# **Statistical Learning with Interaction Data**

Applications to pharmacovigilance and network analysis

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- it allows to model complex phenomena and provides synthetic summaries,
- it provides theoretical guarantees about its performance and behaviour,
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Despite recent impressive results in supervised learning, many situations remain challenging:

- high-dimensional (p large),
- big or as stream (*n* large),
- evolutive (evolving phenomenon),
- heterogeneous (categorical, functional, network, interaction data, ...)

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#### The understanding of the results is essential:

- in many applications, practitioners are very interested in visualizing the processed data,
- and to get a synthetic summary of the data for better interpretation.

### A research team in "core Al", created in 2020:



- 6 permanent researchers, 25 Ph.D. students and postdocs, and 4 engineers,
- located at the Centre Inria of Université Côte d'Azur, in Sophia-Antipolis,

The team focuses on the Models and Algorithms of Artificial Intelligence:



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A summary of our topics:



Figure 1: Scientific objectives of Maasai.

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# Dealing with interaction data

Interaction data become ubiquitous for the analysis of:

 recommendation systems, pharmacovigilance, ...: bipartite networks / incidence matrices.



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Interaction data become ubiquitous for the analysis of:

 recommendation systems, pharmacovigilance, ...: bipartite networks / incidence matrices.  social networks, co-authorship networks, communication networks, ...: networks / adjacency matrices.





Clustering such interaction data is a recurrent task:

- clustering the nodes of a network  $\rightarrow$  detection of influencers, detection of weak signals, ...
- co-clustering of ordinal data  $\rightarrow$  recommendation systems, ...
- co-clustering of count data  $\rightarrow$  bike sharing systems, traffic modeling, ...

With possibilities to apply it to the field of public health:

- detecting safety signals in adverse drug reaction data (pharmacovigilance),
- model and predict the use of health care services of a hospital.
- understand the role of social networks on adverse drug reaction declaration (pharmacovigilance),
- summarizing medical publication networks to help fighting against a pandemic.

**Co-clustering of interaction data streams for Pharmacovigilance** 

#### Pharmacovigilance:

- it is the study of adverse reactions to drugs and vaccines,
- it aims at detecting safety signals about drugs,
- this task is done manually nowadays,
- it can be complicated in case of important media coverage.

The Nice RCPV data:



Figure 4: Evolution of spontaneous reports (extract) to RCPV from 2010 to 2020.

The data we consider are organized as follows:

- rows are indexed by  $i = 1, \dots N$ ;
- columns are indexed by  $j = 1, \ldots, M$ ;
- time instants  $t \in [0, T]$  during which N and M are fixed;
- the  $N \times M \times T$  tensor  $X := \{X_{ij}(t)\}$  contains the number of interactions between any observation and feature pair at any given t.



Figure 5: Data structure.

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#### We aim at estimating:

- The latent variables for the clustering of rows and columns into Q and L groups,
- A latent variable for modeling the evolving sparsity of the data.



Figure 5: Data structure.

#### 1) Modeling of the row and column clusters:

• At each time instant *t*, the *i*th row of *X*(*t*) is assigned to an (unobserved) group among Q(t), according to:

$$Z_i(t) \sim \mathcal{M}(1, \alpha(t) := (\alpha_1(t), \ldots, \alpha_Q(t))),$$

where  $\alpha_q(t) \ge 0$  and  $\sum_{q=1}^{Q} \alpha_q(t) = 1$ , for all t = 0, ..., T.

• similarly, the *j*th column of X(t) is assigned to an (unobserved) group among L(t), according to:

$$W_j(t) \sim \mathcal{M}(1, \beta(t) := (\beta_1(t), \ldots, \beta_L(t))),$$

where  $\beta_{\ell}(t) \ge 0$  and  $\sum_{\ell=1}^{L} \beta_{\ell}(t) = 1$ , for all t = 0, ..., T.

