

Matrix Solvers
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Lecture – 54
Block Relaxation Method

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Block Relaxation Scheme

In ADI, the coefficient matrix A is decomposed as $A=H+V$ so that the pentadiagonal matrix problem can be decomposed into iterative solution of tridiagonal matrices H and V along horizontal and vertical lines.

Pentadiagonal $AU = f$
 Tridiagonal $HU + VU = f$
 $\rightarrow HU = f - VU^k$
 $\rightarrow VU^k = f^x2 - HU^{k+1/2}$

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Welcome, we have been discussing about different relaxation schemes line relaxation and Block Relaxation schemes. In an with a focus to solve the matrix equation not in a point wise manner; that means, each element of the x factor will be updated through the iterative schemes, not going in that way rather trying to identify a part of the solution vector and try to update it at one go and then do it for different parts.

So, we have discussed about alternate direction implicit method which I said is a line relaxation scheme. In ADI the coefficient matrix A is decomposed into two parts H plus V ; H comes from disiteration in x -direction y comes from this V comes from disiteration of in y -direction.

So, that the pin now the pentadiagonal matrix problem can be decomposed as two problems sort of this is ADI is typically for 2-dimensional Laplace equation problem. So, instead of AU is equal to f we can write HU plus Vu is equal to f and HU is has x -directional disiteration, Vu has y -directional disiteration. So, this is pentadiagonal A x is A is pentadiagonal, H is tridiagonal, V is also tridiagonal.

The idea is first solving $HU = f - Vu^*$ and then solving $Vu^k = f^* - HU^k + \frac{1}{2}(f^* - f^*_{old})$. So, this is broken in a pentadiagonal matrix equation is broken into two tri diagonal matrix equations and we know the tridiagonal matrix equations are much faster to solve.

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Block Relaxation Scheme

In ADI, the coefficient matrix A is decomposed as $A=H+V$ so that the pentadiagonal matrix problem can be decomposed into iterative solution of tridiagonal matrices H and V along horizontal and vertical lines.

Similarly, large matrices can be decomposed into small blocks on which direct solution methods can be used and iterations are performed over block of solution vectors.

This can be of help when the matrix is very large, storage and memory will be easier for the small blocks
Direct inversion of block matrices can be possible

Parallel algorithms (Schwarz) can be designed

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Similarly, large matrices can be decomposed into small blocks on which direct solution methods if not direct solution method some of the iterative solution methods which are easy to implement can be used and iterations are performed over blocks of the solution vectors. So, instead of solving the entire matrix equation we will break the matrix into small blocks and try to solve for each of the small blocks.

And while solving for one particular small block we will assume that the Gauss value is obtained for the vectors which are associated with the other small blocks. Only for the diagonal part we will use the we will solve it for off diagonals will assume the Gauss value for the blocks and then iterations are performed over blocks of solution vectors. This can be a fail when the matrix is very large say a million by million matrix which is very often we encounter million by million matrix.

It will be much difficult to store the matrix we need some storage algorithm to consider the sparsity of the matrix and store the right nonzero values. Because storing zero values will be will add a redundancy in the overall algorithm and also when you are using some

storage algorithm that only nonzero values and their right pointers are stored you have to use a memory management methods also.

So, for a large matrix the storage and memory access will be much difficult much complex say and it can be easy if we break down into small blocks we will only store the small blocks and doing any memory management will be easy also in the small blocks. So, this is not only memory management and by the computer, but it is also important when we think of that thing that a part of memory management is actually being done by the processors and the computer architecture also.

So, it is accessing memory at the RAM, putting some of it in cache data from the cache, starting with a memory location and going along the pointer to the new memory the new element of that memory all these things this will be easy when you break it into small pieces of blocks. Direct inversion of block matrices can be possible, you have seen the direct inversion is usually a costly method except a TDMA we if we have to use something like ALU or some algorithm like Cholesky etcetera. These are costly methods and they take almost and if consider an n by n matrix they take almost n^3 operations.

So, you for a million by million matrix it will be extremely large number of operations. So, you probably cannot do it in any practical purpose, but if we break down into small matrix matrices we will may get 100 by 100 matrices, where n^3 operations is still possible. So, direct inversion can be is possible for block matrices. And parallel algorithms like Schwartz method can be designed idea parallel algorithm is that you divide the big matrix intuition equation into number of small matrix blocks and ask different processors in a computer architecture to solve different blocks in parallel.

Nowadays we get computers with dual core 16 core even 32 core processors 32 processors are presented in the computer. So, ask each processor to look into one block. So, that the small 32 small blocks can be done in parallel then you can save a lot of time. Instead of do going from one to million maybe each processor is going 1 to 1000 rows and considering 32 plus 32000 rows are done in the time in which you are supposed to 1000 rows. So, this is the parallel process we will discuss about parallelization in few of the later lectures which also can be done once we think of a block partitioning of the matrix or when you think of block relaxation schemes.

