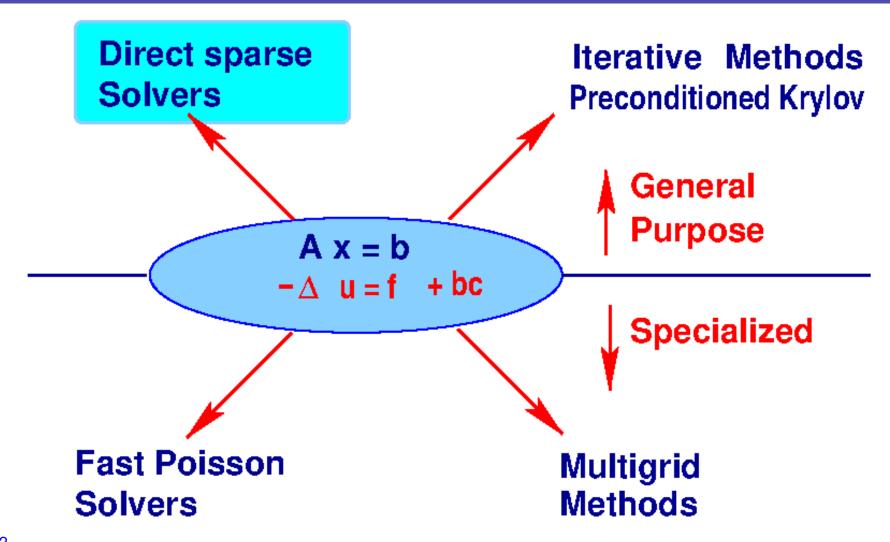
SPARSE GAUSSIAN ELIMINATION

(introduction)

- Gaussian elimination dense case
- Variants
- Main computational kernels
- Gaussian elimination in sparse case: a preview
- Graph model of elimination

Solving sparse systems today



Background. Three types of methods:

- Direct methods: based on sparse Gaussian eimination, sparse Cholesky,...
- ➤ Iterative methods: compute a sequence of iterates which converge to the solution preconditioned Krylov methods..
- > Special purpose methods: Multigrid, Fast-Poisson solvers, ...

Remark: The first 2 classes of methods have always been in competition.

 \succ 40 years ago solving a system with n=10,000 was a challenge [Now you can solve this in a fraction of a second on a laptop]

Quotation from R. Varga's book on iterative methods [1962]

"As an example of the magnitude of problems that have been successfully solved by cyclic iterative methods, the Bettis Atomic Power Laboratory of the Westinghouse Electric Corporation had in daily use in 1960 a 2-dimensional program which would treat as a special case Laplacean-type matrix equations of order 20,000."

He adds in footnote: (paraphrase) the program was written for the Philco-2000 computer which had 32,000 words of core storage (32Kwords!). "Even more staggering": Bettis had a 3-D code which could treat coupled matrix equations of order 108,000.

- > Sparse direct methods made huge gains in efficiency. They are very competitive for 2-D problems.
- > 3-D problems lead to more challenging systems [inherent to the underlying graph]

Difficulty:

No robust 'black-box' iterative solvers.

At issue Robustness in conflict with efficiency.

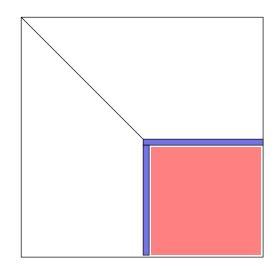
➤ Iterative methods are starting to use some of the tools of direct solvers to gain 'robustness'

Gaussian Elimination. Variants

Recall: Gaussian Elimination has 3 nested loops.

- ➤ The three loop indices are denoted by k, i, j
- We can order each of the loops differently.
- A total of 6 different algorithms [more if we add blocking]: kij, kji, ijk, ikj, jki, jik.
- ➤ IMPORTANT: these algorithms are equivalent. Same operations are done in a different order
- Therefore same operation counts
- Find the analogous algorithms for Matrix-matrix multiplication

```
For k=1:n-1 Do:
a_{ik}:=a_{ik}/a_{kk}
\text{For } \textbf{\textit{j}}:=k+1:n Do
a_{ij}:=a_{ij}-a_{ik}*a_{kj}
End
End
End
```



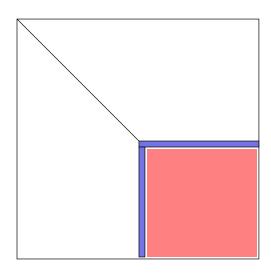
End

- This is the KIJ variant of GE [hint: read the loop indices]
- Also known as the "Outer product" form
- Main drawback: rank-one update matrix at each step.
- Can also have a KJI version [flip loops i and j]

Gauss

```
For k = 1 : n - 1 Do:
   For i := k+1 : n Do
     a_{ik} := a_{ik}/a_{kk}
   EndDo
   For j = k + 1 : n Do:
     For i := k+1 : n Do
        a_{ij} := a_{ij} - a_{ik} * a_{kj}
     End
   End
```





