

lecture_13

March 2, 2017

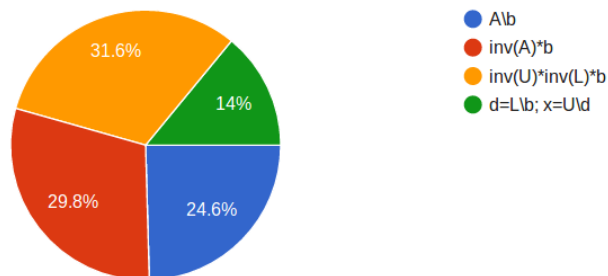
```
In [1]: %plot --format svg
```

```
In [2]: setdefaults
```

0.1 My question from last class

If you are solving the problem $Ax=b$ where $A=LU$ (the lower and upper triangular matrices of A) in matlab or octave what is the most efficient solution?

(57 responses)



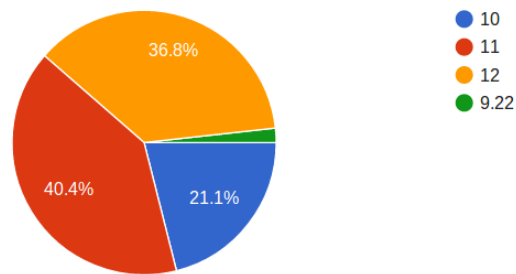
q1

$$A = \begin{bmatrix} 2 & -2 & 0 \\ -1 & 5 & 1 \\ 3 & 4 & 5 \end{bmatrix}$$

0.2 Your questions from last class

1. Do we have to submit a link for HW #4 somewhere or is uploading to Github sufficient?
-no, your submission from HW 3 is sufficient
2. How do I get the formulas/formatting in markdown files to show up on github?
-no luck for markdown equations in github, this is an ongoing request
3. Confused about the $p=1$ norm part and $\|A\|_1$

The $p=1$ norm is the sum of the maximum in each column, what is $\|A\|_1$?
(57 responses)



q2

4. When's the exam?
-next week (3/9)
5. What do you recommend doing to get better at figuring out the homeworks?
-time and experimenting (try going through the lecture examples, verify my work)
6. Could we have an hw or extra credit with a video lecture to learn some simple python?
-Sounds great! how simple?
-Installing Python and Jupyter Notebook (via Anaconda) - <https://www.continuum.io/downloads>
-Running Matlab kernel in Jupyter - <https://anneur.ai.net/2015/11/12/matlab-based-ipython-notebooks/>
-Running Octave kernel in Jupyter - https://anaconda.org/pypi/octave_kernel

1 Markdown examples

" ' ' "

```
x=linspace(0,1);
y=x.^2;
plot(x,y)
for i = 1:10
    fprintf('markdown is pretty')
end
```

1.1 Condition of a matrix

1.1.1 *just checked in to see what condition my condition was in*

1.1.2 Matrix norms

The Euclidean norm of a vector is measure of the magnitude (in 3D this would be: $|x| = \sqrt{x_1^2 + x_2^2 + x_3^2}$) in general the equation is:

$$||x||_e = \sqrt{\sum_{i=1}^n x_i^2}$$

For a matrix, A, the same norm is called the Frobenius norm:

$$||A||_f = \sqrt{\sum_{i=1}^n \sum_{j=1}^m A_{i,j}^2}$$

In general we can calculate any p -norm where

$$||A||_p = \sqrt[p]{\sum_{i=1}^n \sum_{j=1}^m A_{i,j}^p}$$

so the $p=1$, 1-norm is

$$||A||_1 = \sqrt{\sum_{i=1}^n \sum_{j=1}^m A_{i,j}^1} = \sum_{i=1}^n \sum_{j=1}^m |A_{i,j}|$$

$$||A||_\infty = \sqrt{\sum_{i=1}^n \sum_{j=1}^m A_{i,j}^\infty} = \max_{1 \leq i \leq n} \sum_{j=1}^m |A_{i,j}|$$

1.1.3 Condition of Matrix

The matrix condition is the product of

$$Cond(A) = ||A|| \cdot ||A^{-1}||$$

So each norm will have a different condition number, but the limit is $Cond(A) \geq 1$

An estimate of the rounding error is based on the condition of A:

$$\frac{||\Delta x||}{x} \leq Cond(A) \frac{||\Delta A||}{||A||}$$

So if the coefficients of A have accuracy to 10^{-t}

and the condition of A, $Cond(A) = 10^c$

then the solution for x can have rounding errors up to 10^{c-t}

In [7]: $A = [1, 1/2, 1/3; 1/2, 1/3, 1/4; 1/3, 1/4, 1/5]$
 $[L, U] = LU_naive(A)$