• row assignments  $z_i(t)$  are further assumed to be independent from column assignments  $w_j(t)$ , for all i, j;

#### 2) Modeling of a potential extreme sparsity:

• the observed variable X(t) is assumed to be modeled by a mixture of block-conditional Zero-Inflated (ZI) distributions:

$$X_{ij}(t)|Z_i(t) = k, W_j(t) = \ell ~~\sim~~ ZI(\zeta_{k,\ell},\pi(t)),$$

where:

- $\zeta$  is the block-dependent vector of parameters for the distribution  $\phi(X_{ij}(t), .)$ ,
- $\pi(t)$  is the sparsity probability at any given time period t.

The zero-inflated distribution is therefore such that:

$$\begin{cases} X_{ij}(t)|Z_i(t), W_j(t) \sim 0 & \text{with probability } \pi(t) \\ X_{ij}(t)|Z_i(t), W_j(t) \sim \phi(X_{ij}(t); \zeta_{Z_i(t), W_j(t)}) & \text{with probability } 1 - \pi(t) \end{cases}$$
(1)

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(1)

The modeling of the data sparsity can be re-formulated by introducing a latent variable A(t):

- such that  $A_{ij}(t) \sim \mathcal{B}(\pi(t))$ ,
- and we therefore get:

$$\begin{cases} X_{ij}(t)|Z_i(t), W_j(t) \sim 0 & \text{if } A_{ij}(t) = 1 \\ X_{ij}(t)|Z_i(t), W_j(t) \sim \mathcal{P}(\Lambda_{Z_i(t), W_j(t)}) & \text{if } A_{ij}(t) = 0 \end{cases}$$
(2)

3) Modeling of the dynamic of sparsity and cluster proportions:

The evolving mixing proportion and the sparsity parameter are assumed to be generated by three systems of ODEs, respectively:

•  $\frac{d}{dt}a(t) = f_Z(a(t)),$  with  $\alpha_q(t) = \frac{e^{a_q(t)}}{\sum_{q=1}^Q e_q^a(t)},$ •  $\frac{d}{dt}b(t) = f_W(b(t)),$  with  $\beta_\ell(t) = \frac{e^{b_\ell(t)}}{\sum_{\ell=1}^L e_\ell^b(t)},$ •  $\frac{d}{dt}c(t) = f_A(c(t)),$  with  $\pi(t) = \frac{e^{c(t)}}{e^{c(t)} + e^{(1-c(t))}},$ 

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Since we work with discrete time points, the dynamic systems reduce to their Euler schemes:

- $a(t+1) = a(t) + f_Z(a(t)),$
- $b(t+1) = b(t) + f_W(a(t)),$
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We further assume that the functions  $f_Z$ ,  $f_W$  and  $f_A$  can be modeled by three fully connected neural networks.



Figure 6: Graphical representation of the ZIP-dLBM model.

# The joint distribution (Poisson case)

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Given  $\theta = (\Lambda, \alpha(t), \beta(t), \pi(t))$ , we can compute the likelihood of the complete data:

$$p(X, Z, W, A|\theta) = p(X|Z, W, A, \Lambda, \pi)p(A \mid \pi)p(Z|\alpha)p(W|\beta)$$
(3)

where:

$$p(X|A, Z, W, \Lambda, \pi) = \prod_{i=1}^{N} \prod_{j=1}^{M} \prod_{t=1}^{T} \mathbf{1}_{\{X_{ij}(t)=0\}}^{A_{ij}(t)} \left\{ \left( \frac{\Lambda_{Z_{i}(t)W_{j}(t)}^{X_{ij}(t)}}{X_{ij}(t)!} \exp(-\Lambda_{Z_{i}(t)W_{j}(t)}) \right)^{(1-A_{ij}(t))} \right\},$$
(4)

$$p(A|\pi) = \prod_{i=1}^{N} \prod_{j=1}^{M} \prod_{t=1}^{T} \pi(t)^{A_{ij}(t)} (1 - \pi(t)) \Big)^{(1 - A_{ij}(t))},$$
(5)

$$p(Z|\alpha) = \prod_{i=1}^{N} \prod_{q=1}^{Q} \prod_{t=1}^{T} \alpha_q(t)^{Z_{iq}(t)},$$
(6)

$$\rho(W|\beta) = \prod_{j=1}^{M} \prod_{\ell=1}^{L} \prod_{t=1}^{T} \beta_{\ell}(t)^{W_{j\ell}(t)}.$$
(7)

