

10) → KL algorithm.

- Uses difference between number of intracommunity edges & intercommunity edges as the optimisation function.
- A good partition contains more density within community.
- Simple.
- Complexity of  $O(n^2)$
- not suitable for large networks.
- requires prior information so inefficient

→ Faster Newman Algorithm

- Speeds up the community mining process.
- Has a complexity of  $O(mn)$
- Starts off with each node in a separate community.
- Iteratively combines nodes to form bigger communities until some pre-defined halt condition / until one large community has been found. Since only neighbors considered, risk of local optimum
- Constructs the dendrogram in a bottom-up manner;  $a_i = \sum_j c_{ij}$  where  $c_{ij}$  is the edge weight between node  $i$  &  $j$

→ GIA Algorithm

- Uses simulated annealing in order to control local search based optimisation.
- Stochastic Global Search optimisation is used in this method. It finds global optimum.
- Randomness is introduced into this model.

$$P = \begin{cases} 1 & \text{if } C_{t+1} < C_t \\ e^{-(C_{t+1} - C_t)/T} & C_{t+1} \geq C_t \end{cases}$$

where  $P$  is the probability of the node being in the graph.

$C_{t+1}, C_t$  are parameters at time  $t$  &  $t+1$  whereas  $T$  is the temperature.

Initial parameters: Starting configurations, temperature (controls SA)

No prior information is required before starting. The time depends on SA's exec time.

Randomness is implemented by selecting/deleting/merging & splitting node groups.

→ Continuing KL algorithm

In this method, different configurations are tested by randomly picking different nodes & by calculation of objective function.

The algo terminates when the evaluation metric does not improve.

→ Potts Algorithm

A network is considered a multi-step Potts' model, with each node having a spin of value  $q$ .

The best partition is the most stable one with least energy.

In this stable form, the nodes with a

particular spin forms a community inherently.

∴ The aim is to split differently spinning models.

Further approximations to the objective function of the given algorithms by means of Monte Carlo optimisation techniques can lead to improvement in speed, while compromising on the accuracy a little.

(3) Let Alice be 0

Bob	1
Dave	2
Carol	3

Network X code:

```
import networkx as nx.
```

```
G = nx.Graph()
```

```
G.add_weighted_edges_from([(0, 1, 1), (0, 2, 1),  
                           (3, 0, 1)])
```

# (from, to, weighted edge) denoting edge of weight 1

```
pos = nx.ego_graph(G)
```

```
ego-centre = 0
```

```
depth = 3 # friend of a friend of a friend
```

```
nx.draw(G, pos, to with_labels=True, radius=  
        depth, ego-centre =  
        ego-center)
```

# plots the ego-centric graph for Alice.

def snowball(G<sub>1</sub>, centre, max-depth,  
curr-depth, ~~visited~~ = [])

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if max-depth == curr-depth:  
return ~~set~~ friends # reached depth.

if centre in friends:

return friends. # already traversed

for node in G<sub>1</sub>.neighbors(centre):

friends.append(node)

if node not in friends:

snowball(G<sub>1</sub>, node, max-depth)

curr-depth + 1, friends)

# recursive call.

return friends

val =

snowball (G<sub>1</sub>, ego-centre, 3, 0)

# function call.

print (val) # return result of data collection

using same network G<sub>1</sub>.

Data collected in the people in the  
network

## PART-B

7) There are 3 categories for the definition of communities.

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### Local Definition

→ Self referral: Focuses on the vertices of subnetwork & only its neighbors.

Clique: maximally connected subgraph where each node is connected to  $k-1$  other nodes.

$k$ -clique: It is a clique in which the max. distance between nodes is always lesser than  $k$ .

$k$ -plex: Maximally connected graph in which nodes are connected to all but almost ~~no~~  $k$ -of them.

→ Comparative: where edges between mutual connections within network is compared with external edges.

LS set: no. of neighbors inside  $>$  no. of neighbors outside community

Weakly connected graph: no. of connections inside a community  $>$  edge connecting outside.

### Global definition

The community is defined w.r.t the overall network.

Null model: Randomly distributed.

Similar structurally to original w/w but no

properties of its own  
It is created by introducing random changes into the overall network, by reassessing edges to vertices randomly.

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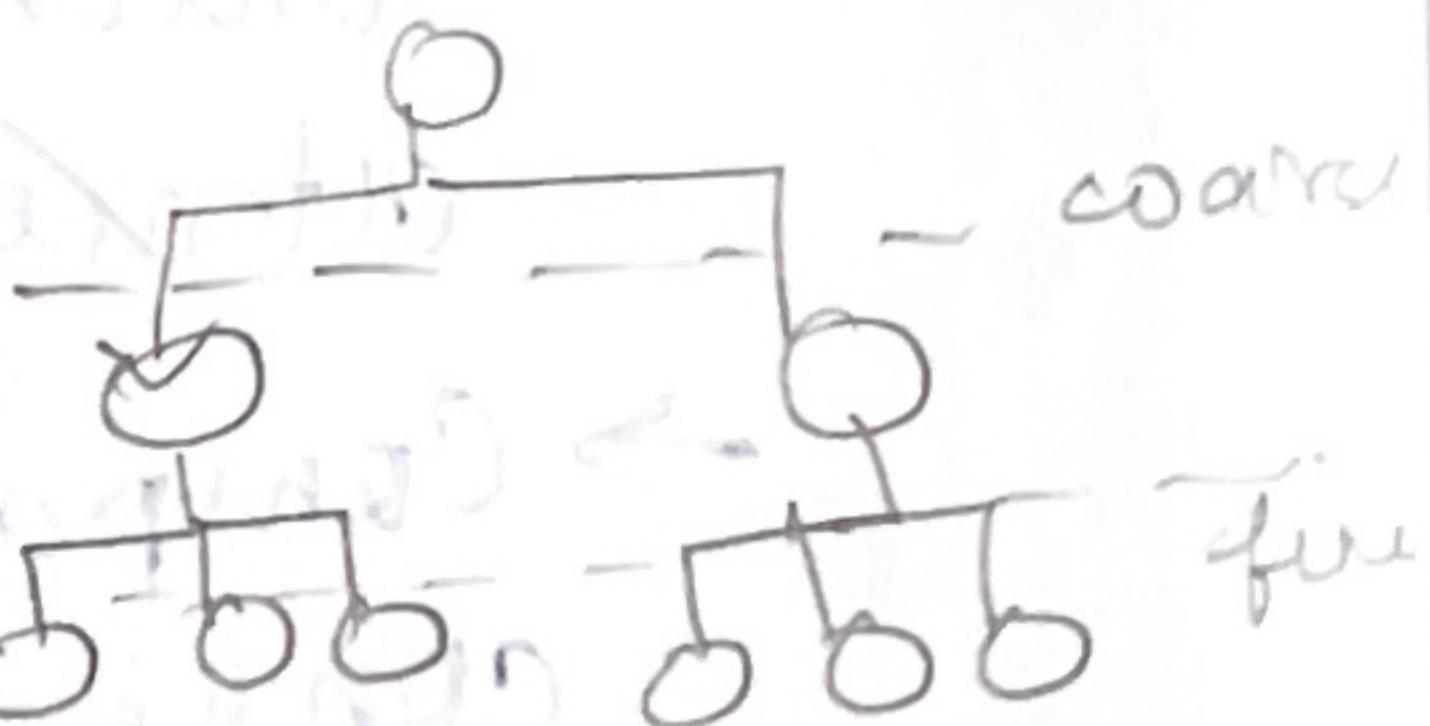
If there is a large difference when compared to original N/W, it is termed a community.

### Based on Vertex Similarity

Similar vertices lie on the same level of the hierarchical dendrogram structure. By taking a horizontal line, various communities can be obtained.

Granularity is fine when the line is low; high coarse when high.

The similarity of vertices is determined quantitatively.



# 8) modularity Quality function

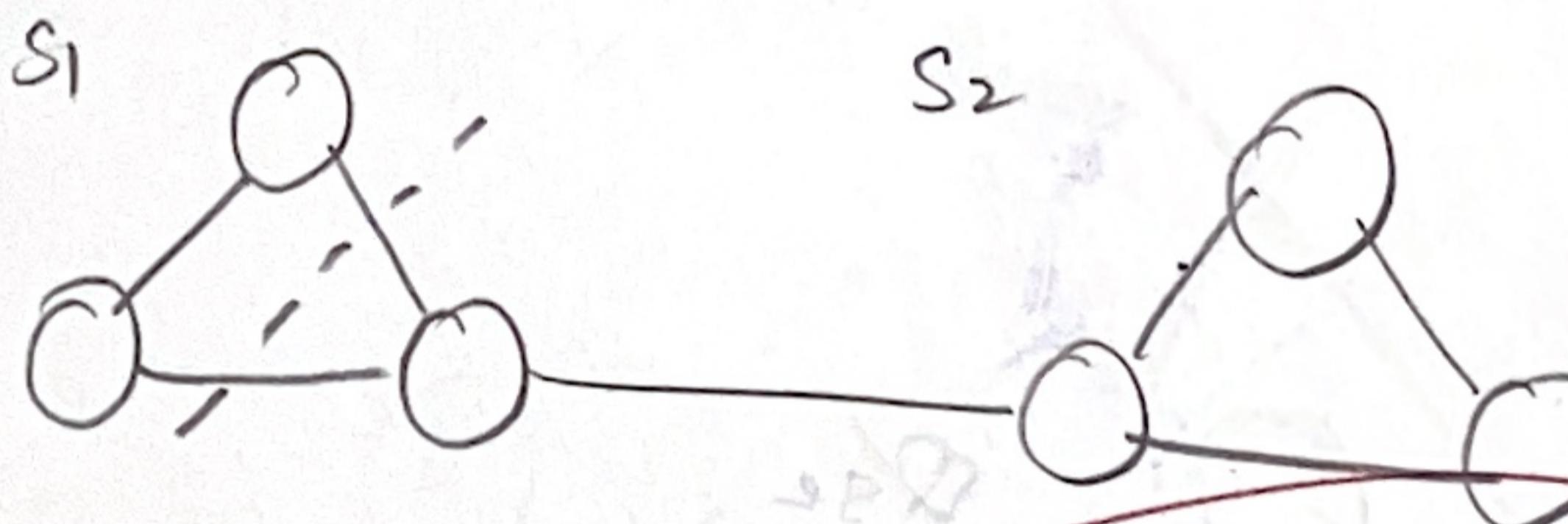
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$$Q = \sum_{i=1}^{nm} \left[ \frac{ds}{L} - \left( \frac{ds}{2L} \right)^2 \right]$$

where L is the total no. of edges in community  
 ds no. of edges within community  
 ds  $\Rightarrow$  degree of nodes' summation

$$Q = \sum \left( \text{Actual fraction of edges inside graph} - \text{Expected Value of edges for graph w/ same vertices} \right)$$

