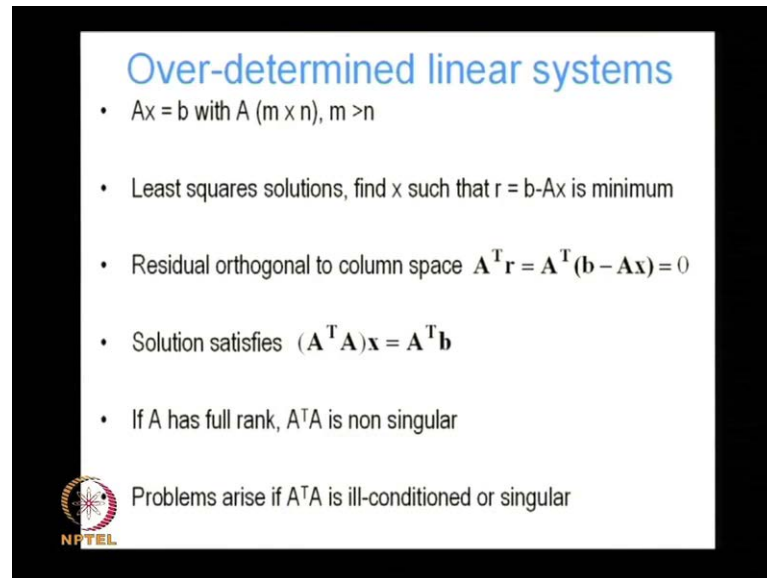


Numerical Methods in Civil Engineering
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Lecture - 10
Iterative Methods - III

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Over-determined linear systems

- $Ax = b$ with A ($m \times n$), $m > n$
- Least squares solutions, find x such that $r = b - Ax$ is minimum
- Residual orthogonal to column space $A^T r = A^T (b - Ax) = 0$
- Solution satisfies $(A^T A)x = A^T b$
- If A has full rank, $A^T A$ is non singular

 Problems arise if $A^T A$ is ill-conditioned or singular

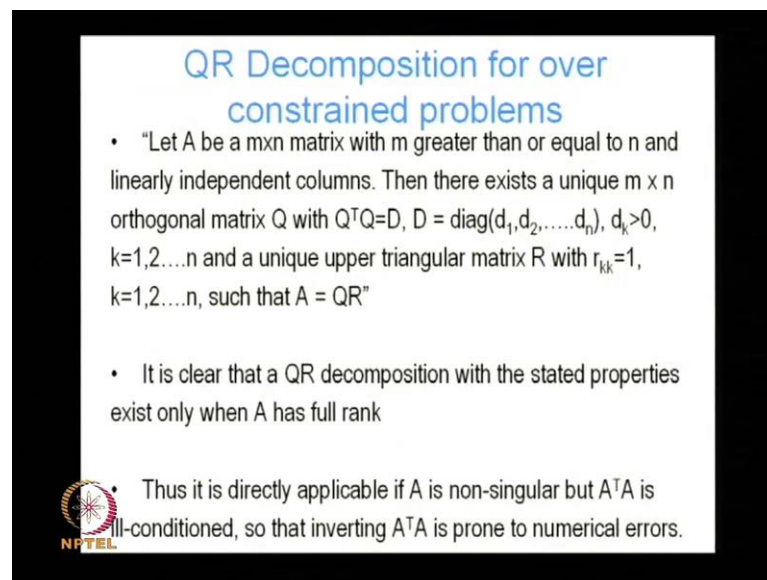
Once again, our series on numerical methods in engineering, recall last time, we talked about iterative method for over determined linear systems. What are over determined linear system? Where there are where there is a co-efficient matrix A , which has size m by n where m is greater than n , basically we have to note. So, those are over determined linear systems. So, for instant for example, for over determine linear system is when we have least when we want to find the least square solution to a problem Ax is equal to b . So, we want to find the solution x , which is the best solution in some norm. So, instance for least square solution we want to find the solution which give the minimum residual in the l_2 norm right.

We found to find x such that r equal to b minus Ax is minimum and the minimum in the l_2 norm and we also recall that the residual if we minimize the function in the and find the minimize residual the l_2 norm of the residual, we find an x and we also find the residual which is equal to b minus Ax and the residual is orthogonal to the space span by column vectors of A . The vector space span by the column of A the residual is orthogonal to that space or $A^T r$ is equal to zero right. So, the residual basically

projects out the component it is basically the component left of the projecting b on to the space which is span by the column vectors of A and the solutions also satisfy.

We also recall this condition this relationship $A^T A x$ is equal to $A^T b$ and if A has full rank, we found last time that $A^T A$ is always going to have full rank two that is a transpose always going to have the positive determine it is going to be non singular. So, by solving this equation by inverting $A^T A$ we can find out a least square solution x . However, problems arise when $A^T A$ is in ill condition or singular if it is ill condition we cannot invert it with appropriate level of accuracy and if it is singular we cannot invert it at all.

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The slide is titled "QR Decomposition for over constrained problems" in blue text. It contains three bullet points. The first bullet point states: "Let A be a $m \times n$ matrix with m greater than or equal to n and linearly independent columns. Then there exists a unique $m \times n$ orthogonal matrix Q with $Q^T Q = D$, $D = \text{diag}(d_1, d_2, \dots, d_n)$, $d_k > 0$, $k=1, 2, \dots, n$ and a unique upper triangular matrix R with $r_{kk}=1$, $k=1, 2, \dots, n$, such that $A = QR$ ". The second bullet point states: "It is clear that a QR decomposition with the stated properties exist only when A has full rank". The third bullet point states: "Thus it is directly applicable if A is non-singular but $A^T A$ is ill-conditioned, so that inverting $A^T A$ is prone to numerical errors." In the bottom left corner of the slide, there is a small circular logo with a star and the text "NPTEL" below it.

In that case, we said that the proper way to go about doing this problem is probably to do QR decomposition QR a decomposition in a QR D decomposition, which says there is the QR theorem with which say, let A be a m by n matrix with m greater to or equal to n and linearly independent column. Then there exist a unique m by n orthogonal matrix Q such that $Q^T Q$ is another diagonal matrix D with all positive entries and an unique upper triangular matrix R with one on all the diagonals right. You need upper triangle matrix R . So, A is equal to QR were Q is orthogonal matrix and R is unit upper triangle matrix.


It is clear that QR decomposition with the stated property exists only when A has full rank why is that because we have said that A has n linearly independent column right.

So, if it has n linearly independent column it is got to have full rank. So, the QR decomposition, assume that A has full rank thus it is applicable if A is non singular, but A transpose is ill condition. So, it is directly applicable to the case where A transpose A is ill condition. So, that a transpose A is prone to numerical errors.

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QR Decomposition

- However a modified form of the decomposition can also be used to find the least squares solution in case A does not possess full rank and consequently $A^T A$ is singular.
- Considering first A with full rank, using QR decomposition, $A^T(b - Ax) = 0$ can be written as: $R^T Q^T (b - Ax) = 0$
- Since R is non-singular (its determinant is one) this condition becomes: $Q^T (b - Ax) = 0 \Rightarrow Q^T b = Q^T Ax = Q^T QRx = DRx$

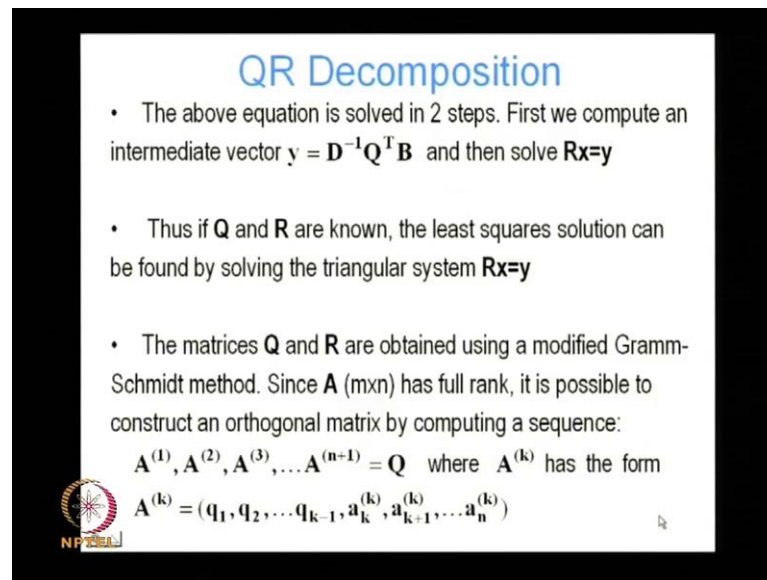


Hence to find x we need to solve the system $x = R^{-1} D^{-1} Q^T b$

However even if A is singular that is a does not have all linear independent column we can use the modified form of QR decomposition to find the least square solution. Let us first consider the case were A has full rank. So, using QR decomposition we can replace A transpose by R transpose Q transpose R transpose Q transpose b minus A x must be equal to zero. Since A is non singular we call R is a unit upper triangular matrix and also recall that upper triangular matrix the determinant is given by the product of its diagonal term and since all it diagonal term is one it is evident that R must be non singular.

