

FACULTY OF MATHEMATICS, COMPUTER SCIENCE AND
NATURAL SCIENCES RESEARCH GROUP COMPUTER SCIENCE 2

RWTH AACHEN UNIVERSITY, GERMANY

Bachelor Thesis

A Term Encoding to Analyze Runtime Complexity via Innermost Runtime Complexity

submitted by

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Coming soon

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Abstract

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Chapter 1

Introduction

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.

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.

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 $\text{irc} = \text{rc}$. 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.

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.

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.

Chapter 2

Preliminaries

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

For now a term can be viewed as some object. Rewriting is then simply a transformation of one term into another. This rewriting is guided by rules, which describe what the input and output of the transformation is. Rules are also associated with some cost, which is considered when analyzing runtime complexity. Unless stated otherwise all rules have a cost of *one*. Rules can also have conditions besides what input they take, however such conditional rules are not in the focus of this thesis. Combining rules together creates a term rewrite system.

Example 1 (Addition).

$$\begin{aligned}\alpha_1 &: \text{add}(x, 0) \rightarrow x \\ \alpha_2 &: \text{add}(x, s(y)) \rightarrow \text{add}(s(x), y)\end{aligned}$$

Example 1. shows a simple term rewrite system (TRS) for the calculation of addition on natural numbers. Numbers here are represented by the so-called successor function and the symbol 0, i.e. 1 would be $s(0)$, 2 would be $s(s(0))$ and so on. For the sake of readability, it is denoted $f^n(x)$, when some function is applied n -many times to an input x .

The base case would be represented in α_1 , where $x + 0 = x$. In α_2 the TRS recursively subtracts 1 from the second position and adds 1 to the first position. This repeats until the second position reaches 0 and thus the first position is output. One could intuitively imagine this as follows:

$$\begin{aligned}3 + 2 &= 4 + 1 = 5 + 0 = 5 \\ \text{add}(s^3(0), s^2(0)) &\rightarrow_{\mathcal{R}} \text{add}(s^4(0), s(0)) \rightarrow_{\mathcal{R}} \text{add}(s^5(0), 0) \rightarrow_{\mathcal{R}} s^5(0)\end{aligned}$$

While not explicitly stated in the example, x and y are variables which can be instantiated with other terms. Without a proper definition, one could assume they are some constants like 0. Therefore each TRS is associated with a set, which holds all function symbols.

Definition 1 (Signature [1]).

A **signature** Σ is a set of function symbols. The number of inputs n of each function symbol is referred to as its arity and it cannot be negative. The subset $\Sigma^{(n)} \subseteq \Sigma$ holds all symbols with arity n . Symbols with arity 0 are called constant.

The signature of example 1. is therefore $\{\text{add}, s, 0\}$ with the accompanying $\Sigma^{(2)} = \{\text{add}\}$, $\Sigma^{(1)} = \{s\}$ and $\Sigma^{(0)} = \{0\}$, 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 $t_1, \dots, t_n \in \mathcal{T}(\Sigma, \mathcal{V})$, we have $f(t_1, \dots, t_n) \in \mathcal{T}(\Sigma, \mathcal{V})$

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, \dots, t_n and a function symbol f with arity n , $f(t_1, \dots, t_n)$ is also a term. $\mathcal{T}(\Sigma, \mathcal{V})$ is denoted as \mathcal{T} , if Σ and \mathcal{V} are irrelevant or clear from the context [2].

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

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 some term in $\mathcal{T}(\Sigma, \mathcal{V})$. The set of **positions** of s contains sequences of natural numbers, denoted $\text{pos}(s)$ and inductively defined as follows:

- If $s = x \in \mathcal{V}$, then $\text{pos}(s) := \{\varepsilon\}$
- If $s = f(s_1, \dots, s_n)$, then

$$\text{pos}(s) := \{\varepsilon\} \cup \bigcup_{i=1}^n \{i.\pi \mid \pi \in \text{pos}(s_i)\}$$

The symbol ε is used to point to the **root position** of a term. The function symbol at that position is called the **root symbol** denoted by $\text{root}(s)$. Positions can be compared to each other

$$\pi \leq \tau, \text{ iff there exists } \pi' \text{ such that } \pi.\pi' = \tau$$

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 $|s| := |\text{pos}(s)|$.
3. For $\pi \in \text{pos}(s)$, the **subterm of s at position π** denoted by $s|_\pi$ is defined as

$$s|_\varepsilon := s$$

$$f(s_1, \dots, s_n)|_{i.\pi} := s_i|_\pi$$

If $\pi \neq \varepsilon$, then $s|_\pi$ is a **proper subterm** of s .