Split 1



$$L = 7$$

$Q_{S_1}$

$$\begin{aligned} ds &= 1 \\ ds &= 4 \end{aligned}$$

$Q_{S_2}$

$$\begin{aligned} ds &= 4 \\ ds &= 10 \end{aligned}$$

$$Q_{S_1} = \frac{1}{7} - \left( \frac{4}{14} \right)^2$$

$$= 0.14 - 0.0816$$

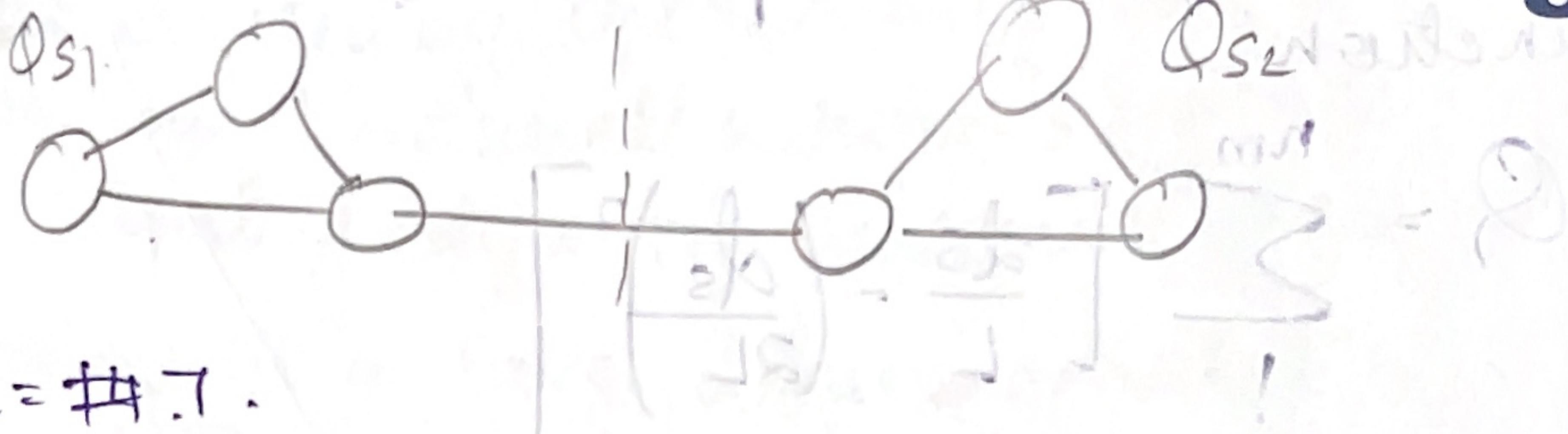
$$\approx 0.061224$$

$$Q_{S_2} = 0.06122$$

$$= 0.5714 - (0.71)^2$$

$$Q = Q_{S_1} + Q_{S_2} = \cancel{0.2244} \cdot 0.12244$$

## Cut 2.



$$L = \#7.$$

$$Q_{S_1}$$

$$ls = 3$$

$$ds = 7$$

$$Q_{S_2}$$

$$ls = 3$$

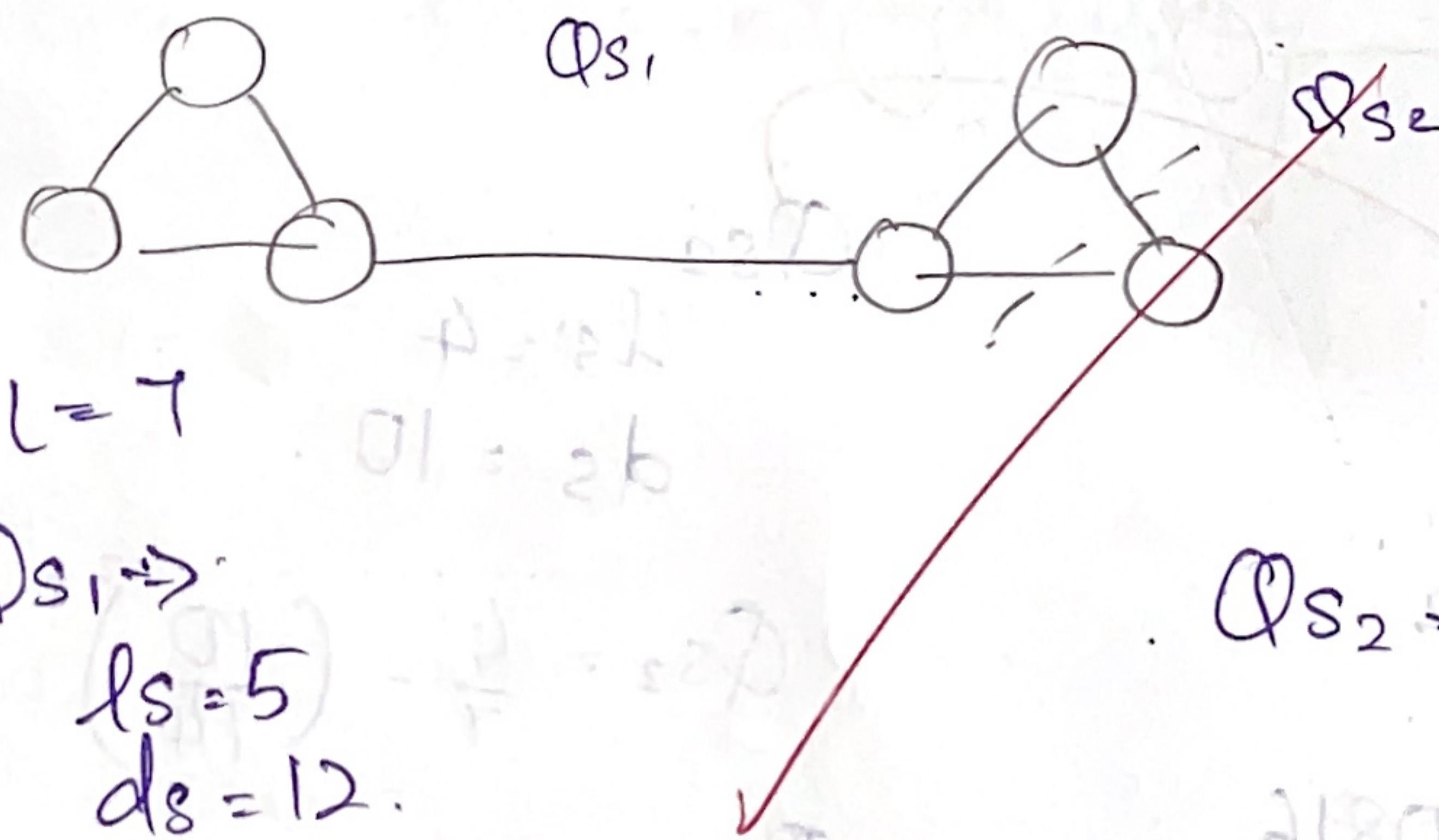
$$ds = 7.$$

$$\begin{aligned} Q_{S_1} &= \frac{3}{7} - \left(\frac{7}{14}\right)^2 \\ &= 0.178 \end{aligned}$$

$$\begin{aligned} Q_{S_2} &= \frac{3}{7} - \left(\frac{7}{14}\right)^2 \\ &= 0.178 \end{aligned}$$

$$\Phi = Q_{S_1} + Q_{S_2} = 0.357$$

## Cut 3



$$Q_{S_1} \rightarrow$$

$$ls = 5$$

$$ds = 12.$$

$$Q_{S_2} \rightarrow$$

$$ls = 0$$

$$ds = 2.$$

$$\begin{aligned} Q_{S_1} &= \frac{5}{7} - \left(\frac{12}{14}\right)^2 \\ &= -0.02 \end{aligned}$$

$$\begin{aligned} Q_{S_2} &= \frac{0}{7} - \left(\frac{2}{14}\right)^2 \\ &= -0.0204. \end{aligned}$$

Negative

## PART-A

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### 1. Non-directional

- friends on facebook
- connections on linkedin
- reddit friendships
- retweeting on twitter.

### 2. The relations for the scenario "children playing together" are.

- Ego-centric network for a particular child
- friendship network for all the children to visualise their mutual friendships

3. It is important because
- Content Filtering: similar people exist in similar networks & share taste.
  - Useful for recommendation systems
  - Visualisation of large scale graphs, for information flow & diffusion
  - Understanding properties and structure of the graphs.