So, since R is the non singular is a non singular upper triangular matrix we know that R is invertible. So, this inflation is effectively becomes Q transpose b minus A x right because R is always invertible. So, Q transpose b minus A x is equal to zero implies Q transpose b is equal to Q transpose A x or Q transpose A x is again is equal to QR using the QR decomposition Q transpose QR x Q transpose Q is equal to the diagonal matrix D . So, Q transpose b is equal to D R x . Hence to find x we have to solve this system we have to invert this D R may product D R and we get x is equal to R inverse D inverse Q transpose b .

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QR Decomposition

- The above equation is solved in 2 steps. First we compute an intermediate vector $y = D^{-1}Q^T B$ and then solve $Rx=y$
- Thus if Q and R are known, the least squares solution can be found by solving the triangular system $Rx=y$
- The matrices Q and R are obtained using a modified Gram-Schmidt method. Since A ($m \times n$) has full rank, it is possible to construct an orthogonal matrix by computing a sequence:
 $A^{(1)}, A^{(2)}, A^{(3)}, \dots, A^{(n+1)} = Q$ where $A^{(k)}$ has the form
 $A^{(k)} = (q_1, q_2, \dots, q_{k-1}, a_k^{(k)}, a_{k+1}^{(k)}, \dots, a_n^{(k)})$

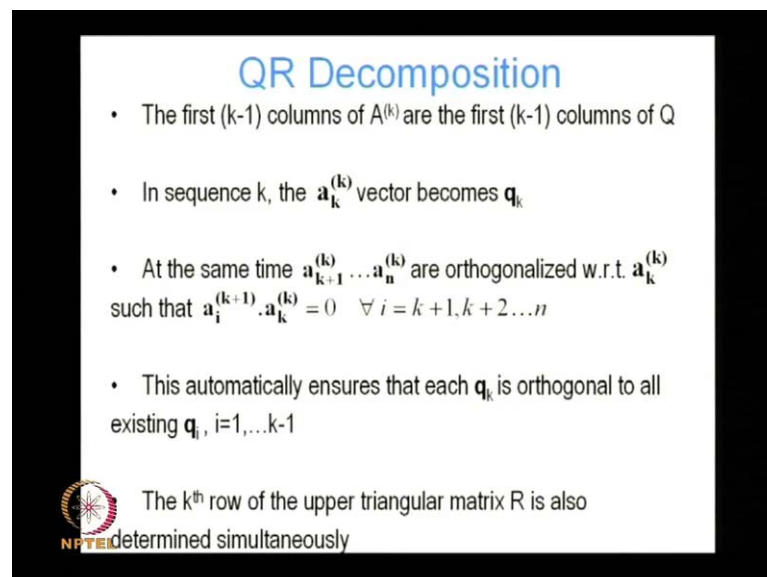
The above equation is solved in two steps, recall whenever we do operations with triangular matrixes are the both upper triangular matrixes then, which we found when we did our decomposition in Gaussian elimination. So, all this sometimes make sense to do this problem in to steps, but in order to do this problem in two steps we have to define the inter mediate vector y is equal to D inverse Q transpose B . So, basically we define intermediate vector comprising these terms right D inverse Q transpose B and then we can calculate and then we solve $R x$ is equal to y .

So, if Q and R is known then least square solution can be found by solving the triangular system $R x$ is equal to y . Off course assuming that, we have already found the intermediate vector y from this operation D inverse Q transpose B . So, again this is a type this is not the V this is because this is that is not the matrix it is the vector right. So, this is $D y$ equal to t inverse t transpose the vector p which is right hand side of my system equations. How do we find the matrix u and r .

Basically, this we can this is the good time to introduce what is known as the method of graham smith, graham smith orthogonalization. So, Q and R are found using a modified graham smith method. Since A has full rank it is possible to construct the orthogonal matrix by computing A sequence A one, A two, A three and so on up till A n . Until we get orthogonal matrix Q were a k has the form a k is equal to q one q two q k minus one and so on and so forth.


So, basically let me try to explain the idea in words first basically we start with initial matrix A and we repeat we change we change one column at a time until we get an orthogonal matrix to start with. So, from A we transform A into Q. So, how do we one column at time? So, first we form the first column of the Q. What is the first column of Q? What is the first column of A? Right, but and then at the same time we make sure that the remaining column of A are orthogonal to the first column of Q. How do we do that well we do that by the orthogonal the remaining column of A with my chosen first column which is Q and we go on and we continue in this fashion. So, the first k if you look at here we are assuming that we have done up to k minus 1 right. So, we have to form the q k minus 1 k minus columns of q and then we are trying to form the k column of q.

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QR Decomposition

- The first $(k-1)$ columns of $A^{(k)}$ are the first $(k-1)$ columns of Q
- In sequence k, the $\mathbf{a}_k^{(k)}$ vector becomes \mathbf{q}_k
- At the same time $\mathbf{a}_{k+1}^{(k)} \dots \mathbf{a}_n^{(k)}$ are orthogonalized w.r.t. $\mathbf{a}_k^{(k)}$ such that $\mathbf{a}_i^{(k+1)} \cdot \mathbf{a}_k^{(k)} = 0 \quad \forall i = k+1, k+2 \dots n$
- This automatically ensures that each \mathbf{q}_k is orthogonal to all existing $\mathbf{q}_i, i=1, \dots, k-1$

 The k^{th} row of the upper triangular matrix R is also determined simultaneously

So, how are we going to do that well what we are going to do is that we will make the a k vector q k. So, this a k k vector right this vector this vector in the k th steps this vector is going to automatically become q k. But in order to become q k we have to make sure that the remaining a k plus 1 k through a n k, they are orthogonal to the vector a k k. How are we going to do that well we are going to do that by taking out the projection the projection of a k with respect of q k a k k, which become q k from all this remaining vector.

So, the same time we make the vector a_k we orthogonalize a_{k+1} through a_n with respect to a_k , such that $a_{k+1} \cdot a_k = 0$. So, that all these vector also now after orthogonalize a_{k+1} become a_{k+1} because that is next iterative right how does a_{k+1} change to a_{k+1} by orthogonalizing it with respect to my new q_k which is a_k .

So, this automatically ensure that each q_k is orthogonal to all the existing q_i . So, basically let me go over it again. So, we start with the first vector of a the first column vector of a and make it first column of q , but in order to do that I have to make sure that all the remaining column vectors of a are orthogonal to that q , how are we going to do that from each of those column vector remaining column vectors, I am going to project out the component of q right.

So, if my second column vector is a_2 right I am going to compute my new a_2 by a_2 minus $a_2 \cdot q_1$ dotted with the unit vector in the q_1 direction. So, I am just projecting out from a_2 the part which is lies along q_1 . So, simultaneously as I compute my orthogonal vector simultaneously as I compute my q_i also compute k th row of my upper triangular matrix r .

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QR Decomposition

- The steps of the algorithm are as follows:

$$q_k = a_k^{(k)} \quad d_k = q_k \cdot q_k \quad r_{kk} = 1$$

$$a_j^{(k+1)} = a_j^{(k)} - \frac{q_k \cdot a_j^{(k)}}{d_k} q_k \quad r_{kj} = \frac{q_k \cdot a_j^{(k)}}{d_k}, \quad j = k+1, \dots, n$$
- We also compute the residual vector as the sequence $b^{(1)} = b, b^{(2)}, b^{(3)}, \dots, b^{(n+1)} = r$ by at each sequence projecting out its components in the newly obtained q_k direction:

$$b^{(k+1)} = b^{(k)} - \frac{q_k \cdot b^{(k)}}{q_k \cdot q_k} q_k = b^{(k)} - r_k q_k$$



