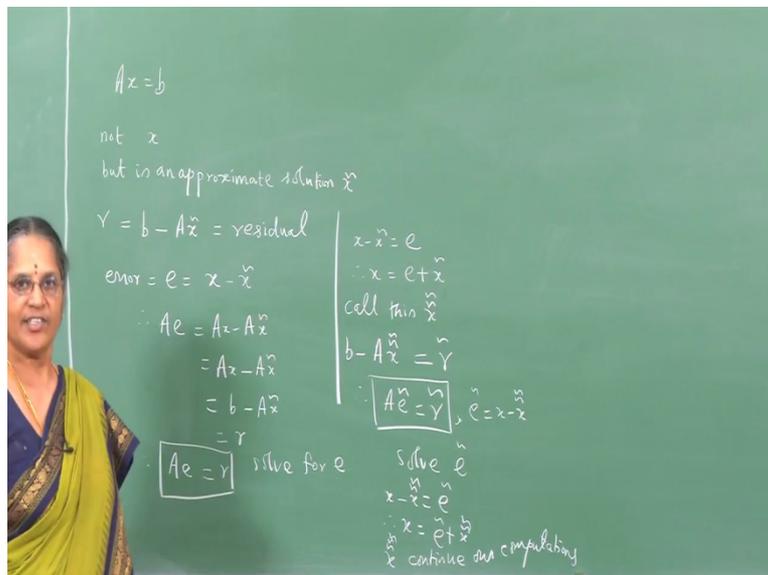


Numerical Analysis.
Professor r. Usha.
Department of Mathematics.
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Lecture-45.
Solution of Linear Systems of equations-8
Iterative Improvement Method Iterative Methods-1.

Good morning everyone, in the last class we discussed the error analysis for the direct methods and we wanted to know whether the ideas that we understood from the error analysis, there is a possibility of improving the accuracy of the approximate solution that we obtained for a system of equations AX equal to b .

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So when we solve the system of equations of the form AX is equal to b by the direct methods, then the solution that we obtained is not the exact solution X but is an approximate solution which we denote by say X tilde due the round-off errors which are incorporated at any stage of our computation. So the question now is can we improve this approximate solution that we obtained using direct methods so that we obtain the solution of the system correct to the desired degree of accuracy. So I compute the A Times X tilde where X tilde is the approximate solution that we obtained using any of the direct methods.

We expect that AX tilde is very close to the right-hand side vector b because we are solving the system AX equal to b . And X tilde being an approximate solution, there will be a difference between b and AX tilde, so I compute this error. Recall b is a vector and A is a

matrix of order $N \times N$, \tilde{X} is an $N \times 1$ vector, so this will be $N \times 1$ vector, b is an $N \times 1$ vector, so $b - A\tilde{X}$ is an $N \times 1$ vector. So I compute this vector and call this as r , namely it is the residual. Now what is the error that has been incorporated? The error if I denote by e is the difference between the exact solution and the approximate solution.

So it is $X - \tilde{X}$ and therefore A times this error is $AX - A\tilde{X}$. And so it is $AX - A\tilde{X}$ is what we have computed using our approximate solution and AX is b , so it is $b - A\tilde{X}$. But what is $b - A\tilde{X}$, that is the residual vector that we obtain. We observe that the equation Ae is equal to r , so there is a relationship connecting the error that one step and the residual at that particular. So e at a particular step which is a vector satisfies the equation Ae is equal to r . So we again have a system of equations in which the unknown vector is e , why is it an unknown e , what is e , e is $X - \tilde{X}$.