A =

1.00000	0.50000	0.33333
0.50000	0.33333	0.25000
0.33333	0.25000	0.20000

L =

1.00000	0.00000	0.00000
0.50000	1.00000	0.00000
0.33333	1.00000	1.00000

U =

1.00000	0.50000	0.33333
---------	---------	---------

```
0.00000    0.08333    0.08333
0.00000   -0.00000    0.00556
```

Then, $A^{-1} = (LU)^{-1} = U^{-1}L^{-1}$

$$Ld_1 = \begin{bmatrix} 1 \\ 0 \\ 0 \end{bmatrix}, Ux_1 = d_1 \dots$$

```
In [8]: invA=zeros(3,3);
        d1=L\[1;0;0];
        d2=L\[0;1;0];
        d3=L\[0;0;1];
        invA(:,1)=U\d1; % shortcut invA(:,1)=A\[1;0;0]
        invA(:,2)=U\d2;
        invA(:,3)=U\d3
        invA*A
```

invA =

```
9.0000    -36.0000    30.0000
-36.0000    192.0000   -180.0000
30.0000   -180.0000    180.0000
```

ans =

```
1.0000e+00    3.5527e-15    2.9976e-15
-1.3249e-14    1.0000e+00   -9.1038e-15
8.5117e-15    7.1054e-15    1.0000e+00
```

Find the condition of A, $cond(A)$

```
In [9]: % Frobenius norm
        normf_A = sqrt(sum(sum(A.^2)))
        normf_invA = sqrt(sum(sum(invA.^2)))

        cond_f_A = normf_A*normf_invA

        norm(A,'fro')

        % p=1, column sum norm
        norm1_A = max(sum(A,2))
        norm1_invA = max(sum(invA,2))
        norm(A,1)

        cond_1_A=norm1_A*norm1_invA
```

```

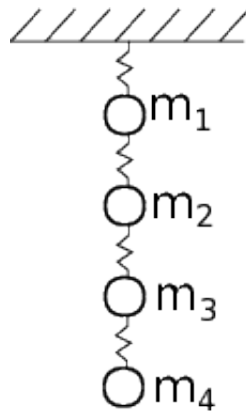
% p=inf, row sum norm
norminf_A = max(sum(A,1))
norminf_invA = max(sum(invA,1))
norm(A,inf)

cond_inf_A=norminf_A*norminf_invA

normf_A = 1.4136
normf_invA = 372.21
cond_f_A = 526.16
ans = 1.4136
norm1_A = 1.8333
norm1_invA = 30.000
ans = 1.8333
cond_1_A = 55.000
norminf_A = 1.8333
norminf_invA = 30.000
ans = 1.8333
cond_inf_A = 55.000

```

Consider the problem again from the intro to Linear Algebra, 4 masses are connected in series to 4 springs with spring constants K_i . What does a high condition number mean for this problem?



Springs-masses

The masses have the following amounts, 1, 2, 3, and 4 kg for masses 1-4. Using a FBD for each mass:

$$m_1 g + k_2(x_2 - x_1) - k_1 x_1 = 0$$

$$m_2 g + k_3(x_3 - x_2) - k_2(x_2 - x_1) = 0$$

$$m_3 g + k_4(x_4 - x_3) - k_3(x_3 - x_2) = 0$$

$$m_4 g - k_4(x_4 - x_3) = 0$$

in matrix form:

$$\begin{bmatrix} k_1 + k_2 & -k_2 & 0 & 0 \\ -k_2 & k_2 + k_3 & -k_3 & 0 \\ 0 & -k_3 & k_3 + k_4 & -k_4 \\ 0 & 0 & -k_4 & k_4 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \\ x_3 \\ x_4 \end{bmatrix} = \begin{bmatrix} m_1 g \\ m_2 g \\ m_3 g \\ m_4 g \end{bmatrix}$$

```

In [10]: k1=10; % N/m
          k2=100000;
          k3=10;
          k4=1;
          m1=1; % kg
          m2=2;
          m3=3;
          m4=4;
          g=9.81; % m/s^2
          K=[k1+k2 -k2 0 0; -k2 k2+k3 -k3 0; 0 -k3 k3+k4 -k4; 0 0 -k4 k4]
          y=[m1*g;m2*g;m3*g;m4*g]

```

K =

```

    100010   -100000         0         0
   -100000    100010        -10         0
         0         -10         11        -1
         0          0         -1         1

```

y =

```

    9.8100
   19.6200
   29.4300
   39.2400

```

```

In [11]: cond(K,inf)
          cond(K,1)
          cond(K,'fro')
          cond(K,2)

```

```

ans =    3.2004e+05
ans =    3.2004e+05
ans =    2.5925e+05
ans =    2.5293e+05

```

```

In [26]: e=eig(K)
          max(e)/min(e)

```

e =

```

    7.9078e-01
    3.5881e+00
    1.7621e+01
    2.0001e+05

```

ans = 2.5293e+05

1.2 P=2 norm is ratio of biggest eigenvalue to smallest eigenvalue!

no need to calculate the inv(K)

2 Iterative Methods

2.1 Gauss-Seidel method

If we have an initial guess for each value of a vector x that we are trying to solve, then it is easy enough to solve for one component given the others.

