Template based Graph Neural Network with Optimal Transport Distances

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Table of content

Optimal Transport and divergences between graphs
  - Gromov-Wasserstein divergence
  - Fused Gromov-Wasserstein and applications on attributed graphs

Template based Graph Neural Network with Optimal Transport Distances
  - Graph Neural Network
  - Template based GNN with FGW
  - Numerical experiments
Optimal Transport and divergences between graphs
Kantorovitch formulation: OT Linear Program

When $\mu_s = \sum_{i=1}^{n_s} a_i \delta_{x_i^s}$ and $\mu_t = \sum_{i=1}^{n_t} b_i \delta_{x_i^t}$

$$W^p_p(\mu_s, \mu_t) = \min_{T \in \Pi(\mu_s, \mu_t)} \left\{ \langle T, C \rangle_F = \sum_{i,j} T_{i,j} c_{i,j} \right\}$$

where $C$ is a cost matrix with $c_{i,j} = c(x_i^s, x_j^t) = \|x_i^s - x_j^t\|^p$ and the constraints are

$$\Pi(\mu_s, \mu_t) = \left\{ T \in (\mathbb{R}^+)^{n_s \times n_t} \mid T \mathbf{1}_{n_t} = a, T^T \mathbf{1}_{n_s} = b \right\}$$

- $W_p(\mu_s, \mu_t)$ is called the Wasserstein distance (EMD for $p = 1$).
- Entropic regularization solved efficiently with Sinkhorn [Cuturi, 2013].
- Classical OT needs distributions lying in the same space $\rightarrow$ Gromov-Wasserstein.
Optimal transport between discrete distributions

**Kantorovitch formulation : OT Linear Program**

When \( \mu_s = \sum_{i=1}^{n_s} a_i \delta_{x_i^s} \) and \( \mu_t = \sum_{i=1}^{n_t} b_i \delta_{x_i^t} \)

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where \( C \) is a cost matrix with \( c_{i,j} = c(x_i^s, x_j^t) = \|x_i^s - x_j^t\|^p \) and the constraints are

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\Pi(\mu_s, \mu_t) = \left\{ T \in (\mathbb{R}^+)^{n_s \times n_t} \mid T1_{n_t} = a, T^T1_{n_s} = b \right\}
\]

- \( W_p(\mu_s, \mu_t) \) is called the Wasserstein distance (EMD for \( p = 1 \)).
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Gromov-Wasserstein divergence

\[ \Omega_s \quad \Omega_t \]

\[ |d_X(x, x') - d_Y(y, y')| \]

Inspired from Gabriel Peyré

GW for discrete distributions [Memoli, 2011]

\[
GW_p(\mu_s, \mu_t) = \left( \min_{T \in \Pi(\mu_s, \mu_t)} \sum_{i,j,k,l} |D_{i,k} - D'_{j,l}|^p T_{i,j} T_{k,l} \right)^{\frac{1}{p}}
\]

with \( \mu_s = \sum_i a_i \delta_{x_i}^s \) and \( \mu_t = \sum_j b_j \delta_{x_j}^t \) and \( D_{i,k} = \|x_i^s - x_k^s\| \), \( D'_{j,l} = \|x_j^t - x_l^t\| \)

- Distance between metric measured spaces: across different spaces.
- Search for an OT plan that preserve the pairwise relationships between samples.
- Invariant to isometry in either spaces (e.g. rotations and translation).
- Entropy regularized GW proposed in [Peyré et al., 2016].
Attributed graphs as distributions

- Joint distribution $\mu$ in the feature/structure space.
  - Nodes are weighted by their mass $h_i$.
  - Structure encoded by $x_i$ ($D$ is adjacency matrix or shortest path).
  - Features values $a_i$ can be compared through the common metric.