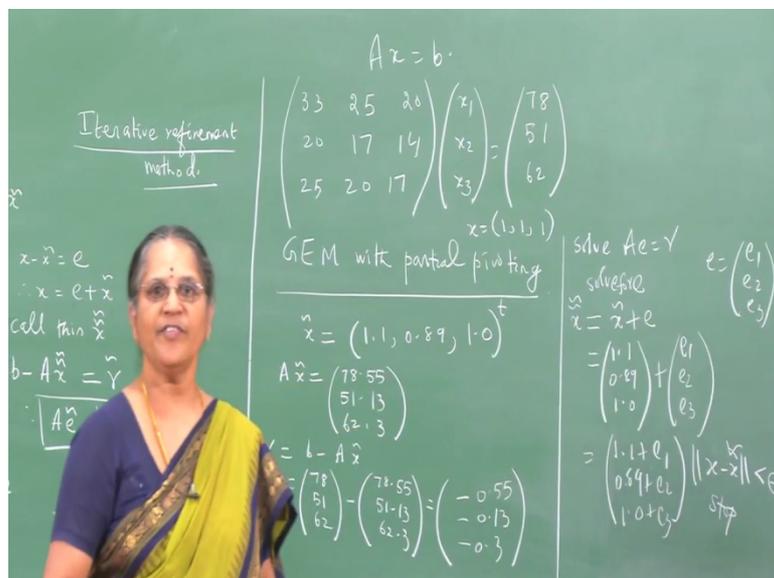
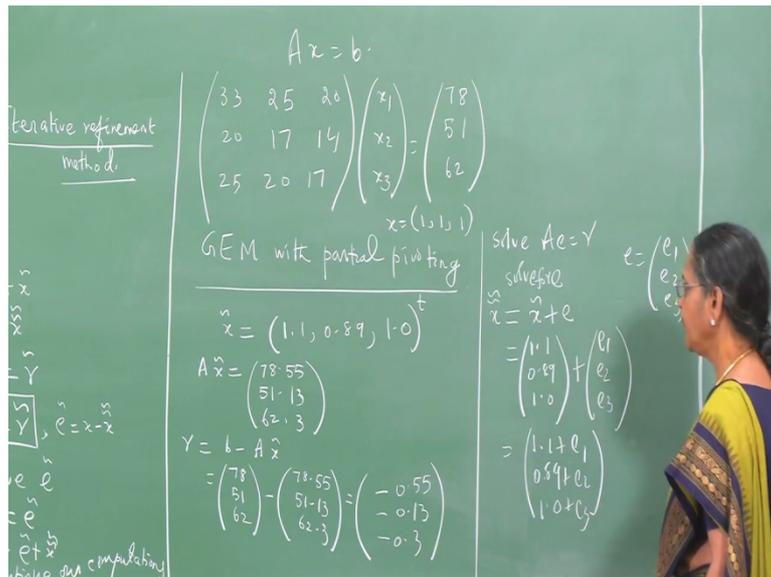
X is not known to us, it is exact solution, that is what we are trying to obtain, so e is an unknown vector, so here e is an unknown vector, so that we solve the system Ae is equal to r . Do we know the right-hand side vector r at this step? Yes, what is it, it is $b - A\tilde{X}$. So we solve the system. How do we solve? Again by any of the direct method that we have learnt and then obtained this error vector e , so once we have e , then we use the fact that the error is $X - \tilde{X}$. So $X - \tilde{X}$ is going to be the e that we have computed and therefore X is equal to $e + \tilde{X}$.

So we call this vector as $\tilde{\tilde{X}}$, we compute what is $b - A\tilde{\tilde{X}}$. If this is such that the $\tilde{\tilde{X}}$ that we computed is within the desired degree of accuracy that we have demanded while computing our solution, then we can take $\tilde{\tilde{X}}$ to be the solution and stop our iterations. If not, what do we do, this is the residual at this step, we call this as \tilde{r} . And therefore we solve the new equation, namely Ae is equal to \tilde{r} . So we compute what is e , what is e , e is $X - \tilde{\tilde{X}}$ which was obtained as the solution at this step.

So again we have a system which satisfies the relationship Ae is equal to \tilde{r} . So we solve for this are known vector again by any of the direct methods that we know. And so we obtain $X - \tilde{\tilde{X}}$ to be e which has been obtained just now, so X is equal to $e + \tilde{\tilde{X}}$. And we check whether the solution that we have obtained at this step is correct to the desired degree of accuracy. If so, we stop our computation, if not, we call this as $\tilde{\tilde{\tilde{X}}}$ and continue our computations till the desired degree of accuracy is achieved.

So the method described here is called the iterative refinement method, which helps us to improve the accuracy of the approximate solution that we have obtained using any of the direct methods. So this method helps us to obtain the solution of the given problem that AX equal to b correct to the desired degree of accuracy. So let us take an example and I will indicate the 1st few steps and you can complete the problem.

