A Recursive Inclusion Checker for Recursively Defined Subtypes

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We present a tableaux procedure that checks logical relations between recursively defined subtypes of recursively defined types and apply this procedure to the problem of resolving ambiguous names in a programming language. This work is part of a project to design a new programming language suitable for efficient implementation of logic. Logical formulas are tree-like structures with many constructors having different arities and argument types. Algorithms that use these structures must perform case analysis on the constructors, and access subtrees whose type and existence depend on the constructor used. In many programming languages, case analysis is handled by matching, but we want to take a different approach, based on recursively defined subtypes. Instead of matching a tree against different constructors, we will classify it by using a set of disjoint subtypes. Subtypes are more general than structural forms based on constructors, we expect that they can be implemented more efficiently, and in addition can be used in static type checking. This makes it possible to use recursively defined subtypes as preconditions or postconditions of functions. We define the types and the subtypes (which we will call adjectives), define their semantics, and give a tableaux-based inclusion checker for adjectives. We show how to use this inclusion checker for resolving ambiguous field references in declarations of adjectives. The same procedure can be used for resolving ambiguous function calls.

Keywords: programming language design; type systems; theorem proving; compiler construction

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Рекурсивная проверка включения для рекурсивно определенных подтипов
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Мы представляем табличную процедуру, которая проверяет логические отношения между рекурсивно определенными подтипами рекурсивно определенных типов, и применяем эту процедуру к проблеме разрешения неоднозначных имен в языке программирования. Эта работа является частью проекта по разработке нового языка программирования, подходящего для эффективной реализации логики. Логические формулы представляют собой древовидные структуры со множеством конструкторов, имеющих различные свойства и типы аргументов. Алгоритмы, использующие эти структуры, должны выполнять анализ вариантов для конструкторов и получать доступ к поддеревьям, тип и существование которых зависят от используемого конструктора. Во многих языках программирования анализ прецедентов обрабатывается путем сопоставления, но мы хотим использовать другой подход, основанный на рекурсивно определенных подтипах. Вместо сопоставления дерева с различными конструкторами мы будем классифицировать его с помощью набора непересекающихся подтипов. Подтипы являются более общими, чем структурные формы, основанные на конструкторах, мы ожидаем, что они могут быть реализованы более эффективно и, кроме того, могут использоваться при статической проверке типов. Это позволяет использовать рекурсивно определенные подтипы в качестве предварительных условий или постусловий функций. Мы определяем типы и подтипы (которые мы будем называть прилагательными), определяем их семантику и даем проверку включения прилагательных на основе таблиц. Мы показываем, как использовать эту проверку включения для разрешения неоднозначных ссылок на поля в объявлениях прилагательных. Та же процедура может быть использована для разрешения неоднозначных вызовов функций.

Ключевые слова: проектирование языков программирования; системы типов; доказательство теорем; построение компилятора

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1. Introduction

Our goal is to develop a programming language for implementation of logic that is convenient to use on one hand, and efficient on the other hand. Traditionally, functional languages like OCaml ([1]), Haskell ([2]) or Scala ([3]) are considered the most suitable for implementation of logic.

Functional languages have inductive types, and use matching for accessing subtrees. Matching can be viewed as simultaneously inspecting a tree and extracting subtrees into local variables. In our proposed language we want to keep inductive types, but add a mechanism for definition of subtypes, which we will use both as a replacement for matching, and as a basis for static overload resolution. The subtypes make it possible to replace matching by usual field access as used in imperative languages like Java or C++. We believe that field access will be more efficient than matching.

We want our language also to support static overloading of function names. In C, defined functions always must have distinct names. In C++ and Java, the same name can be reused for different argument types. The compiler will select the definition that fits best to the arguments. This process is called overload resolution. In this paper, we study overload resolution only for field names (which is already present in C). The same technique is applicable to function calls.

We want some imperative features, because efficient run time handling of containers (also called collections) is difficult in functional setting. In order to obtain this, data must have efficient low level representation in memory. We allow a restricted form of assignment, only to top level variables and substructures at nesting depth one. In this way, side effects due to sharing can be avoided.

We will start by explaining the subtype system. At run time, it is used for inspecting data (formulas) so that we know which fields exist. The advantage of using subtypes for run time classification is that, once one has the mechanism for defining and verifying them, they are more general than structure matching, and can be used at many other points in the programming language, also at compile time.

As an example of our approach, suppose one has a propositional formula \( f \), and one wants to know if it constructed by `and`. One can either match it into `and(F_1, F_2)`, or define a subtype 'constructed by `and'`, and define two fields \( f_1, f_2 \) that can be accessed only on formulas refined by the subtype 'constructed by `and'`. The advantage of subtypes is that they can be easily combined by Boolean operators. For example, it is easy to define the subtype 'constructed by `and` or `or`', or 'atom' and use this as a condition in case analysis.

We will call the subtypes adjectives. In addition to adjectives that correspond to simple structural matchings, our adjective system also allows recursively defined subtypes. This makes it possible to define for example negation normal form or conjunctive normal form as a subtype of formulas. Although such recursively defined subtypes can be used for run time classification, they are not intended for this purpose, because evaluating them at run time would be costly. They are intended to be used for overload resolution at compile time. If one restricts oneself to adjectives of constant depth (like the ones above), they can be efficiently checked in constant time at run time.

Adjectives are somewhat similar to liquid types ([4]) or refinement types ([5]) but they have a different aim. Adjectives are not intended for checking operations whose safety depends on arithmetic. Because of this, adjectives do not need SMT solving. We prefer the term adjective over subtype or refinement type, because adjectives can overlap, and are frequently created for one time use. We also want to use adjectives for static type checking and overload resolution. In order to show that this is possible, we define a procedure that resolves ambiguous field references in the adjectives themselves.

We start by defining the type and adjective system in this section. In Section 2 we give the semantics, which can be used for evaluating adjectives on data, so that they can be used as a replacement for matching. In Section 3, we give a tableaux-based procedure for deciding satisfiability of adjectives. This procedure can be used for checking exhaustiveness of a given case analysis, for checking exclusiveness of cases, and for static type checking. In Section 4 we will apply the tableaux procedure on the problem of resolving
ambiguous field references in the adjective declarations themselves. A similar algorithm could be used for resolving ambiguous function calls. We define primitive types:

**Definition 1.1.** The primitive types are bool, char, nat, double, and selector. We assume that all primitive types, with the exception of selector, have an order < defined on them.

Type selector is a global enumeration type that consists of named constants. It is intended for defining subtypes (see Definition 1.2 below), or representing logical operators. Actually, these uses are the same because the logical operator of a formula defines a subtype at the same time. Enumeration types in Rust play a similar role. We write elements of because the logical operator of a formula defines a subtype at the same time. Enumeration types in Rust have an order < defined on them.

Next we define non-primitive types, and after that follow adjectives. Strictly seen, the definition of adjectives should come first, because types contain adjectives while adjectives do not contain types. We find, however, that this order would be unintuitive, because it would be hard to understand the usefulness of adjectives, without having seen the definition of types. Because of this, we start by introducing types:

**Definition 1.2.** Simple types are recursively defined as follows:

- A primitive type is a simple type.
- An identifier is a simple type (assuming that it has been defined as type, see Definition 1.4 below).
- If T is a simple type, A is an adjective, then T\circ A is also a simple type.

We recursively define compound types:

1. If v₁, ..., vₙ are identifiers, V₁, ..., Vₙ are simple types with n ≥ 0, then \((v₁ : V₁, ..., vₙ : Vₙ)\)' is a compound type.
2. If C is a compound type, v is an identifier and V is a simple type, then v : V, C is a compound type as well.
3. If C₁, ..., Cₘ with m ≥ 1 are compound types, A₁, ..., Aₘ are adjectives, s is an identifier, then s?(A₁ ⇒ C₁, ..., Aₘ ⇒ Cₘ) is a compound type as well.

In case 1, the fields v₁, ..., vₙ are called repeated fields. All other fields defined in cases 2 and 3 are called scalar fields. In case 3, identifier s (which must refer to an earlier defined field) is called the pivot field. Case 3 defines compound types where the existence of later fields depends on the adjective that applies to s.

We distinguish between simple and compound types, because we want every compound type to have a name, which solves some technical problems. It has no consequences for expressiveness, because one can always define a compound type, and use the name as simple type. This can be automatically done by the compiler, if needed. The meaning of T\circ A is: Type T refined by adjective A.

