

Numerical Methods and Computation

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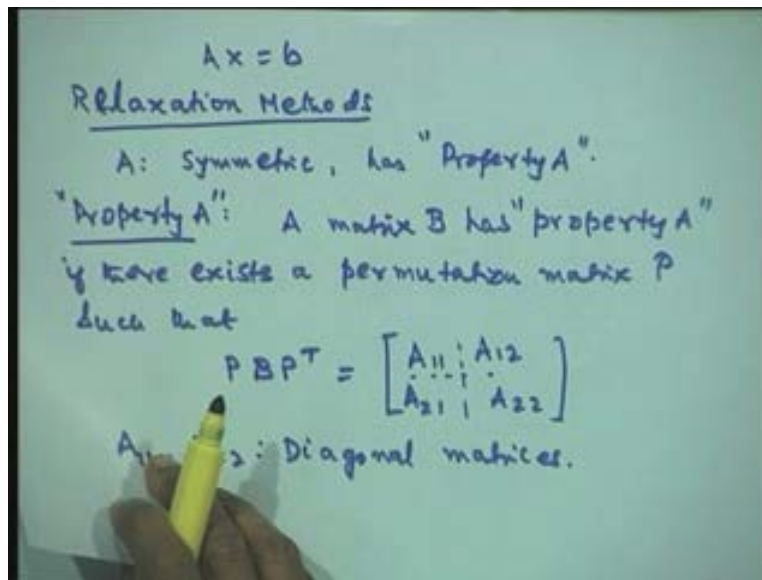
Indian Institute of Technology Delhi

Lecture No # 18

Solution of a System of Linear Algebraic Equations (Continued)

In our last lecture we derived two methods for the solution of the system of linear algebraic equations Ax is equal to b .

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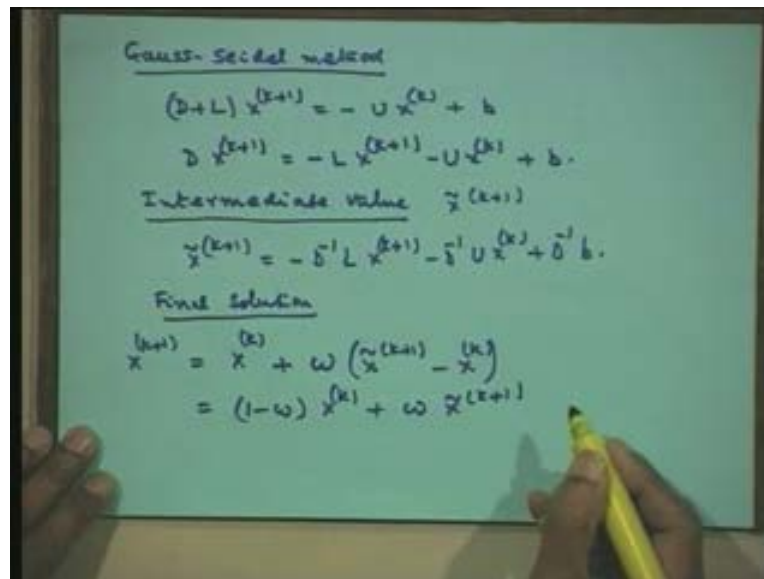
We have we have mentioned that the Gauss Seidel method is at least two times more faster than the Gauss Jacobi method. However in practical applications when we have thousands or lakhs of equations even the convergence of Gauss Seidel is not sufficient, because it will take too much of time. Therefore we need to develop or need to have some methods which works faster than the Gauss Seidel methods. In this direction we have words known as relaxation methods. So we will discuss what is known as the relaxation method. This is a generalization of the Gauss Seidel method. Originally the method was discovered when the matrix has special properties and that is the matrix A has properties which is symmetric and it is has what is known as property A. It's called what is known as property A. This is property A; given the co efficient matrix A , you must able to use inter changes of rows and interchanges of column, that is apply a permutation matrix similar to transformation PAP transpose similarity transformation; such that it can be partitioned as the these two corners. Top and bottom is diagonal matrices and remaining are square matrices or they are any other matrices which is not diagonal.

Let's use another name matrix B as property A. If there exists a permutation matrix P such that PBP transpose can be partitioned as $A_{11} \ A_{12} \ A_{21} \ A_{22}$, where $A_{11} \ A_{22}$ are diagonal matrices, these

are diagonal matrices, these are diagonal matrices; what this really implies is we have pre multiplied by P post multiplied by P transpose. Therefore this gives you interchanges of rows and the corresponding interchanges of columns that is if you interchange column two and five then it will interchange the rows two and five. So it is the corresponding interchanges of this; then I must be able to bring these matrices A_{11} A_{22} as diagonal matrices. Then the matrix is said to have property A, in fact this was done because we will see that the method requires finding the optimal value of a relaxation parameter which gives us fastest convergence for the given method. So that is why we needed this particular formulation to have the analysis of the methods. Therefore without going through whether the matrix property A or now or not even though it is symmetric we use the relaxation methods.

