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## Streszczenie

### Słowa kluczowe

graph queries, Datalog

## Dziedzina pracy (kody wg programu Socrates-Erasmus)

11.3 Informatyka

## Klasyfikacja tematyczna

D. SoftwareD.127. BlabalgorithmsD.127.6. Numerical blabalysis

## Tytuł pracy w języku angielskim

Processing large graphs with Apache Giraph

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## Introduction

In recent years, the humanity has created many graph datasets much larger than those avaiable ever before. Those graphs became a very popular object of research. Most notable examples include the Web graph – a graph of internet websites and links between them, all kinds of social networks. There are also graphs such as transportation routes, similarity of newspaper of scientific articles or citations among them.

With increasing computational power and memory space, we can expect more and more real-life graphs to became subject to computation. We can also expect the existing graphs, such as the Web or social networks, to grow in all aspects.

The graps metioned can be a source of a huge amount of useful information. Hence, there is an increasing number of practical computational problems. Some of the analyses carried out are ranking of the graph nodes (ex. importance of a Web page, determining most influential users in given group of people), clustering (detecting communities), computing metrics for the whole graph or some parts of it (ex. connectivity measures) and connection predictions. Usually, such analyses are built on top of standard graph algorithms, such as PageRank-like procedures, shortest paths or connected components.

In the past, we have seen many tools for efficient distributed large dataset computations, starting from Google's MapReduce [5] and its widely used open source counterpart, Apache's Hadoop, as well as higher-level languages such as PigLatin and Hive. However, those are not well suited for graph computations, as they do not support iteration well.

Recently, there is an outbreak of frameworks and languages for large graphs processing, including industrial systems such as Google's Pregel [6], its open-source version Apache Giraph and Spark GraphX, Giraph++ [7].

The frameworks currently avaiable allow you to implement a graph algorithm in a specified model (for example Pregel's "think like a vertex"). On the other hand, query languages, such as SQL, are a bad fit for graph data since because of limited support of iteration. With the rise of graph computational problems, we need an easier way to extract information from graphs: a query language for effectively expressing data queries typical for graphs.

The Socialite [1] [2] language is one of the most interesting propositions. It is based on a classical language – Datalog. Declatarive semantics makes it easy to distribute the computations, since no execution flow is embedded in the program code. It also gives many possibilities for optimizations and approximate evaluation. At the same time, support for recursion is crucial, since most graph algorithms have iterative nature. However, most of practical graph algorithms can not be expressed efficiently in Datalog because of the language limitiations. With a few extensions to original Datalog, most important of which is recursive aggregation, SociaLite makes it easy to write intuitive programs which can be executed very efficiently.

Unfortunately, there is no solid implementation of SociaLite available. The interpreter published by the authors is undocumented and contains many bugs. It is hard to imagine the language being implemented in the industry in the forseeable future. The papers [1] and [2]

which introduced SociaLite contain certain simplifications and are not specific about some important details in definitions and proofs.

The goal of this work is to bridge the gap between the theoretical idea for SociaLite and a practical implementation and to draw a path towards its usage in the industry. We show how to translate SociaLite declarative programs into Giraph "think-like-a-vertex" programs and introduce a compiler that enables SociaLite programs to be executed on existing Giraph infrastructure. This allows users of Hadoop to write and execute SociaLite programs without any additional effort to build a dedicated server infrastructure for that.

The work consists of six chapters. In 1 we recall definitions of Datalog and its evaluation methods while 2 contains an introduction to the Pregel computation model. In 3 we describe the extensions introduced by SociaLite and provide formal definitions and general-case proofs which the original papers lack. Chapter 4 shows the translation procedure from SociaLite to Giraph programs implemented in the S2G compiler, which is described in 5. In 6 we sketch the possible future work and the path to industrial implementation of the language using the S2G compiler.

## **Datalog**

In this chapter we define the Datalog language.

### 1.1. Definitions

@TODO: te definicje są roboczo przepisane z [4], którzy z kolei podają je za [3].

#### 1.1.1. Queries and instances

Let us assume that **dom** is an infinite set of data values.