# The inference: variational assumptions assumptions

### We rely on the Variational-EM algorithm (VEM) for infering the model:

Given a variational distribution  $q(\cdot)$ :

$$\log p(X|\theta) = \mathcal{L}(q;\theta) + KL(q(.)||p(.|X,\theta)),$$

where:

$$\mathcal{L}(q,\theta) = \sum_{Z} \sum_{W} \sum_{A} q(Z, W, A) \log \frac{p(X, A, Z, W|\theta)}{q(Z, W, A)},$$
$$\mathcal{KL}(q(.)||p(.|X, \theta)) = -\sum_{Z} \sum_{W} \sum_{A} q(Z, W, A) \log \frac{p(Z, W, A|X, \theta)}{q(Z, W, A)}$$

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In order to optimize  $\mathcal{L}(q, \theta)$ , we further assume that q(A, Z, W) can factorize:

$$\begin{aligned} q(Z,W,A) &= q(A)q(Z)q(W) = \prod_{i=1}^{N} \prod_{j=1}^{M} \prod_{t=1}^{T} q(A_{ij}(t)) \prod_{i=1}^{N} \prod_{t=1}^{T} q(Z_{i}(t)) \prod_{j=1}^{M} \prod_{t=1}^{T} q(W_{j}(t)) \\ &= \prod_{i=1}^{N} \prod_{j=1}^{M} \prod_{t=1}^{T} \delta_{ij}(t)^{A_{ij}(t)} (1 - \delta_{ij}(t))^{1 - A_{ij}(t)} \prod_{i=1}^{N} \prod_{q=1}^{Q} \prod_{t=1}^{T} \tau_{iq}(t)^{Z_{iq}(t)} \prod_{j=1}^{M} \prod_{t=1}^{T} \eta_{j\ell}(t)^{W_{j\ell}(t)}. \end{aligned}$$

 $\mathcal{L}(q, \theta)$  can be finally expressed as:

$$\begin{aligned} \mathcal{L}(q,\theta) &= \sum_{t=1}^{T} \sum_{i=1}^{N} \sum_{j=1}^{M} \left\{ \delta_{ij}(t) \log(\pi(t) \mathbf{1}_{\{X_{ij}(t)=0\}}) + (1-\delta_{ij}(t)) \left[ \log(1-\pi(t)) + \right. \\ &+ \sum_{q=1}^{Q} \sum_{\ell=1}^{L} \left\{ \tau_{iq}(t) \eta_{j\ell}(t) X_{ij}(t) \log \Lambda_{q\ell} - \tau_{iq}(t) \eta_{j\ell}(t) \Lambda_{q\ell} \right\} \right] - (1-\delta_{ij}(t)) \log(X_{ij}(t)!) \right\} + \\ &+ \sum_{q=1}^{T} \sum_{i=1}^{N} \sum_{q=1}^{Q} \tau_{iq}(t) \log(\alpha_{q}(t)) + \sum_{t=1}^{T} \sum_{j=1}^{M} \sum_{\ell=1}^{L} \eta_{j\ell}(t) \log(\beta_{\ell}(t)) - \sum_{t=1}^{T} \sum_{i=1}^{N} \sum_{q=1}^{Q} \tau_{iq}(t) \log \tau_{iq}(t) + \\ &- \sum_{t=1}^{T} \sum_{j=1}^{M} \sum_{\ell=1}^{L} \eta_{j\ell}(t) \log(\eta_{j\ell}(t)) - \sum_{t=1}^{T} \sum_{i=1}^{N} \sum_{j=1}^{M} \left( \delta_{ij}(t) \log(\delta_{ij}(t)) + (1-\delta_{ij}(t)) \log(1-\delta_{ij}(t)) \right) \end{aligned}$$

