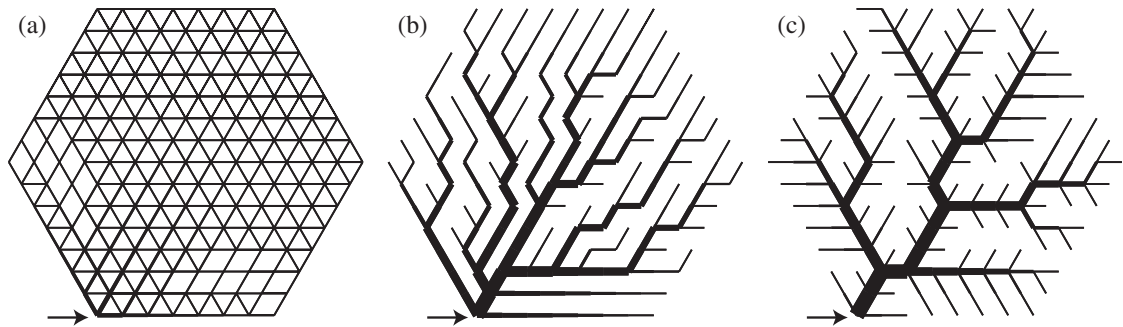




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1. (3 + 3) Reproduce Bohn and Magnasco's Figs. 2a and 2b in [?]:



Steps are given below but please read through the paper to understand how they set things up.

The full team is encouraged to work together on Teams.

- (a) Done (previous assignment): Construct an adjacency matrix \mathbf{A} representing the hexagonal lattice used in [?]. Plot this adjacency matrix.
- (b) Run a minimization procedure to construct Figs. 2a and 2b which correspond to $\gamma = 2$ and $\gamma = 1/2$. Steps:

- i. Set each link's length to 1 (the d_{kl}). The goal then reduces to minimizing the cost

$$C = \sum_{k,l} |I_{kl}|^\Gamma$$

where I_{kl} is the current on link kl and $\Gamma = 2\gamma/(\gamma + 1)$.

- ii. Place a current source of nominal size i_0 at one node (as indicated in Fig. 2 above).

- iii. All other nodes are sinks, drawing a current of

$$i_k = -\frac{i_0}{N_{\text{nodes}} - 1}.$$

- iv. Suggest setting $i_0 = 1000$ (arbitrary but useful value given the size of the network).
- v. Generate an initial set of random conductances for each link, the $\{\kappa_{kl}\}$. From the paper, these must sum to some global constraint as

$$K = \left(\sum_{k,l} \kappa_{kl}^\gamma \right)^{1/\gamma}.$$

This constraint is meant to represent a limitation on the amount of material that can be used to build the network.

Note: There seems to be no reason not to set $K = 1$. However, taking the initial value of K determined by the initial set of random conductances would work.

To our notational peril, we now have a lot of k types on deck.

- vi. Solve the following to determine the potential U at each node, and hence the current on each link using:

$$i_k = \sum_l \kappa_{kl}(U_k - U_l),$$

and then

$$I_{kl} = \kappa_{kl}(U_l - U_k).$$

Note: the paper erroneously has $I_{kl} = R_{kl}(U_l - U_k)$ below equation 4; there are a few other instances of similar miswritings of R_{kl} instead of κ_{kl} .

- vii. Now, use scaling in equation (10) to compute a new set of $\{\kappa_{kl}\}$ from the I_{kl} . Everything boils down to

$$\kappa_{kl} \propto |I_{kl}|^{-(\Gamma-2)},$$

where the constant of proportionality is determined by again making sure $K^\gamma = \sum_{k,l} \kappa_{kl}^\gamma$.

Some help—Let's sort out the key equation:

$$i_k = \sum_l \kappa_{kl}(U_k - U_l).$$

Each time we loop around through this equation, we know the i_k and the κ_{kl} and must determine the U_k . In matrixology, we love $A\vec{x} = \vec{b}$ problems so let's see if we can fashion one:

$$i_k = \sum_l \kappa_{kl}(U_k - U_l)$$

$$\begin{aligned}
&= \sum_l \kappa_{kl} U_k - \sum_l \kappa_{kl} U_l \\
&= U_k \sum_l \kappa_{kl} - \sum_l \mathbf{K}_{kl} U_l \\
&= \lambda_k U_k - [\mathbf{K} \vec{U}]_k
\end{aligned}$$

where we have set $\lambda_k = \sum_l \kappa_{kl}$, the sum of the k th row of the matrix \mathbf{K} . We now construct a diagonal matrix Λ with the λ_k on the diagonal, and obtain:

$$\vec{i} = (\Lambda - \mathbf{K}) \vec{U}.$$

The above is in the form $A\vec{x} = \vec{b}$ so we can solve for \vec{U} using standard features of R, Matlab, Python, ... (hopefully).

Solution:

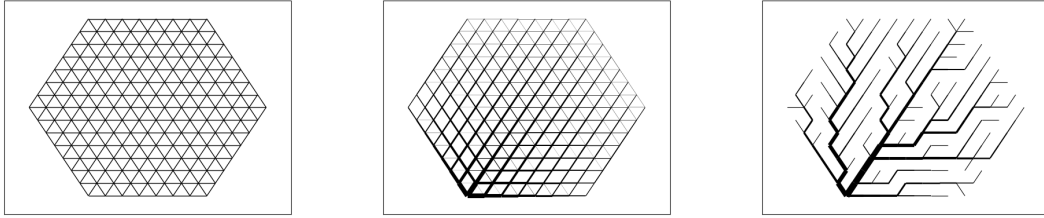


Figure 1: Replication of underlying lattice, and Figs. 2a and 2b from Structure, scaling, and phase transition in the optimal transport network by S. Bohn and M. O. Magnasco

□