End

- \triangleright Can we swap loops k and j ['delay the k-loop']
- Consider first the KIJ (row) version [flip loops k and i]

Gaussian Elimination. IKJ variant

```
For i=2,\ldots,n Do:
a_{ik}:=a_{ik}/a_{kk}
\text{For } j=k+1,\ldots,n \text{ Do:}
a_{ij}:=a_{ij}-a_{ik}*a_{kj}
\text{EndDo}
\text{EndDo}
```

- Also known as the 'up-looking LU'
- Row-oriented 'delayed update' algorithm
- Can also have a column version [very common]

Gaussian Elimination. JKI or 'Left-looking LU'

JKI

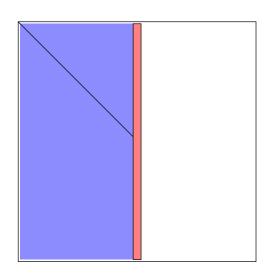
```
for j=1:n

for k=1:j-1

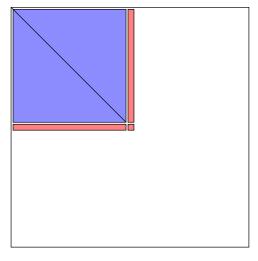
A(k+1:n,j) = A(k+1:n,j) - ...
A(k,j)*A(k+1:n,k);
end

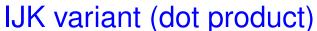
// i loop:

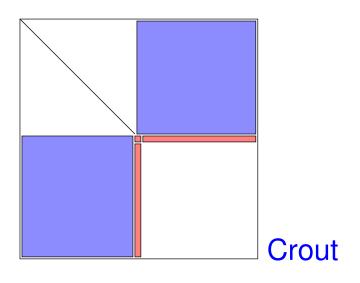
A(j+1:n,j) = A(j+1:n,j) / A(j,j);
end
```



Other variants







➤ Also: corresponding variants of Cholesky – We will only look at the adaptation of the JKI variant of Cholesky..

Column Cholesky (compare with JKI – Gauss)

```
for j=1:n

for k=1:j-1

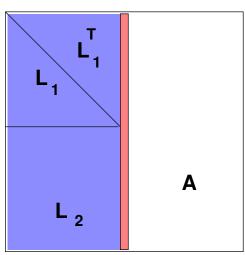
A(j:n,j)=A(j:n,j)-A(j:n,k)*A(j,k);

end

A(j,j)=sqrt(A(j,j));

A(j+1:n,j)=A(j+1:n,j)/A(j,j);

end
```



 \blacktriangleright Column j of A [L-part] becomes column j of L.

- We will often consider Sparse Cholesky because: 1) the SPD case is important; 2) certain aspects are simpler than Gauss; 3) Generalizations are easy...
- Sparse Column Cholesky: same as above algorithm but implemented in sparse mode

Number of operations. The total number of multiplications required to

compute the Cholesky factor L of a matrix A is given by

$$\sum_{k=1}^{n-1} (\eta_k^2 - 1)$$

where η_k is the number of nonzero entries in the k-th column of L

Proof:

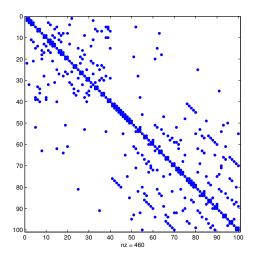
- Consider only the KIJ version of Cholesky which is equivalent.
- ightharpoonup Rank-1 update at each step is $A^{(k)} = A^{(k-1)} v_k v_k^T$, where [using matlab notation]

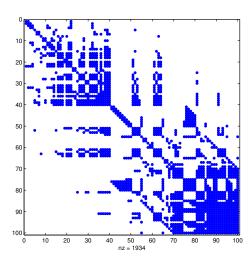
$$m{v}_k = [m{zeros}(1,k), A^{(k-1)}(k,k+1:n)]^T$$

- ► Only lower part is done so cost is $(\eta_k 1)\eta_k$.
- \blacktriangleright Must add η_k-1 scaling operations (mult. by inverse). Total: η_k^2-1
- ightharpoonup OK but η_k 's not known in advance. Dense case $\eta_k = n k + 1$
- \triangleright Storage: $\sum_{k=1}^{n} \eta_k$

Sparse solvers

- Need to develop sparse versions: sparse operations
- Major new consideration: Fill-in.





These variants will lead to different computation kernels

Gaussian Elimination. Graph model

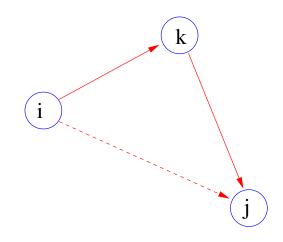
What happens to the graph when a node is eliminated?

