

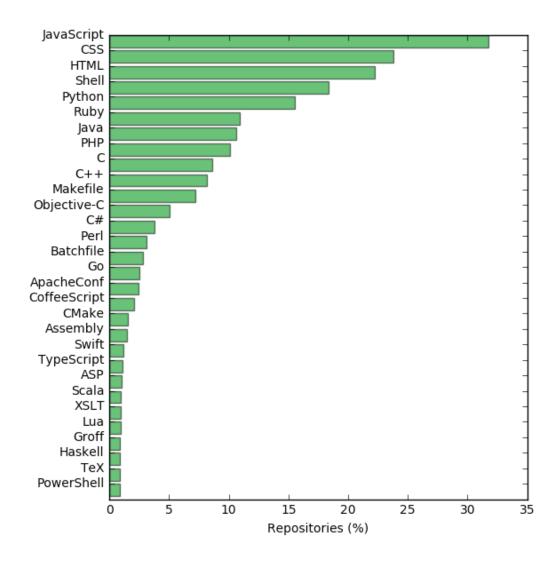
## Matrix Data Visualization

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## GitHub archive dataset

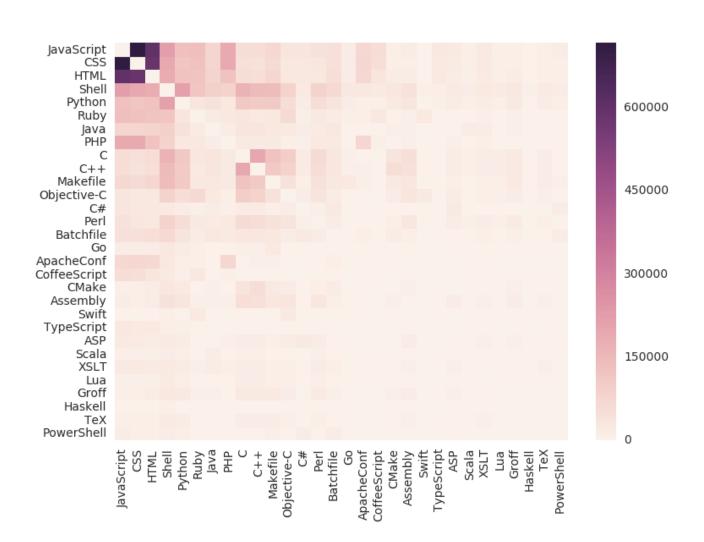
 The plot on this slide shows the percentage of repositories that use specific programming languages

 Suppose our goal is to visualize cooccurrence of pairs of programming languages in different repositories



## Heatmap

- Co-occurrence of pairs of programming languages in different repositories can be represented by a heatmap
- It is important how the rows and columns are sorted to visualize any possibly existing clusters
- In this slide, the rows and columns are sorted in decreasing popularity of programming languages
  - A heuristic revealing some clustering structure, but this is not necessary in general

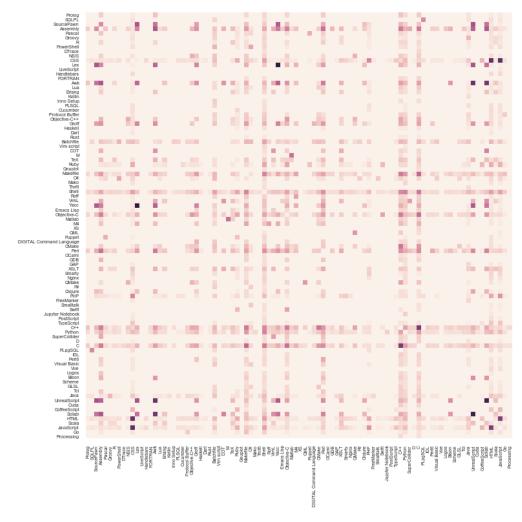


# Use case: similarity matrices

- A similarity matrix quantifies how "similar" pairs of items are
- Similarity may be defined as cosine similarity for items associated with feature vectors: for two non-null feature vectors a and b in  $\mathbf{R}^n$ , the cosine similarity is

