

Symmetric Positive Definite (SPD) Matrix

- An SPD matrix acts sort of like a positive number.
- A is *symmetric* if $a_{ij} = a_{ji}$, for all i and j .
- Several equivalent conditions for A to be *positive definite*:
 - All eigenvalues are > 0
 - LU factorization without pivoting succeeds, and all pivots are > 0
 - For every nonzero vector x , the number $x^T A x > 0$
- SPD matrices come up a lot in scientific computing & data analysis!
- The temperature matrix is SPD.

Conjugate gradient iteration

- An *iterative* algorithm to solve $Ax = b$.
- Start with a guess $x^{(0)}$, then compute $x^{(1)}$, $x^{(2)}$,
- Stop when you think you're close enough.
- In theory, CG can be used to solve *any* system $Ax = b$, provided only that A is SPD.
- In practice, how well CG works depends on specifics of A in subtle ways, involving eigenvalues and condition number.

Conjugate gradient iteration for $Ax = b$

$x^{(0)} = 0$ approximate solution

$r^{(0)} = b$ residual = $b - Ax$

$d^{(0)} = r^{(0)}$ search direction

for $k = 1, 2, 3, \dots :$

$x^{(k)} = x^{(k-1)} + \dots$ new approx solution

$r^{(k)} = \dots$ new residual

$d^{(k)} = \dots$ new search direction

Conjugate gradient iteration for $Ax = b$

$x^{(0)} = 0$ approximate solution

$r^{(0)} = b$ residual = $b - Ax$

$d^{(0)} = r^{(0)}$ search direction

for $k = 1, 2, 3, \dots :$

$\alpha^{(k)} = \dots$ step length

$x^{(k)} = x^{(k-1)} + \alpha^{(k)} d^{(k-1)}$ new approx solution

$r^{(k)} = \dots$ new residual

$d^{(k)} = \dots$ new search direction

Conjugate gradient iteration for $Ax = b$

$x^{(0)} = 0$ approximate solution

$r^{(0)} = b$ residual = $b - Ax$

$d^{(0)} = r^{(0)}$ search direction

for $k = 1, 2, 3, \dots$:

$\alpha^{(k)} = (r^{(k-1)T} r^{(k-1)}) / (d^{(k-1)T} A d^{(k-1)})$ step length

$x^{(k)} = x^{(k-1)} + \alpha^{(k)} d^{(k-1)}$ new approx solution

$r^{(k)} = \dots$ new residual

$d^{(k)} = \dots$ new search direction

Conjugate gradient iteration for $Ax = b$

$x^{(0)} = 0$ approximate solution

$r^{(0)} = b$ residual = $b - Ax$

$d^{(0)} = r^{(0)}$ search direction

for $k = 1, 2, 3, \dots :$

$\alpha^{(k)} = (r^{(k-1)T} r^{(k-1)}) / (d^{(k-1)T} A d^{(k-1)})$ step length

$x^{(k)} = x^{(k-1)} + \alpha^{(k)} d^{(k-1)}$ new approx solution

$r^{(k)} = \dots$ new residual

$\beta^{(k)} = (r^{(k)T} r^{(k)}) / (r^{(k-1)T} r^{(k-1)})$

$d^{(k)} = r^{(k)} + \beta^{(k)} d^{(k-1)}$ new search direction

Conjugate gradient iteration for $Ax = b$

$x^{(0)} = 0$ approximate solution

$r^{(0)} = b$ residual = $b - Ax$

$d^{(0)} = r^{(0)}$ search direction

for $k = 1, 2, 3, \dots$:

$\alpha^{(k)} = (r^{(k-1)T} r^{(k-1)}) / (d^{(k-1)T} A d^{(k-1)})$ step length

$x^{(k)} = x^{(k-1)} + \alpha^{(k)} d^{(k-1)}$ new approx solution

$r^{(k)} = r^{(k-1)} - \alpha^{(k)} A d^{(k-1)}$ new residual

$\beta^{(k)} = (r^{(k)T} r^{(k)}) / (r^{(k-1)T} r^{(k-1)})$

$d^{(k)} = r^{(k)} + \beta^{(k)} d^{(k-1)}$ new search direction

Conjugate gradient iteration to solve $A^*x=b$

$x^{(0)} = 0, r^{(0)} = b, d^{(0)} = r^{(0)}$ (these are all vectors)