4. For $\pi \in \text{pos}(s)$, the **replacement in s at position π with term t** denoted by $s[t]_\pi$ is defined as

$$s[t]_\varepsilon := t$$

$$f(s_1, \dots, s_n)[t]_{i.\pi} := f(s_1, \dots, s_i[t]_\pi, \dots, s_n)$$

5. The set of **variables occurring in s** , denoted $\text{Var}(s)$ is defined as

$$\text{Var}(s) := \{x \in \mathcal{V} \mid \exists \pi \in \text{pos}(s) \text{ such that } s|_\pi = x\}$$

Consider the term $t = \text{add}(\mathbf{x}, \mathbf{s}(\mathbf{y}))$. The set of position is $\text{pos}(t) = \{\varepsilon, 1, 2, 2.1\}$. Some statements that hold for example are $\varepsilon \leq 2 \leq 2.1$ and also $1 \parallel 2$. Some expressions to illustrate are $t|_2 = \mathbf{s}(\mathbf{y})$ and $t[\mathbf{s}(0)]_1 = \text{add}(\mathbf{s}(0), \mathbf{s}(\mathbf{y}))$.

Definition 4 (Substitution [1]).

Let Σ be a signature and \mathcal{V} a finite set of variables. A **substitution** is a function $\sigma : \mathcal{V} \rightarrow \mathcal{T}(\Sigma, \mathcal{V})$ such that $\sigma(x) \neq x$ holds for only finitely many xs . It can be written as

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

The term resulting from $\sigma(s)$ is called an **instance** of s .

Definition 5 (Term rewrite systems [2]).

A **rewrite rule** is a pair of terms $\ell \rightarrow r$ such that ℓ is not a variable and $\mathcal{V}ar(r) \subseteq \mathcal{V}ar(\ell)$. It is referred to ℓ as the left-hand side(lhs) and r as the right-hand side(rhs). A **term rewrite system** is a finite set of such rewrite rules combined with a signature Σ .

1. Given the signature Σ of a TRS \mathcal{R} we have

- the **set of defined symbols** $\Sigma_d := \{root(\ell) \mid \ell \rightarrow r \in \mathcal{R}\}$
- the **set of constructor symbols** $\Sigma_c := \Sigma \setminus \Sigma_d$

2. A term $f(s_1, \dots, s_n)$ is **basic**, if $f \in \Sigma_d$ and $s_1, \dots, s_n \in \mathcal{T}(\Sigma_c, \mathcal{V})$. The **set of all basic terms** over Σ and \mathcal{V} is denoted by $\mathcal{T}_B(\Sigma, \mathcal{V})$.

3. The **number of occurrences of x in term t** is denoted by $\#_x(t)$.

4. A **redex** (**reducible expression**) is an instance of ℓ of some rule $\ell \rightarrow r$.

- 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.

5. Rules are also associated with a **cost** and a **condition**. For the purposes of this thesis all rules have no conditions and a cost of either 0, denoted $\ell \xrightarrow{0} r$, or 1, denoted $\ell \rightarrow r$.

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

Consider the term $s = \text{add}(\text{add}(0, 0), 0)$. With a substitution $\sigma = \{x \mapsto \text{add}(0, 0)\}$ 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 $s|_1 = \text{add}(0, 0)$ can be reduced to 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 $\ell \rightarrow r$ at position π , denoted by $s \rightarrow_{\ell \rightarrow r, \pi} t$ and means that for some substitution σ , $s|_\pi = \sigma(\ell)$ and $t = s[\sigma(r)]_\pi$ holds.

1. Rule and position in the subscript can be omitted, if they are irrelevant. We denote $s \rightarrow_{\mathcal{R}} t$, if it holds for some rule in \mathcal{R} , and in order to distinguish between rules and rewrite steps.

2. A sequence of rewrite steps(or rewrite sequence) $s = t_0 \rightarrow_{\mathcal{R}} t_1 \rightarrow_{\mathcal{R}} \dots \rightarrow_{\mathcal{R}} t_m = t$ is denoted by $s \rightarrow_{\mathcal{R}}^m t$.

3. Only rules with cost 1 contribute to the length of a rewrite sequence, i.e. for $s \rightarrow_{\mathcal{R}}^m t$ it holds that $m = 0$, if it is a single rewrite step that uses a rule with 0 cost.

4. A rewrite step is called **innermost**, if $s|_\pi$ is an innermost redex, and is denoted by $s \xrightarrow{i}_\pi r$

The term $s = \text{add}(\text{add}(0, 0), 0)$ mentioned above can be used to create the following sequence with the TRS from example 1.:

$$\text{add}(\text{add}(0, 0), 0) \rightarrow_{\mathcal{R}} \text{add}(0, 0) \rightarrow_{\mathcal{R}} 0$$

Definition 7 (Derivation height [2]).

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

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

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

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 \rightarrow -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 $rc_{\mathcal{R}} : \mathbb{N} \rightarrow \mathbb{N} \cup \{\omega\}$ and $irc_{\mathcal{R}} : \mathbb{N} \rightarrow \mathbb{N} \cup \{\omega\}$, defined as

$$\begin{aligned} rc_{\mathcal{R}}(n) &= \sup\{dh(t, \rightarrow_{\mathcal{R}}) \mid t \in \mathcal{T}_B, |t| \leq n\} \\ irc_{\mathcal{R}}(n) &= \sup\{dh(t, \overset{i}{\rightarrow}_{\mathcal{R}}) \mid t \in \mathcal{T}_B, |t| \leq n\} \end{aligned}$$

Since every $\overset{i}{\rightarrow}$ -sequence can be viewed as a \rightarrow -sequence, it is clear that $irc_{\mathcal{R}}(n) \leq rc_{\mathcal{R}}(n)$. Therefore, an upper bound for $rc_{\mathcal{R}}$ infers an upper bound for $irc_{\mathcal{R}}$ and a lower bound for $irc_{\mathcal{R}}$ infers a lower bound for $rc_{\mathcal{R}}$. A specific class of TRSs has been previously studied in [2], for which $rc_{\mathcal{R}} = irc_{\mathcal{R}}$ holds. It follows then that the upper bound on irc infers an upper bound on rc and that the lower bound on rc infers a lower bound on irc .