- A database schema  $\sigma$  is a collection of relation names R where every R has arity  $ar(R) \ge 0$ .
- We call  $R(\overline{d})$  a fact when R is a relation name and  $\overline{d}$  is a tuple in **dom**.
- We say that a fact  $R(d_1, \ldots, d_k)$  is over a database schema  $\sigma$  if  $R \in \sigma$  and ar(R) = k.
- A (database) instance I over  $\sigma$  is a finite set of facts over  $\sigma$ .
- We denote by adom(I) the set of all values that occur in facts of I. When  $I = \mathbf{f}$ , we simply write  $adom(\mathbf{f})$  rather than  $adom(\{\mathbf{f}\})$ .
- By |I| we denote the number of facts in I.

#### 1.1.2. Datalog with negation

Let **var** be the universe of variables, disjoint from **dom**. An atom is of the form  $R(u_1, \ldots, u_k)$  where R is a relation name and each  $u_i \in \mathbf{var}$ . We call R the predicate. A *literal* is an atom (positive atom) or a negated atom (negative atom)

We recall Datalog with negation [3], abbreviated Datalog<sup>¬</sup>. Formally, a Datalog<sup>¬</sup> rule  $\phi$  is a quadruple  $(head_{\phi}, pos_{\phi}, neg_{\phi}, ineq_{\phi})$  where  $head_{\phi}$  is an atom;  $pos_{\phi}$  and  $neg_{\phi}$  are sets of atoms;  $ineq_{\phi}$  is a set of inequalities (u=v) with  $u,v \in var$  and the variables of  $\phi$  all occur in  $pos_{\phi}$ . The components  $head_{\phi}$ ,  $pos_{\phi}$  and  $neg_{\phi}$  are called respectively the head, the positive body atoms and the negative body atoms. We refer to  $pos_{\phi} \cup neg_{\phi}$  as the body atoms. Note,  $neg_{\phi}$  contains just atoms, not negative literals. Every Datalog<sup>¬</sup> rule  $\phi$  must have a head,  $pos_{\phi}$  must be non-empty and  $neg_{\phi}$  may be empty. If  $neg_{\phi} = \emptyset$  then  $\phi$  is called positive.

Of course, a rule  $\phi$  may be written in the conventional syntax. For instance, if  $head_{\phi} = T(u, v)$ ,  $pos_{\phi} = \{R(u, v)\}$ ,  $neg_{\phi} = \{S(v)\}$ , and  $ineq_{\phi} = u \neq v$ , with  $u, v \in var$ , then we can write  $\phi$  as:

$$T(u,v): -R(u,v), \neg S(v), u \neq v$$

The set of variables of  $\phi$  is denoted  $vars(\phi)$ . A rule  $\phi$  is said to be over schema  $\sigma$  if for each atom  $R(u_1,\ldots,u_k)\in head\phi\cup pos\phi\cup neg\phi$ , the arity of R in  $\phi$  is k. A Datalog program P over  $\sigma$  is a set of Datalog eg rules over  $\sigma$ . We write sch(P) to denote the (minimal) database schema that P is over. We define  $idb(P)\subset sch(P)$  to be the database schema consisting of all relations in rule-heads of P. We abbreviate edb(P)=sch(P) idb(P). As usual, the abbreviation "idb" stands for "intensional database schema" and "edb" stands for "extensional database schema". A valuation for a rule  $\phi$  in P w.r.t. an instance I over edb(P), is a total function  $V: vars(\phi) \to dom$ . The application of V to an atom  $R(u_1,\ldots,u_k)$  of  $\phi$ , denoted  $V(R(u_1,\ldots,u_k))$ , results in the fact  $R(a_1,\ldots,a_k)$  where  $a_i=V(u_i)$  for each  $i\in 1,\ldots,k$ . This is naturally extended to a set of atoms, which results in a set of facts. The valuation V is said to be satisfying for  $\phi$  on I if  $V(pos_{\phi}) \subset I, V(neg\phi) \cap I = \emptyset$ , and  $V(u) \neq V(v)$  for each  $(u \neq v) \in ineq_{\phi}$ . If so,  $\phi$  is said to derive the fact  $V(head_{\phi})$ .

#### 1.1.3. Positive and semi-positive Datalog

A Datalog  $\neg$  program P is positive if all rules of P are positive. We say that P is semi-positive if for each rule  $\phi \in P$ , the atoms of  $neg_{\phi}$  are over edb(P). We now give the semantics of a semi-positive Datalog  $\neg$  program P. First, let  $T_P$  be the immediate consequence operator that maps each instance J over sch(P) to the instance  $J = J \cup A$  where A is the set of facts derived by all possible satisfying valuations for the rules of P on J. Let I be an instance over edb(P). Consider the infinite sequence  $I_0, I_1, I_2$ , etc, inductively defined as follows:  $I_0 = I$  and  $I_i = T_P(I_{i-1})$  for each  $i \ge 1$ . The output of P on input I, denoted P(I), is defined as  $\bigcup_i I_j$ ; this is the  $minimal\ fixpoint$  of the  $T_P$  operator.

#### 1.1.4. Stratified semantics

We say that P is syntactically stratifiable if there is a function  $\rho: sch(P) \to 1, ..., |idb(P)|$  such that for each rule  $\phi \in P$ , having some head predicate T, the following conditions are satisfied:

- 1.  $\rho(R) \leqslant \rho(T)$  for each  $R(\overline{u}) \in pos_{\phi} \cap idb(P)$
- 2.  $\rho(R) < \rho(T)$  for each  $R(\overline{u}) \in neg_{\phi} \cap idb(P)$ .

For  $R \in idb(P)$ , we call  $\rho(R)$  the stratum number of R. Intuitively,  $\rho$  partitions P into a sequence of semi-positive Datalog  $\neg$  programs  $P_1, ..., P_k$  with  $k \leq |idb(P)|$  such that for each  $i=1,\ldots,k$ , the program  $P_i$  contains the rules of P whose head predicate has stratum number i. This sequence is called a syntactic stratification of P. We can now apply the stratified semantics to P: for an input I over sch(P), we first compute the fixpoint  $P_1(I)$ , then the fixpoint  $P_2(P_1(I))$ , etc. The output of P on input I, denoted P(I), is defined as  $P_k(P_{k-1}(\ldots P_1(I)\ldots))$ . It is well known that the output of P does not depend on the chosen syntactic stratification (if more than one exists). Not all Datalog  $\neg$  programs are syntactically stratifiable. By stratifiedDatalog we refer to all Datalog  $\neg$  programs which are syntactically stratifiable.

## 1.2. Evaluation strategies for Datalog

bottom-up, top-down

# Pregel/Giraph

@TODO: Description of Pregel model and info about the Giraph project, also mention other extensions like Giraph++.

## SociaLite

SociaLite ([1], [2]) is a graph query language based on Datalog. While Datalog allows to express some of graph algorithms in an elegant and succint way, many practical problems cannot be efficiently solved with Datalog programs.

SociaLite allows a programmer to write intuitive queries using declarative semantics, which can often be executed as efficiently as highly optimized dedicated programs. The queries can be executed in a distributed environment with no need for the programmer to worry about distributing the computation or managing the communication protocols.

The language consists of a few extensions to Datalog, most significant of which is the ability to combine recursive rules with aggregation. Under some conditions, such rules can be evaluated incrementally and thus as efficiently as regular recursion in Datalog.

[1] introduces Sequential SociaLite, intended to be executed on one machine, consisting of two main extensions: recursive aggregate functions and tail-nested tables. Recursive aggregate functions are the most important new feature in Socialite – in 3.1 we present a complete definition and proofs of correctness of that extension, which are missing in [1]. Tail-nested tables are a much more straightforward extension – an optimization of data layout in memory. They are described in 3.2

[2] extends Sequential Socialite to *Distributed SociaLite*, executable on a distributed architecture. It introduces a *location operator*, shows how the data and computations can be distributed. The programmer does not have to think about how to distribute the data between machines or manage the communication between them. He only specifies an abstract *location* for each row in the data, and the data and computations are automatically sharded. Distributed SociaLite is covered in section 3.3

Additionally, thanks to the declarative sematics of Datalog and SociaLite, it is possible to provide an important optimization: the *delta stepping* technique, which is an effective way of parallelizing the Dijkstra algorithm [8]. In SociaLite, this technique can be applied automatically to a certain class of recursive aggregate programs.