Now let us define what a relaxation method is. We start from the Gauss Seidel method.

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Handwritten notes on a greenboard showing the Gauss-Seidel method equations:

$$\begin{aligned} \text{Gauss-Seidel method} \\ (D+L)x^{(k+1)} &= -Ux^{(k)} + b \\ Dx^{(k+1)} &= -Lx^{(k+1)} - Ux^{(k)} + b. \\ \text{Intermediate value } \tilde{x}^{(k+1)} \\ \tilde{x}^{(k+1)} &= -D^{-1}Lx^{(k+1)} - D^{-1}Ux^{(k)} + D^{-1}b. \\ \text{Final solution} \\ x^{(k+1)} &= x^{(k)} + \omega(\tilde{x}^{(k+1)} - x^{(k)}) \\ &= (1-\omega)x^{(k)} + \omega\tilde{x}^{(k+1)} \end{aligned}$$

We write the Gauss Seidel method as $Dx^{(k+1)} + Lx^{(k+1)} = -Ux^{(k)} + b$. Now let us take the second term $Lx^{(k+1)}$ plus one to the right hand side. So I will write $Dx^{(k+1)}$ plus one is minus $Lx^{(k+1)}$ plus one minus $Ux^{(k)}$ plus b . Now we shall call this solution of this system as an intermediate value. So let us call it as intermediate value. Let us give a new notation to it. Let's call it is as some $\tilde{x}^{(k+1)}$. Therefore $\tilde{x}^{(k+1)}$ is minus $D^{-1}Lx^{(k+1)}$ plus one minus $D^{-1}Ux^{(k)}$ plus $D^{-1}b$. Now this intermediate value has really no meaning as it is. We shall use this intermediate value to construct the final solution and hence I mean we will not specifically say that this is going to approximate our solution or anything here. Now we shall use this intermediate value and the previous iterated value $x^{(k)}$ to build the final solution. So let us call this as a final solution. Let us first take difference between the two values; that difference between the two values will be $\tilde{x}^{(k+1)} - x^{(k)}$. Therefore this will be the difference of the intermediate value and this is the value at the previous iteration. I multiply this by a parameter ω . I will call this as a relaxation parameter, the value of which we shall determine later on using the property or the condition that this iteration should converge. And the analysis of the convergence would give us what will be the ω that should be used in order that the

solution of the problem can be obtained. Now we shall add this product to the previous iterate x_k , this we shall say is our final solution x_k plus one. Now we can use this as a correction using the parameter ω to the previous iterated value to get the current estimate. I can simplify the right hand side and write it as one minus ω into x_k plus ω $x^{(k)} + 1$. Now I want to simplify this further. Therefore I shall substitute the expression for $x^{(k)} + 1$ from here into this particular equation.

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$$\begin{aligned}
 &= (1-\omega)x^{(k)} + \omega[-D^{-1}Lx^{(k+1)} - D^{-1}Ux^{(k)} + D^{-1}b] \\
 Dx^{(k+1)} &= (1-\omega)Dx^{(k)} + \omega[-Lx^{(k+1)} - Ux^{(k)} + b] \\
 (D+\omega L)x^{(k+1)} &= (1-\omega)Dx^{(k)} + \omega[-Ux^{(k)} + b] \\
 &= [(1-\omega)D - \omega U]x^{(k)} + \omega b \\
 x^{(k+1)} &= (D+\omega L)^{-1}[(1-\omega)D - \omega U]x^{(k)} + \omega(D+\omega L)^{-1}b \\
 &= Hx^{(k)} + C \\
 \text{Iteration matrix} \\
 H &= (D+\omega L)^{-1}[(1-\omega)D - \omega U]
 \end{aligned}$$

