# Graph-Based Weight Cascading Methods for Multisite Time Series Forecasting

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### **Motivations**

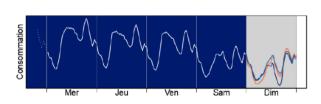
#### **Industrial** context

Anticipation of the consumption of **electricity** and **renewable energy production** is a major challenge for EDF, especially for electricity market operations:

- ▶ Maintaining a balance between electricity supply and demand is important for grid stability;
- Optimizing the production fleet and demand response;
- > **Buying** and **selling** on electricity markets

New **geolocated data** can be exploited by spatial models — such as **GNNs** — and improve forecasts.

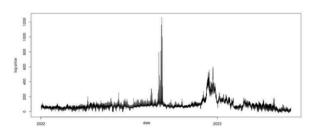
(Obst, Vilmarest, and Goude 2020; Vilmarest and Goude 2021)



**Figure 1** – Electricity consumption.



**Figure 2** – Renewable energy production.

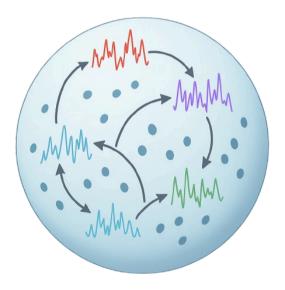


**Figure 3** – Electricity prices.

### **Motivations**

#### *Academic context*

- $\triangleright$  You have a **large dataset** of N time series with a **limited number of observations** T;
- $\triangleright$  You want to have N accurate forecasts but you **do not have a huge budget**.



**Figure 4** – Timeseries may be hard to order.

### **Motivations**

**Approaches** 

We have tested three different approaches:

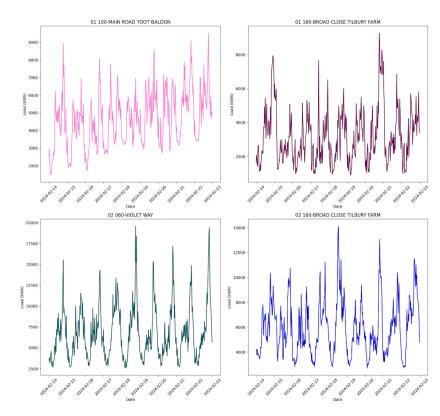
**Approach 1 – Individual**: Train N individual models.

**Approach 2 – Cascade**: Transfer models weights through a tree structure.

**Approach 3 – GNNs**: Train a single Graph Neural Network.

### About the dataset

- ➤ The dataset¹ consists of aggregated half-hourly residential smart meter electricity consumption data collected by four UK Distribution Network Operators (DNOs);
- ▶ 120,000 low voltage feeders;
  → very heterogeneous data;
- Dataset spans January 2024;



**Figure 5** – Subset of 4 nodes of Oxford's urban area.

<sup>&</sup>lt;sup>1</sup>https://weave.energy/

### About the models

#### Feedforward Neural Networks & Graph Neural Networks

#### **Feedforward Neural Networks**

➤ The general update rule of a hidden vector is given by:

$$\boldsymbol{h}^{(\ell+1)} = \sigma \big( \boldsymbol{W}^{(\ell+1)} \boldsymbol{h}^{(\ell)} + \boldsymbol{b}^{(\ell+1)} \big)$$

#### where:

- $\hookrightarrow m{W}^{(\ell+1)} \in \mathbb{R}^{d_{\ell+1} imes d_{\ell}}$  is a learned weight matrix;
- $\rightarrow b^{(\ell+1)} \in \mathbb{R}^{d_{\ell+1}}$  is a learned **bias** vector;
- $\rightarrow \sigma$  is a non-linear **activation function** (e.g. ReLU, tanh).