for $k = 1, 2, 3, \dots :$

$\alpha^{(k)} = (r^{(k-1)T} r^{(k-1)}) / (d^{(k-1)T} A d^{(k-1)})$ step length

$x^{(k)} = x^{(k-1)} + \alpha^{(k)} d^{(k-1)}$ approximate solution

$r^{(k)} = r^{(k-1)} - \alpha^{(k)} A d^{(k-1)}$ residual = $b - A x^{(k)}$

$\beta^{(k)} = (r^{(k)T} r^{(k)}) / (r^{(k-1)T} r^{(k-1)})$ improvement

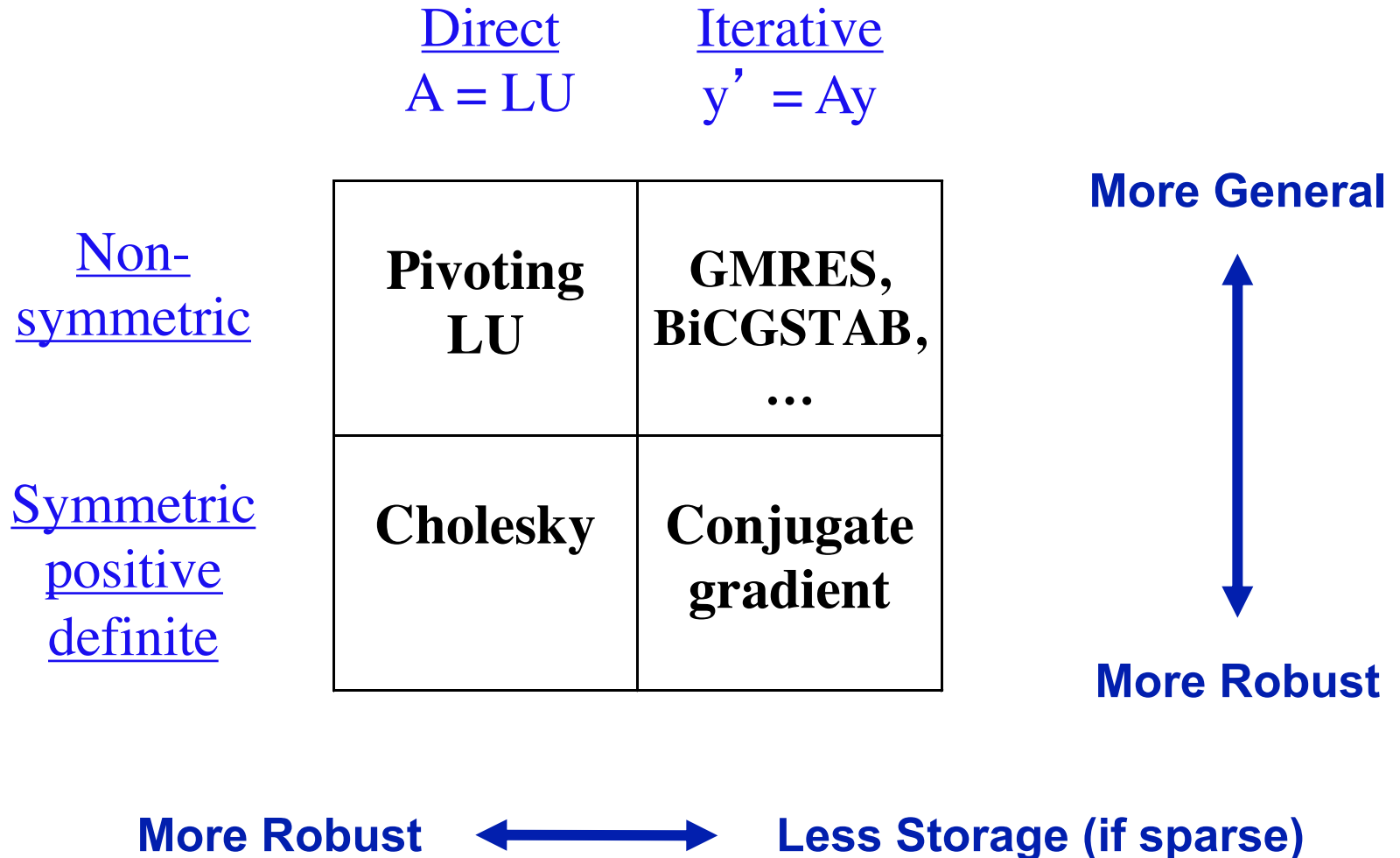
$d^{(k)} = r^{(k)} + \beta^{(k)} d^{(k-1)}$ search direction

