 & V^* C \end{bmatrix} \begin{bmatrix} 1 & 0 \\ 0 & V \end{bmatrix} \\ &= \begin{bmatrix} \lambda & B V \\ 0 & T \end{bmatrix} = (\nabla). \quad \text{QED} \end{aligned}$$

This is an existence proof! it does not tell us how to compute Schur factor!

Note: all eigenvalue algorithms need to be iterative.

(59)

The reason is that finding eigenvalues \Leftrightarrow finding roots of a poly.

(\Rightarrow) Characteristic poly

(\Leftarrow) Companion matrix of poly:

$$p(z) = z^m + a_{m-1}z^{m-1} + \dots + a_1z + a_0$$

$$A = \begin{bmatrix} 0 & -a_0 \\ 1 & -a_1 \\ & \vdots \\ & 0 & -a_{m-2} \\ & & 1 & -a_{m-1} \end{bmatrix}$$

Can verify that, if z is a root of p :
 $v = (1, z, z^2, \dots, z^{m-1})$ is a left eigenvector
of A with eigenvalue z :

$$vA = zv$$

And it is a well known fact that no general formula exists for computing the roots of a polynomial of degree ≥ 5 , so we can only hope to approximate them through an iterative process.

§ 9.2 Power Method

Power method (or iteration) algorithm

$$v^{(0)} = \text{some vector with } \|v^{(0)}\|_{\infty} = 1 = v_p^{(0)}$$

for $k = 1, 2, \dots$

$$w^{(k)} = Av^{(k-1)}$$

$$v^{(k)} = w^{(k)} / \|w^{(k)}\|_{\infty}$$

$$\lambda^{(k)} = w_p^{(k)}, \text{ where } p = \underset{\text{smallest}}{v} \text{ index s.t. } |v_p^{(k-1)}| = \|v^{(k-1)}\|_{\infty} = 1$$

The iteration stops when two successive iterates become close to within some tolerance.

The power method finds the largest eigenvalue (in magnitude) and its associated eigenvector.

Here is a sketch of the convergence proof.

For simplicity we assume A is diagonalizable. We also assume that the largest eigenvalue (in magnitude) of A is simple.

Let $\lambda_1, \lambda_2, \dots, \lambda_n$ be the eigenvalues of A ordered s.t.

$$|\lambda_1| > |\lambda_2| \geq |\lambda_3| \geq \dots \geq |\lambda_n|.$$

Let $q_1, q_2, q_3, \dots, q_n$ the associated eigenvectors (lin indep)

Then given $v^{(0)}$ (The starting vector) $\exists \beta_i$ s.t.

$$v^{(0)} = \sum_{i=1}^n \beta_i q_i.$$

$$A v^{(0)} = \sum_{i=1}^n \beta_i \lambda_i q_i$$

$$A^k v^{(0)} = \sum_{i=1}^n \beta_i \lambda_i^k q_i = \lambda_1^k \beta_1 q_1 + \lambda_1^k \sum_{i=2}^n \beta_i \left(\frac{\lambda_i}{\lambda_1}\right)^k q_i$$

It's not hard to see that when we apply the power method,

we have:

$$v^{(k)} = \frac{A^k v^{(0)}}{\|A^k v^{(0)}\|} = \frac{\lambda_1^k (\beta_1 q_1 + \sum_{i=2}^n \beta_i (\frac{\lambda_i}{\lambda_1})^k q_i)}{|\lambda_1|^k \|\beta_1 q_1 + \sum_{i=2}^n \beta_i (\frac{\lambda_i}{\lambda_1})^k q_i\|_\infty}$$

Now since $\lim_{k \rightarrow \infty} (\frac{\lambda_i}{\lambda_1})^k = 0, i = 2, \dots, n,$

we have $v^{(k)} \rightarrow$ a multiple of q_1 .

and the rate of convergence is linear with ratio $|\frac{\lambda_2}{\lambda_1}|$.