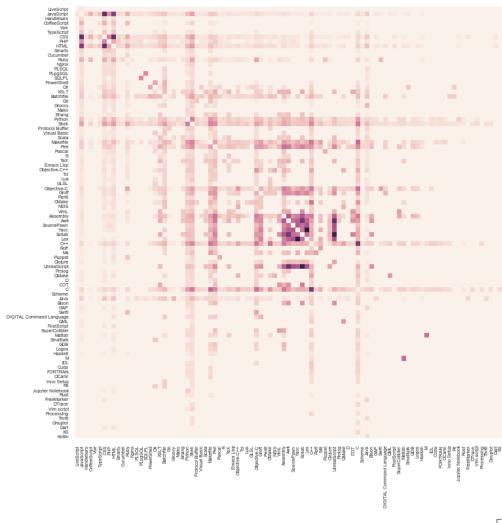
$$sim(a,b) = \frac{\sum_{i=1}^{n} a_i b_i}{\sqrt{\sum_{i=1}^{n} a_i^2} \sqrt{\sum_{i=1}^{n} b_i^2}}$$

 In our example, a feature vector associated with a programming language indicates its usage over different repositories



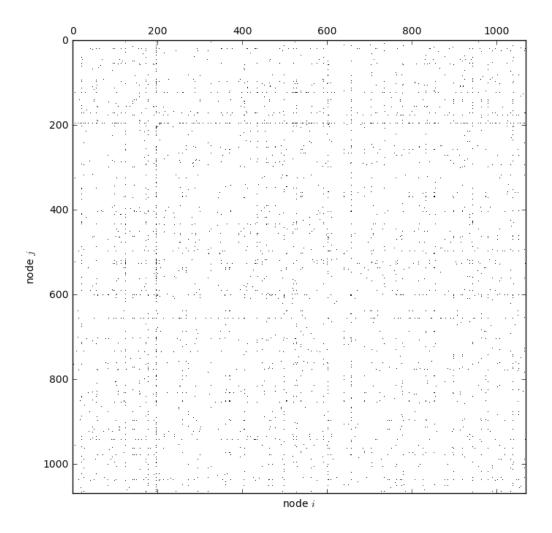
# Use case: similarity matrices (cont'd)

 How can we order rows and columns of a matrix to visualize any possibly existing clusters in matrix data?



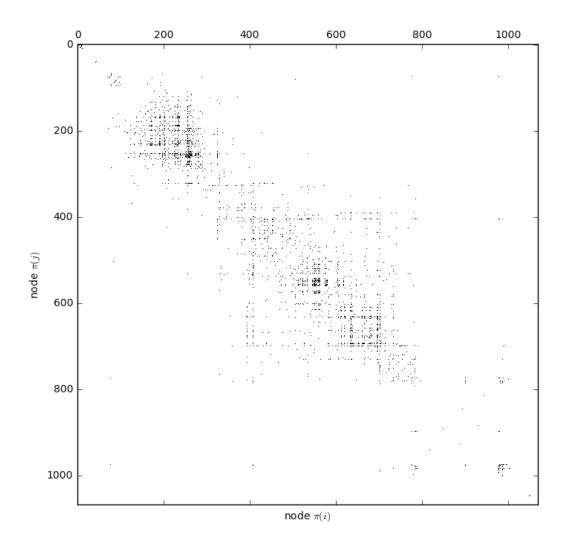
# Use case: adjacency matrices

- Adjacency matrix of a graph indicates whether or not there is an edge between different pairs of nodes of the graph
- The plot in this slide shows the adjacency matrix of a graph that specifies existence of communications between Amazon Mechanical Turk workers
- Dataset source: Yin et al, The Communication Network within the Crowd, WWW 2016



## Use case: adjacency matrices

- Reordered rows and columns
  - i.e. rearranged node identifiers
- The adjacency matrix with reordered rows and columns in this slide suggests existence of different communities of workers