4.

$$\sum_{i=0}^{nm} \left( \frac{\text{actual no. of edges within community}}{\text{Expected number of edges for an equivalent null model graph}} \right)^2$$

$nm \rightarrow \text{no. of communities}$

Originally

$$\sum_{i=0}^{nm} \frac{l_s}{L} \left( \frac{ds}{d_L} \right)^2$$

$\rightarrow \text{total no. of edges}$

$l_s \rightarrow \text{links within } N \text{ communities}$

$ds \rightarrow \text{degree summation}$ :

## Optimization

- 5)  $\rightarrow$  converts the NCMP into an optimization problem; optimise the objective function.
- $\rightarrow$  Obj. function: cut crit. erra, evaluation fns etc
- $\rightarrow$  Further types: local search based  $\leftarrow$  spect  
ral search
- $\rightarrow$  Eg: Faster Newman,  
KL algorithm

## Heuristic

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**SSN**
- $\rightarrow$  Does NCMP as a heuristic based problem taking some assumption/heuristic measure.
- $\rightarrow$  Eg: MFC considers that flow between com. should be greater.
- $\rightarrow$  Eg: WH Algorithm, Gitt Van-Neuman Algorithm

## 6) import networkx.

# initialise:

new\_G = networkx.Graph()  
~~for from, to, edata in G.edges():~~

for node in G.nodes():

'if G.degree(node) > 2: ~~cont~~

continue' # do not add.