Take a 3×3 matrix

$$Ax = b$$
$$\begin{bmatrix} 3 & -0.1 & -0.2 \\ 0.1 & 7 & -0.3 \\ 0.3 & -0.2 & 10 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \\ x_3 \end{bmatrix} = \begin{bmatrix} 7.85 \\ -19.3 \\ 71.4 \end{bmatrix}$$
$$x_1 = \frac{7.85 + 0.1x_2 + 0.2x_3}{3}$$
$$x_2 = \frac{-19.3 - 0.1x_1 + 0.3x_3}{7}$$
$$x_3 = \frac{71.4 + 0.1x_1 + 0.2x_2}{10}$$

```
In [12]: A=[3 -0.1 -0.2;0.1 7 -0.3;0.3 -0.2 10]
        b=[7.85;-19.3;71.4]
```

```
x=A\b
```

A =

```
3.00000    -0.10000    -0.20000
0.10000     7.00000    -0.30000
0.30000    -0.20000    10.00000
```

b =

```
7.8500
-19.3000
71.4000
```

x =

```
3.0000
-2.5000
7.0000
```

2.1.1 Gauss-Seidel Iterative approach

As a first guess, we can use $x_1 = x_2 = x_3 = 0$

$$x_1 = \frac{7.85+0.1(0)+0.3(0)}{3} = 2.6167$$

$$x_2 = \frac{-19.3-0.1(2.6167)+0.3(0)}{7} = -2.7945$$

$$x_3 = \frac{71.4+0.1(2.6167)+0.2(-2.7945)}{10} = 7.0056$$

Then, we update the guess:

$$x_1 = \frac{7.85+0.1(-2.7945)+0.3(7.0056)}{3} = 2.9906$$

$$x_2 = \frac{-19.3-0.1(2.9906)+0.3(7.0056)}{7} = -2.4996$$

$$x_3 = \frac{71.4+0.1(2.9906)+0.2(-2.4996)}{10} = 7.00029$$

The results are converging to the solution we found with \ of $x_1 = 3$, $x_2 = -2.5$, $x_3 = 7$

We could also use an iterative method that solves for all of the x-components in one step:

2.1.2 Jacobi method

$$x_1^i = \frac{7.85+0.1x_2^{i-1}+0.3x_3^{i-1}}{3}$$

$$x_2^i = \frac{-19.3-0.1x_1^{i-1}+0.3x_3^{i-1}}{7}$$

$$x_3^i = \frac{71.4+0.1x_1^{i-1}+0.2x_2^{i-1}}{10}$$

Here the solution is a matrix multiplication and vector addition

$$\begin{bmatrix} x_1^i \\ x_2^i \\ x_3^i \end{bmatrix} = \begin{bmatrix} 7.85/3 \\ -19.3/7 \\ 71.4/10 \end{bmatrix} - \begin{bmatrix} 0 & 0.1/3 & 0.2/3 \\ 0.1/7 & 0 & -0.3/7 \\ 0.3/10 & -0.2/10 & 0 \end{bmatrix} \begin{bmatrix} x_1^{i-1} \\ x_2^{i-1} \\ x_3^{i-1} \end{bmatrix}$$

$x_{\{j\}}$	Jacobi method	vs	Gauss-Seidel
$x_1^i =$	$\frac{7.85+0.1x_2^{i-1}+0.3x_3^{i-1}}{3}$		$\frac{7.85+0.1x_2^{i-1}+0.3x_3^{i-1}}{3}$
$x_2^i =$	$\frac{-19.3-0.1x_1^{i-1}+0.3x_3^{i-1}}{7}$		$\frac{-19.3-0.1x_1^i+0.3x_3^{i-1}}{7}$
$x_3^i =$	$\frac{71.4+0.1x_1^{i-1}+0.2x_2^{i-1}}{10}$		$\frac{71.4+0.1x_1^i+0.2x_2^i}{10}$

```
In [14]: ba=b./diag(A) % or ba=b./[A(1,1);A(2,2);A(3,3)]
        sA=A-diag(diag(A)) % A with zeros on diagonal
        sA(1,:)=sA(1,)/A(1,1);
        sA(2,:)=sA(2,)/A(2,2);
        sA(3,:)=sA(3,)/A(3,3)
        x0=[0;0;0];
        x1=ba-sA*x0
        x2=ba-sA*x1
        x3=ba-sA*x2
        fprintf('solution is converging to [3,-2.5,7]\n')
```