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Block Relaxation Scheme

Let us consider the equation $Ax=b$ and decompose the matrix A and the vectors x and b as:

$$A = \begin{pmatrix} A_{11} & A_{12} & A_{13} & \cdots & A_{1p} \\ A_{21} & A_{22} & A_{23} & \cdots & A_{2p} \\ A_{31} & A_{32} & A_{33} & \cdots & A_{3p} \\ \vdots & \vdots & \vdots & \ddots & \vdots \\ A_{p1} & A_{p2} & \cdots & \cdots & A_{pp} \end{pmatrix}, x = \begin{pmatrix} \xi_1 \\ \xi_2 \\ \xi_3 \\ \vdots \\ \xi_p \end{pmatrix}, b = \begin{pmatrix} \beta_1 \\ \beta_2 \\ \beta_3 \\ \vdots \\ \beta_p \end{pmatrix},$$

LHS of the equation corresponding to any row can be now expressed as:

$$(Ax)_i = \sum_{j=1}^p A_{ij}\xi_j,$$

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So, how will it be let us consider an equation Ax is equal to b the matrix A and the vector x and the solution vector and the and the RHS vector b such that A is divided into number of small matrices all A_{11} , A_{12} , A_{13} , A_{pp} all these are matrices. A is decomposed into number of small matrices each is a block matrix and x is decomposed into p blocks ξ_1 , ξ_2 , ξ_3 , ξ_p each may be excess ten thousand elements and each of them there are 10 blocks each has 1000 elements.

So, all these ξ_i 's will be added up to give us x and similarly b is also decomposed into some small blocks. LHS of the equation corresponding to any row now can be expressed as right hand side is summation of A and multiplication of A and x for one particular row will give Ax of that particular row. Ax is nothing, but $A_{ii}\xi_i$ so, what is the first row of this equation if we go back that.

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Block Relaxation Scheme

Let us consider the equation $Ax=b$ and decompose the matrix A and the vectors x and b as:

$$A = \begin{pmatrix} A_{11} & A_{12} & A_{13} & \cdots & A_{1p} \\ A_{21} & A_{22} & A_{23} & \cdots & A_{2p} \\ A_{31} & A_{32} & A_{33} & \cdots & A_{3p} \\ \vdots & \vdots & \vdots & \ddots & \vdots \\ A_{p1} & A_{p2} & \cdots & \cdots & A_{pp} \end{pmatrix}, \quad x = \begin{pmatrix} \xi_1 \\ \xi_2 \\ \xi_3 \\ \vdots \\ \xi_p \end{pmatrix}, \quad b = \begin{pmatrix} \beta_1 \\ \beta_2 \\ \beta_3 \\ \vdots \\ \beta_p \end{pmatrix},$$

$Ax=b \Rightarrow \sum a_{ij} x_j = b_i \Rightarrow (\sum A_{1i} \xi_i = \beta_1) \rightarrow$ 1st row

LHS of the equation corresponding the any row can be now expressed as: $(Ax)_i = \sum_{j=1}^p A_{ij} \xi_j,$



We are solving Ax is equal to b the first row will be $a_{ij} x_j$ is equal to b_i . If we think of a block partitioning this will be this can be also expressed as summation of $A_{1i} x_i$, so, this is first row $a_{1j} x_j$ is equal to this summation of $A_{1i} x_i$ is equal to β_1 once the entire equations first row of that because this is also a matrix also vector equation β_1 has number of elements, the first row of this particular element. So, we can say that the left hand side of this equation which is $a_{1j} x_j$ or a_{ij} for any a particular y can be expressed as $A_{ij} x_j$ into x_j .

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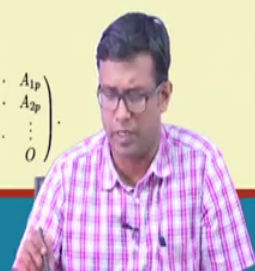
Block Relaxation Scheme- Basic Iterative Schemes

Let us split A similar to the splitting for Jacobi/GS/SOR scheme for the full matrix:

$A=D-E-F$ with D as diagonal, E as a lower triangular and F is an upper triangular matrix

In block decomposition:

$$D = \begin{pmatrix} A_{11} & & & \\ & A_{22} & & \\ & & \ddots & \\ & & & A_{pp} \end{pmatrix},$$

$$E = - \begin{pmatrix} 0 & & & \\ A_{21} & 0 & & \\ \vdots & \vdots & \ddots & \\ A_{p1} & A_{p2} & \cdots & 0 \end{pmatrix}, \quad F = - \begin{pmatrix} 0 & A_{12} & \cdots & A_{1p} \\ 0 & 0 & \cdots & A_{2p} \\ & & \ddots & \vdots \\ 0 & & & 0 \end{pmatrix}.$$



So, let us split a into A is similar to splitting of Jacobi GS or SOR scheme of the full matrix, that is A is equal to D minus E minus F; D is the diagonal, E is the lower triangular, F is upper triangular matrix. Based on the our scheme whether it is Jacobi or Gauss Seidel we will use will take you multiply the last Gauss value with E F or E E plus F we will say that, but we can revise the discussion on basic iterative methods. This is the standard splitting for Jacobi or Gauss Seidel.