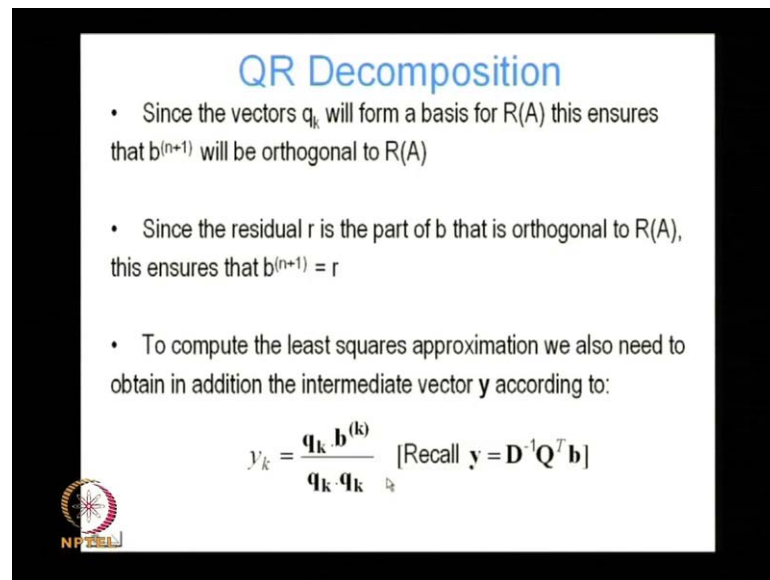
The step of the algorithm is as follows suppose in the q at the k th step first thing that we do is we make the a_k th column a_k to q_k vector right then from all the remaining columns of a from j is equal to $k+1$ to n for all the remaining column. I project out

the component of that vector along the q_k vector and of course, I normalized it by the magnitude of the q_k vector right. So, this make sure that all my remaining columns vector after my k th vector right they all orthogonal to all the previous q_k vector that as been found out till now. So, in the next iteration automatically take the first vector and make it my next q_k vector because i am assured that all those vector are orthogonal to my previously formed q_k vectors. So, at the same time we compute r which is nothing, but r_{kj} is equal to q_k dotted with a_{jk} divided by d_k .

So, this is how we form our qr decomposition. So, we start with the matrix a which is m by n and then we step by step we orthogonalize that matrix may each column orthogonal each of the remaining other columns we do it systematically using gram smith orthogonalization we are going to end up with an orthogonal matrix q and a upper triangular matrix r . At the time we all compute the residual vector as the sequence b_1 b_2 b_3 b_n as we do the as we compute each additional column of q we also change our right hand side b . So, that at.

The end we are going be left with the residual vector r because at the each iteration there are also projecting out from b we start with b_0 b_1 equal to b and each step we project out from b the component its components in the current q_k vector direction. So, we start with my original right hand side at each iteration at each step my qr algorithm i project out from my right hand side the component the projection of the right hand side in the current q_k vector direction. So, at the end i am left with whatever is left is just going to be orthogonal to all the q_k vectors basically since all the q_k vector and my space a ; that means, my residual is going to end up orthogonal to the space span by the column vectors of a . So, this is how i compute the b_k plus one which is b_k minus y_k q_k were i have defined y_k is equal to q_k dotted with b_k divided by q_k dotted with q_k .

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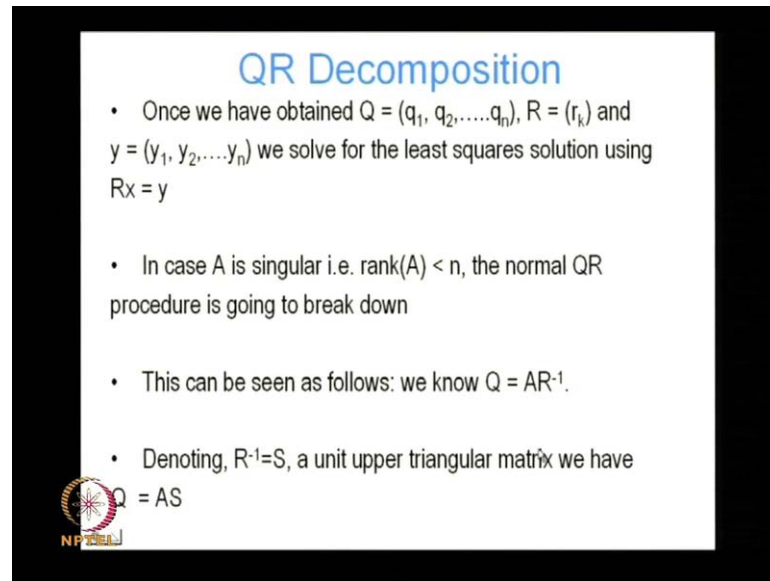


The slide is titled "QR Decomposition" in blue text. It contains three bullet points and a mathematical equation. The first bullet point states that since vectors q_k form a basis for $R(A)$, $b^{(n+1)}$ is orthogonal to $R(A)$. The second bullet point states that since the residual r is the part of b orthogonal to $R(A)$, $b^{(n+1)} = r$. The third bullet point states that to compute the least squares approximation, we need the intermediate vector y . The equation is $y_k = \frac{q_k \cdot b^{(k)}}{q_k \cdot q_k}$ with a note "[Recall $y = D^{-1}Q^T b$]". There is a small NPTEL logo in the bottom left corner of the slide.

Since the vector q_k will form the basis for r a this ensure that b^{n+1} will be orthogonal to r a recall what is r a it is the vector space spanned by all the column vectors of a . So, we since we have to ensure that the b^n are orthogonal to all the q_k s we will ensure that the end my b^{n+1} is orthogonal to all the column.


Vectors of a and therefore, what is that is that is the residual to compute the least square approximation we also need to obtain in addition the intermediate vector y because we need to compute the intermediate vector y which we defined to be y equal to d inverse q transpose b and during the q r process we also compute y like this y is equal to q dotted with b^k divided by q^k dotted with q^k we call it y is equal to d inverse q transpose b . So, this is basically that part q transpose b and then we are dividing by q^k dotted with q^k this is d inverse that operation right.

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QR Decomposition

- Once we have obtained $Q = (q_1, q_2, \dots, q_n)$, $R = (r_k)$ and $y = (y_1, y_2, \dots, y_n)$ we solve for the least squares solution using $Rx = y$
- In case A is singular i.e. $\text{rank}(A) < n$, the normal QR procedure is going to break down
- This can be seen as follows: we know $Q = AR^{-1}$.
- Denoting, $R^{-1} = S$, a unit upper triangular matrix we have $Q = AS$

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So, once we have obtained q_1, q_2, \dots, q_n . So, all these column which are orthogonal to each other right and we have obtained our right angular matrix r and we are obtained the vector intermediate vectors y we solve for the least square solution using $r x$ is equal to y . So, this is what we do when a transpose a is ill condition we can of course, use it a transpose is well condition also dot, but it is particularly situated for when it is ill condition because then with the normal method is going to break down what happens when instead of being ill condition it is acutely singular that is a transpose a is actually singular because a does not have full value what happens then in that case the normal q, r procedure is going to break down how is this going to happen we can see it as follows we know from the q, r decomposition q is equal to $a r^{-1}$ right q is equal to a is equal to $q r$. So, q is equal to $a r^{-1}$ let us denote r^{-1} is equal to s an upper angular matrix we are assuming that since r is the upper triangle its true it is not assuming it is you have to take, but it is true that since r is the upper triangular matrix r^{-1} is also going to be a upper triangular matrix and this also is going to be unit upper triangular matrix right. So, we can write q is equal to a times s a product s matrix product s .

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QR Decomposition

- Thus the k^{th} column of Q can be written as:

$$q_{1k} = a_{11}s_{1k} + a_{12}s_{2k} + \dots + a_{1(k-1)}s_{(k-1)k} + a_{1k}$$

$$q_{2k} = a_{21}s_{1k} + a_{22}s_{2k} + \dots + a_{2(k-1)}s_{(k-1)k} + a_{2k}$$


.....

$$q_{mk} = a_{m1}s_{1k} + a_{m2}s_{2k} + \dots + a_{m(k-1)}s_{(k-1)k} + a_{mk}$$

This can be written concisely as:

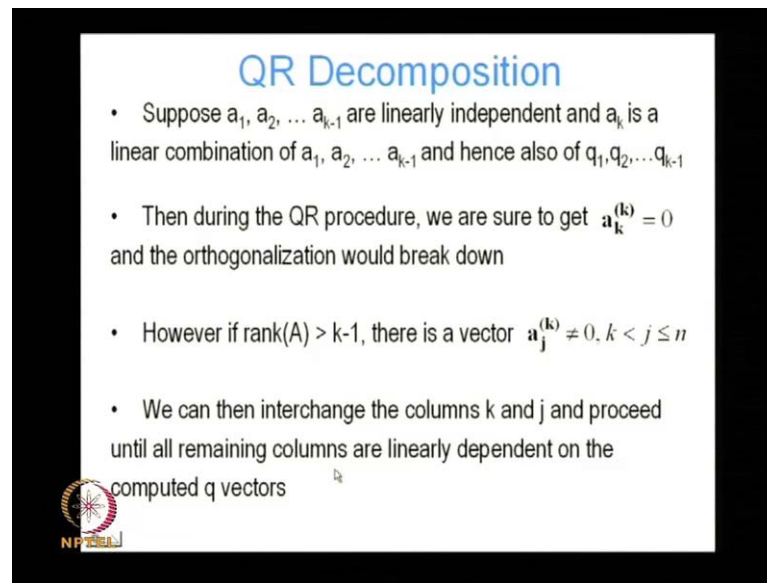
$$q_k = s_{1k}a_1 + s_{2k}a_2 + s_{3k}a_3 + \dots + s_{(k-1)k}a_{k-1} + a_k$$