## Software modules

- Scikit-learn biclustering module
  - Scikit-learn Section 2.4 Biclustering

http://scikit-learn.org/stable/modules/biclustering.html

- from sklearn.cluster.bicluster import SpectralCoclustering
- from sklearn.cluster.bicluster import SpectralBiclustering
- R seriation package
  - https://cran.r-project.org/web/packages/seriation/index.html

# Some linear algebra concepts

# Eigenvalues and eigenvectors

•  $\lambda$  is an eigenvalue of matrix A if for some vector  $x \neq 0$ 

$$Ax = \lambda x$$

A corresponding vector x is called an eigenvector

# Laplacian matrix

- Let A be a real symmetric matrix
- The Laplacian matrix  $L_A$  is defined by

$$L_A = D_A - A$$

where  $D_A$  is a diagonal matrix with  $d_{i,i} = \sum_{j=1}^n a_{i,j}$ 

• A is a real symmetric matrix  $\Rightarrow L_A$  is a real symmetric matrix

# Laplacian matrix (cont'd)

- For any real symmetrix matrix  $A \in \mathbb{R}^{n \times n}$ :
  - A has n eigenvectors
  - All eigenvectors of *A* are pairwise orthogonal
  - All eigenvalues of A are real
- Every Laplacian matrix has all its eigenvalues real and non-negative, which is equivalent to saying that it is positive definite
- Every Laplacian matrix has the vector of all ones  $\boldsymbol{e}$  as an eigenvector corresponding to the eigenvalue zero

# Fiedler value and eigenvector

- For real symmetric matrix  $A \in \mathbb{R}^{n \times n}$ , Fiedler value is defined as the minimum eigenvalue of the Laplacian  $L_A$  that has an eigenvector orthogonal to e
- The corresponding eigenvector is called a Fiedler eigenvector
- The Fiedler value is the optimum value of the optimization problem:

minimize 
$$x^T L_A x$$
  
subject to  $x^T e = 0$   
 $x^T x = 1$ 

#### Historical remarks

- Miroslav Fiedler
- Czech Republic's mathematician
- 1926-2015
- Charles University, Prague



• M. Fiedler, Algebraic connectivity of graphs, Czehoslovak Math. J., 23(98), 1973

# Laplacian matrix (cont'd)

• **Lemma**: For any real, symmetric matrix *A*:

$$x^T L_A x = \sum_{i < j} a_{i,j} (x_i - x_j)^2$$

- Laplacian matrices are closely related to graph cuts
  - Graph G = (V, E) with edge weights: edge (i, j) has weight  $a_{i,j}$
  - A is the adjacency matrix of G
  - Let x takes value in  $\{-1,1\}^n$  defining a vertex cut: partitioning vertices into two sets (negative and positive labeled)
  - Graph cut is defined as the sum of weights of edges whose end vertices belong to different components of the vertex cut

# Seriation

## Seriation

- Input: a real, symmetric matrix  $A \in \mathbb{R}^{n \times n}$ 
  - $a_{i,j}$  interpreted as the similarity between items i and j
- Goal: find a linear ordering (permutation) of items such that similar items are placed nearby, specifically, find a permutation  $\pi^*$  that minimizes

$$c(\pi) = \sum_{i < j} a_{i,j} (\pi_i - \pi_j)^2$$

over the set of all possible permutations of n elements  $\Pi_n$ 

This problem is NP hard

#### Fractional relaxation

• Find optimal solution to the following problem:

minimize 
$$c(x)$$
  
subject to  $x^T e = 0$   
 $x^T x = 1$   
 $x \in \mathbf{R}^n$ 

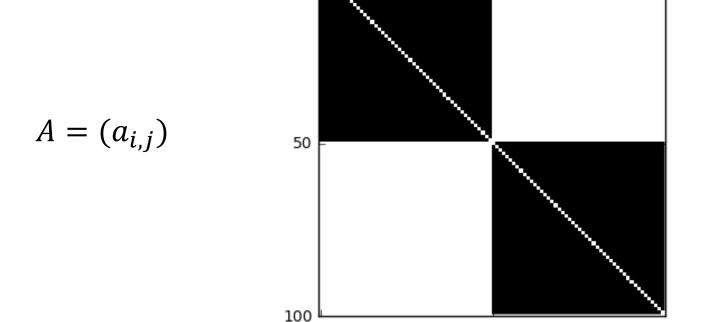
- The first constraint avoids multiplicity of solutions by adding a positive constant to each coordinate of x
- The second constraint avoids vector 0 to be a trivial solution
- The optimum solution is a Fiedler vector of A