Drawback of power method: (for both vector & eigenvalue)

- nobody guarantees initial sol has component in direction of largest eigenvector
- computes only largest eigenvalue & eigenvector
• (we will see how to fix this)
- convergence rate: can be very slow if $\lambda_1 \sim \lambda_2$.

- Advantages
- Algorithm works even when A is nondiagonalizable.
 - Simplicity.
 - Uses only matrix vector prod.
- It's possible to get faster linear convergence of the eigenvalues in the case where A is real & symmetric:

Power method (symmetric case)

$v^{(0)}$ = some vector with $\|v^{(0)}\| = 1$

for $k = 1, 2, \dots$

$$\begin{cases} w^{(k)} = A v^{(k-1)} \\ v^{(k)} = w^{(k)} / \|w^{(k)}\| \\ \lambda^{(k)} = (v^{(k)})^T A v^{(k)} \end{cases}$$

faster than in general case.

In this case we can expect: $|\lambda^{(k)} - \lambda_1| = O\left(\left(\frac{\lambda_2}{\lambda_1}\right)^{2k}\right)$

and $|v^{(k)} - q_1| = O\left(\left(\frac{\lambda_2}{\lambda_1}\right)^k\right)$
 depends on sign of multiple of q_1 we converge to.

The reason behind the update in $\lambda^{(k)}$ is the so called Rayleigh quotient:

$$r(x) = \frac{x^T A x}{x^T x}$$

For real symmetric matrices it can be shown that:

$$\lambda_{\min} \leq r(x) \leq \lambda_{\max}$$

So if $v^{(k)}$ is a good approx to q_1 , then:

$$v^{(k)T} A v^{(k)} = \frac{v^{(k)T} A v^{(k)}}{v^{(k)T} v^{(k)}} \approx \frac{q_1^T A q_1}{q_1^T q_1} = \lambda_1 \frac{q_1^T q_1}{q_1^T q_1}$$

The power method can be modified to zoom in and find the closest eigenvalue to some μ . The idea is that:

if v is an eigenvector of A with eigenvalue λ then $v \frac{(A - \mu I)^{-1}}{(\lambda - \mu)^{-1}}$

So if we apply power method to matrix $(A - \mu I)^{-1}$ the method will converge to the largest eigenvalue $|\lambda_j - \mu|^{-1}$, a λ the closest to μ . This method is:

Inverse iteration

$v^{(0)}$ = some vector with $\|v^{(0)}\| = 1$

for $k = 1, 2, \dots$

solve $(A - \mu I)w^{(k)} = v^{(k-1)}$ for $w^{(k)}$ (= apply $(A - \mu I)^{-1}$)

$v^{(k)} = \frac{w^{(k)}}{\|w^{(k)}\|}$

$\lambda^{(k)} = v^{(k)T} A v^{(k)}$

(here we wrote the version for symmetric matrices, but the same modification can be carried out in the original method)

Convergence is still linear, but once we control μ , we can control the convergence rate.

Suppose λ_{j_1} is the closest eigenvalue of A to μ

λ_{j_2} next to closest

$|\lambda_{j_1} - \mu| < |\lambda_{j_2} - \mu| \leq |\lambda_j - \mu|$ for $j \neq j_1$

then:

$|\lambda^{(k)} - \lambda_{j_1}| = O\left(\left|\frac{\mu - \lambda_{j_1}}{\mu - \lambda_{j_2}}\right|^{2k}\right)$

and

$$\|v^{(k)} - (\pm q_1)\| = O\left(\left|\frac{\mu - \lambda_{j_1}}{\mu - \lambda_{j_2}}\right|^k\right)$$

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So why not use rowable shifts μ that get closer to the eigenvalue of interest?