### The inference: VEM Algorithm

- VE-Step: Lower bound maximization with respect to q(A, Z, W). The optimal sequential updates of the variational distributions are computed through:
  - $\log q^*(A) = E_{W,Z}[\log p(X, A, Z, W \mid \theta)]$
  - $\log q^*(Z) = E_{W,A}[\log p(X, A, Z, W \mid \theta)]$
  - $\log q^*(W) = E_{A,Z}[\log p(X, A, Z, W \mid \theta)]$
- M-Step: Lower bound maximization with respect to  $\theta = (\alpha(t), \beta(t), \pi(t), \Lambda)$ .
  - The derived optimal update of  $\Lambda$  is:

$$\hat{\Lambda}_{q\ell} = rac{\displaystyle \sum_{i=1}^{N} \sum_{j=1}^{M} \sum_{t=1}^{T} \left\{ au_{iq}(t) \eta_{j\ell}(t) \Big( X_{ij}(t) - \delta_{ij}(t) X_{ij}(t) \Big) 
ight\}}{\displaystyle \sum_{i=1}^{N} \sum_{j=1}^{M} \sum_{t=1}^{T} \left\{ au_{iq}(t) \eta_{j\ell}(t) \Big( 1 - \delta_{ij}(t) \Big) 
ight\}}$$

 The optimal updates of α(t), β(t) and π(t) are obtained through a stochastic gradient descent optimization process. 
$$\begin{split} & \overline{\text{Algorithm 1 VEM-SGD Algorithm (for the Zero-Inflated Poisson distribution)} \\ & \overline{\text{Require: } X, Q, L, max.iter, o(l), \beta(l), \tau(l), \Lambda \text{ from Initialization.} \\ & \text{Initialization of } \tau(l) \text{ and } \eta(l): \text{ sampling from } \mathcal{M}(\alpha(l)) \text{ and } \mathcal{M}(\beta(l)), \text{ respectively}; \\ & \text{Initialization of } \tau(l) \text{ matrix of } 1, \text{ then setting } \delta(l) = 0 \text{ when } X > 0; \\ & \text{for it = 1 to max.iter do} \\ & \overline{\text{VE-Step:}} \\ & \text{for } \text{it = 1 to max.iter do} \\ & \overline{\text{VE-Step:}} \\ & \text{for } p = 1 \text{ to Fixed.Point do} \\ & \overline{\text{Update } \delta(l), \tau(l), \eta(l):} \\ & \delta_{\eta}(e) = \frac{\exp(\theta_{\eta}(l))}{(1 + \exp(\theta_{\eta}(q)))}, \\ & \text{where} \\ & R_{ij}(e) = \log(\pi(i) \mathbb{I}_{X_{ij}(e)=0}) + \frac{Q}{2\pi} \sum_{i=1}^{L} \left[ -\pi_{\eta}(l) \eta_{il}(l) \mathbb{I}_{N_{il}(i)} \log \Lambda_{d^{i}} + \tau_{\eta}(l) \eta_{il}(l) \mathbb{I}_{N_{il}(l)} - \log(1 - \pi(l)), \\ & \tau_{\eta}(i) = \frac{L}{D_{i}} \exp\left(\sum_{j=1}^{L} \sum_{i=1}^{L} \left\{ (1 - \delta_{ij}(e)) \left[ \eta_{il}(l) \mathbb{I}_{N_{il}(i)} \log (\Lambda_{q^{i}}) - \eta_{il}(l) \mathbb{A}_{q^{i}} \right] \right\} + \log(\alpha_{\eta}(l)) \right). \\ & \eta_{il}(e) = \sum_{i=1}^{L} \sum_{i=1}^{Q} \left\{ \left[ (1 - \delta_{ij}(e)) \left[ \pi_{il}(l) \mathbb{I}_{N_{il}(i)} \log (\Lambda_{q^{i}}) - \pi_{il}(l) \mathbb{A}_{q^{i}} \right] \right\} + \log(\alpha_{\eta}(t)) \right\}. \\ & \text{ with } D_{q} \text{ and } D_{t} \text{ tormalizing constants.} \end{split}$$

end for

M-Step:

 $\begin{array}{l} \label{eq:constraints} \mathbf{Update} \ \theta = (\lambda, \pi(t), \alpha(t), \beta(t))), \\ \lambda_{\theta'} = & \sum_{i \in \mathcal{A}} \frac{(\lambda_{ij}, \alpha_{ij}, \alpha_$ 

# Introductory example

True alpha







Estimated alpha



Estimated beta



Estimated pi





**Figure 8:** Reorganized incidence matrices at time instants t =10,30, according to the estimates  $\hat{Z}$  and  $\hat{W}$ .