## Robinson matrices

- A matrix A is said to be a Robinson matrix (R-matrix) if, and only if,
  - *A* is symmetric
  - $a_{i,j} \le a_{i,k}$  for j < k < i and  $a_{i,j} \ge a_{i,k}$  for i < j < k

- In other words, a matrix is a R-matrix if it is symmetric and it has off-diagonal elements non-decreasing by moving away from the diagonal
- A matrix A is said to be a pre-R matrix if it can be symmetrically permuted (reordering rows and columns by the same permutation) to become an R-matrix

# An example of a R-matrix



50

100

- Clearly, a symmetric matrix
- The elements are decreasing as we move away from the diagonal along any row

black point if  $a_{i,j} = 1$ white point if  $a_{i,j} = 0$ 

# An example R-matrix: stochastic block model

- Stochastic block model is a random graph model commonly studied in the community detection and graph clustering literature
- A stochastic block model can be defined by
  - Parameters  $0 \le q \le p \le 1$
  - Set of vertices  $V = \{1, 2, \dots, n\}$
  - Hidden bipartition of vertices  $(S, V \setminus S)$
  - Set of edges E: for every pair of vertices (i, j) such that i < j we have

$$(i,j) \in E$$
 independently with probability = 
$$\begin{cases} p & \text{if } i,j \in S \text{ or } i,j \in V \setminus S \\ q & \text{otherwise} \end{cases}$$

- A is defined as the adjacency matrix of G
- The case q=0 and p=1 is trivial when the vertex-set partition is not hidden

## Fiedler vector of R-matrices

• **Theorem**: If A is a R-matrix then it has a monotone Fiedler vector.

• Proof: Atkins et al (1998)

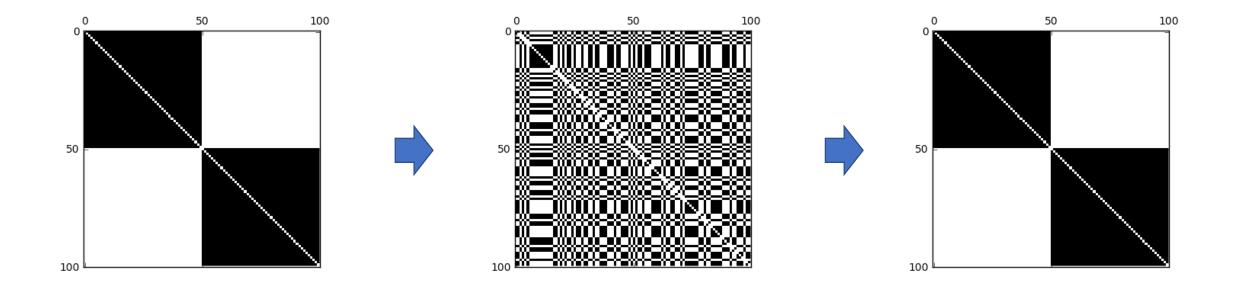
## Fiedler vector of R-matrices

• **Theorem**: Let A be a pre-R matrix with a simple Fiedler value and a Fiedler vector with no repeated elements. Let  $\pi_1$  be the permutation sorting elements of the Fiedler vector in increasing order. Let  $\pi_2$  be the permutation sorting elements of the Fiedler vector in decreasing order. Let  $\Pi_1$  and  $\Pi_2$  be the corresponding permutation matrices.

Then,  $\Pi_1 A \Pi_1$  and  $\Pi_2 A \Pi_2$  are permutation matrices and not other symmetric permutation of A produces an R matrix.