In a block decomposition D will be the blocks containing the diagonal terms. Earlier in a general splitting of a for a basic iterative point relaxation scheme A only contain D only contain the diagonal elements. Here instead of diagonal elements they are containing the diagonal blocks E contain the lower triangular blocks, F contain the upper triangular blocks and E and F both has diagonal element 0. So, instead of splitting it by element we are splitting it by blocks of matrices.

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Block Relaxation Scheme- Basic Iterative Schemes

The Jacobi iteration step will be

$$Dx_{k+1} = (E + F)x_k + b$$

If we consider the block-splitting of the matrix

$$D = \begin{pmatrix} A_{11} & & & \\ & A_{22} & & \\ & & \ddots & \\ & & & A_{pp} \end{pmatrix},$$

$$E = - \begin{pmatrix} 0 & & & \\ A_{21} & 0 & & \\ \vdots & \vdots & \ddots & \\ A_{p1} & A_{p2} & \cdots & 0 \end{pmatrix}, \quad F = - \begin{pmatrix} 0 & A_{12} & \cdots & A_{1p} \\ 0 & 0 & \cdots & A_{2p} \\ \vdots & \vdots & \ddots & \vdots \\ 0 & 0 & \cdots & 0 \end{pmatrix}.$$

$$A_{ii}\xi_i^{(k+1)} = ((E + F)x_k)_i + \beta_i$$

Note the structure of the diagonal matrix D and also using the fact that E and F has diagonal entries=0

Using the last relation: $\xi_i^{(k+1)} = A_{ii}^{-1} ((E + F)x_k)_i + A_{ii}^{-1}\beta_i, \quad i = 1, \dots, p,$

This equation is same as the basic Jacobi iteration step

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So, the Jacobi iteration step is $Dx_{k+1} = (E + F)x_k + b$. If we consider the block splitting of the matrix this will be $A_{11} A_{22}$ etcetera. So, this is one block of matrix is multiplied with the solution value that we have to find out and $E + F$ will be multiplied with x_k which is the last available value. So, this is $A_{ii} x_i^{(k+1)} = (E + F)x_k + \beta_i$. Here we are writing x_k not we are writing x_i because x_k contains the entire solution vector and entire solution vector has to be multiplied with E, E and F and each E

and F has diagonal 0. So, that particular vector which is x_i^k is now being multiplied with E and F .

Note that the structure of the diagonal matrix D and the that has been utilized here that the diagonal matrix is A_{ii} is the structure of diagonal matrix is the diagonal blocks here. This matrix is D is not actually a diagonal matrix rather a block diagonal matrix here and we also use the fact that E and F has diagonal entries 0, so, we can write this. Using the last relation the relation we have developed here we can write x_i^{k+1} which is the new updated block of solution vector. This is not a single solution vector rather this is the block of solution vector a number of x 's; number of x 's are involved in this vector is equal to $A_{ii}^{-1} E + F x^k + A_{ii}^{-1} b_i = 1$ to p .

Now, we can see that this is not a direct solver points or relaxation scheme so, that this is D^{-1} . So, this is a single 1 by the diagonal value this is inverse of a matrix. However, as we think of dividing the matrix into large number of matrices a large number of matrix blocks this matrices are basically smaller matrices and we can easily find out their inverse, it will be not that complex. Complexity is usually in terms of what are the order of operations we need to do to find out in that inverse. Inverse is inverse finding is same like Gauss elimination scheme takes n^3 operations.

As we have broken down the matrix into small blocks the number of rows in each block is very small. If you think of million by million matrix maybe you have 1000 rows in each block. So, this will be much less time consuming number of steps will be less to find out A_{ii}^{-1} because this is a small matrix the number n is smaller here. And we have to do it for all the blocks starting from first block $x_1^1, x_2^1, \dots, x_p^1$ if you do it for all the blocks and then again up iterate it this is an iterative method which is discussing.

The this particular equation is very similar as the basic Jacobi iteration step and if we can think this is actually be Jacobi iteration step. Why because for all off diagonals we are using x^k which is the last iterative value and A_{ii}^{-1} instead of A_{ii} we are solving an equation like that a the instead of directly finding inverse we can use something like a Jacobi iteration also. So, this step can be exactly same as Jacobi iteration step if we use Jacobi iteration for finding A_{ii}^{-1} . But, we can because the A_{ii} matrix is smaller we can use some of the direct solution method or some other method to do it also.

