Faculty of Mathematics, Computer Science and Natural Sciences Research Group Computer Science 2

RWTH Aachen University, Germany

Bachelor Thesis

Analyzing runtime complexity of non-ndg Term Rewrite Systems

submitted by

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Coming soon

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Statutory Declaration in Lieu of an Oath

Abstract

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Contents

Chapter 1

Introduction

Many researchers would disagree that this was the main focus.

Term rewrite systems have been extensively studied, with the main problem concerning their complexity analysis. In recent decades many automated tools have been developed to infer upper bounds on runtime complexity (rc), which considers the longest rewrite sequence starting with a basic term disregarding any evaluation strategy. In contrast to rc, innermost runtime complexity (irc) makes use of an innermost evaluation strategy that can be much easier to analyse. While the analysis of irc has some uses, it is rc that is of greater interest.

Actually, irc is more interesting and thus further developed than rc.

Recently a technique has been developed to overapproximate some TRSs as non-dup generalized (ndg), with a corresponding proof of the equality of rc and irc for such ndg TRSs. The mentioned criteria combined with the much more powerful analysis of irc allowed inferring upper/lower bounds on rc from upper/lower bounds on irc. This method however makes no claims regarding all the TRSs that are non-ndg. It is the goal of this thesis to extend the technique to non-ndg TRSs.

The goal is to transfer the existing results from irc to rc.

Presented here is a transformation of the original non-ndg TRS into an encoded version, that shares the same rc. The goal of the encoding is to create a TRS, for which also holds that irc = rc. While the encoded version may not ndg, what is important is that we can apply the techniques used in the analysis of irc to infer the rc of the original non-ndg TRS.

What does that mean?

?

As it will be shown, not all encodings are equal and some yield substantially better results. Two algorithms have been developed to improve the results of the automated analysis. The first - overapproximating the TRS positions that require encoding and a second - detecting when to alter the encoding to not add any non-terminating rules.

This sounds like there are two approaches to encode the TRS, but there is only one, right?

The structure of the thesis is as follows. Chapter 2 introduces some of the preliminary knowledge regarding TRSs. Chapter 3 contains a description of the first mentioned algorithm, which is used as the foundation, on which the encoding, described in Chapter 4, is built. Chapter 5 is dedicated to the second algorithm and the changes that come up from its result. In Chapter 6 are presented the results from the implementation of the technique in the tool AProVE and it also serves as a summary.

The introduction should be easily understandable for any computer science student that did not yet get to know TRSs.
It should include the following parts, where the first part is mostly missing.
* Why is the topic important?
* What has already been done?
* What is the gap?/ What is missing?
* What does this thesis do to fill the gap?

Chapter 2

Preliminaries

In this chapter is introduced some of the necessary groundwork on term rewrite systems with accompanying examples.

A term rewrite system (TRS) $\mathcal{R}$ is a finite set of rewrite rules $\ell \to r$, where ℓ and r are terms in the set $\mathcal{T}(\Sigma, \mathcal{V})$ over the signature Σ and the set of variables $\mathcal{V}$. In the context of a rule $\ell \to r$, ℓ refers to the left-hand side and r to the right-hand side. The rules describe how terms are rewritten, i.e. the left-hand side of a rule can be matched against some part of a term s, rewrite it to the matching right-hand side and give output t.

Introduce the term before using it.

I like that you start with an example to illustrate very early how a TRS works. You could even pull this into the introduction. However, you should avoid getting into the technical details so early, if you don't want to actually define TRSs at this point. Move these details to def. 5.

Example 1 (Addition)**.**

$$\begin{aligned} \alpha_1 &: \mathsf{add}(\mathsf{x}, \mathsf{0}) &&\to \quad \mathsf{x} \\ \alpha_2 &: \mathsf{add}(\mathsf{x}, \mathsf{s}(\mathsf{y})) &&\to \quad \mathsf{add}(\mathsf{s}(\mathsf{x}), \mathsf{y}) \end{aligned}$$

Example 1. shows a simple TRS for the calculation of addition on natural numbers. Numbers here are represented by the so-called successor function and the base symbol 0, i.e. 1 would be $\mathsf{s}(\mathsf{0})$, 2 - $\mathsf{s}(\mathsf{s}(\mathsf{0}))$ and so on. In this way our base case would be represented in α_1, where $x + 0 = x$. In α_2 the TRS recursively subtracts 1 from the second argument and adds 1 to the first argument. This repeats until the second argument reaches 0 and thus the first argument is output. One could intuitively imagine this as follows:

Avoid unclear notation. Here you could just replace - by "be".

$$3 + 2 \quad = \quad 4 + 1 \quad = \quad 5 + 0 \quad = \quad 5$$

Also denote the rewrite sequence using terms?

