

ΣΥΣΤΗΜΑΤΑ ΓΡΑΜΜΙΚΩΝ ΕΞΙΣΩΣΕΩΝ

Matlab-Octave

```
function x = GaussNaive(A,b)
% GaussNaive: naive Gauss elimination
% A = coefficient matrix
% b = right hand side vector
% x = solution vector
[m,n] = size(A);
if m~=n, error('Matrix A must be square'); end
nb = n+1;
Aug = [A b];
% forward elimination
for k = 1:n-1
    for i = k+1:n
        factor = Aug(i,k)/Aug(k,k);
        Aug(i,k:nb) = Aug(i,k:nb)-factor*Aug(k,k:nb);
    end
end
% back substitution
x = zeros(n,1);
x(n) = Aug(n,nb)/Aug(n,n);
for i = n-1:-1:1
    x(i) = (Aug(i,nb)-Aug(i,i+1:n)*x(i+1:n))/Aug(i,i);
end
```

Παράδειγμα

```
>> a= [2 1 2; 1 2 3; 1 2 2]
```

a =

```
2 1 2  
1 2 3  
1 2 2
```

```
>> b= [ 5; 6; 5]
```

b =

```
5  
6  
5
```

```
>> x = GaussNaive(a,b)
```

x =

```
1  
1  
1
```

```

function x = GaussPivot(A,b)
% x = GaussPivot(A,b): Gauss elimination with pivoting.
% A = coefficient matrix
% b = right hand side vector
% x = solution vector
[m,n]=size(A);
if m~=n, error('Matrix A must be square'); end
nb=n+1;
Aug=[A b];
% forward elimination
for k = 1:n-1
% partial pivoting
[big,i]=max(abs(Aug(k:n,k)));
ipr=i+k-1;
if ipr~=k
Aug([k,ipr],:)=Aug([ipr,k],:);
end
for i = k+1:n
factor=Aug(i,k)/Aug(k,k);
Aug(i,k:nb)=Aug(i,k:nb)-factor*Aug(k,k:nb);
end
end
% back substitution
x=zeros(n,1);
x(n)=Aug(n,nb)/Aug(n,n);
for i = n-1:-1:1
x(i)=(Aug(i,nb)-Aug(i,i+1:n)*x(i+1:n))/Aug(i,i);
end

```

Παράδειγμα

```
>> a=[2 1 2; 1 2 3; 1 2 2]
```

a =

```
2 1 2  
1 2 3  
1 2 2
```

```
>> b= [ 5; 6; 5]
```

b =

```
5  
6  
5
```

```
>> x = GaussPivot(a,b)
```

x =

```
1  
1  
1
```

Η Συνάρτηση rref δίνει την ανηγμένη κλιμακωτή μορφή

```
>> A=[1 2 3;4 5 6;7 8 0]
```

A =

```
1 2 3  
4 5 6  
7 8 0
```

```
>> b=[1;2;3]
```

b =

```
1  
2  
3
```

```
>> rref([A b])
```

ans =

```
1.00000 0.00000 0.00000 -0.33333  
0.00000 1.00000 0.00000 0.66667  
0.00000 0.00000 1.00000 -0.00000
```

Η λύση στην τελευταία
Στήλη του αποτελέσματος

Παράδειγμα (2)

```
>> A=[1 1;2 2]
```

```
A =
```

```
1 1
```

```
2 2
```

```
>> b=[2;3]
```

```
b =
```

```
2
```

```
3
```

```
>> rref([A b])
```

```
ans =
```

```
1 1 0
```

```
0 0 1
```

Το σύστημα είναι ασυμβίβαστο
και δεν υπάρχει λύση

Ενσωματωμένη συνάρτηση επίλυση συστήματος

```
>> A=[1 2 3;4 5 6;7 8 0]
```

```
A =
```

```
1 2 3
```

```
4 5 6
```

```
7 8 0
```

```
>> b=[1;2;3]
```

```
b =
```

```
1
```

```
2
```

```
3
```

```
>> x=A\b
```

```
x =
```

```
-0.3333
```

```
0.6667
```

```
0
```

Παραγοντοποίηση πινάκων DooLittle

```
function [L U] = Doolittle(A)
n = size(A,1);
L = zeros(n,n);           % initilize L and U to be n by n
U = zeros(n,n);           % matrices with all zero entries.
for k = 1:n
    L(k,k) = 1;
    for j = k:n
        U(k,j) = A(k,j);
        for s = 1:(k-1)
            U(k,j) = U(k,j) - L(k,s)*U(s,j);
        end
    end
    for i = (k+1):n
        L(i,k) = A(i,k);
        for s = 1:(k-1)
            L(i,k) = L(i,k) - L(i,s)*U(s,k);
        end
        L(i,k) = L(i,k) / U(k,k);
    end
end
end
```

