

An introduction to graph analysis and modeling

The Stochastic block model

MSc in Statistics for Smart Data – ENSAI

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<http://julien.cremeriefamily.info>



Motivations

Last time: find an underlying organization in a observed network

Spectral or hierachical clustering for network data

↪ Not model-based, thus no statistical inference possible

Today: clustering of network based on a probabilistic model of the graph

Become familiar with

- the stochastic block model, a random graph model tailored for clustering vertices,
- the variational EM algorithm used to infer SBM from network data.

hierarchical clustering ↔ Gaussian mixture models



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Outline

- 1 Background: mixture models and EM
 - Mixture models
 - Expectation-Maximization algorithm
 - Example: mixture of Gaussians
- 2 The Stochastic Block Model (SBM)
 - Some Graphs Models and their limitations
 - Mixture of Erdős-Rényi and the SBM
 - Inference in SBM with variational EM


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Latent variables models

Definition

A **latent variable model** is a statistical model that relates, for $i = 1, \dots, n$ individuals,

- a set of **manifest** (observed) variables $\mathbf{X} = (X_i, i = 1, \dots, n)$ to
- a set of **latent** (unobserved) variables $\mathbf{Z} = (Z_i, i = 1, \dots, n)$.

Common assumption: conditional independence

$$\mathbb{P}((X_1, \dots, X_n) | (Z_1, \dots, Z_n)) = \prod_{i=1}^n \mathbb{P}(X_i | Z_i).$$

Famous examples

- $(Z_i, i \geq 1)$ is Markov chain: **Markov models**
- Z_i categorical and independent: **mixture models**

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- what if $X_i = X_{i,j}$ is a collection of edges in a graph?

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- Z_i categorical and independent: **mixture models**
- what if $X_i = X_{i'j'}$ is a collection of edges in a graph?

Mixture models: the latent variables

When (Z_1, \dots, Z_n) are independent categorical variables, they give a **natural (latent) classification of the observations** (X_1, \dots, X_n) – or **labels**.

Notations

Let (Z_1, \dots, Z_n) be *iid* categorical variables with distribution

$$\mathbb{P}(i \in q) = \mathbb{P}(Z_i = q) = \alpha_q, \quad \text{s.t.} \sum_{q=1}^Q \alpha_q = 1.$$

Alternative (equivalent) notation

Let $Z_i = (Z_{i1}, \dots, Z_{iQ})$ be an indicator vector of label for i :

$$\mathbb{P}(i \in q) = \mathbb{P}(Z_{iq} = 1) = \alpha_q, \quad \text{s.t.} \sum_{q=1}^Q \alpha_q = 1.$$

By definition, $Z_i \sim \mathcal{M}(1, \alpha)$, with $\alpha = (\alpha_1, \dots, \alpha_Q)$.

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Mixture models: the manifest variables

A mixture model represents the **presence of subpopulations** within an overall population as follows:

$$\mathbb{P}(X_i) = \sum_{z_i \in \mathcal{Z}_i} \mathbb{P}(X_i, Z_i) = \sum_{Z_i \in \mathcal{Z}_i} \mathbb{P}(X_i | Z_i) \mathbb{P}(Z_i).$$

Conditional distribution of the manifest variables

We assume a **parametric distribution** of X in each subpopulation

$$X_i | \{Z_i = q\} \sim \mathbb{P}_{\theta_q} \quad \left(\Leftrightarrow X_i | \{Z_{iq}\} = 1 \sim \mathbb{P}_{\theta_q} \right)$$

The specificity of each class is handled by $\{\theta_q\}_{q=1}^Q$.

Mixture models: likelihoods

The complete-data likelihood

It is the joint distribution of (X_i, Z_i) :

$$\mathbb{P}(X_i, Z_i) = \alpha_{Z_i} \mathbb{P}_{\theta_{Z_i}}(X_i)$$

The incomplete-data likelihood

It is the marginal distribution of X_i once Z_i integrated:

$$\mathbb{P}(X_i) = \sum_{q=1}^Q \mathbb{P}(X_i, Z_i = q) = \sum_{q=1}^Q \alpha_q \mathbb{P}_{\theta_q}(X_i)$$

↪ A mixture model is a sum of distributions weighted by the proportion of each subpopulation.

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Intractability of the Likelihood

Maximum Likelihood Estimator

The MLE aims to maximize the (marginal) likelihood of the observations:

$$L(\boldsymbol{\theta}; \mathbf{X}) = \mathbb{P}_{\boldsymbol{\theta}}((X_1, \dots, X_n)) = \int_{\mathbf{Z} \in \mathcal{Z}} \mathbb{P}_{\boldsymbol{\theta}}(\mathbf{X}, \mathbf{Z}) d\mathbf{Z}$$

Integrations are summation over $\{1, \dots, Q\}$: we have Q^n terms !