This gives the following algo:

Rayleigh quotient iteration

$v^{(0)}$ = some vector with $\|v^{(0)}\| = 1$

$\lambda^{(0)} = v^{(0)T} A v^{(0)}$ = Rayleigh quot.

for $k = 1, 2, \dots$

$$\begin{cases} \text{solve } (A - \lambda^{(k-1)} I) w = v^{(k-1)} & (\text{apply } (A - \lambda^{(k-1)} I)^{-1}) \\ v^{(k)} = w / \|w\| \\ \lambda^{(k)} = v^{(k)T} A v^{(k)} & (\text{update guess with Rayleigh quot}) \end{cases}$$

Convergence of RQI is one of the rare cases where one gets cubic convergence!

If the start vector $v^{(0)}$ is sufficiently close to the eigenvector q_j :

$$\|v^{(k+1)} - (\pm q_j)\| = O(\|v^{(k)} - (\pm q_j)\|^3)$$

$$|\lambda^{(k+1)} - \lambda_j| = O(|\lambda^{(k)} - \lambda_j|^3)$$

§ 9.3 Reduction to Hessenberg or tridiagonal form

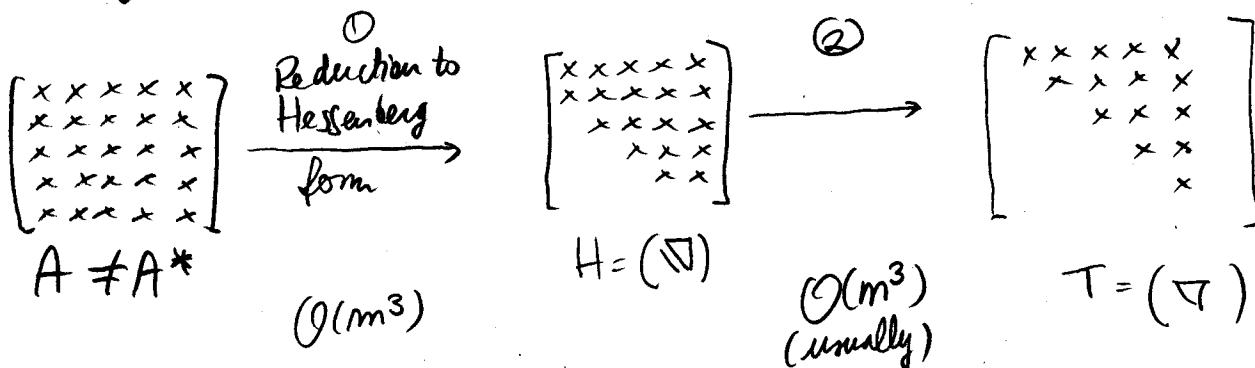
(64)

Motivation: We would like to somehow compute Schur facts of A by successive multiplications by orthogonal matrices: $(AQ = Q^T)$

$$T_j = Q_j^* Q_{j-1}^* \dots Q_2^* Q_1^* A Q_1 Q_2 \dots Q_j$$

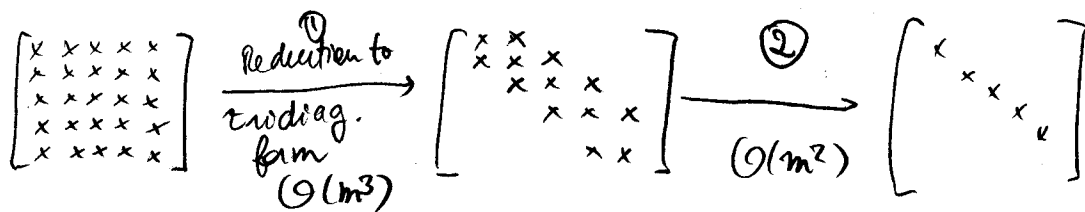
so that $T_j \rightarrow T = (\nabla)$ as $j \rightarrow \infty$.

Usually this is split into 2 phases:



② could run forever, however convergence is usually obtained in m steps and each step in ② takes $O(m^2)$ flops.