Compound types have a tree structure, where every path is a possible realization of the type. The tree branches whenever case 3 is used. In that case, identifier s should refer to a field that occurs before it on the same path. We cannot impose this in Definition 1.2, because trees are constructed bottom up, and the field s must be defined above it in an application of case 2. We give a simple example of a compound type:

**Example 1.3.**

\[ T = s: \text{selector}(\text{?simp}\text{?vect}), \text{?simp} \Rightarrow a: A, b: B, (\text{?vect} \Rightarrow c: C, (d: D, e: E))' \]

Example 1.3 is purely technical. Type T is not intended to be meaningful. Also note that Definition 1.2 is awkward to use when defining types in practice because it is intended as technical representation. It is not intended as the syntax with which the programmer will define types. The first field s is a selector, which must be either ?simp or ?vect. In case it is ?simp, there are two more scalar fields a of type A, and b of type B. In case s equals ?vect, there is one more scalar field c of type C, and an unbounded number of repeated fields...
$d$ of type $D$ and $e$ of type $E$. They must be accessed with array notation $d[i]$ or $e[i]$. They are implemented as vectors in $C^{++}$, which means that the repeated part can dynamically grow and shrink, and an object of type $T$ will be reallocated when it runs out of capacity. Because repeated fields are always last, offsets of scalar fields do not depend on the number of repeated fields present.

For the moment, we assume that different fields of different types have distinct names. This is an unrealistic restriction for a real programming language, because field names can be reused in different types. In Section 4 we will allow reuse of field names between different options of a type, and between different types. We will give a procedure that is able to resolve such ambiguous field references into unambiguous references.

Type definitions need to be stored in a mapping. It is convenient to use two distinct mappings, one for the simple types, and one for the compound types:

**Definition 1.4.** We define two mappings $\Sigma_S$ and $\Sigma_C$. The first mapping maps identifiers to simple types, and the second mapping maps identifiers to compound types. If $\Sigma_S$ contains a value for $v$, we write $\Sigma_S(v)$ for the value. Similarly, if $\Sigma_C$ contains a value for $v$, we write $\Sigma_C(v)$ for the value. We assume that $\Sigma_S$ and $\Sigma_C$ do not contain a value for the same variable $v$. We also assume that $\Sigma_S$ is cycle free.

An example of $\Sigma_S$ not being cycle-free would be when $\Sigma_S(v_1) = v_2 \circ A_2$, and $\Sigma_S(v_2) = v_1 \circ A_1$. Cycles inside $\Sigma_C$ are unproblematic, see for example the definition of prop below in Example 1.10.

This completes the definition of the types, next we define adjectives:

**Definition 1.5.** We recursively define adjectives, starting with primitive adjective constructors:

- If $c$ is a constant of one of the primitive types bool, char, nat, double, then $c$ and $c^A$ are adjectives.
- If $c$ is a constant of type selector, then $c$ is an adjective.
- empty is an adjective.
- If $v$ is an identifier, then $v$ is an adjective.

Next we define recursive adjective constructors:

- If $f$ is an identifier, and $A$ is an adjective, then $f(A)$ is an adjective.
- If $A$ is an adjective, then first($A$) and rest($A$) are adjectives.
- If $A$ is an adjective, then $\forall A$ and $\exists A$ are adjectives.

Adjectives of any type can be combined by propositional operators:

- If $A_1, \ldots, A_n$ are adjectives, then $A_1 \lor \cdots \lor A_n$, and $A_1 \land \cdots \land A_n$ are also adjectives.
- If $A$ is an adjective, then $\neg A$ is also an adjective.

The intuitive meaning of the adjective $c$ is : equal to $c$. The intuitive meaning of $c^A$ is : greater or equal to $c$. The intuitive meaning of $f(A)$ is: whose $f$-field satisfies $A$. If $f$ is a repeated field, then $A$ must be equal to empty or have one of the forms $\forall A$, $\exists A$, first($A$), or rest($A$). The meaning of empty is that no repeated fields are present. The meaning of $f(\forall A)$ is that all repetitions of $f$ must satisfy $A$. Similarly $f(\exists A)$ means that at least one repetition of $f$ must satisfy $A$. The meaning of first($A$) is that $f[1]$ must satisfy $A$. The meaning of rest($A$) is that all repetitions of $f$, except for possibly the first, must satisfy $A$.

**Definition 1.6.** We assume a mapping $\Sigma_A$ that maps identifiers to pairs of form $(T, A)$, where $T$ is a simple type, and $A$ is an adjective. If $\Sigma_A$ contains a value for $v$, we write $\Sigma_A(v)$ for the value.

The intuitive meaning of $\Sigma_A(v) = (T, A)$ is: Identifier $v$ is defined on type $T$ as adjective $A$.

**Definition 1.7.** We write $A[v]$ for an adjective that contains $v$ somewhere not inside a subformula of form $f(\cdot)$. We define a cycle as a sequence of identifiers, s.t.

$$\Sigma_A(v_1) = (T_2, A_2[v_2]), \ldots, \Sigma_A(v_{n-1}) = (T_n, A_n[v_n]), \Sigma_A(v_n) = (T_1, A_1[v_1]).$$
Cycles are logically problematic when they involve negation. Because the intended meaning of an adjective is an inductively defined predicate, an adjective defined through a cycle involving negation, would be ill-defined. This is similar to the situation in logic programming ([7]). Although it would be possible to adapt Definitions 2.2 and 3.8 to monotonic cycles, we are not aware of a meaningful use of it. Therefore, we think it is better to forbid cycles altogether. In the rest of this paper, we will assume that $\Sigma_A$ is cycle free. We give some examples of type and adjective definitions:

**Example 1.8.** The type of complex numbers can be defined as follows:

$$\text{complex} := \text{re: double, im: double, (\_)}.$$  

Complex numbers have two scalar fields re and im, and no repeated fields.

**Example 1.9.** Natural numbers can be defined as follows:

$$\text{nat} := \text{sel: selector, sel(} \begin{cases} \text{?zero} & \Rightarrow (\_), \\
\text{?succ} & \Rightarrow \text{pred: nat, (\_)} \end{cases} \text{)}.$$  

Adjectives odd and even can be defined on nat in mutual recursion:

$$\text{even} := \text{nat(}\text{?zero)} \lor (\text{sel(}\text{?succ)} \land \text{pred(odd)}) \text{)} \text{)}$$  
$$\text{odd} := \text{nat(}\text{?succ)} \land \text{pred(even)}) \text{)}$$

**Example 1.10.** We define propositional logic:

$$\text{prop} := \text{op: selector, op(} \begin{cases} \text{?var} & \Rightarrow (\text{c: char}) \text{),} \\
\text{?not} & \Rightarrow \text{sub: prop, (\_)} \text{),} \\
\text{?implies\lor?equiv} & \Rightarrow \text{sub}_1: \text{prop, sub}_2: \text{prop, (\_)} \text{),} \\
\text{?and\lor?or} & \Rightarrow (\text{sub}_n: \text{prop}) \text{)} \end{cases} \text{)}.$$  

There is no need to define $\top$ or $\bot$ separately, because $\top = \bigwedge \emptyset$, and $\bot = \bigvee \emptyset$. A variable consists of ?var combined with a finite number of characters. A negation consists of ?not combined with a single subformula called sub. A conjunction or disjunction consists of ?and or ?or, followed by an arbitrary number of subformulas called sub$_n$. We define a few adjectives on prop:

**Example 1.11.** 'Constructed by ?and' can be expressed as op(?and). 'Constructed by ?and or ?or' can be expressed as op(?and\lor?or). The property of being an atom can be expressed by

$$\text{atom} := (\text{prop, op(var)} \text{)}.$$  

The property of being a literal can be expressed by

$$\text{literal} := (\text{prop, atom} \lor (\text{op(?not)} \land \text{sub(?atom)})) \text{).}$$

Negation normal form (NNF) can be expressed by:

$$\text{nnf} := (\text{prop, } \bigvee \left\{ \text{literal, op(?and}\lor?or)} \land \text{sub}_n(\forall \text{nnf}) \right\} \text{).}$$

Conjunctive normal form (CNF) can be defined as follows:

$$\text{cnf} := (\text{prop, op(?or)} \land \text{sub}_n(\forall (\text{op(?and)} \land \text{sub}_n(\forall \text{literal}))) \text{).}$$
We define the set of data trees $D$, and the set $\overline{D}$ of finite sequences of data trees, in simultaneous recursion:

- If $d$ is an element of one of the primitive types $T$, defined in Definition 1.1, then $d \in D$.
- If $d_1, \ldots, d_n$ ($n \geq 0$) is a finite sequence of data trees, then $(d_1, \ldots, d_n) \in \overline{D}$. We will write $\|D\|$ for the length $n$ of $\overline{D}$.
- If $f_1, \ldots, f_n$ are pairwise distinct identifiers, and each $d_i \in D \cup \overline{D}$, then $\{(f_1, d_1), \ldots, (f_n, d_n)\} \in \overline{D}$.

Data trees of the third type can be viewed as partial functions mapping each $f_i$ to $d_i$, such that $d_i$ is either a data tree or a sequence of data trees. When $d$ is of the third type, we write $d.f$ for the value attached to $f$ when it exists. If $d.f \in \overline{D}$, we write $d.f[i]$ for its $i$-th element, assuming that $i \leq \|d.f\|$.