Intractable summation

With mixture models, for $\boldsymbol{\theta} = (\boldsymbol{\theta}_1, \dots, \boldsymbol{\theta}_Q)$ we have

$$\log L(\boldsymbol{\theta}; \mathbf{X}) = \sum_{i=1}^n \log \left\{ \sum_{q=1}^Q \alpha_q \mathbb{P}_{\boldsymbol{\theta}_q}(X_i) \right\}.$$

↪ Direct maximization of the likelihood is impossible in practice

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Bayes decision rule / Maximum *a posteriori*

Principle

Affect an individual i to the subpopulation which is the most likely according to the data:

$$\tau_{iq} = \mathbb{P}(Z_{iq} = 1 | X_i = x_i)$$

This is the **posterior probability** for $i \in q$.

Application of the Bayes Theorem

It is straightforward to show that

$$\tau_{iq} = \frac{\alpha_q \mathbb{P}_{\theta_q}(x_i)}{\sum_{q=1}^Q \alpha_q \mathbb{P}_{\theta_q}(x_i)}$$

Principle of the EM algorithm

If θ were known

... estimating the **posterior probability** $\mathbb{P}(Z_i|\mathbf{X})$ of \mathbf{Z} should be easy

By means of the Bayes decision rule

If \mathbf{Z} were known...

... estimating the **best set of parameter** θ should be easy

This is close to usual maximum likelihood estimation

EM principle

Maximize the marginal likelihood iteratively:

- 1 Initialize θ
- 2 Compute the probability of \mathbf{Z} given θ
- 3 Get a better θ with the new \mathbf{Z}
- 4 Iterate until convergence

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Formal algorithm

Initialization: start from a good guess either of \mathbf{Z} or $\boldsymbol{\theta}$, then iterate 1-2

1. Expectation step

Calculate the expected value of the loglikelihood under the current $\boldsymbol{\theta}$

$$Q\left(\boldsymbol{\theta}|\boldsymbol{\theta}^{(t)}\right) = \mathbb{E}_{\mathbf{Z}|\mathbf{X};\boldsymbol{\theta}^{(t)}} \left[\log L(\boldsymbol{\theta}; \mathbf{X}, \mathbf{Z}) \right] \quad (\text{needs } \mathbb{P}_{\boldsymbol{\theta}^{(t)}}(\mathbf{Z}|\mathbf{X}))$$

2. Maximization step

Find the parameters that maximize this quantity

$$\boldsymbol{\theta}^{(t+1)} = \arg \max_{\boldsymbol{\theta}} Q\left(\boldsymbol{\theta}|\boldsymbol{\theta}^{(t)}\right)$$

Stop when $\|\boldsymbol{\theta}^{(t+1)} - \boldsymbol{\theta}^{(t)}\| < \varepsilon$ or $\|Q^{(t+1)} - Q^{(t)}\| < \varepsilon$

(Basic) Convergence analysis

Theorem

At each step of the EM algorithm, the loglikelihood increases. EM thus reaches a local optimum.

Proof.

On board.



Choosing the number of component

Reminder: Bayesian Information Criterion

The BIC is a model selection criterion which penalizes the adjustment to the data by the number of parameter in model \mathcal{M} as follows:

$$\text{BIC}(\mathcal{M}) = \log L(\hat{\boldsymbol{\theta}}; \mathbf{X}) - \frac{1}{2} \log(n) \text{df}(\mathcal{M}).$$

Integrated Classification Criterion

It is an adaptation working with the complete-data likelihood:

$$\begin{aligned} \text{ICL}(\mathcal{M}) &= \log L(\hat{\boldsymbol{\theta}}; \mathbf{X}, \hat{\mathbf{Z}}) + \frac{1}{2} \log(n) \text{df}(\mathcal{M}) \\ &= \text{BIC} - \mathcal{H}(\mathbb{P}(\hat{\mathbf{Z}}|\mathbf{X})), \end{aligned}$$

where the entropy \mathcal{H} measures the separability of the subpopulations.

↪ We choose $\mathcal{M}(Q)$ that maximizes either BIC or ICL

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Mixture of Gaussians

Calculs in the univariate case: complete likelihood

The distribution of X_i conditional on the label of i is assumed to be a univariate Gaussian distribution with unknown parameters:

$$X_i | Z_{iq} = 1 \sim \mathcal{N}(\mu_q, \sigma_q^2)$$

complete Likelihood (\mathbf{X}, \mathbf{Z})

The model complete loglikelihood is

$$\log L(\boldsymbol{\mu}, \boldsymbol{\sigma}^2; \mathbf{X}, \mathbf{Z}) = \sum_{i=1}^n \sum_{q=1}^Q Z_{iq} \left(\log \alpha_q - \log \sigma_q - \log(\sqrt{2\pi}) - \frac{1}{2\sigma_q^2} (x_i - \mu_q)^2 \right)$$

Mixture of Gaussians

Calculs in the univariate case: E-step

E-step

For fixed values of μ_q, σ_q^2 and α_q , the estimates of the posterior probabilities $\hat{\tau}_{iq} = \mathbb{P}(Z_{iq} = 1 | X_i)$ are

$$\hat{\tau}_{iq} = \frac{\alpha_q \mathcal{N}(x_i; \mu_q, \sigma_q^2)}{\sum_{q=1}^Q \alpha_q \mathcal{N}(x_i; \mu_q, \sigma_q^2)},$$

where \mathcal{N} is the density of the normal distribution.