In distributed computations on large graphs, an approximate result is often enough. Usually we can observe the *long tail* phenomenon in the computation, where a good approximate solution is achieved quickly, but it takes a long time to get to the optimal one. In SociaLite, by simply stopping the computation, we can obtain an approximate solution found so far. [2] also shows a method which can significantly reduce memory requirements by storing the intermediate results in an approximate way using Bloom filters. Those topics are covered in section 3.5

## 3.1. Recursive aggregate functions

#### 3.1.1. Motivation

Most graph algorithms are essentially some kind of iteration or recursive computation. Simple recursion can be expressed easily in Datalog. However, in many problems the computation results are gradually refined in each iteration, until the final result is reached. Examples of such algorithms are the Dijkstra algorithm for single source shortest paths or Page Rank. Usually, it is difficult or impossible to express such algorithms in Datalog efficiently, as it would require computing much more intermediate results than it is actually needed to obtain the solution. We will explain that on an example: a simple program that computes shortest paths from a source node.

A straightforward Datalog program for computing single source shortests paths (starting from node 1) is presented below. Due to limitations of Datalog, this program computes all possible path lengths from node 1 (the source) to other nodes in the first place, and after that for each node the minimal distance is chosen. Not only this approach results in bad performance, but causes the program to execute infinitely if a loop in the graph is reachable from the source node.

```
\begin{aligned} & \text{Path}(t,d) & : - & \text{Edge}(1,t,d). \\ & \text{Path}(t,d) & : - & \text{Path}(s,d_1), \text{Edge}(s,t,d_2), d = d_1 + d_2. \\ & \text{MinPath}(t,\text{Min}(d)) & : - & \text{Path}(t,d). \end{aligned}
```

Rysunek 3.1: Datalog query for computing shortest paths from node 1 to other nodes

SociaLite allows aggregation to be combined with recursion under some conditions. This allows us to write staightforward programs for such problems, which finish execution in finite time and often are much more efficient than Datalog programs. An example SociaLite program that computes single source shortest paths is presented below.

```
PATH(t, Min(d)) : - EDGE(1, t, d).: - PATH(s, d_1), EDGE(s, t, d_2), d = d_1 + d_2.
```

Rysunek 3.2: SociaLite query for computing shortest paths from node 1 to other nodes

While being very useful, recursive aggregation rules not always have an unambiguous solution. This is the case only under some conditions on the rules and the aggregation function itself.

Typically, Datalog programs semantics is defined using the least fixed point of instance inclusion. This requires that the subsequent computation interations only add tuples to the database instance, but never remove tuples from the instance. This is the reason for which program?? is inefficient. When recursive aggregate functions are allowed, this is not the case: a tuple in the instance can be replaced with a different one because a new aggregated value appeared. Consequently, in order to define SociaLite programs sematics in terms of least fixed point, we need to use a different order on database instances.

First, we will define a meet operation and show the order that it induces. Then we will show

that if the aggregation function is a meet operation and corresponding rules are monotone with respect to this induced order, then the result of the program is unambiguously defined. We will also show how it can be computed efficiently.

#### 3.1.2. Meet operation and induced ordering

**Definition 3.1.1.** A binary operation is a *meet* operation if it is idempotent, commutative and associative.

@TODO: Maybe remind definitions of semi-lattice and partial order?

#### Order induced by a meet operation

A meet operation  $\sqcap$  defines a semi-lattice: it induces a partial order  $\preceq$  over its domain, such that the result of the operation for any two elements is the least upper bound of those elements with respect to  $\preceq$ 

**Example 3.1.1.**  $\max(a,b)$  for  $a,b \in \mathbb{N}$  is a meet operation; it is:

- idempotent  $-\max(a, a) = a$
- commutative  $-\max(a, b) = \max(b, a)$
- associative  $-\max(a, \max(b, c)) = \max(\max(a, b), c)$

It induces the partial order  $\leq$ : for any two  $a, b \in \mathbb{N}$ ,  $\max(a, b)$  is their least upper bound with respect to  $\leq$ .

On the contrary, + is not a meet operation, since it is not idempotent:  $1+1 \neq 1$ .

#### Order on relation instances

We can extend the partial order  $\leq$  to an order  $\sqsubseteq$  on relation instances as follows:

$$R_1 \sqsubseteq R_2 \iff \forall_{(q_1,\dots,q_{n-1},v)\in g(R_1)} \exists_{(q_1,\dots,q_{n-1},v')\in g(R_2)} v \preceq v'$$

@TODO: Show that this is actually a partial order.

@TODO: komentarz, przykład

We can easily see that an empty relation instance  $\emptyset$  is smaller under  $\sqsubseteq$ 

#### 3.1.3. SociaLite program

A SociaLite program is a Datalog program, with additional aggregation function defined for some of the relations: For each relation R, there can be one column  $agg_R$  chosen for which an aggregation function  $\sqcap_R$  is provided. The rest of the columns are called the *qualifying* columns. Intuitively, after each step of computation, we group the tuples in the relation by the qualifying columns and aggregate the column  $agg_R$  using  $\sqcap_R$ .

For simplicity, we assume that the aggregated column is always the last one.

Syntactically, we require that all rules with a given relation R in the head are defined together, with the head defined once and possibly many rule bodies on the right hand side:

```
P(x_1, \dots, x_k, Agg(y)) : - Q_1(x_1, \dots, x_k, y)
\vdots - Q_m(x_1, \dots, x_k, y).
```

Rysunek 3.3: Definition of multiple rules for one relation.

#### 3.1.4. Semantics and evaluation - one relation case

In this section we will show that the semantics of a SociaLite program can be unambiguously defined using least fixed point, as long as it satisfies some conditions. To simplify the reasoning, we will restrict our attention to programs with only one *idb* relation. In the next section we extend the definitions and theorems from this section to the general case of many *idb* relations.

Let P be a SociaLite program, with only one idb relation R of arity k in the form of:

$$R(x_1, ..., x_{k-1}, F(y))$$
 : -  $Q_1(x_1, ..., x_{k-1}, y)$  ... : -  $Q_m(x_1, ..., x_{k-1}, y)$ .

Let us define:

- $f: \mathbf{dom^k} \to \mathbf{dom^k}$  the function that evaluates the rules  $Q_1, ... Q_m$  based on the contents of the relation given in the argument and returns its input extended with the set of generated tuples
- $f: \mathbf{dom^k} \to \mathbf{dom^k}$  the function that performs grouping and aggregation:  $q(R) = \{(x_1, \dots, x_{k-1}, F(\{y: (x_1, \dots, x_{k-1}, y) \in R\}) : (x_1, \dots, x_{k-1}, \overline{y}) \in R\}$
- $h = g \circ f$

**Theorem 3.1.1.** If f is monotone with respect to  $\sqsubseteq$ , i.e.  $R_1 \subseteq R_2 \to f(R_1) \subseteq f(R_2)$ , and there exists  $n \ge 0$ , such that  $h^n(\emptyset) = h^{n+1}(\emptyset)$ , then  $R^* = h^n(\emptyset)$  is the least fixed point of h, that is:

- 1.  $R^* = h(R^*)$
- 2.  $R^* \sqsubseteq R$  for all R such that R = h(R)

Proof:

g is monotone with respect to  $\sqsubseteq$ . Since we assumed that f is monotone with respect to  $\sqsubseteq$ ,  $h=g\circ f$  is also monotone with respect to  $\sqsubseteq$ .

Let us notice that  $\emptyset$  is greater under  $\sqsubseteq$  than any other element. TODO: show that (very easy)

Let us suppose that R' is any fixpoint of h. We know that  $\emptyset \subseteq R'$ . Applying h n times to both sides of the inequality, we have that  $R^* = h^n(\emptyset) \subseteq h^n(R') = R'$ , thanks to the monotonicity of h with respect to  $\subseteq$ . Therefore, the inductive fixed point  $R^*$  is the least fixed point of h.

@TODO: open questions:

- How we compute recursive functions with non-meet aggregation operators?
- 3.2. Tail-nested tables
- 3.3. Distributed SociaLite
- 3.4. Delta stepping in Distributed SociaLite
- 3.5. Approximate evaluation in Distributed SociaLite

# Translating SociaLite programs to Giraph programs

# Implementation

# Summary

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