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 $rc_{\mathcal{R}} = irc_{\mathcal{R}}$. 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 $rc_{\mathcal{R}} = irc_{\mathcal{R}} = \mathcal{O}(n)$.

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

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

- For all variables x with $\#_x(r) > 1$, $\sigma(x)$ is in normal form, and
- For all $\tau \in pos(\ell) \setminus \{\varepsilon\}$ 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. TRSs and rewrite steps, which do not fit the criteria are therefore called **non-ndg**.

One can now easily observe that the TRS from example 1. is indeed ndg, since none of its rules have duplicated variables on the rhs of rules or nested defined symbols on the lhs of rules. [2] proves that $rc_{\mathcal{R}} = irc_{\mathcal{R}}$ for all ndg TRSs \mathcal{R} .

TO DO: ELABORATE

Example 2.

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

$$\begin{aligned} \alpha_1 : \quad & \mathbf{g} \quad \rightarrow \quad \mathbf{f(a)} \\ \alpha_2 : \quad & \mathbf{a} \quad \rightarrow \quad \mathbf{b} \\ \alpha_3 : \quad & \mathbf{f(a)} \quad \rightarrow \quad \mathbf{f(a)} \end{aligned}$$

The $rc_{\mathcal{R}}$ of this TRS is reached via the infinite sequence $g \rightarrow_{\mathcal{R}} f(a) \rightarrow_{\mathcal{R}} f(a) \rightarrow_{\mathcal{R}} \dots$, while the $irc_{\mathcal{R}}$ corresponds the two step long sequence $g \rightarrow_{\mathcal{R}} f(a) \rightarrow_{\mathcal{R}} f(b)$. 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.

$$\begin{array}{lll} \alpha_1 : & f(x) & \rightarrow \text{dbl}(g(x)) \\ \alpha_2 : & \text{dbl}(x) & \rightarrow d(x, x) \\ \alpha_3 : & g(s(x)) & \rightarrow g(x) \end{array}$$

Here rule α_2 is duplicating. Using the basic term $f(s(0))$ we get the longest sequence

$$f(s(0)) \rightarrow_{\mathcal{R}} \text{dbl}(g(s(0))) \rightarrow_{\mathcal{R}} d(g(s(0)), g(s(0))) \rightarrow_{\mathcal{R}}^2 d(g(0), g(0)),$$

which is just one step longer than the innermost rewrite sequence

$$f(s(0)) \rightarrow_{\mathcal{R}} \text{dbl}(g(s(0))) \rightarrow_{\mathcal{R}} \text{dbl}(g(0)) \rightarrow_{\mathcal{R}} d(g(0), g(0)).$$

While this example shows a difference when exact rc and irc are considered, it is important to note that both are linear in terms of asymptotic notation

This thesis is based on the work in [2] and aims to expand its uses beyond ndg systems. While a non-ndg TRS \mathcal{R} fails to satisfy $rc_{\mathcal{R}} = irc_{\mathcal{R}}$, it is possible that a TRS \mathcal{R}' exists such that $rc_{\mathcal{R}} = irc_{\mathcal{R}'}$. Ideally it would also hold that every full rewrite sequence in \mathcal{R} has an equally long innermost rewrite sequence in \mathcal{R}' and vice versa. The approach taken in this thesis is to transform the original TRS \mathcal{R} by changing some rules and adding new ones, turning it into an encoded version of itself.

Chapter 3

Algorithm for finding non-ndg locations

By definition every TRS \mathcal{R} has finitely many rules and therefore finitely many positions, which can be changed, i.e. encoded. However, for practical purposes encodings that encode every possible position are not useful. Generally speaking, encodings like this introduce new rules, which can produce loops and create infinitely long rewrite sequences. The focus of this chapter lies in the optimization of which positions to encoded(Section 1.) and the limitation of non-terminating rules(Section 2.). To this effort the goal of the encoding is to ensure each innermost rewrite sequence in the encoded TRS \mathcal{R}' has an equally long full rewrite sequence in the TRS \mathcal{R} and vice versa, which is later relevant for the proof of $\text{rc}\mathcal{R} = \text{irc}\mathcal{R}'$.

Before we continue, it is important to clear one misconception. The encoding does not transform a non-ndg TRS \mathcal{R} into an ndg \mathcal{R}' , which we can analyze like in [2]. We can show this best via an example.

Example 4.

