



Department of Electrical Engineering, Computer Science and Mathematics Warburger Straße 100 33098 Paderborn



Master Thesis

Learning to Aggregate on Structured Data

Clemens Damke

1. Reviewer Prof. Dr. Eyke Hüllermeier

Intelligent Systems and Machine Learning Group (ISG)

Paderborn University

2. Reviewer Prof. Dr. Axel-Cyrille Ngonga Ngomo

Data Science Group (DICE) Paderborn University

November 15, 2019

Clemens Damke

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Reviewers: Prof. Dr. Eyke Hüllermeier and Prof. Dr. Axel-Cyrille Ngonga Ngomo

Supervisor: Vitalik Melnikov

Paderborn University

Intelligent Systems and Machine Learning Group (ISG)

Heinz Nixdorf Institute

Department of Electrical Engineering, Computer Science and Mathematics

Warburger Straße 100

33098 Paderborn

Abstract

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Introduction

1.1 Motivation

The field of *machine learning* (ML) on graph-structured data has applications in many domains due to the general expressive power of graphs. Three common types of graph ML problems are

- **1. Link prediction:** A graph with an incomplete edge set is given and the missing edges have to be predicted. The generation of friendship suggestions in a social network is a typical example for this.
- **2. Vertex classification & regression:** Here a class or a score has to be predicted for each vertex of a graph. In social graphs this corresponds to the prediction of properties of individuals. Another example is the prediction of the amount of traffic at the intersections of a street network.
- **3. Graph classification & regression:** In this final problem type a single global class or continuous value has to be predicted for an input graph. The canonical example for this is the prediction of properties of molecule graphs, e.g. the toxicity or solubility of a chemical.

In this thesis we will focus on the last problem type, GC/GR. A ML method for this problem has to accept graphs of varying size and should be permutation invariant wrt. the vertices. Those requirements are not met by the commonly used learners that only accept fixed-size feature vectors as their input, e.g. logistic regression models, support vector machines or multilayer perceptrons.

A GC/GR method has to account for two central aspects of the problem: 1. Local structural analysis and 2. global aggregation. The first aspect is about the extraction of relevant features of substructures of the input graph. The latter is about the way in which the local features are combined into a final class or regression value. The existing GC/GR methods are mostly motivated by local structural graph analysis. The aspect of global aggregation on the other hand is less emphasized by those methods.

There is however a separate branch of research that specifically looks at the problem of learning aggregation functions, called *learning to aggregate* (LTA). Current LTA approaches explicitly learn an aggregation functions for sets which can be interpreted

as graphs without edges. The motivation for this thesis is to generalize LTA from sets to arbitrary graphs. The overall goal is to combine the aggregation learning perspective with existing GC/GR methods.

1.2 Goals

To extend LTA to graphs, three goals have to be achieved:

1. Formalization of LTA: Before LTA can be extended, its essential characteristics have to be defined. Those characteristics should provide the terminology to formally capture the differences and similarities between LTA and existing

GC/GR methods.

2. Give an LTA interpretation of GC/GR methods: Using the LTA formalization,

representative GC/GR approaches should be restated as LTA instances. Currently there is no comprehensive formulation of the relation between both fields

of research; this is addressed by the the second goal.

3. Define an LTA method for graphs: Using the LTA perspective on GC/GR,

hidden assumptions of the existing approaches should become clear and in which way they share the assumptions of LTA. The last goal is to use those

insights to formulate an LTA-GC/GR method that combines ideas from the

existing approaches with the LTA assumptions.

1.3 Structure

Chapter 2: Related Work

Chapter 3: Learning to Aggregate on Graphs

Chapter 4: Evaluation

Chapter 5: Conclusion

Related Work

Before combining LTA and GC/GR as described in section 1.2, we first give an overview of the state-of-the-art in both fields of research. First the existing LTA methods for unstructured set inputs are described. Then we will look at the current approaches to tackle the GC/GR problem.

2.1 Learning to Aggregate

The class of LTA problems was first described by Melnikov and Hüllermeier [MH16]. There an input instance is understood as a composition c of so-called constituents $c_i \in c$, i.e. as a variable-size multiset with n = |c|. The assumption in LTA problems is that for all constituents c_i a local score $y_i \in \mathcal{Y}$ is either given or computable. The set of those local scores should be indicative of the overall score $y \in \mathcal{Y}$ of the composition c. LTA problems typically require two subproblems to be solved:

- **1. Aggregation:** A variadic aggregation function $\mathcal{A}: \mathcal{Y}^* \to \mathcal{Y}$ that estimates composite scores has to be learned, i.e. $y_i \approx \hat{y} = \mathcal{A}(y_1, \dots, y_n)$. Typically the aggregation function \mathcal{A} should be associative and commutative to fit with the multiset-structure of compositions.
- **2. Disaggregation:** In case the constituent scores y_i are not given, they have to be derived from a constituent representation, e.g. a vector $x_i \in \mathcal{X}$. To learn this derivation function $f: \mathcal{X} \to \mathcal{Y}$, only the constituent vectors $\{x_i\}_{i=1}^n$ and the composite score y is given. Thus the constituent scores y_i need to be disaggregated from y in order to learn f.

Overall LTA can be understood as the joint problem of learning the aggregation function A and the local score derivation function f. Two main approaches to

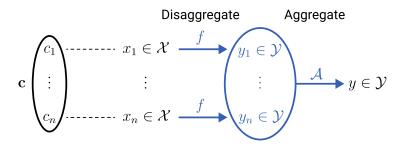


Figure 2.1. Overview of the structure of LTA for multiset compositions.

represent the aggregation function in LTA problems have been explored.

2.1.1 Uninorm-Aggregation

The first approach uses uninorms [MH16] to do so. There the basic idea is to express composite scores as fuzzy truth assignments $y \in [0,1]$. Such a composite assignment y is modeled as the result of a parameterized logical expression of constituent assignments $y_i \in [0,1]$. As the logical expression that thus effectively aggregates the constituents, a uninorm U_{λ} is used. Depending on the parameter $\lambda \in [0,1]$, U_{λ} combines a t-norm T and a t-conorm S which are continuous generalizations of logical conjunction and disjunction respectively. One popular choice of norms are the so-called Łukasiewicz norms:

t-norm
$$T(a,b) := \max\{0, a+b-1\}$$
, t-conorm $S(a,b) := \min\{a+b,1\}$,

uninorm $U_{\lambda}(a,b) := \begin{cases} \lambda T\left(\frac{a}{\lambda}, \frac{b}{\lambda}\right) & \text{if } a,b \in [0,\lambda] \\ \lambda + (1-\lambda)S\left(\frac{a-\lambda}{1-\lambda}, \frac{b-\lambda}{1-\lambda}\right) & \text{if } a,b \in [\lambda,1] \\ \lambda \min\{a,b\} & \text{else} \end{cases}$ (2.1)

At the extreme points (0,0), (0,1), (1,0) and (1,1), T and S coincide with the Boolean operators \wedge and \vee ; the values at all other points are interpolated as shown in fig. 2.2. The uninorm U_{λ} uses the conjunctive t-norm T for values below the threshold λ and the disjunctive t-conorm S for values above the threshold. U_{λ} therefore smoothly interpolates between a conjunctive and disjunctive operator with the extreme points $U_1 = T$ and $U_0 = S$.

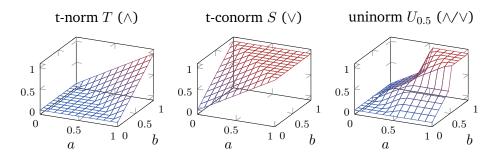


Figure 2.2. The Łukasiewicz norms and the corresponding uninorm for $\lambda=0.5$.

Since t-norms and t-conorms are commutative and associative they can also be applied to non-empty sets of arbitrary size, i.e. $T(\{y_1,\ldots,y_n\})=T(y_1,T(\{y_2,\ldots,y_n\}))$ with fixpoint $T(\{y\})=y$. Using this extension, a uninorm U_λ can be applied to sets which turns it into a parameterized aggregation function $\mathcal{A}_\lambda:[0,1]^*\to[0,1]$. In this simple model the LTA aggregation problem boils down to the optimization of λ . The LTA disaggregation problem is solved by jointly optimizing a logistic regression model, i.e. the constituent scores $\{y_i\in[0,1]\}_{c_i\in c}$ are described by $y_i=\left(1+\exp(-\theta^\top x_i)\right)^{-1}$.

Overall an LTA model is therefore described by the uninorm parameter λ and the regression coefficients θ .

2.1.2 OWA-Aggregation

Recently Melnikov and Hüllermeier [MH19] have looked at an alternative class of aggregation functions. Instead of using fuzzy logic to describe score aggregation, ordered weighted average (OWA) operators were used. OWA aggregators work by sorting the input scores and then weighting them based on their sort position, i.e.

$$\mathcal{A}_{\lambda}(y_1, \dots, y_n) := \sum_{i=1}^n \lambda_i y_{\pi(i)}, \tag{2.2}$$

where $\lambda \in \mathbb{R}^n$ is a weight vector with $\|\lambda\|_1 = 1$ and π is a sorting permutation of the input scores with $y_i < y_j \Rightarrow \pi(i) < \pi(j)$. Depending on the choice of the vector λ , the OWA function \mathcal{A}_{λ} can express common aggregation functions like \min (if $\lambda = (1,0,\ldots,0)$), \max (if $\lambda = (0,\ldots,0,1)$) or the arithmetic mean (if $\lambda = \left(\frac{1}{n},\ldots,\frac{1}{n}\right)$).

To deal with varying composition sizes n, the weights $\{\lambda_i\}_{i=1}^n$ can however not be statically assigned. Instead they are interpolated using a so-called *basic unit interval monotone* (BUM) function $q:[0,1]\to[0,1]$. It takes constituent positions that are normalized to the unit interval, i.e. $\frac{i}{n}\in[0,1]$. The BUM function q is then used to interpolate a weight for any normalized sort position via $\lambda_i:=q\left(\frac{i}{n}\right)-q\left(\frac{i-1}{n}\right)$. Because q is monotone with q(0)=0 and q(1)=1, it always holds that $\|\lambda\|_1=q(1)-q(0)=1$. Using this model, the aggregation problem boils down to optimizing the shape of q.

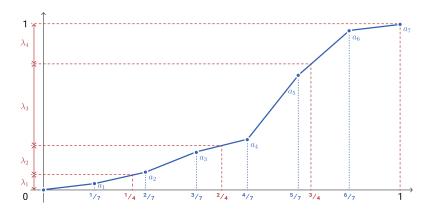


Figure 2.3. Illustration of how *a* describes *q* and its relation to λ (n = 4, m = 7).

In the OWA approach the BUM function q is modeled as a piecewise linear spline. This spline is described by m+1 points $\left\{\left(\frac{j}{m},a_j\right)\right\}_{j=0}^m$, the so-called knots of the spline. The curve of q is obtained by linearly interpolating between neighboring knots as shown in fig. 2.3. If $0=a_0\leq a_1\leq\cdots\leq a_m=1$, q is a BUM function. The LTA

aggregation problem is therefore solved by optimizing $a \in \mathbb{R}^{m+1}$ under this constraint. The disaggregation problem is tackled by adding the scores $y_1, \ldots, y_M \in \mathbb{R}$ to the learnable parameters of the model where M is assumed to be the finite number of constituents. Currently the OWA approach requires all possible constituents to be part of the training dataset since it does not consider constituent features $x_i \in \mathcal{X}$ to predict the scores of previously unseen constituents.

2.2 Graph Regression and Classification

The existing approaches to tackle the GC/GR problem can be categorized into three main families: 1. Explicit graph embeddings, 2. Graph kernels and 3. Graph neural networks. We will now look at the characteristics of those families and give an overview of specific methods.

2.2.1 Explicit Graph Embeddings

The basic idea of graph embedding approaches is to map a graph G to some finite vector space \mathbb{R}^d . By embedding a graph any classification or regression algorithm that works with vectors can then be applied. There are three main types of graph embedding approaches.

Fingerprint embeddings

Unsupervised embeddings

Supervised embeddings

2.2.2 Graph Kernels

2.2.3 Graph Neural Networks

Spatial GNNs

Spectral GNNs

Learning to Aggregate on Graphs

3

- 3.1 Formalization of LTA Characteristics
- 3.2 An LTA Interpretation of GC/GR Methods
- 3.3 LTA on Dynamically Decomposed Graphs

Evaluation 4

Conclusion

- 5.1 Review
- 5.2 Future Directions

Appendix

Bibliography

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Erklärung zur Masterarbeit

Ich, Clemens Damke (Matrikel-Nr. 7011488), versichere, dass ich die Masterarbeit mit dem Thema *Learning to Aggregate on Structured Data* selbstständig verfasst und keine anderen als die angegebenen Quellen und Hilfsmittel benutzt habe. Die Stellen der Arbeit, die ich anderen Werken dem Wortlaut oder dem Sinn nach entnommen habe, wurden in jedem Fall unter Angabe der Quellen der Entlehnung kenntlich gemacht. Das Gleiche gilt auch für Tabellen, Skizzen, Zeichnungen, bildliche Darstellungen usw. Die Masterarbeit habe ich nicht, auch nicht auszugsweise, für eine andere abgeschlossene Prüfung angefertigt. Auf § 63 Abs. 5 HZG wird hingewiesen.

Paderborn, 15. November 2019	
	Clemens Damke