Therefore I would get here one minus ω x_k plus ω into minus $D^{-1}Lx_k$ plus one minus $D^{-1}Ux_k$ plus $D^{-1}b$. Before I simplify it further I would like to pre multiply this equation by D , so that I can remove this D^{-1} from here. Therefore I would get Dx_k plus one is one minus ω Dx_k plus ω minus Lx_k plus one minus Ux_k plus b . I mean we have pre multiplied throughout by D , so that this will simplify and become minus L . This is minus U and this is simply b . Now this x_k plus one belongs to the left hand side because it is a current iterate, it belongs to the left hand side. So let us bring it to the left hand side, then we will have D plus ωL x_k plus one is equal to one minus ω one plus ω one minus ω into Dx_k plus ω minus Ux_k plus b . Now let us combine this x_k term and this x_k term. So I will have one minus ω into D minus ω into U into x_k plus ω into b . Now the solution of the system is our x_k plus one. Therefore let us inverse this matrix and find x_k plus one. Therefore x_k plus one is minus D plus ωL inverse one minus ω into D minus ωU x_k plus ω D plus ωL inverse of b . Now this is our iteration method, therefore we shall write it as Hx_k plus some c , where this H is iteration matrix given by this particular expression.

Therefore our iteration matrix H is equal to minus D plus ωL inverse. There is no minus sign here, there is a plus sign. There is a plus sign D plus ωL inverse one minus ω of D minus ωU .

Now the properties of this iteration matrix, we shall study it later on. If the method is to converge then the Eigen values of this matrix has to be strictly less than one in magnitude. We are going to

prove this in our later lectures. Therefore what we would like to say now here is that the behavior of this iteration matrix will tell us whether the iterations are going to converge or it's not going to converge. However for computation purposes this particular expression is not useful for us because it is going to take a lot of computer time, therefore we would like to use the error format to derive it. Now let us write the error format of this particular method.

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Handwritten derivation of the error format for the Successive Over-Relaxation (SOR) method. The equations are written on a green board:

$$\begin{aligned} \text{Error Format} \\ x^{(k+1)} &= x^{(k)} + [-I + (D + \omega L)^{-1} \{(1 - \omega)D - \omega U\}] x^{(k)} \\ &\quad + \omega (D + \omega L)^{-1} b \\ &= x^{(k)} + (D + \omega L)^{-1} [-(D + \omega L) + (1 - \omega)D - \omega U] x^{(k)} \\ &\quad + \omega (D + \omega L)^{-1} b \end{aligned}$$

I add and subtract x_k on the right hand side, therefore I can write x_{k+1} is equal to x_k plus minus $I D$ plus ωL inverse one minus ω into D minus ω into $U x_k$ plus ωD plus ωL inverse of b . Now I take D plus ωL inverse outside, so I get $x_k D$ plus ωL inverse into minus D plus ωL plus one minus ω into D minus $\omega U x_k$ plus ωD plus ωL inverse of b .

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Handwritten derivation of the error format and definitions for the SOR method. The equations are written on a green board:

$$\begin{aligned} x^{(k+1)} - x^{(k)} &= (D + \omega L)^{-1} [-\omega A] x^{(k)} + \omega (D + \omega L)^{-1} b \\ &= (D + \omega L)^{-1} \omega [b - A x^{(k)}] \end{aligned}$$

Annotations:

- \downarrow $e^{(k)}$: error
- \downarrow $r^{(k)}$: residual vector

$$(D + \omega L) e^{(k)} = \omega r^{(k)}$$

Definitions:

- $\omega = 1$: Gauss-Seidel method
- ω : relaxation factor
- $0 < \omega < 2$: For Convergence
- $1 < \omega < 2$: Over relaxation
- Successive over relaxation
- $0 < \omega < 1$: Under relaxation