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So solve the system AX is equal to b , given that A is a 3 cross 3 matrix having elements as given here, multiplied by the vector X_1, X_2, X_3 and that should be equal to 78, 51, 62, so I have a system AX is equal to b . Say apply Gauss elimination method with partial pivoting, then say I obtain the solution X tilde as 1.1, 0.89 and 1.0 transpose. When I look at the system may immediately see that X_1 equal to 1, X_2 equal to 1 and X_3 equal to 1 is the exact solution,

since $33+25+20$ is 78, $20+17+14$ is 51, $25+20+17$ is 62 and therefore 1, 1, 1 is the exact solution of the system.

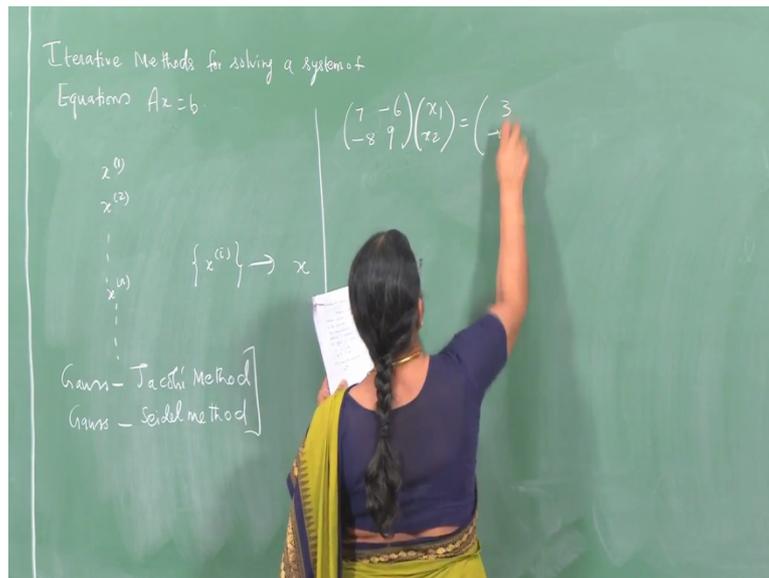
When I apply Gauss elimination method with partial pivoting, I end up with a solution which is this. So it is an approximate solution, so I must improve this approximate solution so that I go closer and closer to the exact solution which is 1, 1, 1. So what should I do according to the iterative refinement method? I must compute A into X tilde. So A is this matrix and X tilde is this vector and so when you compute A into X tilde, that turns out to be vector 78.55, 51.13, 62.3. So I compute what is the residual vector r , what is the residual vector, it is $b - AX$ tilde. And my b is the right-hand side vector 78, 51, 62 - AX tilde is 78.55, 51.13, 62.3. So the residual vector is -0.55, -0.13 and -0.3. So this is my residual vector. So what should I do now?

I should solve the system A into e is equal to this residual vector r . You know what the matrix A is, you also know the right-hand side vector r , so you solve for e , the error vector. So once you get e , then your solution X is going to be this X tilde + e . What is X tilde, X tilde has components 1.1, 0.89 and 1. If suppose you compute your e to have components U_1 , U_2 , U_3 , then add this vector to X tilde, so your new X which I call as X double tilde will be $1.1 + e_1$, $0.89 + e_2$, $1.0 + e_3$. So this is your new solution, this is your new approximation to the solution of the system AX equal to b .

Now check whether $X - X$ double tilde that you have computed is such that the magnitude is less than the prescribed error tolerance. If so, you stop your iterations and you say that X double tilde is your solution correct to the desired degree of accuracy. If not, compute A into X double tilde and then compute $b - AX$ double tilde, call it the residual vector r double tilde and then continue your computations as described in the iterative refinement methods and obtain the solution and see whether that solution is current to the desired degree of accuracy.

So I leave the rest of the computations for you to complete, work out the details. I will give you some more problems on this method in the assignment sheet. So that completes our discussion on the direct methods, so we learnt some direct methods, namely the decomposition methods and Gauss elimination method, Gauss Jordan method and also we incorporated the pivoting strategy in these 2 methods. And then performed some error analysis in the previous class and then using those concepts we have now seen how we can improve the accuracy of the approximate solution that we get using direct methods when we solve a system of equations.

