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LA Lab ESE

Program No.1: Find the Eigen values of the following matrix.

$$A = \begin{bmatrix} 1 & 2 & 2 \\ 2 & 1 & 2 \\ 2 & 2 & 1 \end{bmatrix}$$

Code:-

```
clc; A = [1\ 2\ 2;\ 2\ 1\ 2;\ 2\ 2\ 1]; printf("The Matrix A is: "); disp(A); a = A(1,\ 1) + A(2,\ 2) + A(3,\ 3); b = A(1,\ 1)*A(2,\ 2) - A(1,\ 2)*A(2,\ 1) + A(2,\ 2)*A(3,\ 3) - A(2,\ 3)*A(3,\ 2) + A(1,\ 1)*A(3,\ 3) - A(1,\ 3)*A(3,\ 1); m = det(A); p = [1\ -a\ b\ -m]; m = roots(p); printf("The Eigen Values of Matrix A are:"); disp(m);
```

Output:

```
Scilab 6.1.1 Console

The Matrix A is:

1. 2. 2.

2. 1. 2.

2. 2. 1.

The Eigen Values of Matrix A are:

5. + 0.i

-1. + 0.i

-1. + 0.i

-1. + 0.i
```

Program No. 2:- Solve the given system of equations in terms of x,y,z,w using Gauss Elimination method

```
2x+3y+4z-2w=11
2x+3y-3z-w=1
7x+9y-4z+7z=21
102x-y+70z+301w=72
Code:-
clc;
A= [2 3 4 -2;2 3 -3 -1;7 9 -4 7;102 -1 70 301];
B = [11;1;21;72];
printf("The matrix A is:");
disp(A);
printf("The matrix B is:");
disp(B);
C=[A B];
printf("The matrix C is:");
disp(C);
n=4;
for i=1:n
  if C(i,i) == 0
     C(i,:)=C(i,:)+C(i+1,:);
  end
  if C(i,i) \sim = 0
     C(i,:)=C(i,:)/C(i,i);
  end
  disp(C);
  for j=i+1:n
     C(j,:)=C(j,:)-C(i,:)*C(j,i);
  end
  disp(C);
end
for i=n:-1:2
  for j=1:i-1
     C(j,:)=C(j,:)-C(i,:)*C(j,i);
  end
  disp(C);
printf("x=\%g\n",C(1,5));
printf("y=\%g\n",C(2,5));
printf("z=\% g\n",C(3,5));
printf("w=\%g\n",C(4,5));
```

Output :-

```
The matrix A is:
 2. 3. 4. -2.
 2. 3. -3. -1.
      9. -4. 7.
 7.
 102. -1. 70. 301.
The matrix B is:
 11.
  1.
 21.
 72.
The matrix C is:
 2. 3. 4. -2. 11.
 2. 3. -3. -1.
7. 9. -4. 7.
                   1.
                    21.
 102. -1. 70. 301.
                   72.
      1.5 2. -1.
 1.
                    5.5
                   1.
      3. -3. -1.
 2.
              7.
                    21.
      9. -4.
 102. -1. 70. 301. 72.
 1. 1.5 2. -1.
                    5.5
 0. 0. -7.
              1.
                   -10.
  0. -1.5 -18.
               14. -17.5
  0. -154. -134. 403. -489.
 1. 1.5 2. -1. 5.5
         16.666667 -10. 18.333333
 0. 1.
               14. -17.5
403. -489.
 0. -1.5 -18.
  0. -154. -134.
  1. 1.5 2.
                -1.
                       5.5
  0. 1. 16.666667 -10.
                       18.333333
  0. 0. 7. -1.
                        10.
  0. 0. 2432.6667 -1137. 2334.3333
```

1.5 2. -1. 5.5 0. 1. 16.666667 -10. 18.333333 0. 0. 1. -0.1428571 1.4285714 0. 2432.6667 -1137. 2334.3333 0. 1. 1.5 2. 5.5 -1. 0. 1. 16.666667 -10. 18.333333 1. -0.1428571 1.4285714 0. 0. 0. -789.47619 -1140.9048 0. 0. 1. 1.5 2. 5.5 -1. 0. 1. 16.666667 -10. 18.333333 0. 0. 1. -0.1428571 1.4285714 1. 0. 0. 0. 1.4451414 5.5 1. 1.5 2. -1. 0. 1. 16.666667 -10. 18.333333 0. 0. 1. -0.1428571 1.4285714 1. 0. 0. 0. 1.4451414 1.5 2. 0. 6.9451414 1. 0. 1. 16.666667 0. 32.784748 0. 0. 1. 0. 1.6350202 0. 0. 0. 1. 1.4451414 1. 1.5 0. 0. 3.675101 1. 0. 0. 5.534411 ο. 0. 0. 1. 0. 1.6350202 0. 0. 0. 1. 1.4451414 0. 0. 0. -4.6265155 1. 0. 1. 0. 0. 5.534411 0. 0. 1. 0. 1.6350202 0. 0. 0. 1. 1.4451414 x=-4.62652

y=5.53441 z=1.63502 w=1.44514