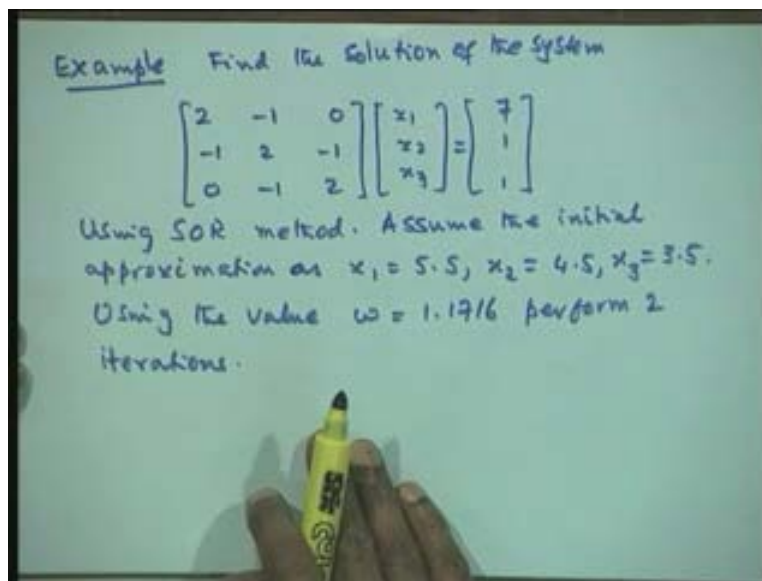
Now let us simplify it further. I can take this x_k to the left hand side and write this as $x_{k+1} - x_k$. Now if you just look at this particular factor there is minus D here and there is plus D here, so minus D cancels with plus D and I have minus ωL minus ωD minus ωU . So what I have here is D plus ωL inverse of minus of ω into A . This is D and D cancels, this is minus ωL minus ωD minus ωU . So I have minus ω common, therefore D plus L plus U that is A x_k plus ωD plus ωL inverse of b ; therefore I can write this as D plus ωL inverse, I can take ω outside and this is $b - Ax_k$. Therefore in the notation that we have used for the Gauss Seidel method we call this as v_k and this as r_k , residual vector. So this is our r_k , the residual vector and this is our error in the solution v_k . Therefore I can write down the method as D plus ωL v_k is equal to ωr_k . So this is our residual vector and this is multiplied by this ω and this D plus ωL , all we have pre multiplied and this is our x_{k+1} . Now you can see that this is in no way computationally expensive than Gauss Seidel except that we have put a factor over here and one more multiplication we put here. Otherwise when ω is equal to one this is same as Gauss Seidel and hence we called this as generalization of Gauss Seidel method.

Now you can see when ω is equal to one we immediately get our Gauss Seidel method here; therefore it is not computationally expensive compared to the Gauss Seidel method at all. Now the ω is called the relaxation factor. The analysis of the method requires that for convergence ω should lie between zero and two. So we should have it such that ω lies between zero and two for convergence. If we use the value greater than one that is above the Gauss Seidel, one less than ω , less than two it is called the over relaxation and it is also called a literature successive over relaxation. And this is one of the most famous and most popular methods. It is simply called SOR, Successive Over Relaxation method. If I use the value between zero and one, it is called under relaxation but it is very rarely used, so this is called the under relaxation. Now therefore this success of the over relaxation procedure would depend on the value relaxation factor that is correctly determined for the properties that I have mentioned earlier. Now symmetric and property A; it is possible for us to say what is the ω optimum value such that this has fastest convergence. It is not unusual to get at least ten times faster than Gauss Seidel in many cases. In fact if you can find the optimum relaxation factor very accurately even getting hundred times faster than Gauss Seidel is not surprising thing for this successive over relaxation.

The successive over relaxation is therefore one of the most important methods for this solution of the linear algebraic equations. There are later on some modifications to make it work better. However the even though the method was derived for the methods having property of A, **later on it has been drop p plus started using** or irrespective of property A, just for symmetric matrices you can use it today even you can try it for the non symmetric matrices also a relaxation

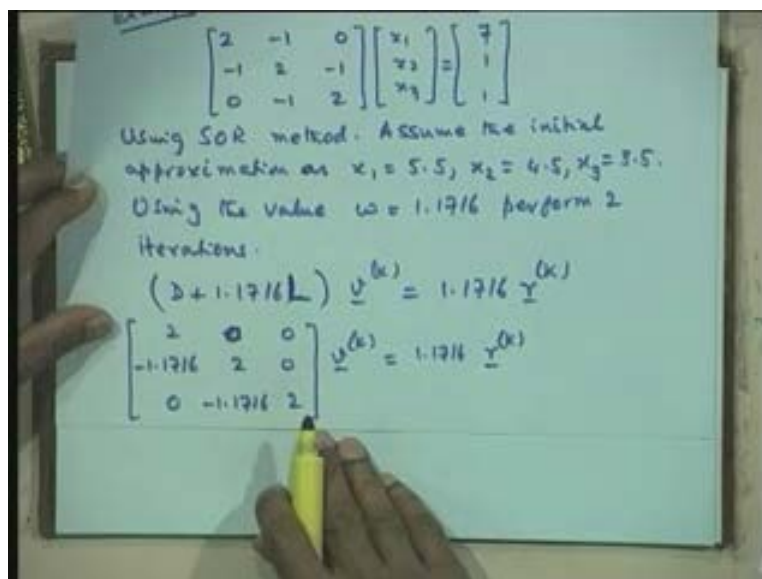
What we are really doing here is we are taking a type of the error between the Gauss Seidel and this one and then multiplying by suitable factor and adding to it in a sense by now improving upon the error in the solution and therefore it would work even for the other cases also. Now before I have a discussion on the error analysis of this, let us take a simple example for the successive over relaxation.