- Importance of the joint modeling:

\[
\begin{align*}
\mu &= \sum_i h_i \delta(x_i, a_i) \\
\mu_A &= \sum_i h_i \delta a_i \\
\mu_X &= \sum_i h_i \delta x_i
\end{align*}
\]
Fused Gromov-Wasserstein distance [Vayer et al., 2020]

\[ \mu_s = \sum_{i=1}^{n} h_i \delta_{x_i, a_i} \quad \text{and} \quad \mu_t = \sum_{j=1}^{m} g_j \delta_{y_j, b_j} \]

\[ \mathcal{FGW}_{p, q, \alpha}(D, D', \mu_s, \mu_t) = \left( \min_{T \in \Pi(\mu_s, \mu_t)} \sum_{i,j,k,l} \left( (1-\alpha)|a|_{C^{q}_{i,j}} + \alpha|D_{i,k}-D'_{j,l}|^q \right)^p T_{i,j} T_{k,l} \right)^{\frac{1}{p}} \]

with \( D_{i,k} = \|x_i - x_k\| \) and \( D'_{j,l} = \|y_i - y_l\| \) and \( C_{i,j} = \|a_i - b_j\| \)

- Parameters \( q > 1, \forall p \geq 1 \).
- \( \alpha \in [0, 1] \) is a trade off parameter between structure and features.
Gromov-Wasserstein distance [Memoli, 2011]

- Divergence between distributions across metric spaces.
- Can be used to measure similarity between graphs seen as distribution their pairwise node relationship.

Fused Gromov-Wasserstein distance [Vayer et al., 2018]

- Model labeled structured data as joint structure/labels distributions.
- New versatile method for comparing structured data based on Optimal Transport
- New notion of barycenter of structured data such as graphs or time series

How to use GW/FGW in graph neural networks?
Template based Graph Neural Network with Optimal Transport Distances
**Graph Neural Networks**

- Each layer of the GNN computes features on graph node using the values from the connected neighbors: message passing principle.
- A step of global aggregation or pooling allows to go from a complex graph object to a vector representation.
- The pooling step must remain invariant to permutations (min, max, mean).

**Principle [Bronstein et al., 2017]**

Can we encode graphs as distributions for pooling in GNN?
Template based Graph Neural Network with OT Distances

- Principle: represent a graph through its distances to learned templates.
- Novel pooling layer derived from OT distances.
- New end-to-end GNN models for graph-level tasks.
- Learnable parameters are illustrated in red above.

Template based FGW layer (TFGW) [Vincent-Cuaz et al., 2022b]
1. Modeling graphs as discrete distributions

- \( D_i \): node relationship matrix e.g. adjacency, shortest-path, laplacian, etc.
- \( F_i \): node feature matrix.
- \( h_i \): nodes relative importance (probabilities).
2. Node embeddings

- \( \phi_u \): GNN of \( L \) layers parameterized by \( u \) e.g. GIN, GAT, etc.
- Promotes discriminant features on the nodes \( \phi_u(F_i) \)
3. Template-based Fused Gromov-Wasserstein (TFGW) pooling

- $\text{FGW}_\alpha$: OT soft graph matching distance.
- $\alpha \in [0; 1]$: relative importance between structure $D_i$ and node features $\phi_u(F_i)$.
- $\{D_k, \bar{F}_k, \bar{h}_k\}$: FGW distances to $K$ templates used as graph representation.
4. Final MLP for predictions

- $\psi_v$: MLP with non-linearities fed with the distance embeddings.
- $\hat{y}_i$: final prediction for graph-level tasks (classification or regression).
- End-to-end optimization of all parameters:
  - $u$ and $v$ parameters of GNN $\phi_u$ and final MLP $\psi_v$.
  - $\{D_k, \overline{F}_k, \overline{h}_k\}$ TFGW graph templates.
### TFGW benchmark