new\_G.add\_node(node) # add in new graph

$$2.8 = 2.0 + 2.1 + 2.1 + 2.3 + 2.0 = 10A$$

$$2.4 = 2.1 + 2.0 + 2.8 + 2.1 + 2.1 = 10A$$

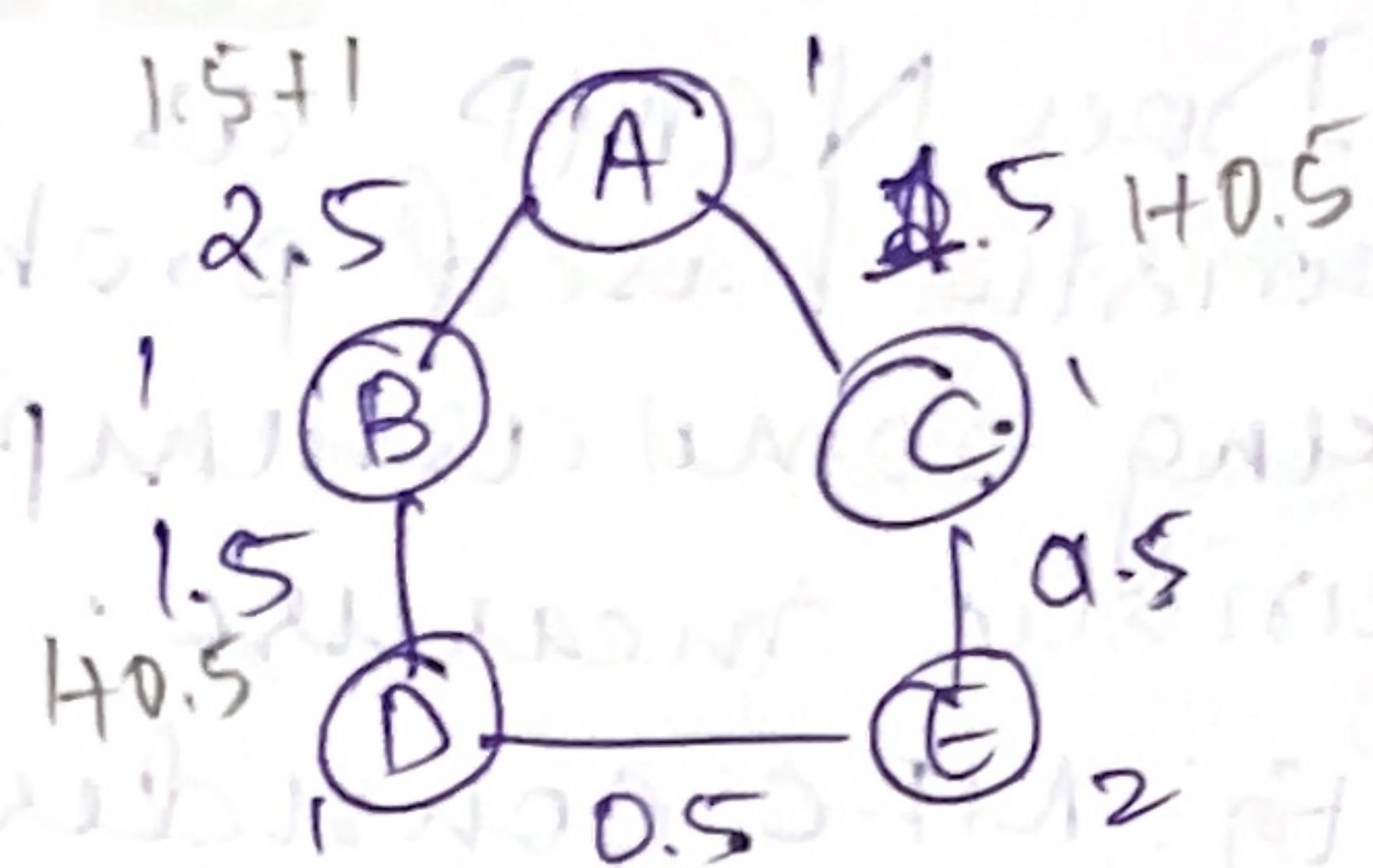
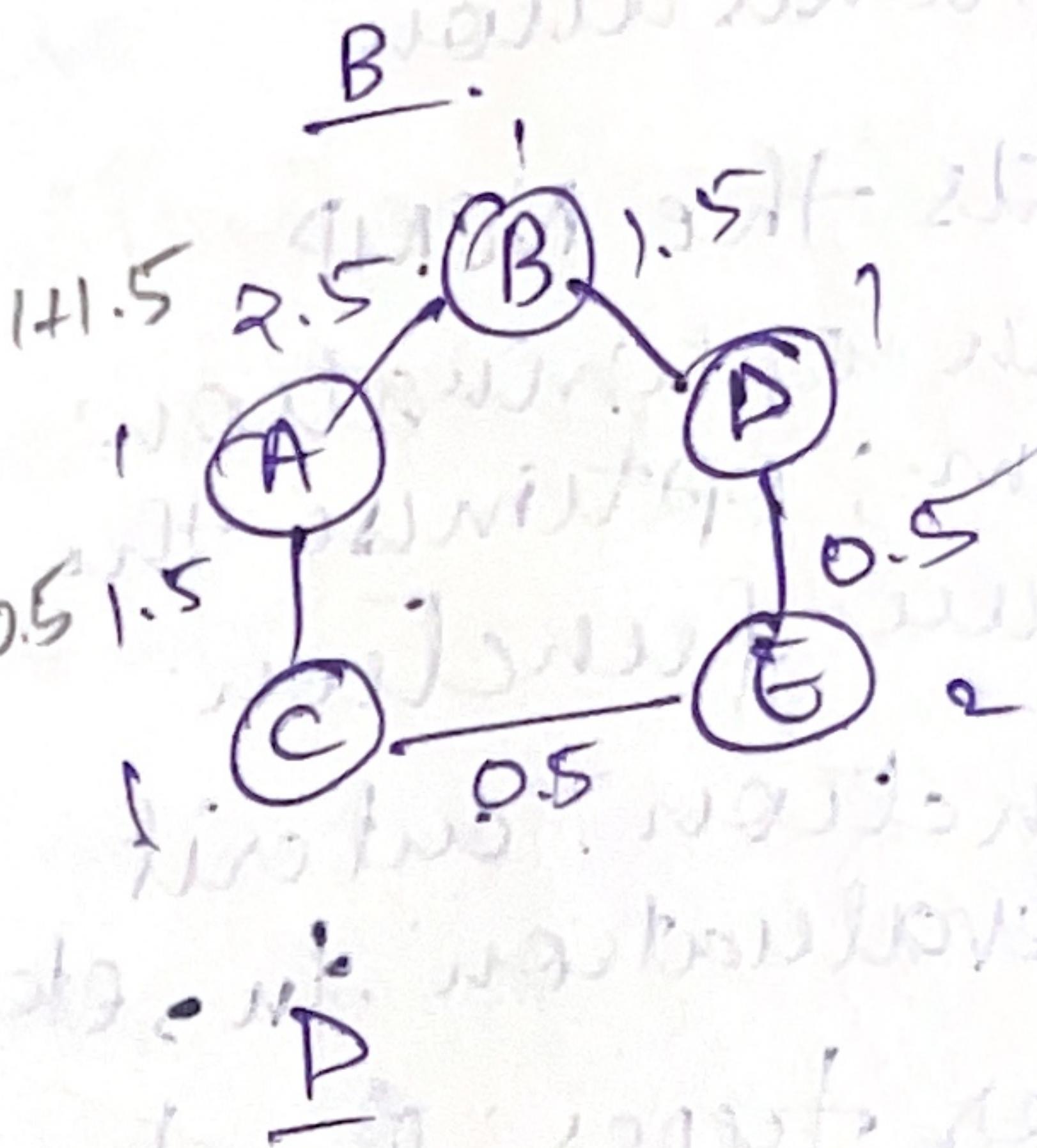
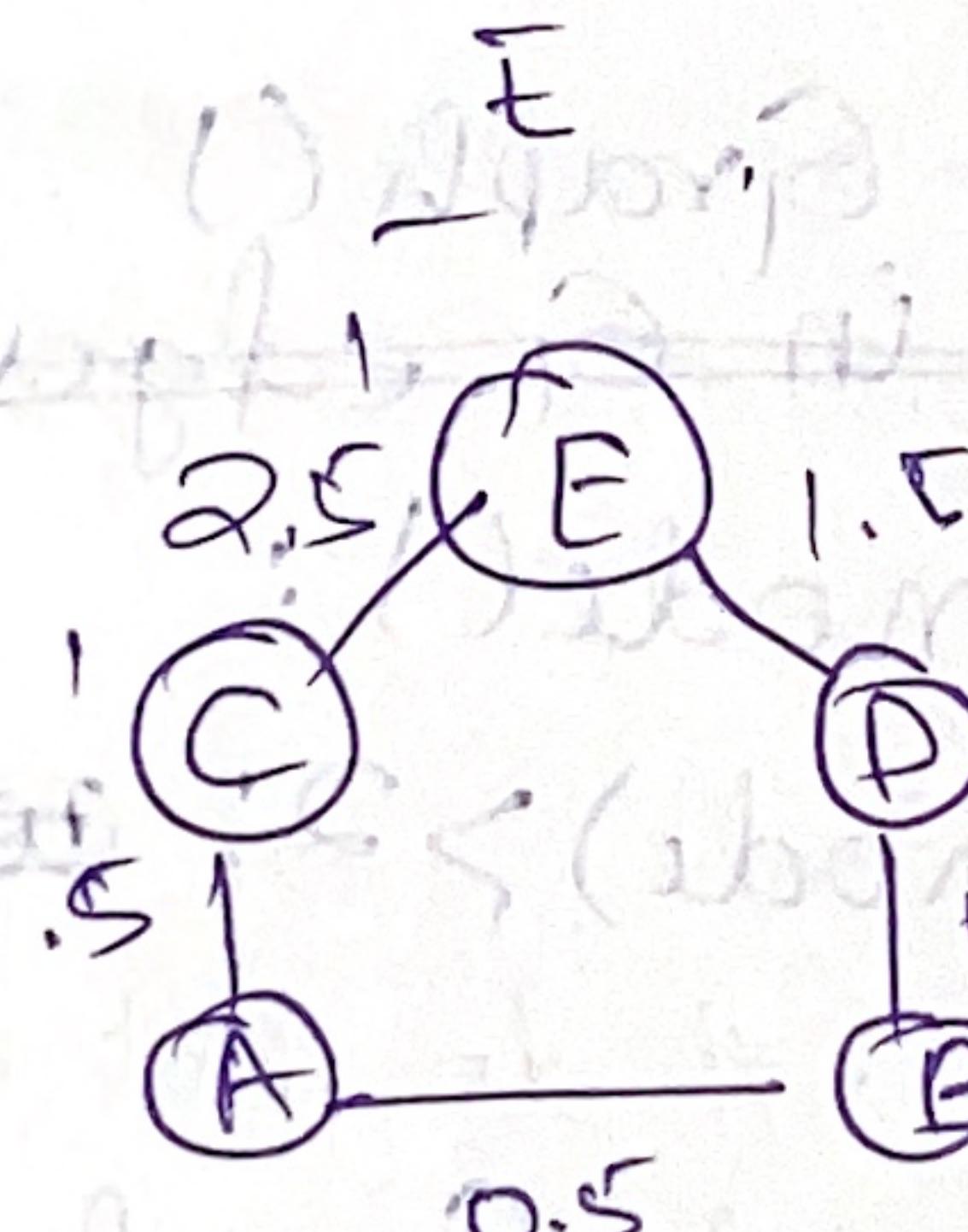
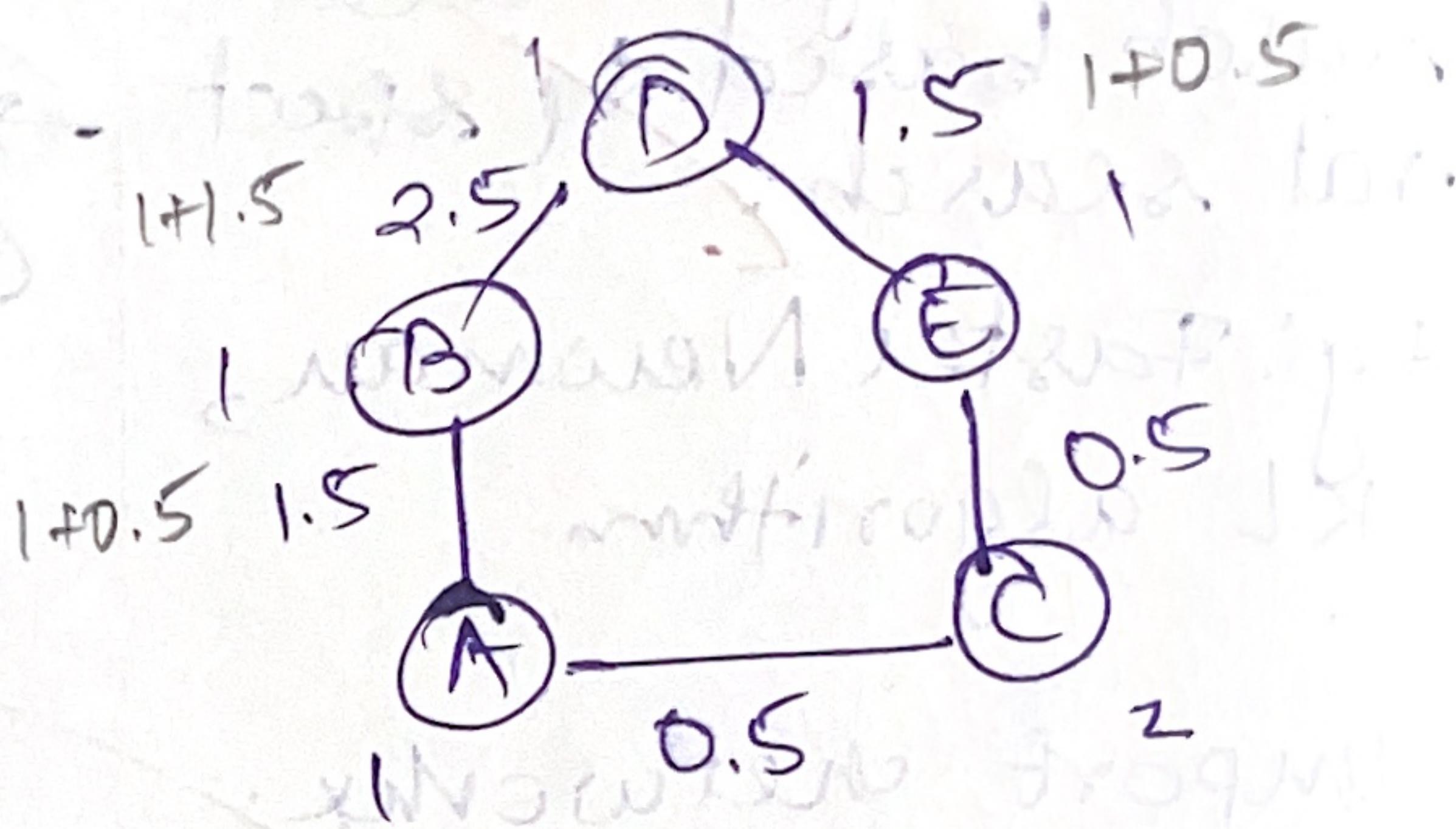
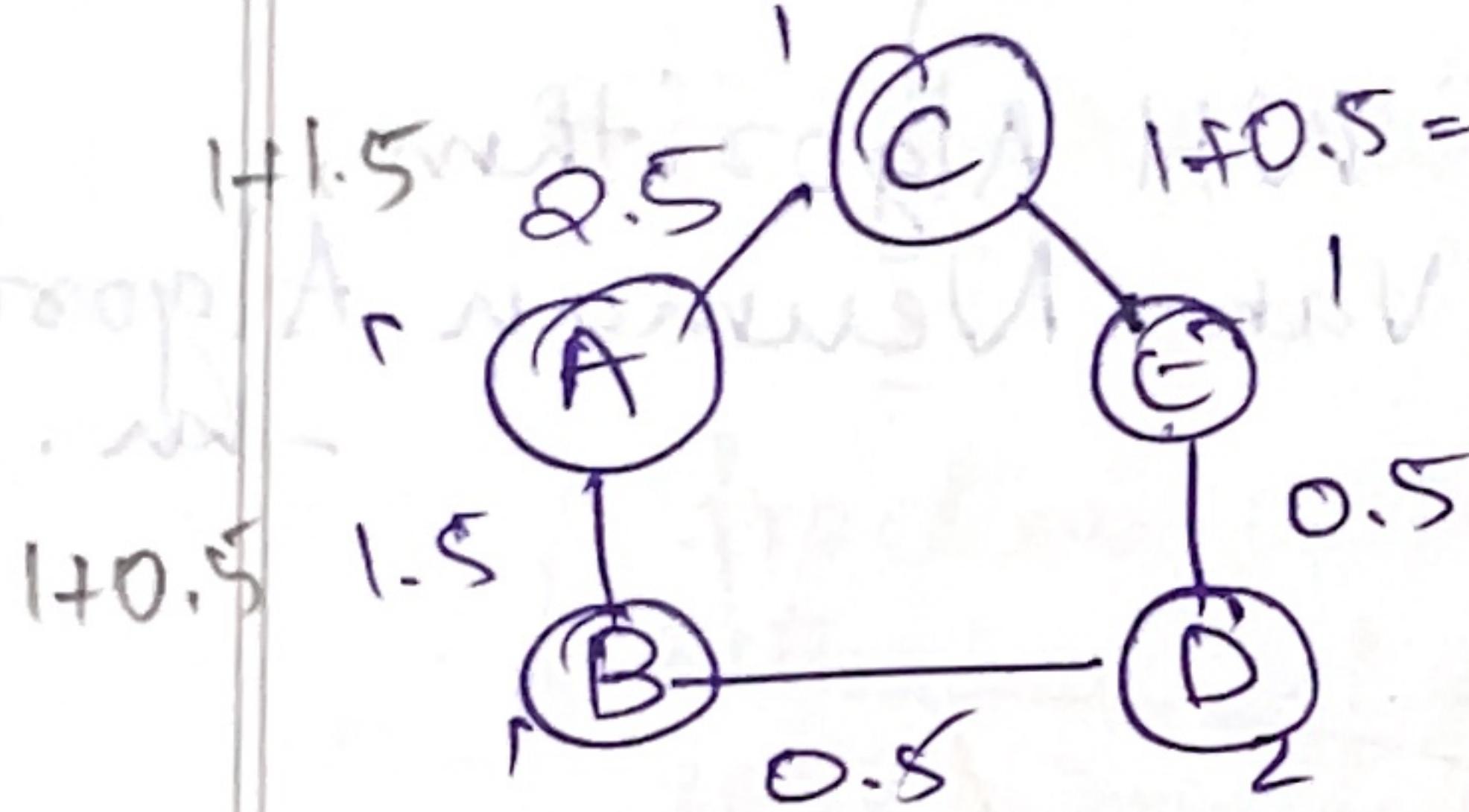
$$2.3 = 2.0 + 2.8 + 2.0 + 2.1 + 2.1 = 10B$$

$$2.2 = 2.8 + 2.0 + 2.1 + 2.0 + 2.0 = 10D$$

$$2.4 = 2.1 + 2.1 + 2.0 + 2.0 + 2.0 = 10D$$

BA extend

9)