ba =

2.6167
-2.7571
7.1400

sA =

```
0.00000  -0.10000  -0.20000
0.10000   0.00000  -0.30000
0.30000  -0.20000   0.00000
```

sA =

```
0.000000  -0.033333  -0.066667
0.014286   0.000000  -0.042857
0.030000  -0.020000   0.000000
```

x1 =

```
2.6167
-2.7571
7.1400
```

x2 =

```
3.0008
-2.4885
7.0064
```

x3 =

```
3.0008
-2.4997
7.0002
```

solution is converging to [3,-2.5,7]

```
In [16]: diag(A)
         diag(diag(A))
```

ans =

```
3
7
10
```

ans =

Diagonal Matrix

```
3  0  0
0  7  0
```

0 0 10

This method works if problem is diagonally dominant,

$$|a_{ii}| > \sum_{j=1, j \neq i}^n |a_{ij}|$$

If this condition is true, then Jacobi or Gauss-Seidel should converge

```
In [15]: A=[0.1,1,3;1,0.2,3;5,2,0.3]
        b=[12;2;4]
        A\b
```

A =

```
0.10000    1.00000    3.00000
1.00000    0.20000    3.00000
5.00000    2.00000    0.30000
```

b =

```
12
 2
 4
```

ans =

```
-2.9393
 9.1933
 1.0336
```

```
In [16]: ba=b./diag(A) % or ba=b./[A(1,1);A(2,2);A(3,3)]
        sA=A-diag(diag(A)) % A with zeros on diagonal
        sA(1,:)=sA(1,+)/A(1,1);
        sA(2,:)=sA(2,+)/A(2,2);
        sA(3,:)=sA(3,+)/A(3,3)
        x0=[0;0;0];
        x1=ba-sA*x0
        x2=ba-sA*x1
        x3=ba-sA*x2
        fprintf('solution is not converging to [-2.93,9.19,1.03]\n')
```

ba =

```
120.000
 10.000
 13.333
```

sA =

```
0   1   3
1   0   3
5   2   0
```

sA =

```
0.00000  10.00000  30.00000
5.00000   0.00000  15.00000
16.66667   6.66667   0.00000
```

x1 =

```
120.000
10.000
13.333
```

x2 =

```
-380.00
-790.00
-2053.33
```

x3 =

```
6.9620e+04
3.2710e+04
1.1613e+04
```

solution is not converging to [-2.93,9.19,1.03]

2.2 Gauss-Seidel with Relaxation

In order to force the solution to converge faster, we can introduce a relaxation term λ .

where the new x values are weighted between the old and new:

$$x^i = \lambda x^i + (1 - \lambda)x^{i-1}$$

after solving for x, lambda weights the current approximation with the previous approximation for the updated x

In [17]: *% rearrange A and b*

```
A=[3 -0.1 -0.2;0.1 7 -0.3;0.3 -0.2 10]
b=[7.85;-19.3;71.4]
```

```
iters=zeros(100,1);
for i=1:100
    lambda=2/100*i;
```

```

        [x,ea,itors(i)]=Jacobi_rel(A,b,lambda);
    end
    plot([1:100]*2/100,itors)