<table>
<thead>
<tr>
<th>category</th>
<th>model</th>
<th>MUTAG</th>
<th>PTC</th>
<th>ENZYMES</th>
<th>PROTEIN</th>
<th>NCI1</th>
<th>IMDB-B</th>
<th>IMDB-M</th>
<th>COLLAB</th>
</tr>
</thead>
<tbody>
<tr>
<td>Ours</td>
<td>TFGW ADJ (L=2)</td>
<td>96.4(3.3)</td>
<td>72.4(5.7)</td>
<td>73.8(4.6)</td>
<td>82.9(2.7)</td>
<td>88.1(2.5)</td>
<td>78.3(3.7)</td>
<td>56.8(3.1)</td>
<td>84.3(2.6)</td>
</tr>
<tr>
<td></td>
<td>TFGW SP (L=2)</td>
<td>94.8(3.5)</td>
<td>70.8(6.3)</td>
<td>75.1(5.0)</td>
<td>82.0(3.0)</td>
<td>86.1(2.7)</td>
<td>74.1(5.4)</td>
<td>54.9(3.9)</td>
<td>80.9(3.1)</td>
</tr>
<tr>
<td>OT emb.</td>
<td>OT-GNN (L=2)</td>
<td>91.6(4.6)</td>
<td>68.0(7.5)</td>
<td>66.9(3.8)</td>
<td>76.6(4.0)</td>
<td>82.9(2.1)</td>
<td>67.5(3.5)</td>
<td>52.1(3.0)</td>
<td>80.7(2.9)</td>
</tr>
<tr>
<td></td>
<td>OT-GNN (L=4)</td>
<td>92.1(3.7)</td>
<td>65.4(9.6)</td>
<td>67.3(4.3)</td>
<td>78.0(5.1)</td>
<td>83.6(2.5)</td>
<td>69.1(4.4)</td>
<td>51.9(2.8)</td>
<td>81.1(2.5)</td>
</tr>
<tr>
<td></td>
<td>WEGL</td>
<td>91.0(3.4)</td>
<td>66.0(2.4)</td>
<td>60.0(2.8)</td>
<td>73.7(1.9)</td>
<td>75.5(1.4)</td>
<td>66.4(2.1)</td>
<td>50.3(1.0)</td>
<td>79.6(0.5)</td>
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<tr>
<td>GNN</td>
<td>PATCHYSAN</td>
<td>91.6(4.6)</td>
<td>58.9(3.7)</td>
<td>55.9(4.5)</td>
<td>75.1(3.3)</td>
<td>76.9(2.3)</td>
<td>62.9(3.9)</td>
<td>45.9(2.5)</td>
<td>73.1(2.7)</td>
</tr>
<tr>
<td></td>
<td>GIN</td>
<td>90.1(4.4)</td>
<td>63.1(3.9)</td>
<td>62.2(3.6)</td>
<td>76.2(2.8)</td>
<td>82.2(0.8)</td>
<td>64.3(3.1)</td>
<td>50.9(1.7)</td>
<td>79.3(1.7)</td>
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<tr>
<td></td>
<td>DropGIN</td>
<td>89.8(6.2)</td>
<td>62.3(6.8)</td>
<td>65.8(2.7)</td>
<td>76.9(4.3)</td>
<td>81.9(2.5)</td>
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<td>80.1(2.8)</td>
</tr>
<tr>
<td></td>
<td>PPGN</td>
<td>90.4(5.6)</td>
<td>65.6(6.0)</td>
<td>66.9(4.3)</td>
<td>77.1(4.0)</td>
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<td>81.0(2.1)</td>
</tr>
<tr>
<td></td>
<td>DIFFPOOL</td>
<td>86.1(2.0)</td>
<td>45.0(5.2)</td>
<td>61.0(3.1)</td>
<td>71.7(1.4)</td>
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</tr>
<tr>
<td>Kernels</td>
<td>FGW - ADJ</td>
<td>82.6(7.2)</td>
<td>55.3(8.0)</td>
<td>72.2(4.0)</td>
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</tr>
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<td></td>
<td>FGW - SP</td>
<td>84.4(7.3)</td>
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<td>47.8(3.8)</td>
<td>77.8(2.4)</td>
</tr>
<tr>
<td></td>
<td>WL</td>
<td>87.4(5.4)</td>
<td>56.0(3.9)</td>
<td>69.5(3.2)</td>
<td>74.4(2.6)</td>
<td>85.6(1.2)</td>
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<tr>
<td></td>
<td>WWL</td>
<td>86.3(7.9)</td>
<td>52.6(6.8)</td>
<td>71.4(5.1)</td>
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</tr>
</tbody>
</table>

- Comparison with state of the art approach from GNN and graph kernel methods.
- Systematic and significant gain of performance with GIN+TFGW.
- Gain independent of GNN architecture (GIN or GAT).
- 1 year after publication world rankings of TFGW on "papers with code": #1 NCI1, #2 COLLAB ENZYMES IMDB-M, #3 MUTAG, PROTEIN.
Going beyond Weisfeiler-Lehman Isomorphism tests