Definition 1 (Signature [1])**.**

A **signature** Σ is a set of function symbols, where each $f \in \Sigma$ is associated with an arity $n \geq 0$. For some valid n, we denote the set of all n-ary elements of Σ as $\Sigma^{(n)}$. The elements of $\Sigma^{(0)}$ are also called constant symbols.

The signature of example 1. is therefore $\{\mathsf{add}, \mathsf{s}, \mathsf{0}\}$ with the accompanying $\Sigma^{(2)} = \{\mathsf{add}\}$, $\Sigma^{(1)} = \{\mathsf{s}\}$ and $\Sigma^{(0)} = \{\mathsf{0}\}$, whereas x and y are variables.

Definition 2 (Terms [1])**.**

Let Σ be a signature and $\mathcal{V}$ a set of variables such that $\Sigma \cap \mathcal{V} = \emptyset$. The set $\mathcal{T}(\Sigma, \mathcal{V})$ of all **terms** over Σ and $\mathcal{V}$ is inductively defined as:

- $\mathcal{V} \subseteq \mathcal{T}(\Sigma, \mathcal{V})$
- for all $n \geq 0$, all $f \in \Sigma^{(n)}$ and all $t_1, ..., t_n \in \mathcal{T}(\Sigma, \mathcal{V})$, we have $f(t_1, ..., t_n) \in \mathcal{T}(\Sigma, \mathcal{V})$

The first item states that each variable on its own is a term and the second item specifies that for arbitrary n many terms $t_1, ..., t_n$ and a function symbol f with arity n, its application $f(t_1, ..., t_n)$ is also a term.

If Σ and $\mathcal{V}$ are irrelevant or clear from the context, only $\mathcal{T}$ is written instead of $\mathcal{T}(\Sigma, \mathcal{V})$ [2] .

Given the signature $\Sigma = \{\mathsf{add}, \mathsf{s}, \mathsf{0}\}$ of example 1, we can construct terms like $\mathsf{s(0)}$ or $\mathsf{add(x, y)}$ or more complex ones like $\mathsf{add(add(x, s(s(0))), s(0))}$.

Definition 3 (Term positions [1, 2])**.**

1. Let Σ be a signature, $\mathcal{V}$ be a set of variables with $\Sigma \cap \mathcal{V} = \emptyset$ and $s \in \mathcal{T}(\Sigma, \mathcal{V})$. The set of **positions** of the term s is a set $pos(s)$ of strings, inductively defined as follows:
   - If $s = x \in \mathcal{V}$, then $pos(s) := \{\epsilon\}$, where ϵ denotes the empty string
   - If $s = f(s_1, ..., s_n)$, then
   $$pos(s) := \{\epsilon\} \cup \bigcup_{i=1}^{n} \{i.\pi \mid \pi \in pos(s_i)\}$$

   The position ϵ of a term s refers to the **root position** of s and the symbol at that position is called the **root symbol** denoted by $root(s)$. The **prefix order** on term positions is defined as
   $$\pi \leq \tau \text{ iff there exists } \pi' \text{ such that } \pi.\pi' = \tau$$
   It is a partial order. Positions for which neither $\pi \leq \tau$, nor $\pi \geq \tau$ hold, are called **parallel positions** denoted by $\pi \parallel \tau$.
2. The **size** of a term s is defined as $|s| := |pos(s)|$.
3. For $\pi \in pos(s)$, the **subterm of s at position π** denoted by $s|_\pi$ is defined inductively as
   $$s|_\epsilon := s$$
   $$f(s_1, ..., s_n)|_{i.\pi} := s_i|_\pi$$
   If $\pi \neq \epsilon$, then $s|_\pi$ is a **proper subterm** of s.
4. For $\pi \in pos(s)$, the **replacement in s at position π with term t** denoted by $s[t]_\pi$ is defined as
   $$s[t]_\epsilon := t$$
   $$f(s_1, ..., s_n)[t]_{i.\pi} := f(s_1, ..., s_i[t]_\pi, ..., s_n)$$
5. By $\mathcal{V}ar(s)$ we denote the set of **variables occurring in** s, i.e.
   $$\mathcal{V}ar(s) := \{x \in \mathcal{V} \mid \exists \pi \in pos(s) \text{ such that } s|_\pi = x\}$$