```

A =

```

    3.00000    -0.10000    -0.20000
    0.10000     7.00000    -0.30000
    0.30000    -0.20000    10.00000

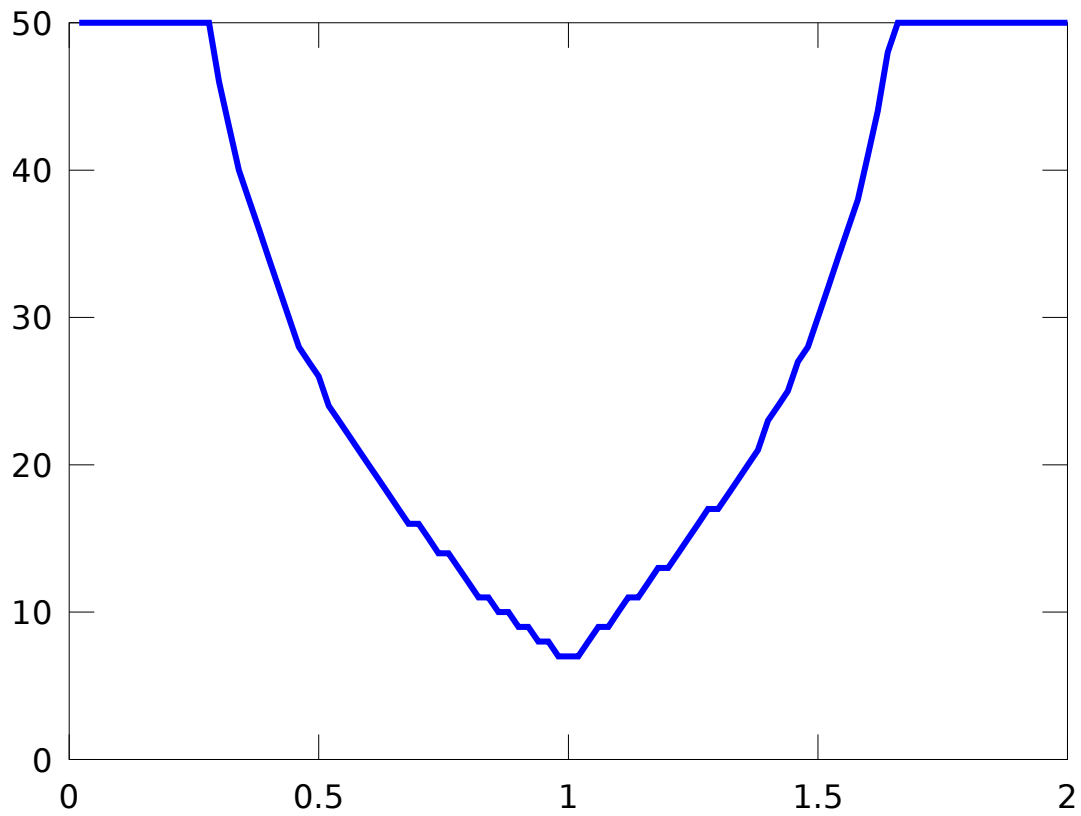
```

b =

```

    7.8500
   -19.3000
    71.4000

```



```

In [107]: l=fminbnd(@(l) lambda_fcn(A,b,l),0.5,1.5)

```

l = 0.99158

```
In [108]: A\b
```

```
ans =
```

```
3.0000  
-2.5000  
7.0000
```

```
In [109]: [x,ea,iter]=Jacobi_rel(A,b,1,0.000001)  
[x,ea,iter]=Jacobi_rel(A,b,1,0.000001)
```

```
x =
```

```
3.0000  
-2.5000  
7.0000
```

```
ea =
```

```
1.8289e-07  
2.1984e-08  
2.3864e-08
```

```
iter = 8
```

```
x =
```

```
3.0000  
-2.5000  
7.0000
```

```
ea =
```

```
1.9130e-08  
7.6449e-08  
3.3378e-08
```

```
iter = 8
```

2.3 Nonlinear Systems

Consider two simultaneous nonlinear equations with two unknowns:

$$x_1^2 + x_1x_2 = 10$$

$$x_2 + 3x_1x_2^2 = 57$$

Graphically, we are looking for the solution:

