# Detection of Duplicates Among Non-structured Data From Different Data Sources

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#### Plan



- 1 Introduction
- 2 Databases
  - Registre des Entreprises du Québec (REQ)
  - Private Dataset
- 3 Similarity Between Two Entities
  - Similarity Algorithm
  - Machine Learning
- 4 Conclusion

#### Situation



When assessing a commercial risk, an insurer needs to gather various information about the risk. This long and complex process implies numerous questions. Thus an insurer is prompt to use an external source to help reduce the number of questions.

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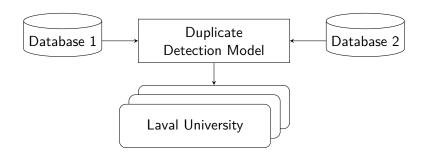
Thus, we want to detect duplicate of commercial risk in another data source using as little as possible information[?].

#### Example

David Beauchemin, the owner of "Beauchemin inc.", calls for insurance. Using minimal information, we want to retrieve as much as possible from an external source to ask him as less than the necessary number of questions.

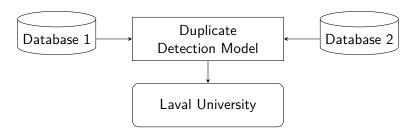
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Moreover, we consider only the top 1 document as the potential candidate.

# How do we detect duplicate?



To detect duplicate we need the following [?]

- databases (at least two) (section 2),
- a way to determine the similarity between two documents (section 3).

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# Registre des Entreprises du Québec (REQ)



- $\sim$  3.5 millions entries [?]
  - Names
  - Address
  - ► Economic activities
  - Administrative informations

#### Private Dataset



- 21,444 enterprises
  - Name
  - Address
  - Economic activity



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- 11,649 commercial risk are in the province of Quebec.
  - ▶ 1706 were annotated
    - ★ 1418 (commercial risk, REQ entity)
    - ★ 288 (commercial risk, None)
- We only use the name and the address.

#### Name



We have used two versions of the name.

Normalize name (NN): lowercase, whitespace and accent trimming.

#### Normalize Name

L'Université Laval ⇒ *l'universite laval* 



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2 No stop words name (NSWN): stop words trimming (i.e *le*, *la*, *de*) [?].

#### No Stop Word Name

l'universite laval ⇒ 'universite laval

#### Address



We also use two versions of the address.

Complete Address Normalize (NA): same as the name.

#### Complete Address Normalize

2325 rue de l'Université, Québec, QC, G1V 0A6

2325 rue de l'universite, quebec, qc, g1v 0a6



We also use two versions of the address.

**1** Complete Address Normalize (NA) : same as the name.

#### Complete Address Normalize

2325 rue de l'Université, Québec, QC, G1V 0A6 ↓

2325 rue de l'universite, quebec, qc, g1v 0a6

2 Address components (AC) : parsed and grouped by address components [?].

#### Address components

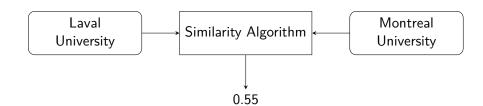
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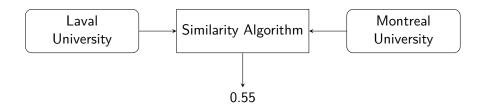
# Similarity Algorithm





# Similarity Algorithm





Their similarity ranks the documents, and we select the best one as the duplicate candidate.

### Similarity Algorithm



Similarities algorithms are one that measures the resemblance between two string base on the distance between their tokens.

10 different algorithms were used.

### Similarity Algorithms



### **Jaccard Similarity**

$$\mathsf{Jaccard}(\mathsf{A},\,\mathsf{B}) = \begin{cases} 0 & \text{if } |\mathsf{A}\cap\mathsf{B}| = 0 \text{ or } |\mathsf{A}\cup\mathsf{B}| = 0 \\ \frac{|\mathsf{A}\cap\mathsf{B}|}{|\mathsf{A}\cup\mathsf{B}|} & \text{otherwise} \end{cases}$$

#### Example

Jaccard(Laval University, Montreal University)

$$\Downarrow$$

$$\frac{|\{\mathsf{University}\}|}{|\{\mathsf{University},\,\mathsf{Montreal},\,\mathsf{Laval}\}|}=\frac{1}{3}=\mathbf{0.333}$$