TRS \mathcal{Q}			encoded TRS \mathcal{Q}'		
$\alpha_1 :$	h	$\rightarrow g(a, a)$	$\alpha'_1 :$	h	$\rightarrow g(i(l_a), i(l_a))$
$\alpha_2 :$	a	$\rightarrow b$	$\alpha'_2 :$	a	$\rightarrow b$
$\alpha_3 :$	$g(x, a)$	$\rightarrow f(x, x)$	$\alpha'_3 :$	$g(x, l_a)$	$\rightarrow f(i(x), i(x))$
			$\alpha'_4 :$	$i(l_a)$	$\xrightarrow{0} i(a)$
			$\alpha'_5 :$	$i(x)$	$\xrightarrow{0} x$

Colored in red are the places where the encoded TRS \mathcal{Q}' differs from \mathcal{Q} . We can observe the introduction of two new symbols, i.e. the defined symbol i and the constructor l_a . Two new rules are also added and their significance will be discussed later. Currently we want to show that the encoded TRS \mathcal{Q}' is non-ndg.

The duplication of variables can not be avoided in \mathcal{Q}' , leaving the possibility of a non-ndg rewrite step. Consider the following sequence

$$h \rightarrow_{\mathcal{R}} g(i(l_a), i(l_a)) \xrightarrow{0}_{\mathcal{R}} g(i(a), i(l_a)) \xrightarrow{0}_{\mathcal{R}} g(i(a), l_a) \xrightarrow{0}_{\mathcal{R}} g(a, l_a) \rightarrow_{\mathcal{R}} f(i(a), i(a)).$$

The last rewrite step duplicates the term a , which is not in normal form. Therefore neither \mathcal{Q} , nor \mathcal{Q}' are ndg.

The new defined symbol i can be more intuitively explained as a helper function, which can choose what part of the term is to be rewritten next. In order to do this, certain positions in the TRS are to be encoded, besides the addition of the new rules. The encoding of these positions usually consists of replacing a defined symbol with its new constructor counterpart, e.g. replacing a with l_a , and/or enclosing the position in i .

The sequence above can be continued to eventually result in the longest sequence in \mathcal{Q}' and it would also be as long as the longest in \mathcal{Q} . It is however not an innermost one. With a slight change of evaluation, the following innermost rewrite sequence is produced:

$$h \rightarrow_{\mathcal{R}} g(i(l_a), i(l_a)) \xrightarrow{0}_{\mathcal{R}} g(l_a, l_a) \rightarrow_{\mathcal{R}} f(i(l_a), i(l_a)) \xrightarrow{0}_{\mathcal{R}} f(i(a), i(a)) \xrightarrow{2}_{\mathcal{R}} f(b, b).$$

This is one of the sequence that the automatic complexity analysis will discover and use for its result, as the rest will either be shorter or equal in length. For example the innermost rewrite sequence

$$h \rightarrow_{\mathcal{R}} g(i(l_a), i(l_a)) \xrightarrow{0}_{\mathcal{R}} g(l_a, i(a)) \rightarrow_{\mathcal{R}} g(l_a, i(b)) \xrightarrow{0}_{\mathcal{R}} g(l_a, b)$$

has only a length of 2, but it can be matched to

$$h \rightarrow_{\mathcal{R}} g(a, a) \rightarrow_{\mathcal{R}} g(a, b),$$

which is exactly one of the properties we desire in our encoding.

3.1 Non-ndg locations

Because we consider rewrite sequences and not just rules, it is important to also ask if certain TRS-positions are reachable in a rewrite sequence and what 'flows' into them. In example 4, it is clear that the rhs of α_1 can be matched against the lhs of α_3 creating a so called **data flow**, i.e. the **a**-symbol at position 1 of the rhs of α_1 flows into the variable **x** at position 1 of the lhs of α_3 . Since it is very confusing to refer to positions in a TRS like that, we can combine this information in a *triple*.

Definition 10 (Location).

Let \mathcal{R} be a TRS. The set of all **locations**(TRS-positions) of \mathcal{R} is defined as

$$\mathcal{L}_{\mathcal{R}} := \{(\alpha, x, \pi) \mid \alpha = \ell \rightarrow r \in \mathcal{R}, x \in \{\mathbf{L}, \mathbf{R}\}, \pi \in \text{pos}(\ell) \text{ if } x = \mathbf{L}, \text{ else } \pi \in \text{pos}(r)\}.$$

- For all $\ell \rightarrow r \in \mathcal{R}$ we write $\mathcal{R}|_{(\ell \rightarrow r, \mathbf{L}, \pi)} = \ell|_{\pi}$ and $\mathcal{R}|_{(\ell \rightarrow r, \mathbf{R}, \pi)} = r|_{\pi}$.
- A location λ is called a **variable location**, if $\mathcal{R}|_{\lambda} \in \mathcal{V}$.
- The function $\text{SUB}_{\mathcal{R}} : \mathcal{L}_{\mathcal{R}} \rightarrow \mathcal{L}_{\mathcal{R}}$ returns the set of all **sub-locations** of a location $\lambda = (\alpha, \mathbf{X}, \tau)$ with $\mathbf{X} \in \{\mathbf{L}, \mathbf{R}\}$ and is defined as

$$\text{SUB}_{\mathcal{R}}(\lambda) = \text{SUB}_{\mathcal{R}}((\alpha, \mathbf{X}, \tau)) := \{(\alpha, \mathbf{X}, \pi) \mid \tau \leq \pi\}.$$

While some work on data flow analysis exists [4], it was deemed not precise enough to use for our encoding. Instead it was preferred to create a separate algorithm, that would return all to-be-encoded positions in a given TRS. After introducing the algorithm, we are going to apply it to the TRS \mathcal{R} from example 4 in order to show how the encoded positions are chosen.