A.B.C.

$$AB = 2.5 + 2.5 + 1.5 + 1.5 + 0.5 = 8.5$$

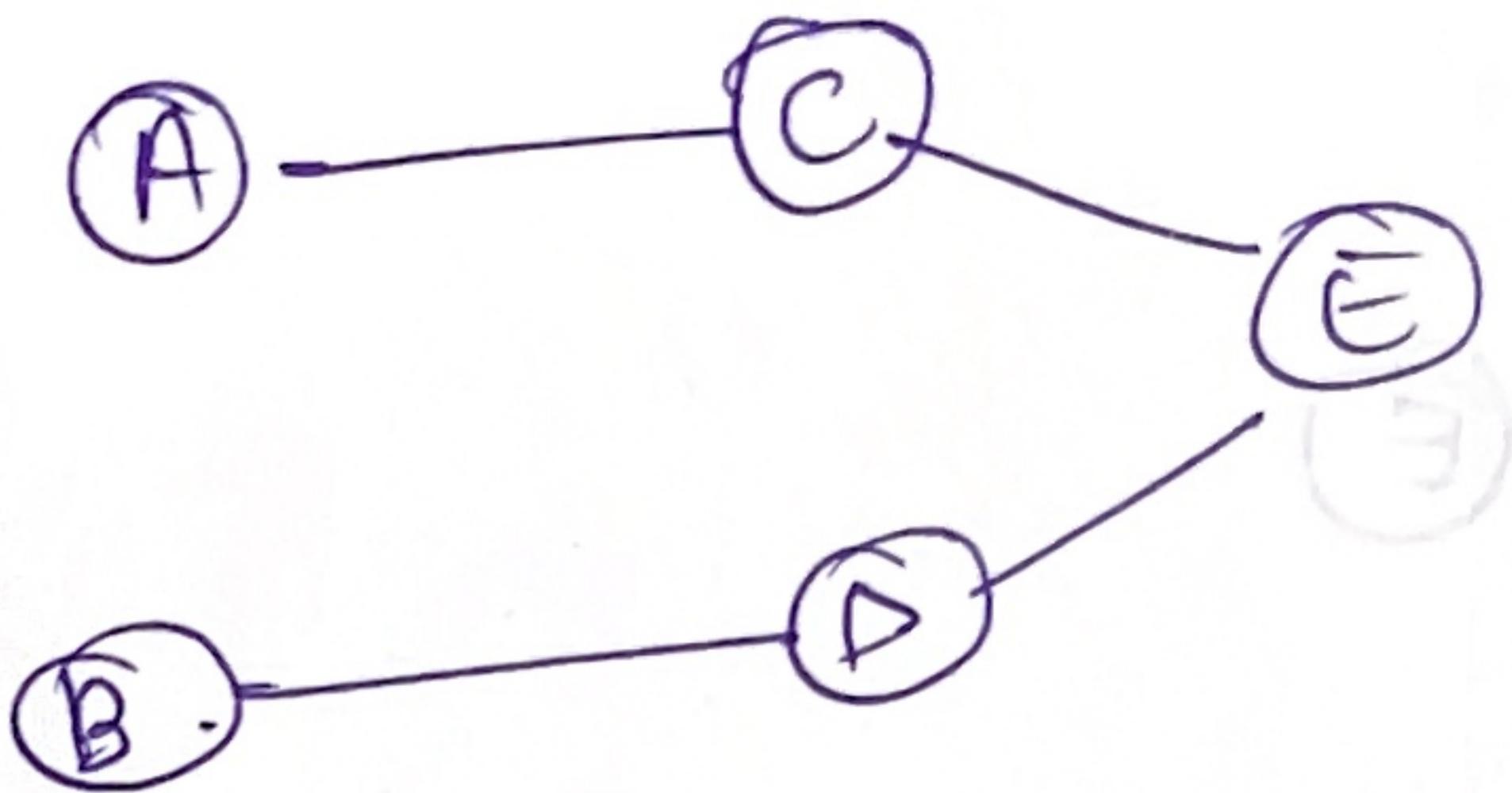
$$AC = 1.5 + 1.5 + 2.5 + 0.5 + 1.5 = 7.5$$

$$BD = 1.5 + 1.5 + 0.5 + 2.5 + 0.5 = 6.5$$

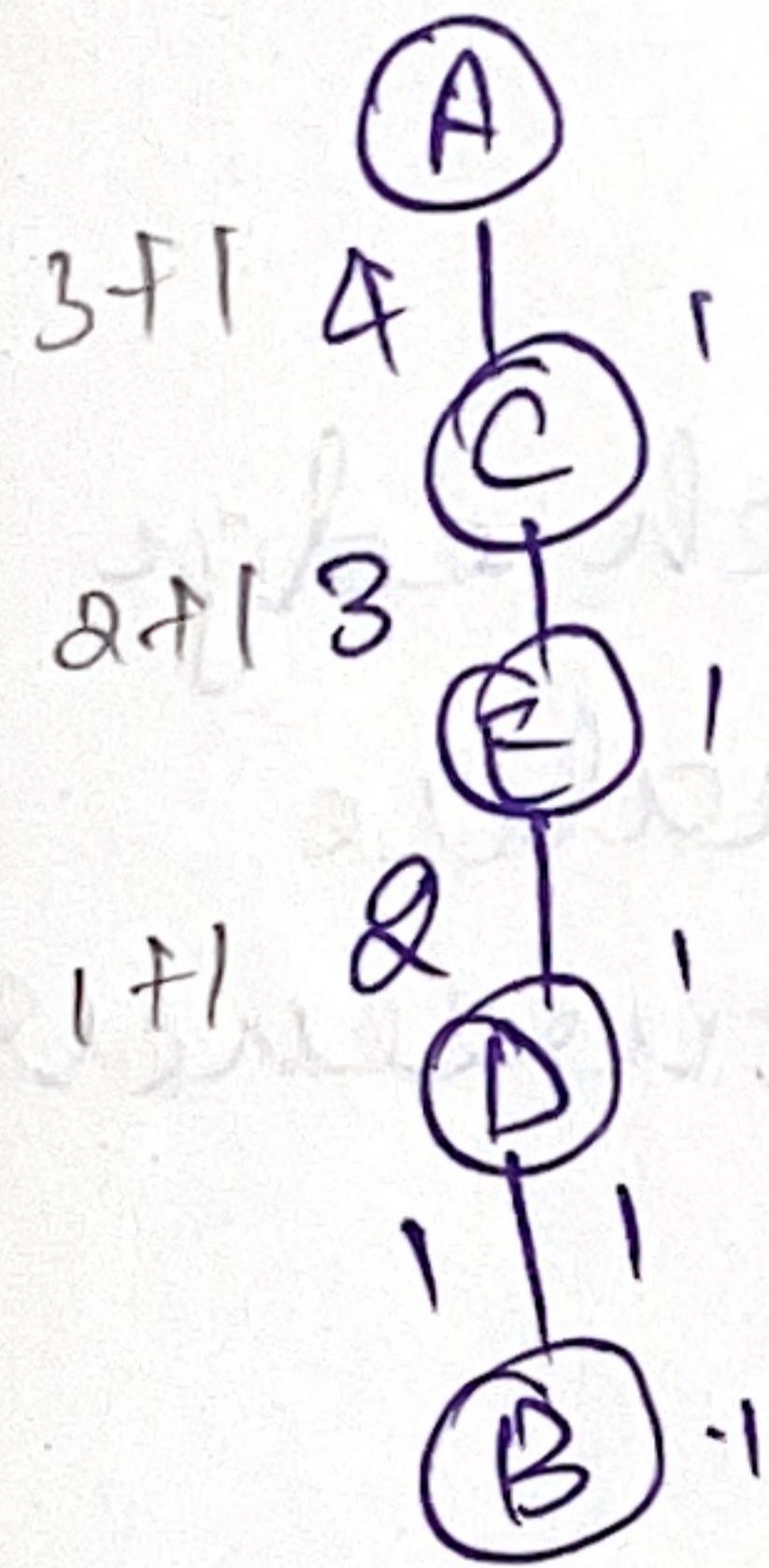
$$CE = 0.5 + 0.5 + 1.5 + 0.5 + 2.5 = 5.5$$