We will define when a data tree has a given type, and when a data tree (or sequence of data trees) satisfies a given adjective. Types may contain adjectives, but adjectives do not contain types. Therefore, we start by defining when a data tree (or sequence of data trees) makes an adjective true:

**Definition 1.12.** Let $f$ be an exact identifier that is declared as field in a compound type $C$. We recursively define the precondition of $f$, written as $\text{PRE}(C, f)$, as follows:

- If $C$ has form $(v_1 : V_1, \ldots, v_n : V_n)$, then $\text{PRE}(f, C) = \top$.
- If $C$ has form $v : V$, then $\text{PRE}(f, C) = \top$. If $f$ is declared in $C$, then $\text{PRE}(f, C) = \text{PRE}(f, C')$.
- If $C$ has form $s?(A_1 \Rightarrow C_1, \ldots, A_m \Rightarrow C_m)$, and $f = s$, then $\text{PRE}(f, C) = A_1 \lor \cdots \lor A_m$. Otherwise, $f$ must be declared in exactly one $C_i$. We define $\text{PRE}(f, C) = s(A_i) \land \text{PRE}(f, C_i)$.

Since $f$ can occur in $\Sigma_C$ only once, it is possible to write $\text{PRE}(f)$ instead of $\text{PRE}(f, C)$, where $C$ is the compound type that contains $f$.

2. **Semantics of Types and Adjectives**

We will describe the semantics of adjectives and types using a simplified high level representation of data. We assume that data are represented by trees whose subtrees are labeled with field names. This representation is convenient for defining the semantics without having to consider low level details like memory lay out. In the implementation we use a low level representation, where scalar fields have a fixed offset, and repeated fields have an offset that can be computed by multiplying the index with the size of the repeated part, and adding a base offset.

**Definition 2.1.** We define the set of data trees $D$, and the set $\overline{D}$ of finite sequences of data trees, in simultaneous recursion:

- If $d$ is an element of one of the primitive types $T$, defined in Definition 1.1, then $d \in D$.
- If $d_1, \ldots, d_n$ ($n \geq 0$) is a finite sequence of data trees, then $(d_1, \ldots, d_n) \in \overline{D}$. We will write $\|D\|$ for the length $n$ of $\overline{D}$.
- If $f_1, \ldots, f_n$ are pairwise distinct identifiers, and each $d_i \in D \cup \overline{D}$, then $\{(f_1, d_1), \ldots, (f_n, d_n)\} \in \overline{D}$.

Data trees of the third type can be viewed as partial functions mapping each $f_i$ to $d_i$, such that $d_i$ is either a data tree or a sequence of data trees. When $d$ is of the third type, we write $d.f$ for the value attached to $f$ when it exists. If $d.f \in \overline{D}$, we write $d.f[i]$ for its $i$-th element, assuming that $i \leq \|d.f\|$.

We will define when a data tree has a given type, and when a data tree (or sequence of data trees) satisfies a given adjective. Types may contain adjectives, but adjectives do not contain types. Therefore, we start by defining when a data tree (or sequence of data trees) makes an adjective true:

**Definition 2.2.** We define in simultaneous recursion when a data tree makes an adjective true, and when a sequence of data trees makes an adjective true. We use the notation $d \vdash A$ for both cases.

We first consider the cases where $d \in D$:

- If $c$ is a constant, then $d \vdash c$ iff $d = c$, and $d \vdash c^e$ if $d \geq c$.
- If $f$ is an identifier, s.t. $d.f$ exists, then
  - if $d.f$ is a single data tree, then $d \vdash f(A)$ iff $d.f \vdash A$.
  - if $d.f$ is a sequence of data trees, $d \vdash f(A)$ iff $d.f \vdash A$.

The two cases appear to be the same, but $d.f$ have different types, hence we prefer to list them separately.

- If $v$ is an identifier, s.t. $\Sigma_A(v)$ is defined, and $d \in D$, then $d \vdash v$ iff $d \vdash \Sigma_A(v)$.

Next we list the cases where $d \in \overline{D}$:

- $(d_1, \ldots, d_n) \vdash \text{empty}$ iff $n = 0$.
- $(d_1, \ldots, d_n) \vdash \forall A$ iff for every $d_i$ we have $d_i \vdash A$. 

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We recursively define when a data tree \( d \) has type \( T \). We use notation \( d : T \). We first list the cases where \( T \) is simple:

- For a primitive type \( T \), we define \( d : T \) as \( d \in T \).
- If \( v \) is an identifier that is defined in \( \Sigma_S \), then \( d : v \) iff \( d : \Sigma_S(v) \).
- If \( v \) is an identifier that is defined in \( \Sigma_C \), then \( d : v \) iff \( d : \Sigma_C(v) \).
- \( d : (T \cdot A) \) iff \( d : T \) and \( d \vdash A \).

Next we list the cases where \( T \) is compound:

- \( d : (\forall \_1, ..., \forall \_n : \forall \_1 \cdot \forall \_n) \) iff either
  - \( n = 0 \), or
  - \( n > 0 \) and \( d.\forall \_1, ..., d.\forall \_n \) are all defined, are all in \( \overline{D} \), have the same length \( L = \|d.\forall \_1\| = \cdots = \|d.\forall \_n\| \),
    and for every \( i (1 \leq i \leq n) \) and \( j (1 \leq j \leq L) \), we have \( d.\forall \_i[j] : \forall \_i \).
- \( d : (v : V, C) \) iff \( d.v \) is defined, \( d.v : V \), and \( d : C \).
- \( d : s ? (A_1 \Rightarrow C_1, ..., A_m \Rightarrow C_m) \) iff \( d.s \) is defined, in \( D \), and there is exactly one \( 1 \leq i \leq m \), s.t. \( d.s \vdash A_i \) and \( d : C_i \).

**Example 2.4.** Following up on Example 1.8, we can see that \( \{ (\text{re}, 1.0), (\text{im}, 2.0) \} \) has type complex.

**Example 2.5.** Using the declaration of prop in Example 1.10, the propositional formula \( p \) will be represented by a data tree \( d_p \) with fields

- \( d_p.\text{op} = ?\text{atom} \)
- \( d_p.\text{c} = (\text{`}p\text{'}\text{)} \)

The propositional formula \( q \) will be represented by \( d_q \) with fields

- \( d_q.\text{op} = ?\text{atom} \)
- \( d_q.\text{c} = (\text{`}q\text{'}\text{)} \)

Similarly, the propositional formula \( r \) will be represented by \( d_r \) with

- \( d_r.\text{op} = ?\text{atom} \)
- \( d_r.\text{c} = (\text{`}r\text{'}\text{)} \)

The formula \( q \lor r \) will be represented by \( d \) with

- \( d.\text{op} = \text{'}\text{or}\text{'} \)
- \( d.\text{sub} = (d_q, d_r) \)

Finally, the formula \( p \rightarrow (q \lor r) \) will be represented by \( d' \) with

- \( d.\text{op} = \text{'}\text{implies}\text{'} \)
- \( d.\text{sub}_1 = d_p \)
- \( d.\text{sub}_2 = d' \)
3. A Tableaux Calculus

In this section we define a calculus for checking satisfiability of adjectives. This is sufficient to answer all logical questions that are needed for the implementation of our programming language. In order to show that adjectives \( A_1, \ldots, A_n \) cover all possible cases in a switch, it is sufficient to show that \( A \land \neg A_1 \land \cdots \land \neg A_n \) is unsatisfiable, where \( A \) is the adjective part of the type of the switch expression. In order to show that a switch has no overlapping cases, it is sufficient to show that \( A \land A_i \land A_j \) is unsatisfiable for all distinct \( i \) and \( j \). In order to show that a function \( f_i \) defined on adjective \( A_1 \) is a better fit than function \( f_2 \) defined on adjective \( A_2 \), it is sufficient to show that \( A_1 \land \neg A_2 \) is unsatisfiable, while \( A_2 \land \neg A_1 \) is satisfiable.

In order to use our calculus, types need to be decomposed into their adjective part, and the part that specifies their implementation. The calculus needs the adjective part of a type, but does not use the implementation. We will define two functions that decompose a simple type into their adjective and implementation components.

**Definition 3.1.** An implementation type is either a primitive type, or an identifier defined in \( \Sigma_C \), i.e. the name of a compound type. For a simple type \( T \), we define \( \text{ADJ}(T) \) and \( \text{IMPL}(T) \) as follows:

- If \( T \) is primitive then \( \text{ADJ}(T) = \top \), and \( \text{IMPL}(T) = t \).
- If \( v \) is an identifier defined in \( \Sigma_C \), then \( \text{ADJ}(v) = \top \), and \( \text{IMPL}(v) = v \).
- \( \text{ADJ}(T \circ A) = \text{ADJ}(T) \land A \), and \( \text{IMPL}(T \circ A) = \text{IMPL}(T) \).

For example, if \( T = \text{prop} \circ \text{nff} \circ \text{cnf} \), then \( \text{ADJ}(T) = \text{nff} \land \text{cnf} \), and \( \text{IMPL}(T) = \text{prop} \).

**Theorem 3.2.** If \( T \) is a simple type, and \( d \) is a data tree, then \( d \models T \iff d \models \text{IMPL}(T) \) and \( d \models \text{ADJ}(T) \).

The theorem can be easily verified by applying the rules of Definition 2.2.

At this point, we can define a simple interface of our tableaux procedure. Its input is just a single type \( T \). The procedure establishes that there exists no data term \( d \) with type \( T \). The applications that we mentioned in the introduction, can be obtained as follows:

- Establish that there exists no data term \( d \) of type \( T \) that satisfies adjective \( A \) : Call the procedure with \( T \circ A \).
- Establish that every data term \( d \) of type \( T \), that satisfies \( A \), must also satisfy \( B \) : Call the procedure with \( T \circ A \circ \neg B \).

The procedure works on the following normal form, which is similar to negation normal form:

**Definition 3.3.** We recursively define when an adjective is in path normal form (PNF):

- If \( c \) is a constant of a primitive type, then \( c, \neg c, c^\circ \) and \( \neg c^\circ \) are in PNF. In the last two cases, \( c \) cannot have type selector.
- empty and \( \neg \text{empty} \) are in PNF.
- If \( v \) is an identifier, then \( v \) and \( \neg v \) are in PNF.
- If \( A \) is in PNF, then \( f(A), \text{first}(A), \text{rest}(A), \forall A, \) and \( \exists A \) are in PNF.
- If \( A_1, \ldots, A_n \) (\( n \geq 0 \)) are in PNF, then \( A_1 \lor \cdots \lor A_n \) and \( A_1 \land \cdots \land A_n \) are in PNF.

**Definition 3.4.** An adjective \( A \) can be brought into PNF by calling \( \text{PNF}(A, \text{pos}) \), where \( \text{PNF}(A, p) \) is defined by cases as follows:

- If \( A \) has form \( c, \neg c^\circ, \text{empty}, \) or \( v \), then
  \[
  \text{PNF}(A, \text{pos}) = A \\
  \text{PNF}(A, \text{neg}) = \neg A
  \]
For the remaining cases:

\[
\begin{align*}
\text{PNF}(A_1 \lor \cdots \lor A_n, \text{pos}) &= \text{PNF}(A_1, \text{pos}) \lor \cdots \lor \text{PNF}(A_n, \text{pos}) \\
\text{PNF}(A_1 \lor \cdots \lor A_n, \text{neg}) &= \text{PNF}(A_1, \text{neg}) \land \cdots \land \text{PNF}(A_n, \text{neg}) \\
\text{PNF}(A_1 \land \cdots \land A_n, \text{pos}) &= \text{PNF}(A_1, \text{pos}) \land \cdots \land \text{PNF}(A_n, \text{pos}) \\
\text{PNF}(A_1 \land \cdots \land A_n, \text{neg}) &= \text{PNF}(A_1, \text{neg}) \lor \cdots \lor \text{PNF}(A_n, \text{neg}) \\
\text{PNF}(f(A), p) &= f(\text{PNF}(A, p)) \text{ for } p \in \{\text{pos}, \text{neg}\} \\
\text{PNF}(\text{first}(A), p) &= \text{first}(\text{PNF}(A, p)) \text{ for } p \in \{\text{pos}, \text{neg}\} \\
\text{PNF}(\text{rest}(A), p) &= \text{rest}(\text{PNF}(A, p)) \text{ for } p \in \{\text{pos}, \text{neg}\} \\
\text{PNF}(\exists A, \text{pos}) &= \exists \text{PNF}(A, \text{pos}) \\
\text{PNF}(\exists A, \text{neg}) &= \forall \text{PNF}(A, \text{neg}) \\
\text{PNF}(\forall A, \text{pos}) &= \forall \text{PNF}(A, \text{pos}) \\
\text{PNF}(\forall A, \text{neg}) &= \exists \text{PNF}(A, \text{neg}) \\
\text{PNF}(\neg A, \text{pos}) &= \text{PNF}(A, \text{neg}) \\
\text{PNF}(\neg A, \text{neg}) &= \text{PNF}(A, \text{pos})
\end{align*}
\]

The following theorem can be straightforwardly proven, using Definition 2.2:

**Theorem 3.5.** Let \( A \) be an adjective. For every data tree \( d \), we have \( d \vdash A \iff d \models \text{PNF}(A, \text{pos}) \). Similarly, \( d \vdash \neg A \iff d \models \text{PNF}(A, \text{neg}) \).

The path normal form of \( \neg (A \lor f(\text{rest}(\forall \neg B))) \) equals \( \neg A \land f(\text{rest}B) \).

**Definition 3.6.** We define a path as a finite sequence \((f_1, \ldots, f_n)\) with \(n \geq 0\), where each \( f_i \) is either an identifier (representing a field name), or an element of \( \{\exists, \forall, \text{first}, \text{rest}\} \). We write \( \emptyset \) for the empty path, and use notation \( \pi. f \) for extending path \( \pi \) with \( f \). If path \( \pi' \) can be obtained from path \( \pi \) by zero or more extensions, we call \( \pi \) a prefix of \( \pi' \). We call \( \pi \) a strict prefix of \( \pi' \) if at least one extension was used.

We assume that there exists a total order \( \prec \) on paths with the property that if \( \pi_1 \) is a strict prefix of \( \pi_2 \), then \( \pi_1 \prec \pi_2 \). Such an order can be easily obtained by fixing a total order on all possible \( f \), and using the alphabetic, lexicographic extension.

**Definition 3.7.** An adjective stack \( S \) is a finite sequence of triples \((\pi_i, A_i, \lambda_i)\), where each \( \pi_i \) is a path, each \( A_i \) is an adjective, and each \( \lambda_i \in \mathcal{N} \). We write \( S[\pi] \) for the set \( \{ A \mid \exists \lambda \text{ s.t. } (\pi, A, \lambda) \text{ occurs in } S \} \) and \( S^\lambda[\pi] \) for the set \( \{ A \mid (\pi, A, \lambda) \text{ occurs in } S \} \).

The attribute \( \lambda_i \) stores the level of \( A_i \) on path \( \pi_i \). More precisely, an adjective that was obtained from a formula on a strict prefix of \( \pi_i \) will have level 0. An adjective that was obtained from an adjective with level \( \lambda \) on the same path, receives level \( \lambda + 1 \).

We define the tableaux procedure. In order to obtain termination, the procedure uses a blocking rule. Whenever a new path \( \pi \) is entered, it checks that there is no strict prefix \( \pi' \) of \( \pi \) containing a set of adjectives that are included in the formulas on the current path. For example, if path \( f \) contains \( A, B \), and path \( f. g \) contains formulas \( A, B, C \), we can close the branch. This rule is correct, because for every data tree \( d \), the subtree \( d. f. g \) is a subtree of \( d. f \). If it is possible that \( d. f. g \models A \land B \land C \), then one can replace \( d. f \) by \( d. f. g \) and obtain a smaller data tree, for which \( d. f \models A \land B \).

The tableaux procedure always tries to extend in deterministic fashion first. If that does not result in a conflict, it selects a path \( \pi \) and tries all possible non-deterministic choices on this path. It keeps on trying until it either has explored all choices, or obtained a consistent stack \( S \).

**Definition 3.8.** We define a procedure that tries to establish that no data term \( d \) can have type \( T \). The procedure starts by creating a stack with one element:

\[
S = ((\epsilon, \text{ADJ}(T), 0)).
\]
After that, it calls $b = \text{deterministic}(0)$. If $b = \perp$, it returns $\perp$. Otherwise, it returns $\text{nondeterministic}(e)$. We list the subprocedures, starting with the procedure $\text{deterministic}(d)$. All subprocedures have access to the stack $S$. We write $|S|$ for the size of the stack, and write the $i$-th element in the form $S_i = (\pi_i, A_i, \lambda_i)$.

Procedure $\text{deterministic}(d)$ must be called with a natural number $1 \leq d \leq |S|$. It returns $\perp$ or $\top$, with $\perp$ indicating contradiction. The implementation is as follows:

1. Set $c = d$.
2. If $c \leq |S|$, then
   • If $A$ has form $v$ with $v$ an identifier that has no definition in $\Sigma_A$, then return $\perp$.
   • If $A$ has form $\neg c^x$ with $c$ the minimal element in its type, then return $\perp$.
   • For all $i < c$ with $\pi_i = \pi_c$, check whether $A_i$ and $A_c$ are in conflict, using the rules listed below. If they are in conflict, return $\perp$. Otherwise try the next $\pi_i$. The rules are:
     - A complementary pair $A, \neg A$ is in conflict.
     - A pair $c_1, c_2$ of any primitive type with $c_1 \neq c_2$ is in conflict.
     - A pair of form $\neg c_1^x, c_2^x$ with $c_1 \leq c_2$ is in conflict.
     - A pair of form $c_1^x, c_2$ with $c_1 > c_2$, is in conflict.
     - A pair of form $\neg c_1^x, c_2$ with $c_1 \leq c_2$ is in conflict.
   • Assign $c = c + 1$ and go back to step 1.
3. If $d \leq |S|$, check if one of the deterministic consequence rules in the table below is applicable on $A_d$. If yes, then for every deterministic consequence $A'$, push $(\pi_d, A', \lambda_d + 1)$ to $S$.

   \[
   \begin{array}{ll}
   \forall A & \Rightarrow \text{PFN}(A, \text{pos}) \quad \text{if } \Sigma_A \text{ contains a definition of form } v := (T, A) \\
   \exists A & \Rightarrow \text{PFN}(A, \text{neg}) \quad \text{if } \Sigma_A \text{ contains a definition of form } v := (T, A) \\
   \forall A & \Rightarrow \text{empty} \lor (\text{first}(A) \land \text{rest}(\forall A)) \\
   \exists A & \Rightarrow \neg \text{empty} \land (\text{first}(A) \lor \text{rest}(\exists A)) \\
   A_1 \land \cdots \land A_n & \Rightarrow A_i \text{ for each } 1 \leq i \leq n
   \end{array}
   \]
4. If $A_d$ has form $f(A')$, then push $(\pi_d, f(A'), 0)$ and $(\pi_d, f, \text{ADJ}(T), 0)$ to $S$, where $\text{ADJ}(T)$ is the adjective component of the type $T$ of $f$.

Procedure $\text{nondeterministic}(\pi)$ must be called with a path $\pi$ that occurs in $S$. It returns $\perp$ or $\top$, with $\perp$ indicating contradiction. The implementation is as follows:

1. For every strict prefix $\pi'$ of $\pi$, do:
   (a) if $S_0[\pi'] \subseteq S[\pi]$, then return $\perp$. (This is the blocking rule, mentioned above.)
2. Call $\text{nondeterministic}(\pi, 1)$.

Procedure $\text{nondeterministic}(\pi, n)$ must be called with a path $\pi$ that occurs in $S$, and with $1 \leq n \leq |S|$. It returns $\perp$ or $\top$, with $\perp$ indicating contradiction. The implementation is as follows:

1. Find the smallest $n' \geq n$, such that $A_{n'}$ has form $A_1 \lor \cdots \lor A_m$ with $m \geq 2$.
2. If $n' = n + 1$ was found in the previous step, then find the next path $\pi' > \pi$ occurring in $S$, using the alphabetic lexicographic order $>$. Call $b = \text{nondeterministic}(\pi')$ and return $b$. If no path could be found, return $\perp$.
3. Set $n = n'$. We know that $S_n$ has form $(\pi, A_1 \lor \cdots \lor A_m, \lambda)$, with $m \geq 2$. For $i$ from $1$ to $n$, do the following:
   • Set $s = |S|$. Push $(\pi, A_i, \lambda + 1)$ on the stack.
   • Call $b = \text{deterministic}(s)$. If $b \neq \perp$, call $b = \text{nondeterministic}(\pi, n + 1)$. If $b = \top$ return $\top$.
   • Restore $S$ to length $s$.

Procedure $\text{nondeterministic}(\pi)$ implements a loop checker which guarantees termination. Its correctness is based on the fact that, if a satisfying data tree can be found, it can be pruned into a data tree in which the calculus does not repeat initial states. We state the following without proof:
**Theorem 3.9.** The tableaux calculus of Definition 3.8 terminates.

In order to establish correctness of Theorem 3.9, it is sufficient establish that every branch is finite. During search, the procedure will only introduce subadjectives of the initial type, combined with subadjectives of identifiers defined in $\Sigma_A$. Therefore, blocking must eventually happen.

**Theorem 3.10.** The procedure of Definition 3.8 is complete.

Доказательство. The proof will be a bit informal. We have to prove that if no data tree $d$ with $d : T$ exists, then the procedure of Definition 3.8 will reject $T$. We will prove the converse: If $T$ is not rejected, then there exists a data tree $d$, s.t. $d : T$. Suppose that $T$ is not rejected. This implies that the procedure terminates with an open branch. Let $S = (\pi_1, A_1, \lambda_1), \ldots, (\pi_n, A_n, \lambda_n)$ be the state of the stack with which the tableaux procedure terminated.

For a path occurring in $S$, let

$$S(\pi) = \{ A \mid S \text{ contains a tuple } (\pi_i, A_i, \lambda_i), \text{s.t. } \pi_i = \pi \text{ and } A_i \text{ has form } c, c^{\varepsilon} \text{ or } \neg c^{\varepsilon} \}.$$ 

For an arbitrary constant $c$, define

$$I(c) = \{ c \} \quad I(c^{\varepsilon}) = \{ c' \mid c' \geq c \}, \quad I(\neg c^{\varepsilon}) = \{ c' \mid c' < c \}.$$ 

It is easily checked that for every $\pi$ occurring in $S$, we have $\bigcap\{I(A) \mid A \in S(\pi)\} \neq \emptyset$, because otherwise the branch would be closed. Hence, we can select a constant $c_\pi$ from each such set. Let $A_S$ be the adjective

$$A_S = \bigwedge\{ \pi(c_\pi) \mid \pi \text{ occurs in } S \}.$$ 

The adjective $A_S$ states that for a given data tree $d$, selecting field sequence $\pi$ will result in constant $c_\pi$. It is easy to construct a data tree $d_S$, s.t. $d_S \vDash A_S$ by starting with the required constants, and combining them as required by $A_S$. We show by backward induction (that is from $n$ towards 1), that

$$d_S \vDash \pi_1(A_1) \land \cdots \land \pi_n(A_n).$$ 

Assume that we already have established that $d_S \vDash \pi_{i+1}(A_{i+1}) \land \cdots \land \pi_n(A_n)$. We proceed by case analysis on the form of $A_i$. We need to consider only the forms of $A_i$ that do not close the branch.

- If $A_i$ is an identifier $v$, then we know that $v$ has a definition $(T, A')$ in $\Sigma_A$, since otherwise the branch would have been closed. Hence we know that PNF$(A', \text{pos})$ occurs among $A_{i+1}, \ldots, A_n$. As a consequence, $d_S \vDash \text{PNF}(A', \text{pos})$. By Theorem 3.5, we know that $d_S \vDash A'$.
- If $A_i$ is a negated identifier $\neg v$, then if $v$ has no definition in $\Sigma_A$, we have $d_S \not\vDash \pi_i(v)$, so that $d_S \vDash \pi_i(\neg v)$. If $v$ has a definition $(T, A')$ in $\Sigma_A$, then PNF$(A', \text{neg})$ occurs among $A_{i+1}, \ldots, A_n$. By induction, $d_S \vDash \text{PNF}(A', \text{neg})$. By Theorem 3.5, we have $d_S \vDash \neg A'$.
- The other cases can be obtained by inspecting the cases in Definition 2.2.

\[ \square \]

**Theorem 3.11.** The procedure of Definition 3.8 is sound.

Доказательство. In order to prove soundness, one must prove that if the procedure rejects a type $T$, then there is no data tree $d$ with $d : T$. We will prove the converse: If there is a data tree $d$ with $d : T$, then $T$ will not be rejected by the procedure.

This would be trivial, if the blocking rule would not exist. It is easy to show that if $d : T$, there exists a stack $S = (\pi_1, A_1, \lambda_1), \ldots, (\pi_n, A_n, \lambda_n)$, with $(\pi_1, A_1, \lambda_1) = (c, \text{ADJ}(T), 0)$, that represents an open branch of the tableaux procedure when the blocking rule is not used.
Now assume that $S$ will be closed when the blocking rule can be used. We will show that there exists a shorter stack $S'$ which also starts with $S_1' = (\varepsilon, \text{ADJ}(T), 0)$, and which also represents an open branch of the tableaux procedure when the blocking rule cannot be used. Repeating this process will result in a stack that does not contain any more applications of the blocking rule.

Let $S$ be a stack with $S_1 = (\varepsilon, \text{ADJ}(T), 0)$, taken from an open branch of the tableaux procedure, when blocking is not used. Suppose that somewhere in $S$, the blocking rule would be applicable. This means that there exist $\pi$ and $\pi'$, s.t. $\pi'$ is a strict prefix of $\pi$, and $S_0[\pi'] \subseteq S[\pi]$. We prune $S$ as follows:

- Remove every $(\pi_i, A_i, \lambda_i)$, s.t. $\pi'$ is a prefix of $\pi_i$ and $\pi_i$ is a strict prefix of $\pi$.
- For every $(\pi_i, A_i, \lambda_i)$, s.t. $\pi$ is a prefix of $\pi_i$, write $\pi_i$ as $\pi \cdot \pi_i'$, and replace it by $\pi' \cdot \pi_i'$.