The CAP function will be used in the algorithm to ease the matching of terms and the detection of data flows. It will be a cause of less precision in our algorithm as will be shown later.

Definition 11 ($\text{CAP}_{\mathcal{R}}$).

For a given TRS \mathcal{R} the function $\text{CAP}_{\mathcal{R}} : \mathcal{T} \rightarrow \mathcal{T}$ replaces all proper subterms with a defined root symbol of a term q with different fresh variables. Multiple occurrences of the same subterm are replaced by pairwise different variables.

$$\text{CAP}_{\mathcal{R}}(q) = \text{CAP}_{\mathcal{R}}(f(t_1, \dots, t_n)) = f(t'_1, \dots, t'_n) \text{ with}$$

- $t'_k = t_k$, if $t_k \in \mathcal{V}$
- $t'_k = \text{CAP}_{\mathcal{R}}(t_k)$, if $\text{root}(t_k) \in \Sigma_c$

- $t'_k = x \in \mathcal{V}$ and $x \notin \text{Var}(t'_j)$ for $1 \leq j < k$, if $\text{root}(t_k) \in \Sigma_d$

Consider the TRS \mathcal{R} from example 1. and the term $t = \text{add}(\text{s}(\text{add}(x, y)), \text{add}(x, y))$. It holds that

$$\text{CAP}_{\mathcal{R}}(t) = \text{add}(\text{s}(x_1), x_2).$$

Algorithm 1.

Let \mathcal{R} be a TRS with signature $\Sigma = \Sigma_d \cup \Sigma_c$ and a set of variables \mathcal{V} .

1. Let $\mathcal{X}_{\mathcal{R}}$ be the smallest set such that

$$\bullet \{(\alpha, L, \tau) \mid \alpha = \ell \rightarrow r \in \mathcal{R}, \tau \in \text{pos}(\ell) \setminus \{\varepsilon\}, \text{root}(\ell|_{\tau}) \in \Sigma_d\} \subseteq \mathcal{X}_{\mathcal{R}} \quad (S1)$$

This set includes all locations of nested defined symbols on the lhs of rules.

$$\bullet \{(\alpha, R, \tau) \mid \alpha = \ell \rightarrow r \in \mathcal{R}, \tau, \pi \in \text{pos}(r), \tau \neq \pi, r|_{\tau} = r|_{\pi} \in \mathcal{V}\} \subseteq \mathcal{X}_{\mathcal{R}} \quad (S2)$$

This set includes all locations of duplicated variables on the rhs of rules.

$$\bullet \{(\alpha, L, \tau) \mid \alpha = \ell \rightarrow r \in \mathcal{R}, \tau \in \text{pos}(\ell), \pi \in \text{pos}(r), \ell|_{\tau} = r|_{\pi} \in \mathcal{V}\} \subseteq \mathcal{X}_{\mathcal{R}}, \text{ if } (\alpha, R, \pi) \in \mathcal{X}_{\mathcal{R}} \quad (S3)$$

This set includes all locations of variables on the lhs of rules, that flow into locations on the rhs of the same rule, which are also included in $\mathcal{X}_{\mathcal{R}}$.

$$\bullet \{(\alpha, R, \tau) \mid \alpha = \ell \rightarrow r \in \mathcal{R}, \tau \in \text{pos}(r), \tau = \pi.v, v \neq \varepsilon, \\ \beta = s \rightarrow t \in \mathcal{R}, \omega \in \text{pos}(s), v \nparallel \omega, \\ \exists \text{ MGU for } \text{CAP}_{\mathcal{R}}(r|_{\pi}) \text{ and } s\} \subseteq \mathcal{X}_{\mathcal{R}}, \text{ if } (\beta, L, \omega) \in \mathcal{X}_{\mathcal{R}} \quad (S4)$$

This set includes all locations on rhs of rules, that flow into locations on lhs of rules, which are also included in $\mathcal{X}_{\mathcal{R}}$.

$$\bullet \{(\alpha, R, \tau) \mid \alpha = \ell \rightarrow r \in \mathcal{R}, \tau \in \text{pos}(r), \\ \beta = s \rightarrow t \in \mathcal{R}, \pi \in \text{pos}(t), \text{root}(t|_{\pi}) \in \Sigma_d, \\ \exists \text{ MGU for } \text{CAP}(t|_{\pi}) \text{ and } \ell\} \subseteq \mathcal{X}_{\mathcal{R}}, \text{ if } (\beta, R, \pi) \in \mathcal{X}_{\mathcal{R}} \quad (S5)$$

This set includes all locations on rhs of rules, whose lhs root symbol is also at a rhs location included in $\mathcal{X}_{\mathcal{R}}$.

2. Let $\mathcal{Y}_{\mathcal{R}}^{\Delta}$ with $\Delta = \{(\alpha, R, \tau) \mid \alpha = \ell \rightarrow r \in \mathcal{R}, \text{root}(r|_{\tau}) \in \Sigma_d\}$ be the smallest set such that

$$\bullet \Delta \subseteq \mathcal{Y}_{\mathcal{R}}^{\Delta} \quad (S6)$$

This set includes all locations on the rhs of rules with a defined symbol.

$$\bullet \{(\beta, L, \omega) \mid \alpha = \ell \rightarrow r \in \mathcal{R}, \tau \in \text{pos}(r), \tau = \pi.v, v \neq \varepsilon, \\ \beta = s \rightarrow t \in \mathcal{R}, \omega \in \text{pos}(s), v \nparallel \omega, \\ \exists \text{ MGU for } \text{CAP}_{\mathcal{R}}(r|_{\pi}) \text{ and } s\} \subseteq \mathcal{Y}_{\mathcal{R}}^{\Delta}, \text{ if } (\alpha, R, \tau) \in \mathcal{Y}_{\mathcal{R}}^{\Delta} \quad (S7)$$

The definition of this subset is similar to (S4). Changed is which location from a data flow is included.

$$\bullet \{(\alpha, R, \tau) \mid \alpha = \ell \rightarrow r \in \mathcal{R}, \pi \in \text{pos}(\ell), \tau \in \text{pos}(r), \ell|_{\pi} = r|_{\tau} \in \mathcal{V}\} \subseteq \mathcal{Y}_{\mathcal{R}}^{\Delta}, \text{ if } (\alpha, L, \pi) \in \mathcal{Y}_{\mathcal{R}}^{\Delta} \quad (S8)$$

The definition of this subset is similar to (S3). Changed is which location from a data flow is included.

3. Let $\mathcal{Z}_{\mathcal{R}} := \{(\alpha, L, \pi) \mid \mathcal{R}|_{(\alpha, L, \pi)} \in \mathcal{V}\}$ be the set of all variable locations on the lhs of rules in \mathcal{R} .

4. The set of all locations to be encoded $\text{ENC}_{\mathcal{R}}$ is defined as

$$\text{ENC}_{\mathcal{R}} := (\mathcal{X}_{\mathcal{R}} \cap \mathcal{Y}_{\mathcal{R}}^{\Delta}) \setminus \mathcal{Z}_{\mathcal{R}}.$$

In order to better illustrate why the subsets ($S1 - S8$) are defined this way, we are going to apply the algorithm to the TRS \mathcal{Q} from example 4. and go into further detail. We call ($S1 - S2$) and ($S6$) *base subsets* as they do not depend on other locations being included in their respective larger set. The rest we call *conditional subsets*. Since the locations we want to encode potentially break ndg criteria, we will call them non-ndg from now on. More specifically, a **non-ndg location** λ is one, for which $\lambda \in \mathcal{X}_{\mathcal{R}}$.

$$\bullet \{(\alpha_3, L, 1)\} \subseteq \mathcal{X}_{\mathcal{Q}} \quad (S1)$$

The first of the two base subsets in $\mathcal{X}_{\mathcal{Q}}$ contains locations of nested defined symbols on lhs of rules. Such locations, if reachable in a sequence, must be replaced by constructors, otherwise the encoded TRS \mathcal{Q}' can not discover an equivalent innermost rewrite sequence, as it will be forced to evaluate the nested defined symbol.

$$\bullet \{(\alpha_3, L, 2), (\alpha_3, R, 1), (\alpha_3, R, 2)\} \subseteq \mathcal{X}_{\mathcal{Q}} \quad (S2)$$

The second of the two base subsets in $\mathcal{X}_{\mathcal{Q}}$ contains locations of duplicated variables on rhs of rules. It is most commonly the duplication of defined symbols that creates a discrepancy between innermost and full runtime complexity. However, it is not always the case that a redex is duplicated and this is what the second part of the algorithm in $\mathcal{V}_{\mathcal{Q}}^{\Delta}$ deals with. For now they are included for having the potential to duplicate redexes.

$$\bullet \{(\alpha_3, L, 2), (\alpha_3, R, 1), (\alpha_3, R, 2), (\alpha_3, L, 1)\} \subseteq \mathcal{X}_{\mathcal{Q}} \quad (S3)$$

The first conditional subset deals with data flows from left- to right-hand side location. We want to include variable locations on the lhs of rules that flow into non-ndg locations. Later these locations included via ($S3$) will be removed, as the encoding does not change them, but for the purposes of building $\mathcal{X}_{\mathcal{Q}}$ now, they must be considered.

$$\bullet \{(\alpha_3, L, 2), (\alpha_3, R, 1), (\alpha_3, R, 2), (\alpha_3, L, 1), (\alpha_1, R, 1), (\alpha_1, R, 2)\} \subseteq \mathcal{X}_{\mathcal{Q}} \quad (S4)$$

The second conditional subset deals with data flows from right- to left-hand side locations

We first split τ in two parts. The symbol at π is the root symbol of the term we try to match against some lhs of any rule. While $\pi = \varepsilon$ is allowed $v = \varepsilon$ is not, because this would include the ε positions of rhs of rules, which can not flow between locations.

TO DO: CAP-FUNCTION ELABORATION

TO DO: σ INSTEAD OF MGU?