$$DE = 0.5 + 0.5 + 0.5 + 1.5 + 1.5 = 4.5$$

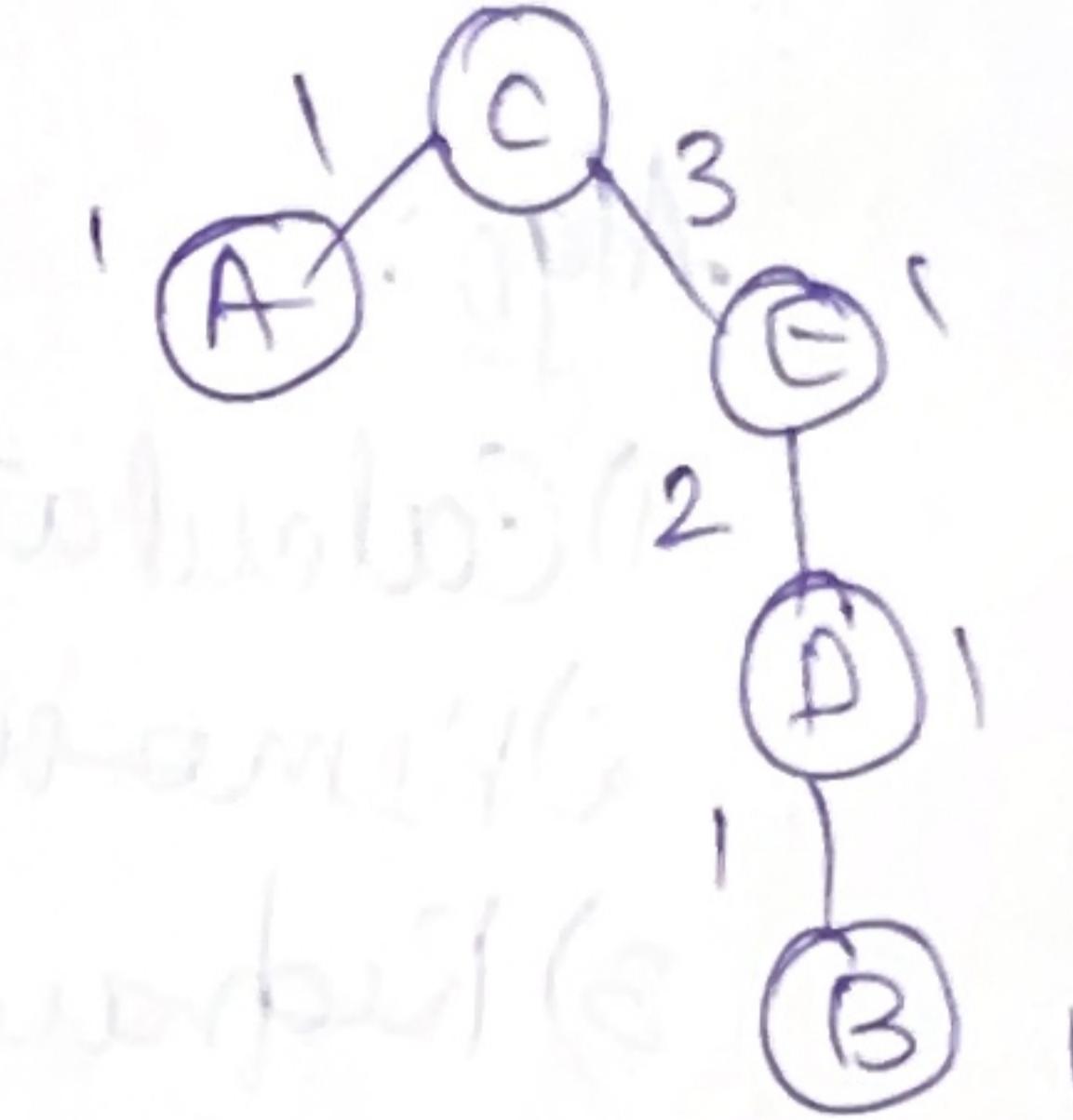
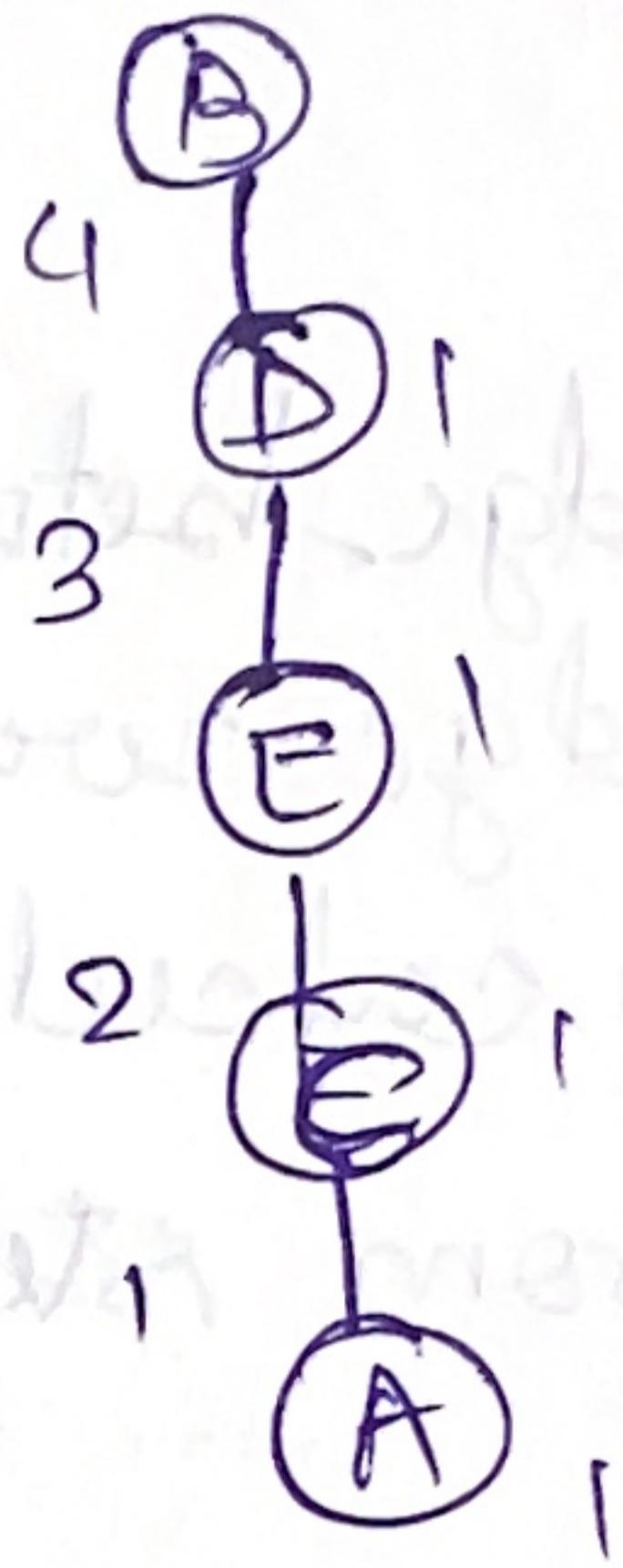
Remove AB .