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So let us take an example, so let's call it as - find the solution of the system of equations. Let me take a simple example, 2, -1, 0; -1, 2, -1; 0, -1, 2 and x_1, x_2, x_3 ; 7, 1, 1. Using SOR method assumes the initial approximation as x_1 one is equal to 5.5, x_2 is equal to 4.5, x_3 is equal to 3.5. Using the value will give the value of omega here. Using the value, omega is 1.1716; perform two iterations. Now here we are given the initial approximation. We have also given the omega value and we need to iterations of this as SOR and we shall see later on how we are going to obtain this value of omega.

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Now let us use this error format of writing the system; that is our D plus omega L D plus omega L into vk is equal to omega 1.1716 rk. So D is the same, L is this part; so I am multiplying by omega, so I can write down 2, -1, 0 omega into this minus 1.1716. Therefore 0, - 1.1716, 2; this is our D plus omega times L. This is our diagonal part. This is the part omega into lower

triangular part, that is D plus omega L, ν_k is 1.1716. Now let's write down r_k in the right, let's write down what is r_k in the next step.

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The whiteboard shows the following handwritten equations:

$$\underline{r}^{(k)} = \underline{b} - A \underline{x}^{(k)}$$

$$= \begin{bmatrix} 7 - 2x_1^{(k)} + x_2^{(k)} \\ 1 + x_1^{(k)} - 2x_2^{(k)} + x_3^{(k)} \\ 1 + x_2^{(k)} - 2x_3^{(k)} \end{bmatrix}$$

$$\underline{x}^{(0)} = [5.5 \quad 4.5 \quad 3.5]^T$$

$$\underline{r}^{(0)} = \begin{bmatrix} 0.5 \\ 1 \\ -0.5 \end{bmatrix}$$

$$\begin{bmatrix} 2 & 0 & 0 \\ -1.1716 & 2 & 0 \\ 0 & -1.1716 & 2 \end{bmatrix} \underline{r}^{(0)} = 1.1716 \quad \underline{x}^{(1)} = 1.1716 \begin{bmatrix} 0.5 \\ 1 \\ -0.5 \end{bmatrix}$$

Now the value of r_k (that is the residual vector) is b minus Ax_k , which is b minus Ax_k . So I can write down b minus Ax_k from here; b is this, entire A is this, so as this goes to the right hand side, I will have this as seven minus two times x_1 plus x_2 at the previous iterate x_k one. Then I am taking this to the right hand side. So plus x_1 of k , two times x_2 of k plus x_3 of k . So I have taken these three values of the right hand side and third the component is one; this is plus x_2 of k two times x_3 of k . Now we are given the initial approximation as our vector x_0 is given as 5.5, 4.5 and 3.5. So we are given this as our initial approximation. So I can just substitute it in r_k . This is the value, so I can substitute x_1 is equal to 5.5, x_2 is 4.5 and x_3 is equal to 3.5. Then I would get r_0 . I will give you the value that I obtained here, this is a very simple thing, 0.5, -0.5. So I just multiplied this, and this gives you minus eleven and this gives four point five; this gives you four point five and that is eleven point five, minus eleven. So point five; similarly I get the other values of r . Once the value of r is obtained, I substitute it in this. In this particular step, so the next step I would have here is 2, 0, 0; -1.1716, 2, 0; 0, -1, 1.1716 and the last element is 2; ν_0 is equal to 1.1716, r_0 is 1.1716, 0.5, 1, -0.5.

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$$x^{(0)} = \begin{bmatrix} 5.5 \\ 4.5 \\ 3.5 \end{bmatrix}$$

$$z^{(0)} = \begin{bmatrix} 0.5 \\ 1 \\ -0.5 \end{bmatrix}$$

$$\begin{bmatrix} 2 & 0 & 0 \\ -1.1716 & 2 & 0 \\ 0 & -1.1716 & 2 \end{bmatrix} v^{(0)} = 1.1716 \quad z^{(0)} = 1.1716 \begin{bmatrix} 0.5 \\ 1 \\ -0.5 \end{bmatrix}$$

Forward Substitution

Now this is a forward substitution method. So I can just find out the first component of this v_1 as this is divided by two and then use this value to get the value of v_2 . Then I can get the value of v_3 , so by forward substitution I can get the solution for this problem.