Component activation: beta



time(t)

## Model selection experiment

- 50 simulated dataset;
- The maximum of the given Q and L is 10;
- ZIP-dLBM succeeds 86% of the time to identify the correct model (Q = 3, L = 2).

Q/L	1	2	3	4	5	6	7	8	9	10
1	0	0	0	0	0	0	0	0	0	0
2	0	0	0	0	0	0	0	0	0	0
3	0	86	0	0	0	0	0	0	0	0
4	0	2	0	0	0	0	0	0	0	0
5	0	2	0	0	0	0	0	0	0	0
6	0	0	0	0	0	0	0	0	0	0
7	0	0	0	0	0	0	0	0	0	0
8	0	4	0	0	0	0	0	0	0	0
9	0	2	0	0	0	0	0	0	0	0
10	0	4	0	0	0	0	0	0	0	0

**Table 1:** Model selection. Percentage of activated components on 50 simulated datasets. Thehighlighted cell corresponds to the actual value of Q and L.

We consider adverse drug reaction (ADR) data collected by the Regional Center of Pharmacovigilance (RCPV), located in the University Hospital of Nice:

- 2.3 million inhabitants;
- several channels (e.g. website form, email, etc);
- time horizon of 7 years (trimester as unity measure);
- 27 754 notifications in the dataset;
- only drugs and ADRs notified more than 20 times are considered;
- the resulting dataset contains 236 drugs, 324 ADRs and 29 trimesters.

Histogram of complete data (2015-2022)



**Figure 9:** Frequency of declarations received by the pharmacovigilance center from January 2015 to March 2022, sorted by month.



Figure 10: Estimated Poisson intensities, each color represents a different drug (ADR) cluster.



**Figure 11:** Evolution of the estimates  $\hat{\pi}$ .



**Figure 12:** Evolution of the estimates  $\hat{\alpha}$  and  $\hat{\beta}$ .

The deep latent position model for network clustering

Networks are specific interaction data where receivers are the same individuals as the senders:

- Nodes (or vertices) represent an object or entity such as a person, an email, etc.
- Links (or edges) represent connections between nodes such as "follow", "send to", etc.



### Statistical models:

- Based on probabilistic generation:
  - SBM (Nowicki et al., 2001),
  - OSBM (Latouche et al., 2011),
  - STBM (Bouveyron et al., 2018), etc.
- Based on latent positions:
  - LPM (Hoff et al., 2002),
  - LPCM (Handcock et al., 2007), etc.
- $\implies$  limitations:
  - challenging inference procedure
  - scaling difficulties

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#### Deep learning approaches:

- Based on VAE architecture:
  - VGAE (Kipf et al., 2016),
  - ARVGA (Pan et al. 2018),
  - DGLFRM (Mehta et al., 2019), etc.

### $\implies$ limitations:

- rely on an external algorithm (e.g. k-means) for clustering;
- without taking into account edge features;
- use an simple inner product as decoder.

### Deep latent position model (deepLPM)

Extending the idea of graph VAE, we propose the deepLPM model here:



Figure 13: A deep-learning-like model view of deepLPM.

- take into account edge features
- use a more general latent position-based decoder
- build an end-to-end clustering model





First, each node is assigned to a cluster via a random variable  $c_i$  encoding its cluster membership:

$$c_i \stackrel{iid}{\sim} \mathcal{M}(1,\pi), \quad \text{with} \quad \pi \in [0,1]^K, \ \sum_{k=1}^K \pi_k = 1.$$
 (8)



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Then, conditionally to its cluster membership, a latent embedding vector  $z_i$  is generated:

$$z_i|(c_{ik}=1) \sim \mathcal{N}(\mu_k, \sigma_k^2 I_P), \quad \text{with} \quad \sigma_k^2 \in \mathbb{R}^{+*}.$$
(9)



Finally, the probability of a connection between nodes i and j is modeled by

$$A_{ij}|z_i, z_j \sim \mathcal{B}(f_{\alpha,\beta}(z_i, z_j)), \tag{10}$$

with

$$f_{\alpha,\beta}(z_i, z_j) = \sigma(\alpha + \beta^T y_{ij} - ||z_i - z_j||^2).$$
(11)

where  $f_{\alpha,\beta}$  is a decoding neural network parametrized by  $\alpha$  and  $\beta$ . Moreover,  $\sigma$  is the logistic sigmoid function and  $y_{ij}$  is the covariate of the edge connecting *i* with *j*.