If $A = A^*$ (A Hermitian)



② each step takes m flops (if no eigenvectors are needed)
 $\sim m$ steps total = $O(m^2)$ flops

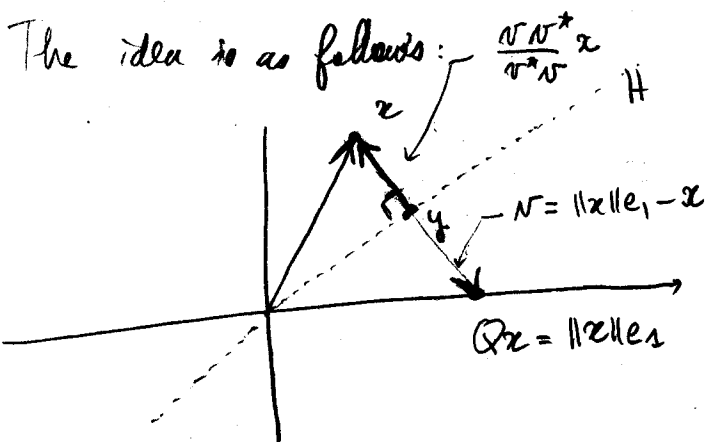
In Burden & Faires: ① = § 9.3 Householder's Method
 ② = § 9.4 The QR algorithm

Phase ② can be made using orthogonal transformations called Householder reflectors (which are extremely useful in linear algebra ...)

Householder reflectors are orthogonal transformations that zero out part of a vector.

Given $x = \begin{bmatrix} x \\ x \\ x \\ \vdots \\ x \end{bmatrix}$ we can build a Q with $Q^*Q = I$ s.t.

$$Qx = \begin{bmatrix} \|x\| \\ 0 \\ \vdots \\ 0 \end{bmatrix} = \|x\| e_1.$$



The \perp proj of x onto v is:

$$\frac{v v^* x}{v^* v}$$

Thus:

$$y = x - \frac{v v^* x}{v^* v}$$

and:

$$Qx = x - 2 \frac{v v^* x}{v^* v}$$

Another geometric interpretation:

y is projection \perp of x onto H

Qx is reflection of x across H .

It's amazing that a simple factor of 2 converts a projector (not full rank) into an orthogonal matrix!

$$Q = I - 2 \frac{v v^*}{v^* v}$$

Q is orthogonal.

$$Q^* Q = \left(I - 2 \frac{v v^*}{v^* v} \right) \left(I - 2 \frac{v v^*}{v^* v} \right)$$

$$= I - 4 \frac{v v^*}{v^* v} + 4 \left(\frac{v v^*}{v^* v} \right) \left(\frac{v v^*}{v^* v} \right) = I.$$

More generally we can introduce 0 anywhere in the vector.

A useful pattern is:

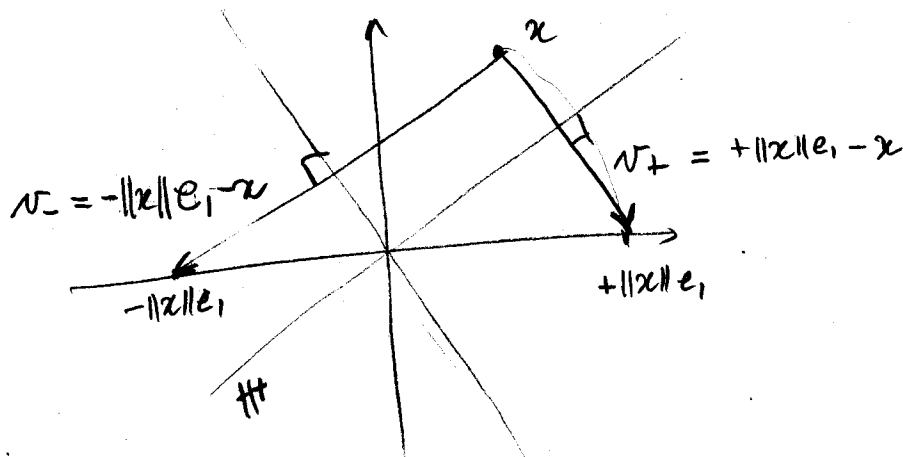
$$Q_k = \begin{bmatrix} 1 & 0 \\ 0 & Q \end{bmatrix}, \text{ where } Q \text{ is a Householder refl.}$$