- It is clear that each q_k is a linear combination of the columns a_1, a_2, \dots, a_k




Then the k^{th} column of q can be returned as q is equal to s . So, that k^{th} column of q is nothing, but q . One k^{th} the k^{th} column all the entries in the k^{th} column q and k q two k through two q m k which is going to be nothing, but a the first row of a multiplied with the k^{th} row of s right. So, this is the first row times the k^{th} row of s second row k^{th} row of s n^{th} row of k a times the k^{th} row of s . So, this gives me the k^{th} column of q . So, we can write this concisely as q_k is equal to some scalar times the first column of a right plus another scalar s_{2k} times the second column of a plus another scalar $s_{k-1,k}$ times the $(k-1)^{\text{th}}$ column of a and finally, we are going to have a one k^{th} right why there is nothing, but because there is unit triangular matrix right. So, s_{kk} is going to be one right. So, $s_{k-1,k}$ $s_{k-2,k}$ $s_{k-3,k}$ is going to be one because it is a unit triangular matrix. So, we can write we can write each q_k as a linear combination of the a vector of the columns of a matrix a one a three up to $A \times$ again this is the typo a_{k-1} plus a_k right. So, it is clear that each q_k is the linear combination of the columns of a one a two through a_k .

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QR Decomposition

- Suppose a_1, a_2, \dots, a_{k-1} are linearly independent and a_k is a linear combination of a_1, a_2, \dots, a_{k-1} and hence also of q_1, q_2, \dots, q_{k-1}
- Then during the QR procedure, we are sure to get $a_k^{(k)} = 0$ and the orthogonalization would break down
- However if $\text{rank}(A) > k-1$, there is a vector $a_j^{(k)} \neq 0, k < j \leq n$
- We can then interchange the columns k and j and proceed until all remaining columns are linearly dependent on the computed q vectors



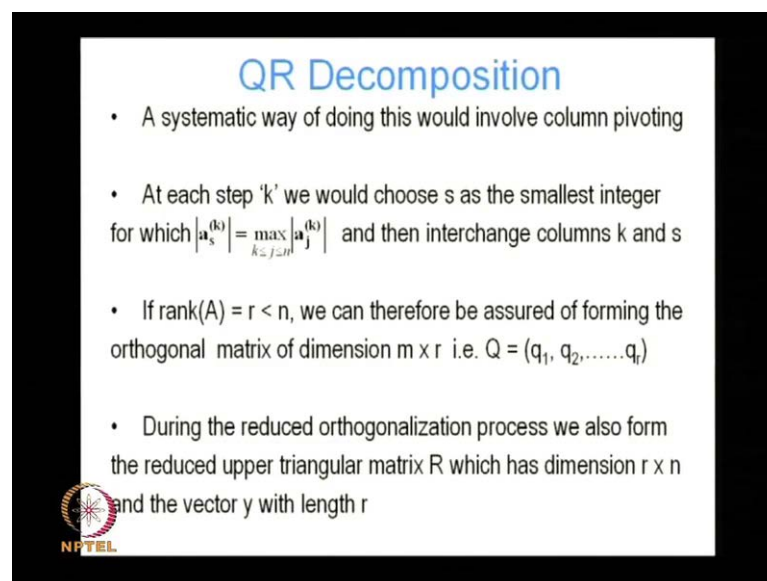
Since suppose let us suppose that a_1, a_2, \dots, a_{k-1} are linearly independent and a_k is a linear combination of a_1, a_2, \dots, a_{k-1} and hence also of q_1, q_2, \dots, q_{k-1} . Now we are considering the case where A does not have full rank. What does that mean? That means that there are columns in A that are linear combinations of the other columns.

Which are linear combinations of the other columns. Right, not all the columns of A are linearly independent. Right, A does not have full rank. So, let us assume that the k th column of A is a linear combination of the first $k-1$ columns. Right, since it is a linear combination of the first $k-1$ columns and we are just shown that each q_k is the linear combination of all the previous $k-1$ columns; that means, that a_k is also a linear combination of all the q_1 through q_{k-1} . Thus during the QR procedure we are sure to get $a_k^{(k)} = 0$ and the orthogonalization is going to break down. Why is that? Because recall at each step at each time we find a new q vector v orthogonalize all the remaining vector with respect to that vector. So, you project out from each remaining vector the q_k vector that have been formed previously. Now since a_k is the linear combination of all of $k-1$ columns of A when we project out from a_k the previous $k-1$ columns we are going to get zero because a_k is basically a linear combination of $k-1$ columns. Right, eventually we are going to land up with an a_k which is zero and then when we do that our orthogonalize procedure breaks down because the QR procedure depends on making the first column the first of the remaining column the next q_k vector, but the first of the

remaining column is now a k, but a k is identically equal to zero. So, our algorithm is going to break down, but; however, if the rank of k is greater than k up till now we.


Have found k minus one independent column, but suppose rank of k rank of a is more than k minus one if the rank of k is just minus well that is it we cannot go any further all the remaining columns are assure to be zero right because they have been all orthogonal with respect to the previous k minus one column right, but if the rank is greater than k minus one there are still some column which are non zero right there are still some of columns of a which are non zero. So, what we do we do the pivoting we do the column interchange we replace the k suppose the j th column is non zero and the j is greater than k, but lesser than or equal n we can then interchange the column k and j and proceed until all the remaining column are linearly independent on the computed k vector. So, we keep on doing this we keep one interchanging until we reach a situation where all the remaining columns are orthogonal to the computed q vectors which means that all the reaming columns are zero right.

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QR Decomposition

- A systematic way of doing this would involve column pivoting
- At each step 'k' we would choose s as the smallest integer for which $|a_s^{(k)}| = \max_{k \leq j \leq n} |a_j^{(k)}|$ and then interchange columns k and s
- If $\text{rank}(A) = r < n$, we can therefore be assured of forming the orthogonal matrix of dimension $m \times r$ i.e. $Q = (q_1, q_2, \dots, q_r)$
- During the reduced orthogonalization process we also form the reduced upper triangular matrix R which has dimension $r \times n$ and the vector y with length r

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So, systematic way of doing this would involved column pivoting at each step we would chose s as a smallest integer for with k s k is equal to max of a j k j varying from k to n and then interchange the.

Column k n s basically we at each step we instead of directly taking the next column and making the q vector we look at all the reaming columns and choose the column which as

the maximum norm maximum norm in the infinite since right infinite norm right is the maximum infinite norm make that the q vector and. So, on and. So, forth right. So, if rank of a is suppose r which is less than n we can therefore, be assure of forming orthogonal matrix of dimension m by r that is q is equal to q one q two to q r. So, that to form r since it has rank r we can form r orthogonal vectors right we can form r orthogonal vectors q is of size m cross r during the reduce orthoganilization process we also form the reduce upper triangle matrix are which has dimension r by n and the vector by the length r using the same procedure that we talked about earlier we also forming r and the y vectors.

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QR Decomposition

- We separate R into two parts, R' and S' with R' of size $r \times r$ and S' of size $r \times (n-r)$
- Similarly, the unknown solution vector x is separated into r and $(n-r)$ components:

$$x_1 = (x_1, x_2, \dots, x_r)$$

$$x_2 = (x_{r+1}, x_{r+2}, \dots, x_n)$$
- Then one can rewrite:

$$Rx = y \text{ as } R'x_1 + S'x_2 = y$$
 which gives $x_1 = (R')^{-1}y - (R')^{-1}S'x_2$ (***)

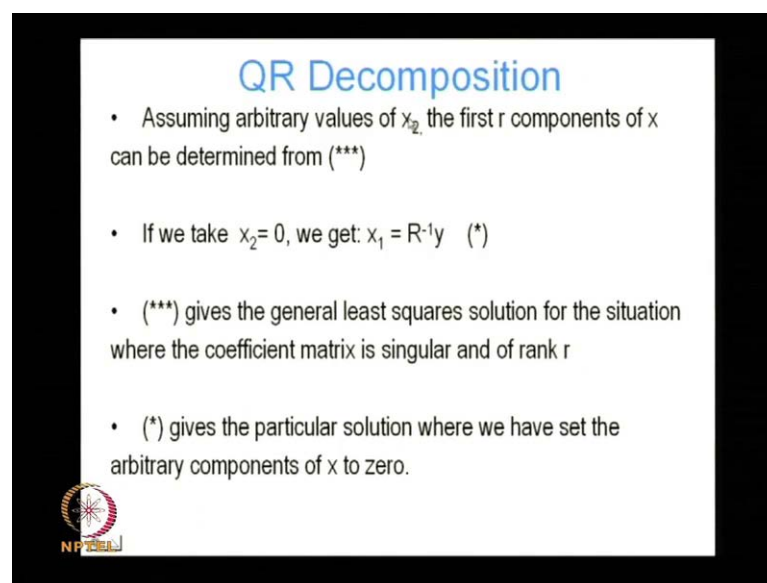


We separate r into two parts r prime and s prime with r prime of size r by remember that q is of size m by r . So, e must be of size r by n because a is of size m by n right. So, m by n is equal to m by r product with r by n . So, now, . So, basically we are saying that we are going to split r into two parts r prime and s prime p prime of size r by r and s prime of size r by n minus r similarly the unknown solution vector x is separated into r n minus r components we separate out the first r component of x like this as x one and r plus one to the n th component as x two then we can write r x is equal to y such that.