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$$v_1^{(0)} = 0.2929, v_2^{(0)} = 0.7574, v_3^{(0)} = -0.435$$

$$x_1^{(1)} = v_1^{(0)} + x_1^{(0)} = 5.5 + 0.2929 = 5.7929$$

$$x_2^{(1)} = v_2^{(0)} + x_2^{(0)} = 4.5 + 0.7574 = 5.2574$$

$$x_3^{(1)} = v_3^{(0)} + x_3^{(0)} = 3.5 - 0.435 = 3.0650$$

$$z^{(1)} = \begin{bmatrix} 0.6716 \\ -0.6569 \\ 0.1274 \end{bmatrix}$$

$$\begin{bmatrix} 2 & 0 & 0 \\ -1.1716 & 2 & 0 \\ 0 & -1.1716 & 2 \end{bmatrix} v^{(1)} = 1.1716 \begin{bmatrix} 0.6716 \\ -0.6569 \\ 0.1274 \end{bmatrix}$$

So if I do that I would get $v_1^{(0)}$ as 0.2929. I have $v_2^{(0)}$ of zero as 0.7574; $v_3^{(0)}$ of zero is -0.435 . This is the solution of this system by just forward substitution. Now I generate my new value, the computed value which will be equal to; $x_1^{(1)}$ of one is equal to $v_1^{(0)}$ of zero plus $x_1^{(0)}$ of zero. So this is the increment, therefore I will write this as the $x_1^{(1)}$ of zero plus $v_1^{(0)}$ of zero and $x_1^{(0)}$ of zero was given as 5.5 and I have this as 0.2929, so I will have 5.7929. Similarly the second component $x_2^{(1)}$ of one is $v_2^{(0)}$ of zero plus $x_2^{(0)}$ of zero. Now the initial approximation is 4.5 plus 0.7574 and that gives you 5.2574. I have added these two to give 5.2574 and the third component is $x_3^{(1)}$ of one is $v_3^{(0)}$ of zero plus $x_3^{(0)}$ of zero and this value was given as 3.5 minus 0.435 that is 3.060 and this is -4.5.

Once I have obtained this value to go to the next iterate I need again this residual vector. So I go back and substitute this in the residual vector and get the next solution for this one. So if I substitute that I would get this particular value, so I get r_1 , which is as same. By substituting this value of the residual vector I get this as 0.6716, -0.6569, and 0.1274. Now I need to solve this system again. So I will again have to solve this system $2, 0, 0; -1.1716, 2, 0; 0, -1.1716, 2$ and v_1 is 1.1716 into r_1 . So I could now write this as 0.6716, -0.6590 and 0.1274.

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$$\begin{aligned}
 u_1^{(1)} &= 0.3934, & u_2 &= -0.1543, & u_3 &= -0.0158 \\
 x_1^{(2)} &= u_1^{(1)} + x_1^{(1)} = 5.7929 + 0.3934 = 6.1863 \\
 x_2^{(2)} &= u_2^{(1)} + x_2^{(1)} = 5.2574 - 0.1543 = 5.1031 \\
 x_3^{(2)} &= u_3^{(1)} + x_3^{(1)} = 3.0650 - 0.0158 = 3.0492 \\
 \text{Exact: } & x_1 = 6, & x_2 &= 5, & x_3 &= 3.
 \end{aligned}$$

Now we have to solve this equation again by forward substitution and I get the solution of this as v_1 of one, that's at this stage it is v_1 of one is 0.3934. I get v_2 is -0.1543, v_3 three is -0.0158, that is from just computation of this value again by forward substitution. Now I will update my value of x , that is x_1 of two is equal to v_1 of one plus x_1 of one. Now let's go back to this slide. So the value of x_1 was 5.7929 plus 0.3934 that is 6.1863. Then x_2 of two is v_2 of one plus x_2 of one, the value of previous iterate was 5.2574. So I will have 5.2574, -0.1543 that is 5.1030, here it is 31 and lastly x_3 of two is v_3 of one plus x_3 of one the value of x_3 is 3.0653, 0.065 and the value of v_3 is -0.0158 that is 3.0492. Now we have completed two iterations and the exact solution for this problem is x_1 is 6, x_2 is 5, x_3 is equal to 3. Now the interesting observation that we would like to do here is that the convergence is so fast here. Look at this particular value r_0 , this is the error in satisfying our given equations, that is 0.51, -0.51. When we have gone to the first step the error in satisfying the equations was 0.6716, -0.569. Now the solution has now moved faster and now the solution has come almost close to the exact solution the second iteration. Now if you find the error r_2 , you will find that this value is very small. Therefore the convergence compared to the Gauss Seidel is much faster. In fact for this particular problem let us take about three or four decimal places accuracy; Gauss Seidel will require almost a hundred iterations to get that accuracy, whereas you will find in SOR that the required is much less almost a factor or so less than the required number of iterations. So that's why the successive over relaxation method is very fast compared to the other methods but the only hitch here is that we have to find suitable value of ω , which gives the fastest convergence.