<table>
<thead>
<tr>
<th>4-Cycles dataset</th>
<th>Skip-Circles dataset</th>
</tr>
</thead>
<tbody>
<tr>
<td>sample $y=0$</td>
<td>sample $y=6$ (11 hops)</td>
</tr>
<tr>
<td></td>
<td>sample $y=1$</td>
</tr>
</tbody>
</table>

TFGW with $K = \#\text{labels}$

<table>
<thead>
<tr>
<th>model</th>
<th>4-Cycles</th>
<th>Skip-Circles</th>
</tr>
</thead>
<tbody>
<tr>
<td>TFGW</td>
<td>0.99(0.03)</td>
<td>1.00(0.00)</td>
</tr>
<tr>
<td>TFGW-fix</td>
<td>0.63(0.11)</td>
<td>1.00(0.00)</td>
</tr>
<tr>
<td>OT-GNN</td>
<td>0.50(0.00)</td>
<td>0.10(0.00)</td>
</tr>
<tr>
<td>GIN</td>
<td>0.50(0.00)</td>
<td>0.10(0.00)</td>
</tr>
<tr>
<td>DropGIN</td>
<td>1.00(0.01)</td>
<td>0.82(0.28)</td>
</tr>
<tr>
<td>PPGN</td>
<td>1.00(0.01)</td>
<td>0.90(0.11)</td>
</tr>
</tbody>
</table>

TFGW with various $K$

- TFGW has good expressivity on problems beyond WL test.
- Learning the templates is important: TFGW $>$ TFGW fix.
Graph OT distance embedding

**TFGW embedding (PCA)**

PCA in the FGW embedding for PTC with no GNN

PCA in the FGW embedding for PTC with 2 GIN layers

PCA in the FGW embedding for PTC with 2 GIN layers

**Learned templates (left) and data samples (right)**

FGW templates for PTC

Samples from PTC
Gromov-Wasserstein family for graph modeling

- Graphs modelled as distributions, $\mathcal{GW}$ can measure their similarity.
- Extensions of GW for labeled graphs and Frechet means can be computed.
- TFGW for graph pooling in GNNs [Vincent-Cuaz et al., 2022b].
- Weights on the nodes are important but rarely available: relax the constraints [Séjourné et al., 2020] or even remove one of them [Vincent-Cuaz et al., 2022a].
- Many applications of FGW from brain imagery [Thual et al., 2022] to Graph Neural Networks [Vincent-Cuaz et al., 2022b].
Python code available on GitHub:

https://github.com/PythonOT/POT

- OT LP solver, Sinkhorn (stabilized, $\epsilon-$scaling, GPU)
- Domain adaptation with OT.
- Barycenters, Wasserstein unmixing.
- Gromov Wasserstein.
- Solvers for Numpy/Pytorch/Jax/tensorflow/Cupy

Tutorial on OT for ML:
http://tinyurl.com/otml-isbi

Papers available on my website:
https://remi.flamary.com/


Sample complexity of sinkhorn divergences.


Gromov wasserstein distances and the metric approach to object matching.


Gromov-wasserstein averaging of kernel and distance matrices.

In *ICML*, pages 2664–2672.


Linear-time gromov wasserstein distances using low rank couplings and costs.


The unbalanced gromov wasserstein distance: Conic formulation and relaxation.


Aligning individual brains with fused unbalanced gromov-wasserstein.

In *Neural Information Processing Systems (NeurIPS).*


Convergence of a block coordinate descent method for nondifferentiable minimization.


Fused gromov-wasserstein distance for structured objects: theoretical foundations and mathematical properties.


