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# Project-I

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## 1 PageRank

### 1.1 Exercise1

i) After randomly initializing  $x_0$  and repeating the experiments. E has the eigenvalues

$$\begin{pmatrix} 1 \\ -0.5000 + 0.2887i \\ -0.5000 - 0.2887i \\ 0 \end{pmatrix}$$

two real and two complex values, the largest eigenvalue is 1. The sequence  $E^{15}x_0$  normalised using l2 norm, as the number of iterations increases tends to the largest eigenvector of E, the eigenvector corresponding to eigenvalue 1. eigenvector

$$\begin{pmatrix} 0.6447 \\ 0.2478 \\ 0.4154 \\ 0.5920 \end{pmatrix}$$

ii) Eigenvector

$$\begin{pmatrix} 0.2230 \\ 0.3755 \\ 0.4201 \\ 0.7955 \end{pmatrix}$$

Eigenvalues

$$\begin{pmatrix} 0 \\ 0.5614 \\ -0.2807 + 0.2640i \\ -0.2807 - 0.2640i \end{pmatrix}$$

The largest eigenvalue is 0.5614 and the sequence converges again to the largest eigenvector, up to a constant, -1 in my case.

iii) Eigenvalues of E

$$\begin{pmatrix} 1.0000 \\ -0.5000 + 0.2887i \\ -0.5000 - 0.2887i \\ -0.0000 \\ 1.0000 \\ -1.0000 \end{pmatrix}$$

There are two eigenvalues of 1. The sequence after 15 steps does not converge to the largest eigenvector (any of them)

$$\begin{pmatrix} 0.3500 \\ 0.1050 \\ 0.2099 \\ 0.3149 \\ 0.6848 \\ 0.5043 \end{pmatrix}$$

## 1.2 Exercise2

i) The eigenvectors are  $-0.6446 - 0.2887 + 0.5000i$   $-0.2887 - 0.5000i$   $0.5345 - 0.2478 - 0.4330 - 0.2500i$   $-0.4330 + 0.2500i$   $-0.8018 - 0.4154$   $0.5774$   $0.5774$   $0.0000$   $-0.5920$   $0.1443 - 0.2500i$   $0.1443 + 0.2500i$   
Eigenvalues are

$$\begin{pmatrix} 1.0000 \\ -0.4250 + 0.2454i \\ -0.4250 - 0.2454i \\ -0.0000 \end{pmatrix}$$

The convergence vector is

$$\begin{pmatrix} 0.6447 \\ 0.2478 \\ 0.4154 \\ 0.5920 \end{pmatrix}$$

ii) GREEN  $0.2769 - 0.3641 - 0.3761i$   $-0.3641 + 0.3761i$   $0.0044$   $0.3984$   $0.5959$   $0.5959$   $-0.1091$   $0.4548$   $-0.3506 + 0.2950i$   $-0.3506 - 0.2950i$   $-0.0657$   $0.7468$   $0.2363 + 0.3242i$   $0.2363 - 0.3242i$   $0.9918$   
Eigenvalues are

$$\begin{pmatrix} 0.6618 \\ -0.2427 + 0.2257i \\ -0.2427 - 0.2257i \\ -0.0264 \end{pmatrix}$$

The convergence vector is

$$\begin{pmatrix} 0.2769 \\ 0.3984 \\ 0.4548 \\ 0.7468 \end{pmatrix}$$

iii) BLUE Eigenvector of M blue

$$\begin{pmatrix} 0.5351 \\ 0.2057 \\ 0.3449 \\ 0.4914 \\ 0.3943 \\ 0.3943 \end{pmatrix}$$

The eigenvalues are

$$\begin{pmatrix} 1.0000 \\ 0.8500 \\ -0.4250 + 0.2454i \\ -0.4250 - 0.2454i \\ 0.0000 \\ -0.8500 \end{pmatrix}$$

The convergence vector is

$$\begin{pmatrix} 0.5200 \\ 0.2012 \\ 0.3358 \\ 0.4778 \\ 0.4257 \\ 0.4085 \end{pmatrix}$$

### 1.3 Exercise3

i) E is a stochastic column matrix for i) and iii). ii) has a column only with 0 elements. M is a stochastic column matrix for i) and iii) again. ii) has a column only of 0.15 values, so it is not .

ii) Let  $A$  be a stochastic column matrix.  $A$  and  $A^\top$  have the same eigenvalues, so it is enough to show that  $A^\top$  has eigenvalue 1, which will imply  $A$  also has eigenvalue of 1.

We use the property that for eigenvalue  $\lambda$  of  $A^\top$   $\det(A^\top - \lambda * I) = 0$ . We compute

$$\det(A^\top - I) = \det \begin{pmatrix} a_{11} - 1 & a_{21} & \dots & a_{n1} \\ a_{12} & a_{22} - 1 & \dots & a_{n2} \\ \dots & \dots & \dots & \dots \\ a_{1n} & a_{2n} & \dots & a_{nn} - 1 \end{pmatrix} \quad (1)$$

$$\det(A^\top - I) = \det \begin{pmatrix} \sum_i a_{i1} - 1 & a_{21} & \dots & a_{n1} \\ \sum_i a_{i2} - 1 & a_{22} - 1 & \dots & a_{n2} \\ \dots & \dots & \dots & \dots \\ \sum_i a_{in} - 1 & a_{2n} & \dots & a_{nn} - 1 \end{pmatrix} \quad (2)$$

$$\det(A^\top - I) = \det \begin{pmatrix} 0 & a_{21} & \dots & a_{n1} \\ 0 & a_{22} - 1 & \dots & a_{n2} \\ \dots & \dots & \dots & \dots \\ 0 & a_{2n} & \dots & a_{nn} - 1 \end{pmatrix} \quad (3)$$

For the first equality we added all the columns 2, 3, ...,  $N - 1$  to the first column. For the second equality, the determinant of a matrix with a column of 0 equals 0. This implies  $A^\top$  and implicitly  $A$  have eigenvalue 1.

Assume that  $A, A^\top$  have an eigenvalue greater than 1,  $\lambda_0$  with eigenvector  $v$ .  $A^\top v = \lambda_0 v$ . Then we have  $\sum_i a_{it} * v_i = \lambda_0 v_t$  for every  $t$  in  $0..N$ . Let  $m$  be the index corresponding to the largest entry in  $v$ . then  $\sum a_{ik} * v_i < \sum a_{ij} v_k = v_k < \lambda v_k$  which is a contradiction, so 1 is the largest eigenvalue.

### 1.4 Exercise4

Gradient of  $f \frac{\partial f}{\partial x} =$

**Acknowledgments**

**References**