

# Network Science Cheatsheet



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## 2 Networks as matrices

### Matrices in short

Matrices are mathematical objects that can be thought as *tables* of numbers. The size of a matrix is expressed as  $m \times n$ , for a matrix with  $m$  rows and  $n$  columns. **The order (row/column) is important.**  $M_{ij}$  is a notation representing the element on **row**  $m$  and **column**  $j$ .

### A - Adjacency matrix

The most natural way to represent a graph as a matrix is called the Adjacency matrix  $A$ . It is defined as a square matrix, such as the number of rows (and the number of columns) is equal to the number of nodes  $N$  in the graph. Nodes of the graph are numbered from 1 to  $N$ , and there is an edge between nodes  $i$  and  $j$  if the corresponding position of the matrix  $A_{ij}$  is not 0.

- A value on the diagonal means that the corresponding node has a **self-loop**
- the graph is **undirected**, the matrix is **symmetric**:  $A_{ij} = A_{ji}$  for any  $i, j$ .
- In an **unweighted** network, and edge is represented by the value 1.
- In a **weighted** network, the value  $A_{ij}$  represents the **weight** of the edge  $(i, j)$

### Typical operations on A

Some operations on Adjacency matrices have straightforward interpretations and are frequently used

**Multiplying A by itself** allows to know the number of walks of a given length that exist between any pair of nodes:  $A_{ij}^2$  corresponds to the number of walks of length 2 from node  $i$  to node  $j$ .  $A_{ij}^3$  to the number of walks of length 3, etc.

**Multiplying A by a column vector W** of length  $1 \times N$  can be thought as setting the  $i$ th value of the vector to the  $i$ th node, and each node *sending* its value to its neighbors (for undirected graphs). The result is a column vector with  $N$  elements, the  $i$ th element corresponding to the sum of the values of its neighbors in  $W$ . This is convenient when working with **random walks** or **diffusion** phenomenon.

### Spectral properties of A

**Spectral Graph Theory** is a whole field in itself, and beyond the scope of this class. A few elements for those with a *linear algebra* background:

- The adjacency matrix of an undirected simple graph is symmetric, and therefore has a complete set of real eigenvalues and an orthogonal eigenvector basis.
- The set of eigenvalues of a graph is the spectrum of the graph.
- Eigenvalues are denoted as  $\lambda_0 \leq \lambda_1 \leq \lambda_2 \leq \dots \lambda_n$
- The largest eigenvalue  $\lambda_0$  lies between the average and maximum degrees
- The number of closed walks of length  $k$  in  $G$  equals  $\sum_i^n \lambda_i^k$
- A graph is bipartite if and only if its spectrum is symmetric (i.e., if  $\lambda$  is an eigenvalue, then so is  $-\lambda$ )
- If  $G$  is connected, then the diameter of  $G$  is strictly less than its number of distinct eigenvalues

### Graph Laplacian

The **Graph Laplacian**, or **Laplacian Matrix** of a graph is a variant of the Adjacency matrix, often used in *Graph theory* and *Spectral Graph Theory*. It is defined as  $D - A$ , with  $D$  the *Degree matrix* of the graph, defined as a  $N \times N$  matrix with  $D_{ii} = k_i$  and zeros everywhere else.

Intuitively, Laplace operator is a generalization of the second derivative, and is defined in discrete situations, for each value, as the sum of differences between the value and its "neighbors". e.g., in time, the  $2^{nd}$  derivative *acceleration* is the difference between current speed and previous speed. In a B&W picture, it's the difference between the greylevel on current pixel and the greylevel of 4 or 8 closest pixels, and perform *edge detection*. On a graph, with  $W$  a column vector representing values on nodes,  $LW$  computes for each node the difference to neighbors.

### Spectral properties of L

Eigenvalues of the Laplacian have many applications, such as *spectral clustering*, *graph matching*, *embedding*, etc. Assuming  $G$  undirected with eigenvalues  $\lambda_0 \leq \lambda_1 \leq \lambda_2 \leq \dots \lambda_n$ , here are some interesting properties:

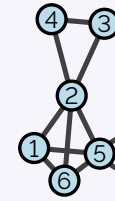
- The smallest eigenvalue  $\lambda_0$  equals 0
- The number of 0 eigenvalues gives the number of connected components

### Random Walk matrix

Another useful matrix of a graph is the **Random Walk Transition Matrix R**. It is the column normalized version of the adjacency matrix.  $R_{ij}$  can be understood as the probability for a random walker located on node  $i$  to move to  $j$ .

### Matrix notation - Example

#### Graph



#### A - Adjacency Mat.

$$\begin{pmatrix} 0 & 1 & 0 & 0 & 1 & 1 \\ 1 & 0 & 1 & 1 & 1 & 1 \\ 0 & 1 & 0 & 1 & 0 & 0 \\ 0 & 1 & 1 & 0 & 0 & 0 \\ 1 & 1 & 0 & 0 & 1 & 1 \\ 1 & 1 & 0 & 0 & 1 & 0 \end{pmatrix}$$

#### D - Degree Matrix

$$\begin{pmatrix} 3 & 0 & 0 & 0 & 0 & 0 \\ 0 & 5 & 0 & 0 & 0 & 0 \\ 0 & 0 & 2 & 0 & 0 & 0 \\ 0 & 0 & 0 & 2 & 0 & 0 \\ 0 & 0 & 0 & 0 & 5 & 0 \\ 0 & 0 & 0 & 0 & 0 & 3 \end{pmatrix}$$

#### L - Laplacian

$$\begin{pmatrix} 3 & -1 & 0 & 0 & -1 & -1 \\ -1 & 5 & -1 & -1 & -1 & -1 \\ 0 & -1 & 2 & -1 & 0 & 0 \\ 0 & -1 & -1 & 2 & 0 & 0 \\ -1 & -1 & 0 & 0 & 4 & -1 \\ -1 & -1 & 0 & 0 & -1 & 3 \end{pmatrix}$$

#### A<sup>2</sup>

$$\begin{pmatrix} 3 & 2 & 1 & 1 & 3 & 2 \\ 2 & 5 & 1 & 1 & 3 & 2 \\ 1 & 1 & 2 & 1 & 1 & 1 \\ 1 & 1 & 1 & 2 & 1 & 1 \\ 3 & 3 & 1 & 1 & 4 & 3 \\ 2 & 2 & 1 & 1 & 3 & 3 \end{pmatrix}$$

#### Random W. mat.

$$\begin{pmatrix} 0 & \frac{1}{5} & 0 & 0 & \frac{1}{4} & \frac{1}{3} \\ \frac{1}{3} & 0 & \frac{1}{2} & \frac{1}{2} & \frac{1}{4} & \frac{1}{3} \\ 0 & \frac{1}{5} & 0 & \frac{1}{2} & 0 & 0 \\ 0 & \frac{1}{5} & \frac{1}{2} & 0 & 0 & 0 \\ \frac{1}{3} & \frac{1}{5} & 0 & 0 & \frac{1}{4} & \frac{1}{3} \\ \frac{1}{3} & \frac{1}{5} & 0 & 0 & \frac{1}{4} & 0 \end{pmatrix}$$