Q_k will leave untouched the first m entries:

$$x = \begin{bmatrix} \left. \begin{matrix} x \\ x \\ x \end{matrix} \right\}^m \\ \left. \begin{matrix} x \\ x \\ x \end{matrix} \right\}^{m-m} \end{bmatrix} \longrightarrow Q_k x = \begin{bmatrix} x \\ x \\ x \\ 0 \\ \vdots \\ 0 \end{bmatrix} \text{ unchanged}$$

There is a slight detail w. th HH reflectors: we could have chosen Q s.t. $Qx = -\|x\|e_1$, so which HH refl is better:

$Qx = +\|x\|e_1$ or the former?



To avoid numerical cancellation in first component one usually takes $v = -\text{sgn}(x_1)\|x\|e_1 - x$

But since $Q = I - 2 \frac{vv^*}{v^*v}$ we can eliminate the $-$ sign:

$$v = \text{sgn}(x_1)\|x\|e_1 - x$$

Note • each application of HH refl Q (or Q^*) is $O(m^2)$.

- we don't need to store Q , only v . Therefore successive applications of different HH refl Q_1, Q_2, \dots, Q_j can be stored with j vectors only.

Back to original problem:

$$\begin{matrix}
 \begin{bmatrix} x & x & x & x & x \\ x & x & x & x & x \\ x & x & x & x & x \\ x & x & x & x & x \\ x & x & x & x & x \end{bmatrix} & \longrightarrow & \begin{bmatrix} x & x & x & x & x \\ x & x & x & x & x \\ & x & x & x & x \\ & & x & x & x \\ & & & x & x \end{bmatrix} \\
 A & & H
 \end{matrix}$$

how do we introduce zeros?

A first idea:

$$Q_1^* A = \begin{bmatrix} x & x & x & x & x \\ 0 & x & x & x & x \\ 0 & x & x & x & x \\ 0 & x & x & x & x \\ 0 & x & x & x & x \end{bmatrix} \xrightarrow{\cdot Q_1} \begin{bmatrix} x & x & x & x & x \\ x & x & x & x & x \\ x & x & x & x & x \\ x & x & x & x & x \\ x & x & x & x & x \end{bmatrix}$$

$$Q_1^* A Q_1$$

each col is lin comb of columns of $Q_1^* A$ so zeros disappear.

does not work!

A better idea: be less ambitious: leaves unchanged

$$\begin{matrix}
 \begin{bmatrix} x & x & x & x & x \\ x & x & x & x & x \\ x & x & x & x & x \\ x & x & x & x & x \\ x & x & x & x & x \end{bmatrix} & \xrightarrow{Q_1^*} & \begin{bmatrix} x & x & x & x & x \\ x & x & x & x & x \\ 0 & x & x & x & x \\ 0 & x & x & x & x \\ 0 & x & x & x & x \end{bmatrix} & \xrightarrow{\cdot Q_1} & \begin{bmatrix} x & x & x & x & x \\ x & x & x & x & x \\ 0 & x & x & x & x \\ 0 & x & x & x & x \\ 0 & x & x & x & x \end{bmatrix} \\
 A & & Q_1^* A & & Q_1^* A Q_1
 \end{matrix}$$

unchanged

$$\begin{matrix}
 \begin{bmatrix} x & x & x & x & x \\ x & x & x & x & x \\ x & & & & \\ & & & & \\ & & & & \end{bmatrix} & \xleftarrow{\cdot Q_2} & \begin{bmatrix} x & x & x & x & x \\ x & x & x & x & x \\ & x & x & x & x \\ & 0 & x & x & x \\ & 0 & x & x & x \end{bmatrix} \\
 & & Q_2^* Q_1^* A Q_1
 \end{matrix}$$

etc...