Now, see we have split out r into two parts r prime which is of size r by r and s prime this is size r n minus r similarly we have split up this vector into this little part x one which is of size r and this part x two which is of size n minus r right. So, r prime x one plus x


prime x_2 must be equal to r which gives me x_1 i can write it has r prime inverse by minus r prime inverse this x_2 are arbitrary right because my matrix has my e matrix has got rank r right we have to assume x_2 arbitrary value x_2 arbitrary remember anytime we have less than full rank we are going to have infinitely many solution right infinitely many solution and we have to assume the as many solution as the null space right in this case the null space is of size n minus r right. So, x_2 is of size n minus r these r arbitrary values and i can write my remaining expand in terms of this x_2 right and my y which i found during the orthogonalisation.

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QR Decomposition

- Assuming arbitrary values of x_2 , the first r components of x can be determined from (***)
- If we take $x_2 = 0$, we get: $x_1 = R^{-1}y$ (*)
- (***) gives the general least squares solution for the situation where the coefficient matrix is singular and of rank r
- (*) gives the particular solution where we have set the arbitrary components of x to zero.

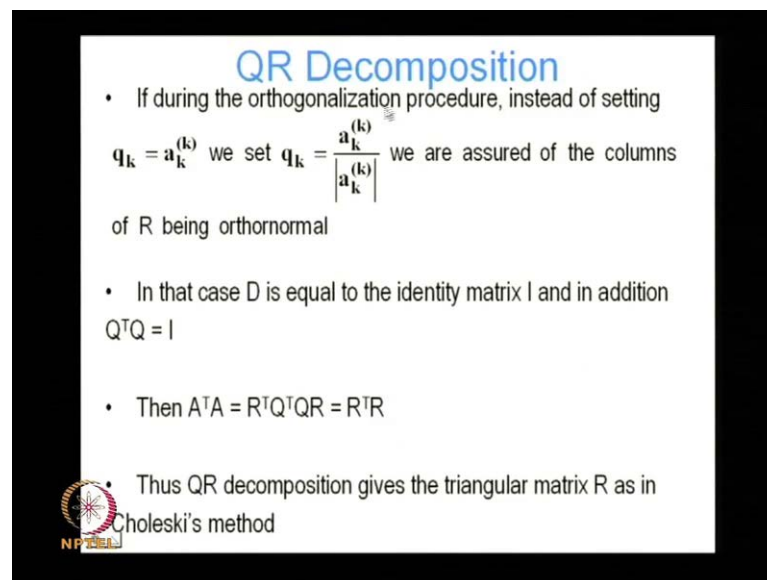


So, assuming the arbitrary value of x_2 the first r components of x can be determined from this expression takes one component similarly if we take x_2 is equal to zero since x_2 is arbitrary we can take any value if we put x_2 is equal to zero we get x_1 is equal to r inverse y .

So, this equation gives the general least square solution for the situation where the coefficient matrix is singular and of rank r right. So, there are infinity many recall now there are infinitely many least square solution right now why because my x_2 are arbitrary right those n minus r components of x are arbitrary i can make them zero i can make them anything i like right, but they all satisfy the least square solution. So, they all minimize the residual they all minimize the residual they all satisfy the least square criteria. So, infinity many least square solution to the problem right unlike the case when

transpose A^T was non singular were we had only one solution in this case the infinite many least square solution, but in all of them make sure that the residual is all them make sure that residual is orthogonal to the to the space span by the column of A in this case the space span by column of Q .

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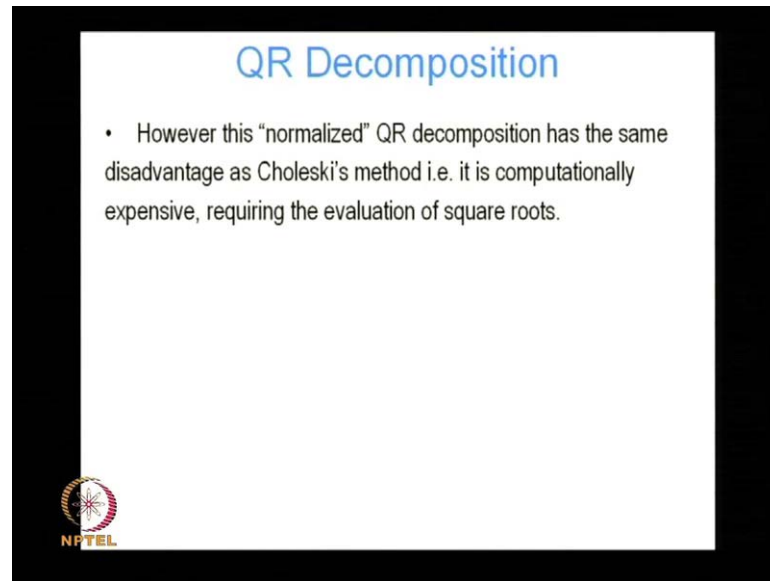
QR Decomposition

- If during the orthogonalization procedure, instead of setting $q_k = a_k^{(k)}$ we set $q_k = \frac{a_k^{(k)}}{\|a_k^{(k)}\|}$ we are assured of the columns of R being orthonormal
- In that case D is equal to the identity matrix I and in addition $Q^T Q = I$
- Then $A^T A = R^T Q^T Q R = R^T R$
- Thus QR decomposition gives the triangular matrix R as in Choleski's method

So, this is the just the diversion, but during the orthogonalization procedure instead of setting q_k is equal to a_k in addition to orthogonalize we also normalize right we ortho-normalize. So, we said the q_k vector. So, that each of them have norm one. So, not only the orthogonal to each other they unit norm also then we are assured that my R my Q matrix is this is again a typo i assured the columns of Q being.

Orthogonal apologize this is Q right the columns of Q are ortho normal right in that case D transpose D^T a sorry $Q^T Q$ is going to be identity matrix. So, D is equal to the identity matrix and in that case $A^T A$ can be written as $R^T Q^T Q R$ which is equal to $R^T R$ we have encountered this decomposition before when we talked about the choleskismethod right. So, this QR decomposition is a triangular matrix R as in choleskis method in case the ortho in addition to orthogonalizing Q we normalize the columns of Q we get a we get the we recover the choleskis method.

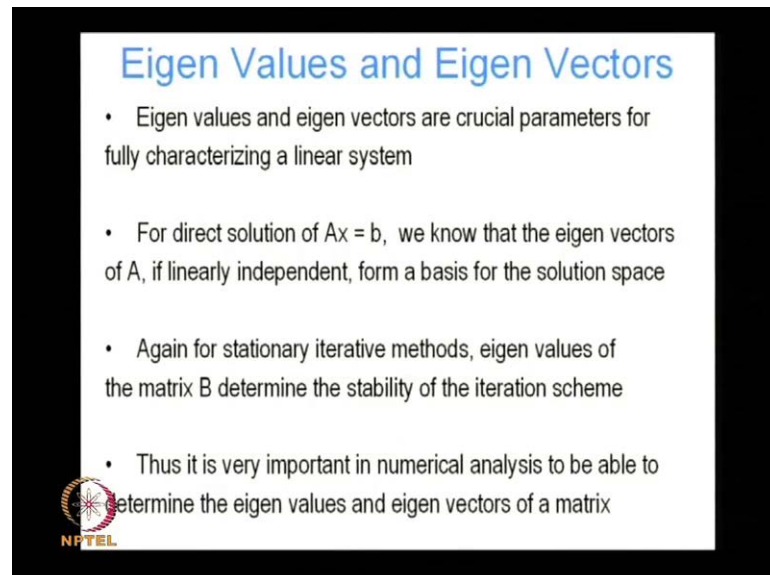
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The slide is titled "QR Decomposition" in blue text. It contains a single bullet point: "• However this 'normalized' QR decomposition has the same disadvantage as Choleski's method i.e. it is computationally expensive, requiring the evaluation of square roots." In the bottom left corner, there is a circular logo with a red and white design and the text "NPTEL" below it.