Mixture of Gaussians

Calculs in the univariate case: M-step

M-step

For fixed values of τ_{iq} , the estimates of the model parameters are

$$\hat{\alpha}_q = \frac{\sum_{i=1}^n \tau_{iq}}{\sum_{i=1}^n \sum_{q=1}^Q \tau_{iq}} \quad \hat{\mu}_q = \frac{\sum_i \tau_{iq} x_i}{\sum_i \tau_{iq}} \quad \hat{\sigma}_q^2 = \frac{\sum_{i=1}^n \tau_{iq} (x_i - \mu_q)^2}{\sum_{i=1}^n \tau_{iq}}$$

R code: auxiliary functions

We start by defining functions to compute the complete model loglikelihood, perform the E step and the M step.

```
get.cloglik <- function(X, Z, theta) {
  alpha <- theta$alpha; mu <- theta$mu; sigma <- theta$sigma
  xs <- scale(matrix(X,length(x),length(alpha)),mu,sigma)
  return(sum(Z*(log(alpha)-log(sigma)-.5*(log(2*pi)+xs^2))))
}

M.step <- function(X, tau) {
  n <- length(X); Q <- ncol(tau)
  alpha <- colMeans(tau)
  mu <- colMeans(tau * matrix(X,n,Q)) / alpha
  sigma <- sqrt(colMeans(tau*sweep(matrix(X,n,Q),2,mu,"-")^2)/alpha)
  return(list(alpha=alpha, mu=mu, sigma=sigma))
}

E.step <- function(X, theta) {
  tau <- mapply(function(alpha, mu, sigma) {
    alpha*dnorm(X,mu,sigma)
  }, theta$alpha, theta$mu, theta$sigma)
  return(tau / rowSums(tau))
}
```

R code: EM for univariate mixture

```
EM.mixture <- function(X, Q,
                       init.cl=sample(1:Q,n,rep=TRUE), max.iter=100, eps=1e-5) {
  n <- length(X); tau <- matrix(0,n,Q); tau[cbind(1:n,init.cl)] <- 1
  Eloglik <- vector("numeric", max.iter)
  iter <- 0; cond <- FALSE

  while (!cond) {
    iter <- iter + 1
    ## M step
    theta <- M.step(X, tau)
    ## E step
    tau <- E.step(X, theta)
    ## check consistency
    Eloglik[iter] <- get.cloglik(X, tau, theta)
    if (iter > 1)
      cond <- (iter>=max.iter) | Eloglik[iter]-Eloglik[iter-1] < eps
  }

  return(list(alpha = theta$alpha, mu = theta$mu, sigma = theta$sigma,
             tau = tau, cl = apply(tau, 1, which.max),
             Eloglik = Eloglik[1:iter]))
}
```

Example: data generation

We first generate data with 4 components:

```
mu1 <- 5    ; sigma1 <- 1; n1 <- 100
mu2 <- 10   ; sigma2 <- 1; n2 <- 200
mu3 <- 15   ; sigma3 <- 2; n3 <- 50
mu4 <- 20   ; sigma4 <- 3; n4 <- 100
c1 <- rep(1:4,c(n1,n2,n3,n4))
x <- c(rnorm(n1,mu1,sigma1),rnorm(n2,mu2,sigma2),
      rnorm(n3,mu3,sigma3),rnorm(n4,mu4,sigma4))
n <- length(x)

## we randomize the class ordering
rnd <- sample(1:n)
c1 <- c1[rnd]
x <- x[rnd]

alpha <- c(n1,n2,n3,n4)/n
```

