

## LARGE SPARSE EIGENVALUE PROBLEMS

- Projection methods
- The subspace iteration
- Krylov subspace methods: Arnoldi and Lanczos
- Golub-Kahan-Lanczos bidiagonalization

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## General Tools for Solving Large Eigen-Problems

- Projection techniques – Arnoldi, Lanczos, Subspace Iteration;
- Preconditionings: shift-and-invert, Polynomials, ...
- Deflation and restarting techniques
- Computational codes often combine these three ingredients

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## A few popular solution Methods

- Subspace Iteration [Now less popular – sometimes used for validation]
- Arnoldi's method (or Lanczos) with polynomial acceleration
- Shift-and-invert and other preconditioners. [Use Arnoldi or Lanczos for  $(A - \sigma I)^{-1}$ .]
- Davidson's method and variants, Jacobi-Davidson
- Specialized method: Automatic Multilevel Substructuring (AMLS).

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## Projection Methods for Eigenvalue Problems

### Projection method onto $K$ orthogonal to $L$

- Given: Two subspaces  $K$  and  $L$  of same dimension.
- Approximate eigenpairs  $\tilde{\lambda}, \tilde{u}$ , obtained by solving:

Find:  $\tilde{\lambda} \in \mathbb{C}, \tilde{u} \in K$  such that  $(\tilde{\lambda}I - A)\tilde{u} \perp L$

- Two types of methods:

Orthogonal projection methods: Situation when  $L = K$ .

Oblique projection methods: When  $L \neq K$ .

- First situation leads to Rayleigh-Ritz procedure

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## Rayleigh-Ritz projection

**Given:** a subspace  $X$  known to contain good approximations to eigenvectors of  $A$ .

**Question:** How to extract 'best' approximations to eigenvalues/eigenvectors from this subspace?

**Answer:** Orthogonal projection method

- Let  $Q = [q_1, \dots, q_m]$  = orthonormal basis of  $X$
- Orthogonal projection method onto  $X$  yields:

$$Q^H(A - \tilde{\lambda}I)\tilde{u} = 0 \rightarrow$$

- $Q^H A Q y = \tilde{\lambda} y$  where  $\tilde{u} = Q y$
- Known as Rayleigh Ritz process

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## Procedure:

1. Obtain an orthonormal basis of  $X$
2. Compute  $C = Q^H A Q$  (an  $m \times m$  matrix)
3. Obtain Schur factorization of  $C$ ,  $C = Y R Y^H$
4. Compute  $\tilde{U} = Q Y$

**Property:** if  $X$  is (exactly) invariant, then procedure will yield exact eigenvalues and eigenvectors.

**Proof:** Since  $X$  is invariant,  $(A - \tilde{\lambda}I)u = Qz$  for a certain  $z$ .  $Q^H Q z = 0$  implies  $z = 0$  and therefore  $(A - \tilde{\lambda}I)u = 0$ .

- Can use this procedure in conjunction with the subspace obtained from subspace iteration algorithm

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## Subspace Iteration

**Original idea:** projection technique onto a subspace of the form  $Y = A^k X$

Practically:  $A^k$  replaced by suitable polynomial

Advantages:

- Easy to implement (in symmetric case);
- Easy to analyze;

Disadvantage: Slow.

- Often used with polynomial acceleration:  $A^k X$  replaced by  $C_k(A)X$ . Typically  $C_k =$  Chebyshev polynomial.

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## Algorithm: Subspace Iteration with Projection

1. **Start:** Choose an initial system of vectors  $X = [x_0, \dots, x_m]$  and an initial polynomial  $C_k$ .
2. **Iterate:** Until convergence do:
  - (a) Compute  $\hat{Z} = C_k(A)X$ . [Simplest case:  $\hat{Z} = AX$ .]
  - (b) Orthonormalize  $\hat{Z}$ :  $[Z, R_Z] = qr(\hat{Z}, 0)$
  - (c) Compute  $B = Z^H A Z$
  - (d) Compute the Schur factorization  $B = Y R_B Y^H$  of  $B$
  - (e) Compute  $X := Z Y$ .
  - (f) Test for convergence. If satisfied stop. Else select a new polynomial  $C_{k'}$  and continue.