The MGU is the matcher between the two terms and it means that a term reached via a rewrite step using rule α , can then potentially be used as input for rule β .

Lastly, we require that $v \not\parallel \omega$. This part ensures that the newly included location in α actually flows into a non-ndg location. In our example $\lambda_1 := (\alpha_1, R, 1)$ flows only into $\lambda_2 := (\alpha_3, L, 1)$ and we would want to include λ_1 , only if λ_2 is non-ndg.

$$\bullet \{(\alpha_3, L, 2), (\alpha_3, R, 1), (\alpha_3, R, 2), (\alpha_3, L, 1), (\alpha_1, R, 1), (\alpha_1, R, 2), (\alpha_2, R, \varepsilon)\} \subseteq \mathcal{X}_{\mathcal{Q}} \quad (S5)$$

The third conditional subset includes the so-called *return positions* of non-ndg locations with defined symbols on the rhs of rules. Because they are non-ndg, it means that they flow into a non-ndg location. But in a rewrite sequence that defined symbol may be rewritten before flowing into said non-ndg location. Therefore we must also include these locations.

While this example was simple enough as to allow for the computation of each of the subsets in a single step, most TRS would require that the subsets ($S3 - S5$) be continuously reevaluated, until no further additions are made.

We can now move onto the second part of the algorithm. It may be the case that only constructor terms flow into some variables, which have no affect on runtime complexity, if duplicated. Similarly, instances of lhs of some rule with a nested defined symbol may never be reachable, when starting with a basic term. In both cases encoding such locations does not in any way better the analysis on the encoded version. Therefore, excluding all locations from $\mathcal{X}_{\mathcal{R}}$ that have no location of a defined symbol flowing into them will create a more precise encoding. The set of all locations that either contain a defined symbol or a location, that contains a defined symbol, flows into them $\mathcal{V}_{\mathcal{Q}}^{\Delta}$ will later be intersected with $\mathcal{X}_{\mathcal{Q}}$ to give us all non-ndg locations, that contain or potentially receive a defined symbol.

$$\bullet \{(\alpha_1 R, \varepsilon), (\alpha_1, R, 1), (\alpha_1, R, 2)\} \subseteq \mathcal{Y}_Q^\Delta \quad (S6)$$

The only base subset of \mathcal{Y}_Q^Δ includes any location of a defined symbol on the rhs of rules. In this example they all occur in rule α_1 .

$$\bullet \{(\alpha_1 R, \varepsilon), (\alpha_1, R, 1), (\alpha_1, R, 2), (\alpha_3, L, 1), (\alpha_3, L, 2)\} \subseteq \mathcal{Y}_Q^\Delta \quad (S7)$$

This and the following conditional subsets are similarly defined to others in \mathcal{X}_Q , but here included are the receivers of the data flow, rather than the source.

As stated previously the CAP_Q has its problems and does not entirely remove overapproximation of non-ndg locations. In order to show this, assume $\mathcal{Q}|_{(\alpha_1, R, 2)} = h$. Then obviously the lhs of α_3 is not reachable, when starting with a basic term and should not be encoded. But since $CAP_Q(g(a, h)) = g(x_1, x_2)$ can be matched with the lhs of α_3 , we include $(\alpha_3, L, 2)$ in \mathcal{Y}_Q^Δ as well. For now this is an acceptable overapproximation.

$$\bullet \{(\alpha_1 R, \varepsilon), (\alpha_1, R, 1), (\alpha_1, R, 2), (\alpha_3, L, 1), (\alpha_3, L, 2), (\alpha_3, R, 1), (\alpha_3, R, 2)\} \subseteq \mathcal{Y}_Q^\Delta \quad (S8)$$

As mentioned above this conditional subset is defined similarly to one in \mathcal{X}_Q . The included location and the location in the condition are swapped just like in (S7). In Q this only happens in α_3 , which adds the last element to \mathcal{Y}_Q^Δ . Analogously to \mathcal{X}_Q , this set is also continuously reevaluated until no further additions are made.

We can now exclude the variable locations on the lhs of rules, since they do not need to be encoded. It is also acceptable to leave them in the set, but to define the encoding in such a way as to ignore them. Here the former approach is taken.

$$\begin{aligned} \bullet \{(\alpha_3 L, 1)\} &= \mathcal{Z}_Q \\ \bullet ENC_Q &= (\mathcal{X}_Q \cap \mathcal{Y}_Q^\Delta) \setminus \mathcal{Z}_Q = \{(\alpha_3, R, 1), (\alpha_3, R, 2), (\alpha_3, L, 2)(\alpha_1, R, 1), (\alpha_1, R, 2)\} \end{aligned}$$

As we can see the non-ndg locations included in ENC_Q correspond exactly to the encoded locations in the TRS Q' . However, the encoding also adds new rules in order to define how the *i*-symbol evaluates. While we have answered the question of what positions we need to encode, we still need to investigate what rules to add to the encoded version of a TRS. This will finally allow us to define the encoding.

3.2 Rules

There are three types of these *i*-rules and they can be summarized as follows:

$$\bullet i(l_a) \rightarrow i(a)$$

The *executing rules* usually take the constructor version of a defined symbol as input and evaluate to it. Since this is happening in the context of an innermost evaluation strategy, this means that the next rewrite step will evaluate the *a*-symbol.