Consider the term $t = \mathsf{add(x, s(y))}$ from example 1. The set of position is $pos(t) = \{\epsilon, 1, 2, 2.1\}$. It holds that $\epsilon \leq 2 \leq 2.1$, $\epsilon \leq 1$ and also $1 \parallel 2$. Some expressions to illustrate are $t|_2 = \mathsf{s(y)}$ and $t[s(0)]_1 = \mathsf{add(s(0), s(y))}$.

Definition 4 (Substitution [1])**.**

Let Σ be a signature and $\mathcal{V}$ a finite set of variables. A **substitution** is a function $\sigma : \mathcal{V} \to \mathcal{T}(\Sigma, \mathcal{V})$ such that $\sigma(x) \neq x$ for only finitely many xs. The **domain** $Dom(\sigma) := \{x \mid \sigma(x) \neq x\}$ is the set of all variables that σ does not map to themselves. The **range** $Ran(\sigma) := \{\sigma(x) \mid x \in Dom(\sigma)\}$ is the set of all terms that σ maps to. If $Dom(\sigma) = \{x_1, ..., x_n\}$, then we may write σ as

Only introduce notation that you actually use in your thesis.

$$\sigma = \{x_1 \mapsto \sigma(x_1) \ , \ \cdots \ , \ x_n \mapsto \sigma(x_n)\} \ .$$

A term t is called an **instance** of a term s, iff there exists a substitution σ such that $\sigma(s) = t$.

Ensure that you explain everything in you own words. There does not have to be a lot of explanation in the preliminaries, but some, which must not be copied word-for-word from another source.

Definition 5 (Term rewrite systems [2])**.**

We don't use TRS rules as identities (i.e., we don't apply them in both directions), so I suggest to omit this wording an notation.

A **rewrite rule** is an identity $\ell \approx r$ such that ℓ is not a variable and $\mathcal{V}ar(r) \subseteq \mathcal{V}ar(\ell)$. In this case it is denoted as $\ell \to r$ instead of $\ell \approx r$. A **term rewrite system** (TRS) is a finite set of such rewrite rules.

1. Given the signature Σ of a TRS $\mathcal{R}$ we have
   - the **set of defined symbols** $\Sigma_d := \{root(\ell) \mid \ell \to r \in \mathcal{R}\}$
   - the **set of constructor symbols** $\Sigma_c := \Sigma \setminus \Sigma_d$
2. A term $f(s_1, ..., s_n)$ is **basic**, if $f \in \Sigma_d$ and $s_1, ..., s_n \in \mathcal{T}(\Sigma_c, \mathcal{V})$. The **set of all basic terms** over Σ and $\mathcal{V}$ is denoted by $\mathcal{T}_B(\Sigma, \mathcal{V})$.
3. Given term t and $x \in \mathcal{V}$ the **number of occurrences of** x **in** t is denoted by $\#_x(t)$.
4. A **redex** (**red**ucible **ex**pression) is an instance of ℓ of some rule $\ell \to r$.
   - A term is an **innermost redex**, if none of its proper subterms is a redex.
   - A term is in **normal form**, if none of its subterms is a redex.

In example 1. we have $\Sigma_d = \{\mathsf{add}\}$ and $\Sigma_c = \{\mathsf{0}, \mathsf{s}\}$.

s is not basic.

Consider the basic term $s = \mathsf{add}(\mathsf{add}(\mathsf{0}, \mathsf{0}), \mathsf{0})$. With a substitution $\sigma = \{x \mapsto \mathsf{add}(\mathsf{0}, \mathsf{0})\}$ it holds that s is an instance of the left-hand side of rule α_1 and therefore s is a redex. But it is not an innermost one as the subterm $s|_1 = \mathsf{add}(\mathsf{0}, \mathsf{0})$ can be reduced to $\mathsf{0}$. Since none of the subterms of $s|_1$ can be reduced, it is an innermost redex.

Definition 6 (Rewrite step [2])**.**

A **rewrite step** is a reduction(or evaluation) of a term s to t by applying a rule $\ell \to r$ at position π, denoted by $s \to_{\ell \to r, \pi} t$ and means that for some substitution σ, $s|_\pi = \sigma(\ell)$ and $t = s[\sigma(r)]_\pi$ holds. A reduction can be expressed as $s \to_{\mathcal{R}, \pi} t$, if it holds for some $\ell \to r \in \mathcal{R}$. Subscripts such as the rule or position of reduction can be omitted, if they are irrelevant.

This should be a = I think.

1. A sequence of rewrite steps(or rewrite sequence) $s \to t_1 \to t_2 \to \cdots \to t_m = t$ is denoted by $s \to^m t$.
2. A rewrite step is called **innermost**, if $s|_\pi$ is an innermost redex, and is denoted by $s \xrightarrow{i}_\pi r$

The basic term $s = \mathsf{add}(\mathsf{add}(\mathsf{0}, \mathsf{0}), \mathsf{0})$ from example 1. can be used to create the following sequence:

$$\mathsf{add}(\mathsf{add}(\mathsf{0}, \mathsf{0}), \mathsf{0}) \to \mathsf{add}(\mathsf{0}, \mathsf{0}) \to \mathsf{0}$$

Again, this is not a basic term. Also, it does not occur in example 1.

Definition 7 (Derivation height [2])**.**

The **derivation height** $dh : \mathcal{T} \times 2^{\mathcal{T}\times\mathcal{T}} \to \mathbb{N}\cup\{\omega\}$ is a function with two inputs:a term t and a binary relation on terms, in this case $\to$. It defines the longest sequence of rewrite steps starting with term t, i.e.

the length of

$$dh(t, \to) = \sup\{m \mid \exists t' \in \mathcal{T}\ ,\ t \to t'\}$$

In the case of an infinitely long rewrite sequence starting with term s, it is denoted as $dh(s, \to) = \omega$.

Definition 8 ((Innermost) Runtime complexity [2])**.**

The **runtime complexity**(rc) of a TRS $\mathcal{R}$ maps any $n \in \mathbb{N}$ to the length of the longest $\to$-sequence (rewrite sequence) starting with a basic term t with $|t| \leq n$. The innermost runtime complexity(irc) is defined analogously, but it only considers innermost rewrite steps. More precisely $\mathrm{rc}_\mathcal{R} : \mathbb{N} \to \mathbb{N}\cup\{\omega\}$ and $\mathrm{irc}_\mathcal{R} : \mathbb{N} \to \mathbb{N}\cup\{\omega\}$, defined as

$$\mathrm{rc}_\mathcal{R}(n) = \sup\{dh(t, \to_\mathcal{R}) \mid t \in \mathcal{T}_B, |t| \leq n\}$$
$$\mathrm{irc}_\mathcal{R}(n) = \sup\{dh(t, \xrightarrow{i}_\mathcal{R}) \mid t \in \mathcal{T}_B, |t| \leq n\}$$

Since every $\xrightarrow{i}$-sequence can be viewed as a $\to$-sequence, it is clear that $\mathrm{irc}_\mathcal{R}(n) \leq \mathrm{rc}_\mathcal{R}(n)$. Therefore, an upper bound for $\mathrm{rc}_\mathcal{R}$ infers an upper bound for $\mathrm{irc}_\mathcal{R}$ and a lower bound for $\mathrm{irc}_\mathcal{R}$ infers a lower bound for $\mathrm{rc}_\mathcal{R}$. A specific class of TRSs has been previously studied in [2], for which $\mathrm{rc}_\mathcal{R} = \mathrm{irc}_\mathcal{R}$ holds. The results from the analysis on runtime complexity of such TRSs can now be used in the analysis on innermost runtime complexity and vice versa, as mentioned above.

Better formulate the "vice versa" case explicitly, since that is the practically relevant case.

Example 1. is sufficiently simple to see that the analysis on rc coincides with the one on irc, because every rewrite sequence starting with a basic term is also an innermost one. This TRS is a part of a class of systems, for which $\mathrm{rc}_\mathcal{R} = \mathrm{irc}_\mathcal{R}$. For now we can say that the length of rewrite sequences in example 1. depends solely on the second argument of the starting basic term. The size of this subterm is decremented by one after each rewrite until it reaches 0, terminating after one more rewrite. Thus it holds that $\mathrm{rc}_\mathcal{R} = \mathrm{irc}_\mathcal{R} \in \mathcal{O}(n)$.

Definition 9 (Non-dup-generalized innermost(ndg) rewrite step [2, 3])**.**

The rewrite step $s \to_{\ell\to r,\pi} t$ with the matching substitution σ is called **non-dup-generalized** innermost(ndg), if

- For all variables x with $\#_x(r) > 1$, $\sigma(x)$ is in normal form, and
- For all $\tau \in pos(\ell)\backslash\{\epsilon\}$ with $root(\ell|_\tau) \in \Sigma_d$, it holds that $\sigma(\ell|_\tau)$ is in normal form.

A TRS $\mathcal{R}$ is called ndg, if every rewrite sequence starting with a basic term consists of only ndg rewrite steps. It is proven in [2] that $\mathrm{rc}_\mathcal{R} = \mathrm{irc}_\mathcal{R}$ for all ndg TRSs $\mathcal{R}$. One can now easily observe that example 1. does fit the criteria, i.e. it has no duplicated variables on the right-hand side of rules and no nested defined symbols on left-hand sides.

Example 2.

This and the following example will present some TRSs, that do not fit one of the criteria stated above, and show why $\mathrm{rc}_\mathcal{R} \neq \mathrm{irc}_\mathcal{R}$ for them.

$$\begin{array}{lccc} \alpha_1: & \mathsf{g} & \to & \mathsf{f(a)} \\ \alpha_2: & \mathsf{a} & \to & \mathsf{b} \\ \alpha_3: & \mathsf{f(a)} & \to & \mathsf{f(a)} \end{array}$$

Always write the TRS in the subscript of the arrow,
to distinguish between the arrow for rules (without the subscript)
and the arrow for rewrite steps (with the subscript).

The $\mathrm{rc}_{\mathcal{R}}$ of this TRS is reached via the infinite sequence $\mathsf{g} \to \mathsf{f}(\mathsf{a}) \to \mathsf{f}(\mathsf{a}) \to \cdots$, while the $\mathrm{irc}_{\mathcal{R}}$ corresponds the two step long sequence $\mathsf{g} \to \mathsf{f}(\mathsf{a}) \to \mathsf{f}(\mathsf{b})$. It is the nested a in rule α_3 that makes this a non-ndg TRS with as shown non-equal rc and irc.

Example 3.

$$\begin{array}{lccc} \alpha_1: & \mathsf{f}(\mathsf{x}) & \to & \mathsf{dbl}(\mathsf{g}(\mathsf{x})) \\ \alpha_2: & \mathsf{dbl}(\mathsf{x}) & \to & \mathsf{d}(\mathsf{x},\mathsf{x}) \\ \alpha_3: & \mathsf{g}(\mathsf{s}(\mathsf{x})) & \to & \mathsf{g}(\mathsf{x}) \end{array}$$

Here rule α_2 is duplicating. Using the basic term $\mathsf{f}(\mathsf{s}(0))$ we get the longest sequence $\mathsf{f}(\mathsf{s}(0)) \to \mathsf{dbl}(\mathsf{g}(\mathsf{s}(0))) \to \mathsf{d}(\mathsf{g}(\mathsf{s}(0)),\mathsf{g}(\mathsf{s}(0))) \to^2 \mathsf{d}(\mathsf{g}(0),\mathsf{g}(0))$, which is just one step longer than the innermost rewrite sequence $\mathsf{f}(\mathsf{s}(0)) \to \mathsf{dbl}(\mathsf{g}(\mathsf{s}(0))) \to \mathsf{dbl}(\mathsf{g}(0)) \to \mathsf{d}(\mathsf{g}(0),\mathsf{g}(0))$.

While the technique used in [2] is helpful in inferring upper and lower bounds for rc and irc on ndg systems, it makes no statements on non-ndg systems. The examples above show that the criteria for ndg systems clearly exclude a lot of TRSs. It is the goal of this thesis to expand the mentioned technique to non-ndg systems by encoding certain positions in the TRS. For the new encoded TRS it will hold that $\mathrm{rc}_{\mathcal{R}} = \mathrm{irc}_{\mathcal{R}}$.

This R is the encoded R.

This R is the original R.

Make sure to clarify this difference.

Chapter 3

Algorithm for finding non-ndg locations

Chapter 4

Encoding

Chapter 5

Encoding with terminating rules

Chapter 6

Results

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