It can be checked that the resulting $S'$ is shorter, because there is at least one $(\pi_i, A_i, \lambda_i)$ with $\pi_i = \pi'$ which will be removed. Moreover, $S'$ still represents an open branch of the tableaux procedure without blocking. Hence we can continue the procedure and obtain an open branch on which the blocking rule is not applicable.

\section{Resolving Overloads in Types and Adjectives}

Until now we have insisted that field and adjective names are always unique. In Example 1.10, we used $\text{sub}_1$, $\text{sub}_2$, and $\text{sub}_3$ as different variations of the name $\text{sub}$, dependent on whether we were taking a subformula of a negated formula, a formula constructed by a binary operator, or a formula constructed by an $n$-ary operator.

Similarly, we did not consider reuse of adjective names between different types. In reality, it is perfectly possible to have different types of formulas, for example modal, first-order and propositional, and define different $\text{nnf}$ adjectives on each of them.

Modern programming languages like $C^+$ or Java allow the use of ambiguous names which are made unambiguous by the compiler. For example, in $C^+$, one can define different print operators $<<$ on different types, and the compiler will pick the right one, when the programmer writes $<<$. This is called \textit{overload resolution}. Without overload resolution, the programmer needs to invent a new name for every type that needs to be printed. For example, in $C$ one has to include the type in the name of a print function, like \text{printfol} or \text{printmodal}.

We want overload resolution in our programming language: it should be possible that the user defines different $\text{nnf}$ adjectives on different types and reuses the same name ‘$\text{nnf}$’ for them. Similarly, we want that the user can call all variations $\text{sub}$, $\text{sub}_1$, $\text{sub}_2$ and $\text{sub}_n$, just ‘$\text{sub}$’.

Concretely, we want overload resolution on field names, adjective names, and function names. There will be no overload on type names, because we think it is unfeasable, and allowing ambiguous type names would result in ambiguous code.

In order to handle overload resolution, we introduce what we call \textit{inexact identifiers}. Inexact identifiers are the identifiers that are used by the programmer in the program. We will call the identifiers that we have been using until now, \textit{exact identifiers}. We assume a function $\nu^?$ that maps exact identifiers $\nu$ to their inexact representations. The inexact representations are used in the program. As an example, one can introduce an inexact name $\text{sub}$ and set $\text{sub}^? = \text{sub}_1^? = \text{sub}_2^? = \text{sub}_n^? = \text{sub}$. Whenever identifier ‘$\text{sub}$’ occurs in the program, the compiler has to find out which of the exact variations of sub is meant. We define inexact identifiers:

\textbf{Definition 4.1.} We assume an infinite set of inexact identifiers. We assume a map $\nu^?$ that maps exact identifiers $\nu$ to inexact identifiers $\nu$ to inexact identifiers.

Now we explain how inexact identifiers are used in the program. In order to do that, we modify the definitions of type and of adjective. In Definition 1.2, in the definition of compound types, we make the following modifications:

1. the identifiers $\nu_1, \ldots, \nu_n$, are inexact.
2. the identifier $v$ is inexact.
3. identifier $s$ is inexact.

In Definition 1.5, we will allow the identifiers $v$ and $f$ to be inexact. Apart from that, there are no changes.

**Example 4.2.** We define propositional and multimodal logic:

\[
\begin{align*}
\text{ident} & := \ (c \colon \text{char})'. \\
\text{prop} & := \text{op} : \text{selector}, \text{op}^? \begin{cases} 
?\text{var} & \Rightarrow \text{ident}, (') \\
?\text{not} & \Rightarrow \text{sub} : \text{prop}, (') \\
?\text{implies}\lor\text{equiv} & \Rightarrow \text{sub}_1 : \text{prop}, \text{sub}_2 : \text{prop}, (') \\
?\text{and}\lor?\text{or} & \Rightarrow (\text{sub} : \text{prop})'
\end{cases}
\end{align*}
\]

Similarly, one can define multimodal logic with inexact identifiers:

\[
\begin{align*}
\text{modal} & := \text{op} : \text{selector}, \text{op}^? \begin{cases} 
?\text{var} & \Rightarrow \text{ident}, (') \\
?\text{not} & \Rightarrow \text{sub} : \text{prop}, (') \\
?\text{implies}\lor\text{equiv} & \Rightarrow \text{sub}_1 : \text{prop}, \text{sub}_2 : \text{prop}, (') \\
?\text{and},?\text{or} & \Rightarrow (\text{sub} : \text{prop})' \\
?\text{box}\lor?\text{dia} & \Rightarrow \text{sub} : \text{prop}, (')
\end{cases}
\end{align*}
\]

In Example 4.2, the inexact field name sub occurs both in prop and in modal. In type prop, it occurs one time as scalar field, and one time as repeated field. If one knows that that $f.\text{op} = ?\text{var}$, one can access $f.\text{sub}$. If one knows that $f.\text{op} \in \{?\text{and},?\text{or}\}$, one can access $f.\text{sub}[i]$. Fieldname sub also occurs three times in type modal, two times as a scalar field, and one time as a repeated field.

**Example 4.3.** The adjectives atom and literal can be defined both on prop and on modal:

\[
\begin{align*}
\text{atom} : \text{prop} & := \text{op}(?\text{var}) \\
\text{atom} : \text{modal} & := \text{op}(?\text{var}) \\
\text{literal} : \text{prop} & := \text{atom} \lor \text{op}(?\text{not}) \land \text{sub}(?\text{atom}) \\
\text{literal} : \text{modal} & := \text{atom} \lor \text{op}(?\text{not}) \land \text{sub}(?\text{atom})
\end{align*}
\]

Similarly, NNF can be defined both on prop and on modal:

\[
\begin{align*}
\text{nnf} : \text{prop} & := \bigvee \left\{ \begin{array}{l}
\text{literal} \\
\text{op}(?\text{and}\lor?\text{or}) \land \text{sub}(?\text{nnf})
\end{array} \right. \\
\text{nnf} : \text{prop} & := \bigvee \left\{ \begin{array}{l}
\text{literal} \\
\text{op}(?\text{and}\lor?\text{or}) \land \text{sub}(?\text{nnf}) \\
\text{op}(?\text{box}\lor?\text{dia}) \land \text{sub}(\text{nnf})
\end{array} \right.
\end{align*}
\]

In the example, there are different occurrences of sub, and it has to be determined which of the possible definitions is being referred to. Before we start discussing the treatment of inexact identifiers, note that defined identifiers in $\Sigma_S$, $\Sigma_C$ and $\Sigma_A$ are always exact. Moreover, we do not allow the use of ambiguous type names. This means that when a defined or built-in type is used, it must be referred to by its exact name.

Resolving of inexact identifiers starts with a preprocessing step on $\Sigma_S$ and $\Sigma_C$. During this step, declared fields are replaced by unique exact identifiers, and any adjectives used in field declarations are moved to $\Sigma_A$. This is done by replacing them with a unique exact identifier, and defining this identifier in $\Sigma_A$. After the preprocessing step, all inexact identifiers are in adjectives in $\Sigma_A$. 

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Adjectives occurring in the types of function declarations will be dealt with in the same way. They are replaced with identifiers, defined in $\Sigma_A$, somewhat similar to the way subformulas are replaced in the CNF transformation ([8]). In our case the goal is not efficiency, but to make overload resolution possible in the first place.

We allow reuse of the same inexact field name in a compound type, as long as the different occurrences of the field name are in different states of the type. For example in prop in Example 4.2, it is possible to reuse fieldname sub between formulas built by ? not and formulas built by ? and or, but renaming both fields sub$_1$ and sub$_2$ into sub at the same time would be impossible. It would be still possible to rename one of them into sub.

We first discuss how type definitions in $\Sigma_S$ and $\Sigma_C$ are processed.

**Definition 4.4.** Let $D = ( w_1 : W_1, \ldots, w_n : W_n )$ be a sequence of declarations. Let $w_{n+1} : W_{n+1}$ be a single declaration. We define

$$D + ( w_{n+1} : W_{n+1} ) = ( w_1 : W_1, \ldots, w_{n+1} : W_{n+1} ).$$

**Definition 4.5.** We define procedure $\text{preprocsimple}(T)$ that preprocesses a simple type $T$. The result is again a simple type. It makes additions to $\Sigma_A$ in the process. The implementation of $\text{preprocsimple}(T)$ is as follows:

- Let $A = \text{ADJ}(T)$, and let $T' = \text{IMPL}(T)$. If $A = \top$, then return $T'$. If $A \neq \top$, then create a new exact identifier $\alpha$, assign $\Sigma_A(\alpha) := (T', A)$, and return $T' \circ \alpha$.

**Definition 4.6.** We define procedure $\text{preproccompound}(C, D)$ that preprocesses a compound type $C$ within declaration context $D$. The second argument is is used for checking that no inexact identifier is reused on a possible realization of $C$.

- If $C$ has form $( v_1 : V_1, \ldots, v_n : V_n )$, then
  - if there exist $(w : W) \in D$ and $1 \leq i \leq n$, s.t. $w^i = v_i$, then create an error.
  - if there exist $1 \leq i < j \leq n$, s.t. $v_i = v_j$, then create an error.
  - Otherwise, create new, exact identifiers $e_i$ for each $v_i$, and set $e_i^0 = v_i$.
  - Set $V'_i = \text{preprocsimple}(V_i)$.
  - Return $(e_1 : V'_1, \ldots, e_n : V'_n)$.

- If $C$ has form $(v : V, C')$ then if there is a $(w : W) \in D$, s.t. $w^0 = v$, create an error. Otherwise, let $e$ be a new, exact identifier for $v$. Set $e^0 = v$. Set $V' = \text{preprocsimple}(V)$. Return $e : V'$, $\text{preproccompound}(C', D + (e : V'))$.

- If $C$ has form $s?(A_1 \Rightarrow C_1, \ldots, A_m \Rightarrow C_m)$, then there must a be unique $(w : W) \in D$, s.t. $w^0 = s$.
  - If no such $w$ exists, or $w$ is not unique, then create an error.
  - Otherwise, create new exact identifiers $\alpha_1, \ldots, \alpha_m$, and set $\Sigma_A(\alpha_i) = (W, A_i)$, for each $1 \leq j \leq m$.
  - After that, return $w?(\alpha_1 \Rightarrow \text{preproccompound}(C_1, D), \ldots, \alpha_m \Rightarrow \text{preproccompound}(C_m, D))$.

We now apply procedures $\text{preprocsimple}$ and $\text{preproccompound}$ as follows:

- For every identifier $v$ in the domain of $\Sigma_S$, replace $\Sigma_S(v)$ by $\text{preprocsimple}(\Sigma_S(v))$.
- For every identifier $v$ in the domain of $\Sigma_C$, replace $\Sigma_C(v)$ by $\text{preproccompound}(\Sigma_C(v), (\ ))$.

**Example 4.7.** After being replaced by $\text{preproccompound}$, the definition of prop will have the following form in $\Sigma_C$:

$$\text{prop} := \text{op}: \text{selector},$$

$$\begin{align*}
\alpha_1 & \Rightarrow \text{ident}, (') \\
\alpha_2 & \Rightarrow \text{sub}: \text{prop}, (') \\
\alpha_3 & \Rightarrow \text{sub}_1 : \text{prop}, \text{sub}_2 : \text{prop}, (') \\
\alpha_4 & \Rightarrow (\text{sub}_n : \text{prop})'
\end{align*}$$
\[
\begin{align*}
\alpha_1 & : \text{selector} := (\text{selector}, ?\text{var}) \\
\alpha_2 & : \text{selector} := (\text{selector}, ?\text{not}) \\
\alpha_3 & : \text{selector} := (\text{selector}, ?\text{implies}?\text{equiv}) \\
\alpha_4 & : \text{selector} := (\text{selector}, ?\text{and}?\text{or}) \\
\end{align*}
\]

At this moment, we have moved all inexactness into \( \Sigma_A \). Adjective definitions in \( \Sigma_A \) have form \( \Sigma_A(\nu) = (T, A) \), where \( \nu \) is an exact identifier, \( T \) is the type on which the defined adjective can be applied, and \( A \) is the adjective. Unfortunately, \( T \) is a simple type, which also may contain inexact adjectives. As an example where this could occur, one could define adjective \( \text{cnf} \) (Example 1.10) on \( \text{prop}\cdot\text{nnf} \) instead of \( \text{prop} \). The definition would have form \( \Sigma_A(\text{cnf}) = (\text{prop}\cdot\text{nnf}, A') \), where \( A' \) is an inexact version of the expression given in Example 1.10, and \( \text{nnf} \) has a definition of form \( \Sigma_A(\text{nnf}) = (\text{prop}, A'') \), where \( A'' \) is the expression of Example 4.3.

In order to solve this, we apply a preprocessing stage on \( \Sigma_A \), similar to the preprocessing of \( \Sigma_S \). We replace all adjectives occurring in domains by exact identifiers, while adding the definitions to \( \Sigma_A \).

- For every identifier \( \nu \) in the domain of \( \Sigma_A \), write \( \Sigma_A(\nu) \) as \( (T, A) \). After that, replace \( \Sigma_A(\nu) \) by \((\text{preprocsimple}(T), A)\).

Note that the call of \text{preprocsimple} may implicitly add new identifiers to \( \Sigma_A \). These new identifiers need not be considered because their domain is an implementation type without adjectives. Hence there is no risk of non-termination. Unfortunately, there potentially exists a circular dependency between preconditions of adjectives that is hard to detect, as illustrated by the following example:

**Example 4.8.** Suppose that \( \Sigma_A \) contains adjective definitions with the following circular structure:

\[
\begin{align*}
v_1 & := (T_1\vee\alpha_1, \cdots) \\
v_2 & := (T_2\vee\alpha_2, \cdots) \\
& \vdots \\
\alpha_1 & := (T_1, \cdots v_2^?) \\
\alpha_2 & := (T_2, \cdots v_1^?)
\end{align*}
\]

If at some other point an occurrence of \( v_1^? \) needs to be resolved, then \( v_1 \) is an overload candidate. In order to decide if \( v_1 \) should be used, one has to resolve the adjective of \( \Sigma_A(\alpha_1) \), which contains \( v_2 \). In order to decide if \( v_2 \) is an overload candidate for this occurrence of \( v_2^? \), one has to resolve the adjective in \( \Sigma(\alpha_2) \), which again contains \( v_1^? \). This circular dependency is not solvable. Note that using \( v_1^? \) or \( v_2^? \) inside the right hand side of \( v_1 \) or \( v_2 \) is unproblematic.

The circular dependency in Example 4.8 is unsolvable, so the compiler has to reject it. Unfortunately, it is difficult to detect, because it only exists if the use of \( v_2^? \) in \( \Sigma(\alpha_1) \) is applied on implementation type \( T_2 \), and the use of \( v_1^? \) in \( \Sigma(\alpha_2) \) is applied on implementation type \( T_1 \). If for example the use of \( v_2^? \) in \( \Sigma(\alpha_1) \) is not on \( T_2 \), then \( v_2 \) is not a candidate for overload, and there is no problem.

We solve this problem by starting to resolve an adjective definition, and whenever we encounter a precondition that has to be resolved first, we recursively try to solve this precondition. If this iteratively results in returning to the original definition, we reject \( \Sigma_A \). In order to detect when we returned to the original definition, we maintain a set \( E \) of adjective definitions that we have encountered already. A circular dependency occurs when we need to resolve the precondition of an identifier that already occurs in \( E \).

At this stage, all inexact identifiers are confined in \( \Sigma_A \), in the second components \( A \) of the definitions \( \Sigma_A(\nu) = (T, A) \). We give the procedure:

**Definition 4.9.** Let \( \Sigma_A \) be the map of inexact adjective definitions. We define procedure \text{RESOLVE}(\nu) \) that tries to resolve the overloads in \( \Sigma_A(\nu) \).

It uses a set \( E \) of identifiers that were already encountered. Initially, \( E = \emptyset \). The purpose of \( E \) is to detect circular dependencies of the type shown in Example 4.8. The implementation is as follows:
• If \( v \) has no definition in \( \Sigma_A(v) \), then the result is an error.
• If \( v \in E \), then the result is also an error. This means that a circular dependency was detected.
• Otherwise, add \( v \) to \( E \).
• Write \( \Sigma_A(v) \) in the form \((U, A)\). If \( U \) has form \((T\circ w)\), then call RESOLVE\((w)\).
• Let \( A' = \text{RESOLVE}_T(\emptyset, A) \), and replace \( \Sigma_A(v) \) by \( A' \). Note that this function calls \( \text{RESOLVE}_T \), that will be defined in Definition 4.10.

Function \( \text{RESOLVE}(v) \) possibly calls itself in order to make \( w \) exact. This will be detected, because we will have \( v \in E \). There is no way to detect such circularity a priori, because its existence depends on the way overloads are resolved.

The second procedure \( \text{RESOLVE}_T(\Gamma, A) \) tries to resolve the fields and adjectives in the non-exact adjective \( A \) in context \( \Gamma \). Context \( \Gamma \) is needed because it may contain preconditions of fields.