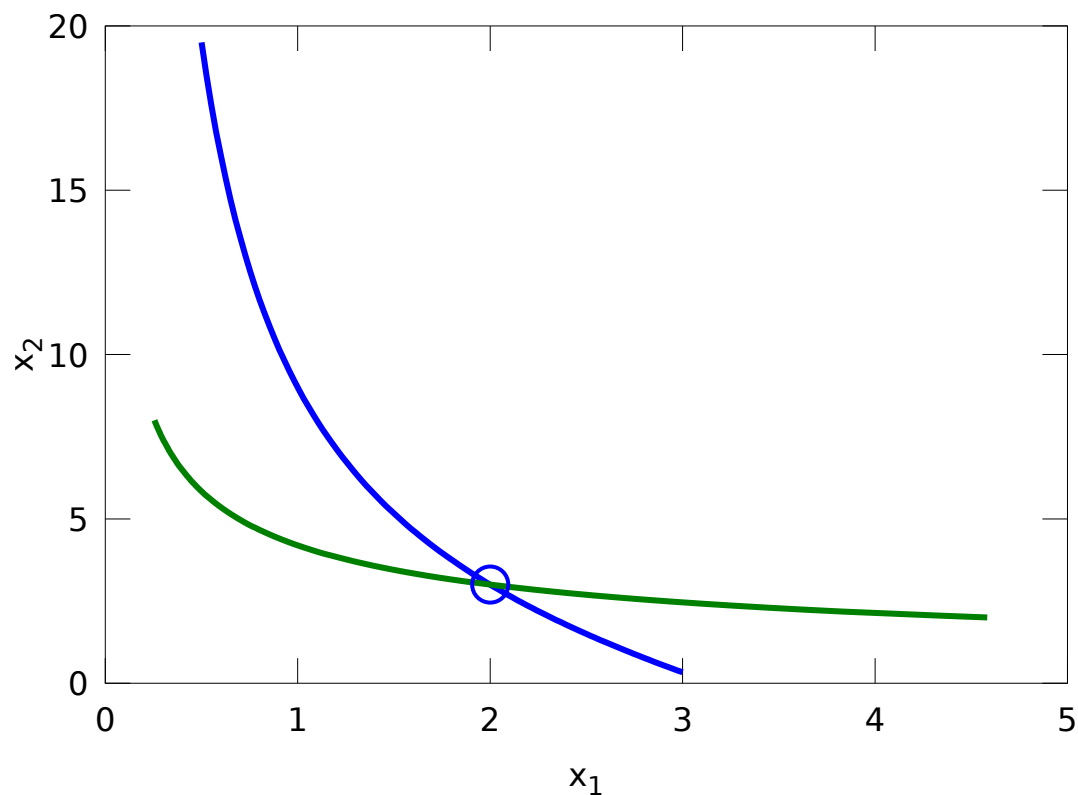
```

In [19]: x11=linspace(0.5,3);
        x12=(10-x11.^2)./x11;

        x22=linspace(2,8);
        x21=(57-x22).*x22.^-2/3;

        plot(x11,x12,x21,x22)
        % Solution at x_1=2, x_2=3
        hold on;
        plot(2,3,'o')
        xlabel('x_1')
        ylabel('x_2')

```



2.4 Newton-Raphson part II

Remember the first order approximation for the next point in a function is:

$$f(x_{i+1}) = f(x_i) + (x_{i+1} - x_i)f'(x_i)$$

then, $f(x_{i+1}) = 0$ so we are left with:

$$x_{i+1} = x_i - \frac{f(x_i)}{f'(x_i)}$$

We can use the same formula, but now we have multiple dimensions so we need to determine the Jacobian

$$[J] = \begin{bmatrix} \frac{\partial f_{1,i}}{\partial x_1} & \frac{\partial f_{1,i}}{\partial x_2} & \dots & \frac{\partial f_{1,i}}{\partial x_n} \\ \frac{\partial f_{2,i}}{\partial x_1} & \frac{\partial f_{2,i}}{\partial x_2} & \dots & \frac{\partial f_{2,i}}{\partial x_n} \\ \vdots & \vdots & & \vdots \\ \frac{\partial f_{n,i}}{\partial x_1} & \frac{\partial f_{n,i}}{\partial x_2} & \dots & \frac{\partial f_{n,i}}{\partial x_n} \end{bmatrix}$$

$$\begin{bmatrix} f_{1,i+1} \\ f_{2,i+1} \\ \vdots \\ f_{n,i+1} \end{bmatrix} = \begin{bmatrix} f_{1,i} \\ f_{2,i} \\ \vdots \\ f_{n,i} \end{bmatrix} + \begin{bmatrix} \frac{\partial f_{1,i}}{\partial x_1} & \frac{\partial f_{1,i}}{\partial x_2} & \dots & \frac{\partial f_{1,i}}{\partial x_n} \\ \frac{\partial f_{2,i}}{\partial x_1} & \frac{\partial f_{2,i}}{\partial x_2} & \dots & \frac{\partial f_{2,i}}{\partial x_n} \\ \vdots & \vdots & & \vdots \\ \frac{\partial f_{n,i}}{\partial x_1} & \frac{\partial f_{n,i}}{\partial x_2} & \dots & \frac{\partial f_{n,i}}{\partial x_n} \end{bmatrix} \left(\begin{bmatrix} x_{i+1} \\ x_{i+1} \\ \vdots \\ x_{i+1} \end{bmatrix} - \begin{bmatrix} x_{1,i} \\ x_{2,i} \\ \vdots \\ x_{n,i} \end{bmatrix} \right)$$

2.4.1 Solution is again in the form $Ax=b$

$$[J]([x_{i+1}] - [x_i]) = -[f]$$

so

$$[x_{i+1}] = [x_i] - [J]^{-1}[f]$$

2.5 Example of Jacobian calculation

2.5.1 Nonlinear springs supporting two masses in series

Two springs are connected to two masses, with $m_1=1$ kg and $m_2=2$ kg. The springs are identical, but they have nonlinear spring constants, of $k_1=100$ N/m and $k_2=-10$ N/m