#### **Graph Neural Networks** (Gori and Monfardini 2005)

 $\triangleright$  The general update rule for a node u is given by:

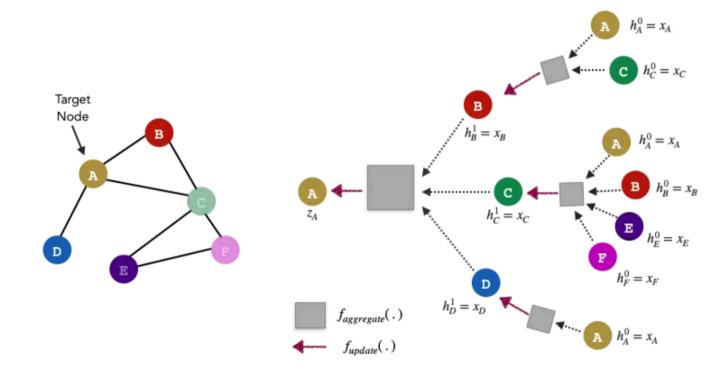
$$\textstyle \boldsymbol{h}_{u}^{(\ell+1)} = \phi\Big(\boldsymbol{h}_{u}^{(\ell)}, \bigoplus_{v \in \mathcal{N}_{u}} \psi\Big(\boldsymbol{h}_{u}^{(\ell)}, \boldsymbol{h}_{v}^{(\ell)}, \boldsymbol{e}_{uv}\Big)\Big)$$

#### where:

- $\hookrightarrow \mathcal{N}_u$  is the set of neighbors of u;
- $\rightarrow$   $\phi$  and  $\psi$  are respectively **update** and **message** functions;
- $\rightarrow$   $\bigoplus$  is the **aggregation** operator;
- $\mapsto e_{uv}$  is the edge representation between u and v.

### About the models

Visual understanding of a GNN



### Two examples of graph convolutions

Graph Convolutional Networks & Graph Attention Networks

# **Graph Convolutional Networks** (Kipf and Welling 2016)

 $\triangleright$  The update rule for a node u is given by:

$$m{h}_u^{(\ell+1)} = \sigma \Big( \sum_{v \in \mathcal{N}_u \cup \{u\}} c_{vu} m{\Theta}^{(\ell)} m{h}_v^{(\ell)} \Big)$$

where

$$c_{vu} = \frac{e_{vu}}{\sqrt{d_v d_u}}.$$

#### **Graph Attention Networks**

(Veličković et al. 2017; Brody, Alon, and Yahav 2022)

 $\triangleright$  The update rule for a node u is given by:

$$\mathbf{h}_u^{(\ell+1)} = \sigma \Big( \sum_{v \in \mathcal{N}_u \cup \{u\}} \alpha_{uv} \mathbf{\Theta}_t^{(\ell)} \mathbf{h}_v^{(\ell)} \Big)$$

where

$$\alpha_{uv} = \frac{\exp\Bigl(\boldsymbol{a}^{\top}\sigma\Bigl(\boldsymbol{\Theta}_{\!s}^{(\ell)}\boldsymbol{h}_{\!u}^{(\ell)} + \boldsymbol{\Theta}_{\!t}^{(\ell)}\boldsymbol{h}_{\!v}^{(\ell)} + \boldsymbol{\Theta}_{\!e}^{(\ell)}\boldsymbol{e}_{uv}\Bigr)\Bigr)}{\sum_{k\in\mathcal{N}_{\!u}\cup\{u\}}\exp\Bigl(\boldsymbol{a}^{\top}\sigma\Bigl(\boldsymbol{\Theta}_{\!s}^{(\ell)}\boldsymbol{h}_{\!u}^{(\ell)} + \boldsymbol{\Theta}_{\!t}^{(\ell)}\boldsymbol{h}_{\!k}^{(\ell)} + \boldsymbol{\Theta}_{\!e}^{(\ell)}\boldsymbol{e}_{uk}\Bigr)\Bigr)}.$$

### About tree algorithms

#### **Minimum Spanning Trees**

- ▷ Apply to undirected graphs;
- Select a subset of edges connecting all nodes with:
  - → Minimum total edge weight;
  - → No cycles;
- ▷ Efficiently computed using **Kruskal**'s or Prim's algorithms;
- Commonly used in network design, clustering, and optimization problems. (Cong and Zhao 2015)

#### **Minimum Cost Arborescences**

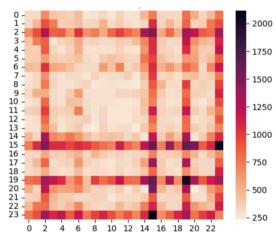
- > Apply to **directed** graphs;
- ▷ Build a rooted spanning tree with:
  - → Minimum total cost of directed edges;
  - → Reachability from the root to all nodes;
- Solved using Chu-Liu Edmonds' algorithm (Chu and Liu 1965; Edmonds 1967);
- □ Useful for hierarchical structures, flow networks, and decision trees.