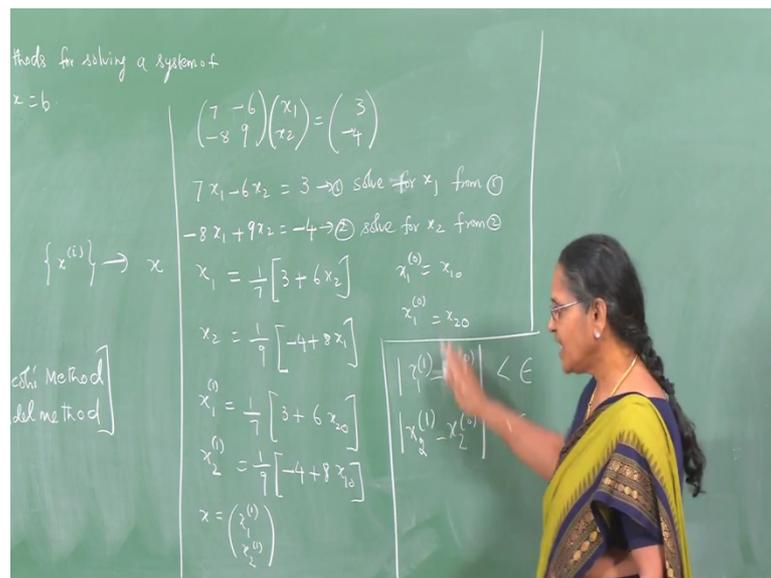
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So now we move onto the iterative methods for solving the system of equations AX equal to b . So our goal now is to develop methods which are iterative in nature for solving a system of equation X equal to b . So what do these methods do? These methods generate successive iterates say X_1, X_2 , etc. X_N and so on for the solution vector X such that this sequence say X_i converges to the solution X . So where do we stop these iterations? We either perform prescribed number of iterations and stop our computations and take the solution obtained at that step to be the solution of the system or we work out the details in such a way that the solution that we obtained at any stage satisfies the required accuracy demand.

So let us see some methods for solving a system of equations iteratively. One of the methods for solving this system is known as Gauss-Jacobi method, the other one that we are going to learn is called Gauss-Seidel method. So we shall describe these methods and see how one can solve a system of equations iteratively. So we shall illustrate it by taking a simple example of a system of 2 equations in 2 unknowns, the procedure is the same for a system of N equations in N unknown.

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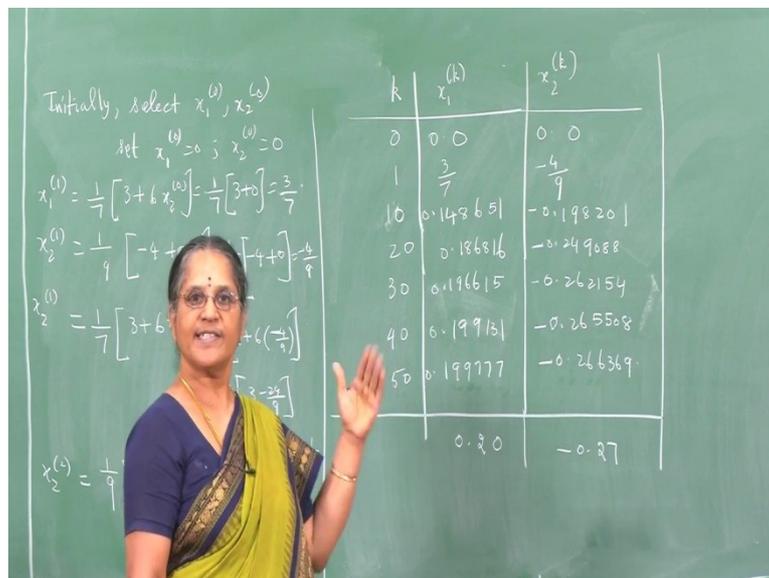
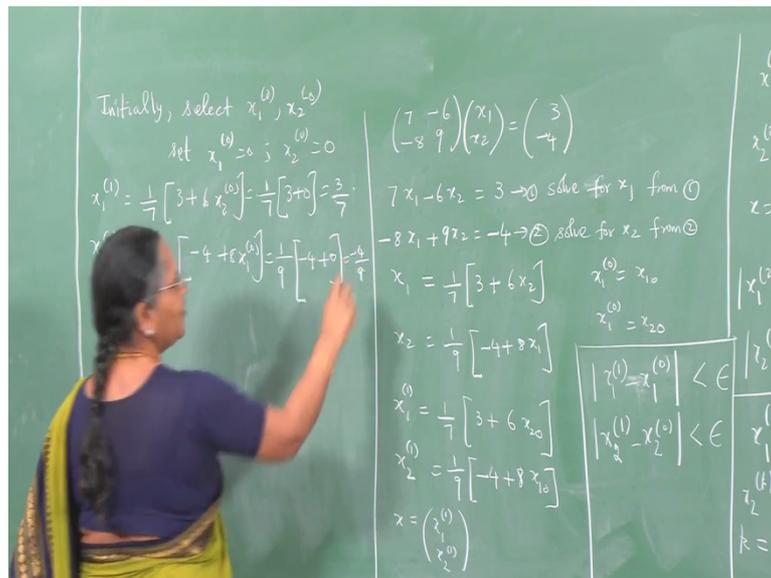