We want to solve for the final position of the masses (x_1 and x_2)

$$m_1 g + k_1(x_2 - x_1) + k_2(x_2 - x_1)^2 + k_1 x_1 + k_2 x_1^2 = 0$$

$$m_2 g - k_1(x_2 - x_1) - k_2(x_2 - x_1)^2 = 0$$

$$J(1,1) = \frac{\partial f_1}{\partial x_1} = -k_1 - 2k_2(x_2 - x_1) + k_1 + 2k_2 x_1$$

$$J(1,2) = \frac{\partial f_1}{\partial x_2} = k_1 + 2k_2(x_2 - x_1)$$

$$J(2,1) = \frac{\partial f_2}{\partial x_1} = k_1 + 2k_2(x_2 - x_1)$$

$$J(2,2) = \frac{\partial f_2}{\partial x_2} = -k_1 - 2k_2(x_2 - x_1)$$

```
In [ ]: m1=1; % kg
        m2=2; % kg
        k1=100; % N/m
        k2=-10; % N/m^2
```

```
In [20]: function [f,J]=mass_spring(x)
        % Function to calculate function values f1 and f2 as well as Jacobian
        % for 2 masses and 2 identical nonlinear springs
        m1=1; % kg
        m2=2; % kg
        k1=100; % N/m
        k2=-10; % N/m^2
        g=9.81; % m/s^2
        x1=x(1);
        x2=x(2);
        J=[-k1-2*k2*(x2-x1)-k1-2*k2*x1,k1+2*k2*(x2-x1);
           k1+2*k2*(x2-x1),-k1-2*k2*(x2-x1)];
```

```

        f=[m1*g+k1*(x2-x1)+k2*(x2-x1).^2-k1*x1-k2*x1^2;
          m2*g-k1*(x2-x1)-k2*(x2-x1).^2];
    end

```

```
In [21]: [f,J]=mass_spring([1,0])
```

```
f =
```

```

-190.19
 129.62

```

```
J =
```

```

-200   120
 120  -120

```

```
In [22]: x0=[3;2];
        [f0,J0]=mass_spring(x0);
        x1=x0-J0\f0
        ea=(x1-x0)./x1
        [f1,J1]=mass_spring(x1);
        x2=x1-J1\f1
        ea=(x2-x1)./x2
        [f2,J2]=mass_spring(x2);
        x3=x2-J2\f2
        ea=(x3-x2)./x3
        x=x3
        for i=1:3
            xold=x;
            [f,J]=mass_spring(x);
            x=x-J\f;
            ea=(x-xold)./x
        end