This normalized q r decomposition has the same disadvantage as Choleski's method it is computationally expensive requiring the valuation of square roots which we find in case of Choleski's method as we have $r^T r$ that going to be square terms and we have to evaluate square roots. So, it become computationally expensive ok.

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The slide is titled "Eigen Values and Eigen Vectors" in blue text. It contains four bullet points: "• Eigen values and eigen vectors are crucial parameters for fully characterizing a linear system", "• For direct solution of $Ax = b$, we know that the eigen vectors of A , if linearly independent, form a basis for the solution space", "• Again for stationary iterative methods, eigen values of the matrix B determine the stability of the iteration scheme", and "• Thus it is very important in numerical analysis to be able to determine the eigen values and eigen vectors of a matrix". In the bottom left corner, there is a circular logo with a red and white design and the text "NPTEL" below it.

So, that was the discussion on over determined linear system u r decomposition to solve the two determined linear system the q r decomposition.

It is very very important tool in the numerical method. So, it is not only usefulness not restricted to over determine linear system it can use it has lots of other uses it can be acutely used to find probably very powerful tool for finding the eigen value also all though we are not talking in detail about use q r in finding eigen value, but it is very use full tool in finding eigen value also it is very important and power full tool, but next lets switch over to finding out how can we calculate how can we solve eigen value problem basically if you have given an n by n matrix and we are interested in finding its eigen value in eigen vector how can we solve it if it is the small matrix if it is the two by two if it is three by three its four by four probably matrix we can do it by hand we can set up the characteristic equation and solve it right, but in case we have very large matrixes then solving that problems is it becomes difficult it requires very sophisticated numerical method and i can assure you that solving eigen value problem are essential if you are going to do any reasonable numerical solution if you want the stability of the analyses we need to find the eigen value countless other applications were eigen value is very important. So, eigen value and eigen vectors are crucial parameter for fully characterizing linear system. So, for direct solution of $Ax = b$ we know that since the eigen vectors of a if linearly independent form basis for the solution space we can use the eigen vectors to solve the problem right again for the stationary iterative method which we have encountered just little while ago we found that the whole stability of the method depends on the eigen value of my b matrix right which was my iteration matrix right. So, the eigen value of the b.

Matrix is determined whether my iteration method is going to be stable or not after i do the n iteration whether i am going to converge to a solution or i am going to divert and get go somewhere totally different right. So, all that depends on my on the spectrum on the spectrum of my b matrix right thus it is very important in numeric analyses to be able to determine the eigen values and eigen vectors of a matrix.

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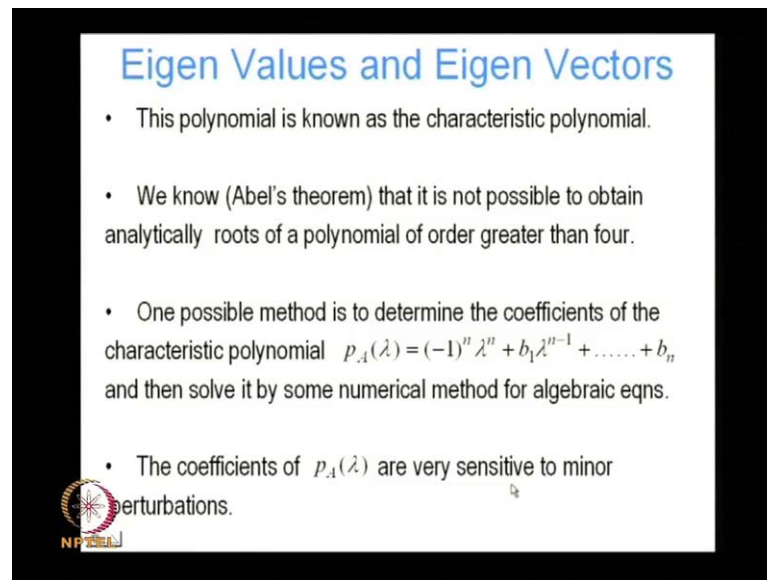
Eigen Values and Eigen Vectors

- The eigen values λ_i and the eigen vectors \mathbf{x}_i of a $n \times n$ matrix \mathbf{A} satisfy: $(\mathbf{A} - \lambda_i \mathbf{I})\mathbf{x}_i = 0, \quad \mathbf{x}_i \neq 0 \quad (*)$
- If the eigen vectors are already known, the corresponding eigen values can be determined as follows:
$$\lambda_i = \frac{\mathbf{x}_i^T \mathbf{A} \mathbf{x}_i}{\mathbf{x}_i^T \mathbf{x}_i}$$
- For the system $(*)$ to have a non-trivial solution, $\det(\mathbf{A} - \lambda_i \mathbf{I}) = 0$
- On expansion we get a polynomial of degree n , whose n roots are the n eigen values of the system.

Eigen values λ_i and the eigen vector \mathbf{x} of n by n matrix satisfy the relation $\mathbf{A} \mathbf{x} - \lambda_i \mathbf{x} = 0$ provided that $\mathbf{x} \neq 0$. If we know the eigen vector we can find the eigen value. We just compute $\mathbf{A} \mathbf{x} = \lambda_i \mathbf{x}$. So, $\mathbf{x}^T \mathbf{A} \mathbf{x} = \lambda_i \mathbf{x}^T \mathbf{x}$. We divide $\mathbf{x}^T \mathbf{A} \mathbf{x}$ by $\mathbf{x}^T \mathbf{x}$ and we get the eigen value right. So, that can be done trivially right provide the eigen vector are known; however, that is not. So, easy finding the eigen vector is not. So, easy. So, we have to find ways of finding the eigen values typically the eigen values are found first and.


Simultaneous eigen vectors are found typically we do not know that eigen vector are known beforehand. So, we know that for this system this is the homogeneous system for it to have nontrivial solution, but meaning that for not all \mathbf{x}_i components of \mathbf{x}_i to be not equal to zero then the determinant $\det(\mathbf{A} - \lambda_i \mathbf{I})$ must be equal to zero on expansion we get the polynomial of degree n . So, if it is three by three matrix we get the polynomial in third degree cubic equation in λ if it is two by two matrix we get the quadratic equation in λ if it is four by four matrix we get the quartic equation in λ . So, we get the polynomial of degree n whose n roots are the eigen values of the systems.

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Eigen Values and Eigen Vectors

- This polynomial is known as the characteristic polynomial.
- We know (Abel's theorem) that it is not possible to obtain analytically roots of a polynomial of order greater than four.
- One possible method is to determine the coefficients of the characteristic polynomial $p_A(\lambda) = (-1)^n \lambda^n + b_1 \lambda^{n-1} + \dots + b_n$ and then solve it by some numerical method for algebraic eqns.
- The coefficients of $p_A(\lambda)$ are very sensitive to minor perturbations.

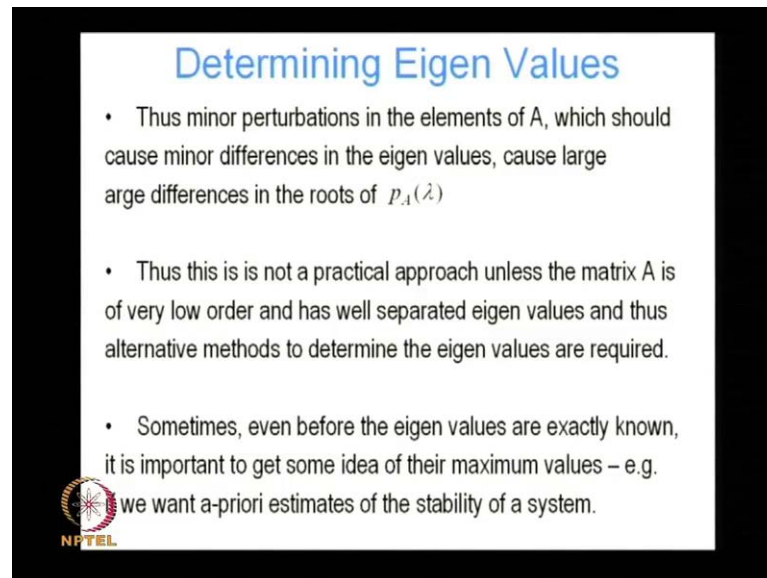


This polynomial is known as the characteristic polynomial again there is theorem which is known as abels theorem it tells that it is not possible to obtain analytically the roots of the polynomial of order greater than four. So, if this polynomial become greater than four it is a five by five matrix as the quintet polynomial right polynomial lambda to the power five in that case there is ables theorem which tells me that you cannot find the roots of this polynomial using any analytical method; however, hardly try right. So, the only way you can find the roots of that equation are through numerical methods right. So, one possible method is to determine the co efficient of the characteristic polynomial p a lambda which is like this lambda to the power n lambda to the power n minus one and. So, on and then solve it by some numerical method algebraic equation why numerical method because for order greater than four we cannot use an analytical methods and even for orders less than four it is not it is not the good idea sometimes to.