Proof: Atkins et al (1998)

# Example 1: p = 1, q = 0

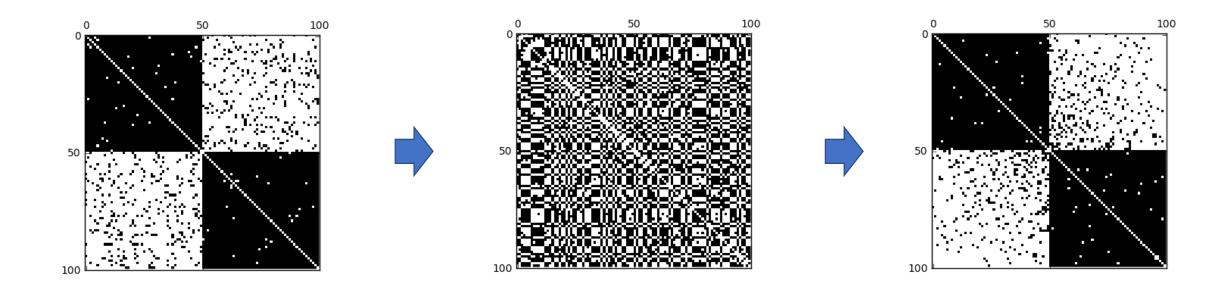


Rows and columns permuted by the same random permutation

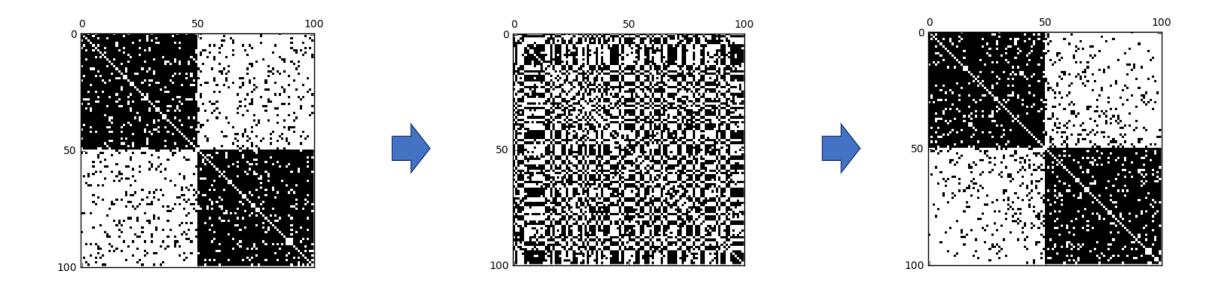
"symmetric permutation"

Rows and columns sorted in a monotonic order of the Fiedler vector elements

# Example 2: p = 0.99, q = 0.01



# Example 3: p = 0.9, q = 0.1



# Spectral co-clustering

# Graph cuts

- Let V be a set of vertices with weights  $w_{i,j}$  for  $i,j \in V$
- k-way partition: let  $P_k(S)$  be the set of all possible partitions of S in k components
  - In particular,  $(S_1, S_2) \in P_2(S)$  is referred as a bipartition
- The cut function is defined as the sum of weights of cut edges:

$$cut(V_1, V_2) = \sum_{i \in V_1, j \in V_2} w_{i,j}$$

More generally, we define

$$\operatorname{cut}(V_1, V_2, \dots, V_k) = \sum_{1 \le i < j \le k} \operatorname{cut}(V_i, V_j)$$

## Graph cuts and the Laplacian matrix

• A bipartition  $(V_1, V_2)$  can be represented by a partition vector x defined as

$$x_i = \begin{cases} 1 & \text{if } i \in V_1 \\ -1 & \text{if } i \in V_2 \end{cases}$$

• Lemma: Any bipartition  $(V_1, V_2)$  and the corresponding partition vector x satisfy

$$\frac{x^T L_A x}{x^T x} = \frac{1}{n} 4 \operatorname{cut}(V_1, V_2)$$

## Normalized cut function

- Define the weight of a set of vertices  $S: w(S) = \sum_{i \in S} w_i$
- The normalized cut function is defined as follows:

$$Q(V_1, V_2) = \frac{\text{cut}(V_1, V_2)}{w(V_1)} + \frac{\text{cut}(V_1, V_2)}{w(V_2)}$$