Suppose saying we are given the system of equations $7x_1 - 6x_2 = 3$, $-8x_1 + 9x_2 = -4$. If I write out this equation, then it is $7x_1 - 6x_2 = 3$, $-8x_1 + 9x_2 = -4$. So the procedure for solving this system using Gauss-Jacobi method is as follows. I take the 1st equation to be associated with the 1st unknown, namely x_1 and I solve for x_1 from the 1st equation. So solve for x_1 from equation 1 and similarly I will associate the 2nd unknown x_2 with the 2nd equation and use the 2nd equation and solve for x_2 using the 2nd equation. Then in that case I get x_1 to be equal to $\frac{1}{7} [3 + 6x_2]$, then x_2 is $\frac{1}{9} [-4 + 8x_1]$. This is the 1st step that I do when I want to apply Gauss-Jacobi method.

Now since it is an iterative method, I must start with an initial approximation. So initially I shall take my solution, right, as say $x_1 = 0$ and $x_2 = 0$. And use that solution in the right-hand side and compute what is the solution that I get for the unknown variable. So when I do that, then I get x_1 to be $\frac{1}{7} [3 + 6 \times 0]$ because my initial approximation for x_2 is $x_2 = 0$, this is what I guess for x_2 . So when I substitute for $x_2 = 0$, I have a value for x_1 , so I say that this is the solution that I obtain of the 1st step, namely $x_1 = 1$. Now I go to the 2nd equation, then that is $\frac{1}{9} [-4 + 8 \times 1]$. So I have a number now which gives me x_2 , so I call this as $x_2 = 1$.

And therefore at the 1st I have my solution for the unknown vector x , namely it is $x_1 = 1$ and $x_2 = 1$ and now I continue my computations. Before continuing my iterations, I check whether the absolute value of the solution that I obtained for x_1 at the 1st step - the solution that I assumed for x_1 at the initial step is less than the prescribed error tolerance and I do that also for the 2nd variable, namely $x_2 = 1 - x_2 = 0$, its absolute value is less than epsilon. If both these

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Then I start with an initial approximation for X_1 0, X_2 0, so this is what is known as Gauss-Jacobi method. So let us apply this technique and solve this system. So we initially select X_1 0 and X_2 0. So we simply set them to be 0. And then use that, so what is X_1 1, it is going to be 1 by 7, 3+6 into X_2 0, so it is 1 by 7, 3+0, so 3 by 7. And what is X_2 1, it is 1 by 9 into $-4+8$ times X_1 0. So it is 1 by 9 into $-4+8$ into 0 and so it is -4 by 9. And suppose say we are asked to do 50 iterations and produce the solution, so we have listed in the table the solutions up to 50 iterations that we have performed and see that X_1 at this step turns out to be 0.199777 and X_2 that we have obtained at the 50th iteration as -0.266369 .

So the 2nd variable X_2 has been determined correct to, say 3 decimal, correct to 2 decimal accuracy and the solution is 0.27 correct up to 2 decimal accuracy. And the solution for X_1 is

obtained in this case again correct to 2 decimal accuracy and that turns out to be 0.20. So at the end of 50 iterations, the solutions have been obtained correct to 2 decimal places using Gauss-Jacobi method. So at this stage we shall see what is Gauss-Seidel method for the same system of equations.