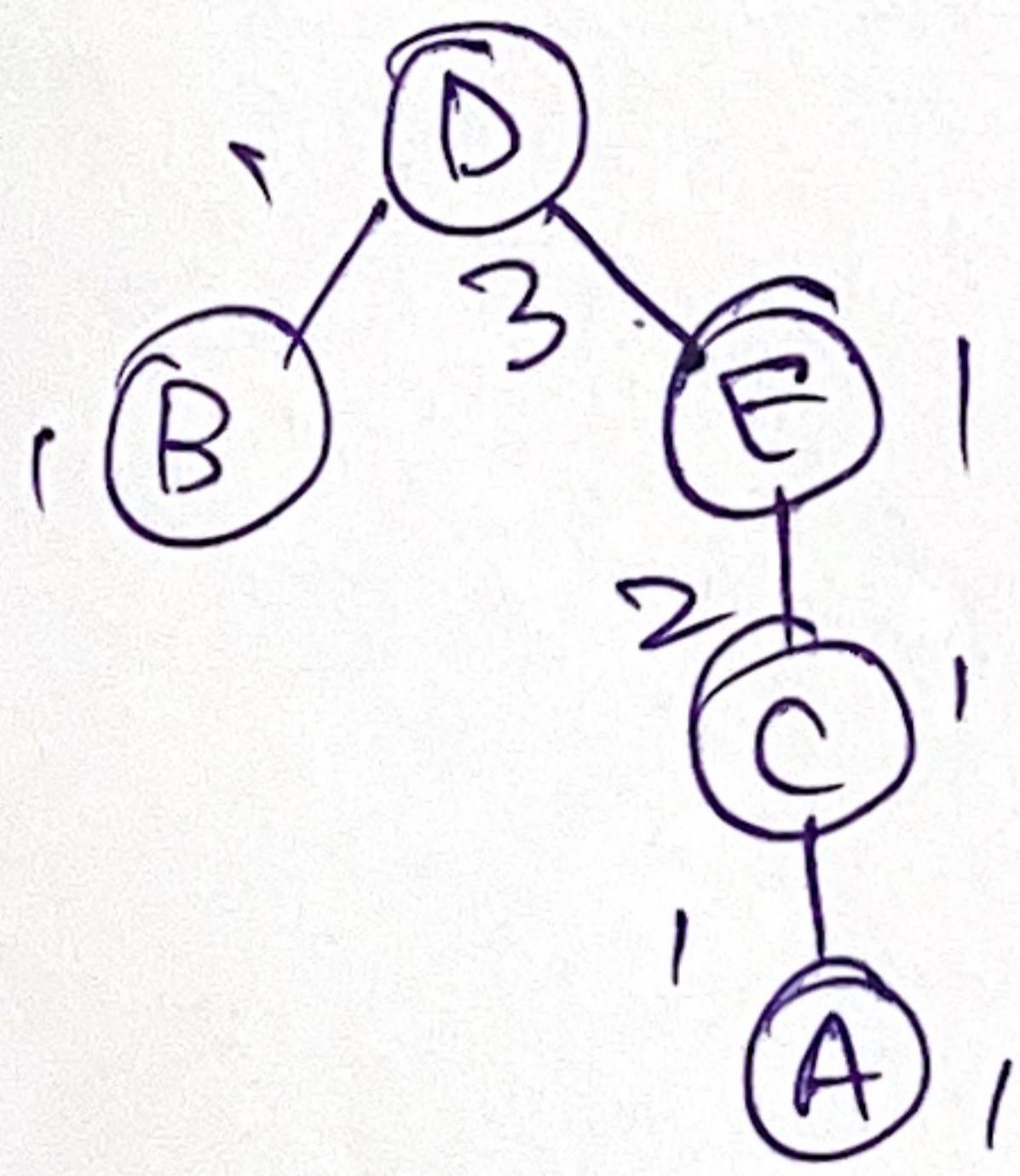
A



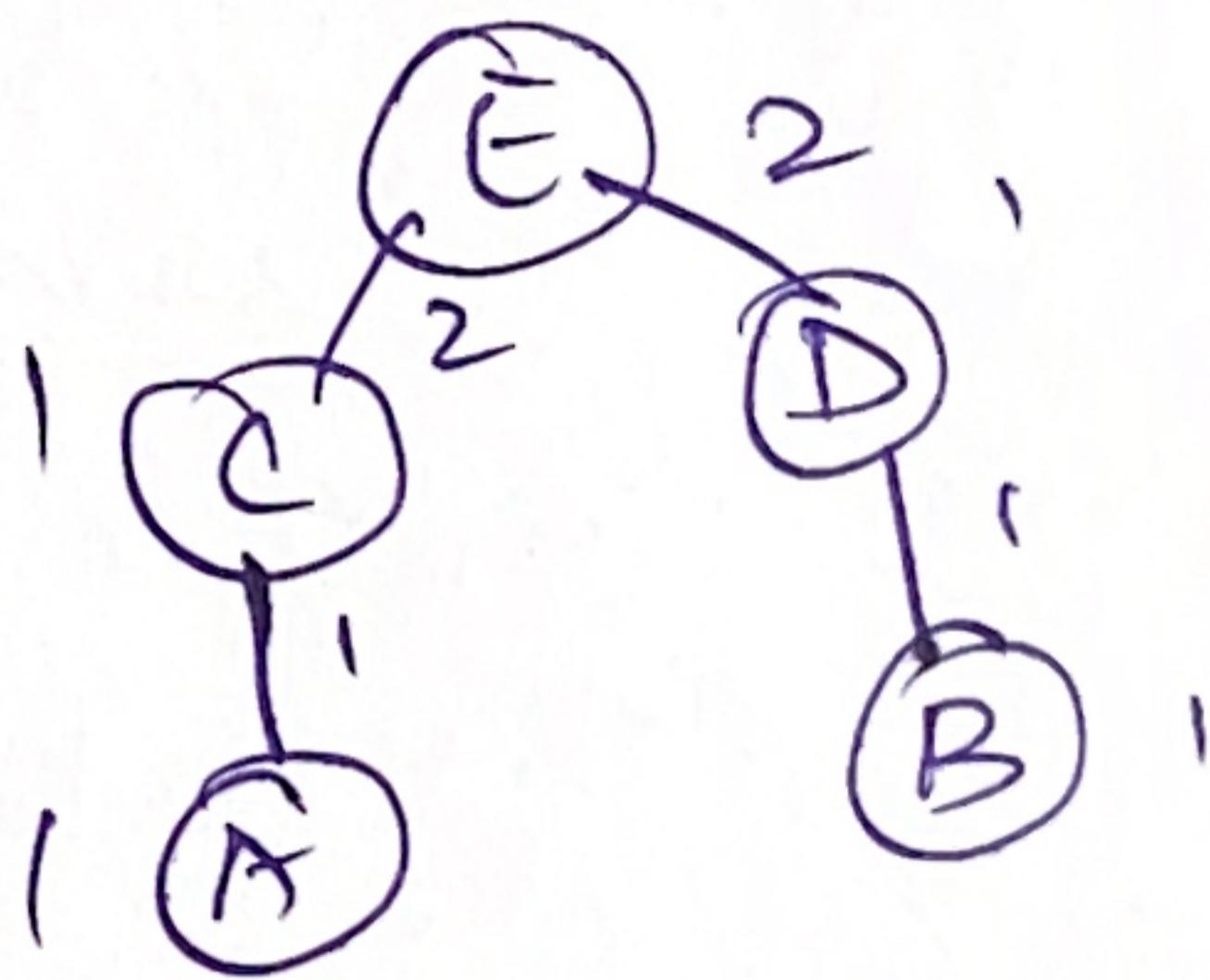
B.



D.



E



$$AC = 4 + 1 + 1 + 1 + 1 = 8$$

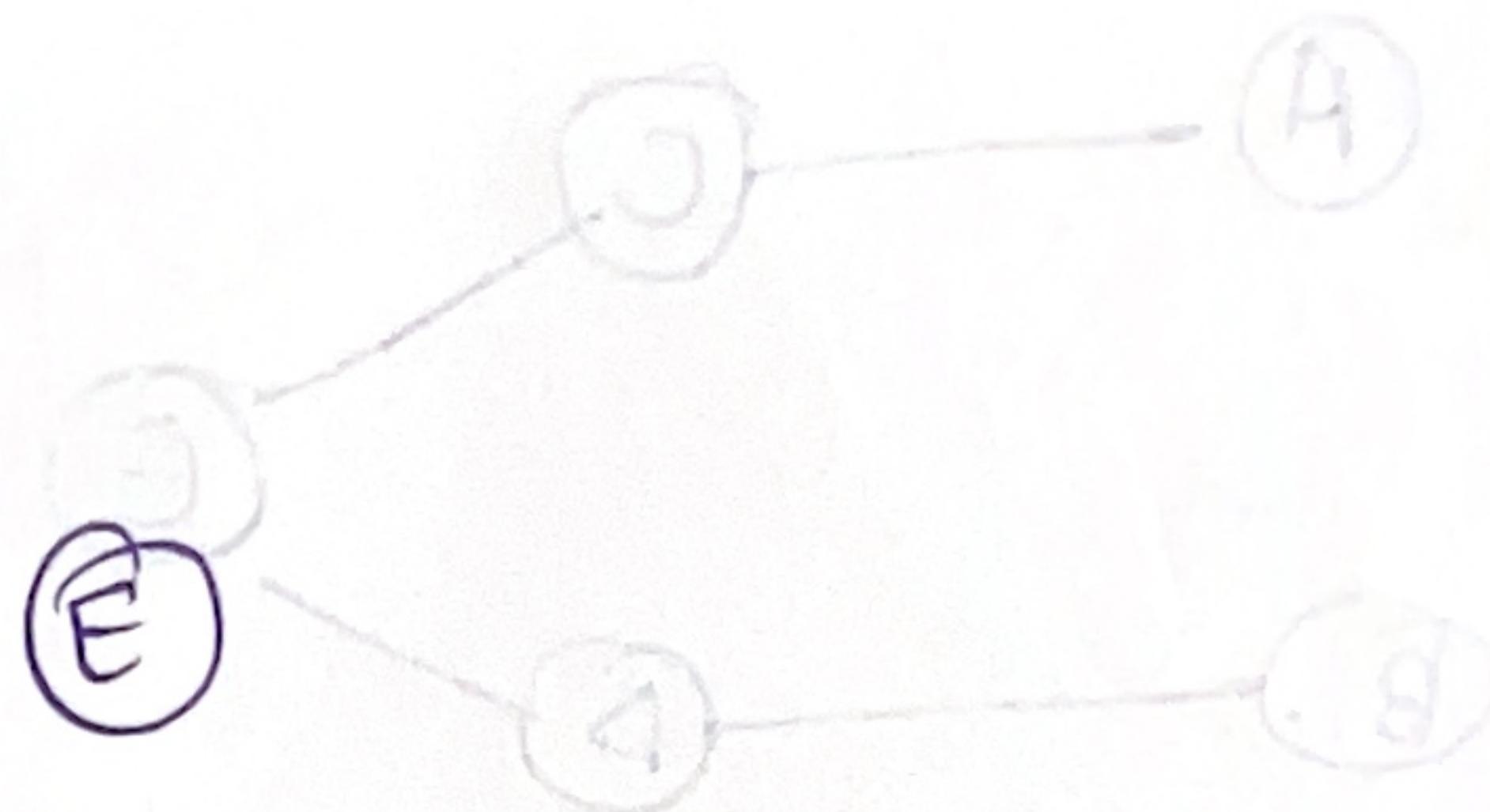
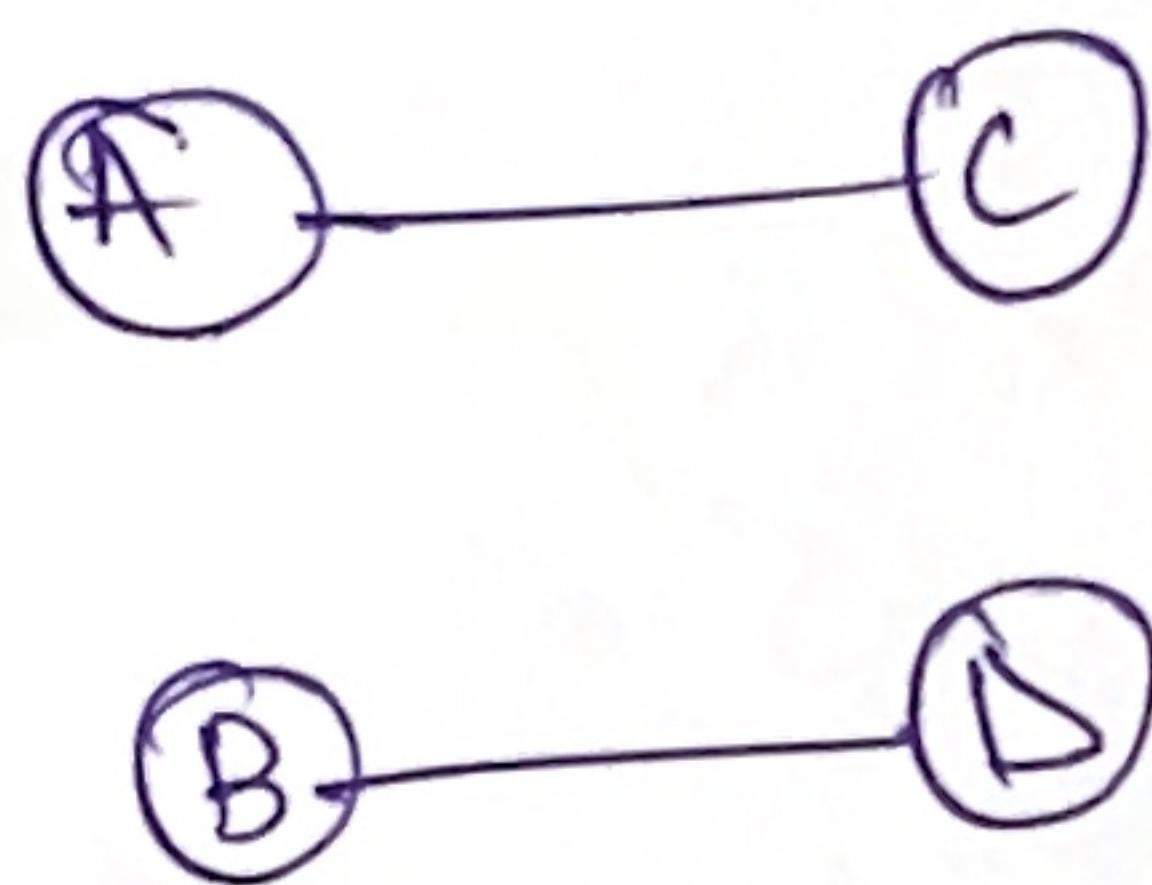
$$CE = 3 + 2 + 3 + 2 + 2 = 12$$