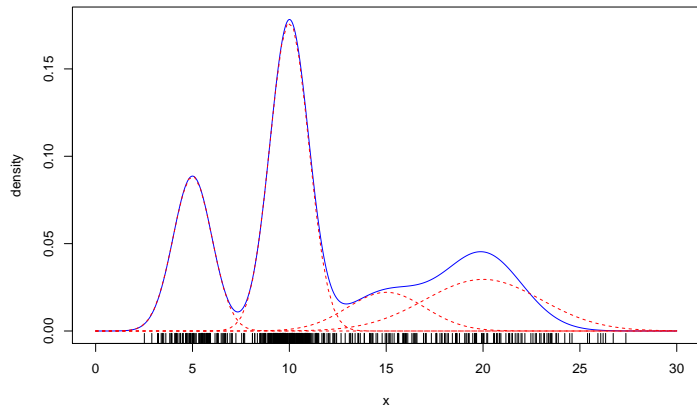
Example: data generation - plot I

Let us plot the data and the theoretical mixture.

```
curve(alpha[1]*dnorm(x,mu1,sigma1) +
       alpha[2]*dnorm(x,mu2,sigma2) +
       alpha[3]*dnorm(x,mu3,sigma3) +
       alpha[4]*dnorm(x,mu4,sigma3),
       col="blue", lty=1, from=0,to=30, n=1000,
       main="Theoretical Gaussian mixture and its components",
       xlab="x", ylab="density")
curve(alpha[1]*dnorm(x,mu1,sigma1), col="red", add=TRUE, lty=2)
curve(alpha[2]*dnorm(x,mu2,sigma2), col="red", add=TRUE, lty=2)
curve(alpha[3]*dnorm(x,mu3,sigma3), col="red", add=TRUE, lty=2)
curve(alpha[4]*dnorm(x,mu4,sigma4), col="red", add=TRUE, lty=2)
rug(x)
```

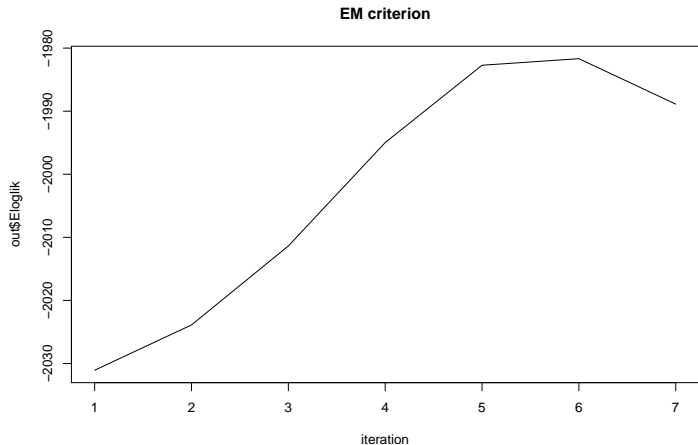
Example: data generation - plot II

Theoretical Gaussian mixture and its components



Example: adjustment

```
out <- EM.mixture(x, Q=4, init.cl=sample(1:4,n,rep=TRUE))  
plot(out$Eloglik, main="EM criterion", type="l", xlab="iteration")
```

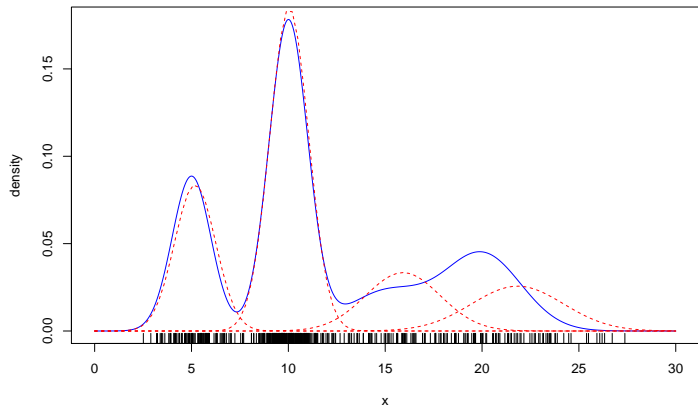


Example: adjustment - plot I

```
out <- EM.mixture(x, Q=4, init.cl=kmeans(x,4)$cl)
curve(alpha[1]*dnorm(x,mu1,sigma1) +
      alpha[2]*dnorm(x,mu2,sigma2) +
      alpha[3]*dnorm(x,mu3,sigma3) +
      alpha[4]*dnorm(x,mu4,sigma3), col="blue",
      lty=1, from=0,to=30, n=1000,
      main="Theoretical Gaussian mixture and estimated components",
      xlab="x", ylab="density")
curve(out$alpha[1]*dnorm(x,out$mu[1],out$sigma[1]), col="red", add=TRUE, lty=2)
curve(out$alpha[2]*dnorm(x,out$mu[2],out$sigma[2]), col="red", add=TRUE, lty=2)
curve(out$alpha[3]*dnorm(x,out$mu[3],out$sigma[3]), col="red", add=TRUE, lty=2)
curve(out$alpha[4]*dnorm(x,out$mu[4],out$sigma[4]), col="red", add=TRUE, lty=2)
rug(x)
```

Example: adjustment - plot II

Theoretical Gaussian mixture and estimated components



Example: adjustment - classification I

```
table(cl, out$cl)
```

```
##  
## cl      1    2    3    4  
##  1      0    2  98    0  
##  2      0 199    1    0  
##  3      1    8    0  41  
##  4     67    1    0  32
```




```
aricode::ARI(cl, out$cl)
```

```
## [1] 0.8449445
```

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References

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Catherine Matias
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A mathematical model: Erdős-Rényi graph

Definition

Let $\mathcal{V} = 1, \dots, n$ be a set of fixed vertices. The (simple) Erdős-Rényi model $\mathcal{G}(n, \pi)$ assumes random edges between pairs of nodes with probability π . In other word, the (random) adjacency matrix \mathbf{X} is such that

$$X_{ij} \sim \mathcal{B}(\pi)$$

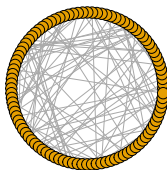
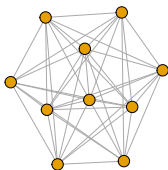
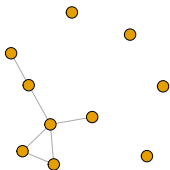
Proposition (degree distribution)