at the end we get (after $m-2$ steps)

$$H = \begin{bmatrix} x & x & x & x & x \\ x & x & x & x & x \\ & x & x & x & x \\ & & x & x & x \\ & & & x & x \end{bmatrix} = Q_{m-2}^* Q_{m-1}^* \dots Q_1^* A Q_1 Q_2 \dots Q_{m-2}$$

Householder reduction to Hessenberg form

for $k = 1 \dots m-2$

$$x = A(k+1:m, k)$$

$$v_k = \text{sign}(x(1)) \|x\|_2 e_1 + x$$

$$v_k = v_k / \|v_k\|_2$$

$$A(k+1:m, k:m) = A(k+1:m, k:m) - 2 v_k (v_k^* A(k+1:m, k:m))$$

$$A(1:m, k+1:m) = A(1:m, k+1:m) - 2 (A(1:m, k+1:m) v_k) v_k^*$$

When $A = A^*$ Hermitian, we can do similar procedure -

$Q_1^* A Q_1$ will also be symmetric so any zeros we introduced in columns of A will also appear at zeros in corresp. row.

Applying HD of l on both sides has exactly same cost, so reduction to tridiagonal form is relatively cheap.

The QR factorization [We assume symm real matrices in following] (69)

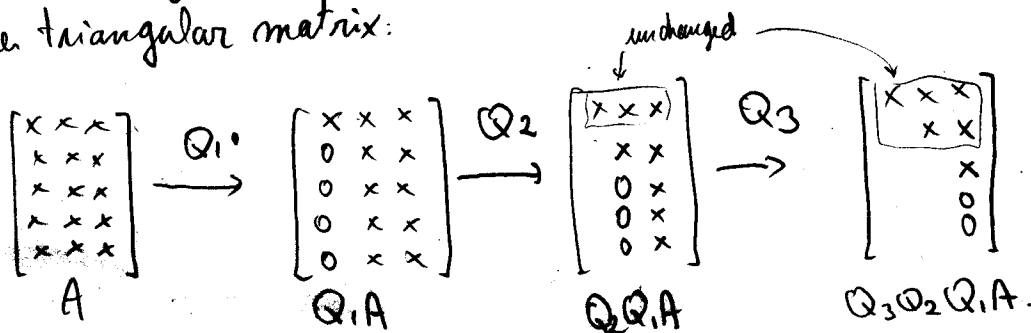
One other important matrix factorization that can be easily obtained with Householder reflectors is the QR factor:

let $A \in \mathbb{R}^{m \times n}$, then $\exists Q$ unitary and $R (\nabla)$ s.t.
 $(Q^*Q = I)$

$$A = QR$$

$m \times m$ $m \times n$ $n \times n$

Idea: apply Householder reflectors on the left of A to obtain upper triangular matrix:



we will not see details. Keep in mind that in Matlab:

$[Q, R] = qr(A, 0)$ gives such a factorization.

Note: if A is full column rank (i.e. columns are lin indep) then $\text{Range}(Q) = \text{Range}(A)$.

Unnormalized Simultaneous iteration

Idea: apply power method to many vectors at once.