In case more than one overload candidate exists, we will take the nearest fit. This approach is used by C++ and Java. For example, if some class \( C \) is a subclass of \( B \), which is a subclass of \( C \), and some function \( f \) has definitions \( f(A) \) and \( f(C) \), then the call \( f(a) \) will be resolved as \( f(A) \), while the call \( f(b) \) will be resolved as \( f(C) \). We have no notion of subclass, but we have implication between adjectives. Using implication between adjectives, the rule becomes as follows: If an application of some inexact identifier \( f \) has different overload candidates \( f_1, \ldots, f_n \), with \( f_1^2 = \ldots = f_n^2 = f \), s.t. each \( f_i \) is defined on adjective \( A_i \), then we resolve \( f \) as follows: If there is a unique \( A_i \), s.t. \( A_i \) implies all of \( A_1, \ldots, A_n \), then \( f \) will be resolved as \( f_i \). In the definition, we will write \( \Gamma \vdash A \), which means that \( \Gamma, \neg A \) is unsatisfiable. It should be noted that this is always in the context of a fixed \( \Sigma_S, \Sigma_C, \) and \( \Sigma_A \). We do not want to use the notation \( \Sigma_S, \Sigma_C, \Sigma_A, \Gamma \vdash A \), because it becomes too long.

**Definition 4.10.** Let \( T \) be an implementation type, let \( \Gamma \) be a set of exact adjectives applicable on \( T \), and let \( A \) be an inexact adjective. \( \text{RESOLVE}_T(\Gamma, A) \) tries to resolve the overloads in \( A \) when applied on \( T \) in context \( \Gamma \). If it succeeds it returns the exact version of \( A \). We define \( \text{RESOLVE}_T(\Gamma, A) \) by cases on the form of \( A \).

- If \( A \) is an (inexact) identifier \( v \), then let \( v_1, \ldots, v_n \) be the adjectives defined in \( \Sigma_A \), that have \( v_i^2 = v \) and for which \( \Sigma_A(v_i) \) has form \((T, A_i)\) or \((T\circ w_i, A_i)\).
  
  Set \( C = \emptyset \). For each \( i \in \{1, \ldots, n\} \), do the following:
  
  - If \( w_i \) is absent, add \( i \) to \( C \).
  - Otherwise, call \( \text{RESOLVE}(w_i) \). (This is the first version, defined in Definition 4.9.) If after that, \( \Gamma \vdash w_i \), then add \( i \) to \( C \).

  At this moment \( C \) is a subset of \( \{1, \ldots, n\} \) containing the candidates that can still be considered as overloads for \( v \). For \( i \in C \) do:
  
  - for \( j \in C \setminus \{i\} \) do:
    
    - \( \text{If } T, w_i \vdash w_j, \text{ then remove } j \text{ from } C. \)

  If \( |C| > 1 \), then create an error message. Otherwise, return \( v_i \) where \( i \) is the unique element of \( C \).
- If \( A \) is a constant \( c \), then if \( c \) has primitive type \( T \), return \( c \). Otherwise create an error.
- If \( A \) has form \( c^0 \), and \( c \) has a primitive type that is not selector, then return \( c^0 \). Otherwise, create an error.
- If \( A \) has form \( f(A') \), then let \( f_1, \ldots, f_n \) be the fields (scalar or repeated) defined on type \( T \), that have \( f_i^2 = f \). Note that \( T \) must be a defined compound type, because primitive types have no fields. Set \( C = \{1, \ldots, n\} \). For each \( i \in C \) do:
  
  - let \( \text{PREC}(f_i) = g_{i,1}(A_{i,1}) \land \cdots \land g_{i,k_i}(A_{i,k_i}) \) with \( k_i \geq 0 \), be the precondition of field \( f_i \) as defined in Definition 1.12. Each \( A_{i,j} \) is either a implementation type or has form \( T_{i,j}\circ w_{i,j} \) with \( w_{i,j} \) an exact identifier defined on implementation type \( T_{i,j} \). For every \( w_{i,j} \) \( (1 \leq j \leq k_i) \) that is present, call \( \text{RESOLVE}(w_{i,j}) \), defined in Definition 4.9.

  After that, check that \( \Gamma \vdash g_{i,1}(A_{i,1}) \land \cdots \land g_{i,k_i}(A_{i,k_i}) \). If not, then remove \( i \) from \( C \).
If \(|C| \neq 1\), then the result is error. Otherwise, let \(i\) be the unique element of \(C\). Assume that the declaration \(f_i\) has form \(f_i : W\). If \(f_i\) is a scalar field, then return
\[
f_i(\text{RESOLVE}_{\text{IMPL}}(W)(\text{ADJ}(W), A')).
\]
if \(f_i\) is a repeated field, then return
\[
f_i(\text{RESOLVE}_{\text{IMPL}}^+(W)(\text{ADJ}(W), A')).
\]
In the latter case, we called \(\text{RESOLVE}^+\) defined below.
- If \(A\) has form \(A_1 \lor \cdots \lor A_n\), then
\[
\text{RESOLVE}_T(\Gamma, A) = \text{RESOLVE}_T(\Gamma, A_1) \lor \cdots \lor \text{RESOLVE}_T(\Gamma, A_n).
\]
- If \(A\) has form \(A_1 \land \cdots \land A_n\), then set \(\Gamma_1 = \Gamma\). For \(i = 1\) to \(n\) do the following:
  - Let \(A'_i = \text{RESOLVE}_T(\Gamma_i, A_i)\).
  - Set \(\Gamma_{i+1} = \Gamma_i \cup \{A'_i\}\).
After that, return \(A'_1 \land \cdots \land A'_n\).

Next we define \(\text{RESOLVE}^+\) which resolves repeated fields.
- If \(A\) has form \(\forall A'\), then return \(\forall \text{RESOLVE}_T(\emptyset, A')\).
- If \(A\) has form \(\exists A'\), then return \(\exists \text{RESOLVE}_T(\emptyset, A')\).
- If \(A = \text{empty}\), and \(T\) is not a compound type, then the result is an error. Otherwise return \(\text{empty}\).

In order to make \(\text{nff}\) of Example 4.3 exact, one has to start by calling \(\text{RESOLVE}(\text{nff})\). Procedure \(\text{RESOLVE}\) will insert \(\text{nff}\) into \(E\), and call \(\text{RESOLVE}_{\text{prop}}(\emptyset, A)\), with \(A\) the expanded definition of \(\text{nff}\). Since \(A\) is a disjunction, \(\text{RESOLVE}_{\text{prop}}(\emptyset, A)\) will process the disjuncts independently.

The first disjunct equals literal. Procedure \(\text{RESOLVE}_{\text{prop}}(\emptyset, \text{literal})\) will establish that literal is the unique overload and return literal. It will not look at the definition of literal. If one wants to resolve the overloads in the definition of literal, one must call \(\text{RESOLVE}(\text{literal})\) separately.

The second disjunct equals \(\text{op}(\text{?and} \lor \text{or}) \land \text{sub}(\text{nff})\). \(\text{RESOLVE}_{\text{prop}}(\emptyset, \text{op}(\text{?and} \lor \text{or}) \land \text{sub}(\text{nff})\) will recursively call \(\text{RESOLVE}_{\text{prop}}(\text{op}(\text{?and} \lor \text{or})\) \), which will resolve \(\text{op}\) (inexact) into \(\text{op}\) (exact) and call \(\text{RESOLVE}_{\text{selector}}(\text{?and} \lor \text{or})\), which will return \(\text{?and} \lor \text{or}\) since both are constants. After that, it will call
\[
\text{RESOLVE}_{\text{prop}}(\text{op}(\text{?and} \lor \text{or}), \text{sub}(\text{nff})\).
\]
There are two possible overloads for \(\text{sub}\), namely \(\text{sub}\) for the \(\text{?not}\) case, and \(\text{sub}_\forall\). We have \(\text{PREC}(\text{sub}) = \text{op}(\text{?not})\), and \(\text{PREC}(\text{sub}_\forall) = \text{op}(\text{?and} \lor \text{or})\). Since only the latter is provable from the premiss, \(\text{sub}_\forall\) will be picked. Since \(\text{sub}_\forall\) is a repeated field, the procedure will recursively call \(\text{RESOLVE}^+_{\text{prop}}(\emptyset, \text{nff})\), which will recursively call \(\text{RESOLVE}_{\text{prop}}(\emptyset, \text{nff})\). This call will return \(\text{nff}\) without expanding it, after which the original call of \(\text{RESOLVE}_{\text{prop}}\) will construct the complete exact overload
\[
\text{RESOLVE}_{\text{prop}}(\text{op}(\text{?and} \lor \text{or}) \lor \text{sub}(\text{nff})\).
\]

\section{Conclusions}
Our goal is to develop and implement an efficient programming language in which it is convenient to implement algorithms on trees whose forms are very different.

In order to obtain this, we have defined a flexible type system together with a way of refining these types by means of adjectives. The adjectives are intended as a replacement for matching in functional languages. In
order to make this replacement possible, we have given a precise semantics for adjectives, so that adjectives can be evaluated on concrete data.

We provided a terminating tableaux calculus for deciding propositional relations between adjectives, and applied this procedure to overload resolution in imprecise formulations of adjectives. The overload resolution procedure replaces ambiguous field references in adjective definitions by exact field references. A similar algorithm can be used for resolving ambiguous overloads in function calls.

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