However, if Jacobi iteration converges this step should also converge. So, the iteration is not in $A^{-1}b$ explicitly here iteration is on finding out x_i^{k+1} . This solution is done some in some other method, but iteration is what that fact that x_i^{k+1} is found out and this is being iterated again and again. So, this iteration must converge if the Jacobi iteration converges.

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The slide illustrates a block relaxation scheme. On the left, a 6x6 grid of points is shown with three vertical columns circled in red, labeled as blocks in the x-vector. The equation $Au=b$ is written in red. On the right, the FDM matrix for Laplace equation is shown as a 6x6 matrix with diagonal blocks A_{11}, A_{22}, A_{33} and off-diagonal blocks $A_{12}, A_{13}, A_{21}, A_{23}, A_{31}, A_{32}$. The matrix is decomposed as $A = \sum_{i=1,2,3} A_i$. The slide also shows the IIT Kharagpur and NPTEL Online Certification Courses logos.

Here we can see an example of block relaxation this is the domain where we have thirty points and a finite difference matrix, of it is a pentadiagonal matrix. So, along any line there are five elements only and they are near that diagonals two super diagonals two sub diagonal and one diagonal line this is the mesh from which we got the finite difference matrix.

Now, we are doing a partitioning here. This is one group one group this is one group this is one group and these are the three diagonal blocks we will get so, block of x vector. So, that that is $x \times 2$ rather instead of x we should write u we are trying to solve Au vectors we will trying to solve Au is equal to b say.

So, $x_1 u_1$ to u_{12} is one group of block of vectors, $x_2 u_2$ to u_{24} is another group I think I should rewrite it this is x_2 , this is x_3 . So, what is that this is one particular block of vector this is one particular block of vector and this is one particular block of vector. So, what the matrix I will get that matrix will be decomposed into several blocks

and these three are the diagonal blocks which will be directly multiplied with x_1 , x_2 , x_3 when we are solving iterating for x_1 , x_2 , x_3 etcetera.

So, these are the A_{11} , A_{22} , A_{33} the diagonal blocks and there are few of diagonal blocks it is important to look that these two of diagonal blocks are basically 0 matrices. So, we can eliminate some of the off diagonal blocks also from the calculations. And, A the net the actual matrix A when we because we are solving the matrix equation $Au = b$, this A matrix is sum of all the small matrix blocks.

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Block Relaxation Scheme- Example of blocks

FDM matrix for Laplace equation

Blocks in x vector

$$A = \sum_{i=1}^3 \sum_{j=1}^3 A_{ij}$$

It is also possible to form blocks with overlap

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It is also possible to form blocks with the overlap; that means, that maybe this line is being shared by two different blocks. So, you can think of another blocking like this so, this is one block maybe this is another block and or rather let us do it like this.

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The slide is titled "Block Relaxation Scheme- Example of blocks". It features a 4x4 grid of nodes. Red circles highlight three overlapping blocks of nodes. To the right, a block-tridiagonal matrix is shown with blocks $A_{11}, A_{12}, A_{13}, A_{21}, A_{22}, A_{23}, A_{31}, A_{32}, A_{33}$ along the diagonal. Below the grid, three vectors are defined: $\vec{u}_1 = \begin{pmatrix} u_1 \\ u_2 \\ \cdot \\ u_{12} \end{pmatrix}$, $\vec{u}_2 = \begin{pmatrix} u_{13} \\ u_{14} \\ \cdot \\ u_{24} \end{pmatrix}$, and $\vec{u}_3 = \begin{pmatrix} u_{25} \\ u_{26} \\ \cdot \\ u_{36} \end{pmatrix}$. The matrix equation $A = \sum_{i=1}^3 \sum_{j=1}^3 A_{ij}$ is also shown. A small video inset of a man in a pink shirt is visible in the bottom right corner of the slide.

Maybe this is one block this is one block and we can say that this is another block this blocks have some points in points common to them. And we will see that block overlapping blocks are important just to maintain continuity of the solution as well as to have some of the parallels algorithms can only work with overlaps so, you will see about the overlaps later.

The best thing of overlap is that when you will try to solve in this domain this point will see the full stencil or full deciduous row for if there is an overlap considered we will come into the overlap later. But, it is also possible to form the blocks with overlap that is what is important to keep in mind here that this is a typical blocking with non-overlapping blocking what we are showing here is a non overlapping blocking, but we can also have overlapping blocking.