Use analytical methods i talk about the reasons just now the coefficient of p a are very sensitive to minor perturbations if you have your original matrix a and i give minor perturbation to the elements of a . So, we add to a matrix epsilon time b were epsilon is the very small number. So, i give perturbations to a . So, you except the eigen values of a to change by epsilon right because i change the components of a by epsilon. So, i would accept the eigen values a also change slightly by epsilon, but it does not happen like that why is that because the coefficient of my characteristic equation are very sensitive to

minor perturbations right. So, they are not very well condition right minor perturbation in a result in large changes in the roots of my polynomials of my characteristic equation.

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Determining Eigen Values

- Thus minor perturbations in the elements of A , which should cause minor differences in the eigen values, cause large differences in the roots of $p_A(\lambda)$
- Thus this is not a practical approach unless the matrix A is of very low order and has well separated eigen values and thus alternative methods to determine the eigen values are required.
- Sometimes, even before the eigen values are exactly known, it is important to get some idea of their maximum values – e.g. we want a-priori estimates of the stability of a system.

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Thus the minor perturbation of elements of A which should cause minor difference in the eigen value cause again there is large differences in the roots of $p_A(\lambda)$ thus this is not the practical approach until the matrix A is a very low order and has well separated eigen values and has the well separate if A is of low order and separate eigen values then this is.

Fine; however, otherwise we need to think of alternative method for determine the eigen values. So, before talking about the alternative method for determining the eigen values we are going to talk about certain important theorem in linear algebra which is known as the Gerschgorin's theorem which is very use full although it is not as use full as (refertime:39:00)it would like to hope it because it is use full because even determine because it allows us to determine the bounds on the eigen we call our stability criteria what was our stability criteria for our jacobians method or for galls syghal method the stability criteria was that the spectral radius must be less than one which means that my largest eigen value must have magnitude must have norm less than one. So, this Gerschgorin's theorem allows us even before actually computing the eigen values to find out bound on the eigen value it tells that what can be the maximum possible eigen values of this matrix given in a matrix without actually computing the eigen values by using Gerschgorin's i can tell i can find out what can be the maximum absolute value $\rho(A)$

beyond which we never be exceeded by the eigen value of that matrix right. So, before that allows us to get some idea of stability of the system even before computing the eigen value it allows us to get half clear estimate of stability of the system which are use full, but not. So, use full because the bounds predicted by the gerschgorins theorem are. So, wide right. So, they are since my eigen value is five suppose and gerschgorins theorem tells me that eigen value will never exceed that value five hundred that bound is not very use full right, but still use full for sometimes, but it is a very very wide bounds. So, narrow bounds are more use full than wide bounds right so.

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Gerschgorin's Theorem

- A theorem due to Gerschgorin provides a bound on the eigen value of a matrix. It states: Let A be a $n \times n$ matrix with eigen values $\lambda_i, i = 1 \dots n$. Then each λ_i lies in the union of the circles:

$$|z - a_{ii}| \leq r_i, \quad r_i = \sum_{\substack{j=1 \\ j \neq i}}^n |a_{ij}|$$

- This can be proved as follows: Let λ be an eigen value of A . Then there is an eigenvector x not equal to zero such that $Ax = \lambda x$

Hence, $(\lambda - a_{ii})x_i = \sum_{\substack{j=1 \\ j \neq i}}^n a_{ij}x_j, \quad i = 1, 2, \dots, n$

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So, let us look at gerschgorins theorem the theorem due to gerschgorins provides a bound of.

The eigen value of a matrix it states that let a be an n by n matrix with eigen value lambda one lambda i i is equal to m then each lambda i lays in the union of the cycle from like that. So, what does; that means, it means that if i form a circle which centre i form a series of circles right and each circles circle one for instants has centre with diagonal element one one right it has center with a diagonal element one one and it has got radius which is some of the absolute values of all the diagonal all the non diagonal elements in that row right if i am looking at the first row alright the radius of the circle is sum of all the non diagonal element in the first row s it tells me that if i compute circles right this look at the first row look at its diagonal element make that the centre of my

circle and draw a circle with a radius which is equal to the sum of norm of my half diagonal elements i do that for the second row again draw a circle centered with the diagonal term and radius equal to the some of the.

Absolute value of diagonal elements i go on doing this then this union of all this circle right the union of all this circle defines my gerschgorins bound. So, if i draw this series of circle and then i draw a circle which circumscise all this circle which is basically the union of all these circle union of all those circle the eigen value will lay within that circle the eigen value cannot lay beyond that circle the maximum eigen value cannot lay beyond that circle this can be proved as follows let lambda be eigen value of a then there is eigen value of x is not equal to zero such that A x is equal to lambda x right hence we can write lambda minus a i i x i is equal to i x a basically i kept the diagonal term on the left hand side and i have moved all of the diagonal term to the right hand side.

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Gerschgorin's Theorem

- Let us consider the equation i corresponding to the component of x with the largest absolute value i.e. $|x_i| = |x|_{\infty}$.

Then
$$|(\lambda - a_{ii})x_i| = \sum_{\substack{j=1 \\ j \neq i}}^n |a_{ij}x_j| \leq \sum_{\substack{j=1 \\ j \neq i}}^n |a_{ij}| |x_j|$$

Dividing throughout by $|x_i|$ we get :

$$\frac{|(\lambda - a_{ii})x_i|}{|x_i|} = |\lambda - a_{ii}| \leq \sum_{\substack{j=1 \\ j \neq i}}^n \frac{|a_{ij}| |x_j|}{|x_i|} \leq \sum_{\substack{j=1 \\ j \neq i}}^n |a_{ij}| = r_i$$

- Thus the largest eigen value lies on or within the **union** of all circles with centers at the diagonal elements of the coefficient matrix A and with radius as given above

So, let us consider. So, if i do that i can do that for each of my equation i equal to one through n assuming that a is the matrix of size m by n. So, i can do it for each of those rows next let us consider the equation i corresponding to the component to the x largest absolute value right. So, let us consider i th row such that i th row x has the maximum absolute value right in that case for that row i can write mode of lambda minus a i i x i is equal to sigma mode of a i j basically am just taking the mode of both sides right am taking the mode on both side in this equation and mode of sum of a i j x j i


can write as this is going to be lesser than or equal to sum of mode of $a_{ij} x_j$ we have seen this many times before and dividing throughout by mode of x_i . So, we divide throughout both side by mode of x_i . So, bound of $\lambda - a_{ii} x_i$ divided by mode of x_i is going to give me mode of $\lambda - a_{ii}$ this is going to be lesser than or equal to mode of $a_{ij} x_j$ by mode of x_i sum from j equal to one to n .

Not equal to i which is going to be lesser than or equal to sum of mode of a_{ij} why because we have chosen i to be the largest component in the x vector. So, all the x_j and these j are expanding from j equal to one to n j not equal to i . So, all these x_j are going to be less than x_i . So, this is going to be this is going to be less than one this factor mode of x_j mode of x_i is going to be less than one. So, this is going to be less than mode of a_{ij} which is equal to r_i right. So, for the row which corresponds to the largest absolute value of x this is true right this is true for that row thus largest eigen values lays on or within the union of all circle with centers at the diagonal elements of the coefficient matrix a and with the radius as given above this is slightly we have to use our imagination basically we are looking for the for the row which has the largest entry in the x this is true right, but we do not know our priority what is my row with the largest entry of x right i do not know what are my eigen vector. So, how do i know which is going to become my row in the eigen vector with the largest entry of x .

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Gerschgorin's Theorem

- The bound above says that the eigen value will lie on or within a circle centered on the diagonal element corresponding to the largest entry in a given eigen vector x_i
- However, we have no way of knowing a priori the location of the entry with the largest magnitude for a certain eigen vector, since the eigen vector itself is unknown
- Hence it is necessary to take the union of all circles centered about each diagonal element with the radius comprising the sum of the absolute values of the off diagonal elements



So, I have to assume that any of those rows in x can be the row with largest entry and have to compute this circle and the radius for each of those rows right and the union of all those circles is going to give me my upper bound right is that clear.

