

Solution Set 5

Problem 1.

We present here an “all-in-one” program that writes out (1) a histogram of the largest eigenvalues, a histogram of the smallest eigenvalues and the cumulative density of states (DOS). This last quantity is the number of eigenvalues smaller than a chosen value.

In order to find the largest eigenvalues, we use the simple iterative algorithm given in the lectures: $x'_k = Ax_{k-1}$ og $x_k = x'_k/|x'_k|$. The largest eigenvalue is then given by $\lambda_m = x_{k+1} \cdot x'_k$.

However, there is a problem. The matrix A has its entries symmetrically distributed about zero. Sometimes the smallest eigenvalue — the most negative — is the largest one in absolute value. The algorithm picks out the eigenvalue with the largest absolute value, try e.g., the matrix $\text{diag}[-1, 1/2]$. (But, with a caveat: See below.) We solve this problem by simply adding a positive constant to the diagonal of the matrix A , and subtracting it from the eigenvalue when done.

By considering *symmetric* matrices, we avoid complex eigenvalues and the problems of having to work with complex eigenvectors.

In order to find the smallest eigenvalue, we define a new matrix

$$A' = \lambda_m - A, \quad (1)$$

and repeat. The smallest eigenvalue in A now becomes the largest eigenvalue in A' .

Lastly, we use the Lambert-Weaire algorithm to map out the distribution of eigenvalues between the smallest and largest eigenvalues of each A . The way I have done it is to choose 100 equally-spaced values for λ and run the Lambert-Weaire algorithm for each. When averaging over many samples, I have lumped together the results for each sample in the corresponding equally-spaced bin. That is, I have lumped the number of eigenvalues less than the 50th λ for sample number one together with the number of eigenvalues less than the 50th λ for sample number two even though they have different largest and smallest eigenvalues. In the end, I have used the *average* smallest and largest eigenvalues to scale the λ -values so that λ number one corresponds to the average smallest eigenvalue and λ number 100 to the average largest eigenvalue.

```

program ranmat
c-----
c Finding the smallest and largest eigenvalues in random
c matrices based on iteration.
c Then it uses the Lambert-Weaire algorithm to map the
c distribution of eigenvalues in between.
```

```

c-----
c n = size of matrix
c nsamp = number of samples
c itemx = max. number of iterations
c nh = number of bins in histograms of largest
c and smallest eigenvalues
c egenmx = value of largest bin in histograms
c egenmn = value of smallest bin in histograms
c nlw = number of bins in Lambert-Weaire calc.
c
      parameter(n=100,nsamp=2000,itemx=2000)
      parameter(nh=100,egenmx=20.,egenmn=-20.)
      parameter(nlw=100)
c-----
      dimension rmat(n,n),smat(n,n),qmat(n,n),vec(n),vecp(n)
6 dimension nhmx(nh),nhmn(nh),nhlw(nlw)
c-----
c Opening files to store output data
c
      open(unit=1,file='ranmat-histogram.dat',status='unknown')
      open(unit=2,file='ranmat-dos.dat',status='unknown')
c-----
c Initializing random number generator
c
      rinv=1./2147483647.
      ibm=4711
      do i=1,1000
      ibm=ibm*16807
      enddo
c-----
c Initializing histograms
c
c nhmx = histogram for largest eigenvalues.
c nhmn = histogram for smallest eigenvalues.
c
      do ih=1,nh
      nhmn(ih)=0
      nhmx(ih)=0
      enddo
c
c egen = bin size
c
      egen=(egenmx-egenmn)/(nh-1)
c

```

```

        avelamsm=0.
        avelambg=0.
c-----
c Initializing Lambert-Weaire result vector
c
        do ilw=1,nlw
            nhlw(ilw)=0
        enddo
c-----
c Loop over samples
c
        do isamp=1,nsamp
c-----
c Generating the matrix
c
c rmat = random matrix
c
        do i=1,n
            do j=1,i
                ibm=ibm*16807
                rmat(i,j)=ibm*rinv
                rmat(j,i)=rmat(i,j)
            enddo
        enddo
c-----
c A necessary trick to stabilize the iteration:
c
        do i=1,n
            rmat(i,i)=rmat(i,i)+1.
        enddo
c-----
c Determining the largest eigenvalue
c
c Initializing the eigenvectors vec.
c
        do i=1,n
            vec( i)=1.
        enddo
c-----
c The iteration:
c
c biglam = biggest eigenvalue
c
        avelam=egenmn