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Block Relaxation Scheme- Formulation

Let V_i be $n \times n_i$ matrix as: $V_i = [e_1, e_2, \dots, e_{n_i}]$ with e_j representing the j -th column of $n \times n$ identity matrix

Let W_i be chosen as another $n \times n_i$ matrix as:

$$W_i = [\eta_1 e_1, \eta_2 e_2, \dots, \eta_{n_i} e_{n_i}]$$

With each η_j representing the weight factor, such that $W_i^T V_i = I_{n \times n_i}$

Handwritten notes in red:
 $\{u\} = \left\{ \begin{matrix} e_1 \\ e_2 \\ \vdots \\ e_{n_i} \end{matrix} \right\}$
 $\left\{ \begin{matrix} \eta_1 e_1 \\ \eta_2 e_2 \\ \vdots \\ \eta_{n_i} e_{n_i} \end{matrix} \right\}$

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Let V we will see the formulation of a block relaxation scheme let V_i be a n into n_i matrix and V_i is basically the columns of identity matrix of up to n_i we will consider in V_i . So, it has it will have n row rows and n_i columns so, there are n_i columns which had e_j 's and they represent the j -th column of an n into n identity matrix. So, it is V_i is a truncated part of an identity matrix.

Where, W_i is chosen and another n into n_i matrix with $\eta_1 e_1, \eta_2 e_2$ etcetera where η_j represents the weight factor such that $W_i^T V_i$ is an identity matrix n_i into n_i . And this is required to map the solutions V_i will be required or as well as W_i to map the solutions x_i to the right x say x_2 I will take which will be after x_1 or we can write that x rather the u matrix sorry u matrix will be x_1, x_2 so on x_1, x_2, \dots, x_p . So, if I find out x_2 that will come in one particular location of u matrix.

So, I have to multiply something with x_2 to make it assembly able to u matrix which comes from these terms $\eta_1 e_1, \eta_2 e_2$. If there is or $e_1 e_2$, if there is overlap then if there is no overlap than $e_1 e_2$ so, x_2 into e_2 will keep x_i here if there is no overlap no overlap it will done there. If there is overlap then maybe this is the location of x_1 and from here we will start the location of x_2 so, you have to multiply something with e_1 .

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Block Relaxation Scheme- Formulation

Let V_i be $n \times n_i$ matrix as: $V_i = [e_1, e_2, \dots, e_{n_i}]$ with e_j representing the j -th column of $n \times n$ identity matrix

Let W_i be chosen as another $n \times n_i$ matrix as:

$$W_i = [\eta_1 e_1, \eta_2 e_2, \dots, \eta_{n_i} e_{n_i}]$$

With each η_j representing the weight factor, such that $W_i^T V_i = I_{n \times n_i}$

For no overlap between the domains, $\eta = 1$

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For no overlap between the domains η is equal to 1 so, the V and W are basically same in these cases.

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Block Relaxation Scheme- Formulation

Now the block A_{ij} , a $n_i \times n_j$ matrix, can be given as: $A_{ij} = W_i^T A V_j$

And similarly; $\xi_i = W_i^T x$, $\beta_i = W_i^T b$

$$x = \sum_{i=1}^{n_{blocks}} V_i \xi_i$$

Each component of Jacobi iteration can be rewritten as projection of the residual in the overlapped domain is zero:

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Now, the block A_{ij} of $n_i \times n_j$ matrix which is an $n_i \times n_j$ matrix can be given as A_{ij} 's W_i transpose $A V_j$. And similarly ξ_i will be given as W_i transpose x i.e. $\xi_i = W_i^T x$. I have used x and u mixed x and u in the discussion. Either x or u is fine, because we started with the problem $A u = b$, but sometimes I confuse myself writing $A x = b$ both are matrix equation be it u be it x is a solution variable.

However, in the present slides that I will upload I will make either x or u uniformly I will make it making by itself. So, nevertheless the xi i will go to the solution vector x we are discussing instead of u using this particular weight matrix W similarly beta will similarly be mapped into b.

So, this is the idea that you have the large matrix b and the part of this the large vector b is say beta q this. So, once you know beta q this will be multiplied with W q transpose b is beta q this will be multiplied like that. So, instead of multiplying W with beta we are doing, but W q is W is transpose multiplied with b that will map b to beta. And, if there are n blocks we can write x is equal to i is equal to one to n V i xi i. So, this is how xi to V x mapping is through V xi 2 x mapping is through V and x to xi mapping is through W they can be same for non overlapping case for overlapping case it will be W will have to carry and weight.

Each component of Jacob iteration can be rewritten as projection of the residual into of the in the overlap domain is equal 0. So, we will get a residual in a in the in a sub domain if there are projections then we will project the if there is overlap then the overlapping of the residual will be obtained. However, some of the residual as projected to the main domain should be 0.