If $A^k v^{(0)} \rightarrow q_1^{(k)}$ we can expect: $\text{span}\{A^k v_1^{(0)}, A^k v_2^{(0)}, \dots, A^k v_n^{(0)}\}$
 $\rightarrow \text{span}\{q_1, q_2, \dots, q_n\}$
 (power method) (simultaneous iteration) (a block power method)
 largest eigenvalue in magnitude largest n eigenvalues in magnitude

In matrix notation:

$$V^{(0)} = \begin{bmatrix} | & | & & | \\ v_1^{(0)} & v_2^{(0)} & \dots & v_m^{(0)} \\ | & | & & | \end{bmatrix}$$

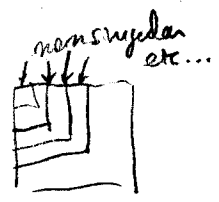
$$V^{(k)} = A^k V^{(0)} = \begin{bmatrix} | & | & & | \\ v_1^{(k)} & v_2^{(k)} & \dots & v_m^{(k)} \\ | & | & & | \end{bmatrix}$$

Since all we want is a basis for Range ($V^{(k)}$) (or column space) we use QR factor:

$$\tilde{Q}^{(k)} \tilde{R}^{(k)} = V^{(k)}$$

If we assume:

- $| \lambda_1 | > | \lambda_2 | > \dots > | \lambda_n | > | \lambda_{n+1} | \geq \dots \geq | \lambda_m |$
 - all leading principal submatrices of $\tilde{Q}^T V^{(0)}$ are non singular
- where $\tilde{Q} = [q_1, q_2, \dots, q_m]$ \uparrow m = matrix with eigenvectors
corresp. to $\lambda_1, \dots, \lambda_m$ as columns.



this condition is similar to condition in power method that initial vector has a component in direction of "largest" eigenvector.

Then it can be shown that:

$$\| q_j^{(k)} - \oplus q_j \| = O(C^k)$$

where $C = \max_{1 \leq k \leq m} \frac{|\lambda_{k+1}|}{|\lambda_k|}$

(convergence ratio is determined by the eigenvalues that are the closest.)

Simultaneous iteration

Obviously if we keep multiplying by A , we can have components that quickly become large in magnitude.

→ as in power method we need to renormalize every time we apply A :

Pick $\tilde{Q}^{(0)} \in \mathbb{R}^{m \times n}$ with orthonormal cols (e.g. from QR factor of a random matrix)

for $k = 1, 2, \dots$

$$\begin{cases} Z = A \tilde{Q}^{(k-1)} \\ \tilde{Q}^{(k)} \tilde{R}^{(k)} = Z \end{cases}$$

It is not hard to show that $\tilde{Q}^{(k)}$ is precisely same matrix for which $\tilde{Q}^{(k)} R = A^k \tilde{Q}^{(0)}$ (provided reuse same initial guess)

Thus convergence study is the same.

QR algorithm (pure version, seldom implemented as is)

$$A^{(0)} = A$$

for $k = 1, 2, \dots$

$$\begin{cases} Q^{(k)} R^{(k)} = A^{(k-1)} \\ A^{(k)} = R^{(k)} Q^{(k)} \end{cases}$$

factors are recombined in reverse order.

Equivalence of QR algorithm and Simultaneous iteration (72)

Of course this equivalence could only happen if we consider simultaneous iteration on all eigenvectors ($m = n$)

To make equivalence obvious we rewrite algorithms:

Simultaneous iteration

$$\begin{aligned} \tilde{Q}^{(0)} &= I; A^{(0)} = A \\ \text{for } k &= 1, 2, \dots \\ Z &= A \tilde{Q}^{(k-1)} \\ \tilde{Q}^{(k)} \tilde{R}^{(k)} &= Z \\ A^{(k)} &= \tilde{Q}^{(k)T} A \tilde{Q}^{(k)} \end{aligned}$$

QR algorithm

$$\begin{aligned} A^{(0)} &= A; \tilde{Q}^{(0)} = I \\ \text{for } k &= 1, 2, \dots \\ Q^{(k)} R^{(k)} &= A^{(k-1)} \\ A^{(k)} &= R^{(k)} Q^{(k)} \\ \tilde{Q}^{(k)} &= Q^{(1)} Q^{(2)} \dots Q^{(k)} \end{aligned}$$