```

```
x1 =
```

```

-1.5142
-1.4341

```

```
ea =
```

```

2.9812
2.3946

```

```
x2 =
```

```
0.049894
```



```

0.248638

ea =

31.3492
6.7678

x3 =

0.29701
0.49722

ea =

0.83201
0.49995

x =

0.29701
0.49722

ea =

0.021392
0.012890

ea =

1.4786e-05
8.9091e-06

ea =

7.0642e-12
4.2565e-12

In [23]: x
X0=fsolve(@(x) mass_spring(x),[3;5])

x =

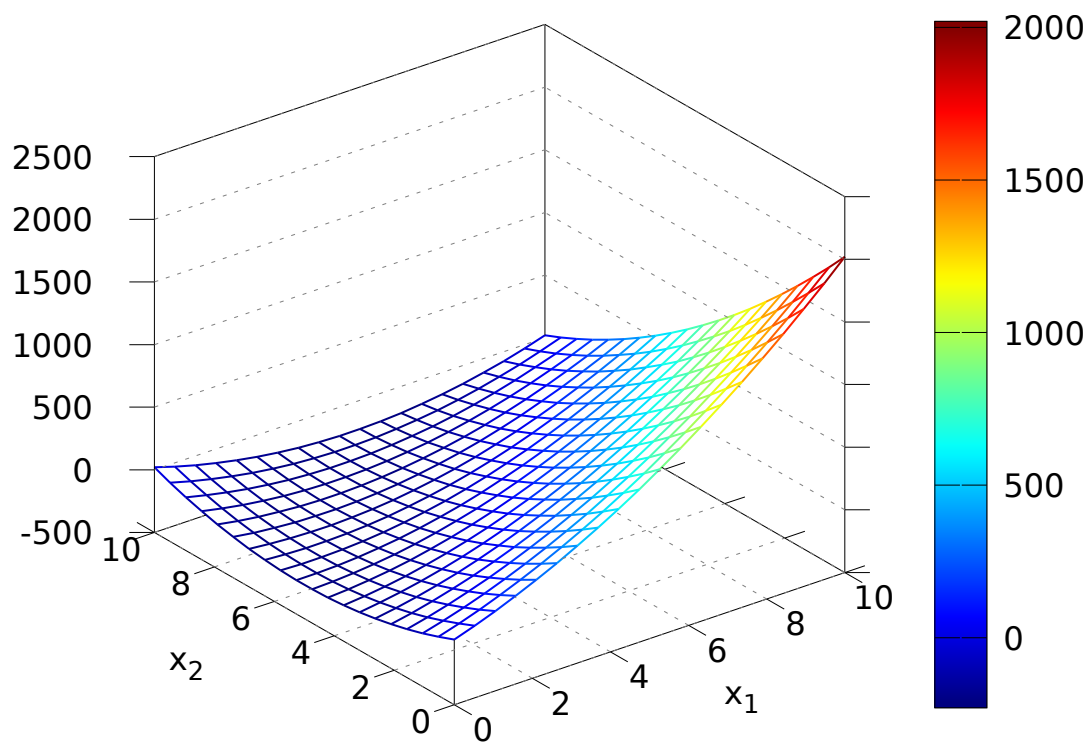
0.30351
0.50372

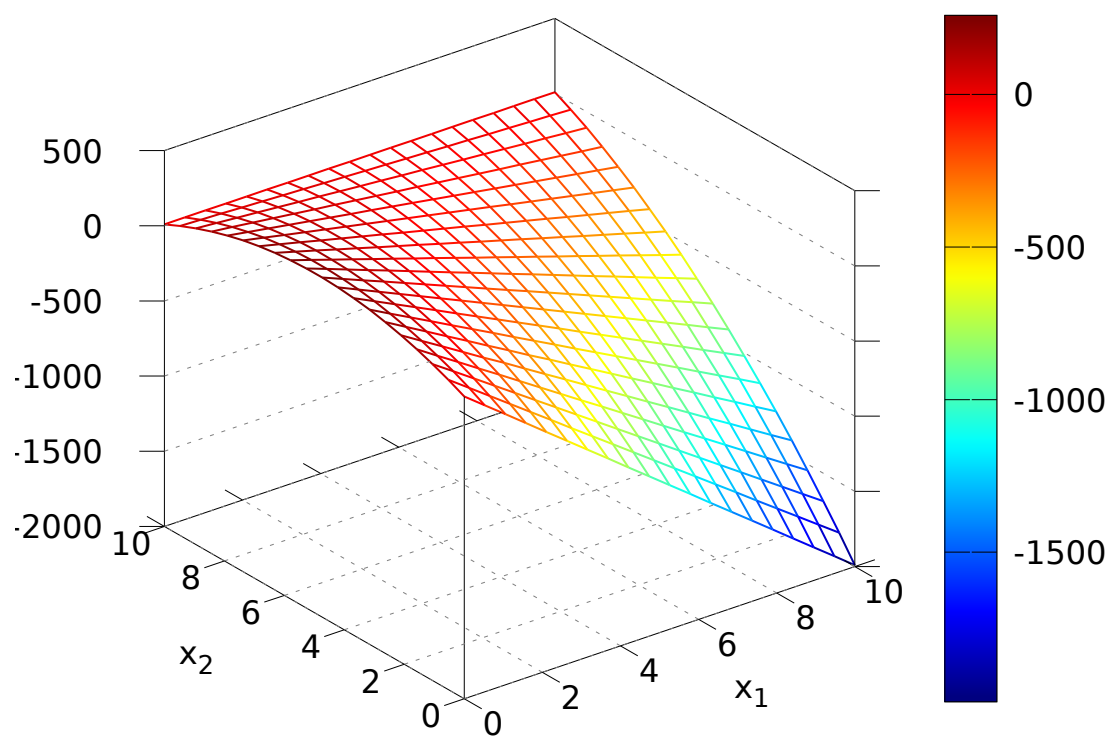
X0 =

```

0.30351
0.50372

```
In [26]: [X,Y]=meshgrid(linspace(0,10,20),linspace(0,10,20));
[N,M]=size(X);
F=zeros(size(X));
for i=1:N
    for j=1:M
        [f,~]=mass_spring([X(i,j),Y(i,j)]);
        F1(i,j)=f(1);
        F2(i,j)=f(2);
    end
end
mesh(X,Y,F1)
xlabel('x_1')
ylabel('x_2')
colorbar()
figure()
mesh(X,Y,F2)
xlabel('x_1')
ylabel('x_2')
colorbar()
```





In []: