Anomaly detection on graphs

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• Problem

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Anomaly detection in graphs

We are trying to solve **unsupervised anomaly detection** on graphs:

- **Our input** is a graph where the nodes have a given set of features.
- **Our output** can either be a *binary labelling* or a *ranking* of the nodes.

The output should reflect some sense of **anomalousness** of the nodes.

Why is this problem interesting?

- Relates to **bot identification** in social networks.
- Can help improve **spam detection**.
- Relates to various types of **fraud**: *insurance*, *healthcare*, *financial*, etc.

→ A good anomaly detection method could also help other graph-based tasks, such as classification or clustering.

Defining anomalies

Defining anomalies on graphs is a **difficult task** since nodes have **two types of information**:

- Personal information, given by their **features**.
- Community information, given by their **neighbours**.

In that context, anomalies can have several forms:

- Feature-based anomalies: an anomaly can be defined uniquely based on its features.
- **Graph-based anomalies:** an anomaly can be defined uniquely *based on its neighbours* and the general graph structure.
- **Combined anomalies:** an anomaly can be defined not in the two previous ways, but by considering *both its features and the general graph structure*.





















































Previous work

The literature in anomaly detection on graphs is very **diverse** and **sparse**.

We did not find **direct baselines** and had to base our work on the following articles.

- Multi-scale Anomaly Detection on Attribute Networks, L. Gutiérrez-Goméz, A. Bovet, J.-C. Delvenne, 2019
- Bayesian Robust Attributed Graph Clustering: Joint Learning of Partial Anomalies and Group Structure, A. Bojchevski, S. Günnemann, 2018
- Maximum Entropy Generators for Energy-Based Models, R. Kumar, S. Ozair, A. Goyal, A. Courville, Y. Bengio, 2019

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General approach

Since we are using an **unsupervised method**, we do not have access to labels or examples of anomalies.

This is a realistic approach since:

- Anomalies are **rare**.
- Anomalies might be **very diverse**.

→ The only assumption we can make is that most nodes are normal.

With these constraints, the **general approach** to anomaly detection is to:

- Define some **energy** of the nodes, related to their *likelihood of normality*.
- Use this energy to identify anomalies:
 - by **ranking** the nodes;
 - by considering the nodes with highest energy;
 - by identifying a **threshold** in the energy distribution.

Our method

The idea:

- We will **reproduce** the original features of the nodes with their expected values.
- We first **hide/mask** some of the features and then try to recreate them using the rest of the information.
- Use *Graph Neural Networks* (**GNNs**) for their high representative power.
- **Compare** the reproduced features with the original ones to identify anomalies.
- → Creation of two algorithms: HideGNN and MaskGNN.

The general method:

- Take the graph as input, with node features.
- Use one of our two algorithms (HideGNN or MaskGNN) to reproduce the node features.
- Use these reproduced features to **augment** the original dataset.
- Apply an **anomaly detection algorithm** (ADA) on the augmented dataset.

Our method



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GNNs

GNNs are **commonly used** in machine learning methods on graphs.

They are based on a **message-passing** algorithm and have the advantage to:

- use **both** the node features and the graph structure;
- and **not depend** on the ordering of the nodes.



HideGNN

The method:

- Start by **hiding** a specific feature.
- Use a GNN on the **other features** to guess the hidden one.
- **Repeat** over all features.
- Eventually obtain **a set of GNNs** able to reproduce the original features.



MaskGNN

The method:

- Start by creating a **random mask** over all features.
- **Replace** the masked features by some value (usually 0).
- Use a GNN to reproduce all features but **only optimize on the masked features**.
- **Resample** the random mask.
- Eventually obtain a **non-trivial GNN** which reproduces the original features.



Pros and Cons



- Pros:
 - Complete reproduction of the features.
 - Less parameters.
- Cons:
 - Heavier to train.
 - Do not use any information from the hidden feature.



- Pros:
 - Faster to train and using a single GNN.
 - Elegant all-in-one tool.
- Cons:
 - More parameters.
 - If not properly parametrized, likely to miss some information.
 - Less stable.