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Block Relaxation Scheme- Formulation

Now the block A_{ij} , a $n_i \times n_j$ matrix, can be given as: $A_{ij} = W_i^T A V_j$

And similarly; $\xi_i = W_i^T x, \beta_i = W_i^T b$

$$x = \sum_{i=1}^{nblocks} V_i \xi_i$$

Each component of Jacobi iteration can be rewritten as projection of residual in the overlapped domain is zero:

$$W_i^T \left[b - A \left(V_i W_i^T x_{k+1} + \sum_{j \neq i} V_j W_j^T x_k \right) \right] = 0.$$

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And that comes as W i transpose b minus V i W i transpose x k plus 1 is if j is not is equal to a n V j W j transpose x k is equal to 0. If there is no overlap V i transpose V i W

if there is no overlap V transpose W is again an identity matrix V and W are same. So, this becomes an identity matrix, but if there is overlap they are not identity matrix. So, this becomes the iteration Jacobi iteration step projecting that the new solutions the residual from the updated solution and non updated solution that residual must be 0 into the overlap domain.

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Block Relaxation Scheme- Formulation

$$W_i^T \left[b - A \left(V_i W_i^T x_{k+1} + \sum_{j \neq i} V_j W_j^T x_k \right) \right] = 0.$$

Using the fact: $\xi_j = W_j^T x$ and rearranging

The iteration step can be rewritten as:

$$\xi_i^{(k+1)} = \xi_i^{(k)} + A_{ii}^{-1} W_i^T (b - Ax_k)$$

The slide also features the IIT Kharagpur and NPTEL Online Certification Courses logos at the bottom, and a small inset image of a lecturer in a pink shirt.

Now, we use the fact that ξ_i is equal to W_j transpose x . So, we multiply W_j W_i transpose with this and write j is equal to i . So, whenever W_i transpose x is there this is ξ_i^k . So, the new if the iteration step will be written as ξ_i^{k+1} which comes from here and this is equal to W_i transpose AV is basically A_{ii} so, A_{ii} . So, this will give us $A_{ii}^{-1} \xi_i^k$ plus 1 is equal to $A_{ij} x^k$ like that.

So, this will be ξ_i^{k+1} is equal to ξ_i^k which will again get here plus $A_{ii}^{-1} W_i^T (b - Ax_k)$. And, this is what we exactly do in a Jacobi iteration step and we get a get the mapping weight function W and then multiply it is transpose with $b - Ax_k$. And, then multiply to A_{ii}^{-1} and add with ξ_i^k and now using this the W is constructed W and V constructed I can get a general algorithm for Jacobi block relaxation scheme.

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Block Iteration Algorithms

General Block Jacobi Iteration

1. For $k = 0, 1, \dots$, until convergence Do:
2. For $i = 1, 2, \dots, p$ Do:
3. Solve $A_{ii}\delta_i = W_i^T(b - Ax_k)$
4. Set $x_{k+1} := x_k + V_i\delta_i$
5. EndDo
6. EndDo

G-S needs less storage and faster

Jacobi can be parallelized due to less data dependency

General Block Gauss-Seidel Iteration

1. Until convergence Do:
2. For $i = 1, 2, \dots, p$ Do:
3. Solve $A_{ii}\delta_i = W_i^T(b - Ax)$
4. Set $x := x + V_i\delta_i$
5. EndDo
6. EndDo

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Which is k is equal to 0 to up to convergence do it up to convergence for each block write $A_{ii}\delta_i = W_i^T(b - Ax_k)$ solve for δ_i and then update $x_{k+1} := x_k + V_i\delta_i$ because δ_i is a solve for $x_{k+1} - x_k$ that is in δ_i . So, you have to multiply V_i with and update x_{k+1} . For better convergence instead of block Jacobi we can think of block Gauss Seidel because each scheme is using the last iteration value of x in block Jacobi in Gauss Seidel instead of using x_k it is using the value x which is the last available value in the neighboring blocks rather.

And, you solve not only neighboring blocks in the in the entire domain the last updated value you solved $A_{ii}\delta_i = W_i^T(b - x)$ and update in each block you update x like this. So, this solution can be done using GS or using Jacobi if we want do a full Jacobi in a block Gauss full Gauss Seidel in a block Gauss Seidel, this solution should also be done using Gauss Seidel.

But, however, when using this particular x it is updated from the neighboring values the neighboring values which the neighboring blocks or the last blocks which are above the x_i 's which are above this particular x_i particular block will have the updated value and the other one will have the last iterate value, this very same as Gauss Seidel.

Gauss Seidel iteration needs less storage because we do not need to storage x_{k+1} in x_k separate leads needs least storage and also faster because, you are using the last

iteration value. However, the Jacobi method can be parallelized due to less data dependency in that sense that once we have the last iteration value we can go to the new Jacobi step. And, if we are doing in different processors none of the processor has to be aware of the fact that what is the updated value in other processors, everybody is using the last iteration value.