- One matrix-vector multiplication per iteration
- Two vector dot products per iteration
- Four n-vectors of working storage

Vector and matrix primitives for CG

- **DAXPY:** $v = \alpha * v + \beta * w$ (vectors v, w ; scalars α, β)
 - Time = $O(n)$
- **DDOT:** $\alpha = v^T * w = \sum_j v[j] * w[j]$ (vectors v, w ; scalar α)
 - Time = $O(n)$
- **Matvec:** $v = A * w$ (matrix A , vectors v, w)
 - This is the hard part!
 - Time = $O(n^2)$ if A is a full matrix stored as a 2-D array
 - But all you need is a subroutine to compute v from w
 - If A is *sparse*, time = $O(\text{\#nonzeros in } A)$

The Landscape of $Ax=b$ Solvers



Optional:

***Analysis of the Conjugate Gradient
Algorithm***

See Shewchuk's paper (linked to course web site) for details.

Conjugate gradient: Krylov subspaces

- Eigenvalues: $Av = \lambda v$ $\{\lambda_1, \lambda_2, \dots, \lambda_n\}$

- Cayley-Hamilton theorem:

$$(A - \lambda_1 I) \cdot (A - \lambda_2 I) \cdot \dots \cdot (A - \lambda_n I) = 0$$

$$\text{Therefore } \sum_{0 \leq i \leq n} c_i A^i = 0 \text{ for some } c_i$$

$$\text{so } A^{-1} = \sum_{1 \leq i \leq n} (-c_i/c_0) A^{i-1}$$

- Krylov subspace:

$$\begin{aligned} &\text{Therefore if } Ax = b, \text{ then } x = A^{-1} b \text{ and} \\ &x \in \text{span}(b, Ab, A^2b, \dots, A^{n-1}b) = K_n(A, b) \end{aligned}$$

Conjugate gradient: Orthogonal sequences

- Krylov subspace: $K_i(A, b) = \text{span}(b, Ab, A^2b, \dots, A^{i-1}b)$
- Conjugate gradient algorithm:
 - for $i = 1, 2, 3, \dots$
 - find $x^{(i)} \in K_i(A, b)$
 - such that $r^{(i)} = (b - Ax^{(i)}) \perp K_i(A, b)$
- Notice $r^{(i)} \in K_{i+1}(A, b)$, so $r^{(i)} \perp r^{(j)}$ for all $j < i$
- Similarly, the “directions” are A -orthogonal:
$$(x^{(i)} - x^{(i-1)})^T \cdot A \cdot (x^{(j)} - x^{(j-1)}) = 0$$
- The magic: Short recurrences. . .
 - A is symmetric \Rightarrow can get next residual and direction from the previous one, without saving them all.

Conjugate gradient: Convergence

- In exact arithmetic, CG converges in n steps
(completely unrealistic!!)
- Accuracy after k steps of CG is related to:
 - consider polynomials of degree k that are equal to 1 at 0.
 - how small can such a polynomial be at all the eigenvalues of A ?
- Thus, eigenvalues close together are good.
- Condition number: $\kappa(A) = \|A\|_2 \|A^{-1}\|_2 = \lambda_{\max}(A) / \lambda_{\min}(A)$
- Residual is reduced by a constant factor by $O(\kappa^{1/2}(A))$ iterations of CG.

Other Krylov subspace methods

- Nonsymmetric linear systems:
 - GMRES:
for $i = 1, 2, 3, \dots$
find $x^{(i)} \in K_i(A, b)$ such that $r^{(i)} = (Ax^{(i)} - b) \perp K_i(A, b)$
But, no short recurrence \Rightarrow save old vectors \Rightarrow lots more space
(Usually “restarted” every k iterations to use less space.)
 - BiCGStab, QMR, etc.:
Two spaces $K_i(A, b)$ and $K_i(A^T, b)$ w/ mutually orthogonal bases
Short recurrences $\Rightarrow O(n)$ space, but less robust
 - Convergence and preconditioning more delicate than CG
 - Active area of current research
- Eigenvalues: Lanczos (symmetric), Arnoldi (nonsymmetric)