Denoting by  $\Theta = \{\pi, \mu_k, \sigma_k^2, \alpha, \beta\}$ , we want to maximize the integrated log-likelihood:

$$\log p(A|\Theta) = \log \int_{Z} \sum_{C} p(A, Z, C|\Theta) dZ, \qquad (12)$$

with respect to  $\Theta$ .

Since Equation (12) is not tractable, we use a variational approach to approximate it

$$\log p(A|\Theta) = \underbrace{\mathcal{L}(q(Z,C);\Theta)}_{ELBO} + D_{KL}(q(Z,C)||p(Z,C|A,\Theta)).$$
(13)

where  $D_{KL}$  denotes the Kullback-Leibler divergence.

### Model inference: variational assumptions



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Assumption 1.

$$q(Z) = \prod_{i=1}^{N} q(z_i) = \prod_{i=1}^{N} \mathcal{N}(z_i; \tilde{\mu}_{\phi}(\overline{A})_i, \tilde{\sigma}_{\phi}^{-2}(\overline{A})_i I_P),$$
(14)

where  $[\tilde{\mu}_{\phi}(\overline{A}), \log \tilde{\sigma}_{\phi}^{2}(\overline{A})] = g_{\phi}(\overline{A}), g_{\phi}$  is a two-layer GCN encoder parametrized by  $\phi$ .  $\overline{A} = D^{-\frac{1}{2}}AD^{-\frac{1}{2}}$  kipf2016semi is the normalized adjacency matrix.

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$$q(Z) = \prod_{i=1}^{N} q(z_i) = \prod_{i=1}^{N} \mathcal{N}(z_i; \tilde{\mu}_{\phi}(\overline{A})_i, \tilde{\sigma}_{\phi}^{-2}(\overline{A})_i I_P),$$
(14)

where  $[\tilde{\mu}_{\phi}(\overline{A}), \log \tilde{\sigma}_{\phi}^{2}(\overline{A})] = g_{\phi}(\overline{A}), g_{\phi}$  is a two-layer GCN encoder parametrized by  $\phi$ .  $\overline{A} = D^{-\frac{1}{2}}AD^{-\frac{1}{2}}$  kipf2016semi is the normalized adjacency matrix.

Assumption 2.

$$q(C) = \prod_{i=1}^{N} \mathcal{M}(c_i; 1, \gamma_i), \quad \text{with} \quad \sum_{k=1}^{K} \gamma_{ik} = 1,$$
(15)

where  $\gamma_{ik}$  represents the variational probability that node *i* is in cluster *k*.

# Model inference: a joint optimization

### With the above assumptions, $\mathcal{L}(\mathsf{ELBO})$ can be further developed as

$$\mathcal{L} = \underbrace{\left[\sum_{i \neq j} A_{ij} \log \eta_{ij} + (1 - A_{ij}) \log(1 - \eta_{ij})\right]}_{implicit optimization} - \underbrace{\sum_{i=1}^{N} \sum_{k=1}^{K} \gamma_{ik} D_{KL}(\mathcal{N}(\tilde{\mu}_{\phi}(\overline{A})_{i}, \tilde{\sigma}_{\phi}^{2}(\overline{A})_{i}l_{P}) ||\mathcal{N}(\mu_{k}, \sigma_{k}^{2}l_{P}))}_{explicit optimization + implicit optimization} + \underbrace{\sum_{i=1}^{N} \sum_{k=1}^{K} \gamma_{ik} \log(\frac{\pi_{k}}{\gamma_{ik}})}_{explicit optimization}, \underbrace{\sum_{i=1}^{N} \sum_{k=1}^{K} \sum_{i=1}^{K} \sum_{k=1}^{K} \gamma_{ik} \log(\frac{\pi_{k}}{\gamma_{ik}})}_{explicit optimization}, \underbrace{\sum_{i=1}^{N} \sum_{i=1}^{K} \sum$$