Gauss Seidel block Gauss Seidel cannot be parallelized in that way because it is data dependent one particular block has to be aware of what happened to the what is the updated value after the previous blocks in that during that particular iteration the blocks which came before this particular block has been updated. So, it is difficult to paralyze it.

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Block Krylov Subspace Methods

Block operations are often performed using Krylov subspace methods
These are equivalents of the point Krylov space methods for matrix blocks

Block Arnoldi Algorithm

1. Choose a unitary matrix V_1 of dimension $n \times p$.
2. For $j = 1, 2, \dots, m$ Do:
3. Compute $H_{ij} = V_i^T AV_j \quad i = 1, 2, \dots, j$
4. Compute $W_j = AV_j - \sum_{i=1}^j V_i H_{ij}$
5. Compute the Q-R factorization of $W_j: W_j = V_{j+1} H_{j+1,j}$
6. EndDo

This can be compared with the original Arnoldi's method which started with a vector v_1 of unit norm and computed h and w by using vector-vector dot product and matrix vector product respectively

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Similarly, Krylov subspace methods can have block operations and they are often performed using block your called block Krylov subspace methods. They are equivalent of Krylov subspace methods, but they are done for matrix blocks and there is a block Arnoldi method which starts with an unitary matrix V 1 dimension n into p computes the Hessenberg block matrix $V_i^T AV_j$.

And then does an orthogonalization of it; that means, it from AV multiplies V with a a and subtract the VH and then does a Q-R orthogonalization of W to find out W . W is $V_{j+1} H_{j+1,j}$ and comes with all the basis functions of W and V . But, these are not basis of the Krylov subspace rather they are basis of the Krylov block subspaces.

This can be compared with original Arnoldi's method where which started with a vector V_1 instead of starting with an unitary matrix V_1 it started with vector V_1 of unit norm and computed Hessen H and Hessenberg matrix and W by using vector vector dot product. Here with instead of matrix here vector vector dot product is used instead of vector and matrix vector dot product and also a matrix vector product has been used respectively. So, the product products are little the processes are little essentially similar, but instead of starting with a vector v instead of starting with a vector v they are starting with a matrix V_1 that is the difference.

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Block Krylov Subspace Methods

Block Arnoldi Algorithm

1. Choose a unitary matrix V_1 of dimension $n \times p$.
2. For $j = 1, 2, \dots, m$ Do:
3. Compute $H_{ij} = V_1^T A V_j$ $i = 1, 2, \dots, j$
4. Compute $W_j = A V_j - \sum_{i=1}^j V_i H_{ij}$
5. Compute the Q-R factorization of W_j : $W_j = V_{j+1} H_{j+1,j}$
6. EndDo

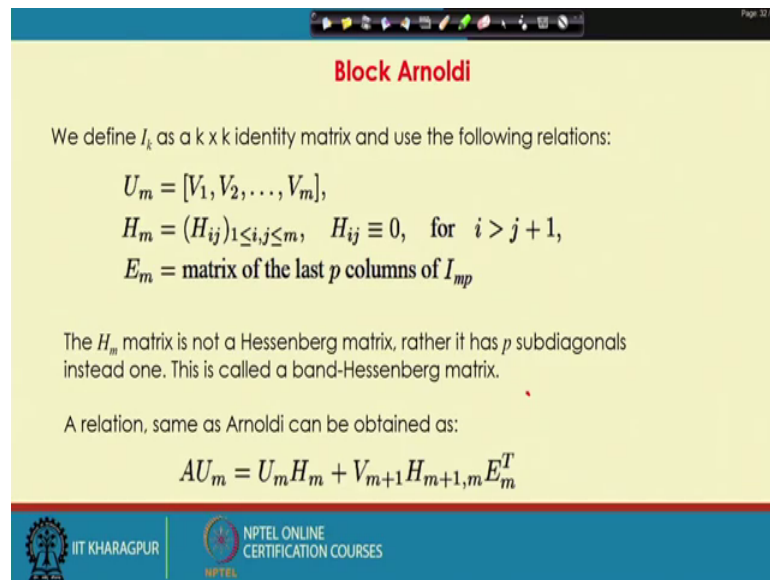
Arnoldi Algorithm

1. Choose a vector v_1 of norm 1
2. For $j = 1, 2, \dots, m$ Do:
3. Compute $h_{ij} = (A v_j, v_i)$ for $i = 1, 2, \dots, j$
4. Compute $w_j := A v_j - \sum_{i=1}^j h_{ij} v_i$
5. $h_{j+1,j} = \|w_j\|_2$
6. If $h_{j+1,j} = 0$ then Stop
7. $v_{j+1} = w_j / h_{j+1,j}$
8. EndDo

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This is block Arnoldi algorithm and this is if we compare with the Arnoldi method it starts with a vector V_1 . So, this is the this these are the difference it starts with a vector. Here it starts with an unitary matrix, here this is compared as a matrix vector product and then it is dot with this. However, there is a matrix product and then it is transpose with H similarly the W_j computing is also W_j compute is also different it blocks subspace. So, we get the basis vectors for a block of Krylov subspace; Krylov subspace of the block matrices.