The (random) degree D_i of vertex i follows a binomial distribution:

$$D_i \sim b(n - 1, \pi).$$

Erdős-Rényi - example

```
G1 <- igraph::sample_gnp(10, 0.1)
G2 <- igraph::sample_gnp(10, 0.9)
G3 <- igraph::sample_gnp(100, .02)
par(mfrow=c(1,3))
plot(G1, vertex.label=NA) ; plot(G2, vertex.label=NA)
plot(G3, vertex.label=NA, layout=layout.circle)
```



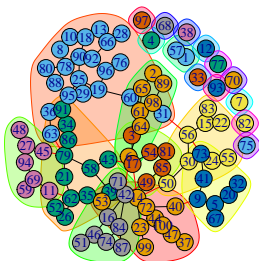
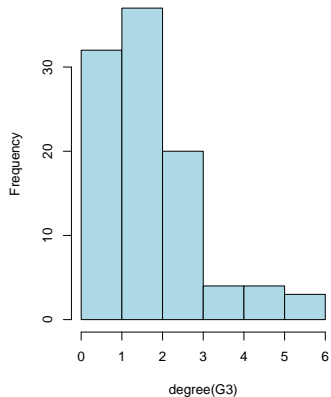
Erdős-Rényi - limitations: very homogeneous

```
average.path.length(G3); diameter(G3)
```

```
## [1] 6.186269
```

```
## [1] 13
```

Histogram of degree(G3)



Mechanism-based model: preferential attachment

The graph is defined dynamically as follows

Definition

Start from a initial graph $\mathcal{G}_0 = (\mathcal{V}_0, \mathcal{E}_0)$, then for each time step,

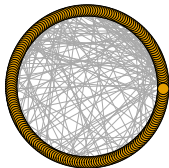
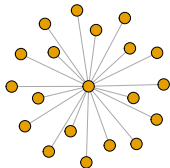
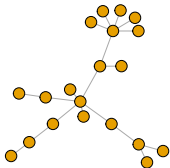
- 1 At t a new node V_t is added
- 2 V_t is connected to $i \in V_{t-1}$ with probability

$$D_i^\alpha + \text{cst.}$$

\rightsquigarrow Nodes with high degree get more connections thus **richers get richers**

Preferential attachment - example

```
G1 <- igraph::sample_pa(20, 1, directed=FALSE)
G2 <- igraph::sample_pa(20, 5, directed=FALSE)
G3 <- igraph::sample_pa(200, directed=FALSE)
par(mfrow=c(1,3))
plot(G1, vertex.label=NA) ; plot(G2, vertex.label=NA)
plot(G3, vertex.label=NA, layout=layout.circle)
```



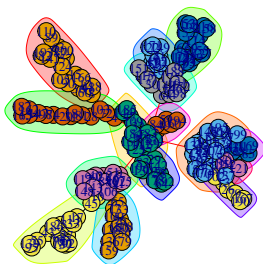
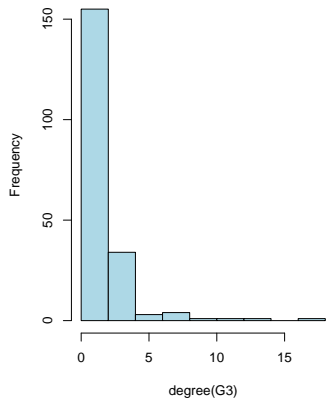
Preferential attachment - limitations

```
average.path.length(G3); diameter(G3)
```

```
## [1] 5.965327
```

```
## [1] 14
```

Histogram of degree(G3)



Limitations

- **Erdős-Rényi**

The ER model does not fit well real world network

- As can be seen from its degree distribution
- ER is generally too homogeneous

- **Preferential attachment**

- Is defined through an algorithm so performing statistics is complicated
- Is stucked to the power-law distribution of degrees

The Stochastic Block Model

The SBM¹ generalizes ER in a mixture framework. It provides

- a statistical framework to adjust and interpret the parameters
- a flexible yet simple specification that fits many existing network data

¹Other models exist (e.g. exponential model for random graphs) but less popular.

Outline

- 1 Background: mixture models and EM
- 2 The Stochastic Block Model (SBM)
 - Some Graphs Models and their limitations
 - Mixture of Erdős-Rényi and the SBM**
 - Inference in SBM with variational EM

Stochastic Block Model: definition

Mixture model point of view: mixture of Erdős-Rényi

Latent structure

Let $\mathcal{V} = \{1, \dots, n\}$ be a fixed set of vertices. We give each $i \in \mathcal{V}$ a **latent label** among a set $\mathcal{Q} = \{1, \dots, Q\}$ such that

- $\alpha_q = \mathbb{P}(i \in q)$, $\sum_q \alpha_q = 1$;
- $Z_{iq} = \mathbf{1}_{\{i \in q\}}$ are independent hidden variables.