Entropic regularized optimal transport

Distributions

Reg. OT matrix with \( \lambda=1e-3 \)

Reg. OT matrix with \( \lambda=1e-2 \)

Entropic regularization [Cuturi, 2013]

\[
W_\epsilon(\mu_s, \mu_t) = \min_{T \in \Pi(\mu_s, \mu_t)} \langle T, C \rangle_F + \epsilon \sum_{i,j} T_{i,j} \log T_{i,j}
\]

- Regularization with the negative entropy \(-H(T)\).
- Looses sparsity, but strictly convex optimization problem [Benamou et al., 2015].
- Can be solved with the very efficient Sinkhorn-Knopp matrix scaling algorithm.
- Loss and OT matrix are differentiable and have better statistical properties [Genevay et al., 2018].
Entropic regularized optimal transport

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GW Upper bond [Vincent-Cuaz et al., 2021]

Let two graphs of order $N$ in the linear embedding $P_s w^1_s$ and $P_s w^2_s$, the $GW$ divergence can be upper bounded by

$$GW^2 \left( \sum_{s \in [S]} w^1_s D_s, \sum_{s \in [S]} w^2_s D_s \right) \leq \|w^1 - w^2\|_M$$  \hspace{1cm} (1)

with $M$ a PSD matrix of components $M_{p,q} = \langle D_h D_p, D_q D_h \rangle_F$, $D_h = diag(h)$.

Discussion

- The upper bound is the value of $GW$ for a transport $T = diag(h)$ assuming that the nodes are already aligned.
- The bound is exact when the weights $w^1$ and $w^2$ are close.
- Solving $GW$ with FW is $O(N^3 \log(N))$ at each iterations.
- Computing the Mahalanobis upper bound is $O(S^2)$: very fast alternative to $GW$ for nearest neighbors retrieval.
Solving the Gromov Wasserstein optimization problem

Optimization problem

$$\mathcal{GW}_p^p(\mu_s, \mu_t) = \min_{T \in \Pi(\mu_s, \mu_t)} \sum_{i,j,k,l} |D_{i,k} - D'_{j,l}|^p T_{i,j} T_{k,l}$$

with $\mu_s = \sum_i a_i \delta_{x^s_i}$ and $\mu_t = \sum_j b_j \delta_{x^t_j}$ and $D_{i,k} = \|x^s_i - x^s_k\|$, $D'_{j,l} = \|x^t_j - x^t_l\|

- Quadratic Program (Wasserstein is a linear program).
- Nonconvex, NP-hard, related to Quadratic Assignment Problem (QAP).
- Large problem and non convexity forbid standard QP solvers.

Optimization algorithms

- Local solution with conditional gradient algorithm (Frank-Wolfe) [Frank and Wolfe, 1956].
- Each FW iteration requires solving an OT problems.
- Gromov in 1D has a close form (solved in discrete with a sort) [Vayer et al., 2019].
- With entropic regularization, one can use mirror descent [Peyré et al., 2016] or fast low rank approximations [Scetbon et al., 2021].
Entropic Gromov-Wasserstein

Optimization Problem

\[
\mathcal{GW}^p_{p,\epsilon}(\mu_s, \mu_t) = \min_{T \in \Pi(\mu_s, \mu_t)} \sum_{i,j,k,l} |D_{i,k} - D'_{j,l}|^p T_{i,j} T_{k,l} + \epsilon \sum_{i,j} T_{i,j} \log T_{i,j} \tag{2}
\]

with \( \mu_s = \sum_i a_i \delta_{x_i^s} \) and \( \mu_t = \sum_j b_j \delta_{x_j^t} \) and \( D_{i,k} = \|x_i^s - x_k^s\|, D'_{j,l} = \|x_j^t - x_l^t\| \)

- Smoothing the original GW with a convex and smooth entropic term.

Solving the entropic \( \mathcal{GW} \) [Peyré et al., 2016]

- Problem (2) can be solved using a KL mirror descent.
- This is equivalent to solving at each iteration \( t \)
  \[
  T^{(t+1)} = \min_{T \in \mathcal{P}} \left\langle T, G^{(t)} \right\rangle_F + \epsilon \sum_{i,j} T_{i,j} \log T_{i,j}
  \]
  Where \( G^{(t)}_{i,j} = 2 \sum_{k,l} |D_{i,k} - D'_{j,l}|^p T^{(t)}_{k,l} \) is the gradient of the GW loss at previous point \( T^{(k)} \).
- Problem above solved using a Sinkhorn-Knopp algorithm of entropic OT.
- Very fast approximation exist for low rank distances [Scetbon et al., 2021].
FGW Properties

\[
\mathcal{FGW}_{p,q,\alpha}(D, D', \mu_s, \mu_t) = \left( \min_{T \in \Pi(\mu_s, \mu_t)} \sum_{i,j,k,l} \left( (1 - \alpha)C_{i,j}^q + \alpha |D_{i,k} - D'_{j,l}|^q \right)^p T_{i,j} T_{k,l} \right)^{\frac{1}{p}}
\]