#### Building a diffusion tree

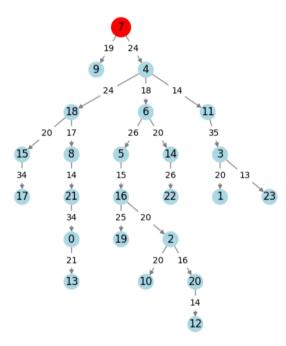
- - → distance-based (e.g. euclidean distance);
  - → spectral-based:

$$oldsymbol{L} = oldsymbol{I} - rac{1}{2} \Big( oldsymbol{\Phi}^{rac{1}{2}} oldsymbol{P} oldsymbol{\Phi}^{-rac{1}{2}} + oldsymbol{\Phi}^{-rac{1}{2}} oldsymbol{P}^ op oldsymbol{\Phi}^{rac{1}{2}} \Big)$$

where P is a transition matrix and  $\Phi$  a matrix with the Perron vector of P in the diagonal and zeros elsewhere (Chung 2005);



**Figure 6** – Distance matrix used to build the diffusion tree.

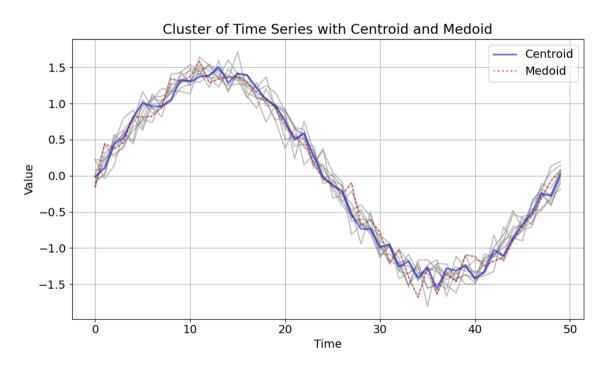


**Figure 7** – Diffusion tree built from the distance matrix.

#### On the prototype selection

- The **prototype** p is the root of the cascade and acts as a source of weight diffusion;
- > Three strategies:

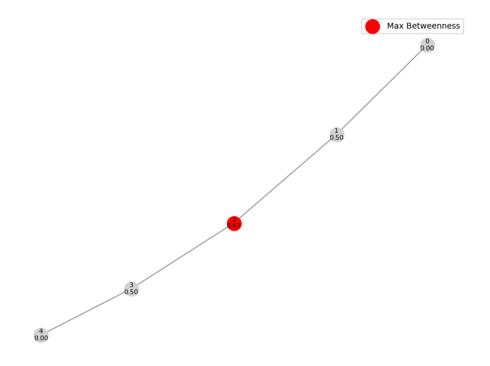
  - → Medoid: most central real point (minimum of pairwise distances); robust to outliers;
  - → **Betweenness centrality**: node with highest betweenness centrality in graph; captures topological importance.



**Figure 8** – Centroid and medoid strategies.

#### On the prototype selection

- The **prototype** p is the root of the cascade and acts as a source of weight diffusion;
- ▷ Three strategies:
  - → **Centroid**: mean of all points; efficient but sensitive to outliers;
  - → **Medoid**: most central real point (minimum of pairwise distances); robust to outliers;
  - → **Betweenness centrality**: node with highest betweenness centrality in graph; captures topological importance.

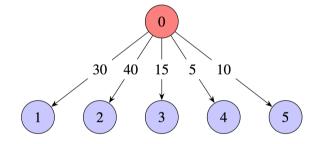


**Figure 9** – Betweenness centrality strategy.

### On the algorithm

#### **Cascading algorithm**

- Consists of 2 stages:
  - $\rightarrow \mathcal{A}_0$ : train prototype model on p;
  - $\hookrightarrow \mathcal{A}_1$ : refine each model using parent weights;
- > **Single-step**: prototype weights broadcast to all cluster members;
- ightharpoonup Multi-step: weights flow through a tree  $\mathcal{T}$  (MST/MCA) from parent to child;
  - → Enables gradual diffusion.

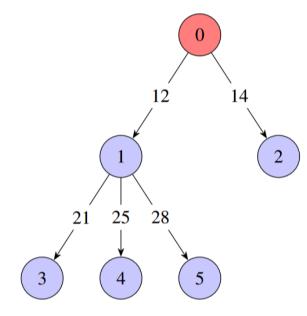


**Figure 10** – Single-step cascade for a total budget of 100.

### *On the algorithm*

#### **Cascading algorithm**

- Consists of 2 stages:
  - $\hookrightarrow \mathcal{A}_0$ : train prototype model on p;
  - $\rightarrow \mathcal{A}_1$ : refine each model using prototype weights;
- ➤ Single-step: prototype weights broadcast to all cluster members;
  - → Can use uniform or distance-based budgets;
- ightharpoonup Multi-step: weights flow through a tree  $\mathcal{T}$  (MST/MCA) from parent to child;



**Figure 11** – Multiple-step cascade for a total budget of 100.

# About the budget

What is "fair"?

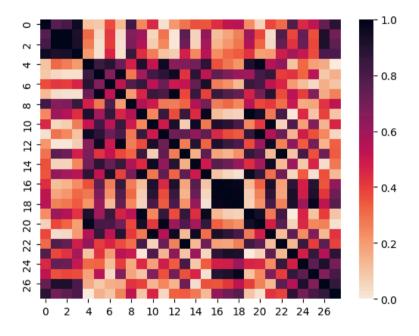
- $\triangleright$  The learning models deployed across these nodes share a **global computational budget** B, and operate using a **fixed batch size**:
  - $\hookrightarrow$  Individual budgets:  $\forall u \in \mathcal{V}, B_u = \frac{B}{N}$  and models weights are randomly initialized;
  - **└→ Cascade budgets** 
    - ightharpoonup Uniform:  $\forall u \in \mathcal{V}, B_u = \frac{B}{N}$ , prototype's model weights are randomly initialized and each child inherits the parent's weights;
    - hinspace > Flexible:  $\forall u,v \in \mathcal{V}, u \to v, \sum_{u \to v} \widetilde{d_{uv}} = 1, B_v = \lceil \widetilde{d_{uv}}B \rceil;$   $\left(\sum_{v \in \mathcal{V}} B_v \simeq B\right)$
  - $\hookrightarrow$  **GNN budget**: *B*.

# Cascade diffusion experiments on synthetic data

- ▶ Both single-step and multi-step cascades outperform individual models especially when the budget is small;
- - → Investigating under **which conditions** multi-step cascades outperform single-step!
- > "Well"-chosen diffusion trees **significantly perform better than random** trees.

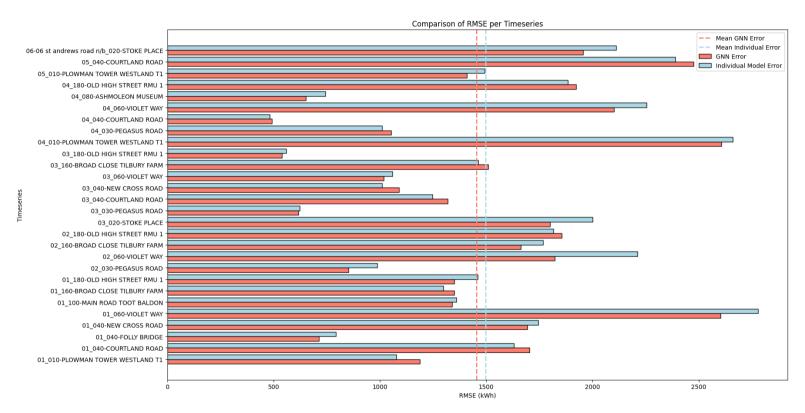
# When Graph Neural Networks Come Into Play

 $\triangleright$  **Designing a diffusion tree**  $\mathcal{T}$ : train a GAT(v2) model and extract **attention weights** to then build a diffusion tree;



▶ A global model: GNNs can also be used as a global model by relying on the spatial links between nodes to efficiently compute representations.

# GNN experiments on real data



**Figure 12** – Comparison of RMSE per timeseries for a **GATv2** and a FFNN with  $B_u=10$ .

### Conclusion

- ▷ Cascading through MSTs or MCAs enables low-cost, scalable model refinement;
- ▷ GNNs can serve as:
  - → a tree generation method to diffuse weights across sites;
  - $\rightarrow$  a **single global model** that captures structural information across sites.

Thanks for your attention!

Feel free to reach out to me at eloi.campagne@ens-paris-saclay.fr.

# **Bibliography**

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# **Appendix**

### Cascading algorithm

#### Algorithm 3 MST Cascade.

- 1: Input: Data  $\{\mathbf{x}_i\}_{i=1}^N$ , distance function between tasks  $\mathtt{dist}()$ , number of clusters K to split the tasks, clustering method  $\mathtt{findClusters}()$ , method that finds a prototype for a given cluster  $\mathtt{findPrototype}()$  budget for prototype training b, total budget per cluster B, training procedures  $A_0$  and  $A_1$ .
- 2: **Output:** Refined models  $\{f_{\widetilde{\theta}_i}\}_{i=1}^N$ .

```
3: ■ (Optional) Partition the problem in a number of non-intersecting clusters.
 4: \{C_k\}_{k=1}^K \leftarrow \text{findClusters}(\{\mathbf{x}_i\}_{i=1}^N, K)
5: \{\mathbf{D}_k\}_{k=1}^K \leftarrow \text{computeDistanceMatrices}(\{C_k\}_{k=1}^K, \text{dist}())
  6: Process each cluster independently.
  7: for each cluster C do
        \mathcal{T} \leftarrow \text{computeMST}(\mathbf{D})
                                                                                                                  > Extract the MST with Kruskal's algorithm
         \mathbf{p} \leftarrow \text{findPrototype}(C, \mathcal{T})
                                                                                                                         \triangleright Compute a cluster prototype within \mathcal{T}
         \mathbf{d} \leftarrow \text{extractTreeWeights}(\mathcal{T})
                                                                                                                              \triangleright Extract the distance vector from \mathcal{T}
11: \mathbf{d}' \leftarrow \text{softmax}(\mathbf{d})
                                                                                                               Normalize the distance vector using softmax
       f_{\widehat{\boldsymbol{\theta}}} \leftarrow \mathcal{A}_0(f_{\boldsymbol{\theta}}, \mathbf{p}, b)
                                                                                              \triangleright Train the prototype model on cluster data with budget b
      for i=1 to |\mathbf{d}'| do
            (\mathbf{b})_i \leftarrow \lceil (\mathbf{d}')_i \cdot B \rceil
                                                                                                                                        15:
          end for
          ■ Refine individual models using allocated budgets.
          Q \leftarrow \emptyset, Q \leftarrow \text{enqueue}(Q, \mathbf{p})
                                                                                             ▶ Initialize a queue with the prototype node for refinement
         while \neg is Empty(Q) do
            \mathbf{x}_{parent} \leftarrow \text{dequeue}(Q)
20:
             for \mathbf{x}_{child} in childrenOf(\mathbf{x}_{parent}) do
                 Q \leftarrow \text{enqueue}(Q, \mathbf{x}_{child})
                                                                                                    > Add the children of the processed node in the queue
21:
                f_{\widetilde{\boldsymbol{	heta}}_{child}} \leftarrow \mathcal{A}_1(f_{\widehat{oldsymbol{	heta}}_{parent}}, \mathbf{x}_{child}, \mathbf{x}_{parent}, (\mathbf{b})_{child})
                                                                                                 > Refine child model using parent model, child data and
22:
23:
              end for
24:
         end while
25: end for
26: return \{f_{\widetilde{\boldsymbol{\theta}}_i}\}_{i=1}^N.
```