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**THEOREM:** Let  $S_0 = \text{span}\{x_1, x_2, \dots, x_m\}$  and assume that  $S_0$  is such that the vectors  $\{Px_i\}_{i=1, \dots, m}$  are linearly independent where  $P$  is the spectral projector associated with  $\lambda_1, \dots, \lambda_m$ . Let  $\mathcal{P}_k$  the orthogonal projector onto the subspace  $S_k = \text{span}\{X_k\}$ . Then for each eigenvector  $u_i$  of  $A$ ,  $i = 1, \dots, m$ , there exists a unique vector  $s_i$  in the subspace  $S_0$  such that  $Ps_i = u_i$ . Moreover, the following inequality is satisfied

$$\|(I - \mathcal{P}_k)u_i\|_2 \leq \|u_i - s_i\|_2 \left( \left| \frac{\lambda_{m+1}}{\lambda_i} \right| + \epsilon_k \right)^k, \quad (1)$$

where  $\epsilon_k$  tends to zero as  $k$  tends to infinity.

## KRYLOV SUBSPACE METHODS

### Krylov subspace methods

**Principle:** Projection methods on Krylov subspaces:

$$K_m(A, v_1) = \text{span}\{v_1, Av_1, \dots, A^{m-1}v_1\}$$

- The most important class of projection methods [for linear systems and for eigenvalue problems]
- Variants depend on the subspace  $L$
- Let  $\mu = \text{deg. of minimal polynom. of } v_1$ . Then:
  - $K_m = \{p(A)v_1 | p = \text{polynomial of degree } \leq m - 1\}$
  - $K_m = K_\mu$  for all  $m \geq \mu$ . Moreover,  $K_\mu$  is invariant under  $A$ .
  - $\dim(K_m) = m$  iff  $\mu \geq m$ .

### Arnoldi's algorithm

- Goal: to compute an orthogonal basis of  $K_m$ .
- Input: Initial vector  $v_1$ , with  $\|v_1\|_2 = 1$  and  $m$ .

#### ALGORITHM : 1. Arnoldi's procedure

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For  $j = 1, \dots, m$  do
  Compute  $w := Av_j$ 
  For  $i = 1, \dots, j$ , do
     $\left\{ \begin{array}{l} h_{i,j} := (w, v_i) \\ w := w - h_{i,j}v_i \end{array} \right.$ 
   $h_{j+1,j} = \|w\|_2$ 
   $v_{j+1} = w/h_{j+1,j}$ 
End
    