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Example Find the solution of the system

$$\begin{bmatrix} 2 & -1 & 0 \\ -1 & 2 & -1 \\ 0 & -1 & 2 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \\ x_3 \end{bmatrix} = \begin{bmatrix} 7 \\ 1 \\ 1 \end{bmatrix}$$

Using SOR method. Assume the initial approximation as $x_1 = 5.5, x_2 = 4.5, x_3 = 3.5$.
Using the value $\omega = 1.1716$ perform 2 iterations.

$(D + \omega L) \underline{y}^{(k)} = \omega \underline{y}^{(k)} + \underline{b}$

$$\begin{bmatrix} 2 & 0 & 0 \\ -1.1716 & 2 & 0 \\ 0 & -1.1716 & 2 \end{bmatrix} \underline{y}^{(k)} = \underline{b}$$

One observation that we should also make here is that when I gave this particular example, which matches exactly over requirements, if I look at this and see its partition, we can apply the permutation matrix such that these are thrown to other way and then we will be able to get the diagonal matrix here and the diagonal matrix here; or you could have another partitions also and then try to attempt that particular way. However as I said it's not necessary nowadays, we can use even if it is a symmetric matrix. Now in order to do the analysis of this even though I will not go deep into how to obtain the omega, or how to obtain for successive over relaxation but we would like to discuss how the Gauss Seidel or the Jacobi or SOR convergence and how do you define by a value that this is the rate of convergence of this particular method; and thereby we can compare the various methods that this is the convergence of the method, therefore it is faster than the other methods. Now to do that I need the background of your matrixes. Let us just revise what we have done earlier and little bit more of the some of the concept about the property of the matrixes. One important property we will need is the norm of a matrix, so let us define the norm of matrix which you might have already done in the course but let's revise it.

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The image shows a whiteboard with handwritten notes in black ink. At the top, it says 'Norm of a matrix $\|A\|$: positive real number' and 'measure of a matrix'. Below this, three properties are listed: $\|A+B\| \leq \|A\| + \|B\|$, $\|AB\| \leq \|A\| \|B\|$, and $\|cA\| = |c| \|A\|$ where c is a real or complex number. The last property is annotated with 'c: real/complex number'. Below these, it says '(i) Euclidean norm' and gives the formula $F(A) = \left[\sum_{i,j=1}^n |a_{ij}|^2 \right]^{1/2}$. A hand holding a yellow marker is visible at the bottom right of the whiteboard.

$$\text{Norm of a matrix } \|A\| : \text{positive real number}$$

measure of a matrix

$$\|A+B\| \leq \|A\| + \|B\|$$
$$\|AB\| \leq \|A\| \|B\|$$
$$\|cA\| = |c| \|A\| \quad c: \text{real/complex number}$$

(i) Euclidean norm

$$F(A) = \left[\sum_{i,j=1}^n |a_{ij}|^2 \right]^{1/2}$$

So I would like to define what's known as norm of a matrix. If given a matrix A, I will denote the norm of the matrix by this, this is a positive real number. This is a positive real number in a sense that it is a measure of a matrix. So it is in a sense actually that we can call this as a measure of a matrix; so we can say given a matrix, this is the norm of this matrix; given another matrix you can say norm of a matrix is this. So you can compare them in terms of norms of the matrixes; so in a sense we can call it as a measure of a matrix. Now if I have two matrixes A and B, then norm of this is your triangular inequality holds here; that is norm of A plus norm of B and if I take the product of two matrixes norm of A into B, this is also norm of A into norm of B. And if I have a complex, any number multiplying, say complex number C into A norm will be where we are talking of C is a real or a complex number. C is a real or a complex number, then this will be equal to magnitude of C into norm of A. This means that if C is a positive or a negative real number or C is complex, then this equality holds. Now let us see what the measures that we can give for this are. The first one norm is called the Euclidean norm. We are talking of a general matrix means real or complex matrix. Therefore what I will do is I will take the magnitude of all the elements, square them, add up and then whatever number that you would get, take the square root of that; that will be called the Euclidean norm. In this you know it is the sort of an Euclidean norm. If you take the distances of points, there is nothing but the distance between the two points is the length. It is the measure, it is Euclidean norm. So that is actually a representation of this in the matrixes. So this is the nothing but I take magnitude of a_{ij} square, then summation over i is equal to j one to n, all of them. Then I take the square root of this and this is called the Euclidean norm. It is usually represented by F of A. So it's a very simple thing that we just take the magnitude of all the elements of the matrix, square them, sum them up and then take the square root of that, a different norm is used in different circumstances.