An example of this is the sequence using TRS Q' from example 4:

$$h \rightarrow_{\mathcal{R}} g(i(l_a), i(l_a)) \xrightarrow{0}_{\mathcal{R}} g(i(a), i(l_a)) \rightarrow_{\mathcal{R}} g(i(b), i(l_a)).$$

$$\bullet i(x) \rightarrow x$$

The *omission rule* is mandatory in all encodings and it effectively means that the next symbol to be evaluated is at a more outer position in the term.

Continuing the sequence from above we get an example of just that:

$$\dots \rightarrow_{\mathcal{R}} g(i(b), i(l_a)) \xrightarrow{0}_{\mathcal{R}} g(b, l_a) \rightarrow_{\mathcal{R}} f(i(b), i(b)).$$

- $i(l_{\text{dbl}}(x)) \rightarrow i(l_{\text{dbl}}(i(x)))$

The *propagation rules* push the i -symbol further inside the term. While this type of rule is missing in example 4, for now we can assume that the dbl -symbol was defined. The usage of these rules means that the next symbol to be evaluated is at a more inner position in the term. These rules have two variations: the first is as shown above (*non-terminating*) and the second omits the i -symbol at position ε on the rhs (*terminating*). We make this distinction due to the necessary omission rule.

A non-terminating rule in combination with $i(x) \rightarrow x$ creates non-terminating rewrite sequences like the following:

$$\cdots \rightarrow_{\mathcal{R}} i(l_{\text{dbl}}(x)) \xrightarrow{0}_{\mathcal{R}} i(l_{\text{dbl}}(i(x))) \xrightarrow{0}_{\mathcal{R}} i(l_{\text{dbl}}(x)) \rightarrow_{\mathcal{R}} \cdots$$

Such sequences could throttle the automatic complexity analysis and therefore it is best to avoid adding non-terminating rules whenever possible. We are now going to focus on why propagating rules are a part of our encoding and how we can detect, which version of them the encoding should add.

Example 5.

TRS \mathcal{P}			encoded		TRS \mathcal{P}'			
$h(x)$	\rightarrow	$f(x, x)$	$h(x)$	\rightarrow	$f(x, a(x))$	$i(s(x))$	\rightarrow	$s(i(x))$
$f(x, s(y))$	\rightarrow	$f(\text{dbl}(g(y, x), y))$	$f(x, s(y))$	\rightarrow	$f(i(l_{\text{dbl}}(l_{\text{lin}}(x))), y)$	$i(d(x, x))$	\rightarrow	$d(i(x), i(x))$
$\text{dbl}(x)$	\rightarrow	$d(x, x)$	$\text{dbl}(x)$	\rightarrow	$d(i(x), i(x))$	$i(l_{\text{dbl}}(x))$	\rightarrow	$i(l_{\text{dbl}}(i(x)))$
$\text{lin}(s(x))$	\rightarrow	$\text{lin}(x)$	$\text{lin}(s(x))$	\rightarrow	$i(l_{\text{lin}}(x))$	$i(l_{\text{lin}}(x))$	\rightarrow	$i(l_{\text{lin}}(i(x)))$
$a(x)$	\rightarrow	$s(x)$	$a(x)$	\rightarrow	$s(x)$	$i(l_{\text{dbl}}(x))$	\rightarrow	$i(\text{dbl}(x))$
						$i(l_{\text{lin}}(x))$	\rightarrow	$i(\text{lin}(x))$
						$i(x)$	\rightarrow	x

There are full rewrite sequences in \mathcal{P} , which result in terms of the shape

$$f(\text{dbl}(\text{lin}(\text{dbl}(\text{lin}(\cdots \text{lin}(x)) \cdots), y),$$

which in order to have an equivalently long innermost rewrite sequence in \mathcal{P}' , the encoding must include some rule that allows for the i -symbol to be moved inside the term. For now this can only be done by the inclusion of the non-terminating propagation rules, colored here in red.

However, this does not necessarily hold for other TRSs, especially for ones that do not use either an outer- or innermost evaluation strategy to reach their worst case time complexity.

The root of this problem is the repeated nesting of defined symbols, which in \mathcal{P} occurs rather obviously in the second rule. Therefore, an algorithm that checks, if a defined symbol f at some location λ eventually flows into a sub-location of λ , then we can conclude that a repeated nesting of this terms occurs and that the encoding must include the non-terminating version of the propagation rule of l_f .

Algorithm 2.

Let \mathcal{R} be a TRS with signature $\Sigma = \Sigma_d \cup \Sigma_c$ and a set of variables \mathcal{V} .

1. Let $\text{INF}_{\mathcal{R}}$ be defined as

$$\text{INF}_{\mathcal{R}} := \{t \in \mathcal{L}_{\mathcal{R}} \mid \text{root}(\mathcal{R}|_t) \in \Sigma_d, \Delta = \{t\}, |\mathcal{Y}_{\mathcal{R}}^{\Delta} \cap \text{SUB}_{\mathcal{R}}(t)| \geq 2\}$$

Chapter 4

Encoding

Chapter 5

Encoding with terminating rules

Chapter 6

Results

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