The conditional distribution of the edges

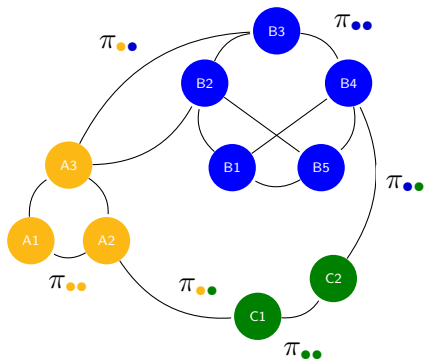
Connexion probabilities depend on the node class belonging:

$$X_{ij} | \{i \in q, j \in \ell\} \sim \mathcal{B}(\pi_{q\ell}) \quad \left(\Leftrightarrow X_{ij} | \{Z_{iq}Z_{j\ell} = 1\} \sim \mathcal{B}(\pi_{q\ell}). \right)$$

The $Q \times Q$ matrix π gives for all couple of labels

$$\pi_{q\ell} = \mathbb{P}(X_{ij} = 1 | i \in q, j \in \ell).$$

Stochastic Block Model: the big picture



Stochastic Block Model

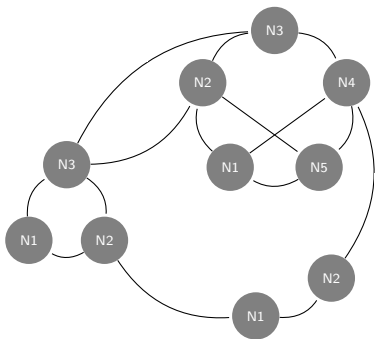
Let n nodes divided into

- $\mathcal{Q} = \{\bullet, \bullet, \bullet\}$ classes
- $\alpha_{\bullet} = \mathbb{P}(i \in \bullet), \bullet \in \mathcal{Q}, i = 1, \dots, n$
- $\pi_{\bullet\bullet} = \mathbb{P}(i \leftrightarrow j | i \in \bullet, j \in \bullet)$

$$Z_i = \mathbf{1}_{\{i \in \bullet\}} \sim^{\text{iid}} \mathcal{M}(1, \alpha), \quad \forall \bullet \in \mathcal{Q},$$

$$X_{ij} | \{i \in \bullet, j \in \bullet\} \sim^{\text{ind}} \mathcal{B}(\pi_{\bullet\bullet})$$

Stochastic Block Model: unknown parameters



Stochastic Block Model

Let n nodes divided into

- $\mathcal{Q} = \{\bullet, \bullet, \bullet\}$, $\text{card}(\mathcal{Q})$ known
- $\alpha_{\bullet} = ?$,
- $\pi_{\bullet\bullet} = ?$

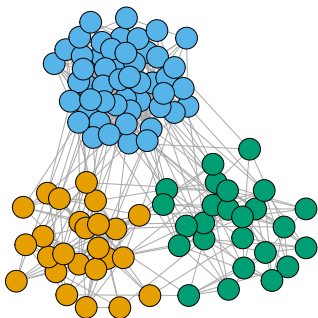
$$Z_i = \mathbf{1}_{\{i \in \bullet\}} \sim^{\text{iid}} \mathcal{M}(1, \alpha), \quad \forall \bullet \in \mathcal{Q},$$

$$X_{ij} \mid \{i \in \bullet, j \in \bullet\} \sim^{\text{ind}} \mathcal{B}(\pi_{\bullet\bullet})$$

Stochastic block models – examples of topology

Community network

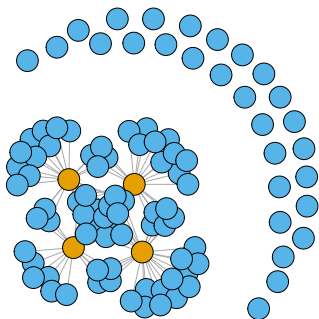
```
pi <- matrix(c(0.3,0.02,0.02,0.02,0.3,0.02,0.02,0.02,0.3),3,3)
communities <- igraph::sample_sbm(100, pi, c(25, 50, 25))
plot(communities, vertex.label=NA, vertex.color = rep(1:3,c(25, 50, 25)))
```



Stochastic block models – examples of topology

Star network

```
pi <- matrix(c(0.05,0.3,0.3,0),2,2)
star <- igraph::sample_sbm(100, pi, c(4, 96))
plot(star, vertex.label=NA, vertex.color = rep(1:2,c(4,96)))
```



Degree distributions

Conditional degree distribution

The conditional degree distribution of a node $i \in q$ is

$$D_i | i \in q \sim \text{b}(n-1, \bar{\pi}) \approx \mathcal{P}(\lambda_q), \quad \bar{\pi}_q = \sum_{\ell=1}^Q \alpha_\ell \pi_{q\ell} \quad \lambda_q = (n-1)\bar{\pi}_q$$