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k	$x_1^{(k)}$	$x_2^{(k)}$
0	0.0	0.0
1	$\frac{3}{7}$	$-\frac{4}{9}$
10	0.148651	-0.198201
20	0.186816	-0.249088
30	0.196615	-0.262154
40	0.199151	-0.265508
50	0.199777	-0.266369
	0.20	-0.27

$$\begin{pmatrix} 7 & -6 \\ -8 & 9 \end{pmatrix} \begin{pmatrix} x_1 \\ x_2 \end{pmatrix} = \begin{pmatrix} 3 \\ -4 \end{pmatrix}$$

$7x_1 - 6x_2 = 3$ solve for x_1 from (1)
 $-8x_1 + 9x_2 = -4$ solve for x_2 from (2)

Gauss-Seidel Method (GSM)

$$x_1 = \frac{1}{7} [3 + 6x_2]$$

$$x_2 = \frac{1}{9} [-4 + 8x_1]$$

$(x_1^{(0)}, x_2^{(0)}) = (0, 0)$

$$x_1^{(1)} = \frac{1}{7} [3 + 6x_2^{(0)}] = \frac{1}{7} [3 + 6(0)] = \frac{3}{7}$$

$$x_2^{(1)} = \frac{1}{9} [-4 + 8x_1^{(1)}] = \frac{1}{9} [-4 + 8(\frac{3}{7})] = \frac{1}{9} [-4 + \frac{24}{7}] = \frac{1}{9} [-\frac{4}{7}] = -\frac{4}{63}$$

$|x_1^{(1)} - x_1^{(0)}| < \epsilon$
 $|x_2^{(1)} - x_2^{(0)}| < \epsilon$

Initially, select $x_1^{(0)}, x_2^{(0)}$
 let $x_1^{(0)} = 0; x_2^{(0)} = 0$

$$x_1^{(1)} = \frac{1}{7} [3 + 6x_2^{(0)}] = \frac{1}{7} [3 + 0] = \frac{3}{7}$$

$$x_2^{(1)} = \frac{1}{9} [-4 + 8x_1^{(1)}] = \frac{1}{9} [-4 + 8(\frac{3}{7})] = \frac{1}{9} [-\frac{4}{7}] = -\frac{4}{63}$$

$$x_1^{(2)} = \frac{1}{7} [3 + 6x_2^{(1)}] = \frac{1}{7} [3 + 6(-\frac{4}{63})] = \frac{1}{7} [3 - \frac{24}{21}] = \frac{1}{7} [\frac{37}{21}] = \frac{37}{147}$$

$$x_2^{(2)} = \frac{1}{9} [-4 + 8x_1^{(2)}] = \frac{1}{9} [-4 + 8(\frac{37}{147})] = \frac{1}{9} [-4 + \frac{296}{147}] = \frac{1}{9} [-\frac{4}{147}] = -\frac{4}{1323}$$

k	$x_1^{(k)}$	$x_2^{(k)}$
0	0.0	0.0
1	$\frac{3}{7}$	$-\frac{4}{9}$
10	0.148651	-0.198201
20	0.186816	-0.249088
30	0.196615	-0.262154
40	0.199151	-0.265508
50	0.199777	-0.266369
	0.20	-0.27

Let us look at the equations, which are given to us, namely 7, -6, -8, 9 is the coefficient matrix, the unknown vector is having 2 components X_1, X_2 , the right-hand side vector is 3, -4. So as before then we write down the system of equations, it is $7X_1 - 6X_2 = 3$ and $-8X_1 + 9X_2 = -4$. As before I associate my 1st equation with the 1st unknown, namely X_1 and I solve for X_1 from the 1st equation and associate the 2nd equation with the 2nd unknown X_2 and

solve for X_2 from the 2nd equation. So what do we get, X_1 is going to be $1/7$ into $3+6 X_2$ and X_2 is $1/9$ into $-4+8 X_1$.