We will show by induction that both algorithms generate the same matrices $\tilde{Q}^{(k)}$, $A^{(k)}$ and $\tilde{R}^{(k)} \equiv R^{(k)} R^{(k-1)} \dots R^{(1)}$,

with:

$$i) \quad A^k = \tilde{Q}^{(k)} \tilde{R}^{(k)}$$

$$ii) \quad A^{(k)} = (\tilde{Q}^{(k)})^T A \tilde{Q}^{(k)}$$

proof: $k=0$ trivial since for both methods $\tilde{Q}^{(0)} = I$, $A^{(0)} = A$, $R^{(0)} = I$
 $I = A^0 = \tilde{Q}^{(0)} \tilde{R}^{(0)} = I I$.

• case $k \geq 1$ for Simult. iter.

$$A^k = A A^{k-1} \underset{\substack{\uparrow \\ \text{induction } i)}}{=} A \tilde{Q}^{(k-1)} \tilde{R}^{(k-1)} \underset{\substack{\uparrow \\ \text{iteration } i}}{=} \tilde{Q}^{(k)} R^{(k)} \tilde{R}^{(k-1)} = \tilde{Q}^{(k)} \tilde{R}^{(k)}$$

• Case $k \geq 1$ for QR algo.

$$A^k = A A^{k-1} \underset{\substack{\uparrow \\ \text{induction } i)}}{=} A \tilde{Q}^{(k-1)} \tilde{R}^{(k-1)} \underset{\substack{\uparrow \\ \text{iteration } i}}{=} \tilde{Q}^{(k-1)} A^{(k-1)} \tilde{R}^{(k-1)} \underset{\substack{\uparrow \\ \text{iteration } ii}}{=} \tilde{Q}^{(k-1)} Q^{(k)} R^{(k)} \tilde{R}^{(k-1)} \underset{\substack{\uparrow \\ \text{iteration } i}}{=} \tilde{Q}^{(k)} \tilde{R}^{(k)}$$

Finally to show ii) for QR algo:

$$\begin{aligned}
A^{(k)} &= R^{(k)} Q^{(k)} \stackrel{\substack{\uparrow \\ \text{QR algo}}}{=} Q^{(k)T} A^{(k-1)} Q^{(k)} \\
&= \underbrace{(Q^{(k)})^T}_{\substack{\uparrow \\ \text{induction}}} (\tilde{Q}^{(k-1)})^T A \tilde{Q}^{(k-1)} Q^{(k)} \\
&= \tilde{Q}^{(k)T} A \tilde{Q}^{(k)} \quad \text{qed.}
\end{aligned}$$

Here is how QR algorithm can be expected to converge.

- QR constructs orthonormal basis of $A^k \rightarrow$ finds eigenvectors
- diagonal elements of $A^{(k)}$ are Rayleigh quotients:

$$A^{(k)} = (Q^{(k)})^T A Q^{(k)}$$

so they should \rightarrow eigenvalues

- strictly triangular part are "generalized" Rayleigh quotients. Since eigenvectors are \perp , these should $\rightarrow 0$.

We are not done with QR. A more practical version would be as follows: Shifted QR

$$A^{(0)} = Q^{(0)} A (Q^{(0)})^T = \left(\begin{array}{c} \equiv \\ \equiv \\ \equiv \end{array} \right) \quad (\text{Reduction to tridiag form})$$

for $k=1, 2, \dots$

pick shift $\mu^{(k)}$ (usually $\mu^{(k)} = A_{mm}^{(k-1)}$)

$$Q^{(k)} R^{(k)} = A^{(k-1)} - \mu^{(k)} I$$

$$A^{(k)} = R^{(k)} Q^{(k)} + \mu^{(k)} I$$

"deflation" to lock in eigenvalues that have converged