Remarks

On our algorithms:

- They are theoretically able to **overfit** and recreate the original features of the dataset.
- They learn from both the graph structure and the features, and their output contains a **lot of information** from the graph.
- They use a novel method combining GNN with methods from **NLP**, which opens the door to new parallel between the two fields.

On our method:

- It allows for **non graph-based algorithms** to be efficiently applied on graphs.
- It could easily be used on **other tasks**.

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Datasets

We tested our models on three datasets:

- **Disney:** 124 nodes and 6 anomalies.
- Books: 1418 nodes and 28 anomalies.
- Enron: 13533 nodes and 5 anomalies.

The inherent **extreme statistics** of these datasets makes this task complicated.

→ Only Books allows an acceptable split into train/valid/test.

	Disney	Books	Enron	
Nodes	124	1418	13533	
Edges	335	3695	176987	
Features	32	21	18	
Outliers	6	28	5	

Experimental setup

- Since only Books can be split, we use this dataset to **optimize** our algorithms.
- Once we finish running the algorithms, we use parameters with the best results on valid of Books.
- We then compute and **report** the results on the test set of Books and on Disney and Enron.

→ We compare the ADAs with or without using our algorithms.

Some problems we encountered:

- None of these datasets were **already split** before our experiments.
- The results previously reported seem to be using the **test set** for hyperparameter tuning.
- We re-computed these results using our framework and found great discrepancies.
 We also did not always manage to recreate their experiments.
- These previous methods also seemed to be using a **modified version** of the datasets, with less features.

Our results

	Books		Disney*		Enron*	
	Original	Reduced	Original	Reduced	Original	Reduced
OCSVM [†]	35.90 [‡]	37.33 [‡]	36.44 [‡]	85.88 [‡]	55.90 [‡]	43.22 [‡]
Hide + OCSVM	49.58 (±1.47)	52.21 (±2.41)	46.97 (±2.26)	$43.16(\pm 1.03)$	32.76 (±3.79)	51.84 (±3.35)
Mask + OCSVM	55.06 (±1.64)	48.49 (±2.07)	47.63 (±1.73)	43.07 (±6.30)	56.28 (±9.73)	50.46 (±6.55)
MEG [†] [24]	63.12 [‡]	56.60 [‡]	50.28 [‡]	39.55 [‡]	26.47^{\ddagger}	63.08 [‡]
Hide + MEG	$60.42 (\pm 7.03)$	65.36 (±3.26)	$46.51 (\pm 6.51)$	69.44 (±3.47)	44.22 (±14.05)	40.12 (±6.54)
Mask + MEG	57.06 (±6.55)	62.17 (±4.67)	51.25 (±3.46)	62.44 (±13.38)	42.36 (±16.82)	40.63 (±12.21)
MADAN [13]	48.80^{\ddagger}	45.54^{\ddagger}	67.51 [‡]	19.49 [‡]	61.81 [‡]	73.64 [‡]
Hide + MADAN	57.21 (±9.26)	52.08 (±7.64)	59.09 (±15.03)	49.82 (±15.02)	54.24 (±14.41)	47.47 (±13.33)
Mask + MADAN	58.43 (±8.87)	52.34 (±7.89)	55.98 (±14.07)	48.87 (±12.54)	60.55 (±14.58)	50.20 (±14.13)
PAICAN [2]	67.76 [‡]	55.38 [‡]	75.14 [‡]	73.45 [‡]	26.94^{\ddagger}	57.77 [‡]
Hide + PAICAN	68.23 (±1.45)	72.82 (±3.38)	62.73 (±8.14)	71.92 (±1.15)	54.59 (±3.31)	$58.65(\pm 10.07)$
Mask + PAICAN	68.90 (±1.15)	61.25 (±3.43)	67.31 (±6.87)	64.76 (±10.82)	$45.92(\pm 8.08)$	65.40 (±5.46)

A synthetic dataset (Thanksgiving)

To try and understand our algorithms and results better, we created a **synthetic dataset**:

- Composed of two communities;
- Each community having a **given set of features**.
- We then modify the features of some members of one community to have the **features of the other community**.

→ This dataset only has combined anomalies and requires to use both the features and the graph structure.



Results on the synthetic dataset

Modifying the difference of **mean** between the two community features:



Modifying the **number of connections** between the two groups:



Thank you!