```

```

c      do ite=1,itemx
c
c      do i=1,n
c      vecp(i)=0.
c      do j=1,n
c      vecp(i)=vecp(i)+rmat(i,j)*vec(j)
c      enddo
c      enddo
c
c      sum=0.
c      do i=1,n
c      sum=sum+vecp(i)*vecp(i)
c      enddo
c      sum=1./sqrt(sum)
c
c      do i=1,n
c      vec(i)=vecp(i)*sum
c      enddo
c
c      biglam=0.
c      do i=1,n
c      biglam=biglam+vec(i)*vecp(i)
c      enddo
c
c      biglam=biglam-1.
c
c      enddo
c-----
c      do i=1,n
c      rmat(i,i)=rmat(i,i)-1.
c      enddo
c-----
c Histogram over biggest eigenvalue
c
c      ih=int((biglam-egenmn)/egen)
c      ih=max(ih,1)
c      ih=min(ih,nh)
c      nhmx(ih)=nhmx(ih)+1
c
c      avelambg=avelambg+biglam
c-----
c smalam = smallest eigenvalue
c

```

```
c Constructing new iteration matrix
```

```
c
```

```
do i=1,n
do j=1,n
smat(i,j)=-rmat(i,j)
enddo
enddo
```

```
c
```

```
do i=1,n
smat(i,i)=smat(i,i)+biglam
enddo
```

```
c
```

```
do i=1,n
vec( i)=1.
enddo
```

```
c
```

```
do ite=1,itemx
```

```
c
```

```
do i=1,n
vecp(i)=0.
do j=1,n
vecp(i)=vecp(i)+smat(i,j)*vec(j)
enddo
enddo
```

```
c
```

```
sum=0.
do i=1,n
sum=sum+vecp(i)*vecp(i)
enddo
sum=1./sqrt(sum)
```

```
c
```

```
do i=1,n
vec(i)=vecp(i)*sum
enddo
smalam=0.
do i=1,n
smalam=smalam+vec(i)*vecp(i)
enddo
```

```
c
```

```
smalam=biglam-smalam
```

```
c
```

```
enddo
```

```
c-----
```

```
c Histogram over smallest eigenvalue
```

```

c
    ih=int((smalam-egenmn)/egen)
    ih=max(ih,1)
    ih=min(ih,nh)
    nhmn(ih)=nhmn(ih)+1
c
    avelamsm=avelamsm+smalam
c-----
c The Lambert-Weaire algorithm
c
    egan=(biglam-smalam)/nlw
c
c Loop over lambda-values
c
    do ilw=1,nlw
c
        alam=smalam+egan*ilw
c
        do i=1,n
            do j=1,n
                smat(i,j)=rmat(i,j)
            enddo
        enddo
c
        nshift=1
c
        do k=n,2,-1
c
            ann=1./(smat(k,k)-alam)
            if(ann.lt.0.) nshift=nshift+1
c
            do i=1,k-1
                do j=1,k-1
                    qmat(i,j)=smat(i,j)-smat(i,k)*smat(k,j)*ann
                enddo
            enddo
c
            do i=1,k-1
                do j=1,k-1
                    smat(i,j)=qmat(i,j)
                enddo
            enddo
c
        enddo

```

```

c-----
c End of Lambert-Weaire iteration
c
    nhlw(ilw)=nhlw(ilw)+nshift
c
    enddo
c-----
c End of loop over samples
c
    enddo
c-----
c Writing the histograms
c
    egenv=egenmn
    do ih=1,nh
        egenv=egenv+egen
        write(1,*) egenv,nhmn(ih),nhmx(ih)
    enddo
c-----
c Writing the DOS
c
    avelambg=avelambg/nsamp
    avelamsm=avelamsm/nsamp
c
    egan=(avelambg-avelamsm)/(nlw-1)
    do ilw=1,nlw
        alam=avelamsm+egan*(ilw-1)
        write(2,*) alam,float(nhlw(ilw))/nsamp
    enddo
c-----
    close(1)
    close(2)
c-----
    end

```

We show the histograms of the smallest and largest eigenvalues in the figure below:
The distribution of eigenvalues found with the Lambert-Weaire algorithm is shown in the figure on the next page.

The data presented here was averaged over 16000 samples. I used 2000 iterations in the iterative algorithms. I checked that this was enough.