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$$\|A \cdot B\| \leq \|A\| \|B\|$$

$$\|c A\| = |c| \|A\| \quad c: \text{real/complex number}$$

(i) Euclidean norm

$$F(A) = \left[\sum_{i,j=1}^n |a_{ij}|^2 \right]^{1/2}$$

(ii) Maximum norm

$$\|A\| = \|A\|_{\infty} = \max_i \sum_k |a_{ik}|$$

(max. absolute row sum)

So let us have another norm called the maximum norm. It is denoted by norm of A with a suffix as infinity. I take any row, take the absolute values of all the elements, sum them up, there will be n such row sums. So I will take the largest of them, that is the maximum absolute row; sum is a measure and that is called the maximum norm. So that is denoted by this; that means this is maximum with respect to row I, summation with respect to k; of all a_{ik} , that is these are all the magnitudes of the elements and this is maximum with respect to I; that is row and I am summing up with respect to k. So that is all the values in the row are added in magnitude and then I maximize it with respect to all the rows and then this will be the maximum norm. So I can call this as maximum absolute row sum.

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$$\|A\| = \|A\|_1 = \max_k \sum_i |a_{ik}|$$

(max. absolute column sum)

(iii) Hilbert norm Spectral norm

$$A, \quad A^* = \bar{A}^T, \quad A^* A$$

Find eigenvalues of $A^* A : \lambda_i$

$$\rho(A^* A) = \max_i |\lambda_i|$$

$$\|A\|_2 = \sqrt{\rho(A^* A)}$$

If I can take maximum absolute row sum, I should be able to take the maximum absolute column sum also. Therefore that would also give me another measure which I would call under the maximum norm only; but with a different notation the notation that is used is this; that is

maximum with respect to k , summation with respect to i , a_{ik} . So now I am taking the maximum with respect to column and now I am adding the rows. Therefore this is the maximum absolute column sum and the third norm that we need is the one that is we use very often called the Hilbert norm or it is also called spectral norm or spectral norm. More often we use it as a spectral norm. What we do here is, given a matrix A , I would find its conjugate transpose and we defined it as A^* ; that is your conjugate transpose. Then I form the matrix $A^* A$, then I will find the Eigen values of $A^* A$. We have defined the largest Eigen value of a matrix as its spectral radius, so I will find the largest Eigen values of this. Therefore I will take this as this, let's call it as λ_i . I find Eigen values of this, that is λ_i ; the maximum Eigen value is the spectral radius of this.

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Handwritten notes on a whiteboard:

Hilbert norm Spectral norm

$$A, A^* = \bar{A}^T, A^* A$$

Find eigenvalues of $A^* A : \lambda_i$

$$\rho(A^* A) = \max_i |\lambda_i|$$

$$\|A\|_2 = \sqrt{\rho(A^* A)}$$

If A is symmetric and real $A = A^*$

$$\rho(A^* A) = \rho(A^2) = [\rho(A)]^2$$

$$\|A\|_2 = \rho(A)$$

The Hilbert norm says that I can define the norm as square root of spectral radius of $A^* A$. That means I find the matrix $A^* A$, find its Eigen value, find its spectral radius, square root of that will be the Hilbert norm. Suppose A is symmetric and real; then look at this definition, A^* is conjugate transpose, it's real, therefore it is same as this called symmetric. Therefore A is equal to A^* transpose, that means A is equal to A^* . Therefore that means this is equal to A , is equal to A^* by definition and whereas this spectral radius of $A^* A$ will become spectral radius of A^2 which is equal to A^* ; therefore it will simply be spectral radius of A^2 . Now we know that if λ_i is an Eigen value of A , then λ_i^2 is the Eigen values of A^2 . Therefore spectral radiuses of A^2 of that Eigen values of A to the power of m is λ_i to the power of m ; λ_i is the Eigen values of A . Therefore in this case this spectral norm will be simply spectral radius of A . Therefore if you have a symmetric matrix and real we are very happy because we can just find the Eigen values of that matrix. The largest Eigen values in magnitude will be simply the spectral norm. Now all these norms that we have defined are all independent norms; the Euclidian norms or the maximum norm or the maximum absolute column sum norm or the Hilbert norm, they are all independent definitions. Therefore if you are given a matrix and if you want to find the norm and if you find all of them, since all are independent you can take the smallest of them as a required norm. Suppose say given a matrix A , what is the largest value of norm this matrix has got. If you are finding the

norm of all the four, the smallest of that will be the required bound; that will give the norm of that particular matrix.