# Similarity Algorithms



### String to String (StoS)

$$\mathsf{StoS}(\mathsf{A},\,\mathsf{B}) = \begin{cases} 1 & \text{if } \mathsf{A}_i = \mathsf{B}_j \,\,\forall i,j \\ 0 & \text{otherwise} \end{cases}$$

#### Example

 $\begin{aligned} \mathsf{StoS}(\mathsf{Laval\ University},\ \mathsf{Montreal\ University}) &\Rightarrow 0 \\ \mathsf{StoS}(\mathsf{Laval\ University},\ \mathsf{Laval\ University}) &\Rightarrow 1 \end{aligned}$ 

# Similarity Algorithms



#### Jaro-Winkler

$$\mathsf{Jaro-Winkler}(\mathsf{A},\,\mathsf{B}) = \mathsf{Jaro}(\mathsf{A},\,\mathsf{B}) + \frac{\min{(P,4)}}{10} \times (1 - \mathsf{Jaro}(\mathsf{A},\,\mathsf{B}))$$

### Example

$$\begin{aligned} & \mathsf{Jaro-Winkler}(\mathsf{David},\,\mathsf{Daniel}) = \\ & \mathsf{Jaro}(\mathsf{David},\,\mathsf{Daniel}) + \tfrac{2}{10} \times (1 - \mathsf{Jaro}(\mathsf{David},\,\mathsf{Daniel})) \\ & \qquad \qquad \downarrow \\ & 0.7 + \tfrac{2}{10} \times 0.7 = 0.84 \end{aligned}$$

# Why Similarity Algorithm?



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- Give good results

# Name Interesting Results <sup>1</sup>



	(%)	StoS	Jaro	Jaro-Winkler	Jaccard
NN	Accuracy	41.47	63.40	63.47	65.73
NSWN	Accuracy	44.15	64.46	65.23	66.50

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- Removing stop words (second line) gives the best results.
- The prefix similarity used by Jaro-Winkler improved results (more when using NSWN).



<sup>1.</sup> Positive examples only.



	(%)	StoS	Jaro	Jaro-Winkler	CSS
CAN	Accuracy	0.00	48.03	48.10	45.91
AC	Accuracy	13.19	51.83	51.83	52.19

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- Using the address components without considering the order of the components improved results (AC).
- The local similarity used by Jaro-Winkler did not improve results this time since an address is rarely defined by is prefix.
- $\blacksquare$  CSS gets the best results (14% below the previous best results using Jaccard).



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## Missing Addresses



Since 16% (270) of the pair (commercial risk, REQ entity) are missing a address, the results are under-evaluated. For example, the CSS accuracy without those pairs is at 62.74% near 10% higher.

Those missing addresses are due to the confidentiality policy of the REQ [?].



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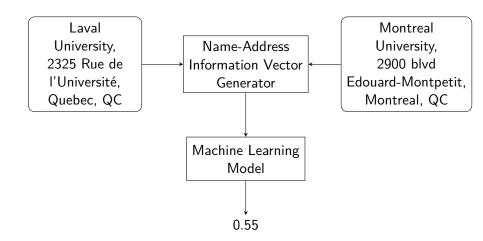
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- Using the NSWN or the AC helps improve the results over the normalized name or address.
- The missing addresses pull down the results.
- To use positive and negative examples, we need to use a decision function. Such as a similarity threshold, were similarity below the threshold are rejected (the results are not shown).

## Machine Learning Similarity Algorithm









We used the previous similarity algorithm to generate an information vector between two entities using the NSWN and the CA.

Example of an information vector						
StoS	Levenshtein	Jaro-Winkler	LCSP	Jaccard	Cosinus	-
0.00	0.15	0.25	0.35	0.15	0.15	-
StoS	Levenshtein	Jaro	LCSP	Jaccard	Cosinus	CSS
0.00	0.16	0.55	0.15	0.45	0.37	0.48

# Machine Learning Algorithm



- Logistic regression
- 2 Random Forest
- Multilayer Perceptron

# Why Machine Learning?



- Allow us to use the name and the address simultaneously
- Generalization capability

### Models' Training Steps



- Data preprocessing
- 2 Hyperparameters grid search
- Model training
- 4 Evaluation of the trained model for the duplicate detection task



■ 80-20 splitting of the data (1706 annoted exemples).