Metric properties [Vayer et al., 2020]

- **FGW** defines a metric over structured data with **measure and features preserving isometries** as invariants.
- **FGW** is a metric for \( q = 1 \) a semi metric for \( q > 1 \), \( \forall p \geq 1 \).
- The distance is nul **iff**:
  - There exists a Monge map \( T \# \mu_s = \mu_t \).
  - Structures are equivalent through this Monge map (isometry).
  - Features are equal through this Monge map.

Bounds and convergence to finite samples [Vayer et al., 2020]

- **FGW**(\( \mu_s, \mu_t \)) is lower bounded by \((1 - \alpha)\mathcal{W}(\mu_A, \mu_B)^q\) and \(\alpha\mathcal{GW}(\mu_X, \mu_Y)^q\)
- Convergence of finite samples when \( \mathcal{X} = \mathcal{Y} \) with \( d = \text{Dim}(\mathcal{X}) + \text{Dim}(\Omega) \):

\[
\mathbb{E} [\mathcal{FGW}(\mu, \mu_n)] = O \left( n^{-\frac{1}{d}} \right)
\]
Solving the unmixing problem

Optimization problem

\[
\min_{\mathbf{w} \in \Sigma_S} \mathcal{GW}_2^2 \left( \sum_{s \in [S]} w_s \overline{D_s}, D \right) - \lambda \| \mathbf{w} \|_2^2
\]

- Non-convex Quadratic Program w.r.t. \( T \) and \( \mathbf{w} \).
- \( \mathcal{GW} \) for fixed \( \mathbf{w} \) already have an existing Frank-Wolfe solver.
- We proposed a Block Coordinate Descent algorithm

**BCD Algorithm for sparse \( \mathcal{GW} \) unmixing [Tseng, 2001]**

1: repeat
2: Compute OT matrix \( T \) of \( \mathcal{GW}_2^2(D, \sum_s w_s \overline{D_s}) \), with FW [Vayer et al., 2018].
3: Compute the optimal \( \mathbf{w} \) given \( T \) with Frank-Wolfe algorithm.
4: until convergence

- Since the problem is quadratic optimal steps can be obtained for both FW.
- BCD convergence in practice in a few tens of iterations.
GDL Extensions

GDL on labeled graphs

- For datasets with labeled graphs, one can learn simultaneously a dictionary of the structure \( \{D_s\}_{s \in [S]} \) and a dictionary on the labels/features \( \{F_s\}_{s \in [S]} \).
- Data fitting is Fused Gromov-Wasserstein distance \( \mathcal{FGW} \), same stochastic algorithm.

Dictionary on weights

\[
\min_{\{(w^{(k)}, v^{(k)})\}_k, \{D_s, h_s\}_s} \sum_{k=1}^{K} GW^2_2 \left( D^{(k)}, \sum_s w_s^{(k)} D_s, h^{(k)}, \sum_s v_s^{(k)} h_s \right) - \lambda \|w^{(k)}\|_2^2 - \mu \|v^{(k)}\|_2^2
\]

- We model the graphs as a linear model on the structure and the node weights

\[
(D^{(k)}, h^{(k)}) \rightarrow \left( \sum_s w_s^{(k)} D_s, \sum_s v_s^{(k)} h_s \right)
\]

- This allows for sparse weights \( h \) so embedded graphs with different order.
- We provide in [Vincent-Cuaz et al., 2021] subgradients of GW w.r.t. the mass \( h \).
Cluster 1

Cluster 2

Cluster 3

Cluster 4

Training dataset examples

Centroids

• Clustering of multiple real-valued graphs. Dataset composed of 40 graphs (10 graphs × 4 types of communities)

• $k$-means clustering using the $FGW$ barycenter
FGW barycenter

Euclidean barycenter

\[ \min_x \sum_k \lambda_k \| x - x_k \|^2 \]