So, bound above says that eigen value will lay on or within the circle centered on the diagonal element corresponding to the largest entry in the given eigen vector x_i ; however, we have no way of knowing our priory the location of the entry with the largest magnitude for the certain eigen vector since the eigen vector itself is unknown hence it is necessary to know union also we have to take all the rows right you have to take the union of all the circle scented about each diagonal element with the radius comprising the sum of absolute values of the of diagonal element.


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Stability of the eigen values

- Let us consider the effect of perturbations in the elements of a matrix A on its eigen values, denoting the perturbed matrix as $A(\varepsilon) = A + \varepsilon B$
- If x_i are the eigen vectors and λ_i are the eigen values of A then we can write the equations $A x_i = \lambda_i x_i, i = 1, 2, \dots, n$ as:

$$A X = X \Lambda, X = (x_1, x_2, \dots, x_n), \Lambda = \text{diag}(\lambda_1, \lambda_2, \dots, \lambda_n)$$
- If A has linearly independent eigen vectors, the matrix X is non-singular. Thus with $C = X^{-1} B X$ we can write:

$$X^{-1} A(\varepsilon) X = X^{-1} A X + \varepsilon X^{-1} B X = \Lambda + \varepsilon C \quad (*)$$



So, that was Gerschgorin's theorem. So, let us use we will use Gerschgorin's theorem look at stability of the eigen value let us consider the affect of perturbation the elements of matrix on its eigen values denoting the perturbed matrix a plus epsilon b . So, I have by original matrix I apply some perturbations on the elements of original matrix by scaling a matrix b with the small number epsilon and adding into my matrix a . So, I am going to know if I know my eigen values of.

A if I know my eigen value of a how what does these tells me about the eigen values of a plus epsilon b first let us assume that a has linearly independent eigen vectors. So, the matrix X is non singular what is the matrix since X I

eigen vectors x_i the eigen values of A we can write the equation $A x_i = \lambda_i x_i$ as A operating on x_i comprises columns n columns each column being in eigen vector of A . So, A operating on x_i is equal to x_i operating on λ_i right basically we are putting all the writing all the eigen value equation together right. So, we are creating a matrix X each column of which is an eigen vector and we are creating the matrix Λ diagonal which is the diagonal matrix and each diagonal matrix is eigen vector we can write it like this and since x_i linearly independent eigen vector the matrix X has full rank because each of its column is a eigen vector in each eigen vector is being linear independent. So, X has got full rank X is invertible right X is non singular. So, we can write. So, we can define a matrix C is equal to $X^{-1} A X$. So, that we can write $X^{-1} A X = C$. So, we are operating on A operating on x_i with A right and then operating on resultant with X^{-1} .

So, $X^{-1} A X$ is going to be $A X^{-1} A X$ right this part plus $\epsilon X^{-1} B X$, but $X^{-1} A X$ is from this $X^{-1} A X$ is nothing, but λ_i right $X^{-1} A X$ is going to give me diagonal matrix Λ . So, $\lambda_i + \epsilon$ $X^{-1} B X$ i define $X^{-1} B X$ is equal to C . So, $X^{-1} A X$ is equal to $\Lambda + \epsilon C$.

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Stability of the eigen values


Suppose $X^{-1}A(\epsilon)X \bar{x}_i = \lambda_i^* \bar{x}_i$ where λ_i^* is an eigen value of $X^{-1}A(\epsilon)X$ and \bar{x}_i is an eigen vector

Then, $A(\epsilon)X \bar{x}_i = \lambda_i^* X \bar{x}_i$

Denoting $X \bar{x}_i = \bar{x}_i$, we have $A(\epsilon) \bar{x}_i = \lambda_i^* \bar{x}_i$

- Thus $A(\epsilon)$ has the same eigen values as $X^{-1}A(\epsilon)X$
- Hence using Gerschgorin's theorem on (*) we have:

$$\left| \lambda_i(\epsilon) - \text{diag}_i(X^{-1}A(\epsilon)X) \right| \leq \sum_{j=1, j \neq i}^n \left| (X^{-1}A(\epsilon)X)_{ij} \right|$$



Suppose $X^{-1} A X$ has got eigen values λ_i^* and eigen vector \bar{x}_i . So, we are now writing the eigen value problem for $X^{-1} A X$ eigen λ_i^* ϵX and we are saying that suppose $X^{-1} A X$ has eigen vector \bar{x}_i and eigen value λ_i^* then $A x_i$ is equal to $\lambda_i^* x_i$ is basically you are bringing this x_i to the

right hand side right $\lambda^* x x^{-1}$ and we have a $\epsilon x x^{-1}$ right. So, if we denote $x x^{-1} \bar{x} x^{-1}$ star as $\bar{x} A x \bar{x}$ is equal to $\lambda^* x \bar{x}$. So, what does this mean this means that a ϵ has the same eigen value as $x^{-1} A x x^{-1} \epsilon$ x why i because the eigen value of $x^{-1} A x$ λ^* and we have just proved that the eigen value of a ϵ are also λ^* thus a ϵ has the same eigen value of $A x^{-1} \epsilon x$ and eigen vector of a ϵ are related to the eigen vector of $x^{-1} A x$ by this operation eigen function of $x^{-1} A x$ by this operation $x x^{-1}$ star. So, if i have the eigen vector $x^{-1} A x$ if i operate that with this matrix x^{-1} am going to get the eigen vector of a ϵ hence using gerschgorins theorem on star basically i am going to use gerschgorins theorem on this equation and what does gerschgorins theorem tell.

Us it tells us that the largest eigen value of this matrix right satisfy this bound right λ_i minus diagonal element of this matrix right must be lesser than or equal to sum of the of diagonal terms right it has to satisfy this equation from gerschgorins theorem diagonal minus of diagonal from which $\lambda_i \epsilon$ which is this right λ ϵ why do you right $\lambda \epsilon$ because we know that λ_i of a λ is nothing, but the eigen values of a ϵ we know that the eigen value of a ϵ as same as the eigen value of $x^{-1} A x$ right.

So, $\lambda_i \epsilon$ is equal minus $\lambda_i \epsilon c_r$ why do we do that well because we call that $x^{-1} A x$ is equal to $\lambda + \epsilon c$ λ is the eigen value of my original a matrix and ϵc where c_i defined as this right. So, we can write this as λ_i minus $\lambda_i \epsilon c_i$ because this is basically the eigen value of diagonal term of this matrix right and this must be lesser than or equal to the sum of the diagonal term which is equal to ϵc_i right because again lets go back again this matrix is diagonal right. So, all the of diagonal term of the $x^{-1} A x$ are given by ϵc right. So, we have this must be lesser than or equal to ϵ mode of c_i j which implies $\lambda_i \epsilon$ minus λ_i must be of order.


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Stability of the eigen values

From which:

$$|\lambda_i(\varepsilon) - \lambda_i - \varepsilon c_{ii}| \leq \sum_{j=1, j \neq i}^n \varepsilon |c_{ij}| \Rightarrow |\lambda_i(\varepsilon) - \lambda_i| = O(\varepsilon)$$

- If A is a symmetric matrix, then X is an orthogonal matrix and can be chosen to be orthonormal. Thus the elements of $C = X^{-1}BX$ are of the same magnitude as B
- Thus the eigen values of the perturbed system are bounded by the perturbations in A . Hence the eigen value problem of a symmetric matrix is always well-conditioned



Epsilon. So, these term are order epsilon. So, different between lambda i epsilon and lambda i must be of order epsilon. So, what does that mean this mean that eigen value of my perturbed matrix lambda i epsilon are different from the eigen value of my original matrix by are by order epsilon right. So, they are of the same order as the perturbations right. So, if lets if a is the symmetric matrix then x is an orthogonal matrix and we can choose x to be a orthogonal matrix by normalizing the eigen vector thus the element of c x inverse b x are of the same magnitude as b why because c is equal to x inverse b x component of column of a each of x each have norm one unit norm.

So, x inverse magnitude norm of x inverse b x of c must be of the same order of b because these thing have each column as unit norm right. So, this c has to have the same norm as must be of the same order as b. So, the eigen values of perturbed system are bounded by the perturbations in a hence the eigen value problem of symmetric matrix is always well condition why do we say it is well condition because i have applied perturbations to a right small perturbations epsilon times b, but that hasn't change my eigen value by a large amount it has only changed it by order epsilon by the order of whatever perturbations i applied right so; that means, the problem is well condition that is the definition of well condition problem a minor change in the input values does not change in does not result in large differences in the solution. So, minor change in my a matrix does not change my eigen value significantly. So, we will continue our discussion of the eigen values next class next lecture. Thank you very much.