```

- Based on Gram-Schmidt procedure

## Result of Arnoldi's algorithm

$$\text{Let: } \overline{H}_m = \begin{pmatrix} x & x & x & x & x \\ x & x & x & x & x \\ & x & x & x & x \\ & & x & x & x \\ & & & x & x \\ & & & & x \end{pmatrix}, H_m = \begin{pmatrix} x & x & x & x & x \\ x & x & x & x & x \\ & x & x & x & x \\ & & x & x & x \\ & & & x & x \end{pmatrix}$$

### Results:

- $V_m = [v_1, v_2, \dots, v_m]$  orthonormal basis of  $K_m$ .
- $AV_m = V_{m+1}\overline{H}_m = V_m H_m + h_{m+1,m}v_{m+1}e_m^T$
- $V_m^T AV_m = H_m \equiv \overline{H}_m$  - last row.

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## Application to eigenvalue problems

- Write approximate eigenvector as  $\tilde{u} = V_m y$

- Galerkin condition:

$$(A - \tilde{\lambda}I)V_m y \perp \mathcal{K}_m \rightarrow V_m^H (A - \tilde{\lambda}I)V_m y = 0$$

- Approximate eigenvalues are eigenvalues of  $H_m$

$$H_m y_j = \tilde{\lambda}_j y_j$$

- Associated approximate eigenvectors are

$$\tilde{u}_j = V_m y_j$$

- Typically a few of the outermost eigenvalues will converge first.

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## Hermitian case: The Lanczos Algorithm

- The Hessenberg matrix becomes tridiagonal :

$$A = A^H \text{ and } V_m^H AV_m = H_m \rightarrow H_m = H_m^H$$

- Denote  $H_m$  by  $T_m$  and  $\overline{H}_m$  by  $\overline{T}_m$ . We can write

$$T_m = \begin{pmatrix} \alpha_1 & \beta_2 & & & & \\ \beta_2 & \alpha_2 & \beta_3 & & & \\ & \beta_3 & \alpha_3 & \beta_4 & & \\ & & & \cdot & \cdot & \cdot \\ & & & & \cdot & \cdot \\ & & & & & \beta_m & \alpha_m \end{pmatrix}$$

- Relation  $AV_m = V_{m+1}\overline{T}_m$

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- Consequence: three term recurrence

$$\beta_{j+1}v_{j+1} = Av_j - \alpha_j v_j - \beta_j v_{j-1}$$

### ALGORITHM : 2. Lanczos

- Choose an initial  $v_1$  with  $\|v_1\|_2 = 1$ ;  
Set  $\beta_1 \equiv 0, v_0 \equiv 0$
- For  $j = 1, 2, \dots, m$  Do:
- $w_j := Av_j - \beta_j v_{j-1}$
- $\alpha_j := (w_j, v_j)$
- $w_j := w_j - \alpha_j v_j$
- $\beta_{j+1} := \|w_j\|_2$ . If  $\beta_{j+1} = 0$  then Stop
- $v_{j+1} := w_j / \beta_{j+1}$
- EndDo

Hermitian matrix + Arnoldi  $\rightarrow$  Hermitian Lanczos

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- In theory  $v_i$ 's defined by 3-term recurrence are orthogonal.
- However: in practice severe loss of orthogonality;

Observation [Paige, 1981]: Loss of orthogonality starts suddenly, when the first eigenpair has converged. It is a sign of loss of linear independence of the computed eigenvectors. When orthogonality is lost, then several the copies of the same eigenvalue start appearing.

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## Reorthogonalization

- Full reorthogonalization – reorthogonalize  $v_{j+1}$  against all previous  $v_i$ 's every time.
- Partial reorthogonalization – reorthogonalize  $v_{j+1}$  against all previous  $v_i$ 's only when needed [Parlett & Simon]
- Selective reorthogonalization – reorthogonalize  $v_{j+1}$  against computed eigenvectors [Parlett & Scott]
- No reorthogonalization – Do not reorthogonalize - but take measures to deal with 'spurious' eigenvalues. [Cullum & Willoughby]

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## Lanczos Bidiagonalization

- We now deal with rectangular matrices. Let  $A \in \mathbb{R}^{m \times n}$ .

### ALGORITHM : 3. Golub-Kahan-Lanczos

1. Choose an initial  $v_1$  with  $\|v_1\|_2 = 1$ ;  
Set  $\beta_0 \equiv 0, u_0 \equiv 0$
2. For  $k = 1, \dots, p$  Do:
3.  $\hat{u} := Av_k - \beta_{k-1}u_{k-1}$
4.  $\alpha_k = \|\hat{u}\|_2$ ;  $u_k = \hat{u}/\alpha_k$
5.  $\hat{v} = A^T u_k - \alpha_k v_k$
6.  $\beta_k = \|\hat{v}\|_2$ ;  $v_{k+1} := \hat{v}/\beta_k$
7. EndDo

Let:  $V_{p+1} = [v_1, v_2, \dots, v_{p+1}] \in \mathbb{R}^{n \times (p+1)}$   
 $U_p = [u_1, u_2, \dots, u_p] \in \mathbb{R}^{m \times p}$

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Let:

$$B_p = \begin{bmatrix} \alpha_1 & \beta_1 & & & & \\ & \alpha_2 & \beta_2 & & & \\ & & \ddots & \ddots & & \\ & & & \ddots & \ddots & \\ & & & & \ddots & \ddots \\ & & & & & \alpha_p & \beta_p \end{bmatrix};$$

- $\hat{B}_p = B_p(:, 1:p)$
- $V_p = [v_1, v_2, \dots, v_p] \in \mathbb{R}^{n \times p}$


Result:

- $V_{p+1}^T V_{p+1} = I$
- $U_p^T U_p = I$
- $AV_p = U_p \hat{B}_p$
- $A^T U_p = V_{p+1} B_p^T$

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- Observe that : 
$$A^T(AV_p) = A^T(U_p\hat{B}_p) = V_{p+1}B_p^T\hat{B}_p$$
  - $B_p^T\hat{B}_p$  is a (symmetric) tridiagonal matrix of size  $(p+1) \times p$
  - Call this matrix  $\bar{T}_p$ . Then:  $(A^T A)V_p = V_{p+1}\bar{T}_p$
  - Standard Lanczos relation !
  - Algorithm is equivalent to standard Lanczos applied to  $A^T A$ .
  - Similar result for the  $u_i$ 's [involves  $AA^T$ ]
-  Work out the details: What are the entries of  $\bar{T}_p$  relative to those of  $B_p$ ?