FGW barycenter

\[ \min_{D \in \mathbb{R}^{n \times n}, \mu} \sum_i \lambda_i \mathcal{FGW}(D_i, D, \mu_i, \mu) \]

**FGW barycenter** \( p = 1, q = 2 \)

- Estimate FGW barycenter using Frechet means (similar to [Peyré et al., 2016]).
- Barycenter optimization solved via block coordinate descent (on \( T, D, \{ a_i \}_i \)).
- Can chose to fix the structure (\( D \)) or the features \( \{ a_i \}_i \) in the barycenter.
- \( a_{ii} \), and \( D \) updates are weighted averages using \( T \).
FGW barycenter on labeled graphs

Barycenter of noisy graphs

- We select a clean graph, change the number of nodes and add label noise and random connections.
- We compute the barycenter on $n = 15$ and $n = 7$ nodes.
- Barycenter graph is obtained through thresholding of the $D$ matrix.
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FGW barycenter for community clustering

Graph approximation and community clustering

\[
\min_{D, \mu} \mathcal{FGW}(D, D_0, \mu, \mu_0)
\]

- Approximate the graph \((D_0, \mu_0)\) with a small number of nodes.
- Can be seen as a FGW (compressed) barycenter for one graph.
- OT matrix give the clustering affectation.
- Works for signle and multiple modes in the clusters.
Experiments - Unsupervised representation learning

Comparison of fixed and learned weights dictionaries

- Graph taken from the IMBD dataset.
- Show original graph and representation after projection on the embedding.
- Uniform weight $h$ has a hard time representing a central node.
- Estimated weights $\tilde{h}$ recover a central node.
- In addition some nodes are discarded with 0 weight (graphs can change order).
Clustering Experiments on real datasets

- Different data fitting losses:
  - Graphs without node attributes: Gromov-Wasserstein.
  - Graphs with node attributes (discrete and real): Fused Gromov-Wasserstein.

- We learn a dictionary on the dataset and perform K-means in the embedding using the Mahalanobis distance approximation.

- Compared to GW Factorization (GWF) [Xu, 2020] and spectral clustering.

- Similar performance for supervised classification (using GW in a kernel).
Aligning individual brains with Fused Unbalanced Gromov-Wasserstein
Inter-subject anatomical and functional variability

- Math-nonmath contrast map from the Mathlang protocol for 3 IBC subjects
- Each subject has different surfaces (mesh) and signal
- Traditional approach maps the signal on an average mesh (fsaverage bottom line) to compute an average.
Two subjects recorded with FMRI doing the same mental task.
We seek an alignment preserving both the local features and the cortex geometry.
Area can have different surface: relax marginal constraints.
Fused Unbalanced Gromov-Wasserstein

\[ L_\theta(P) \triangleq (1 - \alpha) \sum_{0 \leq i < n, 0 \leq j < p} \| F^s_i - F^t_j \|_2^2 P_{i,j} + \alpha \sum_{0 \leq i, k < n, 0 \leq j, l < p} |D^s_{i,k} - D^t_{j,l}|^2 P_{i,j} P_{k,l} + \rho \left( KL(P_{#1} \otimes P_{#1} | w^s \otimes w^s) + KL(P_{#2} \otimes P_{#2} | w^t \otimes w^t) \right) + \varepsilon E(P) \]

Principle [Thual et al., 2022]

- Preserve the features (Wasserstein loss) and the cortex geometry (GW loss).
- Relax the marginal constraints to allows creation/destruction of mass to encore change in surface of the areas in the brain across subjects.
Features VS structure/anatomy preservation

Illustration of the maps (transported atlas) and displacement on the geodesic as a function of $\alpha$. 
Aligning pairs of individuals with FUGW

- Aligning with FUGW leads to significantly increased correlations across subjects.
- Similar gains on other types of stimuli and acquisition time.
Aligning individuals to a FUGW barycenter

\[ \mathbf{X}^B = (F^B, D^B, w^B) \in \arg \min_{\mathbf{X}} \sum_{s \in S} \text{FUGW}(\mathbf{X}^s, \mathbf{X}) \]

Principle

- We compute a barycenter of different subjects (with fixed anatomy).
- FUGW barycenter significantly increases statistical power of group averages.