where  $\eta_{ij} = \sigma(\alpha + \beta^T y_{ij} - ||z_i - z_j||^2).$ 

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#### Algorithm 2 Estimation of deepLPM

```
Input: adjacency matrix A, edge features Y

pretrain_model = pretrain(A, 50 epochs) \triangleright pre-tr

encoder/decoder

while \mathcal{L} increases do

\tilde{\mu}_{\phi}, \tilde{\sigma}_{\phi}^2 = \text{GCN}(A)

explicit optimization (closed formulas):

update \hat{\gamma}_{ik}, \hat{\pi}_k, \hat{\mu}_k, \hat{\sigma}_k^2

calculate loss -\mathcal{L}

implicit optimization (SGD):

update encoder parameter \phi and decoder parameters \alpha, \beta
```

▷ pre-training to save initial weights of

### Numerical experiments: scenario A

Scenario A: 3 communities simulated according to LPCM, the mean of each cluster is set to

$$\begin{cases} \mu_1 = [0,0] \\ \mu_2 = [1.5 * \delta, 1.5 * \delta] \\ \mu_3 = [-1.5 * \delta, 1.5 * \delta] \end{cases}$$

where  $\delta \in [0.2, 0.95]$ .



Figure 14: Network simulated on scenario A.



Figure 15: Clustering ARI with different proximity rate  $\delta$  in Sc.A.

### Numerical experiments: scenario B

**Scenario B**: 1 cluster with large external connectivity and 2 communities with high internal connectivity based on **SBM** 

$$\Pi = \begin{pmatrix} b & a & a \\ a & a & b \\ a & b & a \end{pmatrix}$$

where a = 0.25,  $b = 0.01 + (1 - \delta') * (a - 0.01)$ , with  $\delta' \in [0.2, 1.0]$ 



Figure 16: Network simulated on scenario B.



Figure 17: Clustering ARI with different proximity rate  $\delta'$  in Sc.B.

# Model selection

A key element of our model is to be able to automatically determine P and K thanks to the auto-penalization of the deep encoder (Kingma et al., 2016; Dai et al., 2017).

### Model selection

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1) Fix K = 3, vary  $P \in \{2, 4, 8, 16, 32\}$ :



Figure 18: Averaged training loss (-ELBO) and clustering ARI 50 networks based on scenario B.

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Figure 18: Averaged training loss (-ELBO) and clustering ARI 50 networks based on scenario B. 2) Fix P = 16, vary  $K \in [2, 6]$ :



Figure 19: Averaged training loss (-ELBO) on 50 synthetic data in scenario B.

The considered data (Lamassé et al., 2014) report the ecclesiastical councils that took place in Merovingian Gaul during the 5th and 6th centuries.



**Figure 20:** Visualisation of the ecclesiastical network, highlighting the temporality of the relationships.



Figure 21: Visualisation of 8 cluster partitions with covariates on medieval data.

### Conclusion

#### The ZI-dLBM model:

- a model for the co-clustering of sparse evolving count data matrices,
- an interesting tool to summarize massive pharmacovigilance data and detect patterns,
- we plan to extend this model to the online setup to handle streams of ADR declarations.
- C. Bouveyron, M. Corneli and G. Marchello, A Deep Dynamic Latent Block Model for the Co-clustering of Zero-Inflated Data Matrices, Preprint HAL 03800210, Université Côte d'Azur, 2022
  - G. Marchello, A. Fresse, M. Corneli and C. Bouveyron, *Co-clustering of evolving count matrices in pharmacovigilance with the dynamic latent block model*, Statistics & Computing, in press, 2022.

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#### The deepLPM model:

- a (deep) latent model to represent and cluster network data with covariates,
- we applied it for visualizing and clustering a historical network,
- we plan to extend this model to handle text data as covariates on the nodes.
- C. Bouveyron, M. Corneli, P. Latouche D. Liang, *Clustering by Deep Latent Position Model with Graph Convolutional Network*, Preprint HAL n°03629104, Université Côte d'Azur, 2022.

"Ce qui est simple est toujours faux. Ce qui ne l'est pas est inutilisable."

Paul Valéry





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