Now this is an iterative procedure that I would like to describe, namely it is Gauss-Seidel method. So as before I have to start with an initial approximation for these 2 variables X_1 and X_2 . So suppose you have chosen $X_1 = 0$, $X_2 = 0$ as the initial approximation, then the solution at the 1st step, namely the 1st approximation for X_1 is given by $1/7$ into $3+6$ times $X_2 = 0$, because I have information about X_2 , I have guessed that the solution is $X_2 = 0$. Now I move over to the next iteration, so $X_2 = 1$, which is 1 and into $-4+8$ times, at this stage I see that I have X_1 appearing here but X_1 has already been obtained at this step, namely, the 1st equation gives me X_1 .

So why not I use the currently available solution for X_1 . So I take $X_1 = 1$ for X_1 here and then use that to compute the solution $X_2 = 1$. I recall in Gauss-Jacobi method, we did not use the currently available value for any unknown on the right-hand side. We had any iteration, we always used the solution that was available in the previous step in the previous iteration. Whereas in Gauss-Seidel method, whenever it is possible for us to make use of the currently available solution, it is made use of on the right-hand side to compute the solution at that particular step from the left-hand side. So $X_2 = 1$ is computed as $1/9$ into $-4+8$ times, substitute for X_1 , the currently available value which you have computed just now.

And then now you have to check what is the difference between $X_1 = 1 - X_1 = 0$ in absolute value and $X_2 = 1 - X_2 = 0$ in absolute value. If they satisfy the desired condition for accuracy, then stop your computations and call $X_1 = 1$, $X_2 = 1$ as the solution of the problem, otherwise continue your computations till the desired degree of accuracy is obtained. So let us work out the details for this problem. So if I take the initial approximation as $0, 0$, + I would like to compare my solutions with Gauss-Jacobi method. So at the 1st step what will I get, I get $1/7$ into $3+6$ times $X_2 = 0$ is 0 , so it is $3/7$, that is what we got for $X_1 = 1$ using Gauss-Jacobi method also.

So now I work out $X_2 = 1$ and I see that it is $1/9$ into $-4+8$ times, although initially I have approximated it to be 0 , in the previous step, right, I would like to make use of the solution that is available to me for X_1 currently. Namely I already know that X_1 at this step which is $3/7$, so I would like to make use of this and write down the solution as $1/9$ into $-4+24/7$, so it is $1/9$ into $-28+24/7$, so it is $-4/63$. So you observe that the solution for X_2

even at the 1st step, right, coincides with the solution for X2 that was obtained at the 2nd step using Gauss-Jacobi method.

And so we expect faster convergence using Gauss-Seidel method because it makes use of the currently available solution which is computed at the previous step for an unknown on the right-hand side to compute the solution at this particular step. So we carry out our computations, say we perform 50 iterations and see what our solution is using Gauss-Seidel method. So we present the details in the following table.

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k	$x_1^{(k)}$	$x_2^{(k)}$
0	0	0
1	$\frac{3}{7}$	$-\frac{4}{63}$
2	$\frac{165}{431}$	$\frac{1}{9} \left[8 \left(\frac{165}{431} \right) - 4 \right]$
10	0.219776	-0.249858
20	0.201304	-0.265508
30	0.200086	-0.266590
40	0.200006	-0.266662
50	0.2	-0.266666

k	$x_1^{(k)}$	$x_2^{(k)}$
0	0.0	0.0
1	$\frac{3}{7}$	$-\frac{4}{9}$
10	0.148651	-0.198201
20	0.186816	-0.249888
30	0.196615	-0.262154
40	0.199151	-0.265508
50	0.199777	-0.266369
	0.20	-0.266666

Let us now look at the table for the solutions that we have computed for using Gauss-Seidel method we have performed 50 iterations. And if we still require, say 2 decimal place accuracy for our solution, we observe that this has already been obtained at the 20th iteration. So correct to 2 decimal places, the solution is 0.20 for X1 and 0.27 for X2, whereas for Gauss-Jacobi method, we could achieve that accuracy only at the end of 50 iterations. So Gauss-Seidel method and Gauss-Jacobi method, both generate a sequence of iterates for the solution of the system of equation AX equal to b which converge to the actual solution of the problem.

Gauss-Seidel method converges must faster than Gauss-Jacobi method, the reason being it takes into account the currently available solution at any particular step for the unknowns which have been already computed, as a result the method converges faster. So let us take another example where we have 3 equations into the unknowns and solve the system by both the methods and compare the solutions.