- This cut function captures both sparsity of edge cuts and balancing of the component sizes
- Optimization problem formulation: minimize  $Q(V_1, V_2)$  over  $P_2(V)$

# Special normalized cut functions

#### Ratio-cut

- Each vertex has a unit weight
- Amounts to looking at the eigenvalue problem:  $L_A y = \lambda y$

#### Normalized-cut

- Each vertex weight is equal to the sum of weights of incident edges
- In this case  $w(V_i) = \operatorname{cut}(V_1, V_2) + w^{int}(V_i)$ where  $w^{int}(V_i) = \sum_{(u,v) \in E: u,v \in V_i} w_{u,v}$
- Amounts to the generalized eigenvalue problem :  $L_A y = \lambda D_A y$

## Normalized-cut function

Show that

$$Q(V_1, V_2) = 2 - \left(\frac{w^{int}(V_1)}{w(V_1)} + \frac{w^{int}(V_2)}{w(V_2)}\right)$$

 Minimizing the normalized-cut is equivalent to maximizing the proportion of edge weights that lie within each component of a vertex set partition

# Generalized partition-vector representation

Let y be a generalized partition vector defined as

$$y_{i} = \begin{cases} \sqrt{\frac{w(V_{2})}{w(V_{1})}} & if \ i \in V_{1} \\ -\sqrt{\frac{w(V_{1})}{w(V_{2})}} & if \ i \in V_{2} \end{cases}$$

• Properties: (a)  $y^T D_w y = w(V)$  and (b)  $y^T D_w e = 0$ 

• Lemma:

$$\frac{y^T L_A y}{y^T D_w y} = \frac{\text{cut}(V_1, V_2)}{w(V_1)} + \frac{\text{cut}(V_1, V_2)}{w(V_2)}$$

## Fractional relaxation

minimize 
$$\frac{y^T L_A y}{y^T D_W y}$$
 subject to  $y^T D_W e = 0$ 

• **Theorem**: The solution is the eigenvector corresponding to the second smallest eigenvalue  $\lambda_2$  of the generalized eigenvalue problem:

 $y \in \mathbb{R}^n$ 

 $y \neq 0$ 

$$L_A y = \lambda D_w y$$

• Corollary: The optimum value of the min normalized cut is  $\geq \lambda_2$ 

# Bipartite graph clustering

- Bipartite graph
  - The vertex set consists of disjoint sets of vertices L and R
  - Each edge has its vertices in L and R
- Let  $D_L$  and  $D_R$  be diagonal matrices with diagonal elements

$$(D_L)_{i,i} = \sum_{j \in R} w_{i,j} \text{ and } (D_R)_{i,i} = \sum_{j \in L} w_{j,i}$$

• For graph *G*:

$$A = \begin{pmatrix} 0 & W \\ W^T & 0 \end{pmatrix}$$
,  $D_A = \begin{pmatrix} D_L & 0 \\ 0 & D_R \end{pmatrix}$  and  $L_A = \begin{pmatrix} D_L & -W \\ -W^T & D_R \end{pmatrix}$ 

# The generalized eigenvalue problem

• The generalized eigenvalue problem can be written as

$$D_L x - W y = \lambda D_L x$$
  
-  $A^T x + D_R y = \lambda D_R y$ 

- Assumption: W has a strictly positive element in each row and column
- Note that we can write:

$$D_L^{1/2} x - D_L^{-1/2} W y = \lambda D_L^{1/2} x$$
  
-  $D_R^{-1/2} A^T x + D_R^{1/2} y = \lambda D_R^{1/2} y$ 

# Change of variables

Using the change of variables

$$\widetilde{W} = D_L^{-1/2} W D_R^{-1/2}$$

$$u = D_L^{1/2} x$$

$$v = D_R^{1/2} y \text{ and}$$

$$\sigma = 1 - \lambda$$

we can write

$$\widetilde{W}v = \sigma u$$
 and  $\widetilde{W}^T u = \sigma v$ 

## Eigenvectors

• The eigenvector  $x_2$  corresponding to the second smallest eigenvalue  $\lambda_2$  of the generalized eigenvalue problem can be written as:

$$x_2 = \begin{pmatrix} D_L^{-1/2} u_2 \\ D_R^{-1/2} v_2 \end{pmatrix}$$

where  $u_2$  and  $v_2$  are the left and right singular vectors of  $\widetilde{W}$  corresponding to the singular value  $\sigma_2=1-\lambda_2$ 