Παράδειγμα

```
>> A=[1 2 3;4 5 6;7 8 0]
```

A =

```
1 2 3  
4 5 6  
7 8 0
```

```
>> [L U] = Doolittle(A)
```

L =

```
1 0 0  
4 1 0  
7 2 1
```

U =

```
1 2 3  
0 -3 -6  
0 0 -9
```

```
function x = DoolittleSystemSolution(A,b)
n = size(A,1);
L = zeros(n,n);          % initilize L and U to be n by n
U = zeros(n,n);          % matrices with all zero entries.
for k = 1:n
    L(k,k) = 1;
    for j = k:n
        U(k,j) = A(k,j);
        for s = 1:(k-1)
            U(k,j) = U(k,j) - L(k,s)*U(s,j);
        end
    end
    for i = (k+1):n
        L(i,k) = A(i,k);
        for s = 1:(k-1)
            L(i,k) = L(i,k) - L(i,s)*U(s,k);
        end
        L(i,k) = L(i,k) / U(k,k);
    end
end
end
```

$\Sigma \cup \nabla \epsilon \times \epsilon | \alpha \dots$

% Step 2: Forward subsitution to solve $L * y = b$

y = zeros(n,1); % Initialize y to be a column vector

y(1) = b(1);

for i = 2:n

 y(i) = b(i);

 for j = 1:(i-1)

 y(i) = y(i) - L(i,j)*y(j);

 end

end

% Step 3: Back subsitution to solve $U * x = y$

x = zeros(n,1); % Initialize x to be a column vector

x(n) = y(n) / U(n,n);

for i = (n-1):-1:1

 x(i) = y(i);

 for j = (i+1):n

 x(i) = x(i) - U(i,j)*x(j);

 end

 x(i) = x(i) / U(i,i);

end

end

Παράδειγμα

```
>> A=[1 2 3;4 5 6;7 8 0]
```

```
A =
```

```
1 2 3  
4 5 6  
7 8 0
```

```
>> b=[1;2;3]
```

```
b =
```

```
1  
2  
3
```

```
>> x = DoolittleSystemSolution(A,b)
```

```
x =
```

```
-0.33333  
0.66667  
-0.00000
```

Παραγοντοποίηση πινάκων Crout

```
function [L,U] = crout(A)
[m, n] = size(A);
L=zeros(n,n);
U=zeros(n,n);
for k = 1 : n
    L(k,k)=A(k,k)-L(k,1:k-1)*U(1:k-1,k);
    for j = k : n
        U(k,j) = ( A(k,j) - L(k,1:k-1)*U(1:k-1,j) )/ L(k,k);
    end;
    for i = k+1 : n
        L(i,k) = (A(i,k) - L(i,1:k-1)*U(1:k-1,k) ) / U(k,k);
    end
end
```

Παράδειγμα

```
>> A=[1 2 3;4 5 6;7 8 0]
```

A =

```
1 2 3  
4 5 6  
7 8 0
```

```
>> [L,U] = crout(A)
```

L =

```
1 0 0  
4 -3 0  
7 -6 -9
```

U =

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

Ενσωματωμένη συνάρτηση LU [L, U, P] = lu(A)

A=[1 2 3;4 5 6;7 8 0]

A =

```
1 2 3
4 5 6
7 8 0
```

>> [L, U, P] = lu(A)

L =

```
1.00000 0.00000 0.00000
0.14286 1.00000 0.00000
0.57143 0.50000 1.00000
```

U =

```
7.00000 8.00000 0.00000
0.00000 0.85714 3.00000
0.00000 0.00000 4.50000
```

P =

Permutation Matrix

```
0 0 1
1 0 0
0 1 0
```

Εναλλαγή γραμμής

Παραγοντοποίηση Cholesky

```
>> A = [2,-1,0;-1,2,-1;0,-1,2]
```

```
A =
```

```
 2 -1  0  
-1  2 -1  
 0 -1  2
```

```
>> chol(A)
```

```
ans =
```

```
1.41421 -0.70711  0.00000  
0.00000  1.22474 -0.81650  
0.00000  0.00000  1.15470
```

```
function [x, iflag, itnum] = Jacobi(A,b,x0,delta,max_it)
% x0: initial guess
% delta: error tolerance for the relative difference between two consecutive iterates
% max_it: maximum number of iterations to be allowed
% % iflag: 1 if a numerical solution satisfying the error
% itnum: the number of iterations used to compute x
n = length(b);
iflag = 1;
k = 0;
diagA = diag(A);
A = A-diag(diag(A));
while k < max_it
    k = k+1;
    x = (b-A*x0)./diagA
    relerr = norm(x-x0,inf)/(norm(x,inf)+eps);
    x0 = x;
    if relerr < delta
        break
    end
end
itnum = k;
if (itnum == max_it)
    iflag = -1
end
```

```
function x = GaussSeidel(A,b,es,maxit)
% A = coefficient matrix
% b = right hand side vector
% es = stop criterion (default = 0.00001%)
% maxit = max iterations (default = 50)
if nargin<2,error('at least 2 input arguments required'),end
if nargin<4|isempty(maxit),maxit=50;end
if nargin<3|isempty(es),es=0.00001;end
[m,n] = size(A);
if m~=n, error('Matrix A must be square'); end
C = A;
for i = 1:n
C(i,i) = 0;
x(i) = 0;
end
x = x';
```