Conditional degree distribution

The degree distribution of a node i can be approximated by a mixture of Poisson distributions:

$$\mathbb{P}(D_i = k) = \sum_{q=1}^Q \alpha_q \exp\{-\lambda_q\} \frac{\lambda_q^k}{k!}$$

Likelihoods

Complete-data loglikelihood

$$\log L(\mathbf{X}, \mathbf{Z}) = \sum_{i,q} Z_{iq} \log \alpha_q + \sum_{i < j, q, \ell} Z_{iq} Z_{j\ell} \log \pi_{q\ell}^{X_{ij}} (1 - \pi_{q\ell})^{1 - X_{ij}}.$$

Conditional expectation of the complete-data loglikelihood

$$\mathbb{E}_{\mathbf{Z}|\mathbf{X}}[\log L(\boldsymbol{\theta}; \mathbf{X}, \mathbf{Z})] = \sum_{i,q} \tau_{iq} \log \alpha_q + \sum_{i < j, q, \ell} \eta_{ijq\ell} \log \pi_{q\ell}^{X_{ij}} (1 - \pi_{q\ell})^{1 - X_{ij}}$$

where $\tau_{iq}, \eta_{ijq\ell}$ are the posterior probabilities:

- $\tau_{iq} = \mathbb{P}(Z_{iq} = 1 | \mathbf{X}) = \mathbb{E}[Z_{iq} | \mathbf{X}]$.
- $\eta_{ijq\ell} = \mathbb{P}(Z_{iq} Z_{j\ell} = 1 | \mathbf{X}) = \mathbb{E}[Z_{iq} Z_{j\ell} | \mathbf{X}]$.

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The EM strategy does not apply directly for SBM

Ouch: another intractability problem

- the Z_{iq} are **not independent** in the SBM framework. . .
- we cannot compute $\eta_{ijql} = \mathbb{P}(Z_{iq}Z_{j\ell} = 1|\mathbf{X}) = \mathbb{E}[Z_{iq}Z_{j\ell}|\mathbf{X}]$,
- the conditional expectation $Q(\boldsymbol{\theta})$, i.e. the main EM ingredient, is **intractable**.

Solution: mean field approximation

Approximate η_{ijql} by $\tau_{iq}\tau_{j\ell}$, i.e., **assume independence between Z_{iq}**
 \rightsquigarrow This can be formalized in the variational framework

Revisiting the EM algorithm I

Proposition

Consider a distribution Q for the $\{Z_{iq}\}$. We have

$$\log L(\boldsymbol{\theta}; \mathbf{X}) = \mathbb{E}_Q[\log L(\boldsymbol{\theta}, \mathbf{X}, \mathbf{Z})] + \mathcal{H}(Q) + \text{KL}(Q \mid \mathbb{P}(\mathbf{Z}|\mathbf{X}; \boldsymbol{\theta})),$$

where \mathcal{H} is the entropy and $\text{KL}(\cdot|\cdot)$ is the Kullback-Leibler divergence:

$$\mathcal{H}(Q) = - \sum_z Q(z) \log Q(z) = -\mathbb{E}_Q[\log Q(Z)]$$

$$\text{KL}(Q \mid \mathbb{P}(\mathbf{Z}|\mathbf{X}; \boldsymbol{\theta})) = \sum_z Q(z) \log \frac{Q(z)}{\mathbb{P}(\mathbf{Z}|\mathbf{X}; \boldsymbol{\theta})} = \mathbb{E}_Q \left[\log \frac{Q(z)}{\mathbb{P}(\mathbf{Z}|\mathbf{X}; \boldsymbol{\theta})} \right]$$

Revisiting the EM algorithm II

Let

$$J(\mathbb{Q}, \boldsymbol{\theta}) \triangleq \mathbb{E}_{\mathbb{Q}} (\log L(\boldsymbol{\theta}; \mathbf{X}, \mathbf{Z})) + \mathcal{H}(\mathbb{Q})$$

The steps in the EM algorithm may be viewed as:

Expectation step : choose \mathbb{Q} to maximize $J(\mathbb{Q}; \boldsymbol{\theta}^{(t)})$

The solution is $\mathbb{P}(\mathbf{Z}|\mathbf{X}; \boldsymbol{\theta}^{(t)})$

Maximization step : choose $\boldsymbol{\theta}$ to maximize $J(\mathbb{Q}^{(t)}; \boldsymbol{\theta})$

The solution maximizes $\mathbb{E}_{\mathbf{Z}|\mathbf{X}; \boldsymbol{\theta}^{(t)}} (\log L(\boldsymbol{\theta}; \mathbf{X}, \mathbf{Z}))$

Variational approximation for SBM

Problem for SBM

$\mathbb{P}(\mathbf{Z}|\mathbf{X}; \boldsymbol{\theta}^{(t)})$ cannot be computed thus the E-step cannot be solved.