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- To balance the dataset (1364 positives vs. 241 negatives), we randomly select the first name and address of a REQ entity to create a fake commercial risk and randomly pair it with another REQ entity. Resulting in a training dataset of 2246 (commercial risk, REQ entity) pair.



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- We fit a standard scaler on the training dataset. With that standard scaler, we applied a transformation over all the vectors (train and test).

# Hyperparameters Grid Search & Model Training



- A grid search for the logistic regression (*C* and tolerance) and the random forest (number of estimators).
- A random search for the multilayer perceptron (number of layers and neurons and the tolerance).
- Cross-validation approach using a 5-folds.

After the grid search, we retrain using the best parameters.

#### **Evaluation**



We evaluate our three trained models against the best configuration, Jaccard using the name, but reevaluated with the validation dataset (20% of the 1706 annotated examples).

We evaluated the algorithm with positives and negatives examples (for recall and precision). The decision function is a similarity threshold where a similarity below a numerical threshold (e.g., 0.7) is rejected.

We aim to maximize the recall since our objective is to detect the more duplicate as possible since we can, later on, validate the duplicate manually.



(%)	Logistic Regression	Random Forest	Multilayer perceptron	Jaccard
Precision	89,77	81,06	87,55	81,78
Recall Accuracy	66,67 64,91	73,54 62,87	79,73 73,10	72,51 62,87

 $\blacksquare$  The random forest and the multilayer perceptron achieve better recall than Jaccard. The best being the perceptron with near 80% recall.



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Accuracy	64,91	62,87	73,10	62,87

- $\blacksquare$  The random forest and the multilayer perceptron achieve better recall than Jaccard. The best being the perceptron with near 80% recall.
- The logistic regression achieves the lowest result even if the Jaccard similarity is used in the generation of the information vector.

### Example of an Error



Commercial Risk	Entity		
construction alain cloutier inc.	construction steeve arbourd inc.		
1030 rue de l'ardoise sherbrooke j1c 0j6	2-1822 rue notre-dame, l'ancienne-lorette, g2e 3c7		

The following error was only generated with the logistic regression model. The algorithm matched the two with a similarity of 0.99934, even if the annotated duplicate appear with the same name and a slightly different address (similarity of 0.998).

### Improving the Results



The previous error highlights a problem of our approach; so far, we have tried to match with **the** most similar, and the generated similarity are close to each other, making it restrictive to use only the best similarity.

Since our approach aims to prefill the information of an insurance application, we can return more than one and use a human validation to select the best one from N possibilities.

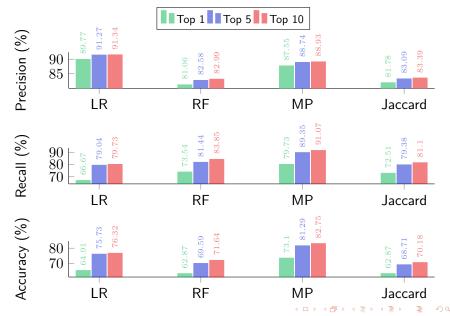
#### N Most Similar



We consider a matching is good when the pair (commercial risk, REQ entity) is included in the N most similar.

#### Results





#### Results



- lacksquare Using a top N approach greatly improved the results.
- Using N=10, we can achieve a near max recall at 91% with the multilayer perceptron (max of 93% with the indexing).

### Inference Times



(second)	Logistic Regression	Random Forest	Multilayer Perceptron	Jaccard
Time	1,32	1,74	1,34	0,25

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- $\blacksquare$  Using a top N approach helps achieve better results when N is greater than 1.
- Inference times (of machine learning models) are similar to using only a similarity algorithm.

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#### Conclusion



- We have shown that using a similarity algorithm can achieve good results.
- Uses of machine learning algorithm (such as multilayer perceptron) can achieve greater results.
- lacktriangle Using a N most similar approach, where N is greater than one, help improved the results, achieving almost the max recall value.

#### **Future Works**



- Word embeddings [?, ?, ?, ?, ?]
- Siamese Network [?, ?]
- Uses of spatial data [?]
- Removal of more specitif stop words using a TF-IDF approach [?]

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### References I