$\Sigma \cup \nabla \epsilon \times \epsilon | \alpha \dots$

```
for i = 1:n
C(i,1:n) = C(i,1:n)/A(i,i);
end
for i = 1:n
d(i) = b(i)/A(i,i);
end
iter = 0;
while (1)
xold = x;
for i = 1:n
x(i) = d(i)-C(i,:)*x;
if x(i) ~= 0
ea(i) = abs((x(i) - xold(i))/x(i)) * 100;
end
end
iter = iter+1;
if max(ea)<=es | iter >= maxit, break, end
end
```

Νόρμες διανύσματος

Η p -νόρμα ($1 \leq p < \infty$) διανύσματος $x \in \mathbb{R}^n$ ορίζεται ως εξής:

$$\|x\|_p = \left(\sum_{i=1}^n |x_i|^p \right)^{1/p}$$

Στην πράξη χρησιμοποιούνται συνήθως οι 1-, η 2- και η ∞ -νόρμα για τις οποίες ισχύουν:

$$\|x\|_1 = \sum_{i=1}^n |x_i|, \quad \|x\|_2 = \left(\sum_{i=1}^n |x_i|^2 \right)^{1/2}, \quad \|x\|_\infty = \max_{1 \leq i \leq n} |x_i|$$

Παράδειγμα

```
>> x=[1 2 3 4]
```

```
x =
```

```
1 2 3 4
```

```
>> norm(x,1)
```

```
ans = 10
```

```
>> norm(x,2)
```

```
ans = 5.4772
```

```
>> norm(x,inf)
```

```
ans = 4
```

Νόρμες Πινάκων

$$\|A\|_1 = \max_{1 \leq j \leq n} \sum_{i=1}^n |a_{ij}|, \quad \|A\|_2 = \sqrt{\textcolor{brown}{r}_\sigma(A^T A)}, \quad \|A\|_\infty = \max_{1 \leq i \leq n} \sum_{j=1}^n |a_{ij}|$$

Παράδειγμα

```
>> A=[ 2 -1 4;-2 3 -1;6 -1 4]
```

```
A =
```

```
 2 -1 4  
-2 3 -1  
 6 -1 4
```

```
>> norm(A,1)
```

```
ans = 10
```

```
>> norm(A,2)
```

```
ans = 8.8141
```

```
>> norm(A,inf)
```

```
ans = 11
```

Δείκτης Κατάστασης

Ο δείκτης κατάστασης (condition number) ενός αντιστρέψιμου πίνακα A ορίζεται ως εξής:

$$\kappa_p(A) = \|A\|_p \|A^{-1}\|_p$$

Αποδεικνύεται ότι

$$\kappa_p(A) \geq 1$$

αν η χρησιμοποιούμενη νόρμα πίνακα είναι **φυσική**. Αν $\kappa_p(A) \gg 1$ λέμε ότι ο A είναι **κακής κατάστασης** (ill or badly conditioned). Διαφορετικά λέμε ότι ο A είναι **καλής κατάστασης** (well conditioned).

Παράδειγμα

Η λύση του γραμμικού συστήματος

$$\begin{bmatrix} .780 & .563 \\ .913 & .659 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \end{bmatrix} = \begin{bmatrix} .217 \\ .254 \end{bmatrix}$$

είναι $x_1 = 1$ και $x_2 = -1$. Πράγματι με τη MATLAB παίρνουμε:

```
>> A=[.780 .563; .913 .659]
```

```
A =
    0.7800    0.5630
    0.9130    0.6590
```

```
>> b=[.217; .254]
```

```
b =
    0.2170
    0.2540
```

```
>> x=A\b
```

```
x =
    1.0000
   -1.0000
```

Παράδειγμα

Αν διαταράξουμε πολύ λίγο τον πίνακα των συντελεστών, και λύσουμε το σύστημα

$$\begin{bmatrix} .780 & .563001 \\ .913 & .659 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \end{bmatrix} = \begin{bmatrix} .217 \\ .254 \end{bmatrix}$$

```
>> format long  
  
>> A(1,2) = .563001  
A =  
    0.780000000000000    0.563001000000000  
    0.913000000000000    0.659000000000000  
  
>> x=A\b  
  
x =  
    8.57471264411556  
   -11.49425287416920  
  
>> cond(A,1)  
  
ans =  
 2.6614e+006
```