Eliminating a node x in Gaussian elimination amounts to

- 1. Deleting node *x* and its adjacent edges from the graph
- 2. Adding edges between any two nodes that were adjacent to x.
- Notation: $A^{(k)} = \text{matrix at } k$ -th step of elimination

Another viewpoint: At step k of Gaussian Elimination, a fill-in is introduced in position (i, j) iff

$$a_{ik}^{(k-1)}
eq 0 \& a_{kj}^{(k-1)}
eq 0$$

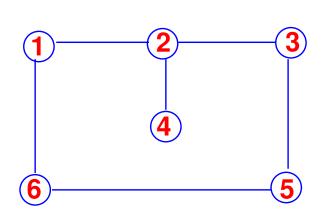


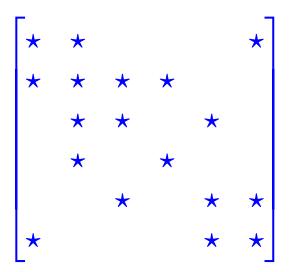
A seminal paper:

- S. Parter, "The use of linear graphs in Gauss elimination", SIAM review, vol. 3, (1961), pp. 119-130.
- ➤ Gave a major impetus to the use of the graph theory approach to sparse matrix techniques.
- Foundation for some of the major ideas (e.g. elimination trees) to come even much later.

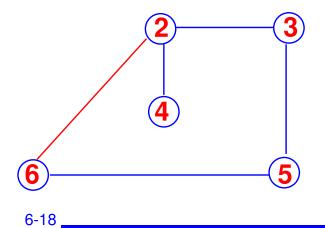
Gauss

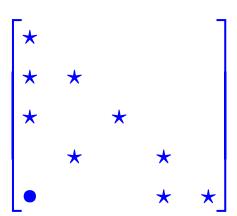
Example: Gaussian elimination steps for following graph



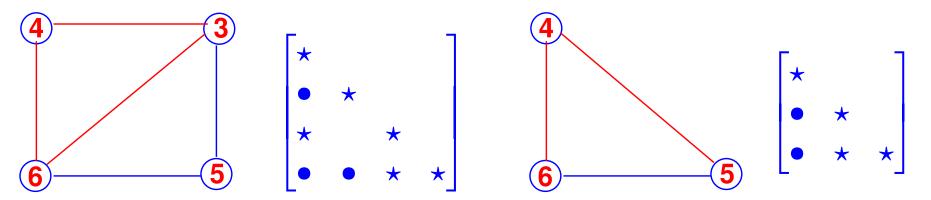


Eliminate 1:

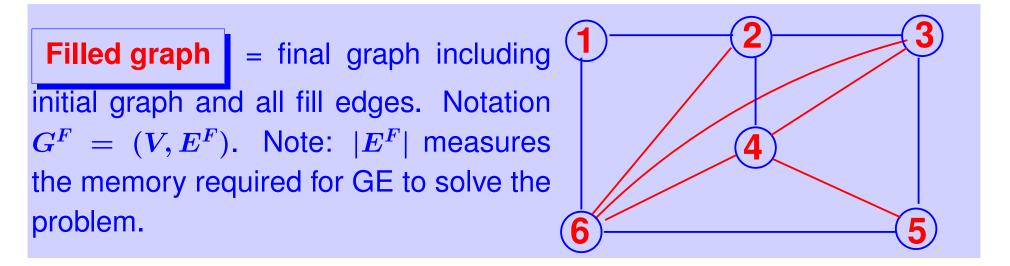




Eliminate 2 and then 3:



Two more steps [omitted - they involve no fill-in]

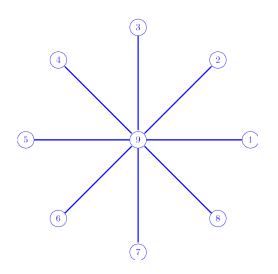


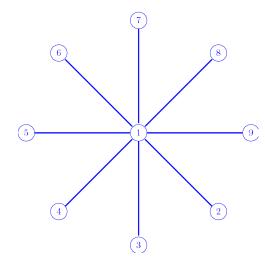
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Gaussian Elimination. The fill path theorem

ightharpoonup A Fill-path is a path between two vertices i and j in the graph of A such that all vertices in the path, except the end points i and j, are numbered less than i and j.

Mhat are all the fill-paths for the two examples below

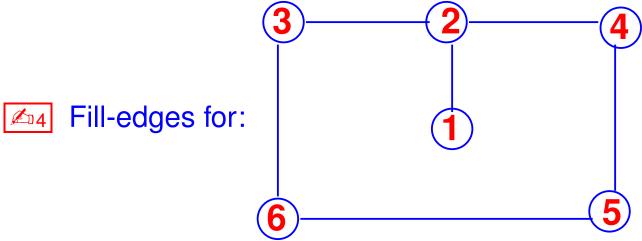




THEOREM [Rose-Tarjan] There is a fill-in in entry (i, j) at the completion of the Gaussian elimination process if and only if, there exists a fill-path between i and j.

Example of application: Separating a graph \equiv finding 3 sets of vertices: V_1, V_2, S such that $V = V_1 \cup V_2 \cup S$ and V_1 and V_2 have no couplings. Labeling nodes of S last prevents fill-ins between nodes of V_1 and V_2 .

What are all the fill-edges for the previous examples (star graphs)



Gauss