• We can think of  $u_2$  to give a partition of the set of left vertices and  $v_2$  to give a partition of the set of right vertices

# Bi-clustering algorithm

- Input: W
- Compute  $\widetilde{W} = D_L^{-1/2} W D_R^{-1/2}$
- Compute the left and right singular vectors  $u_2$  and  $v_2$  of  $\widetilde W$  corresponding to the singular value  $\sigma_2=1-\lambda_2$
- Partition the set of vertices in two components using k-means algorithm for input data points  $x_2$

# k-way clustering algorithm

- Given a positive integer  $k \ge 2$  the goal is to partition the set of left vertices and the set of right vertices in k components
- Let  $U=(u_2,u_3,\dots,u_{\ell+1})$  and  $V=(v_2,v_3,\dots,v_{\ell+1})$  be  $\ell$  the left and right singular vectors

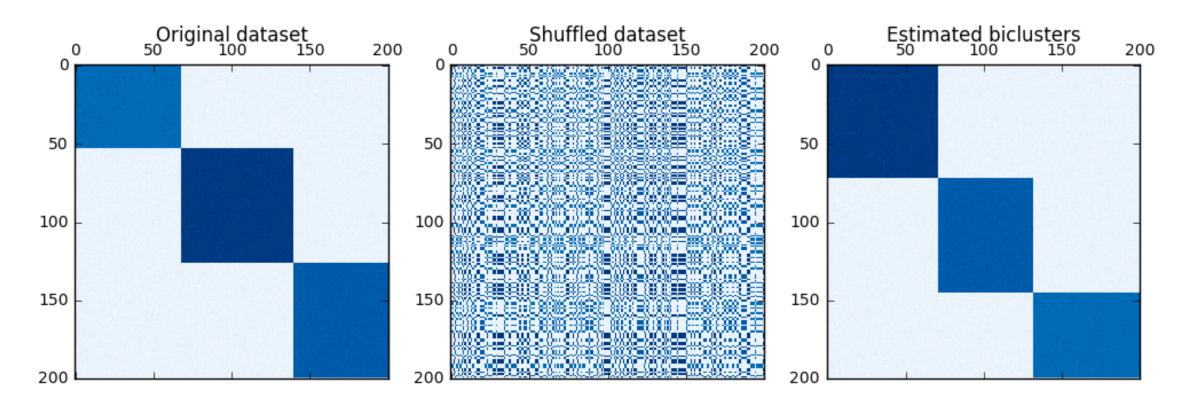
• Let 
$$(x_2, x_3, \dots, x_{\ell+1}) = \begin{pmatrix} D_L^{-1/2} U \\ D_R^{-1/2} V \end{pmatrix}$$