$$ED = 2 + 3 + 2 + 3 + 2 = 12$$

$$DB = 1 + 4 + 1 + 1 + 1 = 8$$

Cut CE & ED.

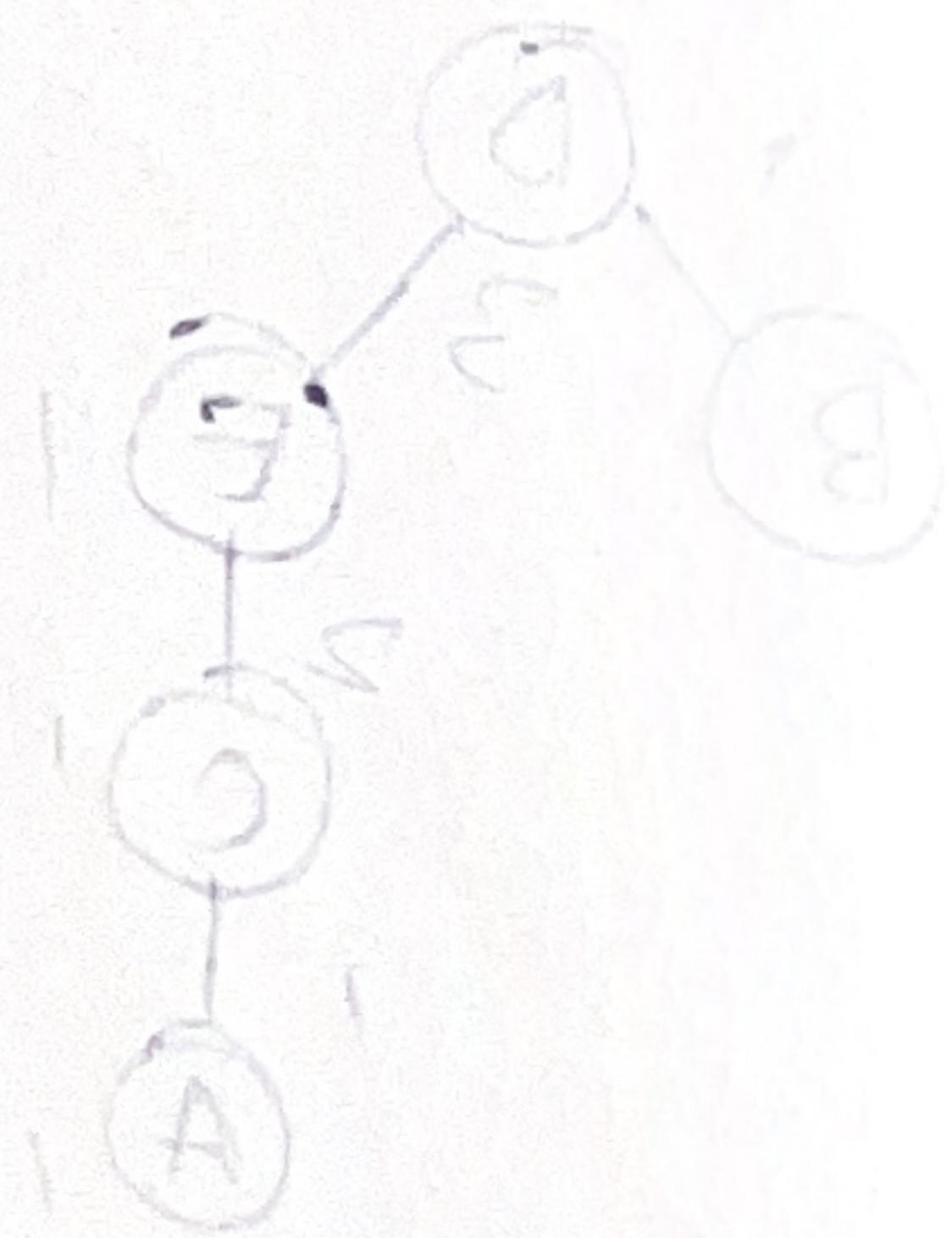
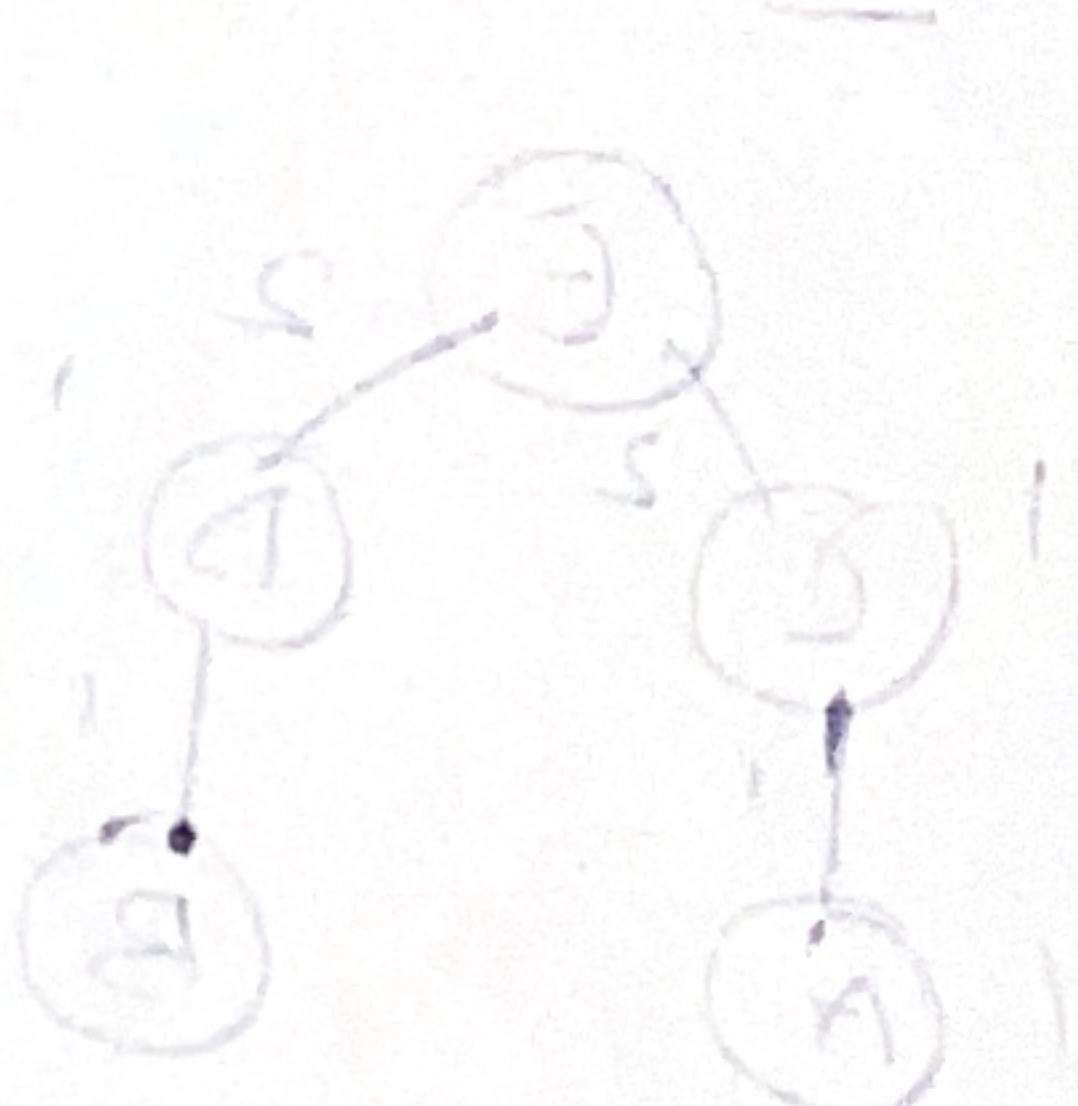
$\therefore$  max. edge centrality



final graph.

Algo:

- 1) Calculate edge betweeness for each edge
- 2) Remove edge with highest value
- 3) Redraw & calculate edge betweeness
- 4) Iterate from step 2.



$$S = 1 + 1 + 1 + 1 + P = 3$$

$$S_1 = 2 + 2 + 8 + 8 + 8 = 30$$

$$S_2 = 8 + 8 + 8 + 8 + 8 = 40$$

$$S = 1 + 1 + 1 + P + 1 = 4$$

GT is 30 true

betweenness of edges.