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Block Arnoldi



We define I_k as a $k \times k$ identity matrix and use the following relations:

$$U_m = [V_1, V_2, \dots, V_m],$$
$$H_m = (H_{ij})_{1 \leq i, j \leq m}, \quad H_{ij} \equiv 0, \quad \text{for } i > j + 1,$$
$$E_m = \text{matrix of the last } p \text{ columns of } I_{mp}$$

The H_m matrix is not a Hessenberg matrix, rather it has p subdiagonals instead one. This is called a band-Hessenberg matrix.

A relation, same as Arnoldi can be obtained as:

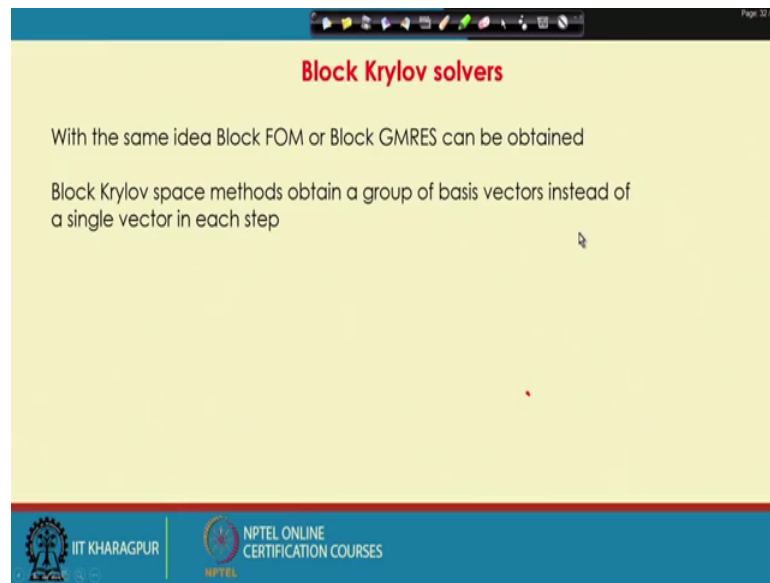
$$AU_m = U_m H_m + V_{m+1} H_{m+1, m} E_m^T$$

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We defined I_k as identity matrix and the and we will see that the following relationships hold same similar type of relationships which we got for simpler or point Arnoldi method. The following relationship shows that Hessenberg matrix H_m is H_m is not a Hessenberg matrix rather it has p sub diagonals and it is called a banded Hessenberg matrix all these are banded matrices.

And, a relationship like this holds which is very similar to our relationship we obtain the to get the Hessenberg relation between the Hessenberg matrix and the Krylov subspace basis vectors. Now, instead of a basis there is a basis matrix instead of the Hessenberg matrix is a band matrix. But, Krylov subspace type of relations are not were applicable for the solution vector x earlier now it is applicable for blocks of the solution vectors x_i 's.

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The slide is titled "Block Krylov solvers" in red text. Below the title, it states: "With the same idea Block FOM or Block GMRES can be obtained". The next line says: "Block Krylov space methods obtain a group of basis vectors instead of a single vector in each step". At the bottom of the slide, there are logos for IIT KHARAGPUR and NPTEL ONLINE CERTIFICATION COURSES.

The same idea with the same idea block full orthogonal method or block GMRES can be obtained. Block Krylov space methods obtain a group of basis vectors instead of a single vector in each step. So, instead of finding out basis vector it gets a basis matrix we can say and the iterations can be done between the blocks of the solution vectors. Within each block I can use even a direct solver, that is how this can give a faster solution or this can be paralyzed also.

So, we have discussed about block relaxation schemes also line relaxation schemes in last two sessions. And, we can see that block relaxation schemes has a potential of paralyzing the solution in a sense that the solution task will be distributed into number of computers which will operate together. And Instead of one computer doing all this n square order or n cube order of operations one computer will do less task and all of the computers we work in parallel, so that the computing tasks can be carried away in less time.

Instead of one people carrying a load if two people instead of one person carrying a load if two people carry the same load it can be much faster. That is the idea block that distribute the matrix into several blocks and use different computers to do it in parallel, that is one great advantage of block partitioning can be this can be exploited for parallel computing and we will look into it in the subsequent classes.

Thank you.