• Apply the k-means algorithm to the input  $\ell$ -dimensional points

# Evaluating a bi-clustering: consensus score

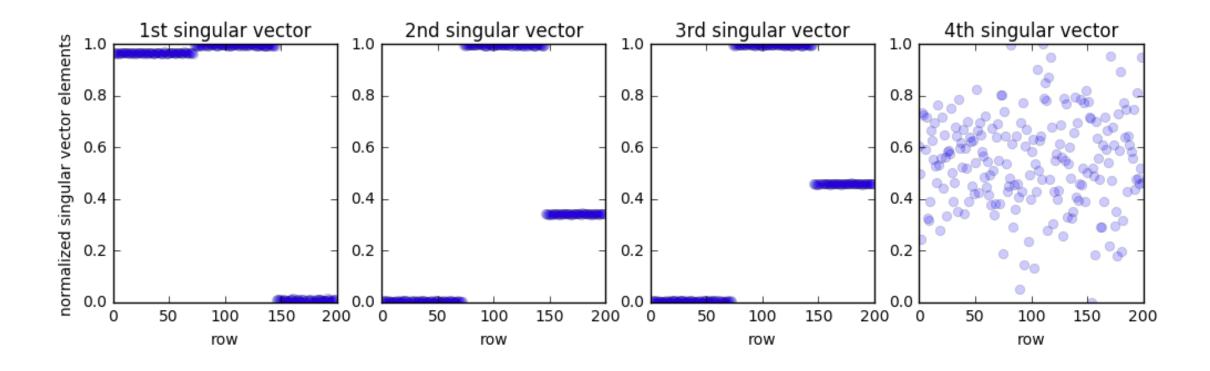
- The definition of the consensus score as implemented in sklearn.metrics.consensus\_score
- Quantifies the similarity of two sets of bi-clusters
- Similarity between individual bi-clusters is computed
- Then the best matching between sets is found using the Hungarian algorithm
- The final score is the sum of similarities divided by the size of the larger set

# Example



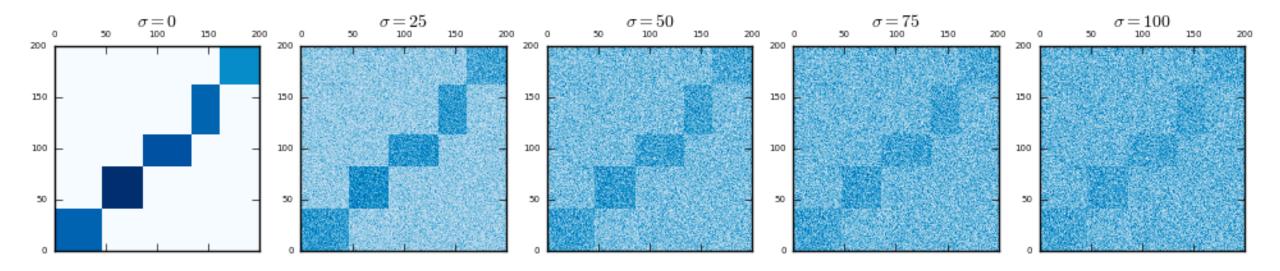
Recovery of hidden co-clusters by spectral co-clustering

# Example: singular vectors



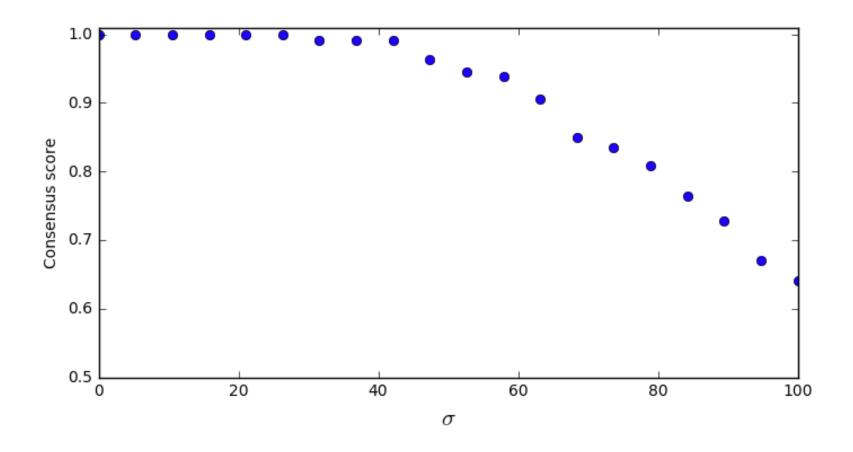
Clusters are clearly indicated by the 2<sup>nd</sup> and 3<sup>rd</sup> singular vectors

## Robustness to noise



• Input matrix corrupted by noise with varying variance  $\sigma$ 

# Robustness to noise (cont'd)



• Consensus score vs standard deviation of noise

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