Idea

Choose \mathbb{Q} in a class of function so that the E-step can be solved.

Family of distribution that factorizes

We chose \mathbb{Q} so as the Z_{iq} are marginally independents:

$$\mathbb{Q}(\mathbf{Z}) = \prod_{i=1}^n \mathbb{Q}_i(Z_i) = \prod_{i=1}^n \prod_{q=1}^Q \tau_{iq}^{Z_{iq}},$$

where $\tau_{iq} = \mathbb{Q}_i(Z_i = q) = \mathbb{E}Q(Z_{iq})$, with $\sum_q \tau_{iq} = 1$ for all $i = 1, \dots, n$.

Variational EM for SBM: the criterion

Lower bound of the loglikelihood

Since \mathbb{Q} is an approximation of $\mathbb{P}(\mathbf{Z}|\mathbf{X})$, the Kullback-Leibler divergence is non-negative and

$$\log L(\boldsymbol{\theta}; \mathbf{X}) \geq \mathbb{E}_{\mathbb{Q}}[\log L(\boldsymbol{\theta}, \mathbf{X}, \mathbf{Z})] + \mathcal{H}(\mathbb{Q}) = J(\mathbb{Q}, \boldsymbol{\theta}).$$

For the SBM,

$$J(\mathbb{Q}, \boldsymbol{\theta}) = \sum_{i,q} \tau_{iq} \log \alpha_q + \sum_{i < j, q, \ell} \tau_{iq} \tau_{j\ell} \log b(X_{ij}; \pi_{q\ell}) - \sum_{i,q} \tau_{iq} \log(\tau_{iq}),$$

\rightsquigarrow we optimize the loglikelihood lower bound $J(\mathbb{Q}, \boldsymbol{\theta}) = J(\boldsymbol{\tau}, \boldsymbol{\theta})$ in $(\boldsymbol{\tau}, \boldsymbol{\theta})$.

E and M steps for SBM

Variational E-step

Maximizing $J(\boldsymbol{\tau})$ for fixed $\boldsymbol{\theta}$, we find a fixed-point relationship:

$$\hat{\tau}_{iq} \propto \alpha_q \prod_j \prod_\ell b(X_{ij}, \pi_{q\ell})^{\hat{\tau}_{j\ell}} \quad (1)$$

M-step

Maximizing $J(\boldsymbol{\theta})$ for fixed $\boldsymbol{\tau}$, we find,

$$\hat{\alpha}_q = \frac{1}{n} \sum_i \hat{\tau}_{iq}, \quad \hat{\pi}_{q\ell} = \frac{\sum_{i \neq j} \hat{\tau}_{iq} \hat{\tau}_{j\ell} X_{ij}}{\sum_{i \neq j} \hat{\tau}_{iq} \hat{\tau}_{j\ell}}. \quad (2)$$

Model selection

We use our lower bound of the loglikelihood to compute an approximation of the ICL

$$\begin{aligned} \text{vICL}(Q) = \mathbb{E}_{\hat{Q}}[\log L(\hat{\boldsymbol{\theta}}); \mathbf{X}, \mathbf{Z}] \\ - \frac{1}{2} \left(\frac{Q(Q+1)}{2} \log \frac{n(n-1)}{2} + (Q-1) \log(n) \right), \end{aligned}$$

where

$$\mathbb{E}_{\hat{Q}}[\log L(\hat{\boldsymbol{\theta}}; \mathbf{X}, \mathbf{Z})] = J(\hat{\boldsymbol{\tau}}, \hat{\boldsymbol{\theta}}) - \mathcal{H}(\hat{Q}).$$

The variational BIC is just

$$\text{vBIC}(Q) = J(\hat{\boldsymbol{\tau}}, \hat{\boldsymbol{\theta}}) - \frac{1}{2} \left(\frac{Q(Q+1)}{2} \log \frac{n(n-1)}{2} + (Q-1) \log(n) \right).$$

Example on the French blogosphere I

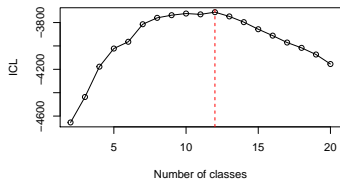
```
library(mixer)
data(blog)
mix.blog <- mixer(x=blog$links,qmin=2,qmax=20)

## Mixer: the adjacency matrix has been transformed in a undirected edge list

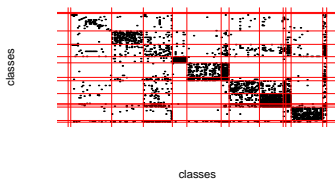
plot(mix.blog)
```

Example on the French blogosphere II

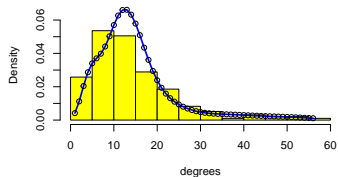
Integrated Classification Likelihood



Reorganized Adjacency matrix



Degree distribution



Inter/intra class probabilities

